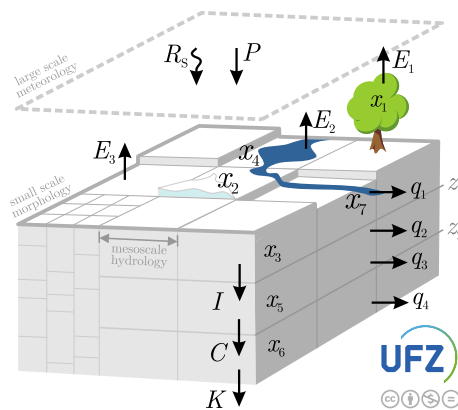


# The mesoscale Hydrologic Model **mHM**



Documentation for version 5.11.2

Presented by the mHM Community

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## The mHM project

[www.ufz.de/mhm](http://www.ufz.de/mhm)

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# Chapter 1

## Introduction to mHM

Welcome to the documentation of mHM version 5.11.2

This chapter is divided in the following sections:

- [Short Description](#)
- [The Grid-based mHM Model](#)
- [Model Formulation](#)
- [The Multiscale Parameter Regionalization Technique](#)
- [The Parameter Estimation Problem](#)
- [Model Calibration](#)
- [Test basin](#)
- [Protocols](#)

### 1.1 Short Description

This document describes the source code for the mesoscale Hydrologic Model mHM. mHM is based on accepted hydrological conceptualizations and is able to reproduce as accurately as possible not only observed discharge hydrographs at any point within a basin but also the distribution of soil moisture among other state variables and fluxes. To achieve these goals and to ensure a reliable performance in ungauged basins, this model employs a multiscale parameter regionalization technique to obtain effective at the scale of interest.

This model is driven by daily or hourly precipitation, temperature fields that are acquired either from satellite products or from observation networks. In the latter case, the driving meteorological data has to be prepared in advance. A module for external drift Kriging is available upon request.

### 1.2 The Grid-based mHM Model

The mHM hydrologic model is based on numerical approximations of dominant hydrological processes that have been tested in various models: HBV [2], [11], [5] and VIC [13]. This model includes also a number of new features that will be described in the next section. In general, this model simulates the following processes: canopy interception, snow accumulation and melting, soil moisture dynamics, infiltration and surface runoff, evapotranspiration,

subsurface storage and discharge generation, deep percolation and baseflow, and discharge attenuation and flood routing (mHM). More information about the model can be found in [16].

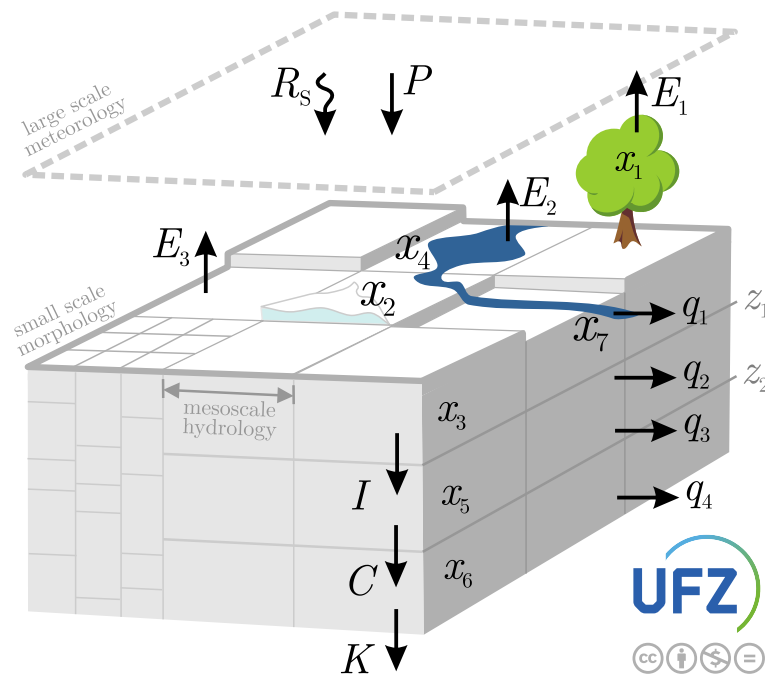


Figure 1.1: Typical mHM cell

Dominant processes of the hydrological cycle at mesoscale span over several orders of magnitude [6]. In this model, three levels (mHM levels) will be differentiated to better represent the spatial variability of state and inputs variables:

- *Level-0*: Spatial discretization suitable to describe the main features of the terrain, the main soil characteristics (pedotop), and the land cover. The cell size at this level is denoted by  $\ell_0$ .
- *Level-1*: Spatial discretization used to describe dominant hydrological processes [4] at the mesoscale as well as the main geological formations of the basin. The cell size at this level is denoted by  $\ell_1$ .
- *Level-2*: Spatial discretization suitable to describe the variability of the meteorological forcings at the mesoscale, for example the formation of convective precipitation. The cell size at this level is denoted by  $\ell_2$ .

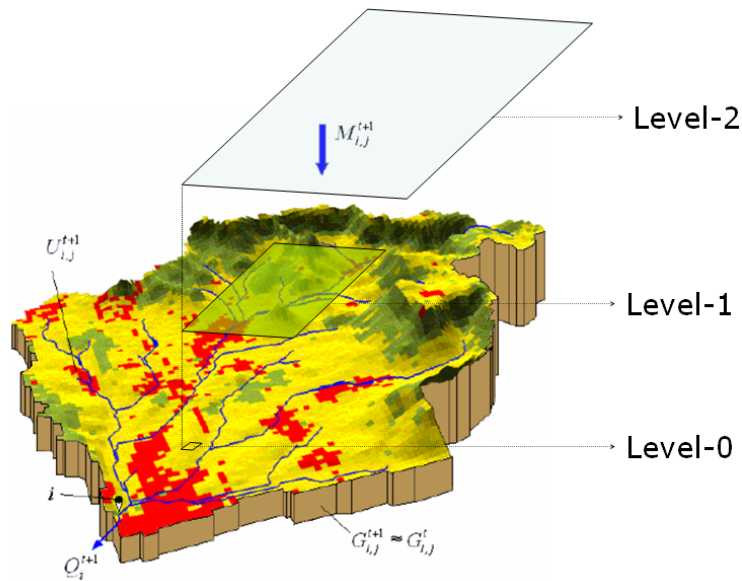


Figure 1.2: Hierarchy of data and modeling levels in mHM

### 1.3 Model Formulation

A mesoscale basin is an open natural system, composed of very heterogeneous materials and having fuzzy boundary conditions. The continuity assumption of the the input variables is quite difficult to justify considering that the spatial heterogeneity of a basin is mostly described by discrete attributes such soil texture types, land cover classes, and geological formations. Due to these reasons, a system of ordinary differential equations ODEs was adopted to describe the evolution of the state variables at a given location  $i$  within the domain  $\Omega$ . This system of ODE is

$$\begin{aligned}
 \dot{x}_{1i} &= P_i(t) - F_i(t) - E_{1i}(t) \\
 \dot{x}_{2i} &= S_i(t) - M_i(t) \\
 \dot{x}_{3i}^l &= (1 - \rho^l) I_i^{l-1}(t) - E_{3i}^l(t) - I_i^l(t) \\
 \dot{x}_{4i} &= \rho^1 (R_i(t) + M_i(t)) - E_{2i}(t) - q_{1i}(t) \\
 \dot{x}_{5i} &= I_i^L(t) - q_{2i}(t) - q_{3i}(t) - C_i(t) \\
 \dot{x}_{6i} &= C_i(t) - q_{4i}(t) \\
 \dot{x}_{7i} &= \hat{Q}_i^0(t) - \hat{Q}_i^1(t)
 \end{aligned}$$

$$\forall i \in \Omega.$$

where

| Inputs | Description  |
|--------|--|
| $P$    | Daily precipitation depth, $\text{mm d}^{-1}$                |
| $E_p$  | Daily potential evapotranspiration (PET), $\text{mm d}^{-1}$ |
| $T$    | Daily mean air temperature, $^{\circ}\text{C}$               |

| Fluxes | Description                                  |
|--------|--|
| $S$    | Snow precipitation depth, $\text{mm d}^{-1}$ |
| $R$    | Rain precipitation depth, $\text{mm d}^{-1}$ |
| $M$    | Melting snow depth, $\text{mm d}^{-1}$       |

| Fluxes | Description   |
|--------|---|
| $E_p$  | Potential evapotranspiration, mm d <sup>-1</sup>                                |
| $F$    | Throughfall, mm d <sup>-1</sup>   |
| $E_1$  | Actual evaporation intensity from the canopy, mm d <sup>-1</sup>                |
| $E_2$  | Actual evapotranspiration intensity, mm d <sup>-1</sup>                         |
| $E_3$  | Actual evaporation from free-water bodies, mm d <sup>-1</sup>                   |
| $I$    | Recharge, infiltration intensity or effective precipitation, mm d <sup>-1</sup> |
| $C$    | Percolation, mm d <sup>-1</sup>   |
| $q_1$  | Surface runoff from impervious areas, mm d <sup>-1</sup>                        |
| $q_2$  | Fast interflow, mm d <sup>-1</sup>  |
| $q_3$  | Slow interflow, mm d <sup>-1</sup>  |
| $q_4$  | Baseflow, mm d <sup>-1</sup>  |

| Outputs | Description   |
|---------|---|
| $Q_i^0$ | Simulated discharge entering the river stretch at cell $i$ , m <sup>3</sup> s <sup>-1</sup> |
| $Q_i^1$ | Simulated discharge leaving the river stretch at cell $i$ , m <sup>3</sup> s <sup>-1</sup>  |

| States | Description   |
|--------|---|
| $x_1$  | Depth of the canopy storage, mm   |
| $x_2$  | Depth of the snowpack, mm   |
| $x_3$  | Depth of soil moisture content in the root zone, mm                       |
| $x_4$  | Depth of impounded water in reservoirs, water bodies, or sealed areas, mm |
| $x_5$  | Depth of the water storage in the subsurface reservoir, mm                |
| $x_6$  | Depth of the water storage in the groundwater reservoir, mm               |
| $x_7$  | Depth of the water storage in the channel reservoir, mm                   |

| Indices  | Description   |
|----------|---|
| $l$      | Index denoting a root zone horizon, $l = 1, \dots, L$ (say $L = 3$ ), in the first layer, $0 \leq z \leq z_1$ |
| $t$      | Time index for each $\Delta t$ interval   |
| $\rho^l$ | Overall influx fraction accounting for the impervious cover within a cell                                     |

## 1.4 The Multiscale Parameter Regionalization Technique

mHM requires at most 28 parameters (depending of the configuration) per cell to account for the spatial variability of the dominant hydrological processes at a mesoscale river basin. These *effective parameters* have to be estimated through calibration. Calibrating this model with a significant number of free parameters for every grid cell would lead to over-parameterization in a mesoscale catchment. This, in turn, would tend to increase the predictive uncertainty of the model due to the *equifinality* [3] of feasible solutions. Moreover, the high dimensionality of this optimization problem is also a daunting task for the state-of-the-art optimization algorithms [14]. To overcome this problem a

multiscale parameter regionalization (MPR) was employed in the mHM model [16].

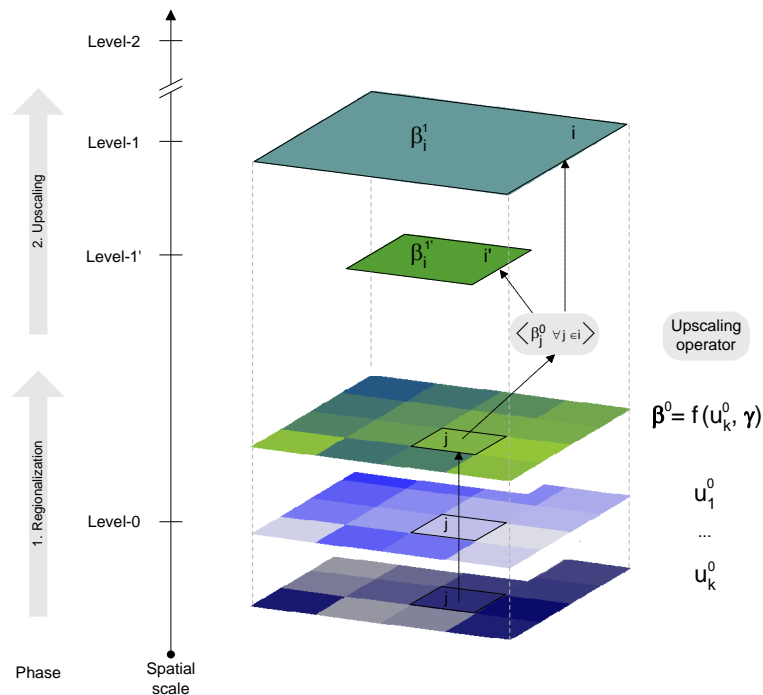


Figure 1.3: MPR

Based on this regionalization method, model parameters at a coarser grid (Level-1) are linked with their corresponding ones at a finer resolution (Level-0) (MPR). The linkage is done with upscaling operators. Model parameters at the finer scale are, in turn, regionalized with nonlinear transfer functions that couple catchment descriptors with the global parameters. Process understanding and empirical evidence are used to define these a priori relationships. The general form of an upscaling operator  $O$  is:

$$\beta_{ki}(t) = O_k \langle \beta_{kj}(t) \quad \forall j \in i \rangle_i$$

where

$$\beta_{kj}(t) = f_k(u_j(t), \gamma).$$

Here  $k = 1, \dots, K$  with  $K$  denoting the number of distributed model parameters.  $u_j$  denotes a  $v$ -dimensional predictor vector for cell  $j$  at level-0 which is contained by cell  $i$  at level-1 (e.g. land cover, elevation, soil texture).

$\gamma$  is a  $s$ -dimensional vector of transfer parameters (super parameters), with  $s$  denoting the number of free parameters to be calibrated or total degrees of freedom.  $O_k \langle \bullet \rangle_i$  denotes the kind of operator applied for the regionalization of the parameter  $k$ . Several types of operators were employed, e.g. majority  $\mathcal{M}$ , arithmetic mean  $\mathcal{A}$ , maximum difference  $\mathcal{D}$ , geometric mean  $\mathcal{G}$ , and harmonic mean  $\mathcal{H}$ . This table also shows the type of relationship employed for the transfer function and the predictors. By establishing such a relationship, the calibration algorithm finds good solutions for the transfer functions parameters ( $s = 45$ ) instead of the model parameters for every grid cell. This, in turn, implies a great reduction of complexity since  $K \times n \gg s$ , where  $n$  denotes the total number of cells of a given basin at level-1.

## 1.5 The Parameter Estimation Problem

Let  $\mathbf{M}\{\mathbf{f}, \mathbf{g}\}$  be a dynamic, spatially distributed, parameter efficient, integrated model that relates a number of state variables  $\mathbf{x}$  with some *observables* called: inputs  $\mathbf{u}$  and outputs  $\mathbf{y}$ . In general, the system is described by the following system of equations

$$\begin{aligned}\dot{\mathbf{x}}(i,t) &= \mathbf{f}(\mathbf{x}, \mathbf{u}, \boldsymbol{\beta}, \boldsymbol{\gamma})(i,t) + \boldsymbol{\eta}(i,t) \\ \mathbf{y}(i,t) &= \mathbf{g}(\mathbf{x}, \mathbf{u}, \boldsymbol{\beta}, \boldsymbol{\gamma})_{\Omega} + \boldsymbol{\varepsilon}(i,t)\end{aligned}$$

where

- $\mathbf{f}$  is a system of functional relationships in mHM (continuous or discrete) that denote the evolution of the system over time.
- $\mathbf{g}$  is a vector of functional relationships used to quantify the expected output (e.g. runoff) of the model denoted by  $\hat{\mathbf{y}}$ .
- $\boldsymbol{\varepsilon}$  denotes the uncertainty of the system originated by defects on measurements of both the inputs and outputs.
- $\boldsymbol{\eta}$  denotes the uncertainty originated by the simplification included during the formulation of  $\mathbf{M}$  or due to the lack of knowledge of the relevant processes (i.e. any kind of model structure deficiency). This term is ignored in the present case.
- $\boldsymbol{\beta}$  denotes the fields of effective mHM parameters at level-1 estimated as described in section (MPR).
- $\boldsymbol{\gamma}$  is a vector of global parameters characterized by a probability density function  $\Phi_{\boldsymbol{\gamma}}$ .
- $i, t$  represent a point in space and time respectively.

In general,  $\boldsymbol{\gamma}$  can be estimated, for example, by

$$\min_{\hat{\boldsymbol{\gamma}}} \|\mathbf{y} - \hat{\mathbf{y}}\|$$

where  $\|\cdot\|$  denotes a robust estimator. Many procedures to estimate global parameters are provided in mHM (e.g. simulated annealing [1], dynamically dimensioned search [17]). Other techniques can be found in the CHS Fortran Library.

## 1.6 Model Calibration

Good parameter sets for  $\boldsymbol{\gamma}$  were identified with a split-sampling technique using an adaptive constrained optimization algorithm based on simulated annealing (SA) [1]. The overall model efficiency was estimated as a weighted combination of four estimators based on the Nash-Sutcliffe efficiency (NSE) between observed and calculated streamflows using three different time scales (daily, monthly and annual) as well as the logarithms of the streamflow to downplay the effects of the peak flows over the low flows [10]. These objective functions are denoted by  $\phi_k$ ,  $k = 1, 4$ . Every objective function should be normalized in the interval  $[0, 1]$ , with 1 representing the best possible solution. The overall objective function to be minimized is then

$$\Phi = \left( \sum_i w_i^p (1 - \phi_i)^p \right)^{\frac{1}{p}}$$

where  $p > 1$ , and  $\sum_{i=1}^4 w_i = 1$ . Here  $p$  is an exponent according to the compromise programming technique [9] and  $w_i$  denote the degree of importance of each objective. High values of  $p$ , say  $p = 6$ , should be chosen to avoid substitution of objective function values at low levels. In general, the estimators related to daily streamflows were twice as important as the long-term ones, thus  $\{w_i\} = \{\frac{2}{4}, \frac{1}{4}, \frac{1}{4}, \frac{2}{4}\}$ . The NSE for a given time interval  $t'$  is given by

$$\phi_k = 1 - \frac{\sum_{t'} (y_k(t') - \hat{y}_k(t'))^2}{\sum_{t'} (y_k(t') - \bar{y}_k(t'))^2}$$



where  $\bar{y}_k(t')$  is the mean value of the observations time series over the calibration period. The index  $k$  denotes here the daily, monthly, yearly, and the transformed  $\ln y(t)$  streamflow discharges.  $y$  and  $\hat{y}$  denote the observed and simulated streamflows at a given time scale. Further details about mHM's calibration options can be found in the section [Calibration Options](#).

## 1.7 Helpful links

Coding and Documentation Style (see [Coding and Documentation Style](#))

Setup netCDF on your MacOS system (see [Install NETCDF](#))

Data Preparation for mHM (see [Data Preparation for mHM](#))

## 1.8 Test basin

mHM comes with a test basin. For details see [The details about the test basin](#).

## 1.9 Protocols

To set up a new input data for mHM see [Protocols for setting up a new mHM basin](#).



## Chapter 2

# Getting Started

This chapter will introduce the structure of mHM technically.

### 2.1 Receive mHM

As from June 2018, the current release of mHM is available through a code repository provided by the version control system GitLab ( <https://git.ufz.de/mhm/mhm>). If you like to contribute to the code development, external accounts be granted upon request through email to [mhm-admin@ufz.de](mailto:mhm-admin@ufz.de).

### 2.2 Install NETCDF

The mHM input and output file format is NETCDF, so the according library is required on your system. Please go to [www.unidata.ucar.edu/software/netcdf](http://www.unidata.ucar.edu/software/netcdf) in order to download the current version.

- **on Linux:** When using ubuntu and gcc, you can use apt to install NETCDF:  

```
sudo apt-get install libnetcdf-dev libnetcdf-fortran-dev
```
- **on MacOS:** You can use [Homebrew](#) to install NETCDF but you need to specify the path in cmake later on:  

```
brew install netcdf
```

### 2.3 Run mHM under CYGWIN on Windows

As from June 2019, we encourage users to make use of cmake guide:

<https://git.ufz.de/mhm/mhm/blob/develop/doc/INSTALL.md>

Further CYGWIN related details might be obsolete.

A NETCDF installation under Windows7 and Windows10 has only been tested using CYGWIN ( <https://www.cygwin.com/>). When executing the CYGWIN setup, the following packages have to be installed (as of May 2018, screenshots kindly provided by Mehmet Cuneyd Demirel):

First, the following netcdf packages have to be installed ([fig\\_cygwin\\_netcdf](#)):

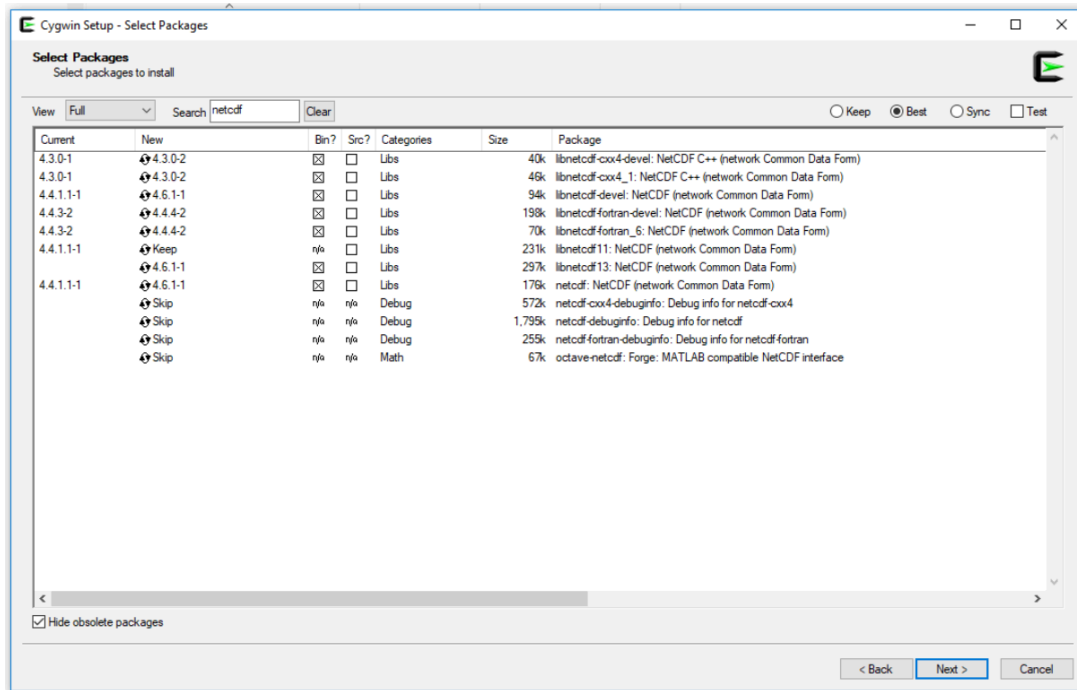


Figure 2.1: netcdf packages for CYGWIN

Second, the following hdf5 version has to be installed for netcdf-4 support.

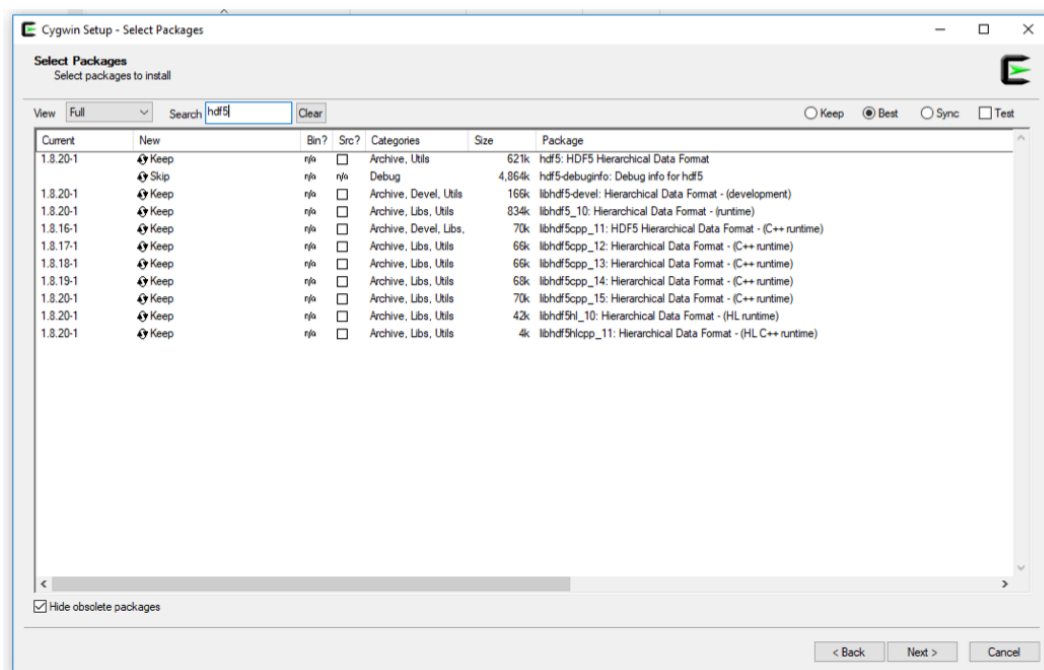


Figure 2.2: hdf5 packages for CYGWIN

Third, the gfortran compiler ([fig\\_cygwin\\_gfortran](#)) and ([fig\\_cygwin\\_libgfortran](#)) have to be installed:

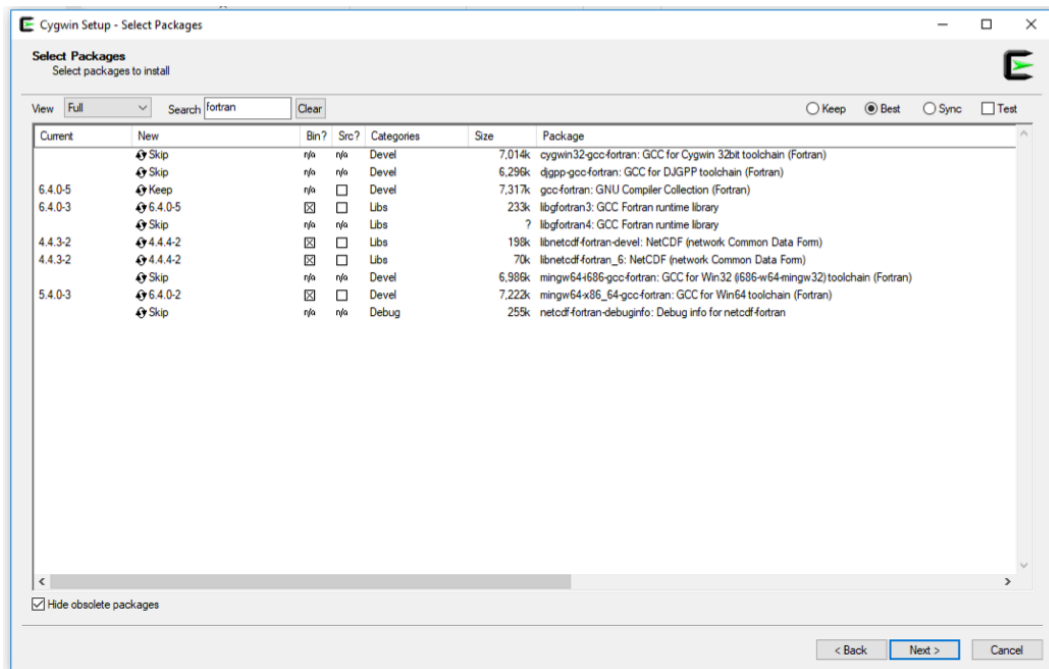


Figure 2.3: gfortran packages for CYGWIN

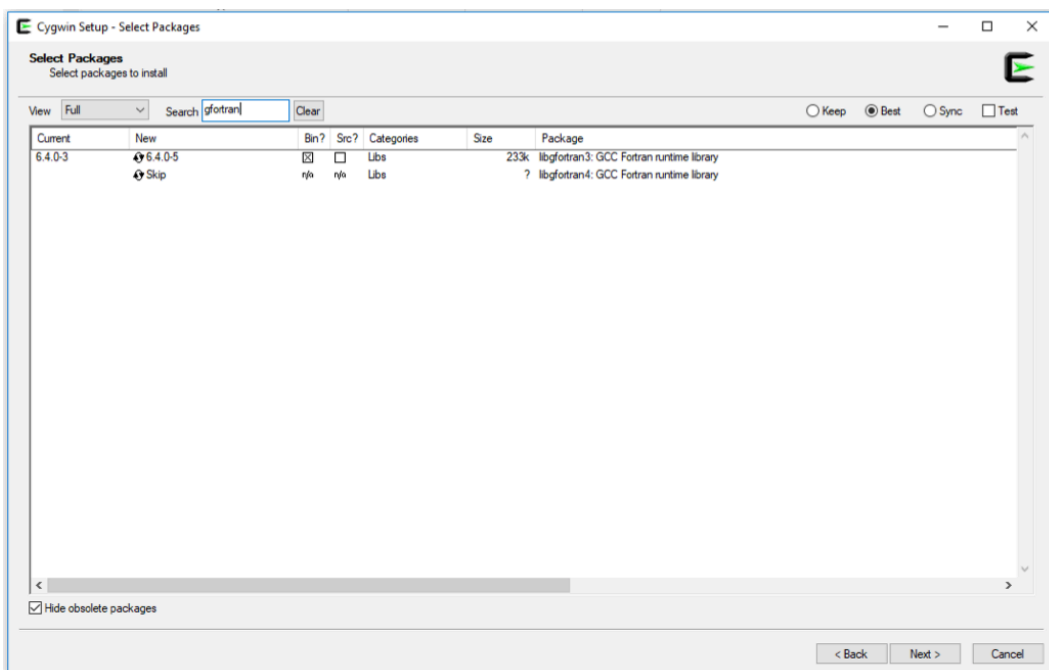


Figure 2.4: libgfortran packages for CYGWIN

Fourth, GNU make has to be made available by selecting the following:

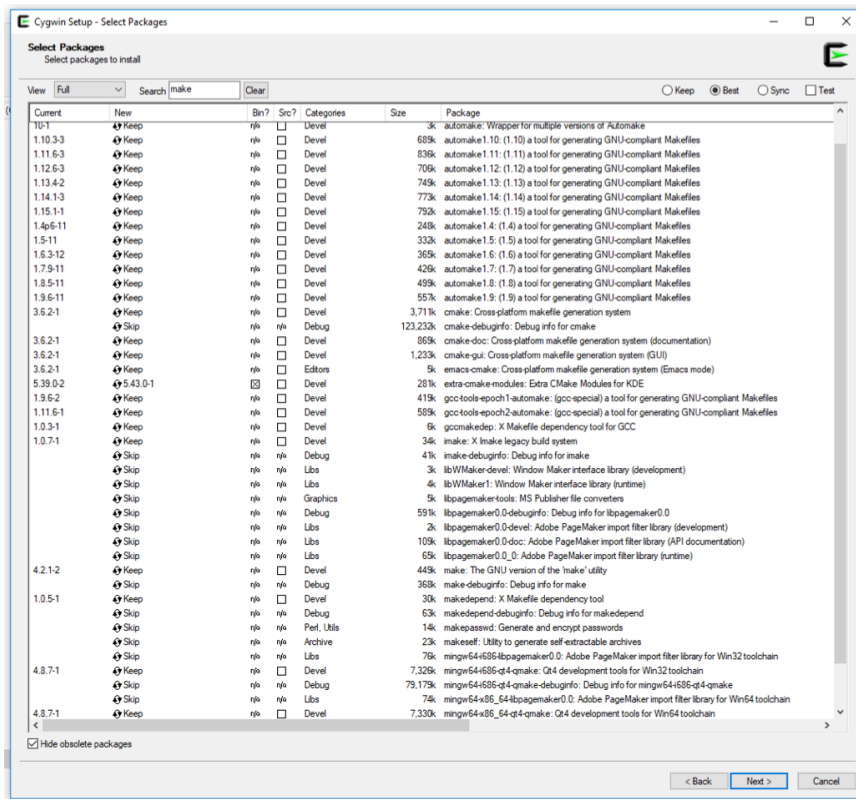


Figure 2.5: make packages for CYGWIN

Note that Cygwin users are advised to define following flag in the Makefile, otherwise compilation won't be successful.

```
EXTRA_LDFLAGS += -Wl,--stack,12485760
EXTRA_CFLAGS += -Wl,--stack,12485760
```

Additionally, make sure to have `\cygwin64\bin` folder in your environmental path ([fig\\_cygwin\\_path](#)).

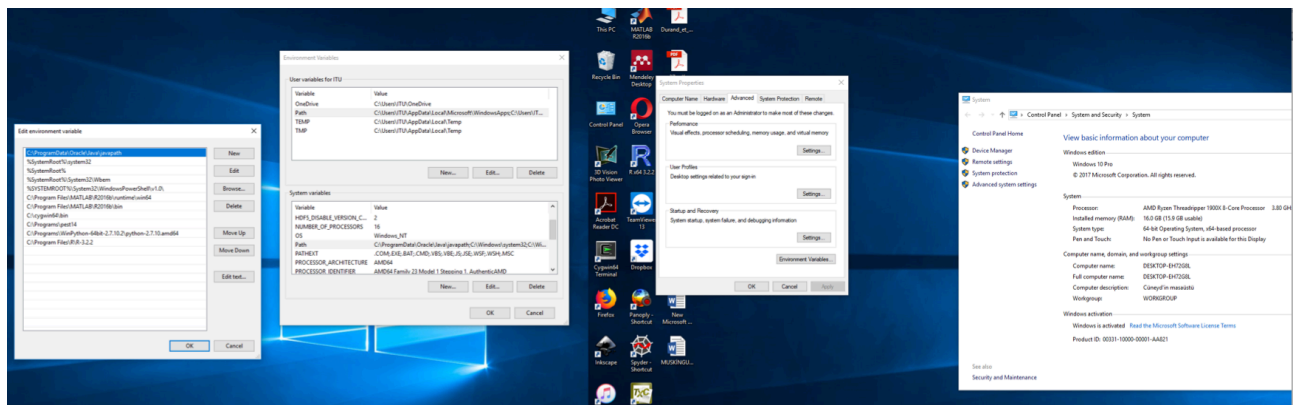


Figure 2.6: setup environmental path CYGWIN

Finally, once the installation of CYGWIN including the packages outlined above is completed, you have to change the system in the Makefile to 'cygwin'. Then, the mhm has proven to compile (simply type `make\bin` on the command line) under Windows. This will produce an executable `mhm\bin` which can be launched by `./mhm\bin` and results for the test basins can be found under `./test_basin/output_b1/` and `./test_basin_2/output_b1/`.

## 2.4 Compile mHM

As from June 2019, we encourage users to make use of cmake guide:

<https://git.ufz.de/mhm/mhm/blob/develop/doc/INSTALL.md>

## 2.5 Test mHM on Example Basin

The mHM distribution usually comes with an example test basin located in `test_basin/`

The directory provides input data for the test basin and output examples. Detailed information about this test basin can be found in the chapter [The details about the test basin](#). All parameters and paths are already set by default to the test case, so you can just start the simulation with the command

```
./mhm
```

This will run mHM on two basins simultaneously and create output files for discharge and interception in `test/output_b*`. The chapter [Visualizing Model Output](#) provides further information on visualising mHM results.

## 2.6 Run your own Simulation

Pretty much the first step mHM takes during runtime is reading the three configuration files:

- `mhm.nml`
- `mhm_output.nml`
- `mhm_parameters.nml`

When editing these files, we recommend to use syntax highlighting for Fortran, which is featured in emacs, for example. This will also guarantee that the file ends with a new line character that is required by the fortran language standard. If you are writing these files with external programs, make sure that the written files comply with Fortran language standard, in particular make sure that the file ends with a new line.

### 2.6.1 Main Configuration: Paths, Periods, Switches

The file `mhm.nml` contains the main configuration for running mHM in your catchments. Since the comments should explain each single setting, this section will only roughly describe the structure of the file. By the way, paths can be relative, absolute and even symbolic links.

- **Common Settings:** Defines output path, input look-up tables and input data format (all "nc" or all "bin") for all basins in your simulation.
- **Basin wise paths:** Set paths for input and output. Create a block for each basin. Remove needless blocks.
- **Resolution:** Hourly or Daily time step. Hydrologic resolution should be factorisable with the input resolutions (e.g. meteo: 4000, hydro: 2000, morph: 100). The routing resolutions determines the velocity of water from cell to cell (keep it greater than the hydro resolution). If you change the routing, remember to recalibrate the model (see [Calibration and Optimization](#)).
- **Restart:** mHM does provide you the option to save the whole model configuration (incl. states, fluxes, routing network, and model parameters) at the end of the simulation period. mHM is then able to restart a new simulation with this model configuration. This reduces the amount of computational time in the newly started simulation because mHM does not have to re-setup the model configuration (e.g., parameter fields and routing network).
- **Periods:** Your actual period of interest should be subsequent of a warming period, where model dynamics can set in properly. Remember to expand your input data by that period, too.

- **Soil Layers:** Soil parameters (not properties) are upscaled vertically in mHM. In this section you specify the number of horizons and depths for the hydrological processing. (This for example you do not find in any other model)
- **Land Cover:** You can provide different land cover files for different years.
- **LAI data:** Switch for choosing LAI option. The user can either select to run mHM with monthly look-up-table LAI values defined for 10 land classes or choose to run mHM using gridded LAI input data (e.g., MODIS). Gridded LAI data must be provided on a daily time-step at the Level-0 resolution. /\*!
- **Process Switches:**
  - Process case 5 - potential evapotranspiration (PET):
    1. PET is input (processCase(5)=0)
    2. PET after Hargreaves-Samani (processCase(5)=1)
    3. PET after Priestley-Taylor (processCase(5)=2)
    4. PET after Penman-Monteith (processCase(5)=3)
  - Process case 8 - routing can be activated (=1) or deactivated (=0).
- **Annual Cycles:** Values for pan evaporation hold in impervious regions only. The meteorological forcing table disaggregates daily input data to hourly values.

## 2.6.2 Output Configuration: Time Steps, States, Fluxes

The file `mhm_output.nml` regulates how often (e.g. `timeStep_model_outputs = 24`) and which variables (fluxes and states) should be written to the final netcdf file `[OUTPUT_DIRECTORY]/mHM_Fluxes_↔ States.nc`. We recommend to only switch on variables that are of actual interest, because the file size will greatly increase with the number of containing variables. During calibration (see [Calibration and Optimization](#)) no output file will be written.

## 2.6.3 Regionalised Parameters: Initial Values and Ranges

The file `mhm_parameters.nml` contains all global parameters and their initial values. They have been determined by calibration in German basins and seem to be transferable to other catchments. If you come up with a very different catchment or routing resolution, these parameters should be recalibrated (see section [Calibration and Optimization](#)).

## 2.7 Calibration and Optimization

By default, mHM runs without caring about observed discharge data. It will use default values of the global regionalised parameters `\gamma`, defined in `mhm_parameters.nml`. In order to fit discharge to observed data, mHM has to be recalibrated. This will optimise the `\gamma` parameters such that mHM arrives at a best fit for discharge. The optimization procedure runs mHM many many times, sampling parameters in their given ranges for each iteration, until the objective function converges to a confident best fit.

### 2.7.1 The Optimization Routines

mHM comes with four available optimization methods:

- **MCMC:** The Monte Carlo Markov Chain sampling of parameter sets is recommended for estimation of parameter uncertainties. Intermediate results are written to `mcmc_tmp_parasets.nc`.
- **DDS:** The Dynamically Dimensioned Search is an optimization routine known to improve the objective within a small number of iterations. However, the result of DDS is not necessarily close to your global optimum. Intermediate results are written to `dds_results.out`.



- **Simulated Annealing:** Simulated Annealing is a global optimization algorithm. SA is known to require a large number of iterations before convergence (e.g. 100 times more than DDS), but finds parameter sets closer to the global minimum like DDS. Intermediate results are written to `anneal_results.out`.
- **SCE:** The Shuffled Complex Evolution is a global optimization algorithm which is based on the shuffling of parameter complexes. It needs more iterations compared to the DDS (e.g. 20 times more), but less compared to Simulated Annealing. The increasing computational effort (i.e. iterations) leads to more reliable estimation of the global optimum compared to DDS. Intermediate results are written to `sce_results.out`.

Objective functions currently implemented in mHM are:

- **NSE:** Nash-Sutcliffe Efficiency, assuming constant + linearly relative errors. Recommended for fitting high flows.
- **InNSE:** logarithmic Nash-Sutcliffe Efficiency, recommended for fitting low flows.
- **0.5\*(NSE+InNSE):** weights both NSE and InNSE by 50%, roughly fits high and low flows.
- **Likelihood:** The confidence bands are probability density functions that capture variable errors, recommended for hydrological discharge.
- **KGE:** Kling Gupta model efficiency: combined measure for variability, bias and correlation
- **PD:** Pattern dissimilarity (PD) of spatially distributed soil moisture
- **ETC:** Combination of other multi-objective functions (streamflow, TWS anomaly, evapotranspiration, soil moisture), see `mhm.nml` for details

## 2.7.2 Calibration Settings for mHM

The following settings in `mhm.nml` are required for calibrating mHM:

- `optimize = .true.`
- `opti_method = 1`  
Choose methods: 0 (MCMC), 1 (DDS), 2 (SA), 3 (SCE)
- `opti_function = 1`  
Choose objective functions: 1 (NSE), 2 (InNSE), 3 (50% NSE+InNSE), 4 (Likelihood)
- `nIterations = 40000`  
Maximum number of iterations (mHM runs), optimisers will exit earlier if convergence criteria are reached.
- `seed = -9`  
The default value -9 will take a seed number from the system clock. Be warned that simulations might not be random if you define a positive seed here.

More specific settings are offered at the very end of the file `mhm.nml`.

In `mhm_parameters.nml` you find the initial values and ranges from which the optimiser will sample parameters. Most ranges are very sensitive and have been determined by detailed sensitivity analysis, so we do not recommend to change them. With the FLAG column you may choose which parameters are allowed to be optimised. The parameter `gain_loss_GWreservoir_karstic` is not meant to be optimised.

Please mind that optimization runs will take long and may demand a huge amount of hardware resources. We recommend to submit those jobs to a cluster computing system.

### 2.7.3 Final Calibration Results

During calibration, mHM does not write out fluxes, states and discharge, so you need to perform a final run with the calibrated parameters. When the simulations are finished, the optimal parameter set is written to

```
[OUTPUT_DIRECTORY]/FinalParams.out  
[OUTPUT_DIRECTORY]/FinalParams.nml
```

You can run mHM directly with the generated namelist (by renaming it) or incorporate the results in `FinalParams.out` as new initial values into `mhm_parameters.nml`. There are two scripts that help you with that:

- `pre-proc/create_multiple_mhm_parameter_nml.sh`  
Useful for many optimisation runs (many lines in `FinalParams.out`)
- `post-proc/opt2nml-params.pl`  
Useful to analyze where the optimisation arrived. Using two arguments, `pathto/FinalParams.out` and `pathto/mhm_parameters.nml`, it will (1) create a new `mhm_parameters.nml.opt` filled with the new values and (2) directly show the new parameters within their ranges and warn if some have been close to their end of range by 1%.

As soon as the new parameters are set, deactivate the `optimize` switch in `mhm.nml` and **rerun** mHM once in order to obtain the final optimised output.

You should also have a look at the parameter evolution (e.g. `sce_results.out`) or final results. If any of the parameters stick to the very end of their allowed range, the result is not optimal and you will be in serious trouble. Possible reasons might be bad parameter ranges (even though they have been optimised mathematically, but in a basin or resolution not comparable to yours) or bad input data.

## Chapter 3

# Data Preparation for mHM

This chapter is divided in the following sections:

- [Getting Started](#)
- [Preparation of the Forcings](#)
- [Preparation of the Morphological Data](#)
- [A possible GIS workflow](#)
- [Table Data](#)
- [Land Cover Data](#)

### 3.1 Getting Started

To run mHM the user requires a number of datasets. The following subsection gives a short overview and references to freely available datasets.

#### 3.1.1 Meteorological variables

Meteorological data is usually available from the weather services of the countries of modelling interest. Freely available alternatives are the EOBS ( <http://www.ecad.eu/download/ensembles/download.php>) and the WATCH datasets ( [http://www.eu-watch.org/gfx\\_content/documents/README-WFDEI.pdf](http://www.eu-watch.org/gfx_content/documents/README-WFDEI.pdf)).

You may have difficulties finding measurement data for Potential Evapotranspiration (PET). One possible solution, would be to calculate PET from the much easier available variables mean, maximum and minimum air temperature, using the Hargreaves-Samani method.

The two meteorological variables which are needed:

| Name                    | Unit | Temporal resolution |
|-------------------------|------|---------------------|
| Precipitation           | mm   | hourly to daily     |
| Average air temperature | °C   | hourly to daily     |

Dependent of the specification for the potential evapotranspiration (processCase(5)) additional meteorological variables may be need:

- processCase(5) = 0 - PET is read from input.

- processCase(5) = 1 - Hargreaves-Samani equation
- processCase(5) = 2 - Priestley-Taylor equation
- processCase(5) = 3 - Penam-Monteith equation

| Name                              | Unit       | Temporal resolution |
|-----------------------------------|------------|---------------------|
| 0 - Potential evapotranspiration  | mm         | hourly to daily     |
| 1 - Minimum air temperature       | °C         | daily               |
| 1 - Maximum air temperature       | °C         | daily               |
| 2 - Net radiation                 | $W m^{-2}$ | daily               |
| 3 - Net radiation                 | $W m^{-2}$ | daily               |
| 3 - Absolut vapur pressure of air | Pa         | daily               |
| 3 - Windspeed                     | $m s^{-1}$ | daily               |

### 3.1.2 Morphological variables

In addition to the Digital Elevation Models (DEM) usually provided by federal authorities, a number of free alternatives exists. SRTM data is available from 60° South to 60° North in a 3" (~ 90 m) resolution (<http://srtm.csi.cgiar.org/>), the ASTER-GDEM covers the entire globe with a resolution of 1" (<http://gdem.ersdac.jspacesystems.or.jp/>). Hydrographically corrected SRTM data and a number of derived products in different resolutions are available from the HydroSHEDS project (<http://hydrosheds.cr.usgs.gov/index.php>).

Soil and hydrogeological data is usually provided by the geological surveys. On the soil side the Harmonized World Soil Database would be a free alternative (<http://webarchive.iiasa.ac.at/Research/LUC/External-World-soil-database/HTML/>).

| Name   | Unit             | Temporal resolution |
|--|------------------|---------------------|
| Digital elevation model  | m                | -                   |
| Soil maps with textural properties (% sand and clay contents, bulk density per horizon, and root depth zone) | -                | -                   |
| Geological maps with aquifer properties (specific yield, permeability, aquifer thickness)                    | -                | -                   |
| Streamflow location  | (m, m) (lat,lon) | -                   |

### 3.1.3 Land Cover

Free land cover data is available from different sources. The Corine programm provides land cover scenes for Europe with resolutions of 100m and 250m (<http://www.eea.europa.eu/publications/COR0-landcover>), the Global Land Cover Map 2000 a dataset covers the entire world in a resolution of 1km.

| Name              | Unit | Temporal resolution |
|-------------------|------|---------------------|
| Land cover scenes | -    | monthly to annual   |
| Leaf area index   | -    | weekly to monthly   |

### 3.1.4 Gauging Station Information

Streamflow measurements should also be available from federal authorities. In addition, the Global Runoff Data Center provides timeseries of varying length for several thousand gauging stations all over the world (<http://www.bafg.de/GRDC>)

| Name                    | Unit                      | Temporal resolution |
|-------------------------|---------------------------|---------------------|
| Streamflow measurements | $\text{m}^3\text{s}^{-1}$ | hourly to daily     |

### 3.1.5 Optional Data

| Name                             | Unit             | Temporal resolution |
|----------------------------------|------------------|---------------------|
| Snow cover                       | -                | daily to weekly     |
| Land surface temperature         | K                | daily to weekly     |
| Groundwater station location     | (m, m) (lat,lon) | -                   |
| Groundwater head measurements    | m                | weekly to monthly   |
| Eddy covariance station location | (m, m) (lat,lon) | -                   |
| Eddy covariance measurements     | m                | hourly              |

## 3.2 Preparation of the Forcings

Forcing data should be available from federal databases like DWD for Germany. Be sure to bring data in the netcdf format, like

```
precipitation-1950-2013.nc
```

These files can be edited with the CDO module package in order to select or change data. Please read the according CDO and NCKS reference sheets for details and make sure to load the CDO modules before starting.

### 3.2.1 Extract Data to the Size of a Catchment

Let FULLMAP.nc be the the forcing data of a region much greater than your catchment. Let X1, X2, Y1, Y2 be the coordinates (lon and lat) of a rectangular overlapping the catchment. The forcing data MAP.nc for your catchment can be extracted by the CDO command:

```
cdo sellonlatbox,X1,X2,Y1,Y2 FULLMAP.nc MAP.nc
```

Optionally, certain pixels can be extracted, too:

```
ncks -d x,1,1 -d y,2,2 MAP.nc PIXEL.nc
```

### 3.2.2 Extact Data to the Time Period of Interest

Let MAP.nc be the forcing data including your period of interest. Let DATE1 and DATE2 be the starting and ending dates of your period, respectively. Their date format is YYYY-MM-DD. The data PERIOD.nc can be extracted by the CDO command:

```
cdo seldate,DATE1,DATE2 MAP.nc PERIOD.nc
```

Please note that a reference date according to DATE1 should be provided in PERIOD.nc:

```
cdo setreftime,DATE1,00:00:00,days PERIOD.nc FINAL.nc
```

### 3.2.3 Data Types

Any data provided for mHM should be of the data type DOUBLE. You should convert all data related to a variable VAR (e.g. tavg) with the following CDO command:

```
cdo -b F64 selname,VAR FINAL.nc FINAL64.nc
```

The only exception here is the time, its data type has to be INTEGER. This can be changed with NCAP2:

```
ncap2 -s 'time=int(time)' STILLNOTFINAL.nc FINAL.nc
```

### 3.2.4 Variable Names

In mHM the variable names of the forcing are hard-coded. They need to be

- "tavg" for average temperature
- "pre" for precipitation
- "pet" for evapotranspiration

For example, you can change the variable name TEMPERATURE in your NETCDF file with the following CDO command:

```
cdo chname,TEMPERATURE,tavg TEMPDATA.nc TEMPDATA-FINAL.nc
```

### 3.2.5 The Header File

Every meteorological data file should come with an additional file "header.txt" in the same directory. In the following example, we have only one pixel of data (ncols=nrows=1) and a cell size of 4km. The easting and northing coordinates of the lower left corner should be similar to the morphological data of your catchment.

```
ncols      1
nrows      1
xllcorner  639357.3
yllcorner  5723706.2
cellsize   4000
NODATA_value -9999
```

### 3.2.6 NODATA values

In order to be consistent in your data, you should specify the same NODATA value everywhere. For example, specify "-9999" in your header file as NODATA value. Usually, this should be changed also in NC files by the `ncatted` command. The following example overwrites or adds (o) the NODATA attribute "\_FillValue" with value "-9999" of type double (d) to the variable "pre":

```
ncatted -O -a _FillValue,pre,o,d,-9999. INOUT.nc
```

## 3.3 Preparation of the LatLon Grid

As input mHM additionally needs a `latlon.nc` file specifying the geographical location of every grid cell in WGS84 coordinates. This file has to be adjusted for every resolution of the level-1 hydrological simulations.

For creating a `latlon` file mHM comes with the python script `create_latlon.py`, which can be found in `pre-proc/`. Detailed information for the usage of this python script can be found in the online help by typing:

```
python [MHM_DIRECTORY]/pre-proc/create_latlon.py -h
```

`create_latlon.py` needs several specifications via command line switches. First, the coordinate system of the morphological and meteorological data have to be specified according to [www.spatialreference.org](http://www.spatialreference.org) by using the switch: `-c`. Second, you need to specify three header files containing the corresponding information for the different spatial resolutions connected to mHM. The three resolutions are the resolution of 1) the morphological input (switch: `-f`), 2) the hydrological simulation (switch: `-g`), and 3) the routing (switch: `-e`).

These header files can be produced by adopting the header file of the meteorological data. Therefore, copy one of these files that you generated for meteorological data (see [Preparation of the Forcings](#)):

```
cp [INPUT_DIRECTORY]/input/meteo/pre/header.txt [INPUT_DIRECTORY]/input/latlon/
```

Edit the new file `header.txt` such that `cellsize` equals your hydrologic resolution. You have to adapt `ncols` and `nrows` to that resolution such that it covers the whole region. For example, reducing the cell size from 2000 to 100, the number of columns and rows should be increased by a factor of 20.

Third, you need to set the path and filename for the resulting output file by using the switch `-o`.

An example for creating a lat-lon-file looks like

```
python [MHM_DIRECTORY]/pre-proc/create_latlon.py -c epsg:31463 -f header_100m.txt -g header_1000m.txt -e header_2000m.txt -o ../latlon/latlon.nc
```

Keep in mind that you need one `latlon` file for each hydrologic resolution in mHM. Since the filename `latlon.nc` is hard-coded in mHM, we recommend to create different directories for each resolution you need, containing the related header and `latlon` file.

## 3.4 Preparation of the Morphological Data

MHM needs several morphological input datasets. All these have to be provided as raster maps in the ArcGis ascii-format, which stores the above header and the actual data as plain text. Some of the raster files need to be complemented by look-up tables providing additional information. The tables and their structure are described in more detail in subsection [Table Data](#). Take care of the following limitations to mHM input data during data processing:

- All gridded input, i.e. your morphological and your meteorological data, needs to cover the same spatial domain. That means, that the values `xllcorner`, `yllcorner`, `xllcorner+ncols*cellsize` and `yllcorner+nrows*cellsize` have to be identical for all files!
- MHM allows you to provide your meteorological forcing in a different horizontal resolution than the morphological data. The larger cellsize however needs to be a multiple of the smaller.

The required datasets and their corresponding filenames:

| Description                                 | Raster file name  | Table file name             |
|---|-------------------|-----------------------------|
| Sink filled Digital Elevevation Model (DEM) | dem.asc           | -                           |
| Slope map                                   | slope.asc         | -                           |
| Aspect Map                                  | aspect.asc        | -                           |
| Flow Direction map                          | fdir.asc          | -                           |
| Flow Accumulation map                       | facc.asc          | -                           |
| Gauge(s) position map                       | idgauges.asc      | [gauge-id].txt              |
| Soil map                                    | soil_class.asc    | soil_classdefinition.txt    |
| Hydrogeological map                         | geology_class.asc | geology_classdefinition.txt |
| Leaf Area Index (LAI) map                   | LAI_class.asc     | LAI_classdefinition.txt     |
| Land use map                                | your choice       | -                           |

## 3.5 A possible GIS workflow

In the following paragraphs a possible GIS workflow is outlined using the software ArcMAP 10 with the Spatial Analyst Extension and optionally the Arc Hydro Tools ( [http://downloads.esri.com/blogs/hydro/AH2/ArcHydroTools\\_2\\_0.zip](http://downloads.esri.com/blogs/hydro/AH2/ArcHydroTools_2_0.zip))

### 3.5.1 General considerations

- As the spatial discretizations (i.e. resolutions, origin) of your datasets will most likely differ, it is recommended to set the following environmental settings on every processing step outlined in the following paragraphs.

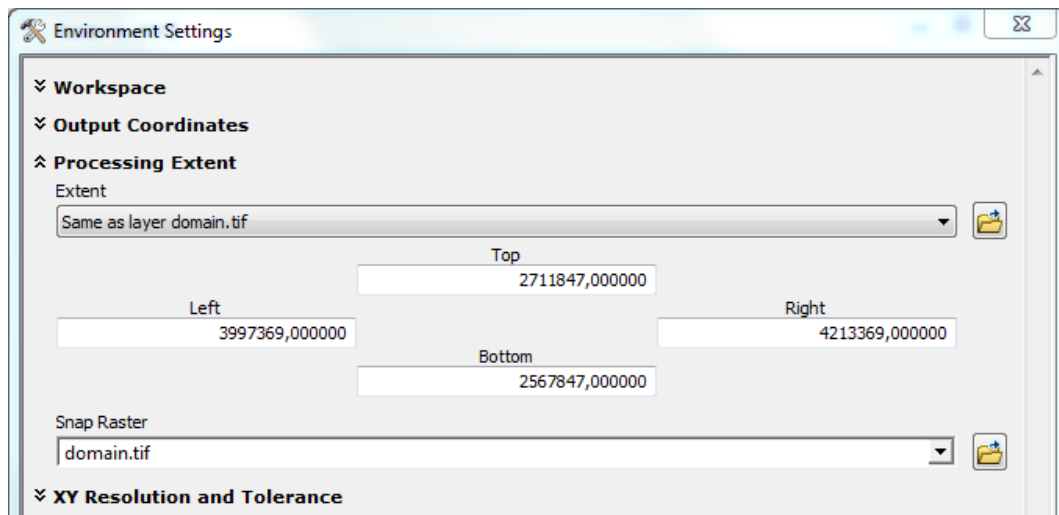


Figure 3.1: Button 'Environments...' in all Toolbox windows

Point to the largest of your input data grids in 'Extent' and 'Snap Raster', which is usually the dataset with the coarsest horizontal resolution. In the likely case, that this is your meteorological input, create a grid from any of your netcdf-files first.

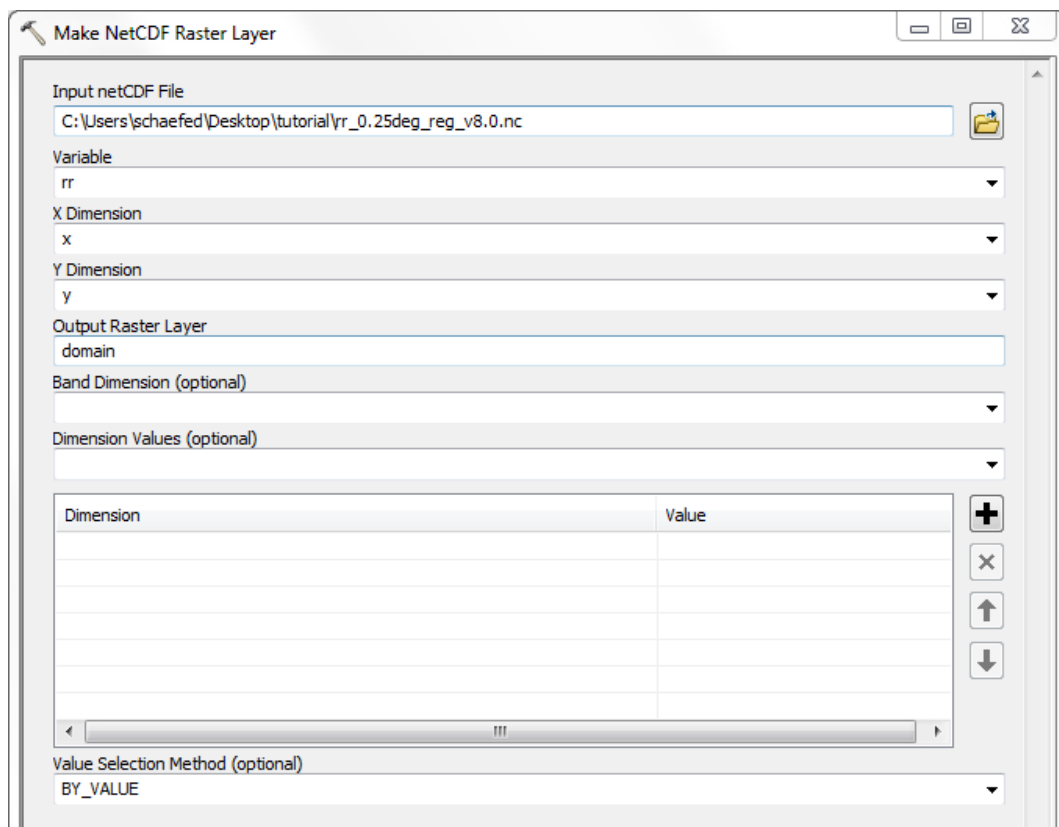


Figure 3.2: System Toolboxes -> Multidimension Tools -> Make NetCDF Raster Layer

Save the output grid (right click the raster in the 'Table of Contents' -> 'Data' -> 'Export Data...').

- If the map projections of your datasets differ, a harmonization of these becomes necessary. Repeat the depicted step to convert all your input files. Which projection to choose is highly dependent on your simulation domain and size. For the current model version an equal-area projection is strongly recommended.



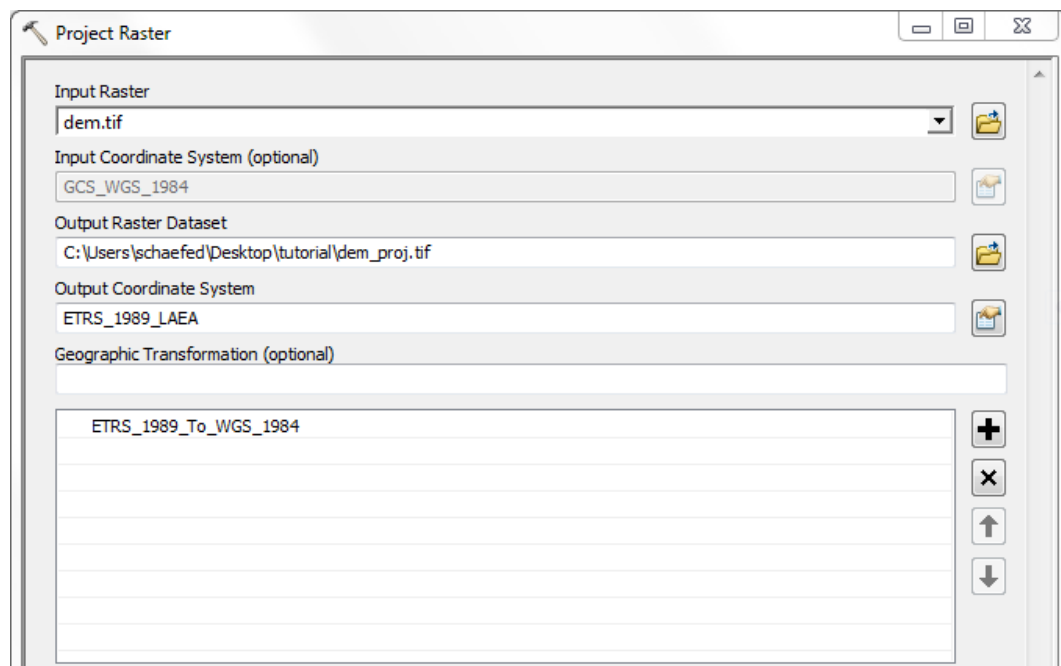


Figure 3.3: System Toolboxes -> Data Management Tools -> Projections and Transformations -> Raster -> Project Raster

- If your input datasets are not already in the desired level-0 resolution, resample the DEM, the hydrogeological, LAI, soil and land use maps. Choosing an appropriate resolution depends on data quality and needed level of simulation detail, but keep in mind that:
  1. Your different input resolution levels must be multiples of each other. E.g. you should choose a level-0 resolution of 100m (instead of 90m in case you are using SRTM data) if your meteorological input resolution is 4km.
  2. Model runtime directly depends on the number of grid cells.

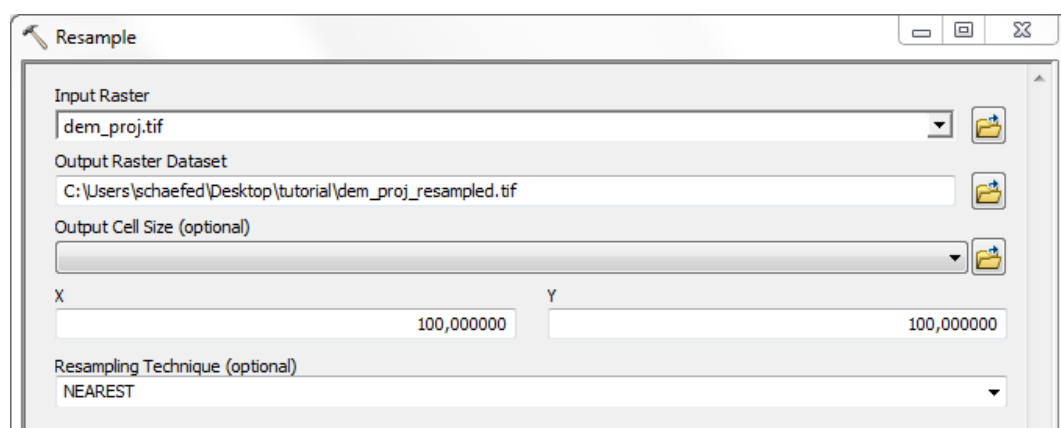


Figure 3.4: System Toolboxes -> Data Management Tools -> Raster -> Raster Processing -> Resample

- It is important that all your morphological input files exactly cover the same spatial domain. That also means that if a cell contains valid data in any one of the datasets, the very same cell must also be defined in all the others. One possibility to solve this typical problem would be to set such 'doubtful' cells to the corresponding NODATA\_value. Therefore create a mask as depicted below, which only contains cells, that are defined everywhere.

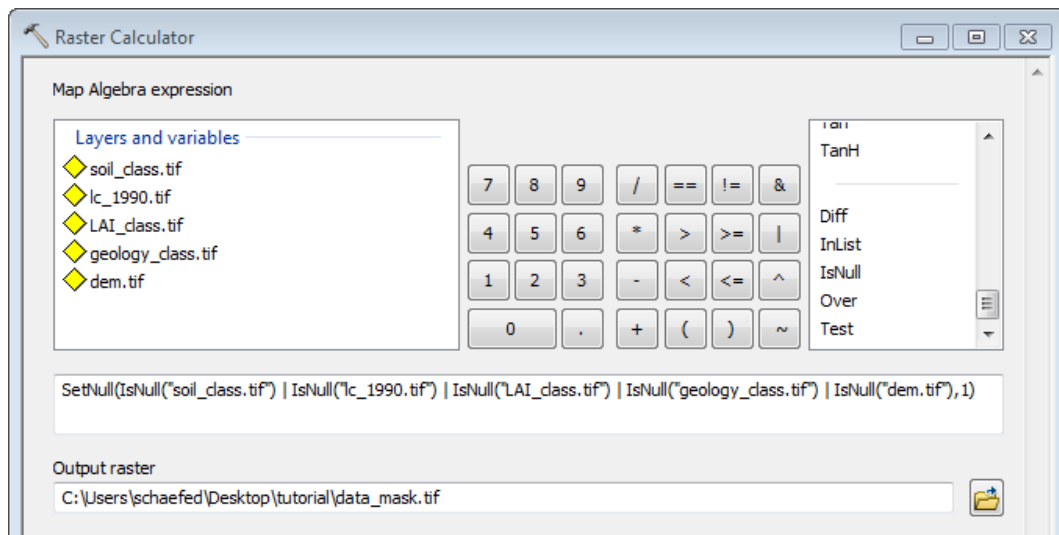


Figure 3.5: System Toolboxes -> Spatial Analyst Tools -> Map Algebra -> Raster Calculator

- Mask all the mentioned datasets with the output of the 'Raster Calculator' following the procedure described in [Mask the datasets](#) of this tutorial. In case the described processing step is necessary, accomplish it **before** you reach subsection [Flow direction and flow accumulation](#) ! Masking these maps would most likely disturb the hydrological properties of your catchment data and result in unexpected model behaviour.

### 3.5.2 Slope map

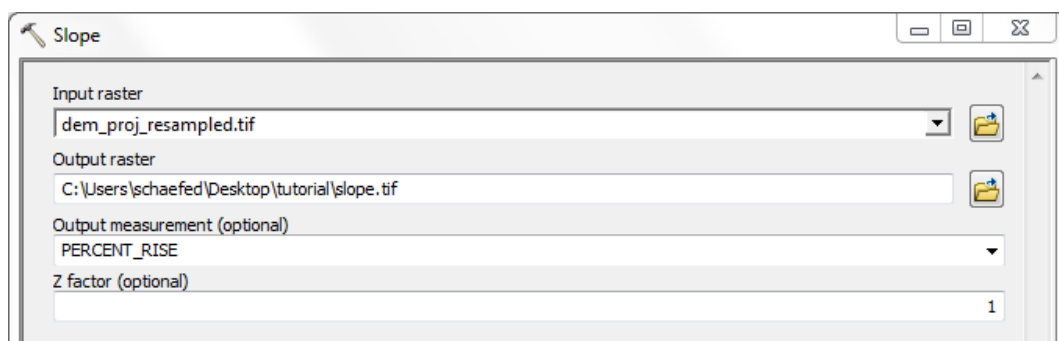


Figure 3.6: System Toolboxes -> Spatial Analyst Tools -> Surface -> Slope

### 3.5.3 Aspect map

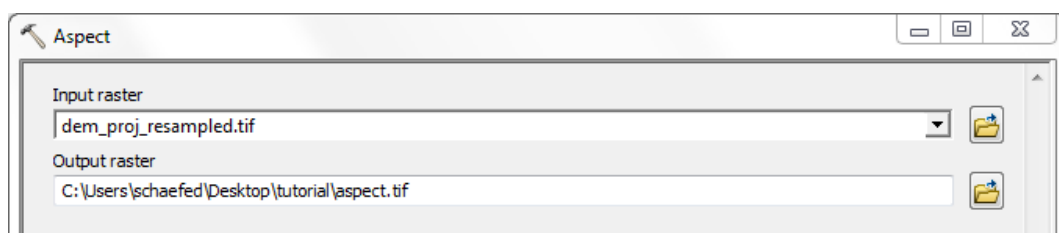


Figure 3.7: System Toolboxes -> Spatial Analyst Tools -> Surface -> Aspect

### 3.5.4 Fill DEM sinks

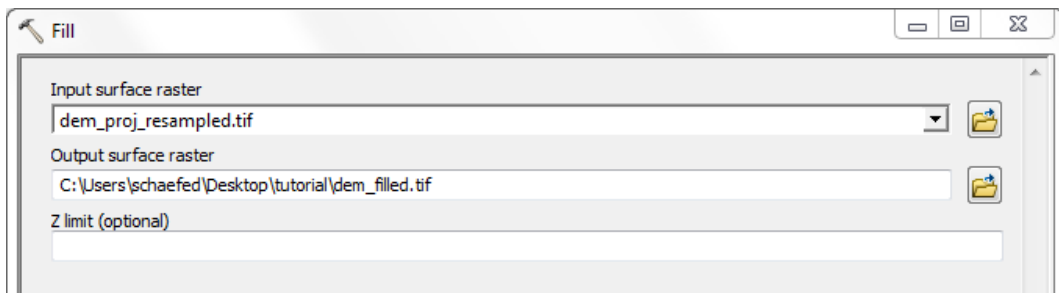


Figure 3.8: System Toolboxes -> Spatial Analyst Tools -> Hydrology -> Fill

### 3.5.5 Flow direction and flow accumulation

Depending on quality and resolution of the DEM map, these steps can be done with the respective tools from the Spatial Analyst Extension or by using the Arc Hydro Tools.

#### 3.5.5.1 Spatial Analyst

If a high quality DEM, with a resolution fine enough to represent small scale river morphology is available, you may calculate flow direction and flow accumulation directly.

- Flow Direction

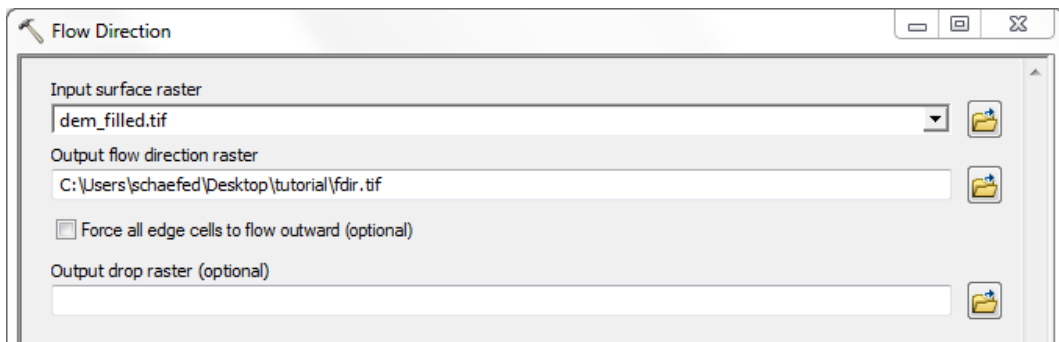


Figure 3.9: System Toolboxes -> Spatial Analyst Tools -> Hydrology -> Flow Direction

- Flow Accumulation

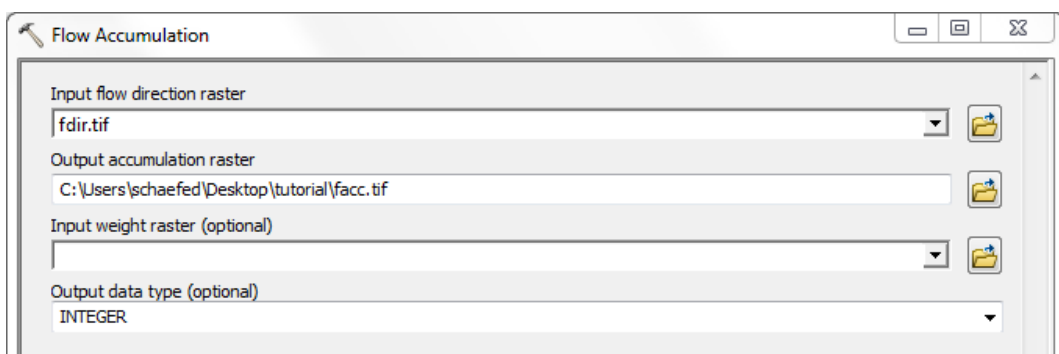


Figure 3.10: System Toolboxes -> Spatial Analyst Tools -> Hydrology -> Flow Accumulation

### 3.5.5.2 Arc Hydro Tools

Using the Arc Hydro Tools is recommended in case of doubtful DEM quality and/or coarse map resolutions.

- In a first step the original DEM must be reconditioned, i.e. the given altitudes will be reassigned in dependence of a stream network. The latter must be given as a Line Shapefile. In case the necessary stream network file is not available, you can get one from the USGS HydroSHEDS download portal <http://hydrosheds.cr.usgs.gov/dataavail.php>.



Figure 3.11: System Toolboxes -> Arc Hydro Tools -> Terrain Preprocessing -> DEM Reconditioning

- Fill the sinks in the resulting reconditioned DEM

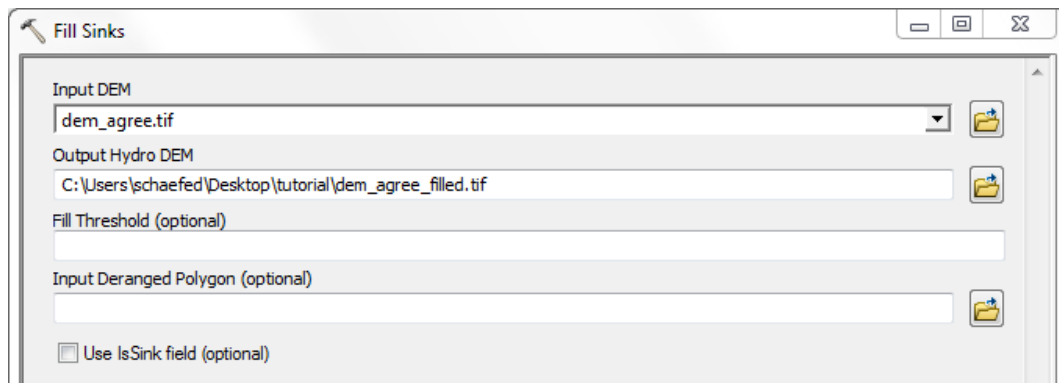


Figure 3.12: System Toolboxes -> Arc Hydro Tools -> Terrain Preprocessing -> Fill Sinks

- Calculate Flow Direction

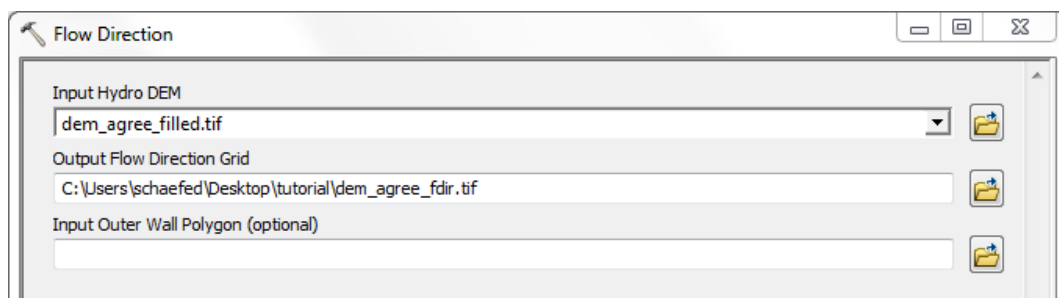


Figure 3.13: System Toolboxes -> Arc Hydro Tools -> Terrain Preprocessing -> Flow Direction

- Create a Flow Accumulation map

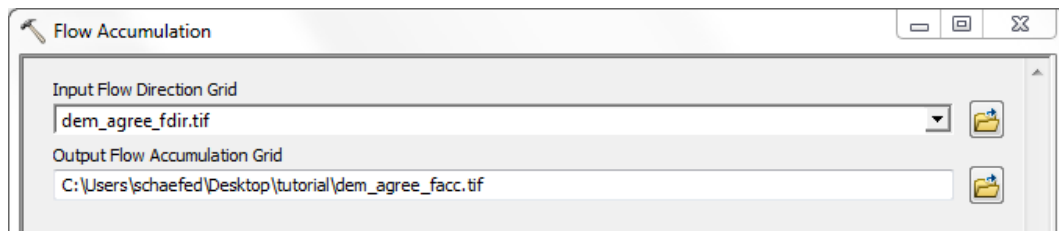


Figure 3.14: System Toolboxes -> Arc Hydro Tools -> Terrain Preprocessing -> Flow Accumulation

### 3.5.6 Gauges map

- Assuming that you have the positions of your gauges in a table, which has at least the columns x, y, id (in any order and/or amongst other columns), you are able to convert your data into a Point Shapefile. Choose the appropriate fields and do not forget to set the coordinate system information.

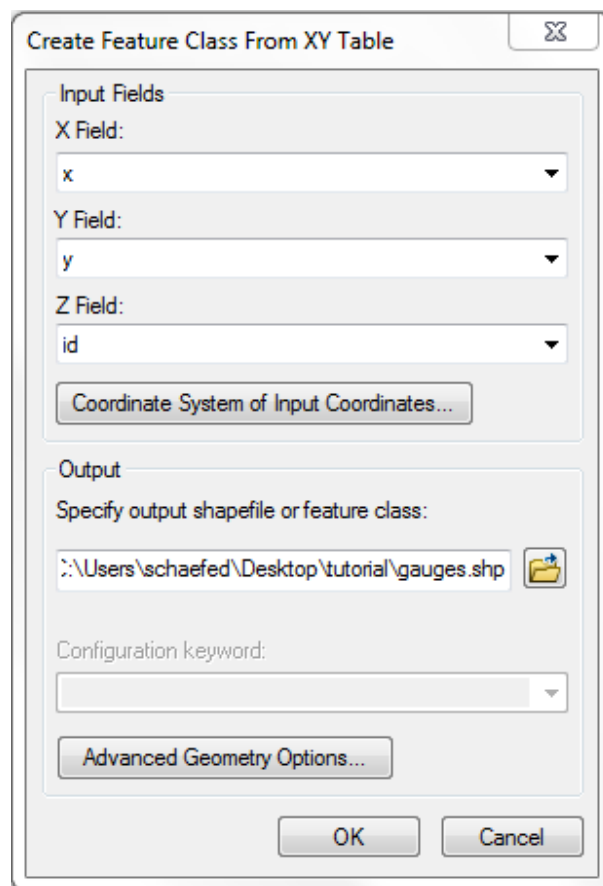


Figure 3.15: Right click on the table in Arc Catalog -> Create Feature Class -> From XY Table

- Check the positions of your gauges against the previously created Flow Accumulation map. If the points do not exactly match the stream-like features on the Flow Accumulation grid, edit the Shapefile (right click on the Point Feature in the Table of Contents -> Edit Features -> Start Editing) and move the gauges to do so. Changing the depiction of the Flow Accumulation map might facilitate this step (i.e. right click the Flow Accumulation map in the Table of Contents -> Properties -> Tab 'Symbology' -> Choose 'Stretched' in the 'Show' box on the left hand side and Type 'Standard Deviations' in the center box. It might also be necessary to invert the color ramp by (un-)checking the 'Invert' check box). Remember to stop the editing process and save your edits (Editor Toolbar -> Editor -> Stop Editing).

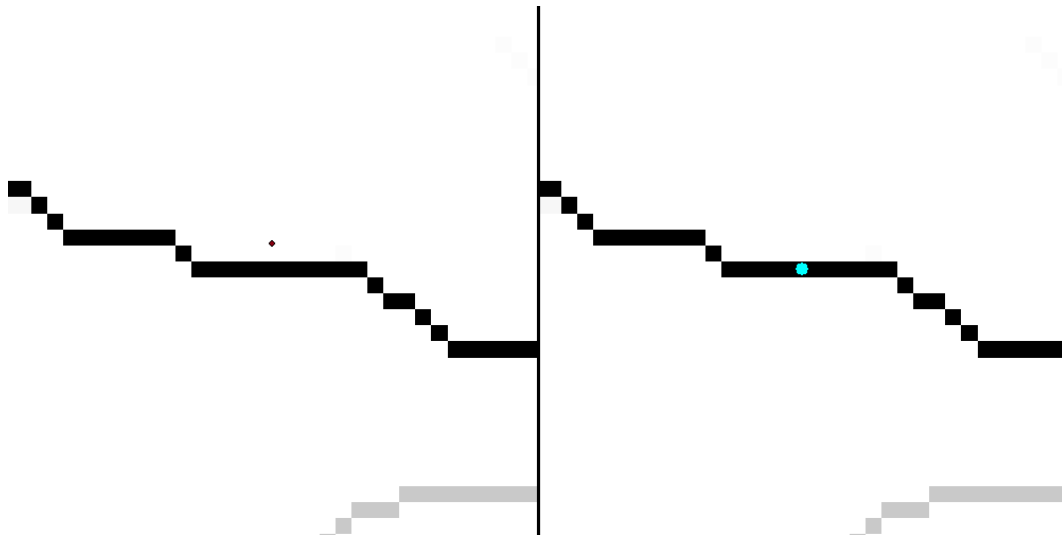


Figure 3.16: Mismatching gauges (left) should be corrected (right)

- Convert the resulting shapefile into a raster map with level-0 resolution.

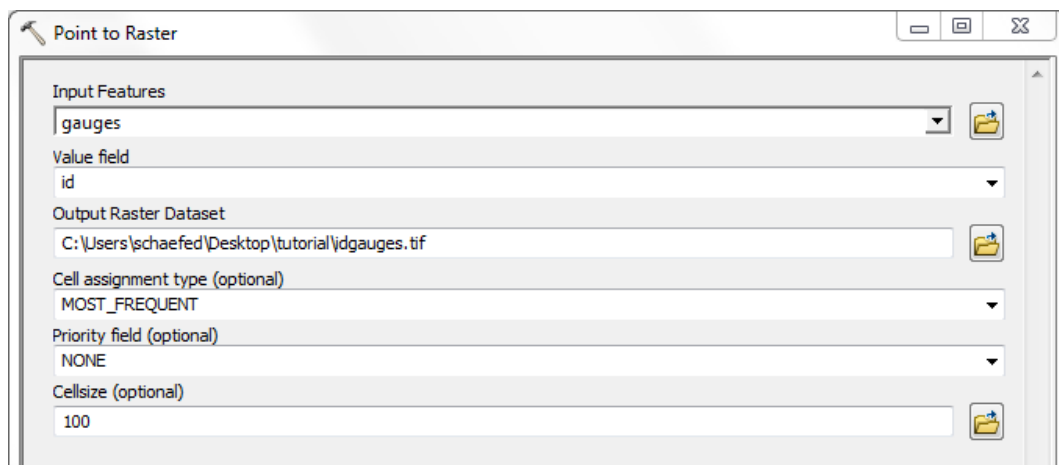


Figure 3.17: System Toolboxes -> Conversion Tools -> To Raster -> Point to Raster

### 3.5.7 Watershed delineation

- Edit the (possibly corrected) gauges shapefile created during the last processing step again and remove all but the outlet gauge. If you need the unedited file, create a copy of the original Shapefile before editing it.
- Delineate the basin

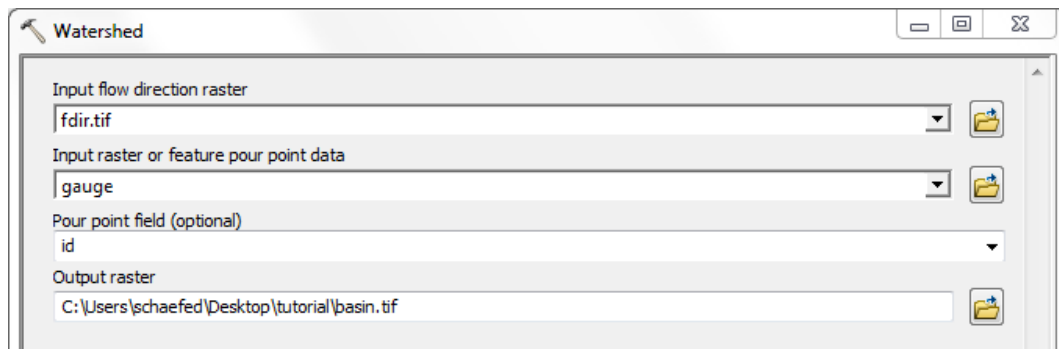


Figure 3.18: System Toolboxes -&gt; Spatial Analyst -&gt; Hydrology -&gt; Watershed

### 3.5.8 Mask the datasets

Mask all mHM input raster files with the created watershed mask

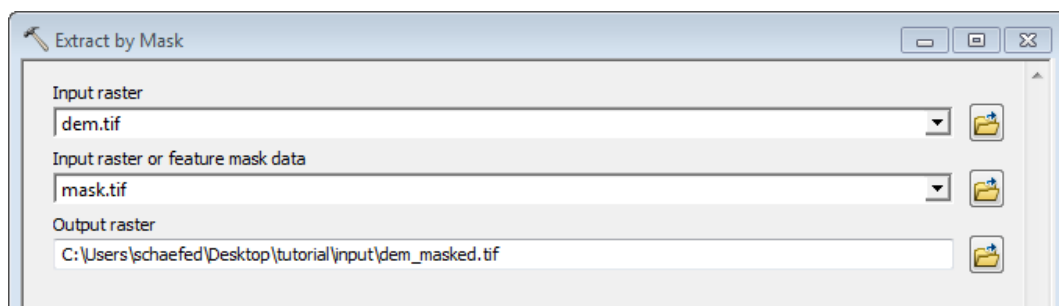


Figure 3.19: System Toolboxes -&gt; Spatial Analysis Tools &gt; Extraction &gt; Extract by Mask

### 3.5.9 Write the ascii grids

Convert the processed and masked raster maps into ASCII files

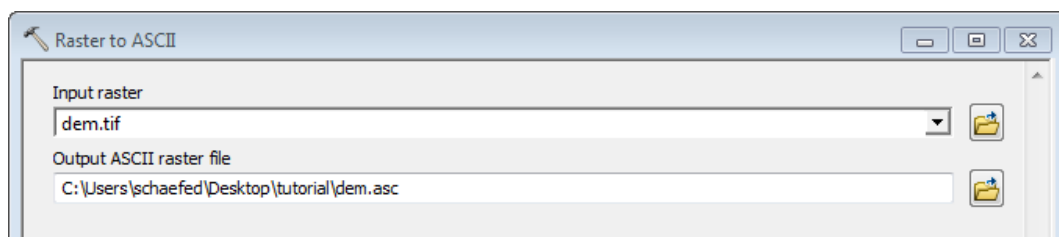


Figure 3.20: System Toolboxes -&gt; Conversion Tools -&gt; From Raster -&gt; Raster to ASCII

## 3.6 Land Cover Data

Unlike the other classified datasets (soils, hydrogeology, LAI), where as much information as available can be added, the land use data is restricted to three different classes, which are listed below:

| Class | Description |
|-------|-------------|
| 1     | Forest      |
| 2     | Impervious  |
| 3     | Pervious    |

### 3.7 Table Data

As previously stated, the input datasets 'soil\_class.asc', 'geology\_class.asc', 'idgauges.asc' and 'LAI\_class.asc' need to be complemented by look-up-tables. All these look-up tables specify the total number of classes/units to read in the very first line, following the keywords 'nSoil\_Types', 'nGeo\_Formations' and 'NoLAIclasses' respectively. The second line acts as a header describing the contents of the following data. In the subsection [The soil look-up table](#) hydrogeo\_table and [The LAI look-up table](#) screenshots of shorted, but sufficient table files are presented. All fields are further listed and described in the respective tables.

#### 3.7.1 The soil look-up table

```
nSoil_Types 2
MU_GLOBAL  HORIZON UD[mm] LD[mm] CLAY[%] SAND[%] BD[gcm-3]
1           1         0       300   23.0   38.0   1.2
1           2       300    1000  31.0   25.0   1.4
2           1         0        50   19.85  53.45  1.41
2           2         50       300   31.33  45.22  1.5
2           3       300    1450  35.42  39.8   1.67
```

Figure 3.21: A possible 'soil\_classdefinon.txt' file

| Field     | Description  |
|-----------|--|
| MU_GLOBAL | Soil id as given in the corresponding 'soil_class.asc' map. All raster map ids must be given here                                    |
| HORIZON   | Horizon number starting at the top of the soil column. You are free to provide any number of horizons for each soil individually     |
| UD[mm]    | Upper depth of the horizon in millimeter. Should be 0 for the first, and the lower depth of the superimposing horizon for all others |
| LD[mm]    | Lower depth of the horizon in millimeter   |
| CLAY[%]   | Clay content in percent  |
| SAND[%]   | Sand content in percent  |
| BD[gcm-3] | Mineral bulk density in grams per cubic centimeter   |

#### 3.7.2 The hydrogeology look-up table

```
nGeo_Formations 2
GeoParam(i)  ClassUnit  Karstic  Description
1           1         0       GeoUnit-1
2           2         0       GeoUnit-2
```

Figure 3.22: A possible 'geology\_classdefinon.txt' file

| Field       | Description  |
|-------------|--|
| GeoParam(i) | Parameter number of the formation, the link to 'mhm_parameter.nml' |
| ClassUnit   | Class id as present in the 'hydrogeology_class.asc' raster map     |
| Karstic     | Presence of karstic formation (0: False, 1: True)                  |
| Description | Text description of the unit                                       |



### 3.7.3 The LAI look-up table

| NoLAIclasses |                     | 10   |      |      |      |      |      |      |      |      |      |      |      |
|--------------|---------------------|------|------|------|------|------|------|------|------|------|------|------|------|
| ID           | LAND-USE            | Jan. | Feb. | Mar. | Apr. | May  | Jun. | Jul. | Aug. | Sep. | Oct. | Nov. | Dec. |
| 1            | Coniferous-forest   | 11   | 11   | 11   | 11   | 11   | 11   | 11   | 11   | 11   | 11   | 11   | 11   |
| 2            | Deciduous-forest    | 0.5  | 0.5  | 1.5  | 4.0  | 7.0  | 11   | 12   | 12   | 11   | 8.0  | 1.5  | 0.5  |
| 3            | Mixed-forest        | 3.0  | 3.0  | 4.0  | 6.0  | 8.0  | 11   | 11.5 | 11.5 | 11   | 9.0  | 4.0  | 3.0  |
| 4            | Sparse-forest       | 2.0  | 2.0  | 3.0  | 5.5  | 6.5  | 7.5  | 7.5  | 7.5  | 6.5  | 4.0  | 2.5  | 2.0  |
| 5            | Sealed-Water-bodies | 0.01 | 0.01 | 0.02 | 0.04 | 0.06 | 0.09 | 0.09 | 0.07 | 0.06 | 0.04 | 0.02 | 0.02 |
| 6            | Viniculture         | 1.0  | 1.0  | 1.0  | 1.5  | 2.0  | 3.5  | 4.0  | 4.0  | 4.0  | 1.5  | 1.0  | 1.0  |
| 7            | Intensive-orchards  | 2.0  | 2.0  | 2.0  | 2.0  | 3.0  | 3.5  | 4.0  | 4.0  | 4.0  | 2.5  | 2.0  | 2.0  |
| 8            | Pasture             | 2.0  | 2.0  | 3.0  | 4.0  | 5.0  | 5.0  | 5.0  | 5.0  | 5.0  | 3.0  | 2.5  | 2.0  |
| 9            | Fields              | 0.4  | 0.4  | 0.3  | 0.7  | 3.0  | 5.2  | 4.6  | 3.1  | 1.3  | 0.2  | 0.0  | 0.0  |
| 10           | Wetlands            | 2.0  | 2.0  | 3.0  | 4.0  | 5.0  | 5.0  | 5.0  | 5.0  | 5.0  | 3.0  | 2.5  | 2.0  |

Figure 3.23: A possible 'LAI\_classdefinion.txt' file

| Field        | Description                                     |
|--------------|---|
| ID           | LAI class id as given in the file LAI_class.asc |
| LAND-USE     | Description of the LAI class                    |
| Jan. to Dec. | Monthly LAI values                              |

### 3.7.4 The gauge files

```
00139:BRUGG/AARE (Abfluss)    DAILY
nodata -9999
n      1      measurements per day [1, 1440]
start 1996 01 01 00 00 (YYYY MM DD HH MM)
end   1996 01 06 00 00 (YYYY MM DD HH MM)
1996 01 01 00 00      368.500
1996 01 02 00 00      408.100
1996 01 03 00 00      493.600
1996 01 04 00 00      502.500
1996 01 05 00 00      453.100
1996 01 06 00 00      439.600
```

Figure 3.24: A possible gauge file

The structure of the gauge files is different from the look-up tables listed so far. The following table describes its content.

| Line     | Description   |
|----------|---|
| 1        | Gives basic information about the gauging station (Station-Id:Station-Name/River-Name)  |
| 2        | Specifies the nodata value  |
| 3        | The number of measurements per day  |
| 4        | The start date of the time series in the format given in parentheses  |
| 5        | The end date of the time series in the format given in parentheses  |
| 6 to end | The measurement data in cubic meter per second preceded by the actual date in the same format as the dates of lines four and five |

For every gauge id given in 'idgauges.asc' one gauge file has to be created as outlined above. The file name has to **exactly** reflect the gauge id and carry a '.txt' extension. The table data file for a gauge with an id of 0343 in 'idgauges.asc' should therefore be named '0343.txt'.

## 3.8 Post-GIS preparation

Make sure that all decimals are indicated with dots, not commas:

```
perl -i.bak -pe 's/[,]/./g' *.asc
```

Make sure that accuracy of your llcorners is reasonable. For example, the following code removes all but one digit after the dot:

```
perl -i.bak -pe 's/([xy].+)\.(\d)\d+/$1\.$2/g' *.asc
```

Make sure that your files cover the same spatial domain. In case your grid origins match, but you are missing rows and columns, the perl script `pre-proc/enlargegrid.pl` will pad your data at the top and the right end of the grid.

```
/basin/input/morph/$ ~/mhm/pre-proc/enlargegrid.pl aspect.pl
31 rows with 47 cols.
/basin/input/morph/$ ~/mhm/pre-proc/enlargegrid.pl 50 40
aspect.asc
dem.asc
facc.asc
fdir.asc
/basin/input/morph/$ ~/mhm/pre-proc/enlargegrid.pl aspect.pl
40 rows with 50 cols.
```

If your data differs more substantially the program 'mHM/pre-proc/resize\_grid.py' will harmonize your grids.

```
python match_grids.py source_grid target_grid out_grid
```

The script takes three Input arguments:

1. The source grid to be enlarged
2. The target grid which could also be an header file as described in subsection [The Header File](#)
3. An output file

The program will fail, if your target grid is smaller than the source grid or if the cellsizes of both grids are not divisable. If necessary the source grid will be shifted in order to make both origins match. This 'shift' is only accomplished by changing the values of the corner coordinates, no real interpolation will be done.

## Chapter 4

# Visualizing Model Output

mHM usually writes two files into the output directory specified in mhm.nml:

```
ConfigFile.log
daily_discharge.out
discharge.nc
mHM_Fluxes_States.nc
mRM_Fluxes_States.nc
```

In addition to these, three restart files are saved into the restart directory also specified in mhm.nml:

```
xxx_config.out
xxx_L11_config.nc
001_states.nc
```

The first file ConfigFile.log is ASCII type and summarizes the main configuration of mHM. The second file daily\_discharge.out is ASCII type and can be read by any standard editor or ASCII interpreting scripts. It contains the simulated and observed discharge. The same timeseries are stored for all gauges in discharge.nc (the third file), but in NETCDF format. In mHM\_Fluxes\_States.nc all the spatial and temporal output is written, with the exception of routed discharge which can be found in mRM\_Fluxes\_States.nc (also a NETCDF file). These binary NETCDF files can be visualised with NCVIEW (part of the NETCDF library), VISIT (LLNL Software) or analysis scripts based on Python (PyNGL). Please note that daily\_discharge.out, discharge.nc and mRM\_Fluxes\_States will only be written if the routing is switched on (i.e., process\_case 8 equals one).

The restarting files are mainly intended for improving model performance (i.e., to avoid spin-up time) rather than for visualization purposes. The xxx prefix indicate the basin number. Since these files are also written into a binary NETCDF file, some remarks will be made below for the right interpretation of their content.

For a quick analysis we recommend NCVIEW since it quickly visualises those files via command-line and X-forwarding. Make sure to load the NC module before starting.

```
ncview FILE.nc
```

In order to hide unnecessary output messages, you may pipe them to a log file:

```
ncview FILE.nc 1> ~/log/ncview.out 2> ~/log/ncview.err
```

If you want to transfer large nc files through servers or visualise them locally, it might be useful to compress the data before. The featured nc file deflation with the ncks module will decrease the file size usually by 30-60%. Choose the deflation level from minimum (0) to maximum (9). The -4 switch in the following command will convert netcdf3 files to version 4 simultaneously.

```
ncks -4 -L 9 IN.nc OUT.nc
```

### Using X-Forwarding under Windows

On Windows we recommend CYGWIN including MINTTY, OPENSSSH and XINIT, XMING as the X server. An alternative is PUTTY using X-forwarding. The workflow in MINTTY is as follows:

```
/bin/XWin.exe -multiwindow
ssh -Y SERVERNAME
module load ncview
ncview FILE.nc
```

In some cases it has proven to be necessary to add the following line to /etc/profile

```
export DISPLAY=:0
```

### Exploring the contents of the Restarting files

Restarting files are simple binary dumps of arrays and vectors of all constants, parameters, state variables (1D, 2D) and fluxes at a given point in time of a simulation that are needed for executing the subsequent mHM time step in a new instance of this model without performing spin-out simulations and additional run time up to the previous time point. The stored information is divided into three categories: 1) Configuration variables at L0 and L1 levels (`xxx_config.nc`), 2) configuration variables at L11 level (i.e., routing) (`xxx_L11_config.nc`), and 3) and effective parameters, state variables and fluxes at L1 and L11 levels (`xxx_states.nc`).

#### WARNINGS

1) Results may appear inverted in NCVIEW with respect to the geographical coordinates used to describe the basin domain. This is due to the lack of geographical coordinates in this NetCDF visualizer and the fact that it simply dumps the content of a 2D array in C convention (row first, DIM2). In addition to that, mHM 2D arrays are stored in (lon, lat) or (X,Y) ordering, which is the transpose of the ESRI convention (lat, lon) or (Y,X).

2) Caution should be taken to interpret flow direction variables (i.e., variable `L11_fDir` in `xxx_L11_config.nc`) if visualized with NCVIEW because of its implicit 90 ° counter-clockwise rotation. Hence the values depicted NCVIEW using 8-flow direction convention (1,2,4,8,16,32,64,128) (i.e., 1 pointing to the east and then moving clockwise every 45 °) have to be rotated accordingly. In other words, direction 1 should be interpreted as 64, 2 as 128, and so on. All 2D variables, however, included the previous one, will produce consistent maps as that shown in ([Moselle Basin](#)) if plotted with appropriate routines (e.g., Python or NCL) that take into account the geographic coordinates stored within the NetCDF file.

## Chapter 5

# Calibration Options

### 5.1 Calibration of Parameters on Observations

In order to use the calibration feature of mHM, turn the `optimize` switch in `mhm.nml` to `.true.`.

#### 5.1.1 Parameters

MHM calibrates all parameters in `mhm_parameters.nml` which are flagged for optimization (column "FLAG"). Parameters can be forced to stay constant during calibration by setting the flag to 0. The upper and lower bounds define the ranges in which the optimization methods vary the parameters. Although permitted, we recommend not to change the bounds, since they are the result of intensive sensitivity studies covering a wide range of basins.

#### 5.1.2 Methods

Different optimization methods are available to find the best configuration of parameters for the selected objective function. You may choose between Simulated Annealing (SA), Dynamically Dimensioned Search (DDS) and Shuffled Complex Evolution algorithm (SCE). Details about these methods can be found in the module description part of this manual. At the very end of `mhm.nml` additional settings are offered for optimization.

#### 5.1.3 Functions

The measure to define how well a quantity fits to the observations depends strongly on the specific type of application. mHM offers a variety of options for different quantities like discharge, soil moisture, or total water storage. For example, an NSE type of function will not be able to catch low values as well as  $\ln(\text{NSE})$  would. In `mhm.nml`, individual objective functions are defined for specific quantities.

#### 5.1.4 Data

Discharge and total water storage data should be provided in ASCII file format for each gauge/basin. Soil moisture and neutron data need to be provided as spatio-temporal netcdf files. Example files can be found in the folder `optional_data` in the `test_basin`. Furthermore, a `optional_data` namelist in `mhm.nml` provides several options regarding the handling of such data. For example, for soil moisture data, the temporal resolution needs to be set in `mhm.nml`. However, neutron data is restricted to be daily in the current release. The spatial resolution of soil moisture and neutrons data needs to match the hydrological resolution of mHM.

For correct preparation of input data, please take a look at the input data in the `test_basin`, which will serve as a good example. Furthermore, it could be helpful to generate optional data files from precedent mHM output, because then the temporal resolution and the spatial grid is guaranteed to match the needs of mHM. The files then can be modified with tools like `cdo` or `python`.

### 5.1.5 Output

Calibration runs result in a parameter file called `FinalParams.nml` which contains the optimized parameter set. Replace `mhm_parameters.nml` with `FinalParams.nml`, then run mHM in forward mode in order to obtain hydrologic predictions.

## Chapter 6

# Coding and Documentation Style

The coding and documentation style guide has the following sections:

[General](#)

[In and Out of Files](#)

[Modules, Subroutines and Functions](#)

[Variables](#)

[Documentation](#)

### General

- Follow the coding and documentation style of template [mo\\_template.f90](#) very closely.
- Enforce SI units in all code, with the only exception of  $\text{mms}^{-1}$  for  $\text{kg m}^{-2}\text{s}^{-1}$ .
- mHM is computed in double precision (dp).  

```
real(dp) :: wind_speed ! wind speed in [m s-1]
```
- New routines should be tested with at least two different compilers (of different vendors).

### In and Out of Files

- Filenames are all lower case.
- Module names and filenames are the same and start with mo\_
- Filenames should be descriptive of the content and use as little abbreviations as possible.
- Fortran filenames end with .f90  

```
mo_string_utils.f90
```
- Never use tabs.
- Break lines at column 130, at most. This includes comments  

```
sum_of_all = a + b + c + d + e + f + g + h + i + j + k + l + m + n + o + p + q + r + s + &  
           t + u + v + w + x + y + z  
zero      = 0.0_dp ! This literal can be used whenever something has to be initialised, so that  
               ! it will be initialised with the right number precision
```
- Indent all code blocks.  

```
if (abs(a-b)>epsilon(1.0_dp)) then  
  if (a >= b) then  
    a = b  
  else  
    b = a  
  endif  
endif  
endif
```

## Modules, Subroutines and Functions

- All modules and programs have IMPLICIT NONE.
- All USE statements have the only clause.
- Modules are PRIVATE by default.
- Module routines are made available explicitly, i.e. PUBLIC.

- Give 1-line descriptions after the public definition.

```

module mo_calc
  use mo_kind, only: i4, dp
  implicit none
  ...
  public :: calc ! Calculates the thing
contains
  function calc(x)
  ...

```

- Try to write elemental pure subroutines whenever possible (see below).
- Extremely short subroutines should be avoided.
- Avoid very long subroutines (>500 lines).

## Variables

- Use expressive and descriptive variable names for all variables in the namelist and for all variables that have a larger scope, i.e. that are used in more than a screen full. This means that counter variables can still be i, ii, j, ... and local variables can also be called short names such as tmp or itmp. But variables, which are important to read the formula, should have descriptive names such as ido\_that or iDoThat. Variable names should not be too long either. The agreed maximum numbers are

- filenames < 30 characters
- variable names < 20 characters

- All variables and literals use explicit type declarations from mo\_kind.

```

elemental pure function circum(radius)
  use mo_kind, only: dp
  use mo_consts, only: pi_dp
  implicit none
  real(dp), intent(in) :: radius
  real(dp) :: circum
  circum = 2.0_dp * pi_dp * radius
end function circum

```

- Array dimensions are in the following order: lon (east-west), lat (north-south), z (vertical), t (time level). Additional dimensions can be used and are placed in front of t.
- Always pass arrays as sub-program arguments. This means: use as little global variables as possible.
- All input/output arguments have an intent.

```

subroutine calc(x,y)
  use mo_kind, only: i4, dp
  implicit none
  real(dp), dimension(:, :), intent(in) :: x
  integer(i4), dimension(size(x,1),size(x,2)), intent(out) :: y
  ...

```



- Use intuitive names for optional arguments.

```
function calc(x, inverse)
  use mo_kind, only: i4, dp
  implicit none
  real(dp), dimension(:, :), intent(in) :: x
  logical, optional, intent(in) :: inverse
  integer(i4), dimension(size(x,1), size(x,2)) :: calc
  logical :: iinverse
  iinverse = .false.
  if (present(inverse)) then
    iinverse = .false.
    if (inverse) iinverse = .true.
  endif
  ...
end function
```

- Global variables are initialised at the setup stage.
- Never use hardcoded number literals such as 5., 3.14, etc. Exceptions might be -1, 0, 1, and 2. But all literals are appended by `_dp`, `_i4`, etc.

```
elemental pure function circum(radius)
  use mo_kind, only: dp
  use mo_consts, only: pi_dp
  implicit none
  real(dp), intent(in) :: radius
  real(dp) :: circum
  circum = 2.0_dp * pi_dp * radius
end function circum
```

- Avoid pointers whenever possible.
- Index notation is encouraged whenever whole arrays are treated such as `a(:) = b(:)*c(:)`. It is still very good for contiguous subarrays such as `a(1:n-1) = b(1:n-1)*c(1:n-1)`. It is discouraged, though, for mixed indexing such as `a(1:n-1) = b(2:n)*c(1:n-1)`. This is (1) prone to coding error, (2) not easy to read by other programmers, and (3) hard to parallelise with OpenMP or MPI. Use `forall` or `do` instead.

## Documentation

Follow the documentation structure before the interface mean in [mo\\_template.f90](#).

Documentation uses the automatic documentation system doxygen ( <http://www.doxygen.org/>). See the manual for the complete list of documentation tags: <http://www.doxygen.nl/manual.html>

- The documentation of a routine is in front of the routine definition.
- Documentation for a module interface must be in front of the interface definition. The individual routines do not need extra documentation.
- Only public elements will be considered by the automatic documentation system doxygen.
- Make your routines public at the beginning of the file. Append a one-line descriptions to the public statement.



## Chapter 7

# The details about the test basin

The test data coming with the mHM source code are for the Moselle River basin upstream of Perl, a place, where the Moselle River leaves France and enters Luxembourg and Germany ([Moselle Basin](#)). The catchment area is approximately 11,500 km<sup>2</sup>, altitude ranges between 150 and 1300 m. a.m.s.l. The Moselle River originates from the Vosges Mountains and is a tributary of the Rhine River. The origin of data used in the test example is listed below.

### **Meteorological forcings (temperature and precipitation):**

We acknowledge the E-OBS dataset from the EU-FP6 project ENSEMBLES ( <http://ensembles-eu.metoffice.com>) and the data providers in the ECA&D project ( <http://www.ecad.eu>).

"Haylock, M.R., N. Hofstra, A.M.G. Klein Tank, E.J. Klok, P.D. Jones, M. New. 2008: A European daily high-resolution gridded dataset of surface temperature and precipitation. J. Geophys. Res (Atmospheres), 113, D20119, doi:10.1029/2008JD10201".

( <http://www.ecad.eu/download/ensembles/ensembles.php> 02.04.2014)

### **Hydrological observations:**

We acknowledge the Global Runoff Data Centre (GRDC), a repository for the world's river discharge data and associated metadata for providing the discharge data.

( [http://www.bafg.de/GRDC/EN/01\\_GRDC/12\\_plcy/policy\\_guidelines.pdf;jsessionid=AA96F6F200B3880575879F15534B6669.live2052?\\_\\_blob=publicationFile](http://www.bafg.de/GRDC/EN/01_GRDC/12_plcy/policy_guidelines.pdf;jsessionid=AA96F6F200B3880575879F15534B6669.live2052?__blob=publicationFile) 02.04.2014)

### **SRTM (Shuttle Radar Topography Mission):**

We acknowledge the U.S. Geological Survey's Earth Resources Observation and Science (EROS) Center or NASA's Land Processes Distributed Active Archive Center (LP DAAC) for providing the digital elevation model.

( <http://www2.jpl.nasa.gov/srtm/cbanddataproducts.html> 15.05.2014)

### **Harmonized world soil database:**

We acknowledge the use of Harmonized World Soil Database dataset of the Food and Agriculture Organization of the United Nations (FAO) the International Institute for Applied Systems Analysis (IIASA), International Soil Reference and Information Centre (ISRIC), Institute of Soil Science – Chinese Academy of Sciences (ISSCAS) and Joint Research Centre of the European Commission (JRC).

( <http://webarchive.iiasa.ac.at/Research/LUC/External-World-soil-database/HTML/> 02.04.2014)

### **European soil database:**

We acknowledge the European Commission for providing the European soil database.

( [http://ec.europa.eu/geninfo/legal\\_notices\\_en.htm](http://ec.europa.eu/geninfo/legal_notices_en.htm) 02.04.2013)

### **Land cover:**

We acknowledge the European Environment Agency for providing the CORINE land cover dataset.

( <http://www.eea.europa.eu/publications/COR0-landcover> – 15.04.2014)

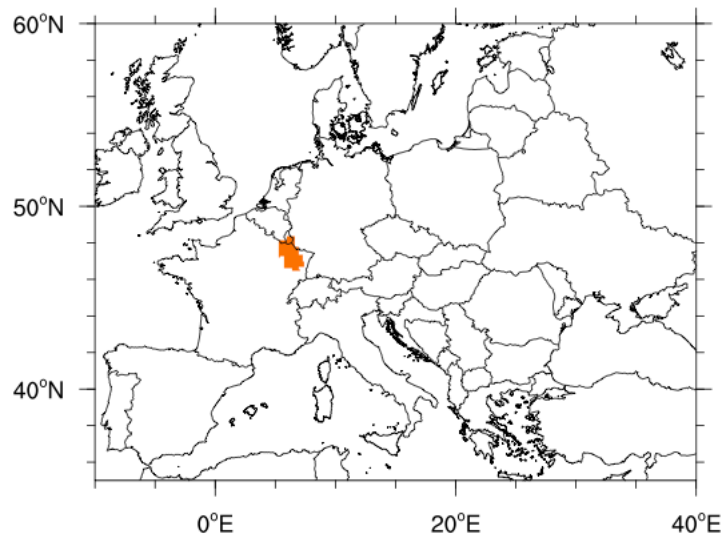


Figure 7.1: Location of the Moselle River basin upstream of Perl within the European domain

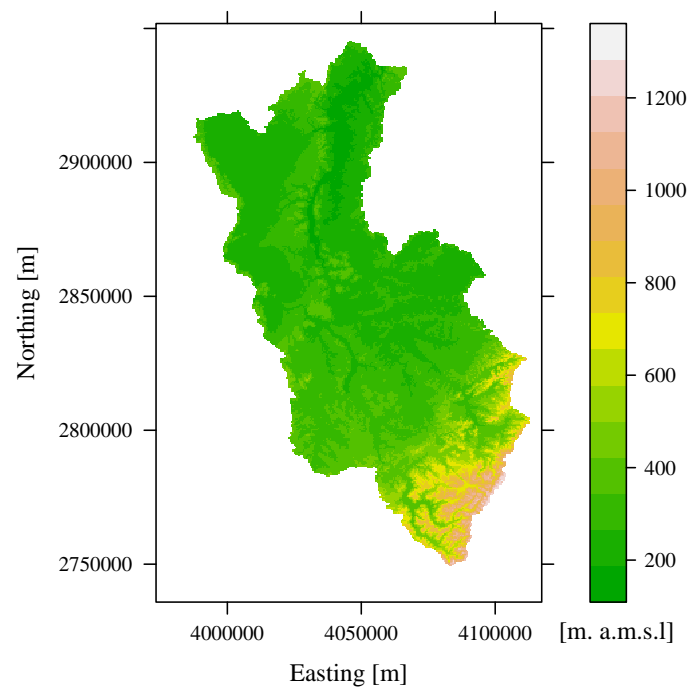


Figure 7.2: Digital elevation model for the Moselle River basin upstream of Perl

## Chapter 8

# Protocols for setting up a new mHM basin

The following checks and protocols describe some standard procedures for the setup of a new basin in mHM. They are based on the personal experience of the authors and do not cover all possible subjects. They are to be seen as a first guideline when working with a new basin in mHM but do not represent a complete instruction.

This chapter is divided into the following sections:

- check for input data errors
- protocol for generating a restart file
- protocol for determining the warm-up period for a new basin

### 8.1 Check for possible input data errors using mHM outputs

Besides mHM check for the possible input data errors (in the [mo\\_startup](#) routine), there could still be room for the data errors. Before you make the exhaustive model runs for calibration, make sure that you have processed your input data correctly. Below are some tips for detecting data problems using the mHM outputs.

1. Make a default mHM simulation for a relatively longer time-period, as supported by the input data sets (say 20-30 years).
2. Print out distributed fluxes/states, for example, at a monthly time scale. This feature is supported in the current mHM version using the name list file "mhm\_outputs.nml". You just need to tick ".TRUE." to the fluxes/ states you want to save as a netcdf file.

Check at-least the following model outputs:

- (a) height of snow pack (L1\_snowpack) [mm] – case 2 outputFlxState(2)=.TRUE.
  - This variable should be check in those basins which have snow covers/glaciers.
  - Snowpacks in the mountainous/glacier regions should be higher than those of other areas.
  - Snowpack during winter should be higher than that during summer.
- (b) mean volumetric soil moisture averaged over all soil layers [mm/mm] – case 5 ( = L1\_soilMoist / L1\_↔ soilMoistSat ) outputFlxState(5)=.TRUE.
  - Soil moist. during winter should be higher than that of during summer.
  - In general, soil moist. in hilly areas is higher compared to that in flatter areas
- (c) actual evapotranspiration aET [mm/T] – case 9 outputFlxState(9)=.TRUE.
  - Evapotrans. during winter should be lower than that of during summer.
  - In general, evapotrans. in hilly areas is lower compared to that in flatter areas
- (d) total discharge generated per cell (L1\_total\_runoff) [mm/T] – case 10 outputFlxState(10)=.TRUE.

- In general, runoff in hilly areas is higher compared to that in flatter areas
3. Check the temporal dynamics of modeled fluxes/states across several locations (i.e., at several grid cells).
    - They should, in general, exhibit distinct behavior in the temporal dynamics.
  4. You could calculate a long-term mean annual dynamics of variables like
    - (a) actual evapotranspiration
    - (b) total runoff
    - (c) precipitation (use values provided in the meteo-input file)
    - This could fairly give you an idea about the respective water balance regime of the modeled basin. For example, hilly areas should exhibit a higher runoff, as a result of a relatively higher precipitation and lower ET values, compared to those of the flat areas.
    - Compute the values of runoff coefficient (long-term mean annual runoff / long-term mean annual precipitation) These values should be lesser than 1.0 everywhere.
    - Compute the values of long-term mean annual evapotranspiration / long-term mean annual pot. evapotranspiration These values should be also lesser than 1.0 everywhere.
  5. Compare the respective spatial patterns of each variable with some known literature results (google them for your respective basins)

Above check list are just few (based on the personal experience of the authors) among many possible ones, and the list will be updated as soon as we find new ways to deal with data problems.

## 8.2 Protocol for generating a restart file

Not specified yet.

## 8.3 Protocol for determining the warm-up period for a new basin

Not specified yet.

## Chapter 9

# mRM - the Multiscale Routing Model: Running the stand-alone version of the routing model

### 9.1 Introduction to mRM

This section provides information on how to run the routing model mRM as a stand-alone program outside of mHM. The general philosophy is as follows: mRM can be compiled and run in a similar way as mHM. Specific files for mRM can be found in the main path and are denoted as <filename>\_mRM.<file\_type>, where <filename>.<file\_type> is the corresponding mHM file.

This Subsection provides a short description of the method used in mRM. For further details and an analysis over the European domain, please refer to Thober et al. 2019.

#### 9.1.1 Finite Difference Approximation of Kinematic Wave Equation

The multiscale Routing Model mRM uses the kinematic wave equation to describe the water flow within a stream as

$$\frac{\partial Q}{\partial t} + c \frac{\partial Q}{\partial x} = 0 \quad (9.1)$$

where  $Q$  ( $m^3, s^{-1}$ ) is streamflow,  $x$  (m) the space dimension,  $t$  (s) the time dimension, and  $c$  ( $m, s^{-1}$ ) the celerity. The kinematic wave equation is a simplification of the Saint-Venant equations. The derivation is based on the assumption that the continuity equation is sufficient to describe the movement of flood waves. In detail, a constant river bed slope and time-invariant celerity  $c$  have to be assumed. mRM employs the classical finite difference weighted approximation on a four points scheme to solve the equation above. A short derivation is as follows:

The partial derivatives within the above equation are represented as finite differences between four values, that means on two locations at two points in time:

$$\begin{aligned} \frac{\partial Q}{\partial t} &\approx \frac{\varepsilon(Q(x_j, t_{i+1}) - Q(x_j, t_i)) + (1 - \varepsilon)(Q(x_{j+1}, t_{i+1}) - Q(x_{j+1}, t_i))}{\Delta t}, \\ \frac{\partial Q}{\partial x} &\approx \frac{\varphi(Q(x_{j+1}, t_{i+1}) - Q(x_j, t_{i+1})) + (1 - \varphi)(Q(x_{j+1}, t_i) - Q(x_j, t_i))}{\Delta x}, \end{aligned}$$

where  $j$  denotes the spatial location (i.e., reach id) and  $i$  denotes the timestep.  $\varepsilon$  is a space-weighting factor and  $\varphi$  is a time-weighting factor. mRM uses a rectangular grid to represent the river network with a river reach in the model connecting two center grid locations. Different spatial locations are separated by  $\Delta x$  and time steps by  $\Delta t$ . The two weighting factors,  $\varepsilon$  and  $\varphi$ , can be chosen between 0 and 1, but the numerical solution becomes unstable for  $\varepsilon > 0.5$ . The numerical diffusion depends linearly on  $\varepsilon$ , with 0 implying full numerical diffusion and 0.5 no numerical diffusion, respectively.

Setting  $\varphi$  to 0.5, which represents a time-centered scheme, and substituting the above equation into equation (9.1) results in the classical linear equation:

$$Q(x_{j+1}, t_{i+1}) = C_1 Q(x_j, t_{i+1}) + C_2 Q(x_j, t_i) + C_3 Q(x_{j+1}, t_i), \quad (9.2)$$

with the coefficients  $C_1$ ,  $C_2$  and  $C_3$  being:

$$\begin{aligned} C_1 &= \frac{-2\Delta x \varepsilon + c \Delta t}{2\Delta x(1-\varepsilon) + c \Delta t}, \\ C_2 &= \frac{2\Delta x \varepsilon + c \Delta t}{2\Delta x(1-\varepsilon) + c \Delta t}, \\ C_3 &= \frac{2\Delta x(1-\varepsilon) - c \Delta t}{2\Delta x(1-\varepsilon) + c \Delta t}. \end{aligned} \quad (9.3)$$

The coefficients  $C_1$ ,  $C_2$  and  $C_3$  add up to one. The spatial resolution at which mRM is applied is called "routing" resolution in the following.

### 9.1.2 Stream Celerity Parametrization based on Terrain Slope

Two parametrizations of equation (9.3) are available in mRM: firstly the regionalized Muskingum-Cunge (rMC) parametrization with a fixed time step, and secondly a parametrization using spatially varying celerities in combination with an adaptive time step (aTS). A short summary of the former is presented in Subsection [Regionalized Muskingum-Cunge parametrization](#) and is referred to as rMC in the following. The latter is described in detail in this and the next section and is referred to as aTS scheme.

The aTS calculates stream celerity as a function of terrain slope using a simple relationship:

$$c_i = \gamma \sqrt{s_i}, \quad (9.4)$$

where  $c_i$ ,  $s_i$  and  $\gamma$  are celerity, terrain slope and a free model parameter, respectively, and  $i$  is the grid cell index.

The above equation is applied at the resolution of the Digital Elevation Model (DEM), from which terrain slope is derived. A median absolute deviation (MAD) filter is applied to the high resolution slope data along the path of the main river with a threshold value of 2.25 to correct for outliers. Large slope outliers happen easily in DEMs, for example, when the river flows in a valley and one grid cell represents the valley while the next grid cell represents the hill top. A minimum river bed slope of 0.1% is further assumed. The celerities are then upscaled to the routing resolution, i.e., the resolution at which the kinematic wave equation is solved. The upscaling is by averaging with the harmonic mean, the correct averaging operator for celerities (or speed). The upscaling considers also only those high-resolution grid cells that align with the path of the main river because aTS only considers the flow in the main river reach, assuming that travel times in the main reach dominate the routing process in tributaries.

### 9.1.3 Adaptive Time Step (aTS) Implementation

The aTS scheme uses an adaptive time step to calculate the linear coefficients in equation (9.2). The basic idea is that the time step should be such that water has not been transported further than  $\Delta x$  during a single step. This condition is generally known as Courant-Friedrich-Lewy criterium, which is a necessary condition for numerical stability of finite difference schemes [8]. The condition can be expressed as:

$$C_r = \frac{c \Delta t}{\Delta x} \leq C_{max} = 1 \quad (9.5)$$

where  $C_{max}$  is the Courant number. aTS uses a Courant number of 1. The Courant condition couples the applied spatial resolution with the integration time step of the finite difference scheme. Celerities  $c_i$  are typically in the order of a few  $\text{m s}^{-1}$ . The time step  $\Delta t$  is chosen such that it does not deviate too much from the Courant number  $C_{max}$  to keep computational demand to a minimum. This implies that  $\Delta t$  ranges from a few minutes for high-resolution grids to a few hours for continental scale applications.

In detail, aTS chooses  $\Delta t$  from a set of prescribed values such that  $c_i \Delta t / \Delta x$  is close to but less than 1 for all celerities  $c_i$ . The prescribed values range from one minute to one day, namely: 1 min, 2 min, 3 min, 4 min, 5 min, 6 min, 10 min, 12 min, 15 min, 20 min, 30 min, 1 h, 2 h, 3 h, 4 h, 6 h, 8 h, 12 h, and 1 day. The choice of these values is



motivated from the fact that these represent multiples or dividers of hourly and daily time steps. These time steps allow in principle model applications from 100 m to 100 km, for typical celerities around  $1.5 \text{ m s}^{-1}$ .

Note that the chosen time step depends only on the spatial resolution and is independent of the time resolution of the provided forcing. For example, applying aTS at 12 km resolution using a celerity of  $c \sim 1.5 \text{ m s}^{-1}$  gives  $\Delta x/c$  of 2.2 hours and, hence, a time step of two hours will be chosen. If mRM is forced with hourly input, it will aggregate the input over two consecutive time steps prior to the routing. The calculated streamflow is then distributed to the previous two time steps because these represent the mean flow over this period. If aTS is forced with daily input, it will use internally 12 iterations of 2-hour time steps to route the water through the river network. In this case, aTS will also return the average of the calculated 12 streamflow values at each reach.

### 9.1.4 Regionalized Muskingum-Cunge parametrization

The regionalized Muskingum-Cunge (rMC) parametrization implemented in the mesoscale Hydrologic Model mHM calculates the Muskingum coefficients  $C_1$ ,  $C_2$ , and  $C_3$  in equation (9.2) as a function of high-resolution river network properties. The coefficients  $C_1$ ,  $C_2$ , and  $C_3$  are parametrized as follows

$$C_1 = v_2; C_2 = v_1 - v_2; C_3 = 1 - v_1, \quad (9.6)$$

where the parameters  $v_1$  and  $v_2$  are given as

$$\begin{aligned} v_1 &= \frac{\Delta t}{\beta(1-\varepsilon) + \frac{\Delta t}{2}}; \\ v_2 &= \frac{\frac{\Delta t}{2} - \beta\varepsilon}{\beta(1-\varepsilon) + \frac{\Delta t}{2}} \end{aligned} \quad (9.7)$$

following the nomenclature of appendix A2 in [16]. This formulation is identical to equation (9.3) of the present study, using  $\beta = \Delta x/c$  in the equation above and substituting this equation into two equations above. The parameters  $\beta$  and  $\varepsilon$  are then conceptualized as

$$\begin{aligned} \beta &= \gamma_1 + \gamma_2 L + \gamma_3 S + \gamma_4 C; \\ \varepsilon &= \gamma_5 \frac{S}{\max(S)}, \end{aligned} \quad (9.8)$$

where  $L$  is the length of the reach,  $S$  is the slope of the reach, and  $C$  is the fraction of impervious land cover within the floodplains (see table 4 in [12]). Overall, there are five global parameters  $\gamma_1$  to  $\gamma_5$  in the above equation that can be chosen by the user. The integration time step is fixed at one hour. To guarantee the numerical stability of the parameterization, the following upper and lower bounds are applied

$$0 < \varepsilon \leq 0.5, \quad (9.9)$$

$$\frac{\Delta t}{2(1-\varepsilon)} < \beta \leq \frac{\Delta t}{2\varepsilon}, \quad (9.10)$$

where  $\Delta t$  is set to one hour.

## 9.2 Getting Started

For receiving mRM, installing NETCDF, and running mRM under CYGWIN on Windows, please follow Sections [Install NETCDF](#), [Run mHM under CYGWIN on Windows](#), and [Compile mHM](#).

## 9.2.1 Compiling mRM

The compilation of mRM can be achieved by running the *make* command on the **Makefile\_mRM** file. This can be done in two ways:

1. `make -f Makefile_mRM`
2. renaming of `Makefile_mRM` to `Makefile` via

```
mv Makefile_mRM Makefile
```

## 9.2.2 Test mRM on Example Basin

The received mHM distribution contains an example test basin in the folder `test_basin/`

This folder also contains all necessary data to run mRM. After you compiled mRM, simply run the executable via `./mrm`

Output will be written in the `test_basin/output_b1`

## 9.2.3 Run your own Simulation

As for mHM, there are three mRM configuration files. These are:

- `mrm.nml`
- `mrm_output.nml`
- `mrm_parameter.nml`

Please note that the information contained in these files can also be found in the respective files for mHM because mRM is part of mHM.

### 9.2.3.1 Main configuration: Path, Periods, Switches

The file `mrm.nml` contains the main configuration for running mRM in your catchments. Since the comments should explain each single setting, this section will only roughly describe the structure of the file. All mandatory namelists have to be specified explicitly for both mHM and mRM. Namelists that are specific for mRM are denoted by `_mRM`. Please keep in mind, that paths can be relative, absolute and even symbolic links.

- **project\_description:** Defines meta data that is written in the output netcdf files.
- **mainconfig:** Defines number of basins and hydrology resolution.
- **mainconfig\_mhm\_mrm:** Defines input paths for reading restart files, routing resolution, and optimization.
- **mainconfig\_mrm:** Defines file name, variable name, and data convention for runoff input field for mRM.
- **directories\_general:** Defines common input and output paths for mHM and mRM.
- **directories\_mRM:** Defines paths where gauge information and input runoff fields can be found.
- **processSelection:** Only processCase seven is used for mRM, all other options are ignored.
- **LCover:** defines information on LAI data.
- **time\_periods:** Defines simulation periods.
- **evaluation\_gauges:** Defines number of gauges per basin and file names containing gauge observations.
- **Optimization:** Defines parameters for optimization.

### 9.2.4 Output Configuration: Time Steps, States, Fluxes

The file `mrm_outputs.nml` regulates how often (e.g. `timeStep_model_outputs_mrm = 24`) and which variables (fluxes and states) should be written to the final netcdf file `[OUTPUT_DIRECTORY]/mRM_Fluxes←_States.nc`. We recommend to only switch on variables that are of actual interest, because the file size will greatly increase with the number of containing variables. During calibration (see [Calibration and Optimization](#)) no output file will be written.

### 9.2.5 Calibration and Optimization

The parameter calibration can be conducted in the same way as for mHM. Please see Section [Calibration and Optimization](#) and Section [Calibration of Parameters on Observations](#) for details. Parameters contained in the `mrm_parameter.←nml` are considered during the optimization.

## 9.3 Data preparation for mRM

The data preparation for the static data for mRM is identical to that of mHM. It is described in the Chapter "Data preparation for mHM" in subsection [Extract Data to the Size of a Catchment](#), [Extact Data to the Time Period of Interest](#), [Preparation of the LatLon Grid](#), [Preparation of the Morphological Data](#) (with the exception of soil map, hydrogeological map, LAI map), [A possible GIS workflow](#) (with the exception of aspect), [Land Cover Data](#), and [Table Data](#).

Additionally, forcing data is required. The forcing for mRM is grid-scale runoff. This runoff needs to be provided in a netcdf file that follows the input format of mHM for forcing variables. The only difference to the forcing variables is that the variable and file name can be both provided in the `mrm.nml` file in the namelist `mainconfig_mrm`.

Please see other details in Section [Preparation of the Forcings](#).



## Chapter 10

# mHM Dependencies

### 10.1 NetCDF4

Reading and writing of input and output files requires the NetCDF4 library to be present. For example, version 4.1.2 can be [downloaded from here](#). See also the [NetCDF 4.1.2 release notes](#).



# Chapter 11

## Install Instruction

The section 'Dependencies' lists the general requirements for the installation. The section 'System-dependent installation instructions' gives some instructions on how to install these dependencies on Windows, some Linux-distributions and soon MacOS.

The section 'Specific setups' can be skipped in most cases. If you are using a module system or MacOS there are some additional hints for the installation process. It might be helpful to first read the section 'Installation'. The section 'Specific setups' then helps to adjust the commands for a build on the specific system.

The section 'Installation' then is a step-by-step guide to install mHM in the command line.

### 11.1 Dependencies

For Windows, some Linux distributions and soon also MacOS specific installation instructions for the following list can be found below.

- a fortran compiler
- make (a tool to compile a program)
- cmake (version  $\geq 3.5$ ) (a tool to create a system-dependent makefile)
- fitting netcdf-fortran libraries (libraries for the use of the data format netcdf on which mhm depends)
- (optional, but makes things much easier) git

Git is a version-control system. If you want to contribute to a project, it is highly recommended to use Git. You can use Git to download (clone) the project to your local pc and have a look at the history or synchronize it without copying the whole repository again. You can also download the project folder without Git, but this would not allow you to pull updates from and push changes to our repository.

### 11.2 System-dependent installation instructions

#### Windows:

**Cygwin** is an environment with a terminal that allows to compile and run programs of Unix-like systems. You can find further instructions to install cygwin on the webpage, as well as instructions on how to install further dependencies after the installation.

After the installation of cygwin and its dependencies mHM will be installed using cygwin. All commands and the execution of mHM only run in that environment.

Install cygwin by executing the cygwin setup and choose the following dependencies:

- [ ] gcc-fortran (the fortran compiler)
- [ ] make
- [ ] cmake (version  $\geq 3.12$ )
- [ ] libnetcdf-fortran-devel
- [ ] libhdf5-devel
- [ ] libgfortran
- [ ] gfortran

While installing cygwin you will have to choose a mirror. A mirror is a server on the internet where the files for the installation come from. Choose any server located near your city and when in doubt, choose the first one in the list. In the next step you can find all available packages provided by cygwin, set the view to "full". In the search panel you can filter the packages by the dependencies listed above (e.g. make). When you choose a version, the newest one is usually a good choice if not marked as experimental.

*Note for UFZ members:* Install cygwin locally (on your own computer), do not choose a location on the network for the installation.

Some cygwin versions create a new home directory for you. You may check e.g. here:

```
C:\cygwin64\home\$username
```

As from December 2019, step-by-step guidelines, how to install all cygwin libraries can be viewed in [this youtube video](#) created by Mehmet Cüneyd Demirel (Istanbul Technical University).

Step-by-step mHM compilation in CYGWIN platform

- 1) Change directory to mHM folder. Use single quote e.g. 'D:/mhm-v5.11.1/' if there is space in the path. `cd mhm`
- 2) Make a sub-directory inside mHM folder e.g. `build mkdir build`
- 3) Change directory to build subfolder `cd build`
- 4) Execute `cmake` with the path to the Git source directory as parameter. (If you followed the instructions above, the path is `..`) `cmake ..`

Other `cmake` options:

To avoid memory issues, allocate stack memory during `cmake`

```
cmake -DCMAKE_Fortran_FLAGS="-Wl,--stack,12485760" ..
```

If you will run mHM in parallel using OpenMP then you will need Microsoft MPI installed in your PC. Search for "Download Microsoft MPI" on internet.

Then use `cmake` option below. Note that memory dump is a common issue for cygwin users when compiling with OpenMP. For memory allocation please also use this line below.

```
cmake -DCMAKE_Fortran_FLAGS="${CMAKE_Fortran_FLAGS} -Wl,--stack,12485760" -DCMAKE_WITH_OpenMP=ON -DCMAKE_BUILD
```

- 4) Execute `make make`
- 5) If all went well then `mhm.exe` must be created inside build folder. If you are in build folder then copy `mhm.exe` to upper folder. `cp ./mhm.exe ..`
- 6) Change directory to upper level and then call `mhm cd .. ./mhm`

## Ubuntu, Mint and other apt-get based systems with matching repositories

```
sudo apt-get install git # (optional)
sudo apt-get install gfortran netcdf-bin libnetcdf-dev libnetcdf-dev cmake
```



## Archlinux

```
sudo pacman -S git # (optional)
sudo pacman -S gcc-libs netcdf-fortran cmake
```

## Module systems

If you are on a module system, load the modules gcc or intel depending on your favorite compiler. Then, load the modules netcdf-fortran and cmake.

These modules will have system specific names, environments, etc. You may use `module spider` to find the right packages and the right dependencies, potentially use corresponding wiki pages.

### On eve (the cluster at the UFZ)

A set of load-scripts is provided in `moduleLoadScripts`, to load all need modules for specific compilers:

- GNU 7.3 compiler (`foss/2018b Toolchain`):  
`source moduleLoadScripts/eve.gcc73`  
 or (MPI support)  
`source moduleLoadScripts/eve.gcc73MPI`
- GNU 8.3 compiler (`foss/2019b Toolchain`):  
`source moduleLoadScripts/eve.gcc83`  
 or (MPI support)  
`source moduleLoadScripts/eve.gcc83MPI`
- Intel 18 compiler (`iomkl/2018b Toolchain`):  
`source moduleLoadScripts/eve.intel18`  
 or (MPI support)  
`source moduleLoadScripts/eve.intel18MPI`
- Intel 20 compiler (`iomkl/2020a Toolchain`):  
`source moduleLoadScripts/eve.intel20`  
 or (MPI support)  
`source moduleLoadScripts/eve.intel20MPI`
- NAG 6.2 compiler:  
`source moduleLoadScripts/eve.nag62`

Then you can compile mHM with cmake. We prepared a set of scripts, to automatize the build and compilation process to generate an executable in the root directory with the following naming scheme:

- Release version mhm:  
`source CI-scripts/compile`
- Debug version mhm\_debug:  
`source CI-scripts/compile_debug`
- Release version with MPI support mhm\_mpi:  
`source CI-scripts/compile_MPI`
- Debug version with MPI support mhm\_mpi\_debug:  
`source CI-scripts/compile_MPI_debug`
- Release version with OpenMP support mhm\_openmp:  
`source CI-scripts/compile_OpenMP`
- Debug version with OpenMP support mhm\_openmp\_debug:  
`source CI-scripts/compile_OpenMP_debug`

## MacOS

\*(to be added)\*

## 11.3 Specific setups

The following hints can replace the step `cmake ..` in the installation instruction.

You can skip this part and continue with "Installation", if you do not have a module system setup (like on clusters) or if you have not installed all packages with a package manager, such as `cygwin` or `apt-get`.

### Module systems

If you are okay with loading the needed modules as described in the section 'Module systems' above before executing the program you can skip this step and continue with 'Installation'.

The executable can be built in such a way that it does not require any modules to be loaded. The module system, though, adds system paths in the background the user should not care about too much, so the setup is a workaround. (This would be the case with any other building tool aswell.) It should be stable, anyway.

In case you want to have a module-independent build, instead of just executing `cmake ..`, either run

```
cmake -DCMAKE_BUILD_MODULE_SYSTEM_INDEPENDENT=ON ..
```

or

```
cmake -C ../CMakeCacheFiles/eve ..
```

or change the variable `CMAKE_BUILD_MODULE_SYSTEM_INDEPENDENT` with `ccmake` to `ON` after running `cmake ...`

### Non-standard locations for the netcdf-library (e.g. standard setup Macs in CHS):

Find the location of the `nf-config` file, for example by using:

```
find / -iname "*nf-config*" 2>/dev/null
```

This searches the root directory `/` for a file with a name containing the string "nf-config" (case-insensitive). It writes error messages like "permission denied" into the void.

Then, instead of running `cmake ..` if not using the standard compiler, set the fortran compiler variable to the desired compiler, e.g.

```
export FC=gfortran
```

then either run

```
cmake -DCMAKE_NETCDF_DIR=/path/to/nf-config/of/used/compiler
```

or copy the file `specificSetup` to some other place:

```
cp ../CMakeCacheFiles/specificSetup .
```

However, in case you want to keep it, you should choose a place outside the build repository. Edit the file as follows: add the path to your `nf-config` file, and after editing, run:

```
cmake -C specificSetup ..
```

or change the variable `CMAKE_NETCDF_DIR` to the path to the `nf-config` file with `ccmake` after running `cmake ...`

## 11.4 Installation

1. Change to a directory where you want to store the source code.
2. Clone the corresponding mHM repository into a folder, either using Git (if installed):

```
git clone --recurse-submodules https://git.ufz.de/mhm/mhm.git mhm/
```

for cloning it into a folder `mhm`, or download and unpack it using the download link on <https://git.ufz.de/mhm/mhm> (the cloud symbol with the arrow on it).

3. Create and change to a build directory where you want to store the build, e.g. inside the Git source directory

```
cd mhm
mkdir build
```

Change into the build directory:

```
cd build
```

4. Generate a system-dependent makefile

Execute `cmake` with the path to the Git source directory as parameter. (If you followed the instructions above, the path is `..` )

```
cmake ..
```

If everything worked well a Makefile was created with the corresponding paths.

*Note: have a look at "Specific setups" above in case you are using module systems, or when the netcdf libraries are not located where the package manager usually installs libraries, or when they are not saved in environment variables (i.e., classical MacOS setups at CHS).*

5. Make the build:

Execute make:

```
make
```

If this also worked fine, an executable was created, which has to be moved or copied to the Git source directory.

6. Execute the file:

```
cd ..
cp build/mhm .
```

On Windows the executable is called `mhm.exe` instead of `mhm`. In that case instead of `cp build/mhm .` execute

```
cp build/mhm.exe .
```

Now you might execute mHM:

```
./mhm
```

*Note concerning the development of the cmake setup: one could automatically link the executable with the cmake code inside the Git source directory which is not done for two reasons:*

- *The executable depends on your local system, so it should never be committed and pushed to other users. Nothing should be built inside the source directory automatically.*
- *The directory where mHM is executed usually is not the source directory but the directory where you want to run your tests. In case of the test setup it is the same, usually it is not.*

## 11.5 Building Release or Debug versions

If you want to set up specific versions of the build, you can create different folders for that. In case a release and a debug version need to be set up, it makes sense to create two folders named `debug` and `release`.

```
mkdir release
```

```
mkdir debug
```

inside the `release` folder one would execute

```
cmake -DCMAKE_BUILD_TYPE=Release ..
```

and inside the `debug` folder

```
cmake -DCMAKE_BUILD_TYPE=Debug ..
```

Executing

```
make
```

in the corresponding folder would then always result in a release build or respectively in a debug build.

## 11.6 Troubleshooting

**cmake** is far from being my main task, so it will probably take a while until I can track a problem. Nonetheless, I would be happy having bug reports. Feel free to write an email to [maren.kaluza@ufz.de](mailto:maren.kaluza@ufz.de) if there are any bugs, remarks or questions.

### MacOS, homebrew

On brew/homebrew setup MacOS systems a working `nf-config` is not yet available at this time. Execute:

```
nf-config --all
```

and if it says something like "is not implemented yet" the issue is not solved yet. But it is on my tracklist.

### cygwin

If libraries are not found, the problem can be:

- you accidentally tried to use commands within the `cmd` shell and not within the `cygwin` shell
- your `cygwin` setup might be broken. Try deinstalling, following the instructions <https://cygwin.com/faq/faq.html#faq.setup.uninstall-service>, and reinstalling again. During reinstallation only install the required dependencies.

## Chapter 12

### Publications using mHM

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## Chapter 13

# mHM Release Notes

### 13.1 mHM v5.11.2 (Jul 2021)

#### Enhancements

- documentation modernized with `doxygen-awesome-css` (!86)
- cmake update to be able to install mHM (`cmake --install`) (!85)
- added pFUnit tests thanks to Nicola Döring (!76)
- link to a new [YouTube tutorial](#) for compiling mHM with cygwin by Mehmet Cüneyd Demirel added to the documentation (!74)
- NetCDF output: add deflate and precision option to namelists (!73)
- refactor cmake workflow (!72)

#### Bugfixes

- fixed: `mrm` tried to write output even if routing was switched off (!82)
- unreachable `else` branch in `feddes_et_reduction` removed (!77)
- unnecessary `inout` variable intent in `soil_moisture` removed (!77)

### 13.2 mHM v5.11.1 (Mar 2021)

#### Enhancements

- added compile information for cygwin (!68)

#### Bugfixes

- removed note about mHM 5.10 from the README
- `smhorizon`: `tmp_rootfraccoef` was corrected directly if it is not between 0 and 1, but actually `FCnorm` should always be between 0 and 1 (!67)

## 13.3 mHM v5.11 (Feb 2021)

### Experimental Features

- river temperature routing was implemented in an alpha version 0.1 (!37)
  - this feature is in an experimental stage and should not be considered stable!

### Enhancements

- introduced central version files `version.txt` and `version_data.txt` (!51)
- added Feddes and global FC dependency on root fraction coef. at SM process(3)=4 (!43)
- Online documentation generated with doxygen: <https://mhm.pages.ufz.de/mhm/develop/> (!44)
- CI/CD with GitLab Runner (!11, !13, !14, !28, !32, !48, !50)
  - building on EVE for multiple compiler (GNU 7.3/8.3, Intel 18/19, NAG 6.2)
  - building debug/release serial/parallel
  - memcheck with valgrind
  - running all check-cases with all compiled versions
  - calculation of coverage
  - new checking script `run_mhm_checks.py`
- the domain loop is now parallelized with MPI
- objective function for boxcox-transformed streamflow
- post processing script for probabilistic forecasts
- different module load scripts for EVE
- Objective function from separate mhm calls (!7)
- new data type for simulated gridded optidata (!10)
- new datatype `datetimeinfo` (!16)
- added module `mo_os` to check files and directories (!41, !57)

### Changes

- internal: "basin" renamed to "domain"
- TWS input file changed from ascii to netCDF (!9)
- Switched to cell wise kge of et and tws in `opti_function` 33 (!12)
- restart files are now given by name (!34)
- removed mRM standalone and statically integrate mRM into mHM (!53)
- removed the old makefile and legacy checking scripts (!55)
- minimal Cmake version is now 3.12 (!58)

## Bugfixes

- Finalparam.nml is now written with specific format (Intel/GNU compatibility) (#40)
- FinalParam.nml routing section bug fixed (#49, !25)
- dirEvapotranspiration is now allocated before writing
- cmake: netcdf link flags where separated by ";"
- sharing of L0 domain now working
- added L1\_jarvis\_thresh\_c1 to restart file for process id 2 AND 3 (#29, !15)
- allowing higher routing resolution than hydrology (!21)
- domainID not set correctly for mRM if restart is activated (!30)
- mHM states\_fluxes netCDF output was curvilinear even if coordinate system is set to regular latlon (#98, !31)
- mismatch in messages about written mhm fluxes (!42)
- Fixing wrongly matched IDs from L1 to L11 when routing resolution (L11) is finer than L1 resolution (!45)
- The length in net\_startup was only cut in case there are less then 2 lengths (!46)
- corrected unit attributes for lat lon variables (!47)
- Allow run mHM and mRM without any observed gauge for processCase(8) = 2 / 3 (#27, !52)

## 13.4 mHM v5.10 (June 2019)

### New Features:

- New routing process introduced `processCase(8) = 3` [see Thober et al, 2019, GMDD, in press](#) for more details
- mRM is decoupled from mHM and mRM now resides in `deps/mrm` as an independent submodule [more information on handling submodules](#)
- New option to compile mHM with cmake is provided, see more details under [cmake manual](#).
- Visualization/animation R script (producing PDF and GIF) of mHM netcdf files included under `post-proc animatel.R`
- New option for coupling of mRM to a groundwater model (`gw_coupling = .true.`). The river head can be computed based on the Manning equation.

### Bugs resolved from release 5.9:

- Enable use of i8 for `time_data` in `common/mo_read_forcing_nc.f90`, otherwise netcdf time stamps with initial dates prior to 1900 were wrong (due to overflow)

**Known bugs:**

1. Adaptive routing does not allow to run without at least 1 gauge specification
2. Incompatibility of Finalparam.nml format between Intel and GNU
3. If ProcessOption(3) is set to 3 and optimization is activated, the created FinalParam.nml misses the header for the namelist of the soil moisture parameters.
4. Land cover scenes cannot be changed between the run generating the restart file and the run using the restart file. This causes unpredictable behaviour by the model.
5. Simulation period must span overall land cover scenes specified in the namelist.
6. Cut-off for link length is calculated with missing values, but those should be neglected.
7. Using a higher routing resolution than hydrology resolution may cause segmentation faults because mapping from L1id on L11 is not working correctly
8. If ProcessOption(5) is set to -1 and optimization is activated, the created FinalParam.nml misses the header for the namelist of the PET process.

**Restrictions:**

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).

**13.5 mHM v5.9 (July 2018)****New Features:**

- Major restructuralization of the mHM code.
- MPR is now executed before mHM is run.
- MPR can be compiled as a standalone tool.
- mHM without mRM can be compiled.
- The code in general is strictly reorganized into modules that belong to their respective processes (MPR, mHM, mRM). This is not only done for constants, global variables and so on, but also for every module and the reading of input as well as the namelists.
- This leads to the creation of those folders:
  - `./common` (code shared by MPR, mHM and mRM) included for every compilation option
  - `./lib` (code shared by MPR, mHM and mRM) included for every compilation option
  - `./MPR` (code for MPR), not included for mRM standalone
  - `./common_mHM_mRM` (code shared by mHM and mRM), not included for MPR standalone
  - `./mHM` (code for mHM), not included for MPR and mRM standalone
  - `./mRM` (code for mRM), not included for MPR standalone and mHM without routing
- Code is reformatted (indentation=2, spacing unified)
- Removed many duplicate code parts (e.g. shared between mHM and mRM)
- Check cases are minimized (reduced output, shorter time periods ( $\leq 2$  years), less basins)

- Check cases can now be set up more easily (by use of `model_wrapper` for automatic creation of new `nml` files)
- Check cases now run a python script for output comparison, advantage: tolerance now also allowed for `ascii-output` and support of 4-D `netCDF` files without time dimension
- `fSealed` is now an effective parameter
- mHM effective parameters now all have three dimensions internally (`nCells_L1`, [`iHorizon`, `LAI-Time`], `nLCoverScenes`)
- Introduction of new derived types:
  - `Grid` (merge of `basin_info`, `basin_info_mrm`, `gridGeoRef`, `nCells`, `longitude`, `latitude`, `Id`) used for each level (0, 1, 11, 2) individually
  - `GridRemapper` (merge of `lower_bound`, `upper_bound`, etc.)
- `mhm_eval` and `mrm_eval` now have common procedure interface (needed for fully flexible optimisation)
- A new post-processor can check and adapt some fields for doxygen generation
- Changes that are not backwards-compatible:
  - Restart files are restructured (now contain only the minimum required for restart):
  - MPR: effective parameters at L1 + grid information
  - mHM: effective parameters, states and fluxes at L1 + grid information
  - mRM: routing-specific parameters, states, fluxes, configs at L1/L11 + grid information
  - `mhm.nml` is restructured and not backwards compatible mainly due to the reorganization of modules in dependency of their processes
  - gridded LAI values are now used for all effective parameters (no fallback to LAIclasses anymore)
  - canopy height used for aerodynamic resistance is now scaled with the actual LAI timeseries and not with a dummy timeseries of intensive orchard
- Considerable improvements and reduced redundancy in estimation of an empirical distribution of slopes (sort function in the [mo\\_startup](#)). Example for the Australian domain (180 million L0 cells), it reduced time for sorting slope from 32hours to only 1 minute.
- `mtCLIM` preprocessor in `pre-proc/mtCLIM`, based on [mtCLIM v4.3](#). This code is able to estimate humidity (vapore pressure or vapore pressure deficit) and incoming shortwave radiation based on meteorological variables (minimum and maximum air temperature, precipitation) and morphological characteristics of the underlying terrain (digital elevation model, slope, aspect).
- `mHM2OGS` preprocessor in `pre-proc/GIS2FEM3`. This code converts the triangular-wise or quadrilateral-wise recharge data from mHM into the nodal source terms of a three dimensional finite element model for [OGS](#).
- Updated documentation for CYGWIN installations under Windows 7 and 10.
- Added new objective function number 31: weighted NSE (NSE is weighted with observed discharge)
- Removal of `bin` files and related code (only `nc` and `ascii` files are used)

### Bugs resolved from release 5.8:

- Enable use of `i8` for `time_data` in `common/mo_read_forcing_nc.f90`, otherwise `netcdf` time stamps with initial dates prior to 1900 were wrong (due to overflow)

### Known bugs:

None.

**Restrictions:**

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).
- If you wish to use the multi-scale Routing Model as stand-alone version, please contact [Stephan Thober](#).

## 13.6 mHM v5.8 (Dec 2017)

**New Features:**

- Implementation of a new process for PET correction based on LAI at PET `process(5)=-1` (Cuneyd Demirel + *GEUS* colleagues);
- Pre-processor code for SOILGRIDS data as used for the EDgE project (Rohini Kumar);
- Reduced computational time of the neutron forward model COSMIC by factors of 30–100 (Maren Kaluza);
- Compression of the netCDF output files (David Schaefer);
- Optional project description added into the `mhm.nml`

**Bugs resolved from release 5.7:**

- `processCase(3)=3` did not work when compiled with openMP.
- openMP declarations missing in `mo_mpr_smhorizons.f90` for the case of `iFlag_soil=1`

**Known bugs:**

None.

**Restrictions:**

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use a special process description of evapotranspiration (`processCase(4)`) please contact [Matthias Zink](#).
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).
- If you wish to use the multi-scale Routing Model as stand-alone version, please contact [Stephan Thober](#).

## 13.7 mHM v5.7 (Jun 2017)

**New Features:**

- New process descriptions for the soil evapotranspiration module:
  - Field capacity dependency to root fraction coefficient (`processCase(3)=2`) – implemented by *GEUS*.

- Jarvis (1989, J. Hydrol.) evapotranspiration reduction (`processCase(3)=3`) – implemented by *GEUS*.
- Use local, monthly LAI climatology instead of look-up table, i.e., `LAI_classdefinition.txt` (`timeStep_LAI←_input=1`).
- New objective functions for model calibration
  - Calibration of mHM using catchment average evapotranspiration (`opti_function=27`).
  - Calibration of mHM using soil moisture and streamflow simultaneously (`opti_function=28`).

### Bugs resolved:

- Calibration using `processCase(8)=2` (routing with adaptive timestep) does properly work now.
- Streamflow output is now properly written to the NetCDF.

### Known bugs:

None

### Restrictions:

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use a special process description of evapotranspiration (`processCase(4)`) please contact [Matthias Zink](#).
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).
- If you wish to use the multi-scale Routing Model as stand-alone version, please contact [Stephan Thober](#).

## 13.8 mHM v5.6 (Dec 2016)

### New Features:

- **Routing extended:** Implementation of a new parametrization for the routing model (`processCase(8)=2`). This routing option is based on an adaptive time step to improve the scalability and transferability of the model as well as a significant reduction in run time. The adaptive time step is calculated as ratio of routing resolution and celerity, the latter can be given as parameter in `mhm_parameter.nml`.

### Bugs resolved:

- Any model time step from 1 h to 24 h can be chosen (in releases v5.4 and v5.5 only 1 h worked properly).
- Estimation of the Hargreaves-Samani PET for high altitudes works properly now (there have been numerical issues for high latitude values).
- Reading catchment outlets from the `restart` file works now (bug appeared in v5.5).

### Known bugs:

- Calibration using `processCase(8)=2` (adaptive timestep) does not work, please use `processCase(8)=1`.

**Restrictions:**

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use a special process description of evapotranspiration (`processCase(4)`) please contact [Matthias Zink](#).
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).
- If you wish to use the multi-scale Routing Model as stand-alone version, please contact [Stephan Thober](#).

## 13.9 mHM v5.5 (Jun 2016)

**New Features:**

- Routing works on domains with multiple outlets (e.g., continental level).
- New option for providing soil data. They can be provided as predefined layers (one map per layer).
- Speed up of mHM for big domains, due to reformulations in the model start up.
- Pre-processing: new tools for i) cutting out a catchment from a existing dataset, ii) estimation of Hargreaves-Samani evapotranspiration, and iii) enlarging the grids of the input files.

**Bugs resolved:**

- Assigning routing parameters is done properly now.

**Known bugs:**

- Specifying a model time step of 24h (in `mhm.nml`) does not work, please stick with the default time step (1h)

**Restrictions:**

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use a special process description of evapotranspiration (`processCase(4)`) please contact [Matthias Zink](#).
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).
- If you wish to use the multi-scale Routing Model as stand-alone version, please contact [Stephan Thober](#).

## 13.10 mHM v5.4 (Dec 2015)

**New Features:**

- The routing of mHM can be used as a stand alone version/independent software called *multiscale Routing Model mRM*, e.g. for coupling to other environmental models.



- A new output, i.e. fields of routed discharge, is now available. They are stored in `mRM_fluxes_and_states.nc` and are controlled by a new namelist contained in the `mrm_outputs.nml` file, which is an optional file.
- New calibration objectives have been incorporated. It is now possible, additionally to the former objectives, to calibrate mHM against additional input data:
  - total water storage (e.g. GRACE) and discharge simultaneously (`opti_function=15`), and/or
  - cosmic ray neutron counts (`opti_function=17`).
- New post-processing: a mHM python class for reading all inputs and outputs of a model run can be found in `post-proc/`.
- Reorganization of the NetCDF writing in mHM to simplify future implementations of additional outputs.

### Bugs resolved:

- Calibration with catchment average soil moisture (`opti_function=10`) works properly now.
- Discharge output for multi basin runs with different time periods for each basin works properly now.
- Hargreaves-Samani PET calculation (`processCase(5)=1`) is valid on southern hemisphere too now.

### Known bugs:

- None.

### Restrictions:

- For `gfortran` compilers mHM supports only v4.8 and higher.
- If you wish to use a special process description of evapotranspiration (`processCase(4)`) please contact [Matthias Zink](#).
- If you wish to use features connected to ground albedo neutrons (`processCase(9)`), please contact [Martin Schrön](#).
- If you wish to use the multi-scale Routing Model as stand-alone version, please contact [Stephan Thober](#).

## 13.11 mHM v5.3 (Jun 2015)

### New Features:

- Simulation period and warming days can be now given per basin (see `time_periods` in `mhm.nml`)
- Enabling use of MPI (set `mpi=true` in `[Makefile](Makefile)` and use `#ifdef MPI` for MPI specific code)
- Optional input data can be loaded for example to calibrate against soil moisture (see `optional_data` in `mhm.nml`)
- Generation of ground albedo cosmic-ray neutrons (see `processCase(10)` in `mhm.nml`); these calculations are based on the `COSMIC` code, which was originally written by Rafael Rosolem. Please contact [Martin Schrön](#) if you like to use this new feature.
- Several new objective functions, e.g. calibrating the Kling-Gupta efficiency of catchment's average soil moisture (`opti_function=10`) or calibrating multiple basins regarding Kling-Gupta efficiency of discharge (`opti_function=14`) among others; calibration against soil moisture is still purpose to research (`opti_function=10-14`). The interested user may contact [Matthias Zink](#) for further details.

**Bugs resolved:**

- Calibration using potential evapotranspiration from input file (i.e. `processCase(4)=0`) is now working properly

**Known bugs:**

- Compiling mHM with the recent Cygwin version under Windows is leading to an error message indicating circular dependencies. The reason for this is Unicode characters in some source code files. Please contact [mhm-admin@ufz.de](mailto:mhm-admin@ufz.de), if you get this error message. We will provide you the files with cleaned characters..

**Restrictions:**

- If you wish to use a special process description of evapotranspiration (i.e. Hargreaves-samani, Priestley-Taylor, or Penman-Monteith) please contact [Matthias Zink](#). The special cases are set in `mhm.nml` (see `processCase(4)`).
- If you wish to use the new feature of calculating neutron counts please contact [Martin Schrön](#). The feature can be enabled in `mhm.nml` (see `processCase(8)`).

**13.12 mHM 5.2 (Dec 2014)****New Features:**

- Chunk-wise reading of input data (see `timestep_model_inputs`)
- Complete revision of writing netCDF files
- Possibility to discard multi-scale parameter regionalization (MPR) calculations (see `perform_mpr`)
- Several process descriptions of evapotranspiration implemented (see `processCase(4)`): Read PET, Hargreaves-Samani, Priestley-Taylor, Penman-Monteith. Please contact [Matthias Zink](#) if you use one of the last three options, since the code is not under GNU Public license up to now.
- Adding routines for signature calculations of time series (see `mo_signatures`)
- New objective function for calibrating discharge with Kling-Gupta efficiency measure (KGE, see `opti_function`)
- New output variables (see `mhm_outputs.nml`)
- Sorting algorithm changed to public available library orderpack (see `mo_orderpack`)

**Bugs resolved:**

- Some variables in restart file where not assigned correctly
- Variables not initialized correctly

**Known bugs:**

- Calibration using PET values read from the input file (`processCase(5)=0`) is running, but yields wrong results due to a wrong initialization of variables. **The bug is resolved and will be released with version 5.3.**

## 13.13 mHM 5.1 (Jun 2014)

### New Features:

- OpenMP handling of routines such as the multi-scale parameter regionalization (MPR)
- Multi-scale implementation, i.e. running mHM simultaneously in several basins with different resolutions
- Automatic check case framework, i.e. testing new implementations on their validity and back-compatibility
- Implementation of inflow gauges, i.e. feeding discharge time series from upstream areas at catchment boundaries
- File `gaugeinfo.txt` specifying gauging stations is now part of namelist `mhm.nml`
- Code is now free of Numerical Recipes proprietary code
- Can now run on a single cell (no routing performed) Hydrological modelling resolution (L1) equal to morphological input data resolution (L0) possible
- Windows compatible (with Cygwin)
- Support of regular geographic coordinate systems (e.g lat-lon) in addition to equal-area coordinate systems (UTM)

### Bugs resolved:

- Initialization of states was not correct when running mHM in calibration mode.
- Calculated parameter values ( `mhm_parameters.nml`) not necessarily in bound (check added).
- Aggregation/Disaggregation of meteorological data corrected.
- Forecast with mHM did not work because modelling period was restricted to discharge data period.
- Wrong mapping of evaluation discharge gauges for runs involving multiple gauges.

### Known bugs:

- Print out of River network in Config File is wrong for Multi-Basin setup, i.e., the River network is always properly written for the first basin, but not properly for subsequent basins when these are either different ones or the same one with a different Hydrology or Routing resolution.
- mHM does not abort if x-axis of L0 (morphological data) and L2 (meteorological data) do not span over exactly the same range.

## 13.14 mHM 5.0 (Dec 2013)

### New Features:

- Full modular version
- Automatic documentation by doxygen
- Running mHM for multiple basin simultaneously
- Definition of 8 major processes:
  - interception,
  - snow,

- soil moisture,
  - direct runoff,
  - evapotranspiration,
  - interflow,
  - percolation,
  - routing
- Choice of different descriptions of processes possible
  - Input in binary `*.bin` or netcdf `*.nc` format
  - Various calibration routines and objective functions
  - Consistent numerical precision handling of variables

**Known bugs:**

- None.

# Chapter 14

## Modules Index

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# Chapter 16

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# Chapter 17

## Module Documentation

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### 17.2 mo\_canopy\_interc Module Reference

Canopy interception.

#### Functions/Subroutines

- elemental pure subroutine, public [canopy\\_interc](#) (pet, interc\_max, precip, interc, throughfall, evap\_canopy)  
*Canopy interception.*

#### 17.2.1 Detailed Description

Canopy interception.

This module deals with processes related to canopy interception, evaporation and throughfall.

#### Authors

Vladyslav Prykhodko

#### Date

Dec 2012

#### 17.2.2 Function/Subroutine Documentation

##### 17.2.2.1 canopy\_interc()

```
elemental pure subroutine, public mo_canopy_interc::canopy_interc (  
    real(dp), intent(in) pet,  
    real(dp), intent(in) interc_max,  
    real(dp), intent(in) precip,  
    real(dp), intent(inout) interc,  
    real(dp), intent(out) throughfall,  
    real(dp), intent(out) evap_canopy )
```

Canopy interception.

Calculates throughfall. Updates interception and evaporation intensity from canopy. Throughfall ( $F$ ) is estimated as a function of the incoming precipitation ( $P$ ), the current status of the canopy water content ( $C$ ), and the max. water

$$F = \text{Max}((P + C - C_{max}), 0)$$

Evaporation ( $E$ ) from canopy is estimated as a fraction of the potential evapotranspiration ( $E_p$ ) depending on the current status of the canopy water content ( $C$ ) and the max. water content ( $C_{max}$ ) that can be intercepted by the vegetation.

$$E = E_p (C/C_{max})^{2/3}$$

ADDITIONAL INFORMATION content ( $C_{max}$ ) that can be intercepted by the vegetation. canopy\_interc(pet, interc\_← month\_max, interc\_max, precip, throughfall, evap\_canopy, interc)

#### Parameters

|         |                         |  |
|---------|-------------------------|--|
| in      | REAL(dp) :: pet         | Potential evapotranspiration [mm TS-1]           |
| in      | REAL(dp) :: interc_max  | Maximum interception [mm]                        |
| in      | REAL(dp) :: precip      | Daily mean precipitation [mm]                    |
| in, out | REAL(dp) :: interc      | Interception [mm]                                |
| out     | REAL(dp) :: throughfall | Throughfall [mm TS-1]                            |
| out     | REAL(dp) :: evap_canopy | Real evaporation intensity from canopy [mm TS-1] |

#### Authors

Vladyslav Prykhodko

#### Date

Dec 2012

Definition at line 74 of file mo\_canopy\_interc.f90.

References mo\_common\_constants::eps\_dp.

Referenced by mo\_mhm::mhm().

Here is the caller graph for this function:



## 17.3 mo\_check Module Reference

### Functions/Subroutines

- subroutine, public [check\\_dir](#) (path, text\_, throwError\_, tab\_, text\_length\_)  
Check if a given directory exists.

#### 17.3.1 Function/Subroutine Documentation

### 17.3.1.1 check\_dir()

```
subroutine, public mo_check::check_dir (
    character(len=*), intent(in) path,
    character(len=*), intent(in), optional text_,
    logical, intent(in), optional throwError_,
    integer(i4), intent(in), optional tab_,
    integer(i4), intent(in), optional text_length_ )
```

Check if a given directory exists.

Check if a given directory exists and write out a message about it. Will also give potential information about prefixes given with the path

#### Authors

Sebastian Mueller

#### Date

Nov 2020

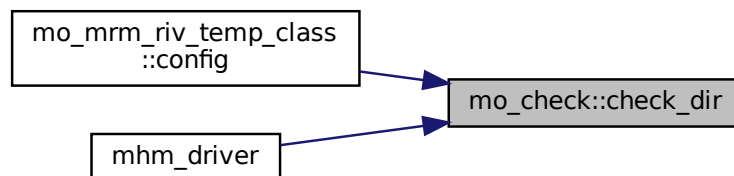
#### Parameters

|    |                                |   |
|----|--------------------------------|---|
| in | <i>path</i>                    | input path to check                               |
| in | <i>text_</i>                   | text to write out                                 |
| in | <i>throwerror_</i><br>_        | wheather to throw an error if folder not existing |
| in | <i>tab_</i>                    | tab-depth   |
| in | <i>text_</i><br><i>length_</i> | maximal text length (for aligning)                |

Definition at line 25 of file mo\_check.f90.

Referenced by mo\_mrm\_riv\_temp\_class::config(), and mhm\_driver().

Here is the caller graph for this function:



## 17.4 mo\_common\_constants Module Reference

Provides constants commonly used by mHM, mRM and MPR.

## Variables

- real(dp), parameter, public `eps_dp` = `epsilon(1.0_dp)`  
*epsilon(1.0) in double precision*
- real(sp), parameter, public `eps_sp` = `epsilon(1.0_sp)`  
*epsilon(1.0) in single precision*
- integer(i4), parameter, public `nodata_i4` = -9999\_i4
- real(dp), parameter, public `nodata_dp` = -9999.\_dp
- real(dp), parameter, public `p1_initstatefluxes` = 0.00\_dp
- integer(i4), parameter, public `ncolpars` = 5\_i4
- integer(i4), parameter, public `maxnodomains` = 50\_i4
- integer(i4), parameter, public `maxnlcovers` = 50\_i4
- character(64), parameter, public `soilhorizonsvarname` = "L1\_SoilHorizons"
- character(64), parameter, public `landcoverperiodsvarname` = "L1\_LandCoverPeriods"
- character(64), parameter, public `laivarname` = "L1\_LAITimesteps"

### 17.4.1 Detailed Description

Provides constants commonly used by mHM, mRM and MPR.

Provides commonly used by mHM, mRM and MPR such as no\_data values and eps

#### Authors

Robert Schweppe

#### Date

Dec 2017

### 17.4.2 Variable Documentation

#### 17.4.2.1 eps\_dp

```
real(dp), parameter, public mo_common_constants::eps_dp = epsilon(1.0_dp)
```

*epsilon(1.0) in double precision*

Definition at line 24 of file mo\_common\_constants.f90.

Referenced by mo\_multi\_param\_reg::aerodynamical\_resistance(), mo\_canopy\_interc::canopy\_interc(), mo\_mpr←  
\_startup::l0\_check\_input(), mo\_mpr\_read\_config::mpr\_read\_config(), mo\_objective\_function::objective\_kge\_q←  
rmse\_et(), mo\_objective\_function::objective\_kge\_q\_rmse\_tws(), mo\_soil\_database::read\_soil\_lut(), mo\_runoff←  
::runoff\_unsat\_zone(), and mo\_soil\_moisture::soil\_moisture().

#### 17.4.2.2 eps\_sp

```
real(sp), parameter, public mo_common_constants::eps_sp = epsilon(1.0_sp)
```

*epsilon(1.0) in single precision*

Definition at line 26 of file mo\_common\_constants.f90.



### 17.4.2.3 laivarname

```
character(64), parameter, public mo_common_constants::laivarname = "L1_LAItimeSteps"
```

Definition at line 42 of file mo\_common\_constants.f90.

Referenced by mo\_restart::read\_restart\_states(), and mo\_mpr\_restart::write\_mpr\_restart\_files().

### 17.4.2.4 landcoverperiodsvarname

```
character(64), parameter, public mo_common_constants::landcoverperiodsvarname = "L1_LandCover←  
Periods"
```

Definition at line 41 of file mo\_common\_constants.f90.

Referenced by mo\_mrm\_restart::mrm\_write\_restart(), mo\_restart::read\_restart\_states(), and mo\_mpr\_restart←  
::write\_mpr\_restart\_files().

### 17.4.2.5 maxnlcovers

```
integer(i4), parameter, public mo_common_constants::maxnlcovers = 50_i4
```

Definition at line 38 of file mo\_common\_constants.f90.

Referenced by mo\_common\_read\_config::common\_read\_config().

### 17.4.2.6 maxnodomains

```
integer(i4), parameter, public mo_common_constants::maxnodomains = 50_i4
```

Definition at line 37 of file mo\_common\_constants.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_read←  
\_config::common\_read\_config(), mo\_mrm\_riv\_temp\_class::config(), mo\_mhm\_read\_config::mhm\_read\_config(),  
mo\_mpr\_read\_config::mpr\_read\_config(), and mo\_mrm\_read\_config::mrm\_read\_config().

### 17.4.2.7 ncolpars

```
integer(i4), parameter, public mo_common_constants::ncolpars = 5_i4
```

Definition at line 36 of file mo\_common\_constants.f90.

Referenced by mo\_mpr\_read\_config::mpr\_read\_config(), and mo\_mrm\_read\_config::read\_mrm\_routing\_params().

### 17.4.2.8 nodata\_dp

```
real(dp), parameter, public mo_common_constants::nodata_dp = -9999._dp
```

Definition at line 30 of file mo\_common\_constants.f90.

Referenced by mo\_mrm\_river\_head::avg\_and\_write\_timestep(), mo\_multi\_param\_reg::baseflow\_param(), mo\_←  
meteo\_forcings::chunk\_config(), mo\_objective\_function::create\_domain\_avg\_et(), mo\_objective\_function::create←  
\_domain\_avg\_tws(), mo\_mrm\_river\_head::create\_output(), mo\_write\_fluxes\_states::createoutputfile(), mo\_mrm←

`_write_fluxes_states::createoutputfile()`, `mo_soil_database::generate_soil_database()`, `mo_grid::init_lowres_level()`, `mo_mrm_river_head::init_masked_zeros_l0()`, `mo_multi_param_reg::karstic_layer()`, `mo_mrm_net_startup::l11_calc_celerity()`, `mo_mrm_pre_routing::l11_e_acc()`, `mo_mrm_net_startup::l11_flow_accumulation()`, `mo_mrm_net_startup::l11_fraction_sealed_floodplain()`, `mo_mrm_pre_routing::l11_meteo_acc()`, `mo_mrm_pre_routing::l11_runoff_acc()`, `mo_mrm_net_startup::l11_stream_features()`, `mo_meteo_forcings::meteo_forcings_wrapper()`, `mo_meteo_forcings::meteo_weights_wrapper()`, `mo_multi_param_reg::mpr()`, `mo_mpr_read_config::mpr_read_config()`, `mo_mpr_runoff::mpr_runoff()`, `mo_mpr_soilmoist::mpr_sm()`, `mo_mpr_smhorizons::mpr_smhorizons()`, `mo_mrm_init::mrm_init()`, `mo_mrm_read_data::mrm_read_discharge()`, `mo_mrm_restart::mrm_read_restart_config()`, `mo_mrm_read_data::mrm_read_total_runoff()`, `mo_mrm_restart::mrm_write_restart()`, `mo_objective_function::objective()`, `mo_objective_function::objective_et_kge_catchment_avg()`, `mo_objective_function::objective_kge_q_rmse_et()`, `mo_objective_function::objective_kge_q_rmse_tws()`, `mo_objective_function::objective_master()`, `mo_objective_function::objective_neutrons_kge_catchment_avg()`, `mo_objective_function::objective_q_et_tws_kge_catchment_avg()`, `mo_objective_function::objective_sm_kge_catchment_avg()`, `mo_objective_function::objective_sm_pd()`, `mo_objective_function::objective_subprocess()`, `mo_read_wrapper::read_data()`, `mo_common_read_data::read_dem()`, `mo_soil_database::read_soil_lut()`, `mo_read_optional_data::readoptdataobs()`, `mo_spatial_agg_disagg_forcing::spatial_aggregation::spatial_aggregation_3d()`, `mo_spatial_agg_disagg_forcing::spatial_aggregation::spatial_aggregation_4d()`, `mo_spatial_agg_disagg_forcing::spatial_disaggregation::spatial_disaggregation_3d()`, `mo_spatial_agg_disagg_forcing::spatial_disaggregation::spatial_disaggregation_4d()`, `mo_write_ascii::write_configfile()`, `mo_mrm_write::write_configfile()`, `mo_mrm_write::write_daily_obs_sim_discharge()`, `mo_mpr_restart::write_eff_params()`, `mo_common_restart::write_grid_info()`, `mo_restart::write_restart_files()`, `mo_write_fluxes_states::writevariableattributes()`, `mo_mrm_write_fluxes_states::writevariableattributes()`, `mo_write_fluxes_states::writevariabletimestep()`, and `mo_mrm_write_fluxes_states::writevariabletimestep()`.

#### 17.4.2.9 nodata\_i4

```
integer(i4), parameter, public mo_common_constants::nodata_i4 = -9999_i4
```

Definition at line 29 of file `mo_common_constants.f90`.

Referenced by `mo_mrm_river_head::calc_channel_elevation()`, `mo_mrm_riv_temp_class::config()`, `mo_soil_database::generate_soil_database()`, `mo_grid::init_lowres_level()`, `mo_multi_param_reg::karstic_layer()`, `mo_mrm_init::l0_check_input_routing()`, `mo_upscaling_operators::l0_fractionalcover_in_lx()`, `mo_mrm_net_startup::l11_calc_celerity()`, `mo_mrm_net_startup::l11_flow_accumulation()`, `mo_mrm_net_startup::l11_flow_direction()`, `mo_mrm_net_startup::l11_l1_mapping()`, `mo_mrm_net_startup::l11_link_location()`, `mo_mrm_net_startup::l11_routing_order()`, `mo_mrm_net_startup::l11_set_drain_outlet_gauges()`, `mo_mrm_net_startup::l11_set_network_topology()`, `mo_mrm_net_startup::l11_stream_features()`, `mo_mhm_read_config::mhm_read_config()`, `mo_multi_param_reg::mpr()`, `mo_mpr_runoff::mpr_runoff()`, `mo_mpr_soilmoist::mpr_sm()`, `mo_mrm_init::mrm_init()`, `mo_mrm_read_config::mrm_read_config()`, `mo_mrm_read_data::mrm_read_l0_data()`, `mo_mrm_restart::mrm_write_restart()`, `mo_read_wrapper::read_data()`, `mo_common_read_data::read_lcover()`, `mo_soil_database::read_soil_lut()`, `mo_mrm_read_data::rotate_fdir_variable()`, `mo_common_read_config::set_land_cover_scenes_id()`, `mo_mpr_restart::write_eff_params()`, and `mo_common_restart::write_grid_info()`.

#### 17.4.2.10 p1\_initstatefluxes

```
real(dp), parameter, public mo_common_constants::p1_initstatefluxes = 0.00_dp
```

Definition at line 33 of file `mo_common_constants.f90`.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_init_states::variables_alloc()`, `mo_init_states::variables_default_init()`, and `mo_mrm_init::variables_default_init_routing()`.

### 17.4.2.11 soilhorizonsvarname

character(64), parameter, public mo\_common\_constants::soilhorizonsvarname = "L1\_SoilHorizons"

Definition at line 40 of file mo\_common\_constants.f90.

Referenced by mo\_restart::read\_restart\_states(), and mo\_mpr\_restart::write\_mpr\_restart\_files().

## 17.5 mo\_common\_datetime\_type Module Reference

### Data Types

- type [datetimeinfo](#)

### Functions/Subroutines

- subroutine [datetimeinfo\\_init](#) (this, iDomain)
- subroutine [datetimeinfo\\_increment](#) (this)
- subroutine [datetimeinfo\\_update\\_lai\\_timestep](#) (this)
- logical function [datetimeinfo\\_writeout](#) (this, timeStep\_model\_outputs, tt)

### 17.5.1 Function/Subroutine Documentation

#### 17.5.1.1 datetimeinfo\_increment()

```
subroutine mo_common_datetime_type::datetimeinfo_increment (
    class(datetimeinfo), intent(inout) this )
```

Definition at line 105 of file mo\_common\_datetime\_type.f90.

References mo\_common\_mhm\_mrm\_variables::timestep.

#### 17.5.1.2 datetimeinfo\_init()

```
subroutine mo_common_datetime_type::datetimeinfo_init (
    class(datetimeinfo), intent(inout) this,
    integer(i4), intent(in) iDomain ) [private]
```

Definition at line 70 of file mo\_common\_datetime\_type.f90.

References mo\_common\_mhm\_mrm\_variables::lcyyearid, mo\_common\_mhm\_mrm\_variables::ntstepday, mo\_↔  
common\_mhm\_mrm\_variables::simper, and mo\_common\_mhm\_mrm\_variables::timestep.

#### 17.5.1.3 datetimeinfo\_update\_lai\_timestep()

```
subroutine mo_common_datetime_type::datetimeinfo_update_lai_timestep (
    class(datetimeinfo), intent(inout) this )
```

Definition at line 132 of file mo\_common\_datetime\_type.f90.

References mo\_mpr\_global\_variables::timestep\_lai\_input.

### 17.5.1.4 datetimeinfo\_writeout()

```
logical function mo_common_datetime_type::datetimeinfo_writeout (
    class(datetimeinfo), intent(in) this,
    integer(i4), intent(in) timeStep_model_outputs,
    integer(i4), intent(in) tt )
```

Definition at line 154 of file mo\_common\_datetime\_type.f90.

## 17.6 mo\_common\_file Module Reference

Provides file names and units for mRM.

### Variables

- character(len= \*), parameter `file_dem` = 'dem.asc'  
*DEM input data file.*
- integer, parameter `udem` = 53  
*Unit for DEM input data file.*
- integer, parameter `ulcoverclass` = 61  
*Unit for LCover input data file.*
- character(len= \*), parameter `file_config` = 'ConfigFile.log'  
*file defining mHM's outputs*
- integer, parameter `uconfig` = 68  
*Unit for file defining mHM's outputs.*

### 17.6.1 Detailed Description

Provides file names and units for mRM.

Provides all filenames as well as all units used for the multiscale Routing Model mRM.

#### Authors

Matthias Cuntz, Stephan Thober

#### Date

Aug 2015

### 17.6.2 Variable Documentation

#### 17.6.2.1 file\_config

```
character(len=*), parameter mo_common_file::file_config = 'ConfigFile.log'
```

file defining mHM's outputs

Definition at line 26 of file mo\_common\_file.f90.

Referenced by mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.6.2.2 file\_dem

```
character(len=*), parameter mo_common_file::file_dem = 'dem.asc'
```

DEM input data file.

Definition at line 19 of file mo\_common\_file.f90.

Referenced by mo\_common\_read\_data::read\_dem().

### 17.6.2.3 uconfig

```
integer, parameter mo_common_file::uconfig = 68
```

Unit for file defining mHM's outputs.

Definition at line 28 of file mo\_common\_file.f90.

Referenced by mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.6.2.4 udem

```
integer, parameter mo_common_file::udem = 53
```

Unit for DEM input data file.

Definition at line 21 of file mo\_common\_file.f90.

Referenced by mo\_common\_read\_data::read\_dem().

### 17.6.2.5 ulcoverclass

```
integer, parameter mo_common_file::ulcoverclass = 61
```

Unit for LCover input data file.

Definition at line 23 of file mo\_common\_file.f90.

Referenced by mo\_common\_read\_data::read\_lcover().

## 17.7 mo\_common\_functions Module Reference

Provides small utility functions used by multiple parts of the code (mHM, mRM, MPR)

### Functions/Subroutines

- logical function, public [in\\_bound](#) (params)

*TODO: add description.*

#### 17.7.1 Detailed Description

Provides small utility functions used by multiple parts of the code (mHM, mRM, MPR)

Provides the functions `in_bound` used to check `global_parameter` ranges

**Authors**

Robert Schweppe

**Date**

Dec 2017

**17.7.2 Function/Subroutine Documentation****17.7.2.1 in\_bound()**

```
logical function, public mo_common_functions::in_bound (
    real(dp), dimension(:, :) :: params )
```

TODO: add description.

TODO: add description

**Parameters**

|                 |  |   |
|-----------------|--|---|
| <code>in</code> | <code>real(dp), dimension(:, :) :: params</code> | parameter: col_1=Lower bound, col_2=Upper bound col_3=initial |
|-----------------|--|---|

**Authors**

Robert Schweppe

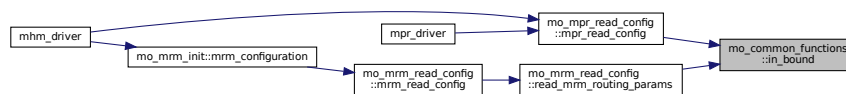
**Date**

Jun 2018

Definition at line 46 of file mo\_common\_functions.f90.

Referenced by mo\_mpr\_read\_config::mpr\_read\_config(), and mo\_mrm\_read\_config::read\_mrm\_routing\_params().

Here is the caller graph for this function:

**17.8 mo\_common\_mhm\_mrm\_file Module Reference**

Provides file names and units for mHM.

**Variables**

- character(len=\*), parameter `file_opti` = 'FinalParam.out'  
*file defining optimization outputs (objective and parameter set)*
- integer, parameter `uopti` = 72

*Unit for file optimization outputs (objective and parameter set)*

- character(len=\*), parameter `file_opti_nml` = 'FinalParam.nml'

*file defining optimization outputs in a namelist format (parameter set)*

- integer, parameter `uopti_nml` = 73

*Unit for file optimization outputs in a namelist format (parameter set)*

### 17.8.1 Detailed Description

Provides file names and units for mHM.

Provides all filenames as well as all units used for the hydrologic model mHM.

#### Authors

Matthias Cuntz

#### Date

Jan 2012

### 17.8.2 Variable Documentation

#### 17.8.2.1 file\_opti

```
character(len = *), parameter mo_common_mhm_mrm_file::file_opti = 'FinalParam.out'
```

file defining optimization outputs (objective and parameter set)

Definition at line 18 of file mo\_common\_mHM\_mRM\_file.f90.

Referenced by mo\_mrm\_write::mrm\_write\_optifile(), and mo\_write\_ascii::write\_optifile().

#### 17.8.2.2 file\_opti\_nml

```
character(len = *), parameter mo_common_mhm_mrm_file::file_opti_nml = 'FinalParam.nml'
```

file defining optimization outputs in a namelist format (parameter set)

Definition at line 22 of file mo\_common\_mHM\_mRM\_file.f90.

Referenced by mo\_mrm\_write::mrm\_write\_optinamelist(), and mo\_write\_ascii::write\_optinamelist().

#### 17.8.2.3 uopti

```
integer, parameter mo_common_mhm_mrm_file::uopti = 72
```

Unit for file optimization outputs (objective and parameter set)

Definition at line 20 of file mo\_common\_mHM\_mRM\_file.f90.

Referenced by mo\_mrm\_write::mrm\_write\_optifile(), and mo\_write\_ascii::write\_optifile().

### 17.8.2.4 uopti\_nml

integer, parameter mo\_common\_mhm\_mrm\_file::uopti\_nml = 73

Unit for file optimization outputs in a namelist format (parameter set)

Definition at line 24 of file mo\_common\_mHM\_mRM\_file.f90.

Referenced by mo\_mrm\_write::mrm\_write\_optinamelist(), and mo\_write\_ascii::write\_optinamelist().

## 17.9 mo\_common\_mhm\_mrm\_mpi\_tools Module Reference

tools for MPI communication that are mHM or mRM specific

### Functions/Subroutines

- subroutine, public [distribute\\_parameterset](#) (parameterset)
- subroutine, public [get\\_parameterset](#) (parameterset)

### 17.9.1 Detailed Description

tools for MPI communication that are mHM or mRM specific

This module contains sending and receiving subroutines for data that are specific for mHM or mRM

#### Authors

Maren Kaluza

#### Date

Jul 2019

### 17.9.2 Function/Subroutine Documentation

#### 17.9.2.1 distribute\_parameterset()

```
subroutine, public mo_common_mhm_mrm_mpi_tools::distribute_parameterset (
    real(dp), dimension(:), intent(in) parameterset )
```

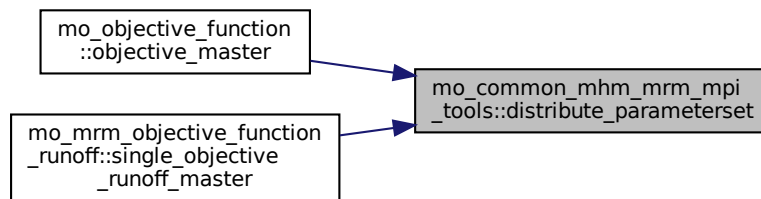
Definition at line 30 of file mo\_common\_mHM\_mRM\_MPI\_tools.f90.

References mo\_common\_variables::domainmeta.

Referenced by mo\_objective\_function::objective\_master(), and mo\_mrm\_objective\_function\_runoff::single\_↔objective\_runoff\_master().



Here is the caller graph for this function:



### 17.9.2.2 get\_parameterset()

```

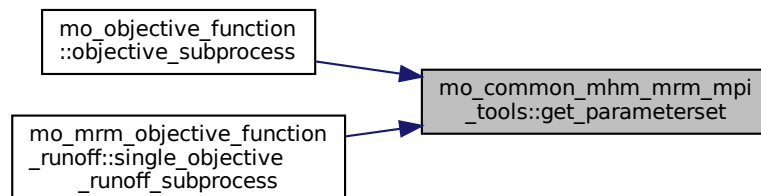
subroutine, public mo_common_mhm_mrm_mpi_tools::get_parameterset (
    real(dp), dimension(:), intent(inout), allocatable parameterset )
  
```

Definition at line 48 of file `mo_common_mHM_mRM_MPI_tools.f90`.

References `mo_common_variables::domainmeta`.

Referenced by `mo_objective_function::objective_subprocess()`, and `mo_mrm_objective_function_runoff::single_objective_runoff_subprocess()`.

Here is the caller graph for this function:



## 17.10 mo\_common\_mhm\_mrm\_read\_config Module Reference

Reading of main model configurations.

### Functions/Subroutines

- subroutine, public `common_mhm_mrm_read_config` (`file_namelist`, `unamelist`)  
*Read main configurations for common parts.*
- subroutine, public `check_optimization_settings`  
*TODO: add description.*
- subroutine, public `common_check_resolution` (`do_message`, `allow_subgrid_routing`)

*TODO: add description.*

- subroutine [period\\_copy\\_period\\_data](#) (toPeriod, fromPeriod)

### 17.10.1 Detailed Description

Reading of main model configurations.

This routine reads the configurations of common program parts

#### Authors

Matthias Zink

#### Date

Dec 2012

### 17.10.2 Function/Subroutine Documentation

#### 17.10.2.1 check\_optimization\_settings()

subroutine, public mo\_common\_mhm\_mrm\_read\_config::check\_optimization\_settings

*TODO: add description.*

*TODO: add description*

#### Authors

Robert Schweppe

#### Date

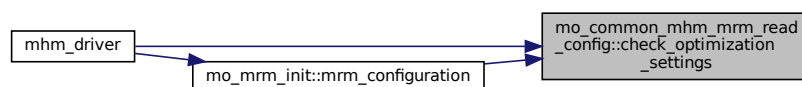
Jun 2018

Definition at line 233 of file mo\_common\_mHM\_mRM\_read\_config.f90.

References [mo\\_common\\_mhm\\_mrm\\_variables::dds\\_r](#), [mo\\_common\\_variables::global\\_parameters](#), [mo\\_common\\_mhm\\_mrm\\_variables::niterations](#), [mo\\_common\\_mhm\\_mrm\\_variables::sce\\_ngs](#), [mo\\_common\\_mhm\\_mrm\\_variables::sce\\_npg](#), and [mo\\_common\\_mhm\\_mrm\\_variables::sce\\_nps](#).

Referenced by [mhm\\_driver\(\)](#), and [mo\\_mrm\\_init::mrm\\_configuration\(\)](#).

Here is the caller graph for this function:



### 17.10.2.2 common\_check\_resolution()

```
subroutine, public mo_common_mhm_mrm_read_config::common_check_resolution (
    logical, intent(in) do_message,
    logical, intent(in) allow_subgrid_routing )
```

TODO: add description.

TODO: add description

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>logical :: do_message</i>            |  |
| in | <i>logical :: allow_subgrid_routing</i> |  |

#### Authors

Robert Schwappe

#### Date

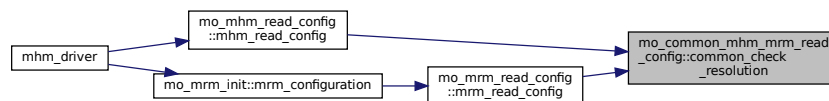
Jun 2018

Definition at line 294 of file mo\_common\_mHM\_mRM\_read\_config.f90.

References mo\_common\_variables::domainmeta, mo\_common\_variables::resolutionhydrology, and mo\_↔  
common\_mhm\_mrm\_variables::resolutionrouting.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config(), and mo\_mrm\_read\_config::mrm\_read\_config().

Here is the caller graph for this function:



### 17.10.2.3 common\_mhm\_mrm\_read\_config()

```
subroutine, public mo_common_mhm_mrm_read_config::common_mhm_mrm_read_config (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist )
```

Read main configurations for common parts.

TODO: add description

#### Parameters

|    |                                      |  |
|----|--------------------------------------|--|
| in | <i>character(*) :: file_namelist</i> |  |
| in | <i>integer :: unamelist</i>          |  |

**Authors**

Matthias Zink

**Date**

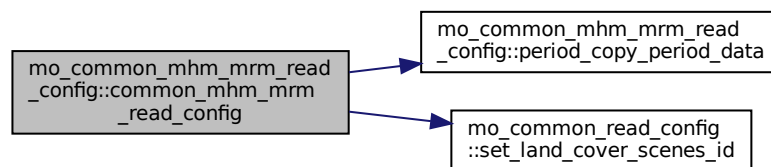
Dec 2012

Definition at line 49 of file mo\_common\_mHM\_mRM\_read\_config.f90.

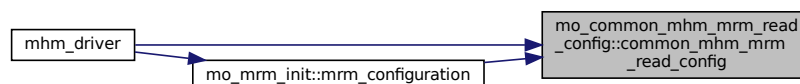
References mo\_common\_mhm\_mrm\_variables::dds\_r, mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_variables::evalper, mo\_common\_variables::lcfilename, mo\_common\_mhm\_mrm\_variables::lcyetid, mo\_common\_constants::maxnodomains, mo\_common\_mhm\_mrm\_variables::mcmc\_error\_params, mo\_common\_mhm\_mrm\_variables::mcmc\_opti, mo\_common\_mhm\_mrm\_variables::mhmfilerestartin, mo\_common\_mhm\_mrm\_variables::mrm\_read\_river\_network, mo\_common\_mhm\_mrm\_variables::mrmfilerestartin, mo\_common\_mhm\_mrm\_variables::niterations, mo\_common\_mhm\_mrm\_variables::ntstepday, mo\_common\_mhm\_mrm\_variables::opti\_function, mo\_common\_mhm\_mrm\_variables::opti\_method, mo\_common\_mhm\_mrm\_variables::optimize, mo\_common\_mhm\_mrm\_variables::optimize\_restart, period\_copy\_period\_data(), mo\_common\_variables::processmatrix, mo\_common\_mhm\_mrm\_variables::read\_restart, mo\_common\_mhm\_mrm\_variables::resolutionrouting, mo\_common\_mhm\_mrm\_variables::sa\_temp, mo\_common\_mhm\_mrm\_variables::sce\_ngs, mo\_common\_mhm\_mrm\_variables::sce\_npg, mo\_common\_mhm\_mrm\_variables::sce\_nps, mo\_common\_mhm\_mrm\_variables::seed, mo\_common\_read\_config::set\_land\_cover\_scenes\_id(), mo\_common\_mhm\_mrm\_variables::simper, mo\_common\_mhm\_mrm\_variables::timestep, mo\_common\_mhm\_mrm\_variables::warmingdays, and mo\_common\_mhm\_mrm\_variables::warmper.

Referenced by mhm\_driver(), and mo\_mrm\_init::mrm\_configuration().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.10.2.4 period\_copy\_period\_data()**

```

subroutine mo_common_mhm_mrm_read_config::period_copy_period_data (
    type(period), intent(inout) toPeriod,
    type(period), intent(in) fromPeriod )
  
```

Definition at line 354 of file mo\_common\_mHM\_mRM\_read\_config.f90.

Referenced by common\_mhm\_mrm\_read\_config().

Here is the caller graph for this function:



## 17.11 mo\_common\_mhm\_mrm\_restart Module Reference

TODO: add description.

### Data Types

- interface [check\\_consistency\\_element](#)

### Functions/Subroutines

- subroutine, public [check\\_dimension\\_consistency](#) (iBasin, nSoilHorizons\_temp, soilHorizonBoundaries\_temp, nLAIs\_temp, LAIBoundaries\_temp, nLandCoverPeriods\_temp, landCoverPeriodBoundaries\_temp)  
*checks dimension configurations read from restart file*
- subroutine [check\\_consistency\\_element\\_dp](#) (item1, item2, name, iBasin)
- subroutine [check\\_consistency\\_element\\_i4](#) (item1, item2, name, iBasin)

#### 17.11.1 Detailed Description

TODO: add description.

TODO: add description

##### Authors

Robert Schweppe

##### Date

Aug 2019

#### 17.11.2 Function/Subroutine Documentation

##### 17.11.2.1 check\_consistency\_element\_dp()

```
subroutine mo_common_mhm_mrm_restart::check_consistency_element_dp (
    real(dp), intent(in) item1,
    real(dp), intent(in) item2,
    character(*), intent(in) name,
    integer(i4), intent(in) iBasin )
```

Definition at line 90 of file mo\_common\_mHM\_mRM\_restart.f90.

### 17.11.2.2 check\_consistency\_element\_i4()

```
subroutine mo_common_mhm_mrm_restart::check_consistency_element_i4 (
    integer(i4), intent(in) item1,
    integer(i4), intent(in) item2,
    character(*), intent(in) name,
    integer(i4), intent(in) iBasin )
```

Definition at line 109 of file mo\_common\_mHM\_mRM\_restart.f90.

### 17.11.2.3 check\_dimension\_consistency()

```
subroutine, public mo_common_mhm_mrm_restart::check_dimension_consistency (
    integer(i4), intent(in) iBasin,
    integer(i4), intent(in) nSoilHorizons_temp,
    real(dp), dimension(:), intent(inout) soilHorizonBoundaries_temp,
    integer(i4), intent(in) nLAIs_temp,
    real(dp), dimension(:), intent(inout) LAIBoundaries_temp,
    integer(i4), intent(in) nLandCoverPeriods_temp,
    real(dp), dimension(:), intent(inout) landCoverPeriodBoundaries_temp )
```

checks dimension configurations read from restart file

#### Authors

Robert Schweppe

#### Date

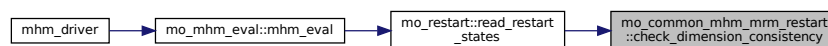
Aug 2019

Definition at line 46 of file mo\_common\_mHM\_mRM\_restart.f90.

References mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_mpr\_global\_variables::laiboundaries, mo\_↔  
common\_variables::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_mpr\_global\_variables::nlai, mo\_↔  
common\_variables::nlcoverscene, and mo\_mpr\_global\_variables::nsoilhorizons\_mhm.

Referenced by mo\_restart::read\_restart\_states().

Here is the caller graph for this function:



## 17.12 mo\_common\_mhm\_mrm\_variables Module Reference

Provides structures needed by mHM, mRM and/or mpr.

## Variables

- integer(i4) [mrm\\_coupling\\_mode](#)
- integer(i4), public [timestep](#)
- real(dp), public [c2tstu](#)
- real(dp), dimension(:), allocatable, public [resolutionrouting](#)
- logical, public [read\\_restart](#)
- logical, public [mrm\\_read\\_river\\_network](#)
- type([period](#)), dimension(:), allocatable, public [warmper](#)
- type([period](#)), dimension(:), allocatable, public [evalper](#)
- type([period](#)), dimension(:), allocatable, public [simper](#)
- type([period](#)), public [readper](#)
- integer(i4), dimension(:), allocatable, public [warmingdays](#)
- integer(i4), dimension(:, :), allocatable, public [lcyearid](#)
- integer(i4), public [ntstepday](#)
- character(256), dimension(:), allocatable, public [mhmfilerestartin](#)
- character(256), dimension(:), allocatable, public [mrmfilerestartin](#)
- integer(i4), public [opti\\_method](#)
- integer(i4), public [opti\\_function](#)
- logical, public [optimize](#)
- logical, public [optimize\\_restart](#)
- integer(i8), public [seed](#)
- integer(i4), public [niterations](#)
- real(dp), public [dds\\_r](#)
- real(dp), public [sa\\_temp](#)
- integer(i4), public [sce\\_ngs](#)
- integer(i4), public [sce\\_npg](#)
- integer(i4), public [sce\\_nps](#)
- logical, public [mcmc\\_opti](#)
- integer(i4), parameter, public [nerror\\_model](#) = 2
- real(dp), dimension([nerror\\_model](#)), public [mcmc\\_error\\_params](#)

### 17.12.1 Detailed Description

Provides structures needed by mHM, mRM and/or mpr.

Provides the global structure `period` that is used by both mHM and mRM.

#### Authors

Stephan Thober

#### Date

Sep 2015

### 17.12.2 Variable Documentation

#### 17.12.2.1 c2tstu

```
real(dp), public mo_common_mhm_mrm_variables::c2tstu
```

Definition at line 32 of file `mo_common_mHM_mRM_variables.f90`.

Referenced by `mo_startup::constants_init()`, and `mo_mhm_eval::mhm_eval()`.

### 17.12.2.2 dds\_r

```
real(dp), public mo_common_mhm_mrm_variables::dds_r
```

Definition at line 75 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::check\_optimization\_settings(), mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().

### 17.12.2.3 evalper

```
type(period), dimension(:), allocatable, public mo_common_mhm_mrm_variables::evalper
```

Definition at line 38 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mrm\_river\_head::create\_output(), mo\_write\_fluxes\_states::createoutputfile(), mo\_mrm\_write\_fluxes\_states::createoutputfile(), mo\_mrm\_objective\_function\_runoff::extract\_runoff(), mo\_mrm\_read\_data::mrm\_read\_discharge(), mo\_mrm\_write::mrm\_write(), mo\_mrm\_objective\_function\_runoff::multi\_objective\_ae\_fdc\_lsv\_nse\_djf(), mo\_objective\_function::objective\_kge\_q\_rmse\_et(), mo\_objective\_function::objective\_kge\_q\_rmse\_tws(), mo\_read\_optional\_data::readoptidataobs(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), and mo\_mrm\_write::write\_daily\_obs\_sim\_discharge().

### 17.12.2.4 lcyearid

```
integer(i4), dimension(:, :), allocatable, public mo_common_mhm_mrm_variables::lcyearid
```

Definition at line 42 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_datetime\_type::datetimeinfo\_init(), mo\_mhm\_eval::mhm\_eval(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.12.2.5 mcmc\_error\_params

```
real(dp), dimension(nerror_model), public mo_common_mhm_mrm_variables::mcmc_error_params
```

Definition at line 89 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().

### 17.12.2.6 mcmc\_opti

```
logical, public mo_common_mhm_mrm_variables::mcmc_opti
```

Definition at line 85 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().



### 17.12.2.7 mhmfilerestartin

```
character(256), dimension(:), allocatable, public mo_common_mhm_mrm_variables::mhmfilerestartin
```

Definition at line 53 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mhm\_eval::mhm←\_eval(), and mo\_startup::mhm\_initialize().

### 17.12.2.8 mrm\_coupling\_mode

```
integer(i4) mo_common_mhm_mrm_variables::mrm_coupling_mode
```

Definition at line 26 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_mrm\_init::mrm\_configuration(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_write::mrm\_write(), and mo\_mrm\_write::write\_configfile().

### 17.12.2.9 mrm\_read\_river\_network

```
logical, public mo_common_mhm_mrm_variables::mrm_read_river_network
```

Definition at line 35 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_mrm\_init←::mrm\_init().

### 17.12.2.10 mrmfilerestartin

```
character(256), dimension(:), allocatable, public mo_common_mhm_mrm_variables::mrmfilerestartin
```

Definition at line 54 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mhm\_eval::mhm←\_eval(), and mo\_mrm\_init::mrm\_init().

### 17.12.2.11 nerror\_model

```
integer(i4), parameter, public mo_common_mhm_mrm_variables::nerror_model = 2
```

Definition at line 87 of file mo\_common\_mHM\_mRM\_variables.f90.

### 17.12.2.12 niterations

```
integer(i4), public mo_common_mhm_mrm_variables::niterations
```

Definition at line 74 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::check\_optimization\_settings(), mo\_common\_mhm\_mrm←read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().

### 17.12.2.13 ntstepday

```
integer(i4), public mo_common_mhm_mrm_variables::ntstepday
```

Definition at line 47 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_meteo\_forcings::chunk\_size(), mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_datetime\_type::datetimeinfo\_init(), mo\_mrm\_objective\_function\_runoff::extract\_runoff(), mo\_write\_fluxes\_states::fluxesunit(), mo\_meteo\_forcings::is\_read(), mhm\_driver(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_read\_data::mrm\_read\_discharge(), and mo\_mrm\_write::mrm\_write().

### 17.12.2.14 opti\_function

```
integer(i4), public mo_common_mhm_mrm_variables::opti_function
```

Definition at line 63 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_discharge(), mo\_mrm\_objective\_function\_runoff::multi\_objective\_runoff(), mo\_objective\_function::objective(), mo\_objective\_function::objective\_master(), mo\_objective\_function::objective\_subprocess(), mo\_optimization::optimization(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_master(), and mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_subprocess().

### 17.12.2.15 opti\_method

```
integer(i4), public mo_common_mhm_mrm_variables::opti_method
```

Definition at line 59 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_optimization::optimization(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_master(), and mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_subprocess().

### 17.12.2.16 optimize

```
logical, public mo_common_mhm_mrm_variables::optimize
```

Definition at line 67 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_read\_data::mrm\_read\_discharge(), and mo\_mrm\_mpr::mrm\_update\_param().

### 17.12.2.17 optimize\_restart

```
logical, public mo_common_mhm_mrm_variables::optimize_restart
```

Definition at line 69 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().

### 17.12.2.18 read\_restart

logical, public mo\_common\_mhm\_mrm\_variables::read\_restart

Definition at line 34 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mhm\_eval::mhm\_eval(), mo\_startup::mhm\_initialize(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.12.2.19 readper

type(period), public mo\_common\_mhm\_mrm\_variables::readper

Definition at line 40 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_meteo\_forcings::meteo\_forcings\_wrapper(), mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

### 17.12.2.20 resolutionrouting

real(dp), dimension(:), allocatable, public mo\_common\_mhm\_mrm\_variables::resolutionrouting

Definition at line 33 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_check\_resolution(), mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_mpr::mrm\_update\_param(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.12.2.21 sa\_temp

real(dp), public mo\_common\_mhm\_mrm\_variables::sa\_temp

Definition at line 77 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().

### 17.12.2.22 sce\_ngs

integer(i4), public mo\_common\_mhm\_mrm\_variables::sce\_ngs

Definition at line 79 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::check\_optimization\_settings(), mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), and mo\_optimization::optimization().

### 17.12.2.23 sce\_npg

integer(i4), public mo\_common\_mhm\_mrm\_variables::sce\_npg

Definition at line 81 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by `mo_common_mhm_mrm_read_config::check_optimization_settings()`, `mo_common_mhm_mrm_read_config::common_mhm_mrm_read_config()`, and `mo_optimization::optimization()`.

#### 17.12.2.24 `sce_nps`

```
integer(i4), public mo_common_mhm_mrm_variables::sce_nps
```

Definition at line 83 of file `mo_common_mHM_mRM_variables.f90`.

Referenced by `mo_common_mhm_mrm_read_config::check_optimization_settings()`, `mo_common_mhm_mrm_read_config::common_mhm_mrm_read_config()`, and `mo_optimization::optimization()`.

#### 17.12.2.25 `seed`

```
integer(i8), public mo_common_mhm_mrm_variables::seed
```

Definition at line 72 of file `mo_common_mHM_mRM_variables.f90`.

Referenced by `mo_common_mhm_mrm_read_config::common_mhm_mrm_read_config()`, and `mo_optimization::optimization()`.

#### 17.12.2.26 `simper`

```
type(period), dimension(:), allocatable, public mo_common_mhm_mrm_variables::simper
```

Definition at line 39 of file `mo_common_mHM_mRM_variables.f90`.

Referenced by `mo_meteo_forcings::chunk_size()`, `mo_common_mhm_mrm_read_config::common_mhm_mrm_read_config()`, `mo_common_datetime_type::datetimeinfo_init()`, `mo_write_fluxes_states::fluxesunit()`, `mo_meteo_forcings::is_read()`, `mo_mhm_eval::mhm_eval()`, `mo_mrm_read_data::mrm_read_discharge()`, `mo_mrm_read_data::mrm_read_total_runoff()`, `mo_mrm_write::mrm_write()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

#### 17.12.2.27 `timestep`

```
integer(i4), public mo_common_mhm_mrm_variables::timestep
```

Definition at line 31 of file `mo_common_mHM_mRM_variables.f90`.

Referenced by `mo_common_mhm_mrm_read_config::common_mhm_mrm_read_config()`, `mo_startup::constants_init()`, `mo_common_datetime_type::datetimeinfo_increment()`, `mo_common_datetime_type::datetimeinfo_init()`, `mo_write_fluxes_states::fluxesunit()`, `mo_meteo_forcings::is_read()`, `mo_mhm_eval::mhm_eval()`, `mo_mrm_mpr::mrm_init_param()`, `mo_mrm_read_data::mrm_read_total_runoff()`, `mo_mrm_mpr::mrm_update_param()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

#### 17.12.2.28 `warmingdays`

```
integer(i4), dimension(:), allocatable, public mo_common_mhm_mrm_variables::warmingdays
```

Definition at line 41 of file `mo_common_mHM_mRM_variables.f90`.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_mrm\_objective\_↔  
function\_runoff::extract\_runoff(), mo\_mhm\_eval::mhm\_eval(), and mo\_mrm\_write::mrm\_write().

### 17.12.2.29 warmper

type(period), dimension(:), allocatable, public mo\_common\_mhm\_mrm\_variables::warmper

Definition at line 37 of file mo\_common\_mHM\_mRM\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_write\_ascii::write\_↔  
\_configfile(), and mo\_mrm\_write::write\_configfile().

## 17.13 mo\_common\_read\_config Module Reference

Reading of main model configurations.

### Functions/Subroutines

- subroutine, public [common\\_read\\_config](#) (file\_namelist, unamelist)  
*Read main configurations commonly used by mHM, mRM and MPR.*
- subroutine, public [set\\_land\\_cover\\_scenes\\_id](#) (sim\_Per, LCyear\_Id)  
*Read main configurations commonly used by mHM, mRM and MPR.*
- subroutine [init\\_domain\\_variable](#) (nDomains, optiData, domainMeta)
- subroutine [init\\_domain\\_variable\\_for\\_master](#) (domainMeta, colMasters, colDomain)
- subroutine [distributeddomainsroundrobin](#) (nproc, rank, domainMeta)
- subroutine [distribute\\_processes\\_to\\_domains\\_according\\_to\\_role](#) (optiData, rank, domainMeta, colMasters, colDomain)

### 17.13.1 Detailed Description

Reading of main model configurations.

This routine reads the configurations of namelists commonly used by mHM, mRM and MPR

#### Authors

Matthias Zink

#### Date

Dec 2012

### 17.13.2 Function/Subroutine Documentation

#### 17.13.2.1 common\_read\_config()

```
subroutine, public mo_common_read_config::common_read_config (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist )
```

Read main configurations commonly used by mHM, mRM and MPR.

Read the main configurations commonly used by mHM, mRM and MPR, namely: project\_description, directories←  
\_general, mainconfig, processSelection, LCover

#### Parameters

|    |                                      |              |
|----|--------------------------------------|--------------|
| in | <i>character(*) :: file_namelist</i> | name of file |
| in | <i>integer :: unamelist</i>          | id of file   |

#### Authors

Matthias Zink

#### Date

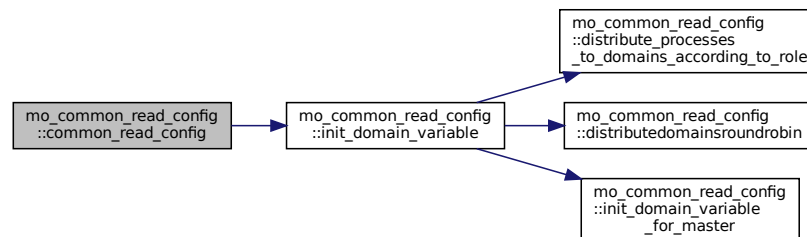
Dec 2012

Definition at line 50 of file mo\_common\_read\_config.f90.

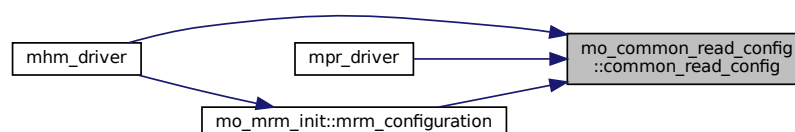
References mo\_common\_variables::contact, mo\_common\_variables::conventions, mo\_common\_variables←  
::dircommonfiles, mo\_common\_variables::dirconfigout, mo\_common\_variables::dircover, mo\_common\_variables←  
::dirmorpho, mo\_common\_variables::dirout, mo\_common\_variables::domainmeta, mo\_common\_variables←  
::filelatlon, mo\_common\_variables::history, mo\_common\_variables::iflag\_cordinate\_sys, init\_domain\_variable(),  
mo\_common\_variables::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::lcfilename,  
mo\_common\_constants::maxnlcovers, mo\_common\_constants::maxnodomains, mo\_common\_variables::mhm←  
\_details, mo\_common\_variables::mhmfilerestartout, mo\_common\_variables::mrmfilerestartout, mo\_common←  
\_variables::nlcoversscene, mo\_common\_variables::nprocesses, mo\_common\_variables::nunique0domains,  
mo\_common\_variables::processmatrix, mo\_common\_variables::project\_details, mo\_common\_variables←  
::resolutionhydrology, mo\_common\_variables::setup\_description, mo\_common\_variables::simulation\_type, and  
mo\_common\_variables::write\_restart.

Referenced by mhm\_driver(), mpr\_driver(), and mo\_mrm\_init::mrm\_configuration().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.13.2.2 distribute\_processes\_to\_domains\_according\_to\_role()

```

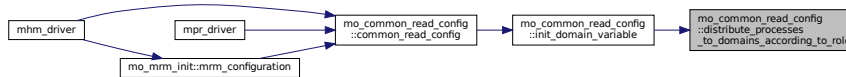
subroutine mo_common_read_config::distribute_processes_to_domains_according_to_role (
  integer(i4), dimension(:), intent(in) optiData,
  integer(i4), intent(in) rank,
  type(domain_meta), intent(inout) domainMeta,
  integer(i4), intent(out) colMasters,
  integer(i4), intent(out) colDomain )

```

Definition at line 478 of file mo\_common\_read\_config.f90.

Referenced by init\_domain\_variable().

Here is the caller graph for this function:



### 17.13.2.3 distributeddomainsroundrobin()

```

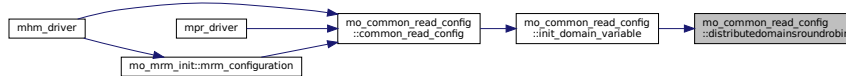
subroutine mo_common_read_config::distributeddomainsroundrobin (
  integer(i4), intent(in) nproc,
  integer(i4), intent(in) rank,
  type(domain_meta), intent(inout) domainMeta )

```

Definition at line 455 of file mo\_common\_read\_config.f90.

Referenced by init\_domain\_variable().

Here is the caller graph for this function:



### 17.13.2.4 init\_domain\_variable()

```

subroutine mo_common_read_config::init_domain_variable (
  integer(i4), intent(in) nDomains,
  integer(i4), dimension(:), intent(in) optiData,
  type(domain_meta), intent(inout) domainMeta )

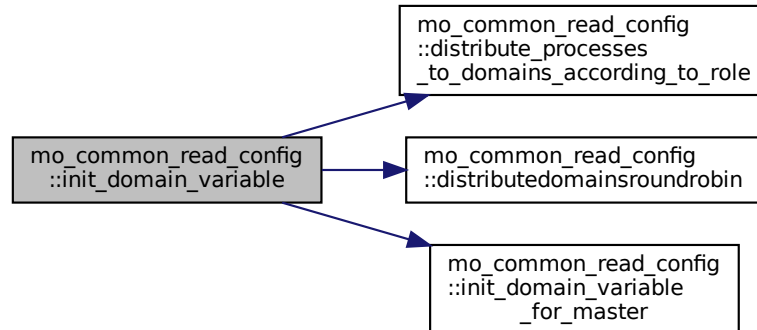
```

Definition at line 358 of file mo\_common\_read\_config.f90.

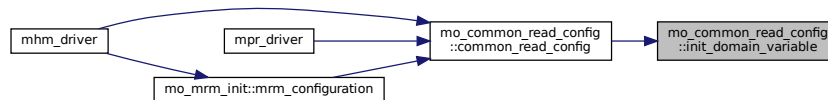
References mo\_common\_variables::comm, distribute\_processes\_to\_domains\_according\_to\_role(), distributeddomainsroundrobin(), and init\_domain\_variable\_for\_master().

Referenced by `common_read_config()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.13.2.5 `init_domain_variable_for_master()`

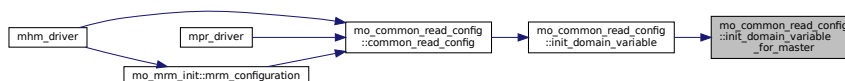
```

subroutine mo_common_read_config::init_domain_variable_for_master (
    type(domain_meta), intent(inout) domainMeta,
    integer(i4), intent(out) colMasters,
    integer(i4), intent(out) colDomain )
  
```

Definition at line 435 of file `mo_common_read_config.f90`.

Referenced by `init_domain_variable()`.

Here is the caller graph for this function:





### 17.13.2.6 set\_land\_cover\_scenes\_id()

```
subroutine, public mo_common_read_config::set_land_cover_scenes_id (
    type(period), dimension(:), intent(in) sim_Per,
    integer(i4), dimension(:, :), intent(inout), allocatable LCyear_Id )
```

Read main configurations commonly used by mHM, mRM and MPR.

Read the main configurations commonly used by mHM, mRM and MPR, namely: project\_description, directories\_↵\_general, mainconfig, processSelection, LCover

#### Parameters

|         |   |  |
|---------|---|--|
| in      | <i>type(period), dimension(:) :: sim_Per</i>      |  |
| in, out | <i>integer(i4), dimension(:, :) :: LCyear_Id</i>  |  |
| in, out | <i>character(256), dimension(:) :: LCfilename</i> |  |

#### Authors

Matthias Zink

#### Date

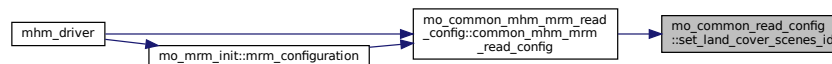
Dec 2012

Definition at line 270 of file mo\_common\_read\_config.f90.

References mo\_common\_variables::domainmeta, mo\_common\_variables::lc\_year\_end, mo\_common\_variables\_↵::lc\_year\_start, mo\_common\_variables::nlcoverscene, and mo\_common\_constants::nodata\_i4.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config().

Here is the caller graph for this function:



## 17.14 mo\_common\_read\_data Module Reference

TODO: add description.

### Functions/Subroutines

- subroutine, public [read\\_dem](#)  
*TODO: add description.*
- subroutine, public [read\\_lcover](#)  
*TODO: add description.*

#### 17.14.1 Detailed Description

TODO: add description.

TODO: add description

**Authors**

Robert Schweppe

**Date**

Jun 2018

**17.14.2 Function/Subroutine Documentation****17.14.2.1 read\_dem()**

```
subroutine, public mo_common_read_data::read_dem
```

TODO: add description.

TODO: add description

**Authors**

Robert Schweppe

**Date**

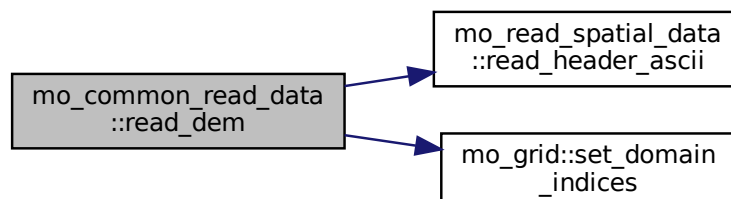
Jun 2018

Definition at line 41 of file mo\_common\_read\_data.f90.

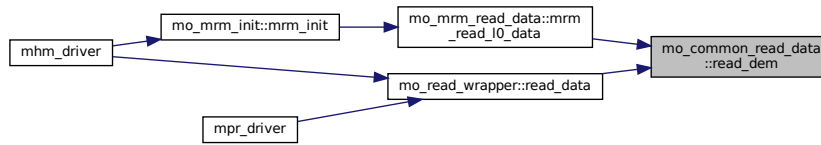
References mo\_common\_variables::dirmorpho, mo\_common\_variables::domainmeta, mo\_common\_file::file\_dem, mo\_common\_variables::l0\_elev, mo\_common\_variables::level0, mo\_common\_constants::nodata\_dp, mo\_read\_spatial\_data::read\_header\_ascii(), mo\_common\_variables::resolutionhydrology, mo\_grid::set\_domain\_indices(), and mo\_common\_file::udem.

Referenced by mo\_mrm\_read\_data::mrm\_read\_l0\_data(), and mo\_read\_wrapper::read\_data().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.14.2.2 read\_lcover()

subroutine, public mo\_common\_read\_data::read\_lcover

TODO: add description.

TODO: add description

#### Authors

Robert Schweppe

#### Date

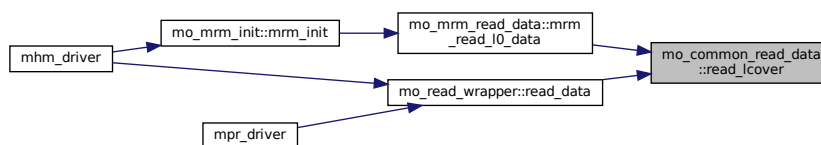
Jun 2018

Definition at line 143 of file mo\_common\_read\_data.f90.

References mo\_common\_variables::dirlcover, mo\_common\_variables::domainmeta, mo\_common\_variables::l0\_lcover, mo\_common\_variables::lcfilename, mo\_common\_variables::level0, mo\_common\_variables::nlcoverscene, mo\_common\_constants::nodata\_i4, and mo\_common\_file::ulcoverclass.

Referenced by mo\_mrm\_read\_data::mrm\_read\_io\_data(), and mo\_read\_wrapper::read\_data().

Here is the caller graph for this function:



## 17.15 mo\_common\_restart Module Reference

TODO: add description.

### Functions/Subroutines

- subroutine, public [write\\_grid\\_info](#) (grid\_in, level\_name, nc)  
*write restart files for each domain*

- subroutine, public `read_grid_info` (domainID, InFile, level\_name, new\_grid)  
*reads configuration apart from Level 11 configuration from a restart directory*

### 17.15.1 Detailed Description

TODO: add description.

TODO: add description

#### Authors

Robert Schweppe

#### Date

Jun 2018

### 17.15.2 Function/Subroutine Documentation

#### 17.15.2.1 `read_grid_info()`

```
subroutine, public mo_common_restart::read_grid_info (
    integer(i4), intent(in) domainID,
    character(256), intent(in) InFile,
    character(*), intent(in) level_name,
    type(grid), intent(inout) new_grid )
```

reads configuration apart from Level 11 configuration from a restart directory

read configuration variables from a given restart directory and initializes all configuration variables, that are initialized in the subroutine initialise, contained in module `mo_startup`.

#### Parameters

|         |                                   |                                     |
|---------|-----------------------------------|-------------------------------------|
| in      | <i>integer(i4) :: iDomain</i>     | number of domain                    |
| in      | <i>character(256) :: InFile</i>   | Input Path including trailing slash |
| in      | <i>character(*) :: level_name</i> | level_name (id)                     |
| in, out | <i>type(Grid) :: new_grid</i>     | grid to save information to         |

#### Authors

Stephan Thober

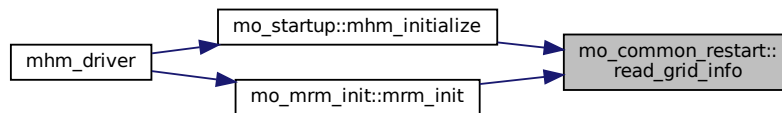
**Date**

Apr 2013

Definition at line 162 of file mo\_common\_restart.f90.

Referenced by mo\_startup::mhm\_initialize(), and mo\_mrm\_init::mrm\_init().

Here is the caller graph for this function:

**17.15.2.2 write\_grid\_info()**

```

subroutine, public mo_common_restart::write_grid_info (
    type(Grid), intent(in) grid_in,
    character(*), intent(in) level_name,
    type(NcDataset), intent(inout) nc )
  
```

write restart files for each domain

write restart files for each domain. For each domain three restart files are written. These are `xxx_states.nc`, `xxx_L11_config.nc`, and `xxx_config.nc` (`xxx` being the three digit domain index). If a variable is added here, it should also be added in the read restart routines below.

**Parameters**

|         |   |                                   |
|---------|---|-----------------------------------|
| in      | <code>type(Grid) :: grid_in</code>      | level to be written               |
| in      | <code>character(*) :: level_name</code> | level_id                          |
| in, out | <code>type(NcDataset) :: nc</code>      | NcDataset to write information to |

**Authors**

Stephan Thober

**Date**

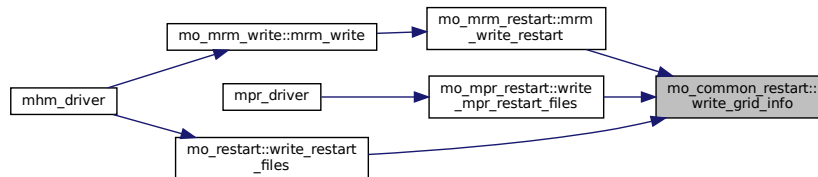
Jun 2014

Definition at line 62 of file mo\_common\_restart.f90.

References mo\_common\_constants::nodata\_dp, and mo\_common\_constants::nodata\_i4.

Referenced by mo\_mrm\_restart::mrm\_write\_restart(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

Here is the caller graph for this function:



## 17.16 mo\_common\_variables Module Reference

Provides structures needed by mHM, mRM and/or mpr.

### Data Types

- type [period](#)
- type [grid](#)
- type [gridremapper](#)
- type [domain\\_meta](#)

### Variables

- character(1024), public [project\\_details](#)
- character(1024), public [setup\\_description](#)
- character(1024), public [simulation\\_type](#)
- character(256), public [conventions](#)
- character(1024), public [contact](#)
- character(1024), public [mhm\\_details](#)
- character(1024), public [history](#)
- integer(i4), public [iflag\\_cordinate\\_sys](#)
- real(dp), dimension(:), allocatable, public [resolutionhydrology](#)
- integer(i4), dimension(:), allocatable, public [IO\\_domain](#)
- logical, public [write\\_restart](#)
- character(256), dimension(:), allocatable, public [mhmfilerrestartout](#)
- character(256), dimension(:), allocatable, public [mrmfilerrestartout](#)
- character(256), public [dirconfigout](#)
- character(256), public [dircommonfiles](#)
- character(256), dimension(:), allocatable, public [dirmorpho](#)
- character(256), dimension(:), allocatable, public [dirlcover](#)
- character(256), dimension(:), allocatable, public [dirout](#)
- character(256), dimension(:), allocatable, public [filelatlon](#)

- type([grid](#)), dimension(:), allocatable, target, public [level0](#)
- type([grid](#)), dimension(:), allocatable, target, public [level1](#)
- type([gridremapper](#)), dimension(:), allocatable, public [l0\\_l1\\_remap](#)
- real(dp), dimension(:), allocatable, public [l0\\_elev](#)
- integer(i4), dimension(:, :), allocatable, public [l0\\_lcover](#)
- type(mpi\_comm) [comm](#)
- type([domain\\_meta](#)), public [domainmeta](#)
- integer(i4), public [nunique0domains](#)
- integer(i4), public [nlcoverscene](#)
- character(256), dimension(:), allocatable, public [lcfilename](#)
- integer(i4), dimension(:), allocatable, public [lc\\_year\\_start](#)
- integer(i4), dimension(:), allocatable, public [lc\\_year\\_end](#)
- integer(i4), parameter, public [nprocesses](#) = 11
- integer(i4), dimension([nprocesses](#), 3), public [processmatrix](#)
- real(dp), dimension(:, :), allocatable, target, public [global\\_parameters](#)
- character(256), dimension(:), allocatable, public [global\\_parameters\\_name](#)
- logical [alma\\_convention](#)

### 17.16.1 Detailed Description

Provides structures needed by mHM, mRM and/or mpr.

Provides the global structure period that is used by both mHM and mRM.

#### Authors

Stephan Thober

#### Date

Sep 2015

### 17.16.2 Variable Documentation

#### 17.16.2.1 alma\_convention

```
logical mo_common_variables::alma_convention
```

Definition at line 209 of file mo\_common\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), and mo\_mrm\_read\_data::mrm\_read\_total\_runoff().

#### 17.16.2.2 comm

```
type(mpi_comm) mo_common_variables::comm
```

Definition at line 131 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::init\_domain\_variable().

**17.16.2.3 contact**

```
character(1024), public mo_common_variables::contact
```

Definition at line 33 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_mrm\_river\_head::create\_output().

**17.16.2.4 conventions**

```
character(256), public mo_common_variables::conventions
```

Definition at line 32 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_mrm\_river\_head::create\_output().

**17.16.2.5 dircommonfiles**

```
character(256), public mo_common_variables::dircommonfiles
```

Definition at line 52 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_read\_wrapper::read\_data().

**17.16.2.6 dirconfigout**

```
character(256), public mo_common_variables::dirconfigout
```

Definition at line 51 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_write::mrm\_write\_optifile(), mo\_mrm\_write::mrm\_write\_optinamelist(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), mo\_write\_ascii::write\_optifile(), and mo\_write\_ascii::write\_optinamelist().

**17.16.2.7 dirlcover**

```
character(256), dimension(:), allocatable, public mo_common_variables::dirlcover
```

Definition at line 54 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_init::config\_output(), mo\_common\_read\_data::read\_lcover(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.16.2.8 dirmorpho**

```
character(256), dimension(:), allocatable, public mo_common_variables::dirmorpho
```

Definition at line 53 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_init::config\_output(), mo\_mrm\_read\_data::mrm\_read\_IO\_data(), mo\_read\_wrapper::read\_data(), mo\_common\_read\_data::read\_dem(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().



### 17.16.2.9 dirout

```
character(256), dimension(:), allocatable, public mo_common_variables::dirout
```

Definition at line 55 of file mo\_common\_variables.f90.

Referenced by mo\_write\_fluxes\_states::close(), mo\_mrm\_write\_fluxes\_states::close(), mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_init::config\_output(), mo\_write\_fluxes\_states::createoutputfile(), mo\_mrm\_write\_fluxes\_states::createoutputfile(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), and mo\_mrm\_write::write\_daily\_obs\_sim\_discharge().

### 17.16.2.10 domainmeta

```
type(domain_meta), public mo_common_variables::domainmeta
```

Definition at line 161 of file mo\_common\_variables.f90.

Referenced by mo\_mrm\_river\_head::calc\_channel\_elevation(), mo\_mrm\_net\_startup::celllength(), mo\_common\_mhm\_mrm\_read\_config::common\_check\_resolution(), mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_riv\_temp\_class::config(), mo\_mrm\_init::config\_output(), mo\_common\_mhm\_mrm\_mpi\_tools::distribute\_parameterset(), mo\_common\_mhm\_mrm\_mpi\_tools::get\_parameterset(), mo\_mrm\_net\_startup::l11\_calc\_celerity(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mhm\_eval::mhm\_eval(), mo\_startup::mhm\_initialize(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_mpr\_eval::mpr\_eval(), mo\_mpr\_startup::mpr\_initialize(), mo\_mpr\_read\_config::mpr\_read\_config(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_discharge(), mo\_mrm\_read\_data::mrm\_read\_l0\_data(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_write::mrm\_write(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_objective\_function::objective\_et\_kge\_catchment\_avg(), mo\_objective\_function::objective\_kge\_q\_et(), mo\_objective\_function::objective\_kge\_q\_rmse\_et(), mo\_objective\_function::objective\_kge\_q\_rmse\_tws(), mo\_objective\_function::objective\_kge\_q\_sm\_corr(), mo\_objective\_function::objective\_master(), mo\_objective\_function::objective\_neutrons\_kge\_catchment\_avg(), mo\_objective\_function::objective\_q\_et\_tws\_kge\_catchment\_avg(), mo\_objective\_function::objective\_sm\_corr(), mo\_objective\_function::objective\_sm\_kge\_catchment\_avg(), mo\_objective\_function::objective\_sm\_pd(), mo\_objective\_function::objective\_sm\_sse\_standard\_score(), mo\_objective\_function::objective\_subprocess(), mo\_optimization::optimization(), mo\_meteo\_forcings::prepare\_meteo\_forcings\_data(), mo\_read\_wrapper::read\_data(), mo\_common\_read\_data::read\_dem(), mo\_common\_read\_data::read\_lcover(), mo\_common\_read\_config::set\_land\_cover\_scenes\_id(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_master(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_subprocess(), mo\_mrm\_init::variables\_alloc\_routing(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), mo\_mrm\_write::write\_daily\_obs\_sim\_discharge(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

### 17.16.2.11 filelatlon

```
character(256), dimension(:), allocatable, public mo_common_variables::filelatlon
```

Definition at line 56 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_read\_latlon::read\_latlon().

### 17.16.2.12 global\_parameters

```
real(dp), dimension(:, :), allocatable, target, public mo_common_variables::global_parameters
```

Definition at line 197 of file mo\_common\_variables.f90.

Referenced by `mo_common_mhm_mrm_read_config::check_optimization_settings()`, `mo_mrm_objective_↔function_runoff::loglikelihood_evin2013_2()`, `mo_multi_param_reg::mpr()`, `mo_mpr_read_config::mpr_read_↔_config()`, `mo_mrm_init::mrm_init()`, `mo_optimization::optimization()`, `mo_read_wrapper::read_data()`, `mo_↔mrm_read_config::read_mrm_routing_params()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_↔configfile()`.

### 17.16.2.13 global\_parameters\_name

```
character(256), dimension(:), allocatable, public mo_common_variables::global_parameters_name
```

Definition at line 201 of file `mo_common_variables.f90`.

Referenced by `mo_mpr_read_config::mpr_read_config()`, `mo_mrm_read_config::read_mrm_routing_params()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

### 17.16.2.14 history

```
character(1024), public mo_common_variables::history
```

Definition at line 35 of file `mo_common_variables.f90`.

Referenced by `mo_common_read_config::common_read_config()`, and `mo_mrm_river_head::create_output()`.

### 17.16.2.15 iflag\_cordinate\_sys

```
integer(i4), public mo_common_variables::iflag_cordinate_sys
```

Definition at line 40 of file `mo_common_variables.f90`.

Referenced by `mo_mrm_river_head::calc_slope()`, `mo_common_read_config::common_read_config()`, `mo_↔write_fluxes_states::createoutputfile()`, `mo_mrm_write_fluxes_states::createoutputfile()`, `mo_grid::l0_grid_setup()`, `mo_mrm_net_startup::l11_stream_features()`, `mo_mrm_mpr::mrm_init_param()`, `mo_mrm_mpr::mrm_update_↔_param()`, `mo_mrm_write_fluxes_states::newoutputdataset()`, `mo_write_fluxes_states::newoutputdataset()`, and `mo_write_ascii::write_configfile()`.

### 17.16.2.16 l0\_domain

```
integer(i4), dimension(:), allocatable, public mo_common_variables::l0_domain
```

Definition at line 42 of file `mo_common_variables.f90`.

### 17.16.2.17 l0\_elev

```
real(dp), dimension(:), allocatable, public mo_common_variables::l0_elev
```

Definition at line 122 of file `mo_common_variables.f90`.

Referenced by `mo_mrm_river_head::calc_channel_elevation()`, `mo_mpr_startup::l0_check_input()`, `mo_mrm_net_↔_startup::l11_stream_features()`, and `mo_common_read_data::read_dem()`.

**17.16.2.18 l0\_l1\_remap**

```
type(gridremapper), dimension(:), allocatable, public mo_common_variables::l0_l1_remap
```

Definition at line 115 of file mo\_common\_variables.f90.

Referenced by mo\_mpr\_eval::mpr\_eval(), mo\_mpr\_startup::mpr\_initialize(), and mo\_mrm\_init::mrm\_init().

**17.16.2.19 l0\_lcover**

```
integer(i4), dimension(:, :), allocatable, public mo_common_variables::l0_lcover
```

Definition at line 124 of file mo\_common\_variables.f90.

Referenced by mo\_mpr\_startup::l0\_check\_input(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mpr\_eval::mpr\_eval(), mo\_mrm\_read\_data::mrm\_read\_l0\_data(), and mo\_common\_read\_data::read\_lcover().

**17.16.2.20 lc\_year\_end**

```
integer(i4), dimension(:), allocatable, public mo_common_variables::lc_year_end
```

Definition at line 171 of file mo\_common\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_restart::read\_restart\_states(), mo\_common\_read\_config::set\_land\_cover\_scenes\_id(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), mo\_mpr\_restart::write\_eff\_params(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

**17.16.2.21 lc\_year\_start**

```
integer(i4), dimension(:), allocatable, public mo_common_variables::lc_year_start
```

Definition at line 170 of file mo\_common\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_restart::read\_restart\_states(), mo\_common\_read\_config::set\_land\_cover\_scenes\_id(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), mo\_mpr\_restart::write\_eff\_params(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

**17.16.2.22 lcfilename**

```
character(256), dimension(:), allocatable, public mo_common_variables::lcfilename
```

Definition at line 169 of file mo\_common\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_read\_config::common\_read\_config(), mo\_common\_read\_data::read\_lcover(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.16.2.23 level0**

```
type(grid), dimension(:), allocatable, target, public mo_common_variables::level0
```

Definition at line 98 of file mo\_common\_variables.f90.

Referenced by mo\_mrm\_river\_head::avg\_and\_write\_timestep(), mo\_mrm\_river\_head::calc\_river\_head(), mo\_mrm\_net\_startup::celllength(), mo\_mrm\_river\_head::init\_masked\_zeros\_l0(), mo\_mpr\_startup::l0\_check\_input(), mo\_mrm\_init::l0\_check\_input\_routing(), mo\_mpr\_startup::l0\_variable\_init(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_startup::mhm\_initialize(), mo\_mpr\_eval::mpr\_eval(), mo\_mpr\_startup::mpr\_initialize(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_read\_data::mrm\_read\_l0\_data(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_read\_wrapper::read\_data(), mo\_common\_read\_data::read\_dem(), mo\_common\_read\_data::read\_lcover(), mo\_mrm\_river\_head::reset\_sum(), mo\_mrm\_init::variables\_alloc\_routing(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.16.2.24 level1**

```
type(grid), dimension(:), allocatable, target, public mo_common_variables::level1
```

Definition at line 99 of file mo\_common\_variables.f90.

Referenced by mo\_write\_fluxes\_states::createoutputfile(), mo\_mrm\_net\_startup::l11\_l1\_mapping(), mo\_meteo\_forcings::meteo\_forcings\_wrapper(), mo\_meteo\_forcings::meteo\_weights\_wrapper(), mo\_mhm\_eval::mhm\_eval(), mo\_startup::mhm\_initialize(), mo\_mpr\_eval::mpr\_eval(), mo\_mpr\_startup::mpr\_initialize(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_read\_data::mrm\_read\_total\_runoff(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_objective\_function::objective\_kge\_q\_et(), mo\_objective\_function::objective\_kge\_q\_sm\_corr(), mo\_objective\_function::objective\_sm\_corr(), mo\_objective\_function::objective\_sm\_kge\_catchment\_avg(), mo\_objective\_function::objective\_sm\_pd(), mo\_objective\_function::objective\_sm\_sse\_standard\_score(), mo\_restart::read\_restart\_states(), mo\_read\_optional\_data::readoptidataobs(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

**17.16.2.25 mhm\_details**

```
character(1024), public mo_common_variables::mhm_details
```

Definition at line 34 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_mrm\_river\_head::create\_output().

**17.16.2.26 mhmfilerestartout**

```
character(256), dimension(:), allocatable, public mo_common_variables::mhmfilerestartout
```

Definition at line 49 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), mpr\_driver(), and mo\_write\_ascii::write\_configfile().

**17.16.2.27 mrmfilerestartout**

```
character(256), dimension(:), allocatable, public mo_common_variables::mrmfilerestartout
```

Definition at line 50 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), mo\_mrm\_write::mrm\_write(), and mo\_mrm\_write::write\_configfile().

### 17.16.2.28 nlcoverscene

```
integer(i4), public mo_common_variables::nlcoverscene
```

Definition at line 168 of file mo\_common\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_common\_read\_config::common\_read\_config(), mo\_mpr\_startup::init\_eff\_params(), mo\_mpr\_startup::io\_check\_input(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_common\_read\_data::read\_lcover(), mo\_restart::read\_restart\_states(), mo\_common\_read\_config::set\_land\_cover\_scenes\_id(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

### 17.16.2.29 nprocesses

```
integer(i4), parameter, public mo_common_variables::nprocesses = 11
```

Definition at line 176 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_write\_ascii::write\_optinamelist().

### 17.16.2.30 nunique10domains

```
integer(i4), public mo_common_variables::nunique10domains
```

Definition at line 162 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config().

### 17.16.2.31 processmatrix

```
integer(i4), dimension(nprocesses, 3), public mo_common_variables::processmatrix
```

Definition at line 188 of file mo\_common\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_read\_config::common\_read\_config(), mo\_startup::constants\_init(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_multi\_param\_reg::mpr(), mo\_mpr\_read\_config::mpr\_read\_config(), mo\_mrm\_init::mrm\_configuration(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_io\_data(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_mpr::mrm\_update\_param(), mo\_mrm\_write::mrm\_write\_optinamelist(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_meteo\_forcings::prepare\_meteo\_forcings\_data(), mo\_read\_wrapper::read\_data(), mo\_mrm\_read\_config::read\_mrm\_routing\_params(), mo\_restart::read\_restart\_states(), mo\_write\_ascii::write\_configfile(), mo\_mrm\_write::write\_configfile(), and mo\_mpr\_restart::write\_eff\_params().

### 17.16.2.32 project\_details

```
character(1024), public mo_common_variables::project_details
```

Definition at line 29 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_mrm\_river\_head::create\_output().

### 17.16.2.33 resolutionhydrology

```
real(dp), dimension(:), allocatable, public mo_common_variables::resolutionhydrology
```

Definition at line 41 of file mo\_common\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_read\_config::common\_check\_resolution(), mo\_common\_read\_config::common\_read\_config(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_startup::mpr\_initialize(), mo\_mrm\_init::mrm\_init(), mo\_common\_read\_data::read\_dem(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.16.2.34 setup\_description

```
character(1024), public mo_common_variables::setup_description
```

Definition at line 30 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_mrm\_river\_head::create\_output().

### 17.16.2.35 simulation\_type

```
character(1024), public mo_common_variables::simulation_type
```

Definition at line 31 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), and mo\_mrm\_river\_head::create\_output().

### 17.16.2.36 write\_restart

```
logical, public mo_common_variables::write_restart
```

Definition at line 43 of file mo\_common\_variables.f90.

Referenced by mo\_common\_read\_config::common\_read\_config(), mhm\_driver(), mpr\_driver(), mo\_mrm\_write::mrm\_write(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

## 17.17 mo\_file Module Reference

Provides file names and units for mHM.

### Variables

- character(len=\*), parameter `version` = '5.11.2'  
*Current mHM model version (will be set to 5.11.2)*
- character(len=\*), parameter `version_date` = 'Jul 2021'

- Time of current mHM model version release.*

  - character(len=\*), parameter `file_main` = 'mhm\_driver.f90'  
*Driver file.*
  - character(len=\*), parameter `file_namelist_mhm` = 'mhm.nml'  
*Namelist file name.*
  - integer, parameter `unamelist_mhm` = 30  
*Unit for namelist.*
  - character(len=\*), parameter `file_namelist_mhm_param` = 'mhm\_parameter.nml'  
*Parameter namelists file name.*
  - integer, parameter `unamelist_mhm_param` = 31  
*Unit for namelist.*
  - character(len=\*), parameter `file_defoutput` = 'mhm\_outputs.nml'  
*file defining mHM's outputs*
  - integer, parameter `undefoutput` = 67  
*Unit for file defining mHM's outputs.*

### 17.17.1 Detailed Description

Provides file names and units for mHM.

Provides all filenames as well as all units used for the hydrologic model mHM. The `version` parameter will be set during compilation to " 5.11.2". The `version_date` parameter will be set during compilation to " Jul 2021", if it is a release version, otherwise it will be the current date.

#### Authors

Matthias Cuntz  
Sebastian Mueller

#### Date

Jan 2012

### 17.17.2 Variable Documentation

#### 17.17.2.1 file\_defoutput

```
character(len = *), parameter mo_file::file_defoutput = 'mhm_outputs.nml'
```

file defining mHM's outputs

Definition at line 48 of file mo\_file.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config().

#### 17.17.2.2 file\_main

```
character(len = *), parameter mo_file::file_main = 'mhm_driver.f90'
```

Driver file.

Definition at line 38 of file mo\_file.f90.

Referenced by mhm\_driver().

### 17.17.2.3 file\_namelist\_mhm

```
character(len = *), parameter mo_file::file_namelist_mhm = 'mhm.nml'
```

Namelist file name.

Definition at line 40 of file mo\_file.f90.

### 17.17.2.4 file\_namelist\_mhm\_param

```
character(len = *), parameter mo_file::file_namelist_mhm_param = 'mhm_parameter.nml'
```

Parameter namelists file name.

Definition at line 44 of file mo\_file.f90.

Referenced by mo\_startup::constants\_init().

### 17.17.2.5 udefoutput

```
integer, parameter mo_file::udefoutput = 67
```

Unit for file defining mHM's outputs.

Definition at line 50 of file mo\_file.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config().

### 17.17.2.6 unamelist\_mhm

```
integer, parameter mo_file::unamelist_mhm = 30
```

Unit for namelist.

Definition at line 42 of file mo\_file.f90.

### 17.17.2.7 unamelist\_mhm\_param

```
integer, parameter mo_file::unamelist_mhm_param = 31
```

Unit for namelist.

Definition at line 46 of file mo\_file.f90.

### 17.17.2.8 version

```
character(len = *), parameter mo_file::version = '5.11.2'
```

Current mHM model version (will be set to 5.11.2)

Definition at line 31 of file mo\_file.f90.

Referenced by mo\_mrm\_river\_head::create\_output(), mo\_write\_fluxes\_states::createoutputfile(), mhm\_driver(), and mo\_write\_ascii::write\_configfile().



### 17.17.2.9 version\_date

```
character(len = *), parameter mo_file::version_date = 'Jul 2021'
```

Time of current mHM model version release.

Definition at line 34 of file mo\_file.f90.

Referenced by mhm\_driver().

## 17.18 mo\_global\_variables Module Reference

Global variables ONLY used in reading, writing and startup.

### Variables

- integer(i4) [output\\_deflate\\_level](#)
- logical [output\\_double\\_precision](#)
- integer(i4) [timestep\\_model\\_outputs](#)
- logical, dimension(noutflxstate) [outputflxstate](#)
- integer(i4), dimension(:), allocatable, public [timestep\\_model\\_inputs](#)
- logical, public [read\\_meteo\\_weights](#)
- character(256), public [inputformat\\_meteo\\_forcings](#)
- character(256), dimension(:), allocatable, public [dirprecipitation](#)
- character(256), dimension(:), allocatable, public [dirtemperature](#)
- character(256), dimension(:), allocatable, public [dirminttemperature](#)
- character(256), dimension(:), allocatable, public [dirmaxtemperature](#)
- character(256), dimension(:), allocatable, public [dirnetradiation](#)
- character(256), dimension(:), allocatable, public [dirabsvappressure](#)
- character(256), dimension(:), allocatable, public [dirwindspeed](#)
- character(256), dimension(:), allocatable, public [dirreferencecet](#)
- character(256), dimension(:), allocatable, public [dirradiation](#)
- integer(i4), parameter, public [routingstates](#) = 2
- type([grid](#)), dimension(:), allocatable, public [level2](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_temp\\_weights](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_pet\\_weights](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_pre\\_weights](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_pre](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_temp](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_pet](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_tmin](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_tmax](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_netrad](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_absvappress](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_windspeed](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_ssrd](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_strd](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_tann](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_sm](#)
- logical, dimension(:, :), allocatable, public [l1\\_sm\\_mask](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_neutronsdata](#)
- logical, dimension(:, :), allocatable, public [l1\\_neutronsdata\\_mask](#)
- integer(i4) [nsoilhorizons\\_sm\\_input](#)
- type(optidata), dimension(:), allocatable, public [l1\\_smobs](#)
- type(optidata), dimension(:), allocatable, public [l1\\_neutronsobs](#)

- type(optidata), dimension(:), allocatable, public [l1\\_etobs](#)
- type(optidata), dimension(:), allocatable, public [l1\\_twsaobs](#)
- real(dp), dimension(:), allocatable, public [l1\\_inter](#)
- real(dp), dimension(:), allocatable, public [l1\\_snowpack](#)
- real(dp), dimension(:), allocatable, public [l1\\_sealstw](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_soilmoist](#)
- real(dp), dimension(:), allocatable, public [l1\\_unsatstw](#)
- real(dp), dimension(:), allocatable, public [l1\\_satstw](#)
- real(dp), dimension(:), allocatable, public [l1\\_neutrons](#)
- real(dp), dimension(:), allocatable, public [l1\\_pet\\_calc](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_aetsoil](#)
- real(dp), dimension(:), allocatable, public [l1\\_aetcanopy](#)
- real(dp), dimension(:), allocatable, public [l1\\_aetsealed](#)
- real(dp), dimension(:), allocatable, public [l1\\_baseflow](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_infilsoil](#)
- real(dp), dimension(:), allocatable, public [l1\\_fastrunoff](#)
- real(dp), dimension(:), allocatable, public [l1\\_melt](#)
- real(dp), dimension(:), allocatable, public [l1\\_percol](#)
- real(dp), dimension(:), allocatable, public [l1\\_preeffect](#)
- real(dp), dimension(:), allocatable, public [l1\\_rain](#)
- real(dp), dimension(:), allocatable, public [l1\\_runoffseal](#)
- real(dp), dimension(:), allocatable, public [l1\\_slowrunoff](#)
- real(dp), dimension(:), allocatable, public [l1\\_snow](#)
- real(dp), dimension(:), allocatable, public [l1\\_throughfall](#)
- real(dp), dimension(:), allocatable, public [l1\\_total\\_runoff](#)
- real(dp), dimension(yearmonths\_i4), public [evap\\_coeff](#)
- real(dp), dimension(yearmonths\_i4), public [fdlay\\_prec](#)
- real(dp), dimension(yearmonths\_i4), public [fnight\\_prec](#)
- real(dp), dimension(yearmonths\_i4), public [fdlay\\_pet](#)
- real(dp), dimension(yearmonths\_i4), public [fnight\\_pet](#)
- real(dp), dimension(yearmonths\_i4), public [fdlay\\_temp](#)
- real(dp), dimension(yearmonths\_i4), public [fnight\\_temp](#)
- real(dp), dimension(yearmonths\_i4), public [fdlay\\_ssr](#)
- real(dp), dimension(yearmonths\_i4), public [fnight\\_ssr](#)
- real(dp), dimension(yearmonths\_i4), public [fdlay\\_strd](#)
- real(dp), dimension(yearmonths\_i4), public [fnight\\_strd](#)
- real(dp), dimension(:), allocatable, public [neutron\\_integral\\_afast](#)

### 17.18.1 Detailed Description

Global variables ONLY used in reading, writing and startup.

TODO: add description

#### Authors

Luis Samaniego

#### Date

Dec 2012

### 17.18.2 Variable Documentation

### 17.18.2.1 dirabsvappressure

`character(256), dimension(:), allocatable, public mo_global_variables::dirabsvappressure`

Definition at line 76 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_read_config::mhm_read_config()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.2 dirmaxtemperature

`character(256), dimension(:), allocatable, public mo_global_variables::dirmaxtemperature`

Definition at line 74 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_read_config::mhm_read_config()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.3 dirmintemperature

`character(256), dimension(:), allocatable, public mo_global_variables::dirmintemperature`

Definition at line 73 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_read_config::mhm_read_config()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.4 dirnetradiation

`character(256), dimension(:), allocatable, public mo_global_variables::dirnetradiation`

Definition at line 75 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_read_config::mhm_read_config()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.5 dirprecipitation

`character(256), dimension(:), allocatable, public mo_global_variables::dirprecipitation`

Definition at line 71 of file `mo_global_variables.f90`.

Referenced by `mo_startup::l2_variable_init()`, `mhm_driver()`, `mo_mhm_read_config::mhm_read_config()`, `mo_meteo_forcings::prepare_meteo_forcings_data()`, and `mo_write_ascii::write_configfile()`.

### 17.18.2.6 dirradiation

`character(256), dimension(:), allocatable, public mo_global_variables::dirradiation`

Definition at line 80 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_read_config::mhm_read_config()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.7 dirreferenceet

```
character(256), dimension(:), allocatable, public mo_global_variables::dirreferenceet
```

Definition at line 78 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config(), mo\_meteo\_forcings::prepare\_meteo\_forcings\_data(), and mo\_write\_ascii::write\_configfile().

### 17.18.2.8 dirtemperature

```
character(256), dimension(:), allocatable, public mo_global_variables::dirtemperature
```

Definition at line 72 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config(), mo\_meteo\_forcings::prepare\_meteo\_forcings\_data(), and mo\_write\_ascii::write\_configfile().

### 17.18.2.9 dirwindspeed

```
character(256), dimension(:), allocatable, public mo_global_variables::dirwindspeed
```

Definition at line 77 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

### 17.18.2.10 evap\_coeff

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::evap_coeff
```

Definition at line 173 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

### 17.18.2.11 fday\_pet

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fday_pet
```

Definition at line 176 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

### 17.18.2.12 fday\_prec

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fday_prec
```

Definition at line 174 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.18.2.13 fday\_ssrđ**

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fday_ssrđ
```

Definition at line 180 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.18.2.14 fday\_strđ**

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fday_strđ
```

Definition at line 182 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.18.2.15 fday\_temp**

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fday_temp
```

Definition at line 178 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.18.2.16 fnight\_pet**

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fnight_pet
```

Definition at line 177 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.18.2.17 fnight\_prec**

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fnight_prec
```

Definition at line 175 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.18.2.18 fnight\_ssrđ**

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fnight_ssrđ
```

Definition at line 181 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

### 17.18.2.19 `fnight_strd`

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fnight_strd
```

Definition at line 183 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, and `mo_mhm_read_config::mhm_read_config()`.

### 17.18.2.20 `fnight_temp`

```
real(dp), dimension(yearmonths_i4), public mo_global_variables::fnight_temp
```

Definition at line 179 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, and `mo_mhm_read_config::mhm_read_config()`.

### 17.18.2.21 `inputformat_meteo_forcings`

```
character(256), public mo_global_variables::inputformat_meteo_forcings
```

Definition at line 65 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_read_config::mhm_read_config()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.22 `l1_absvappress`

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_absvappress
```

Definition at line 107 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

### 17.18.2.23 `l1_aetcanopy`

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_aetcanopy
```

Definition at line 153 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_restart::read_restart_states()`, `mo_init_states::variables_alloc()`, `mo_init_states::variables_default_init()`, and `mo_restart::write_restart_files()`.

### 17.18.2.24 `l1_aetsealed`

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_aetsealed
```

Definition at line 154 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_restart::read_restart_states()`, `mo_init_states::variables_alloc()`, `mo_init_states::variables_default_init()`, and `mo_restart::write_restart_files()`.

### 17.18.2.25 l1\_aetsoil

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_aetsoil
```

Definition at line 152 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

### 17.18.2.26 l1\_baseflow

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_baseflow
```

Definition at line 155 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

### 17.18.2.27 l1\_etobs

```
type(optidata), dimension(:), allocatable, public mo_global_variables::l1_etobs
```

Definition at line 129 of file mo\_global\_variables.f90.

Referenced by mo\_objective\_function::create\_domain\_avg\_et(), mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_objective\_function::objective\_kge\_q\_et(), mo\_objective\_function::objective\_kge\_q\_rmse\_et(), and mo\_objective\_function::objective\_q\_et\_tws\_kge\_catchment\_avg().

### 17.18.2.28 l1\_fastrunoff

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_fastrunoff
```

Definition at line 157 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

### 17.18.2.29 l1\_infilsoil

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_infilsoil
```

Definition at line 156 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

### 17.18.2.30 l1\_inter

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_inter
```

Definition at line 140 of file mo\_global\_variables.f90.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_restart::read_restart_states()`, `mo_init_states::variables_alloc()`, `mo_init_states::variables_default_init()`, and `mo_restart::write_restart_files()`.

#### 17.18.2.31 `l1_melt`

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_melt
```

Definition at line 158 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_restart::read_restart_states()`, `mo_init_states::variables_alloc()`, `mo_init_states::variables_default_init()`, and `mo_restart::write_restart_files()`.

#### 17.18.2.32 `l1_netrad`

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_netrad
```

Definition at line 106 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, and `mo_meteo_forcings::prepare_meteo_forcings_data()`.

#### 17.18.2.33 `l1_neutrons`

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_neutrons
```

Definition at line 146 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_init_states::variables_alloc()`, and `mo_init_states::variables_default_init()`.

#### 17.18.2.34 `l1_neutronsdata`

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_neutronsdata
```

Definition at line 119 of file `mo_global_variables.f90`.

#### 17.18.2.35 `l1_neutronsdata_mask`

```
logical, dimension(:, :), allocatable, public mo_global_variables::l1_neutronsdata_mask
```

Definition at line 120 of file `mo_global_variables.f90`.

#### 17.18.2.36 `l1_neutronsobs`

```
type(optidata), dimension(:), allocatable, public mo_global_variables::l1_neutronsobs
```

Definition at line 127 of file `mo_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_mhm_read_config::mhm_read_config()`, and `mo_objective_function::objective_neutrons_kge_catchment_avg()`.



**17.18.2.37 l1\_percol**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_percol
```

Definition at line 159 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.38 l1\_pet**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_pet
```

Definition at line 103 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.39 l1\_pet\_calc**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_pet_calc
```

Definition at line 151 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_init\_states::variables\_alloc(), and mo\_init\_states::variables\_default\_init().

**17.18.2.40 l1\_pet\_weights**

```
real(dp), dimension(:, :, :), allocatable, public mo_global_variables::l1_pet_weights
```

Definition at line 99 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.41 l1\_pre**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_pre
```

Definition at line 101 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.42 l1\_pre\_weights**

```
real(dp), dimension(:, :, :), allocatable, public mo_global_variables::l1_pre_weights
```

Definition at line 100 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

#### 17.18.2.43 l1\_preeffect

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_preeffect
```

Definition at line 160 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

#### 17.18.2.44 l1\_rain

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_rain
```

Definition at line 161 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

#### 17.18.2.45 l1\_runoffseal

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_runoffseal
```

Definition at line 162 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

#### 17.18.2.46 l1\_satstw

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_satstw
```

Definition at line 145 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

#### 17.18.2.47 l1\_sealstw

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_sealstw
```

Definition at line 142 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

#### 17.18.2.48 l1\_slowrunoff

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_slowrunoff
```

Definition at line 163 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.49 l1\_sm**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_sm
```

Definition at line 116 of file mo\_global\_variables.f90.

**17.18.2.50 l1\_sm\_mask**

```
logical, dimension(:, :), allocatable, public mo_global_variables::l1_sm_mask
```

Definition at line 117 of file mo\_global\_variables.f90.

**17.18.2.51 l1\_slobs**

```
type(optidata), dimension(:), allocatable, public mo_global_variables::l1_slobs
```

Definition at line 125 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_objective\_function↔::objective\_kge\_q\_sm\_corr(), mo\_objective\_function::objective\_sm\_corr(), mo\_objective\_function↔kge\_catchment\_avg(), mo\_objective\_function::objective\_sm\_pd(), and mo\_objective\_function::objective\_sm\_sse↔\_standard\_score().

**17.18.2.52 l1\_snow**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_snow
```

Definition at line 164 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.53 l1\_snowpack**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_snowpack
```

Definition at line 141 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.54 l1\_soilmoist**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_soilmoist
```

Definition at line 143 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.55 l1\_ssrđ**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_ssrđ
```

Definition at line 110 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.56 l1\_strđ**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_strđ
```

Definition at line 111 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.57 l1\_tann**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_tann
```

Definition at line 112 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.58 l1\_temp**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_temp
```

Definition at line 102 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.59 l1\_temp\_weights**

```
real(dp), dimension(:, :, :), allocatable, public mo_global_variables::l1_temp_weights
```

Definition at line 98 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.60 l1\_throughfall**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_throughfall
```

Definition at line 165 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.61 l1\_tmax**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_tmax
```

Definition at line 105 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.62 l1\_tmin**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_tmin
```

Definition at line 104 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.63 l1\_total\_runoff**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_total_runoff
```

Definition at line 166 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.64 l1\_twsaobs**

```
type(optidata), dimension(:), allocatable, public mo_global_variables::l1_twsaobs
```

Definition at line 131 of file mo\_global\_variables.f90.

Referenced by mo\_objective\_function::create\_domain\_avg\_tws(), mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_objective\_function::objective\_kge\_q\_rmse\_tws(), and mo\_objective\_function::objective\_q\_et\_tws\_kge\_catchment\_avg().

**17.18.2.65 l1\_unsatstw**

```
real(dp), dimension(:), allocatable, public mo_global_variables::l1_unsatstw
```

Definition at line 144 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), and mo\_restart::write\_restart\_files().

**17.18.2.66 l1\_windspeed**

```
real(dp), dimension(:, :), allocatable, public mo_global_variables::l1_windspeed
```

Definition at line 108 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

### 17.18.2.67 level2

```
type(grid), dimension(:), allocatable, public mo_global_variables::level2
```

Definition at line 90 of file mo\_global\_variables.f90.

Referenced by mo\_meteo\_forcings::meteo\_forcings\_wrapper(), mo\_meteo\_forcings::meteo\_weights\_wrapper(), and mo\_startup::mhm\_initialize().

### 17.18.2.68 neutron\_integral\_afast

```
real(dp), dimension(:), allocatable, public mo_global_variables::neutron_integral_afast
```

Definition at line 190 of file mo\_global\_variables.f90.

Referenced by mo\_startup::constants\_init(), and mo\_mhm\_eval::mhm\_eval().

### 17.18.2.69 nsoilhorizons\_sm\_input

```
integer(i4) mo_global_variables::nsoilhorizons_sm_input
```

Definition at line 123 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

### 17.18.2.70 output\_deflate\_level

```
integer(i4) mo_global_variables::output_deflate_level
```

Definition at line 55 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config(), and mo\_write\_fluxes\_states::newoutputvariable().

### 17.18.2.71 output\_double\_precision

```
logical mo_global_variables::output_double_precision
```

Definition at line 56 of file mo\_global\_variables.f90.

Referenced by mo\_write\_fluxes\_states::createoutputfile(), and mo\_mhm\_read\_config::mhm\_read\_config().

### 17.18.2.72 outputflxstate

```
logical, dimension(noutflxstate) mo_global_variables::outputflxstate
```

Definition at line 58 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_write\_fluxes\_states::newoutputdataset(), and mo\_write\_fluxes\_states::updatedataset().

**17.18.2.73 read\_meteo\_weights**

```
logical, public mo_global_variables::read_meteo_weights
```

Definition at line 64 of file mo\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.74 routingstates**

```
integer(i4), parameter, public mo_global_variables::routingstates = 2
```

Definition at line 85 of file mo\_global\_variables.f90.

**17.18.2.75 timestep\_model\_inputs**

```
integer(i4), dimension(:), allocatable, public mo_global_variables::timestep_model_inputs
```

Definition at line 63 of file mo\_global\_variables.f90.

Referenced by mo\_meteo\_forcings::chunk\_size(), mo\_meteo\_forcings::is\_read(), mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), and mo\_meteo\_forcings::prepare\_meteo\_forcings\_data().

**17.18.2.76 timestep\_model\_outputs**

```
integer(i4) mo_global_variables::timestep_model_outputs
```

Definition at line 57 of file mo\_global\_variables.f90.

Referenced by mo\_write\_fluxes\_states::fluxesunit(), mo\_mhm\_eval::mhm\_eval(), and mo\_mhm\_read\_config::mhm\_read\_config().

**17.19 mo\_grid Module Reference**

TODO: add description.

**Functions/Subroutines**

- subroutine, public [init\\_lowres\\_level](#) (highres, target\_resolution, lowres, highres\_lowres\_remap)  
*Level-1 variable initialization.*
- subroutine, public [set\\_domain\\_indices](#) (grids, indices)  
*TODO: add description.*
- subroutine, public [IO\\_grid\\_setup](#) (new\_grid)  
*level 0 variable initialization*
- subroutine, public [mapcoordinates](#) (level, y, x)  
*Generate map coordinates.*
- subroutine, public [geocoordinates](#) (level, lat, lon)  
*Generate geographic coordinates.*
- subroutine [calculate\\_grid\\_properties](#) (nrowsIn, ncolsIn, xllcornerIn, yllcornerIn, cellsizeIn, aimingResolution, nrowsOut, ncolsOut, xllcornerOut, yllcornerOut, cellsizeOut)

*Calculates basic grid properties at a required coarser level using information of a given finer level. Calculates basic grid properties at a required coarser level (e.g., L11) using information of a given finer level (e.g., L0). Basic grid properties such as nrows, ncols, xllcorner, yllcorner cellsize are estimated in this routine.*

### 17.19.1 Detailed Description

TODO: add description.

TODO: add description

#### Authors

Robert Schweppe

#### Date

Jun 2018

### 17.19.2 Function/Subroutine Documentation

#### 17.19.2.1 calculate\_grid\_properties()

```
subroutine mo_grid::calculate_grid_properties (
    integer(i4), intent(in) nrowsIn,
    integer(i4), intent(in) ncolsIn,
    real(dp), intent(in) xllcornerIn,
    real(dp), intent(in) yllcornerIn,
    real(dp), intent(in) cellsizeIn,
    real(dp), intent(in) aimingResolution,
    integer(i4), intent(out) nrowsOut,
    integer(i4), intent(out) ncolsOut,
    real(dp), intent(out) xllcornerOut,
    real(dp), intent(out) yllcornerOut,
    real(dp), intent(out) cellsizeOut )
```

Calculates basic grid properties at a required coarser level using information of a given finer level. Calculates basic grid properties at a required coarser level (e.g., L11) using information of a given finer level (e.g., L0). Basic grid properties such as nrows, ncols, xllcorner, yllcorner cellsize are estimated in this routine.

TODO: add description

#### Parameters

|     |                                     |                                |
|-----|-------------------------------------|--------------------------------|
| in  | <i>integer(i4) :: nrowsIn</i>       | no. of rows at an input level  |
| in  | <i>integer(i4) :: ncolsIn</i>       | no. of cols at an input level  |
| in  | <i>real(dp) :: xllcornerIn</i>      | xllcorner at an input level    |
| in  | <i>real(dp) :: yllcornerIn</i>      | yllcorner at an input level    |
| in  | <i>real(dp) :: cellsizeIn</i>       | cell size at an input level    |
| in  | <i>real(dp) :: aimingResolution</i> | resolution of an output level  |
| out | <i>integer(i4) :: nrowsOut</i>      | no. of rows at an output level |
| out | <i>integer(i4) :: ncolsOut</i>      | no. of cols at an output level |
| out | <i>real(dp) :: xllcornerOut</i>     | xllcorner at an output level   |
| out | <i>real(dp) :: yllcornerOut</i>     | yllcorner at an output level   |
| out | <i>real(dp) :: cellsizeOut</i>      | cell size at an output level   |



**Authors**

Matthias Zink &amp; Rohini Kumar

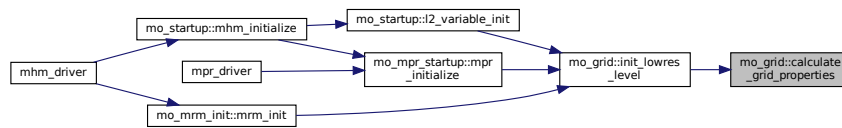
**Date**

Feb 2013

Definition at line 489 of file mo\_grid.f90.

Referenced by init\_lowres\_level().

Here is the caller graph for this function:

**17.19.2.2 geocoordinates()**

```

subroutine, public mo_grid::geocoordinates (
    type(grid), intent(in) level,
    real(dp), dimension(:, :) , intent(out), allocatable lat,
    real(dp), dimension(:, :) , intent(out), allocatable lon )
  
```

Generate geographic coordinates.

Generate geographic coordinate arrays for given domain and level

**Parameters**

|     |  |                   |
|-----|--|-------------------|
| in  | <i>type(Grid) :: level</i>                   | -> grid reference |
| out | <i>real(dp), dimension(:, :) :: lat, lon</i> |                   |
| out | <i>real(dp), dimension(:, :) :: lat, lon</i> |                   |

**Authors**

Matthias Zink

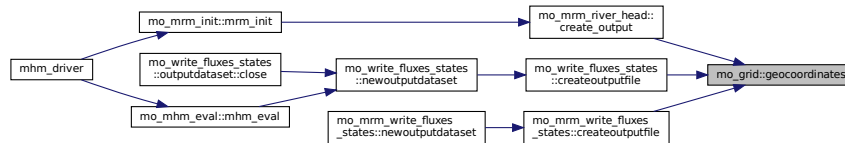
## Date

Apr 2013

Definition at line 433 of file mo\_grid.f90.

Referenced by mo\_mrm\_river\_head::create\_output(), mo\_write\_fluxes\_states::createoutputfile(), and mo\_mrm\_↔ write\_fluxes\_states::createoutputfile().

Here is the caller graph for this function:



### 17.19.2.3 init\_lowres\_level()

```

subroutine, public mo_grid::init_lowres_level (
    type(Grid), intent(in), target highres,
    real(dp), intent(in) target_resolution,
    type(Grid), intent(inout), target lowres,
    type(GridRemapper), intent(inout), optional highres_lowres_remap )
  
```

Level-1 variable initialization.

following tasks are performed for L1 datasets

- cell id & numbering
- mask creation
- storage of cell coordinates (row and column id)
- storage of four corner L0 coordinates If a variable is added or removed here, then it also has to be added or removed in the subroutine config\_variables\_set in module [mo\\_restart](#) and in the subroutine set\_config in module mo\_set\_netcdf\_restart

#### Parameters

|         |   |  |
|---------|---|--|
| in      | <i>type(Grid) :: highres</i>                                |  |
| in      | <i>real(dp) :: target_resolution</i>                        |  |
| in, out | <i>type(Grid) :: lowres</i>                                 |  |
| in, out | <i>type(GridRemapper), optional :: highres_lowres_remap</i> |  |

#### Authors

Rohini Kumar

#### Date

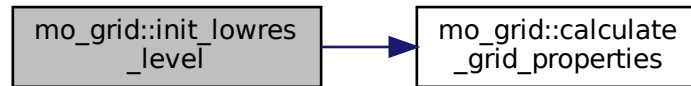
Jan 2013

Definition at line 59 of file mo\_grid.f90.

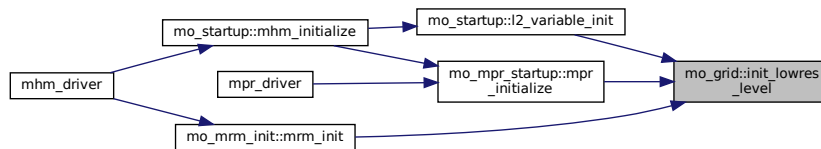
References `calculate_grid_properties()`, `mo_common_constants::nodata_dp`, and `mo_common_constants::nodata_i4`.

Referenced by `mo_startup::l2_variable_init()`, `mo_mpr_startup::mpr_initialize()`, and `mo_mrm_init::mrm_init()`.

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.19.2.4 l0\_grid\_setup()

```

subroutine, public mo_grid::l0_grid_setup (
    type(Grid), intent(inout) new_grid )
  
```

level 0 variable initialization

following tasks are performed for L0 data sets

- cell id & numbering
- storage of cell coordinates (row and column id)
- empirical dist. of terrain slope
- flag to determine the presence of a particular soil id in this configuration of the model run If a variable is added or removed here, then it also has to be added or removed in the subroutine `config_variables_set` in module `mo_restart` and in the subroutine `set_config` in module `mo_set_netcdf_restart`

#### Parameters

|         |                                     |  |
|---------|-------------------------------------|--|
| in, out | <code>type(Grid) :: new_grid</code> |  |
|---------|-------------------------------------|--|

#### Authors

Rohini Kumar

**Date**

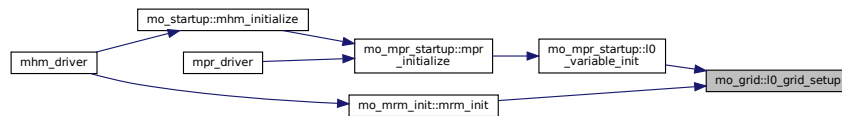
Jan 2013

Definition at line 268 of file mo\_grid.f90.

References mo\_common\_variables::iflag\_cordinate\_sys.

Referenced by mo\_mpr\_startup::!0\_variable\_init(), and mo\_mrm\_init::mrm\_init().

Here is the caller graph for this function:

**17.19.2.5 mapcoordinates()**

```

subroutine, public mo_grid::mapcoordinates (
    type(grid), intent(in) level,
    real(dp), dimension(:), intent(out), allocatable y,
    real(dp), dimension(:), intent(out), allocatable x )
  
```

Generate map coordinates.

Generate map coordinate arrays for given domain and level

**Parameters**

|     |                                       |                   |
|-----|---------------------------------------|-------------------|
| in  | <i>type(Grid) :: level</i>            | -> grid reference |
| out | <i>real(dp), dimension(:) :: x, y</i> |                   |
| out | <i>real(dp), dimension(:) :: x, y</i> |                   |

**Authors**

Matthias Zink

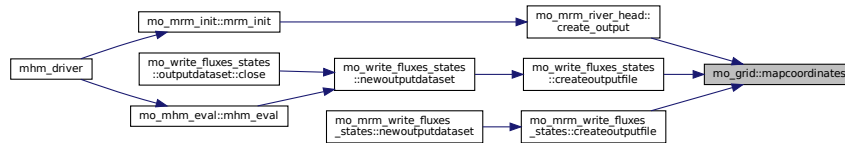
## Date

Apr 2013

Definition at line 370 of file mo\_grid.f90.

Referenced by mo\_mrm\_river\_head::create\_output(), mo\_write\_fluxes\_states::createoutputfile(), and mo\_mrm\_↔write\_fluxes\_states::createoutputfile().

Here is the caller graph for this function:



## 17.19.2.6 set\_domain\_indices()

```

subroutine, public mo_grid::set_domain_indices (
    type(grid), dimension(:), intent(inout) grids,
    integer(i4), dimension(:), intent(in), optional indices )

```

TODO: add description.

TODO: add description

## Parameters

|         |  |
|---------|--|
| in, out | <i>type(Grid), dimension(:) :: grids</i> |
|---------|--|

## Authors

Robert Schwappe

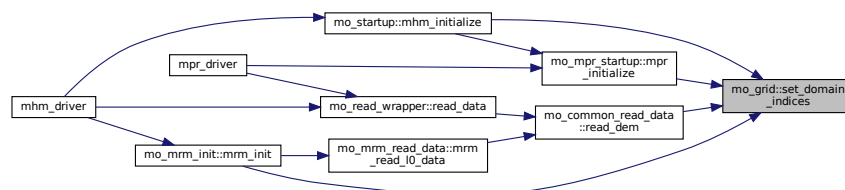
## Date

Jun 2018

Definition at line 205 of file mo\_grid.f90.

Referenced by mo\_startup::mhm\_initialize(), mo\_mpr\_startup::mpr\_initialize(), mo\_mrm\_init::mrm\_init(), and mo\_↔\_common\_read\_data::read\_dem().

Here is the caller graph for this function:



## 17.20 mo\_init\_states Module Reference

Initialization of all state variables of mHM.

### Functions/Subroutines

- subroutine, public [variables\\_alloc](#) (ncells1)  
*Allocation of space for mHM related L1 and L11 variables.*
- subroutine, public [variables\\_default\\_init](#)  
*Default initialization mHM related L1 variables.*

### 17.20.1 Detailed Description

Initialization of all state variables of mHM.

This module initializes all state variables required to run mHM. Two options are provided:

- (1) default values
- (2) from nc file

#### Authors

Luis Samaniego & Rohini Kumar

#### Date

Dec 2012

### 17.20.2 Function/Subroutine Documentation

#### 17.20.2.1 variables\_alloc()

```
subroutine, public mo_init_states::variables_alloc (
    integer(i4), intent(in) ncells1 )
```

Allocation of space for mHM related L1 and L11 variables.

Allocation of space for mHM related L1 and L11 variables (e.g., states, fluxes, and parameters) for a given domain. Variables allocated here is defined in them [mo\\_global\\_variables.f90](#) file. After allocating any variable in this routine, initialize them in the following variables\_default\_init subroutine:

#### Parameters

|    |                        |  |
|----|------------------------|--|
| in | integer(i4) :: ncells1 |  |
|----|------------------------|--|

#### Authors

Rohini Kumar

**Date**

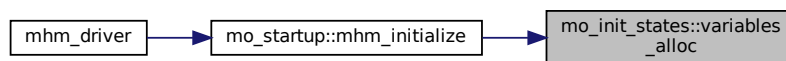
Jan 2013

Definition at line 62 of file mo\_init\_states.f90.

References mo\_mpr\_constants::c1\_initstatesm, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_global\_variables::l1\_aetcanopy, mo\_global\_variables::l1\_aetsealed, mo\_global\_variables::l1\_aetsoil, mo\_global\_variables::l1\_baseflow, mo\_global\_variables::l1\_fastrunoff, mo\_global\_variables::l1\_infilsoil, mo\_global\_variables::l1\_inter, mo\_global\_variables::l1\_melt, mo\_global\_variables::l1\_neutrons, mo\_global\_variables::l1\_percol, mo\_global\_variables::l1\_pet\_calc, mo\_global\_variables::l1\_preeffect, mo\_global\_variables::l1\_rain, mo\_global\_variables::l1\_runoffseal, mo\_global\_variables::l1\_satstw, mo\_global\_variables::l1\_sealstw, mo\_global\_variables::l1\_slowrunoff, mo\_global\_variables::l1\_snow, mo\_global\_variables::l1\_snowpack, mo\_global\_variables::l1\_soilmoist, mo\_global\_variables::l1\_throughfall, mo\_global\_variables::l1\_total\_runoff, mo\_global\_variables::l1\_unsatstw, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_common\_constants::p1\_initstatefluxes, mo\_mpr\_constants::p2\_initstatefluxes, mo\_mpr\_constants::p3\_initstatefluxes, mo\_mpr\_constants::p4\_initstatefluxes, and mo\_mpr\_constants::p5\_initstatefluxes.

Referenced by mo\_startup::mhm\_initialize().

Here is the caller graph for this function:

**17.20.2.2 variables\_default\_init()**

```
subroutine, public mo_init_states::variables_default_init
```

Default initialization mHM related L1 variables.

Default initialization of mHM related L1 variables (e.g., states, fluxes, and parameters) as per given constant values given in [mo\\_mhm\\_constants](#). Variables initialized here is defined in the [mo\\_global\\_variables.f90](#) file. Only Variables that are defined in the variables\_alloc subroutine are initialized here. If a variable is added or removed here, then it also has to be added or removed in the subroutine state\_variables\_set in the module [mo\\_restart](#) and in the subroutine set\_state in the module [mo\\_set\\_netcdf\\_restart](#).

**Authors**

R. Kumar &amp; J. Mai

**Date**

Sep 2013

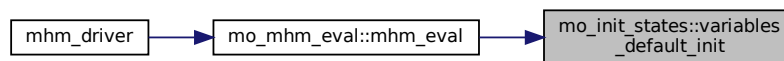
Definition at line 192 of file mo\_init\_states.f90.

References mo\_mpr\_constants::c1\_initstatesm, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_mpr\_global\_variables::l1\_aeroresist, mo\_global\_variables::l1\_aetcanopy, mo\_global\_variables::l1\_aetsealed, mo\_global\_variables::l1\_aetsoil, mo\_mpr\_global\_variables::l1\_alpha, mo\_global\_variables::l1\_baseflow, mo\_mpr\_global\_variables::l1\_degday, mo\_mpr\_global\_variables::l1\_degdayinc, mo\_mpr\_global\_variables::l1\_degdaymax, mo\_mpr\_global\_variables::l1\_degdaynopre, mo\_mpr\_global\_variables::l1\_fasp, mo\_global\_variables::l1\_fastrunoff,

mo\_mpr\_global\_variables::l1\_froots, mo\_mpr\_global\_variables::l1\_fsealed, mo\_mpr\_global\_variables::l1\_harsamcoeff, mo\_global\_variables::l1\_infilsoil, mo\_global\_variables::l1\_inter, mo\_mpr\_global\_variables::l1\_jarvis\_thresh\_c1, mo\_mpr\_global\_variables::l1\_karstloss, mo\_mpr\_global\_variables::l1\_kbaseflow, mo\_mpr\_global\_variables::l1\_kfastflow, mo\_mpr\_global\_variables::l1\_kperco, mo\_mpr\_global\_variables::l1\_kslowflow, mo\_mpr\_global\_variables::l1\_maxinter, mo\_global\_variables::l1\_melt, mo\_global\_variables::l1\_neutrons, mo\_global\_variables::l1\_percol, mo\_global\_variables::l1\_pet\_calc, mo\_mpr\_global\_variables::l1\_petlaicorfactor, mo\_global\_variables::l1\_preeffect, mo\_mpr\_global\_variables::l1\_prietayalpha, mo\_global\_variables::l1\_rain, mo\_global\_variables::l1\_runoffseal, mo\_global\_variables::l1\_satstw, mo\_mpr\_global\_variables::l1\_sealedthresh, mo\_global\_variables::l1\_sealstw, mo\_global\_variables::l1\_slowrunoff, mo\_global\_variables::l1\_snow, mo\_global\_variables::l1\_snowpack, mo\_global\_variables::l1\_soilmoist, mo\_mpr\_global\_variables::l1\_soilmoistexp, mo\_mpr\_global\_variables::l1\_soilmoistfc, mo\_mpr\_global\_variables::l1\_soilmoistsat, mo\_mpr\_global\_variables::l1\_surfresist, mo\_mpr\_global\_variables::l1\_tempthresh, mo\_global\_variables::l1\_throughfall, mo\_global\_variables::l1\_total\_runoff, mo\_global\_variables::l1\_unsatstw, mo\_mpr\_global\_variables::l1\_unsatthresh, mo\_mpr\_global\_variables::l1\_wiltingpoint, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_common\_constants::p1\_initstatefluxes, mo\_mpr\_constants::p2\_initstatefluxes, mo\_mpr\_constants::p3\_initstatefluxes, mo\_mpr\_constants::p4\_initstatefluxes, and mo\_mpr\_constants::p5\_initstatefluxes.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the caller graph for this function:



## 17.21 mo\_meteo\_forcings Module Reference

Prepare meteorological forcings data for mHM.

### Functions/Subroutines

- subroutine, public [prepare\\_meteo\\_forcings\\_data](#) (iDomain, domainID, tt)
 

*Prepare meteorological forcings data for a given variable.*
- subroutine [meteo\\_forcings\\_wrapper](#) (iDomain, dataPath, inputFormat, dataOut1, lower, upper, nvarName)
 

*Prepare meteorological forcings data for mHM at Level-1.*
- subroutine [meteo\\_weights\\_wrapper](#) (iDomain, read\_meteo\_weights, dataPath, dataOut1, lower, upper, nvarName)
 

*Prepare weights for meteorological forcings data for mHM at Level-1.*
- subroutine [chunk\\_config](#) (iDomain, tt, read\_flag, readPer)
 

*determines the start date, end date, and read\_flag given Domain id and current timestep*
- logical function [is\\_read](#) (iDomain, tt)
 

*evaluate whether new chunk should be read at this timestep*
- subroutine [chunk\\_size](#) (iDomain, tt, readPer)
 

*calculate beginning and end of read Period, i.e. that is length of current chunk to read*

### 17.21.1 Detailed Description

Prepare meteorological forcings data for mHM.

Prepare meteorological forcings data for mHM.



**Authors**

Rohini Kumar

**Date**

Jan 2012

**17.21.2 Function/Subroutine Documentation****17.21.2.1 chunk\_config()**

```

subroutine mo_meteo_forcings::chunk_config (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) tt,
    logical, intent(out) read_flag,
    type(period), intent(out) readPer )

```

determines the start date, end date, and read\_flag given Domain id and current timestep

TODO: add description

**Parameters**

|     |                                |  |
|-----|--------------------------------|--|
| in  | <i>integer(i4) :: iDomain</i>  | current Domain                               |
| in  | <i>integer(i4) :: tt</i>       | current timestep                             |
| out | <i>logical :: read_flag</i>    | indicate whether reading data should be read |
| out | <i>type(period) :: readPer</i> | start and end dates of reading Period        |

**Authors**

Stephan Thober

**Date**

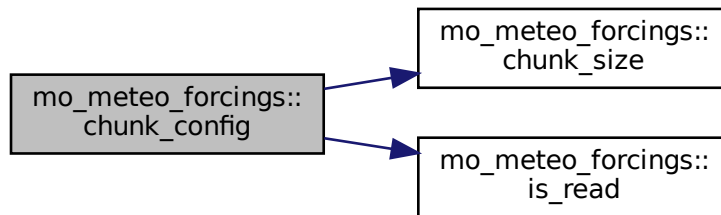
Jun 2014

Definition at line 584 of file mo\_meteo\_forcings.f90.

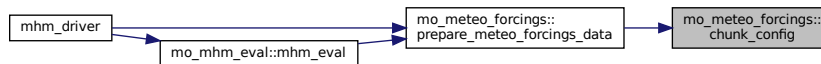
References chunk\_size(), is\_read(), and mo\_common\_constants::nodata\_dp.

Referenced by prepare\_meteo\_forcings\_data().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.21.2.2 chunk\_size()

```

subroutine mo_meteo_forcings::chunk_size (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) tt,
    type(period), intent(out) readPer )
  
```

calculate beginning and end of read Period, i.e. that is length of current chunk to read

TODO: add description

#### Parameters

|     |                                |                                    |
|-----|--------------------------------|------------------------------------|
| in  | <i>integer(i4) :: iDomain</i>  | current Domain to process          |
| in  | <i>integer(i4) :: tt</i>       | current time step                  |
| out | <i>type(period) :: readPer</i> | start and end dates of read Period |

#### Authors

Stephan Thober

#### Date

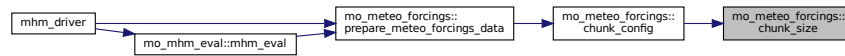
Jun 2014

Definition at line 759 of file mo\_meteo\_forcings.f90.

References mo\_common\_mhm\_mrm\_variables::ntstepday, mo\_common\_mhm\_mrm\_variables::simper, and mo←\_global\_variables::timestep\_model\_inputs.

Referenced by chunk\_config().

Here is the caller graph for this function:



### 17.21.2.3 is\_read()

```

logical function mo_meteo_forcings::is_read (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) tt )
  
```

evaluate whether new chunk should be read at this timestep

TODO: add description

#### Parameters

|    |                               |                   |
|----|-------------------------------|-------------------|
| in | <i>integer(i4) :: iDomain</i> | current Domain    |
| in | <i>integer(i4) :: tt</i>      | current time step |

#### Authors

Stephan Thober

#### Date

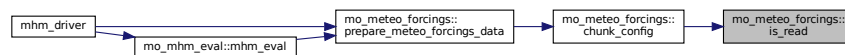
Jun 2014

Definition at line 648 of file mo\_meteo\_forcings.f90.

References mo\_common\_mhm\_mrm\_variables::ntstepday, mo\_common\_mhm\_mrm\_variables::simper, mo\_↔  
common\_mhm\_mrm\_variables::timestep, and mo\_global\_variables::timestep\_model\_inputs.

Referenced by chunk\_config().

Here is the caller graph for this function:



### 17.21.2.4 meteo\_forcings\_wrapper()

```

subroutine mo_meteo_forcings::meteo_forcings_wrapper (
    integer(i4), intent(in) iDomain,
    character(len = *), intent(in) dataPath,
  
```

```

character(len = *), intent(in) inputFormat,
real(dp), dimension(:, :), intent(inout), allocatable dataOut1,
real(dp), intent(in), optional lower,
real(dp), intent(in), optional upper,
character(len = *), intent(in), optional nvarName )

```

Prepare meteorological forcings data for mHM at Level-1.

Prepare meteorological forcings data for mHM, which include 1) Reading meteo. datasets at their native resolution for every Domain 2) Perform aggregation or disaggregation of meteo. datasets from their native resolution (level-2) to the required hydrologic resolution (level-1) 3) Pad the above datasets of every Domain to their respective global ones

#### Parameters

|         |   |  |
|---------|---|--|
| in      | <i>integer(i4) :: iDomain</i>                   | Domain Id  |
| in      | <i>character(len = *) :: dataPath</i>           | Data path where a given meteo. variable is stored              |
| in      | <i>character(len = *) :: inputFormat</i>        | only 'nc' possible at the moment                               |
| in, out | <i>real(dp), dimension(:, :) :: dataOut1</i>    | Packed meteorological variable for the whole simulation period |
| in      | <i>real(dp), optional :: lower</i>              | Lower bound for check of validity of data values               |
| in      | <i>real(dp), optional :: upper</i>              | Upper bound for check of validity of data values               |
| in      | <i>character(len = *), optional :: nvarName</i> | name of the variable (for .nc files)                           |

#### Authors

Rohini Kumar

#### Date

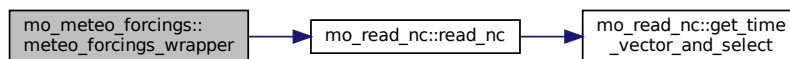
Jan 2013

Definition at line 267 of file mo\_meteo\_forcings.f90.

References mo\_common\_variables::level1, mo\_global\_variables::level2, mo\_common\_constants::nodata\_dp, mo\_read\_nc::read\_nc(), and mo\_common\_mhm\_mrm\_variables::readper.

Referenced by prepare\_meteo\_forcings\_data().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.21.2.5 meteo\_weights\_wrapper()

```

subroutine mo_meteo_forcings::meteo_weights_wrapper (
    integer(i4), intent(in) iDomain,
    logical, intent(in) read_meteo_weights,
    character(len = *), intent(in) dataPath,
    real(dp), dimension(:, :, :), intent(inout), allocatable dataOut1,
    real(dp), intent(in), optional lower,
    real(dp), intent(in), optional upper,
    character(len = *), intent(in), optional nvarName )

```

Prepare weights for meteorological forcings data for mHM at Level-1.

Prepare meteorological weights data for mHM, which include 1) Reading meteo. weights datasets at their native resolution for every Domain 2) Perform aggregation or disaggregation of meteo. weights datasets from their native resolution (level-2) to the required hydrologic resolution (level-1) 3) Pad the above datasets of every Domain to their respective global ones

## Parameters

|         |   |  |
|---------|---|--|
| in      | <i>integer(i4) :: iDomain</i>                   | Domain Id  |
| in      | <i>logical :: read_meteo_weights</i>            | Flag for reading meteo weights                                 |
| in      | <i>character(len = *) :: dataPath</i>           | Data path where a given meteo. variable is stored              |
| in, out | <i>real(dp), dimension(:, :, :) :: dataOut1</i> | Packed meteorological variable for the whole simulation period |
| in      | <i>real(dp), optional :: lower</i>              | Lower bound for check of validity of data values               |
| in      | <i>real(dp), optional :: upper</i>              | Upper bound for check of validity of data values               |
| in      | <i>character(len = *), optional :: nvarName</i> | name of the variable (for .nc files)                           |

## Authors

Stephan Thober & Rohini Kumar

## Date

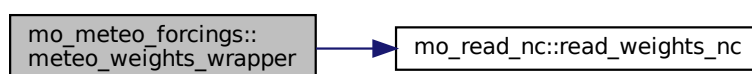
Jan 2017

Definition at line 432 of file mo\_meteo\_forcings.f90.

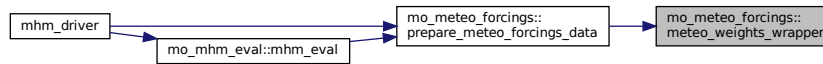
References mo\_common\_variables::level1, mo\_global\_variables::level2, mo\_common\_constants::nodata\_dp, and mo\_read\_nc::read\_weights\_nc().

Referenced by prepare\_meteo\_forcings\_data().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.21.2.6 prepare\_meteo\_forcings\_data()

```

subroutine, public mo_meteo_forcings::prepare_meteo_forcings_data (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) domainID,
    integer(i4), intent(in) tt )
  
```

Prepare meteorological forcings data for a given variable.

Prepare meteorological forcings data for a given variable. Internally this subroutine calls another routine `meteo_weights_wrapper` for different meteorological variables

#### Parameters

|    |                               |                  |
|----|-------------------------------|------------------|
| in | <i>integer(i4) :: iDomain</i> | Domain Id        |
| in | <i>integer(i4) :: tt</i>      | current timestep |

#### Authors

Rohini Kumar

#### Date

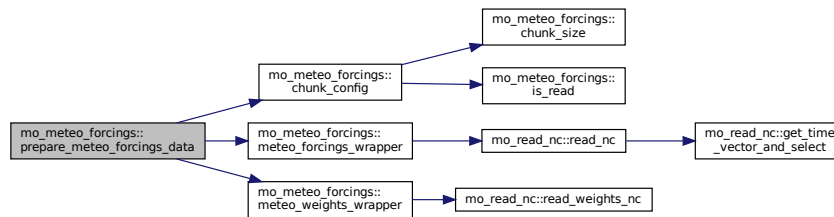
Jan 2013

Definition at line 63 of file `mo_meteo_forcings.f90`.

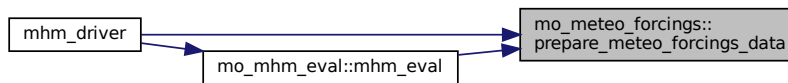
References `chunk_config()`, `mo_global_variables::dirabsvappressure`, `mo_global_variables::dirmaxtemperature`, `mo_global_variables::dirmintemperature`, `mo_global_variables::dirnetradiation`, `mo_global_variables::dirprecipitation`, `mo_global_variables::dirradiation`, `mo_global_variables::dirreferenceet`, `mo_global_variables::dirtemperature`, `mo_global_variables::dirwindspeed`, `mo_common_variables::domainmeta`, `mo_global_variables::inputformat_meteo_forcings`, `mo_global_variables::l1_absvappress`, `mo_global_variables::l1_netrad`, `mo_global_variables::l1_pet`, `mo_global_variables::l1_pet_weights`, `mo_global_variables::l1_pre`, `mo_global_variables::l1_pre_weights`, `mo_global_variables::l1_ssr`, `mo_global_variables::l1_strd`, `mo_global_variables::l1_tann`, `mo_global_variables::l1_temp`, `mo_global_variables::l1_temp_weights`, `mo_global_variables::l1_tmax`, `mo_global_variables::l1_tmin`, `mo_global_variables::l1_windspeed`, `meteo_forcings_wrapper()`, `meteo_weights_wrapper()`, `mo_common_variables::processmatrix`, `mo_global_variables::read_meteo_weights`, `mo_common_mhm_mrm_variables::readper`, and `mo_global_variables::timestep_model_inputs`.

Referenced by `mhm_driver()`, and `mo_mhm_eval::mhm_eval()`.

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.22 mo\_mhm Module Reference

Call all main processes of mHM.

### Functions/Subroutines

- subroutine, public **mhm** (read\_states, tt, time, processMatrix, horizon\_depth, nCells1, nHorizons\_mHM, nimesteps\_day, c2TSTu, neutron\_integral\_AFast, global\_parameters, latitude, evap\_coeff, fday\_prec, fnight\_prec, fday\_pet, fnight\_pet, fday\_temp, fnight\_temp, temp\_weights, pet\_weights, pre\_weights, read\_meteo\_weights, pet\_in, tmin\_in, tmax\_in, netrad\_in, absvappres\_in, windspeed\_in, prec\_in, temp\_in, fSealed1, interc, snowpack, sealedStorage, soilMoisture, unsatStorage, satStorage, neutrons, pet\_calc, aet\_soil, aet\_canopy, aet\_sealed, baseflow, infiltration, fast\_interflow, melt, perc, prec\_effect, rain, runoff\_sealed, slow\_interflow, snow, throughfall, total\_runoff, alpha, deg\_day\_incr, deg\_day\_max, deg\_day\_noprec, deg\_day, fAsp, petLAIcorFactorL1, HarSamCoeff, PrieTayAlpha, aeroResist, surfResist, frac\_roots, interc\_max, karst\_loss, k0, k1, k2, kp, soil\_moist\_FC, soil\_moist\_sat, soil\_moist\_exponen, jarvis\_thresh\_c1, temp\_thresh, unsat\_thresh, water\_thresh\_sealed, wilting\_point)

*Pure mHM calculations.*

### 17.22.1 Detailed Description

Call all main processes of mHM.

This module calls all processes of mHM for a given configuration. The configuration of the model is stored in the a process matrix. This configuration is specified in the namelist mhm.nml. The processes are executed in ascending order. At the moment only process 5 and 8 have options. Currently the following processes are implemented:

| Process | Name             | Flag | Description   |
|---------|------------------|------|---|
| 1       | interception     | 1    | Maximum interception                                |
| 2       | snow and melting | 1    | Degree-day  |
| 3       | soil moisture    | 1    | Feddes equation for ET reduction, Brooks-Corey like |

| Process | Name                      | Flag | Description  |
|---------|---------------------------|------|--|
| 3       | soil moisture             | 2    | Jarvis equation for ET reduction, Brooks-Corey like        |
| 3       | soil moisture             | 3    | Jarvis eq. for ET red. + FC dependency on root frac. coef. |
| 4       | direct runoff             | 1    | Linear reservoir exceedance                                |
| 5       | PET                       | -1   | PET is input, LAI based correction, dynamic scaling func.  |
| 5       | PET                       | 0    | PET is input, Aspect based correction                      |
| 5       | PET                       | 1    | Hargreaves-Samani  |
| 5       | PET                       | 2    | Priestley-Taylor   |
| 5       | PET                       | 3    | Penman-Monteith  |
| 6       | interflow                 | 1    | Nonlinear reservoir with saturation excess                 |
| 7       | percolation and base flow | 1    | GW linear reservoir  |
| 8       | routing                   | 0    | no routing   |
| 8       | routing                   | 1    | use mRM i.e. Muskingum                                     |
| 8       | routing                   | 2    | use mRM i.e. adaptive timestep                             |

### Authors

Luis Samaniego

### Date

Dec 2012

## 17.22.2 Function/Subroutine Documentation

### 17.22.2.1 mhm()

```

subroutine, public mo_mhm::mhm (
    logical, intent(in) read_states,
    integer(i4), intent(in) tt,
    real(dp), intent(in) time,
    integer(i4), dimension(:, :), intent(in) processMatrix,
    real(dp), dimension(:), intent(in) horizon_depth,
    integer(i4), intent(in) nCells1,
    integer(i4), intent(in) nHorizons_mHM,
    real(dp), intent(in) ntimesteps_day,
    real(dp), intent(in) c2TSTu,
    real(dp), dimension(:), intent(in) neutron_integral_AFast,
    real(dp), dimension(:), intent(in) global_parameters,
    real(dp), dimension(:), intent(in) latitude,
    real(dp), dimension(:), intent(in) evap_coeff,
    real(dp), dimension(:), intent(in) fday_prec,
    real(dp), dimension(:), intent(in) fnight_prec,
    real(dp), dimension(:), intent(in) fday_pet,
    real(dp), dimension(:), intent(in) fnight_pet,
    real(dp), dimension(:), intent(in) fday_temp,
    real(dp), dimension(:), intent(in) fnight_temp,
    real(dp), dimension(:, :, :), intent(in) temp_weights,
    real(dp), dimension(:, :, :), intent(in) pet_weights,
    real(dp), dimension(:, :, :), intent(in) pre_weights,
    logical, intent(in) read_meteo_weights,
    real(dp), dimension(:), intent(in) pet_in,

```



```

real(dp), dimension(:), intent(in) tmin_in,
real(dp), dimension(:), intent(in) tmax_in,
real(dp), dimension(:), intent(in) netrad_in,
real(dp), dimension(:), intent(in) absvappres_in,
real(dp), dimension(:), intent(in) windspeed_in,
real(dp), dimension(:), intent(in) prec_in,
real(dp), dimension(:), intent(in) temp_in,
real(dp), dimension(:), intent(inout) fSealed1,
real(dp), dimension(:), intent(inout) interc,
real(dp), dimension(:), intent(inout) snowpack,
real(dp), dimension(:), intent(inout) sealedStorage,
real(dp), dimension(:, :), intent(inout) soilMoisture,
real(dp), dimension(:), intent(inout) unsatStorage,
real(dp), dimension(:), intent(inout) satStorage,
real(dp), dimension(:), intent(inout) neutrons,
real(dp), dimension(:), intent(inout) pet_calc,
real(dp), dimension(:, :), intent(inout) aet_soil,
real(dp), dimension(:), intent(inout) aet_canopy,
real(dp), dimension(:), intent(inout) aet_sealed,
real(dp), dimension(:), intent(inout) baseflow,
real(dp), dimension(:, :), intent(inout) infiltration,
real(dp), dimension(:), intent(inout) fast_interflow,
real(dp), dimension(:), intent(inout) melt,
real(dp), dimension(:), intent(inout) perc,
real(dp), dimension(:), intent(inout) prec_effect,
real(dp), dimension(:), intent(inout) rain,
real(dp), dimension(:), intent(inout) runoff_sealed,
real(dp), dimension(:), intent(inout) slow_interflow,
real(dp), dimension(:), intent(inout) snow,
real(dp), dimension(:), intent(inout) throughfall,
real(dp), dimension(:), intent(inout) total_runoff,
real(dp), dimension(:), intent(inout) alpha,
real(dp), dimension(:), intent(inout) deg_day_incr,
real(dp), dimension(:), intent(inout) deg_day_max,
real(dp), dimension(:), intent(inout) deg_day_noprec,
real(dp), dimension(:), intent(inout) deg_day,
real(dp), dimension(:), intent(inout) fAsp,
real(dp), dimension(:), intent(inout) petLAIcorFactorL1,
real(dp), dimension(:), intent(inout) HarSamCoeff,
real(dp), dimension(:), intent(inout) PrieTayAlpha,
real(dp), dimension(:), intent(inout) aeroResist,
real(dp), dimension(:), intent(inout) surfResist,
real(dp), dimension(:, :), intent(inout) frac_roots,
real(dp), dimension(:), intent(inout) interc_max,
real(dp), dimension(:), intent(inout) karst_loss,
real(dp), dimension(:), intent(inout) k0,
real(dp), dimension(:), intent(inout) k1,
real(dp), dimension(:), intent(inout) k2,
real(dp), dimension(:), intent(inout) kp,
real(dp), dimension(:, :), intent(inout) soil_moist_FC,
real(dp), dimension(:, :), intent(inout) soil_moist_sat,
real(dp), dimension(:, :), intent(inout) soil_moist_exponen,
real(dp), dimension(:), intent(inout) jarvis_thresh_c1,
real(dp), dimension(:), intent(inout) temp_thresh,
real(dp), dimension(:), intent(inout) unsat_thresh,
real(dp), dimension(:), intent(inout) water_thresh_sealed,
real(dp), dimension(:, :), intent(inout) wilting_point )

```

Pure mHM calculations.

Pure mHM calculations. All variables are allocated and initialized. They will be local variables within this call.

### Parameters

|         |   |   |
|---------|---|---|
| in      | <i>logical :: read_states</i>                           | indicated whether states have been read from file                       |
| in      | <i>integer(i4) :: tt</i>                                | simulation time step  |
| in      | <i>real(dp) :: time</i>                                 | current decimal Julian day  |
| in      | <i>integer(i4), dimension(:, :) :: processMatrix</i>    | mHM process configuration matrix  |
| in      | <i>real(dp), dimension(:) :: horizon_depth</i>          | Depth of each horizon in mHM  |
| in      | <i>integer(i4) :: nCells1</i>                           | number of cells in a given domain at level L1                           |
| in      | <i>integer(i4) :: nHorizons_mHM</i>                     | Number of Horizons in mHM   |
| in      | <i>real(dp) :: ntimesteps_day</i>                       | number of time intervals per day, transformed in dp                     |
| in      | <i>real(dp), dimension(:) :: neutron_integral_AFast</i> | tabular for neutron flux approximation                                  |
| in      | <i>real(dp), dimension(:) :: global_parameters</i>      | global mHM parameters   |
| in      | <i>real(dp), dimension(:) :: latitude</i>               | latitude on level 1   |
| in      | <i>real(dp), dimension(:) :: evap_coeff</i>             | Evaporation coefficient for free-water surface of that current month    |
| in      | <i>real(dp), dimension(:) :: fday_prec</i>              | [-] day ratio precipitation < 1   |
| in      | <i>real(dp), dimension(:) :: fnight_prec</i>            | [-] night ratio precipitation < 1                                       |
| in      | <i>real(dp), dimension(:) :: fday_pet</i>               | [-] day ratio PET < 1   |
| in      | <i>real(dp), dimension(:) :: fnight_pet</i>             | [-] night ratio PET < 1   |
| in      | <i>real(dp), dimension(:) :: fday_temp</i>              | [-] day factor mean temp  |
| in      | <i>real(dp), dimension(:) :: fnight_temp</i>            | [-] night factor mean temp  |
| in      | <i>real(dp), dimension(:, :, :) :: temp_weights</i>     | multiplicative weights for temperature (deg K)                          |
| in      | <i>real(dp), dimension(:, :, :) :: pet_weights</i>      | multiplicative weights for potential evapotranspiration                 |
| in      | <i>real(dp), dimension(:, :, :) :: pre_weights</i>      | multiplicative weights for precipitation                                |
| in      | <i>logical :: read_meteo_weights</i>                    | flag whether weights for tavg and pet have read and should be used      |
| in      | <i>real(dp), dimension(:) :: pet_in</i>                 | [mm d-1] Daily potential evapotranspiration (input)                     |
| in      | <i>real(dp), dimension(:) :: tmin_in</i>                | [degc] Daily minimum temperature  |
| in      | <i>real(dp), dimension(:) :: tmax_in</i>                | [degc] Daily maximum temperature  |
| in      | <i>real(dp), dimension(:) :: netrad_in</i>              | [w m2] Daily average net radiation                                      |
| in      | <i>real(dp), dimension(:) :: absvappres_in</i>          | [Pa] Daily average absolute vapour pressure                             |
| in      | <i>real(dp), dimension(:) :: windspeed_in</i>           | [m s-1] Daily average wind speed  |
| in      | <i>real(dp), dimension(:) :: prec_in</i>                | [mm d-1] Daily mean precipitation                                       |
| in      | <i>real(dp), dimension(:) :: temp_in</i>                | [degc] Daily average temperature  |
| in, out | <i>real(dp), dimension(:) :: fSealed1</i>               | fraction of sealed area at scale L1                                     |
| in, out | <i>real(dp), dimension(:) :: interc</i>                 | Interception  |
| in, out | <i>real(dp), dimension(:) :: snowpack</i>               | Snowpack  |
| in, out | <i>real(dp), dimension(:) :: sealedStorage</i>          | Retention storage of impervious areas                                   |
| in, out | <i>real(dp), dimension(:, :) :: soilMoisture</i>        | Soil moisture of each horizon   |
| in, out | <i>real(dp), dimension(:) :: unsatStorage</i>           | Upper soil storage  |
| in, out | <i>real(dp), dimension(:) :: satStorage</i>             | Groundwater storage   |
| in, out | <i>real(dp), dimension(:) :: neutrons</i>               | Ground albedo neutrons  |
| in, out | <i>real(dp), dimension(:) :: pet_calc</i>               | [mm TS-1] estimated PET (if PET is input = corrected values (fAsp*PET)) |

## Parameters

|         |  |   |
|---------|--|---|
| in, out | <i>real(dp), dimension(:, :) :: aet_soil</i>           | actual ET   |
| in, out | <i>real(dp), dimension(:) :: aet_canopy</i>            | Real evaporation intensity from canopy                                      |
| in, out | <i>real(dp), dimension(:) :: aet_sealed</i>            | Actual ET from free-water surfaces  |
| in, out | <i>real(dp), dimension(:) :: baseflow</i>              | Baseflow  |
| in, out | <i>real(dp), dimension(:, :) :: infiltration</i>       | Recharge, infiltration intensity or effective precipitation of each horizon |
| in, out | <i>real(dp), dimension(:) :: fast_interflow</i>        | Fast runoff component   |
| in, out | <i>real(dp), dimension(:) :: melt</i>                  | Melting snow depth  |
| in, out | <i>real(dp), dimension(:) :: perc</i>                  | Percolation   |
| in, out | <i>real(dp), dimension(:) :: prec_effect</i>           | Effective precipitation depth (snow melt + rain)                            |
| in, out | <i>real(dp), dimension(:) :: rain</i>                  | Rain precipitation depth  |
| in, out | <i>real(dp), dimension(:) :: runoff_sealed</i>         | Direct runoff from impervious areas   |
| in, out | <i>real(dp), dimension(:) :: slow_interflow</i>        | Slow runoff component   |
| in, out | <i>real(dp), dimension(:) :: snow</i>                  | Snow precipitation depth  |
| in, out | <i>real(dp), dimension(:) :: throughfall</i>           | Throughfall   |
| in, out | <i>real(dp), dimension(:) :: total_runoff</i>          | Generated runoff  |
| in, out | <i>real(dp), dimension(:) :: alpha</i>                 | Exponent for the upper reservoir  |
| in, out | <i>real(dp), dimension(:) :: deg_day_incr</i>          | Increase of the Degree-day factor per mm of increase in precipitation       |
| in, out | <i>real(dp), dimension(:) :: deg_day_max</i>           | Maximum Degree-day factor   |
| in, out | <i>real(dp), dimension(:) :: deg_day_noprec</i>        | Degree-day factor with no precipitation                                     |
| in, out | <i>real(dp), dimension(:) :: deg_day</i>               | Degree-day factor   |
| in, out | <i>real(dp), dimension(:) :: fAsp</i>                  | [1] PET correction for Aspect at level 1                                    |
| in, out | <i>real(dp), dimension(:) :: petLAIcorFactorL1</i>     | PET correction factor based on LAI at level 1                               |
| in, out | <i>real(dp), dimension(:) :: HarSamCoeff</i>           | [1] PET Hargreaves Samani coefficient at level 1                            |
| in, out | <i>real(dp), dimension(:) :: PrieTayAlpha</i>          | [1] PET Priestley Taylor coefficient at level 1                             |
| in, out | <i>real(dp), dimension(:) :: aeroResist</i>            | [s m <sup>-1</sup> ] PET aerodynamical resistance at level 1                |
| in, out | <i>real(dp), dimension(:) :: surfResist</i>            | [s m <sup>-1</sup> ] PET bulk surface resistance at level 1                 |
| in, out | <i>real(dp), dimension(:, :) :: frac_roots</i>         | Fraction of Roots in soil horizon   |
| in, out | <i>real(dp), dimension(:) :: interc_max</i>            | Maximum interception  |
| in, out | <i>real(dp), dimension(:) :: karst_loss</i>            | Karstic percolation loss  |
| in, out | <i>real(dp), dimension(:) :: k0</i>                    | Recession coefficient of the upper reservoir, upper outlet                  |
| in, out | <i>real(dp), dimension(:) :: k1</i>                    | Recession coefficient of the upper reservoir, lower outlet                  |
| in, out | <i>real(dp), dimension(:) :: k2</i>                    | Baseflow recession coefficient  |
| in, out | <i>real(dp), dimension(:) :: kp</i>                    | Percolation coefficient   |
| in, out | <i>real(dp), dimension(:, :) :: soil_moist_FC</i>      | Soil moisture below which actual ET is reduced                              |
| in, out | <i>real(dp), dimension(:, :) :: soil_moist_sat</i>     | Saturation soil moisture for each horizon [mm]                              |
| in, out | <i>real(dp), dimension(:, :) :: soil_moist_exponen</i> | Exponential parameter to how non-linear is the soil water retention         |
| in, out | <i>real(dp), dimension(:) :: jarvis_thresh_c1</i>      | jarvis critical value for normalized soil water content                     |
| in, out | <i>real(dp), dimension(:) :: temp_thresh</i>           | Threshold temperature for snow/rain   |
| in, out | <i>real(dp), dimension(:) :: unsat_thresh</i>          | Threshold water depth in upper reservoir                                    |
| in, out | <i>real(dp), dimension(:) :: water_thresh_sealed</i>   | Threshold water depth in impervious areas                                   |
| in, out | <i>real(dp), dimension(:, :) :: wilting_point</i>      | Permanent wilting point for each horizon                                    |

## Authors

Luis Samaniego & Rohini Kumar

## Date

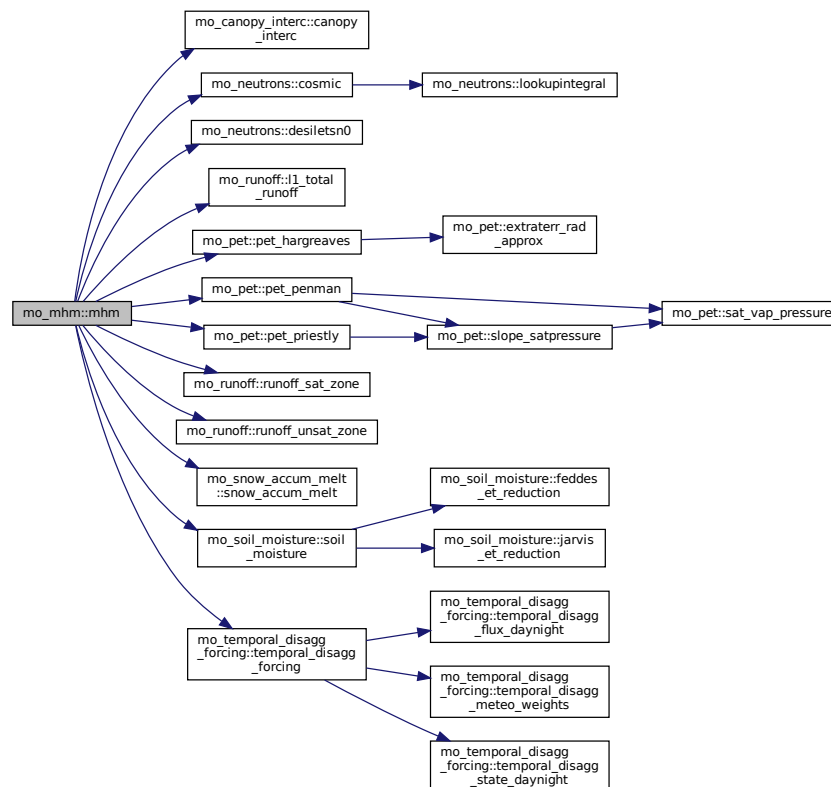
Dec 2012

Definition at line 200 of file mo\_mhm.f90.

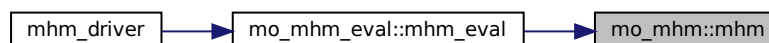
References mo\_canopy\_interc::canopy\_interc(), mo\_neutrons::cosmic(), mo\_neutrons::desiletsn0(), mo\_mhm↔\_constants::harsamconst, mo\_runoff::l1\_total\_runoff(), mo\_pet::pet\_hargreaves(), mo\_pet::pet\_penman(), mo↔\_pet::pet\_priestly(), mo\_runoff::runoff\_sat\_zone(), mo\_runoff::runoff\_unsat\_zone(), mo\_snow\_accum\_melt::snow↔\_accum\_melt(), mo\_soil\_moisture::soil\_moisture(), and mo\_temporal\_disagg\_forcing::temporal\_disagg\_forcing().

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.23 mo\_mhm\_constants Module Reference

Provides mHM specific constants.

### Variables

- real(dp), parameter, public `h2odens` = 1000.0\_dp
- real(dp), parameter, public `p2_initstatefluxes` = 15.00\_dp
- real(dp), parameter, public `p3_initstatefluxes` = 10.00\_dp
- real(dp), parameter, public `p4_initstatefluxes` = 75.00\_dp
- real(dp), parameter, public `p5_initstatefluxes` = 1500.00\_dp
- real(dp), parameter, public `c1_initstatesm` = 0.25\_dp
- integer(i4), parameter, public `noutfluxstate` = 20\_i4
- real(dp), parameter, public `harsamconst` = 17.800\_dp  
*Hargreaves-Samani ref. ET formula [deg C].*
- real(dp), parameter, public `duffiedr` = 0.0330\_dp
- real(dp), parameter, public `duffiedelta1` = 0.4090\_dp
- real(dp), parameter, public `duffiedelta2` = 1.3900\_dp
- real(dp), parameter, public `tetens_c1` = 0.6108\_dp  
*Tetens's formula to calculate saturated vapour pressure.*
- real(dp), parameter, public `tetens_c2` = 17.270\_dp
- real(dp), parameter, public `tetens_c3` = 237.30\_dp
- real(dp), parameter, public `satpressureslope1` = 4098.0\_dp  
*calculation of the slope of the saturation vapour pressure curve following Tetens*
- real(dp), parameter, public `desilets_a0` = 0.0808\_dp  
*Neutrons and moisture: N0 formula, Desilets et al. 2010.*
- real(dp), parameter, public `desilets_a1` = 0.372\_dp
- real(dp), parameter, public `desilets_a2` = 0.115\_dp
- real(dp), parameter, public `cosmic_bd` = 1.4020\_dp  
*Neutrons and moisture: COSMIC, Shuttleworth et al. 2013.*
- real(dp), parameter, public `cosmic_vwclat` = 0.0753\_dp
- real(dp), parameter, public `cosmic_n` = 348.33\_dp
- real(dp), parameter, public `cosmic_alpha` = 0.2392421548\_dp
- real(dp), parameter, public `cosmic_l1` = 161.98621864\_dp
- real(dp), parameter, public `cosmic_l2` = 129.14558985\_dp
- real(dp), parameter, public `cosmic_l3` = 107.82204562\_dp
- real(dp), parameter, public `cosmic_l4` = 3.1627190566\_dp

### 17.23.1 Detailed Description

Provides mHM specific constants.

Provides mHM specific constants such as flood plain elevation.

#### Authors

Matthias Cuntz

#### Date

Nov 2011

## 17.23.2 Variable Documentation

### 17.23.2.1 c1\_initstatesm

```
real(dp), parameter, public mo_mhm_constants::c1_initstatesm = 0.25_dp
```

Definition at line 30 of file mo\_mhm\_constants.f90.

### 17.23.2.2 cosmic\_alpha

```
real(dp), parameter, public mo_mhm_constants::cosmic_alpha = 0.2392421548_dp
```

Definition at line 59 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.3 cosmic\_bd

```
real(dp), parameter, public mo_mhm_constants::cosmic_bd = 1.4020_dp
```

Neutrons and moisture: COSMIC, Shuttleworth et al. 2013.

Definition at line 56 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.4 cosmic\_l1

```
real(dp), parameter, public mo_mhm_constants::cosmic_l1 = 161.98621864_dp
```

Definition at line 60 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.5 cosmic\_l2

```
real(dp), parameter, public mo_mhm_constants::cosmic_l2 = 129.14558985_dp
```

Definition at line 61 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.6 cosmic\_l3

```
real(dp), parameter, public mo_mhm_constants::cosmic_l3 = 107.82204562_dp
```

Definition at line 62 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.7 cosmic\_l4

```
real(dp), parameter, public mo_mhm_constants::cosmic_l4 = 3.1627190566_dp
```

Definition at line 63 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.8 cosmic\_n

```
real(dp), parameter, public mo_mhm_constants::cosmic_n = 348.33_dp
```

Definition at line 58 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.9 cosmic\_vwclat

```
real(dp), parameter, public mo_mhm_constants::cosmic_vwclat = 0.0753_dp
```

Definition at line 57 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic().

### 17.23.2.10 desilets\_a0

```
real(dp), parameter, public mo_mhm_constants::desilets_a0 = 0.0808_dp
```

Neutrons and moisture: N0 formula, Desilets et al. 2010.

Definition at line 51 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::desiletsn0().

### 17.23.2.11 desilets\_a1

```
real(dp), parameter, public mo_mhm_constants::desilets_a1 = 0.372_dp
```

Definition at line 52 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::desiletsn0().

### 17.23.2.12 desilets\_a2

```
real(dp), parameter, public mo_mhm_constants::desilets_a2 = 0.115_dp
```

Definition at line 53 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::desiletsn0().

### 17.23.2.13 duffiedelta1

```
real(dp), parameter, public mo_mhm_constants::duffiedelta1 = 0.4090_dp
```

Definition at line 40 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::extraterr\_rad\_approx().

### 17.23.2.14 duffiedelta2

```
real(dp), parameter, public mo_mhm_constants::duffiedelta2 = 1.3900_dp
```

Definition at line 41 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::extraterr\_rad\_approx().

### 17.23.2.15 duffiedr

```
real(dp), parameter, public mo_mhm_constants::duffiedr = 0.0330_dp
```

Definition at line 39 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::extraterr\_rad\_approx().

### 17.23.2.16 h2odens

```
real(dp), parameter, public mo_mhm_constants::h2odens = 1000.0_dp
```

Definition at line 23 of file mo\_mhm\_constants.f90.

Referenced by mo\_neutrons::cosmic(), mo\_mrm\_riv\_temp\_class::finalize\_source\_e(), and mo\_mrm\_riv\_temp\_class::get\_e\_io().

### 17.23.2.17 harsamconst

```
real(dp), parameter, public mo_mhm_constants::harsamconst = 17.800_dp
```

Hargreaves-Samani ref. ET formula [deg C].

Definition at line 36 of file mo\_mhm\_constants.f90.

Referenced by mo\_mhm::mhm().

### 17.23.2.18 noutflxstate

```
integer(i4), parameter, public mo_mhm_constants::noutflxstate = 20_i4
```

Definition at line 33 of file mo\_mhm\_constants.f90.



**17.23.2.19 p2\_initstatefluxes**

```
real(dp), parameter, public mo_mhm_constants::p2_initstatefluxes = 15.00_dp
```

Definition at line 26 of file mo\_mhm\_constants.f90.

**17.23.2.20 p3\_initstatefluxes**

```
real(dp), parameter, public mo_mhm_constants::p3_initstatefluxes = 10.00_dp
```

Definition at line 27 of file mo\_mhm\_constants.f90.

**17.23.2.21 p4\_initstatefluxes**

```
real(dp), parameter, public mo_mhm_constants::p4_initstatefluxes = 75.00_dp
```

Definition at line 28 of file mo\_mhm\_constants.f90.

**17.23.2.22 p5\_initstatefluxes**

```
real(dp), parameter, public mo_mhm_constants::p5_initstatefluxes = 1500.00_dp
```

Definition at line 29 of file mo\_mhm\_constants.f90.

**17.23.2.23 satpressureslope1**

```
real(dp), parameter, public mo_mhm_constants::satpressureslope1 = 4098.0_dp
```

calculation of the slope of the saturation vapour pressure curve following Tetens

Definition at line 48 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::slope\_satpressure().

**17.23.2.24 tetens\_c1**

```
real(dp), parameter, public mo_mhm_constants::tetens_c1 = 0.6108_dp
```

Tetens's formula to calculate saturated vapour pressure.

Definition at line 44 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::sat\_vap\_pressure().

**17.23.2.25 tetens\_c2**

```
real(dp), parameter, public mo_mhm_constants::tetens_c2 = 17.270_dp
```

Definition at line 45 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::sat\_vap\_pressure().

### 17.23.2.26 tetens\_c3

```
real(dp), parameter, public mo_mhm_constants::tetens_c3 = 237.30_dp
```

Definition at line 46 of file mo\_mhm\_constants.f90.

Referenced by mo\_pet::sat\_vap\_pressure(), and mo\_pet::slope\_satpressure().

## 17.24 mo\_mhm\_eval Module Reference

Runs mhm with a specific parameter set and returns required variables, e.g. runoff.

### Functions/Subroutines

- subroutine, public [mhm\\_eval](#) (parameterset, opti\_domain\_indices, runoff, smOptiSim, neutronsOptiSim, et←OptiSim, twsOptiSim)

*Runs mhm with a specific parameter set and returns required variables, e.g. runoff.*

### 17.24.1 Detailed Description

Runs mhm with a specific parameter set and returns required variables, e.g. runoff.

Runs mhm with a specific parameter set and returns required variables, e.g. runoff.

#### Authors

Juliane Mai, Rohini Kumar

#### Date

Feb 2013

### 17.24.2 Function/Subroutine Documentation

#### 17.24.2.1 mhm\_eval()

```
subroutine, public mo_mhm_eval::mhm_eval (
    real(dp), dimension(:), intent(in) parameterset,
    integer(i4), dimension(:), intent(in), optional opti_domain_indices,
    real(dp), dimension(:, :), intent(out), optional, allocatable runoff,
    type(optidata_sim), dimension(:), intent(inout), optional smOptiSim,
    type(optidata_sim), dimension(:), intent(inout), optional neutronsOptiSim,
    type(optidata_sim), dimension(:), intent(inout), optional etOptiSim,
    type(optidata_sim), dimension(:), intent(inout), optional twsOptiSim )
```

Runs mhm with a specific parameter set and returns required variables, e.g. runoff.

Runs mhm with a specific parameter set and returns required variables, e.g. runoff.

## Parameters

|     |   |   |
|-----|---|---|
| in  | <i>real(dp), dimension(:) :: parameterset</i>               | a set of global parameter (gamma) to run mHM, DIMENSION [no. of global_Parameters]  |
| out | <i>real(dp), dimension(:, :), optional :: runoff</i>        | returns runoff time series, DIMENSION [nTimeSteps, nGaugesTotal]  |
| out | <i>real(dp), dimension(:, :), optional :: sm_opti</i>       | returns soil moisture time series for all grid cells (of multiple Domains concatenated), DIMENSION [nCells, nTimeSteps] time series, DIMENSION [nTimeSteps, nDomains] |
| out | <i>real(dp), dimension(:, :), optional :: neutrons_opti</i> | dim1=ncells, dim2=time  |
| out | <i>real(dp), dimension(:, :), optional :: et_opti</i>       | returns evapotranspiration time series for  |
| out | <i>real(dp), dimension(:, :), optional :: tws_opti</i>      | returns tws time series all grid cells (of multiple Domains concatenated), DIMENSION [nCells, nTimeSteps]   |

## Authors

Juliane Mai, Rohini Kumar

## Date

Feb 2013

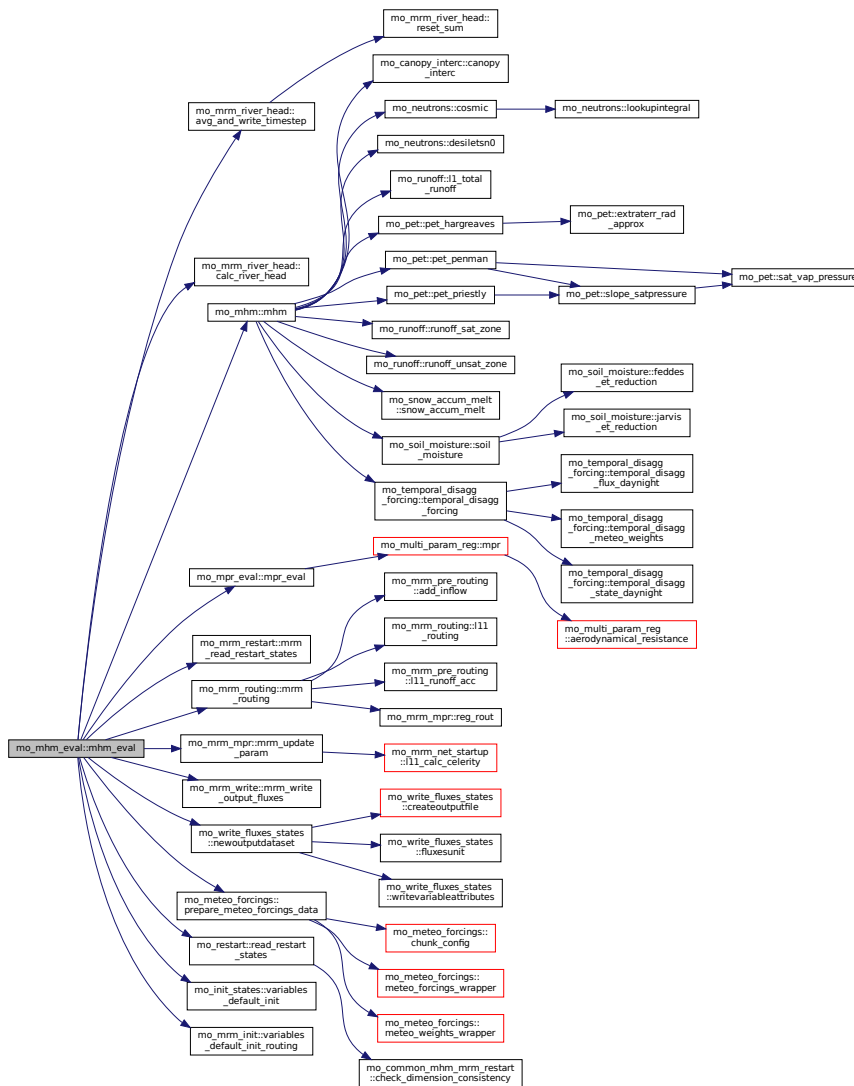
Definition at line 83 of file mo\_mhm\_eval.f90.

References mo\_mrm\_river\_head::avg\_and\_write\_timestep(), mo\_common\_mhm\_mrm\_variables::c2tstu, mo\_mrm\_river\_head::calc\_river\_head(), mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_global\_variables::evap\_coeff, mo\_global\_variables::fday\_pet, mo\_global\_variables::fday\_prec, mo\_global\_variables::fday\_ssr, mo\_global\_variables::fday\_strd, mo\_global\_variables::fday\_temp, mo\_global\_variables::fnight\_pet, mo\_global\_variables::fnight\_prec, mo\_global\_variables::fnight\_ssr, mo\_global\_variables::fnight\_strd, mo\_global\_variables::fnight\_temp, mo\_mrm\_global\_variables::gw\_coupling, mo\_mpr\_global\_variables::horizontdepth\_mhm, mo\_mrm\_global\_variables::inflowgauge, mo\_mrm\_global\_variables::l0\_river\_head\_mon\_sum, mo\_mrm\_global\_variables::l11\_c1, mo\_mrm\_global\_variables::l11\_c2, mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_global\_variables::l11\_l1\_id, mo\_mrm\_global\_variables::l11\_length, mo\_mrm\_global\_variables::l11\_netperm, mo\_mrm\_global\_variables::l11\_nlinkfracpimp, mo\_mrm\_global\_variables::l11\_noutlets, mo\_mrm\_global\_variables::l11\_qmod, mo\_mrm\_global\_variables::l11\_qout, mo\_mrm\_global\_variables::l11\_qtin, mo\_mrm\_global\_variables::l11\_qtr, mo\_mrm\_global\_variables::l11\_slope, mo\_mrm\_global\_variables::l11\_ton, mo\_mrm\_global\_variables::l11\_tsout, mo\_global\_variables::l1\_absvappress, mo\_mpr\_global\_variables::l1\_aeroresist, mo\_global\_variables::l1\_aetcanopy, mo\_global\_variables::l1\_aetsealed, mo\_global\_variables::l1\_aetsoil, mo\_mpr\_global\_variables::l1\_alpha, mo\_global\_variables::l1\_baseflow, mo\_mpr\_global\_variables::l1\_degday, mo\_mpr\_global\_variables::l1\_degdayinc, mo\_mpr\_global\_variables::l1\_degdaymax, mo\_mpr\_global\_variables::l1\_degdaynopre, mo\_global\_variables::l1\_etobs, mo\_mpr\_global\_variables::l1\_fasp, mo\_global\_variables::l1\_fastrunoff, mo\_mpr\_global\_variables::l1\_froots, mo\_mpr\_global\_variables::l1\_fsealed, mo\_mpr\_global\_variables::l1\_harsamcoeff, mo\_global\_variables::l1\_infilsoil, mo\_global\_variables::l1\_inter, mo\_mpr\_global\_variables::l1\_jarvis\_thresh\_c1, mo\_mpr\_global\_variables::l1\_karstloss, mo\_mpr\_global\_variables::l1\_kbaseflow, mo\_mpr\_global\_variables::l1\_kfastflow, mo\_mpr\_global\_variables::l1\_kperco, mo\_mpr\_global\_variables::l1\_kslowflow, mo\_mrm\_global\_variables::l1\_l11\_id, mo\_mpr\_global\_variables::l1\_maxinter, mo\_global\_variables::l1\_melt, mo\_global\_variables::l1\_netrad, mo\_global\_variables::l1\_neutrons, mo\_global\_variables::l1\_neutronsobs, mo\_global\_variables::l1\_percol, mo\_global\_variables::l1\_pet, mo\_global\_variables::l1\_pet\_calc, mo\_global\_variables::l1\_pet\_weights, mo\_mpr\_global\_variables::l1\_pettaicorfactor, mo\_global\_variables::l1\_pre, mo\_global\_variables::l1\_pre\_weights, mo\_global\_variables::l1\_preeffect, mo\_mpr\_global\_variables::l1\_prietaryalpha, mo\_global\_variables::l1\_rain, mo\_global\_variables::l1\_runoffseal, mo\_global\_variables::l1\_satstw, mo\_mpr\_global\_variables::l1\_sealedthresh, mo\_global\_variables::l1\_sealstw, mo\_global\_variables::l1\_slowrunoff, mo\_global\_variables::l1\_smobs, mo\_global\_variables::l1\_snow, mo\_global\_variables::l1\_snowpack, mo\_global\_variables::l1\_soilmoist, mo\_mpr\_global\_variables::l1\_soilmoistexp, mo\_mpr\_global\_variables::l1\_soilmoistfc, mo\_mpr\_global\_variables::l1\_soilmoistsat, mo\_global\_variables::l1\_ssr, mo\_global\_variables::

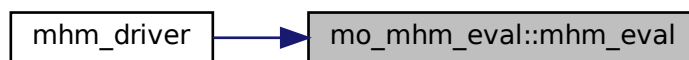
::l1\_strd, mo\_mpr\_global\_variables::l1\_surfresist, mo\_global\_variables::l1\_tann, mo\_global\_variables::l1\_↵  
temp, mo\_global\_variables::l1\_temp\_weights, mo\_mpr\_global\_variables::l1\_tempthresh, mo\_global\_variables↵  
::l1\_throughfall, mo\_global\_variables::l1\_tmax, mo\_global\_variables::l1\_tmin, mo\_global\_variables::l1\_total↵  
\_runoff, mo\_global\_variables::l1\_twsaobs, mo\_global\_variables::l1\_unsatstw, mo\_mpr\_global\_variables::l1\_↵  
unsatthresh, mo\_mpr\_global\_variables::l1\_wiltingpoint, mo\_global\_variables::l1\_windspeed, mo\_common\_↵  
mhm\_mrm\_variables::lcyyearid, mo\_common\_variables::level1, mo\_mrm\_global\_variables::level11, mo\_mhm\_↵  
::mhm(), mo\_common\_mhm\_mrm\_variables::mhmfilerestartin, mo\_mpr\_eval::mpr\_eval(), mo\_mrm\_restart::mrm\_↵  
\_read\_restart\_states(), mo\_mrm\_routing::mrm\_routing(), mo\_mrm\_global\_variables::mrm\_runoff, mo\_mrm\_↵  
mpr::mrm\_update\_param(), mo\_mrm\_write::mrm\_write\_output\_fluxes(), mo\_common\_mhm\_mrm\_variables↵  
::mrmfilerestartin, mo\_global\_variables::neutron\_integral\_afast, mo\_write\_fluxes\_states::newoutputdataset(), mo\_↵  
\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_global\_variables::nsoilhorizons\_sm\_input, mo\_common\_mhm\_↵  
\_mrm\_variables::ntstepday, mo\_common\_mhm\_mrm\_variables::optimize, mo\_global\_variables::outputflxstate,↵  
mo\_mrm\_global\_variables::outputflxstate\_mrm, mo\_meteo\_forcings::prepare\_meteo\_forcings\_data(), mo\_↵  
common\_variables::processmatrix, mo\_global\_variables::read\_meteo\_weights, mo\_common\_mhm\_mrm\_↵  
variables::read\_restart, mo\_restart::read\_restart\_states(), mo\_common\_mhm\_mrm\_variables::readper, mo\_↵  
\_common\_variables::resolutionhydrology, mo\_common\_mhm\_mrm\_variables::resolutionrouting, mo\_mrm\_↵  
global\_variables::riv\_temp\_pcs, mo\_common\_mhm\_mrm\_variables::simper, mo\_common\_mhm\_mrm\_variables↵  
::timestep, mo\_global\_variables::timestep\_model\_inputs, mo\_global\_variables::timestep\_model\_outputs, mo\_↵  
\_mrm\_global\_variables::timestep\_model\_outputs\_mrm, mo\_init\_states::variables\_default\_init(), mo\_mrm\_init\_↵  
::variables\_default\_init\_routing(), and mo\_common\_mhm\_mrm\_variables::warmingdays.

Referenced by mhm\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.25 mo\_mhm\_read\_config Module Reference

Reading of main model configurations.

## Functions/Subroutines

- subroutine, public `mhm_read_config` (`file_namelist`, `unamelist`)  
*Read main configurations for mHM.*

### 17.25.1 Detailed Description

Reading of main model configurations.

This routine reads the configurations of mHM including, input and output directories, module usage specification, simulation time periods, global parameters, ...

#### Authors

Matthias Zink

#### Date

Dec 2012

### 17.25.2 Function/Subroutine Documentation

#### 17.25.2.1 `mhm_read_config()`

```
subroutine, public mo_mhm_read_config::mhm_read_config (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist )
```

Read main configurations for mHM.

The main configurations in mHM are read from three files:

1. `mhm.nml`
2. `mhm_parameters.nml`
3. `mhm_outputs.nml`

For details please refer to the above mentioned namelist files.

#### Parameters

|    |                                      |  |
|----|--------------------------------------|--|
| in | <i>character(*) :: file_namelist</i> |  |
| in | <i>integer :: unamelist</i>          |  |

#### Authors

Matthias Zink

#### Date

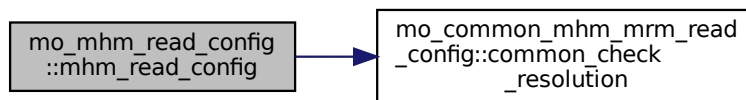
Dec 2012

Definition at line 89 of file `mo_mhm_read_config.f90`.

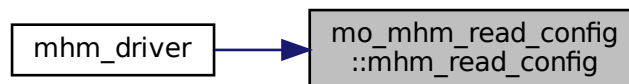
References mo\_common\_mhm\_mrm\_read\_config::common\_check\_resolution(), mo\_global\_variables::dirabsvappressure, mo\_global\_variables::dirmaxtemperature, mo\_global\_variables::dirminttemperature, mo\_global\_variables::dirnetradiation, mo\_global\_variables::dirprecipitation, mo\_global\_variables::dirradiation, mo\_global\_variables::dirreferenceet, mo\_global\_variables::dirtemperature, mo\_global\_variables::dirwindspeed, mo\_common\_variables::domainmeta, mo\_global\_variables::evap\_coeff, mo\_global\_variables::fday\_pet, mo\_global\_variables::fday\_prec, mo\_global\_variables::fday\_ssr, mo\_global\_variables::fday\_strd, mo\_global\_variables::fday\_temp, mo\_file::file\_defoutput, mo\_global\_variables::fnight\_pet, mo\_global\_variables::fnight\_prec, mo\_global\_variables::fnight\_ssr, mo\_global\_variables::fnight\_strd, mo\_global\_variables::fnight\_temp, mo\_global\_variables::inputformat\_meteo\_forcings, mo\_global\_variables::l1\_etobs, mo\_global\_variables::l1\_neutronsobs, mo\_global\_variables::l1\_smobs, mo\_global\_variables::l1\_twsaobs, mo\_common\_constants::maxnodomains, mo\_mpr\_constants::maxnosoilhorizons, mo\_common\_constants::nodata\_i4, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_global\_variables::nsoilhorizons\_sm\_input, mo\_common\_mhm\_mrm\_variables::opti\_function, mo\_common\_mhm\_mrm\_variables::optimize, mo\_global\_variables::output\_deflate\_level, mo\_global\_variables::output\_double\_precision, mo\_global\_variables::outputflxstate, mo\_common\_variables::processmatrix, mo\_global\_variables::read\_meteo\_weights, mo\_global\_variables::timestep\_model\_inputs, mo\_global\_variables::timestep\_model\_outputs, and mo\_file::udefoutput.

Referenced by mhm\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.26 mo\_mpr\_constants Module Reference

Provides MPR specific constants.

### Variables

- integer(i4), parameter, public `nlcover_class` = 3\_i4
- integer(i4), parameter, public `maxgeounit` = 25\_i4
- integer(i4), parameter, public `maxnosoilhorizons` = 10\_i4
- real(dp), parameter, public `p2_initstatefluxes` = 15.00\_dp
- real(dp), parameter, public `p3_initstatefluxes` = 10.00\_dp

- real(dp), parameter, public `p4_initstatefluxes` = 75.00\_dp
- real(dp), parameter, public `p5_initstatefluxes` = 1500.00\_dp
- real(dp), parameter, public `c1_initstatesm` = 0.25\_dp
- real(dp), parameter, public `bulkdens_orgmatter` = 0.224\_dp
- real(dp), parameter, public `field_cap_c1` = -0.60\_dp
- real(dp), parameter, public `field_cap_c2` = 2.0\_dp
- real(dp), parameter, public `vgenuchten_sandfresh` = 66.5\_dp
- real(dp), parameter, public `vgenuchtenn_c1` = 1.392\_dp
- real(dp), parameter, public `vgenuchtenn_c2` = 0.418\_dp
- real(dp), parameter, public `vgenuchtenn_c3` = -0.024\_dp
- real(dp), parameter, public `vgenuchtenn_c4` = 1.212\_dp
- real(dp), parameter, public `vgenuchtenn_c5` = -0.704\_dp
- real(dp), parameter, public `vgenuchtenn_c6` = -0.648\_dp
- real(dp), parameter, public `vgenuchtenn_c7` = 0.023\_dp
- real(dp), parameter, public `vgenuchtenn_c8` = 0.044\_dp
- real(dp), parameter, public `vgenuchtenn_c9` = 3.168\_dp
- real(dp), parameter, public `vgenuchtenn_c10` = -2.562\_dp
- real(dp), parameter, public `vgenuchtenn_c11` = 7.0E-9\_dp
- real(dp), parameter, public `vgenuchtenn_c12` = 4.004\_dp
- real(dp), parameter, public `vgenuchtenn_c13` = 3.750\_dp
- real(dp), parameter, public `vgenuchtenn_c14` = -0.016\_dp
- real(dp), parameter, public `vgenuchtenn_c15` = -4.197\_dp
- real(dp), parameter, public `vgenuchtenn_c16` = 0.013\_dp
- real(dp), parameter, public `vgenuchtenn_c17` = 0.076\_dp
- real(dp), parameter, public `vgenuchtenn_c18` = 0.276\_dp
- real(dp), parameter, public `ks_c` = 10.0\_dp
- real(dp), parameter, public `pwp_c` = 1.0\_dp
- real(dp), parameter, public `pwp_matpot_thetar` = 15000.0\_dp
- real(dp), parameter, public `windmeasheight` = 10.0\_dp
  - *assumed meteorol. measurement hight for estimation of aeroResist and surfResist*
- real(dp), parameter, public `karman` = 0.41\_dp
  - *von karman constant*
- real(dp), parameter, public `lai_factor_surfresi` = 0.3\_dp
  - *LAI factor for bulk surface resistance formulation.*
- real(dp), parameter, public `lai_offset_surfresi` = 1.2\_dp
  - *LAI offset for bulk surface resistance formulation.*
- real(dp), parameter, public `max_surfresist` = 250.0\_dp
  - *maximum bulk surface resistance*

### 17.26.1 Detailed Description

Provides MPR specific constants.

Provides MPR specific constants such as flood plain elevation.

#### Authors

Matthias Cuntz

#### Date

Nov 2011



## 17.26.2 Variable Documentation

### 17.26.2.1 bulkdens\_orgmatter

```
real(dp), parameter, public mo_mpr_constants::bulkdens_orgmatter = 0.224_dp
```

Definition at line 35 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::mpr\_sm().

### 17.26.2.2 c1\_initstatesm

```
real(dp), parameter, public mo_mpr_constants::c1_initstatesm = 0.25_dp
```

Definition at line 31 of file mo\_mpr\_constants.f90.

Referenced by mo\_init\_states::variables\_alloc(), and mo\_init\_states::variables\_default\_init().

### 17.26.2.3 field\_cap\_c1

```
real(dp), parameter, public mo_mpr_constants::field_cap_c1 = -0.60_dp
```

Definition at line 37 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::field\_cap().

### 17.26.2.4 field\_cap\_c2

```
real(dp), parameter, public mo_mpr_constants::field_cap_c2 = 2.0_dp
```

Definition at line 38 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::field\_cap().

### 17.26.2.5 karman

```
real(dp), parameter, public mo_mpr_constants::karman = 0.41_dp
```

von karman constant

Definition at line 68 of file mo\_mpr\_constants.f90.

Referenced by mo\_multi\_param\_reg::aerodynamical\_resistance().

### 17.26.2.6 ks\_c

```
real(dp), parameter, public mo_mpr_constants::ks_c = 10.0_dp
```

Definition at line 60 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::hydro\_cond().

### 17.26.2.7 lai\_factor\_surfresi

```
real(dp), parameter, public mo_mpr_constants::lai_factor_surfresi = 0.3_dp
```

LAI factor for bulk surface resistance formulation.

Definition at line 71 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_pet::bulksurface\_resistance().

### 17.26.2.8 lai\_offset\_surfresi

```
real(dp), parameter, public mo_mpr_constants::lai_offset_surfresi = 1.2_dp
```

LAI offset for bulk surface resistance formulation.

Definition at line 73 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_pet::bulksurface\_resistance().

### 17.26.2.9 max\_surfresist

```
real(dp), parameter, public mo_mpr_constants::max_surfresist = 250.0_dp
```

maximum bulk surface resistance

Definition at line 75 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_pet::bulksurface\_resistance().

### 17.26.2.10 maxgeounit

```
integer(i4), parameter, public mo_mpr_constants::maxgeounit = 25_i4
```

Definition at line 23 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_read\_config::mpr\_read\_config().

### 17.26.2.11 maxnosoilhorizons

```
integer(i4), parameter, public mo_mpr_constants::maxnosoilhorizons = 10_i4
```

Definition at line 24 of file mo\_mpr\_constants.f90.

Referenced by mo\_mhm\_read\_config::mhm\_read\_config(), and mo\_mpr\_read\_config::mpr\_read\_config().

### 17.26.2.12 nlcover\_class

```
integer(i4), parameter, public mo_mpr_constants::nlcover_class = 3_i4
```

Definition at line 22 of file mo\_mpr\_constants.f90.

Referenced by mo\_soil\_database::read\_soil\_lut().

#### 17.26.2.13 p2\_initstatefluxes

```
real(dp), parameter, public mo_mpr_constants::p2_initstatefluxes = 15.00_dp
```

Definition at line 27 of file mo\_mpr\_constants.f90.

Referenced by mo\_init\_states::variables\_alloc(), and mo\_init\_states::variables\_default\_init().

#### 17.26.2.14 p3\_initstatefluxes

```
real(dp), parameter, public mo_mpr_constants::p3_initstatefluxes = 10.00_dp
```

Definition at line 28 of file mo\_mpr\_constants.f90.

Referenced by mo\_init\_states::variables\_alloc(), and mo\_init\_states::variables\_default\_init().

#### 17.26.2.15 p4\_initstatefluxes

```
real(dp), parameter, public mo_mpr_constants::p4_initstatefluxes = 75.00_dp
```

Definition at line 29 of file mo\_mpr\_constants.f90.

Referenced by mo\_init\_states::variables\_alloc(), and mo\_init\_states::variables\_default\_init().

#### 17.26.2.16 p5\_initstatefluxes

```
real(dp), parameter, public mo_mpr_constants::p5_initstatefluxes = 1500.00_dp
```

Definition at line 30 of file mo\_mpr\_constants.f90.

Referenced by mo\_init\_states::variables\_alloc(), and mo\_init\_states::variables\_default\_init().

#### 17.26.2.17 pwp\_c

```
real(dp), parameter, public mo_mpr_constants::pwp_c = 1.0_dp
```

Definition at line 62 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::pwp().

#### 17.26.2.18 pwp\_matpot\_thetar

```
real(dp), parameter, public mo_mpr_constants::pwp_matpot_thetar = 15000.0_dp
```

Definition at line 63 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::pwp().

**17.26.2.19 vgenuchten\_sandtresh**

```
real(dp), parameter, public mo_mpr_constants::vgenuchten_sandtresh = 66.5_dp
```

Definition at line 40 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.20 vgenuchtenn\_c1**

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c1 = 1.392_dp
```

Definition at line 41 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.21 vgenuchtenn\_c10**

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c10 = -2.562_dp
```

Definition at line 50 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.22 vgenuchtenn\_c11**

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c11 = 7.0E-9_dp
```

Definition at line 51 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.23 vgenuchtenn\_c12**

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c12 = 4.004_dp
```

Definition at line 52 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.24 vgenuchtenn\_c13**

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c13 = 3.750_dp
```

Definition at line 53 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.25 vgenuchtenn\_c14**

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c14 = -0.016_dp
```

Definition at line 54 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

#### 17.26.2.26 vgenuchtenn\_c15

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c15 = -4.197_dp
```

Definition at line 55 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

#### 17.26.2.27 vgenuchtenn\_c16

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c16 = 0.013_dp
```

Definition at line 56 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

#### 17.26.2.28 vgenuchtenn\_c17

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c17 = 0.076_dp
```

Definition at line 57 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

#### 17.26.2.29 vgenuchtenn\_c18

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c18 = 0.276_dp
```

Definition at line 58 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

#### 17.26.2.30 vgenuchtenn\_c2

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c2 = 0.418_dp
```

Definition at line 42 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

#### 17.26.2.31 vgenuchtenn\_c3

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c3 = -0.024_dp
```

Definition at line 43 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

### 17.26.2.32 vgenuchtenn\_c4

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c4 = 1.212_dp
```

Definition at line 44 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

### 17.26.2.33 vgenuchtenn\_c5

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c5 = -0.704_dp
```

Definition at line 45 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

### 17.26.2.34 vgenuchtenn\_c6

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c6 = -0.648_dp
```

Definition at line 46 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

### 17.26.2.35 vgenuchtenn\_c7

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c7 = 0.023_dp
```

Definition at line 47 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

### 17.26.2.36 vgenuchtenn\_c8

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c8 = 0.044_dp
```

Definition at line 48 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

### 17.26.2.37 vgenuchtenn\_c9

```
real(dp), parameter, public mo_mpr_constants::vgenuchtenn_c9 = 3.168_dp
```

Definition at line 49 of file mo\_mpr\_constants.f90.

Referenced by mo\_mpr\_soilmoist::genuchten().

**17.26.2.38 windmeasheight**

```
real(dp), parameter, public mo_mpr_constants::windmeasheight = 10.0_dp
```

assumed meteorol. measurement hight for estimation of aeroResist and surfResist

Definition at line 66 of file mo\_mpr\_constants.f90.

Referenced by mo\_multi\_param\_reg::aerodynamical\_resistance().

**17.27 mo\_mpr\_eval Module Reference**

Runs MPR and writes to global effective parameters.

**Functions/Subroutines**

- subroutine, public [mpr\\_eval](#) (parameterset, opti\_domain\_indices)  
*Runs MPR and writes to global effective parameters.*

**17.27.1 Detailed Description**

Runs MPR and writes to global effective parameters.

Runs MPR and writes to global effective parameters

**Authors**

Robert Schweppe

**Date**

Feb 2018

**17.27.2 Function/Subroutine Documentation****17.27.2.1 mpr\_eval()**

```
subroutine, public mo_mpr_eval::mpr_eval (
    real(dp), dimension(:), intent(in), optional parameterset,
    integer(i4), dimension(:), intent(in), optional opti_domain_indices )
```

Runs MPR and writes to global effective parameters.

Runs MPR and writes to global effective parameters

**Parameters**

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:), optional :: parameterset</i> | a set of global parameter (gamma) to run mHM, DIMENSION [no. of global_Parameters] |
|----|---|--|

**Authors**

Juliane Mai, Rohini Kumar

Date

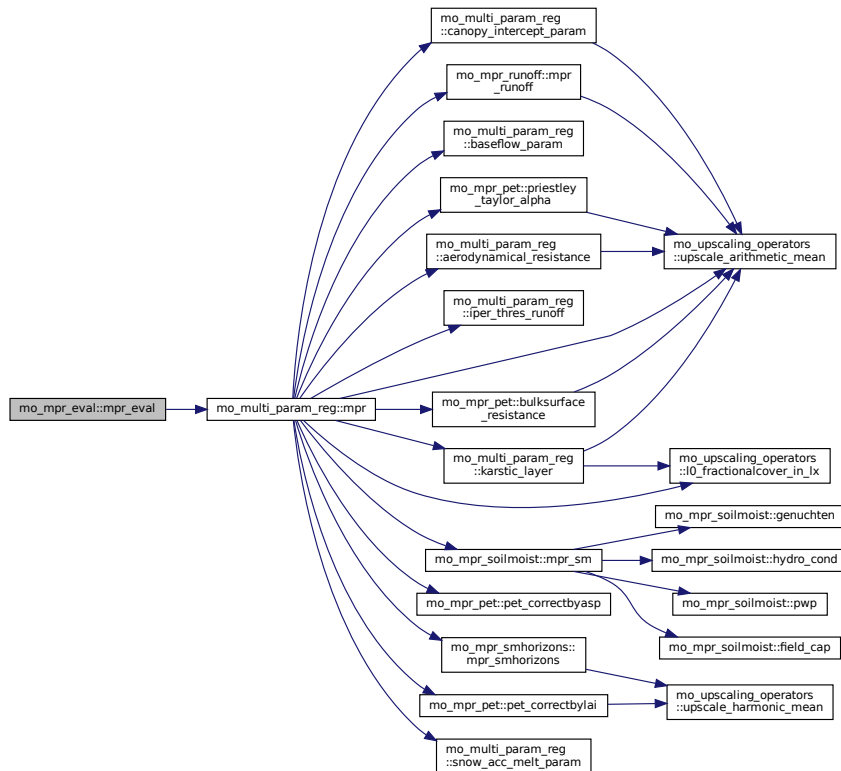
Feb 2013

Definition at line 71 of file mo\_mpr\_eval.f90.

References mo\_common\_variables::domainmeta, mo\_mpr\_global\_variables::l0\_esp, mo\_mpr\_global\_variables::l0\_geounit, mo\_mpr\_global\_variables::l0\_gridded\_lai, mo\_common\_variables::l0\_l1\_remap, mo\_common\_variables::l0\_lcover, mo\_mpr\_global\_variables::l0\_slope\_emp, mo\_mpr\_global\_variables::l0\_soilid, mo\_mpr\_global\_variables::l1\_aeroresist, mo\_mpr\_global\_variables::l1\_alpha, mo\_mpr\_global\_variables::l1\_degdayinc, mo\_mpr\_global\_variables::l1\_degdaymax, mo\_mpr\_global\_variables::l1\_degdaynopre, mo\_mpr\_global\_variables::l1\_fasp, mo\_mpr\_global\_variables::l1\_froots, mo\_mpr\_global\_variables::l1\_fsealed, mo\_mpr\_global\_variables::l1\_harsamcoeff, mo\_mpr\_global\_variables::l1\_jarvis\_thresh\_c1, mo\_mpr\_global\_variables::l1\_karstloss, mo\_mpr\_global\_variables::l1\_kbaseflow, mo\_mpr\_global\_variables::l1\_kfastflow, mo\_mpr\_global\_variables::l1\_kperco, mo\_mpr\_global\_variables::l1\_kslowflow, mo\_mpr\_global\_variables::l1\_maxinter, mo\_mpr\_global\_variables::l1\_petlaicorfactor, mo\_mpr\_global\_variables::l1\_prietayalpha, mo\_mpr\_global\_variables::l1\_sealedthresh, mo\_mpr\_global\_variables::l1\_soilmoistexp, mo\_mpr\_global\_variables::l1\_soilmoistfc, mo\_mpr\_global\_variables::l1\_soilmoistsat, mo\_mpr\_global\_variables::l1\_surfresist, mo\_mpr\_global\_variables::l1\_tempthresh, mo\_mpr\_global\_variables::l1\_unsatthresh, mo\_mpr\_global\_variables::l1\_wiltingpoint, mo\_common\_variables::level0, mo\_common\_variables::level1, and mo\_multi\_param\_reg::mpr().

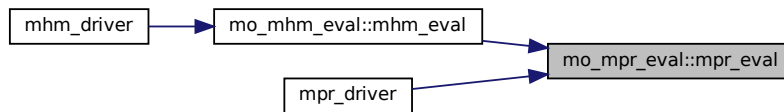
Referenced by mo\_mhm\_eval::mhm\_eval(), and mpr\_driver().

Here is the call graph for this function:





Here is the caller graph for this function:



## 17.28 mo\_mpr\_file Module Reference

Provides file names and units for mRM.

### Variables

- character(len=\*), parameter `version` = '0.1'  
*Current mHM model version.*
- character(len=\*), parameter `version_date` = 'Jun 2019'  
*Time of current mHM model version release.*
- character(len=\*), parameter `file_main` = 'mpr\_driver.f90'  
*Driver file.*
- character(len=\*), parameter `file_namelist_mpr` = 'mpr.nml'  
*Namelist file name.*
- integer, parameter `unamelist_mpr` = 80  
*Unit for namelist.*
- character(len=\*), parameter `file_namelist_mpr_param` = 'mpr\_parameter.nml'  
*Parameter namelists file name.*
- integer, parameter `unamelist_mpr_param` = 31  
*Unit for namelist.*
- character(len=\*), parameter `file_soil_database` = 'soil\_classdefinition.txt'  
*Soil database file (iFlag\_soilDB = 0) = classical mHM format.*
- character(len=\*), parameter `file_soil_database_1` = 'soil\_classdefinition\_iFlag\_soilDB\_1.txt'  
*Soil database file (iFlag\_soilDB = 1)*
- integer, parameter `usoil_database` = 52  
*Unit for soil data base.*
- character(len=\*), parameter `file_slope` = 'slope.asc'  
*slope input data file*
- integer, parameter `uslope` = 54  
*Unit for slope input data file.*
- character(len=\*), parameter `file_aspect` = 'aspect.asc'  
*aspect input data file*
- integer, parameter `uaspect` = 55  
*Unit for aspect input data file.*
- character(len=\*), parameter `file_hydrogeoclass` = 'geology\_class.asc'  
*hydrogeological classes input data file*
- integer, parameter `uhydrogeoclass` = 58  
*Unit for hydrogeological classes input data file.*
- character(len=\*), parameter `file_soilclass` = 'soil\_class.asc'

- soil classes input data file*

  - integer, parameter `usoilclass` = 59

*Unit for soil classes input data file.*
- character(len=\*), parameter `file_laiclass` = 'LAI\_class.asc'

*LAI classes input data file.*
- integer, parameter `ulaiclass` = 60

*Unit for LAI input data file.*
- character(len=\*), parameter `file_geolut` = 'geology\_classdefinition.txt'

*geological formation lookup table file*
- integer, parameter `ugeolut` = 64

*Unit for geological formation lookup table file.*
- character(len=\*), parameter `file_lailut` = 'LAI\_classdefinition.txt'

*LAI classes lookup table file.*
- integer, parameter `ulailut` = 65

*Unit for LAI classes lookup table file.*
- character(len=\*), parameter `file_meteo_header` = 'header.txt'

*Input nCols and nRows of binary meteo and LAI files are in header file.*
- integer, parameter `umeteo_header` = 50

*Unit for meteo header file.*
- character(len=\*), parameter `file_meteo_binary_end` = '.bin'

*File ending of meteo files.*
- integer, parameter `umeteo` = 51

*Unit for meteo files.*

### 17.28.1 Detailed Description

Provides file names and units for mRM.

Provides all filenames as well as all units used for the multiscale Routing Model mRM.

#### Authors

Matthias Cuntz, Stephan Thober

#### Date

Aug 2015

### 17.28.2 Variable Documentation

#### 17.28.2.1 file\_aspect

```
character(len = *), parameter mo_mpr_file::file_aspect = 'aspect.asc'
```

aspect input data file

Definition at line 44 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

### 17.28.2.2 file\_geolut

```
character(len = *), parameter mo_mpr_file::file_geolut = 'geology_classdefinition.txt'
```

geological formation lookup table file

Definition at line 61 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

### 17.28.2.3 file\_hydrogeoclass

```
character(len = *), parameter mo_mpr_file::file_hydrogeoclass = 'geology_class.asc'
```

hydrogeological classes input data file

Definition at line 48 of file mo\_mpr\_file.f90.

Referenced by mo\_startup::constants\_init(), and mo\_read\_wrapper::read\_data().

### 17.28.2.4 file\_laiclass

```
character(len = *), parameter mo_mpr_file::file_laiclass = 'LAI_class.asc'
```

LAI classes input data file.

Definition at line 56 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

### 17.28.2.5 file\_lailut

```
character(len = *), parameter mo_mpr_file::file_lailut = 'LAI_classdefinition.txt'
```

LAI classes lookup table file.

Definition at line 66 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

### 17.28.2.6 file\_main

```
character(len = *), parameter mo_mpr_file::file_main = 'mpr_driver.f90'
```

Driver file.

Definition at line 22 of file mo\_mpr\_file.f90.

### 17.28.2.7 file\_meteo\_binary\_end

```
character(len = *), parameter mo_mpr_file::file_meteo_binary_end = '.bin'
```

File ending of meteo files.

Definition at line 75 of file mo\_mpr\_file.f90.

### 17.28.2.8 file\_meteo\_header

```
character(len = *), parameter mo_mpr_file::file_meteo_header = 'header.txt'
```

Input nCols and nRows of binary meteo and LAI files are in header file.

Definition at line 71 of file mo\_mpr\_file.f90.

Referenced by mo\_startup::l2\_variable\_init().

### 17.28.2.9 file\_namelist\_mpr

```
character(len = *), parameter mo_mpr_file::file_namelist_mpr = 'mpr.nml'
```

Namelist file name.

Definition at line 24 of file mo\_mpr\_file.f90.

### 17.28.2.10 file\_namelist\_mpr\_param

```
character(len = *), parameter mo_mpr_file::file_namelist_mpr_param = 'mpr_parameter.nml'
```

Parameter namelists file name.

Definition at line 28 of file mo\_mpr\_file.f90.

Referenced by mpr\_driver().

### 17.28.2.11 file\_slope

```
character(len = *), parameter mo_mpr_file::file_slope = 'slope.asc'
```

slope input data file

Definition at line 40 of file mo\_mpr\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_I0\_data(), and mo\_read\_wrapper::read\_data().

### 17.28.2.12 file\_soil\_database

```
character(len = *), parameter mo_mpr_file::file_soil_database = 'soil_classdefinition.txt'
```

Soil database file (iFlag\_soilDB = 0) = classical mHM format.

Definition at line 34 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

### 17.28.2.13 file\_soil\_database\_1

```
character(len = *), parameter mo_mpr_file::file_soil_database_1 = 'soil_classdefinition_iFlag←  
_soilDB_1.txt'
```

Soil database file (iFlag\_soilDB = 1)

Definition at line 36 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.14 file\_soilclass

```
character(len = *), parameter mo_mpr_file::file_soilclass = 'soil_class.asc'
```

soil classes input data file

Definition at line 52 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.15 uaspect

```
integer, parameter mo_mpr_file::uaspect = 55
```

Unit for aspect input data file.

Definition at line 46 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.16 ugeolut

```
integer, parameter mo_mpr_file::ugeolut = 64
```

Unit for geological formation lookup table file.

Definition at line 63 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.17 uhydrogeoclass

```
integer, parameter mo_mpr_file::uhydrogeoclass = 58
```

Unit for hydrogeological classes input data file.

Definition at line 50 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.18 ulaiclass

```
integer, parameter mo_mpr_file::ulaiclass = 60
```

Unit for LAI input data file.

Definition at line 58 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.19 ulailut

```
integer, parameter mo_mpr_file::ulailut = 65
```

Unit for LAI classes lookup table file.

Definition at line 68 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.28.2.20 umeteo

```
integer, parameter mo_mpr_file::umeteo = 51
```

Unit for meteo files.

Definition at line 77 of file mo\_mpr\_file.f90.

#### 17.28.2.21 umeteo\_header

```
integer, parameter mo_mpr_file::umeteo_header = 50
```

Unit for meteo header file.

Definition at line 73 of file mo\_mpr\_file.f90.

Referenced by mo\_startup::l2\_variable\_init().

#### 17.28.2.22 unamelist\_mpr

```
integer, parameter mo_mpr_file::unamelist_mpr = 80
```

Unit for namelist.

Definition at line 26 of file mo\_mpr\_file.f90.

#### 17.28.2.23 unamelist\_mpr\_param

```
integer, parameter mo_mpr_file::unamelist_mpr_param = 31
```

Unit for namelist.

Definition at line 30 of file mo\_mpr\_file.f90.

Referenced by mpr\_driver().

#### 17.28.2.24 uslope

```
integer, parameter mo_mpr_file::uslope = 54
```

Unit for slope input data file.

Definition at line 42 of file mo\_mpr\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_IO\_data(), and mo\_read\_wrapper::read\_data().

### 17.28.2.25 usoil\_database

```
integer, parameter mo_mpr_file::usoil_database = 52
```

Unit for soil data base.

Definition at line 38 of file mo\_mpr\_file.f90.

Referenced by mo\_soil\_database::read\_soil\_lut().

### 17.28.2.26 usoilclass

```
integer, parameter mo_mpr_file::usoilclass = 59
```

Unit for soil classes input data file.

Definition at line 54 of file mo\_mpr\_file.f90.

Referenced by mo\_read\_wrapper::read\_data().

### 17.28.2.27 version

```
character(len = *), parameter mo_mpr_file::version = '0.1'
```

Current mHM model version.

Definition at line 18 of file mo\_mpr\_file.f90.

### 17.28.2.28 version\_date

```
character(len = *), parameter mo_mpr_file::version_date = 'Jun 2019'
```

Time of current mHM model version release.

Definition at line 20 of file mo\_mpr\_file.f90.

## 17.29 mo\_mpr\_global\_variables Module Reference

Global variables for mpr only.

### Data Types

- type [soiltype](#)

### Variables

- real(dp), public [tillagedepth](#)
- integer(i4), public [nsoiltypes](#)
- integer(i4), public [iflag\\_soildb](#)
- integer(i4), public [nsoilhorizons\\_mhm](#)

- real(dp), dimension(:), allocatable, public [horizondepth\\_mhm](#)
- type([soiltype](#)), public [soildb](#)
- integer(i4), public [ngeounits](#)
- integer(i4), dimension(:), allocatable, public [geounitlist](#)
- integer(i4), dimension(:), allocatable, public [geounitkar](#)
- character(256), public [inputformat\\_gridded\\_lai](#)
- integer(i4), public [timestep\\_lai\\_input](#)
- integer(i4), public [nlaiclass](#)
- integer(i4), public [nlai](#)
- real(dp), dimension(:), allocatable, public [laiboundaries](#)
- integer(i4), dimension(:), allocatable, public [laiunitlist](#)
- real(dp), dimension(:, :), allocatable, public [lailut](#)
- type([period](#)), dimension(:), allocatable, public [laiper](#)
- real(dp), public [fracsealed\\_cityarea](#)
- real(dp), dimension(:), allocatable, public [l0\\_slope\\_emp](#)
- real(dp), dimension(:, :), allocatable, public [l0\\_gridded\\_lai](#)
- real(dp), dimension(:), allocatable, public [l0\\_slope](#)
- real(dp), dimension(:), allocatable, public [l0\\_asp](#)
- integer(i4), dimension(:, :), allocatable, public [l0\\_soilid](#)
- integer(i4), dimension(:), allocatable, public [l0\\_geounit](#)
- character(256), dimension(:), allocatable, public [dirgridded\\_lai](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_fsealed](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_alpha](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_degdayinc](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_degdaymax](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_degdaynopre](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_degday](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_karstloss](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_fasp](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_petlaicorfactor](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_harsamcoeff](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_prietayalpha](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_aeroresist](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_surfresist](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_froots](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_maxinter](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_kfastflow](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_kslowflow](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_kbaseflow](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_kperco](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_soilmoistfc](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_soilmoistsat](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_soilmoistexp](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_jarvis\\_thresh\\_c1](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_tempthresh](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_unsatthresh](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_sealedthresh](#)
- real(dp), dimension(:, :, :), allocatable, public [l1\\_wiltingpoint](#)



## 17.29.1 Detailed Description

Global variables for mpr only.

TODO: add description

### Authors

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### Date

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## 17.29.2 Variable Documentation

### 17.29.2.1 dirgridded\_lai

```
character(256), dimension(:), allocatable, public mo_mpr_global_variables::dirgridded_lai
```

Definition at line 115 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_read\_config::mpr\_read\_config(), mo\_prepare\_gridded\_lai::prepare\_gridded\_daily\_lai\_data(), and mo\_prepare\_gridded\_lai::prepare\_gridded\_mean\_monthly\_lai\_data().

### 17.29.2.2 fracsealed\_cityarea

```
real(dp), public mo_mpr_global_variables::fracsealed_cityarea
```

Definition at line 92 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_multi\_param\_reg::mpr(), and mo\_mpr\_read\_config::mpr\_read\_config().

### 17.29.2.3 geounitkar

```
integer(i4), dimension(:), allocatable, public mo_mpr_global_variables::geounitkar
```

Definition at line 76 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_multi\_param\_reg::karstic\_layer(), and mo\_read\_wrapper::read\_data().

### 17.29.2.4 geounitlist

```
integer(i4), dimension(:), allocatable, public mo_mpr_global_variables::geounitlist
```

Definition at line 75 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_multi\_param\_reg::baseflow\_param(), mo\_startup::constants\_init(), mo\_multi\_param\_reg::karstic\_layer(), and mo\_read\_wrapper::read\_data().

### 17.29.2.5 horizondepth\_mhm

`real(dp), dimension(:), allocatable, public mo_mpr_global_variables::horizondepth_mhm`

Definition at line 31 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_common_mhm_mrm_restart::check_dimension_consistency()`, `mo_soil_database::generate_soil_database()`, `mo_mhm_eval::mhm_eval()`, `mo_multi_param_reg::mpr()`, `mo_mpr_read_config::mpr_read_config()`, `mo_soil_database::read_soil_lut()`, `mo_init_states::variables_alloc()`, `mo_init_states::variables_default_init()`, `mo_mpr_restart::write_mpr_restart_files()`, and `mo_restart::write_restart_files()`.

### 17.29.2.6 iflag\_soildb

`integer(i4), public mo_mpr_global_variables::iflag_soildb`

Definition at line 29 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_soil_database::generate_soil_database()`, `mo_mpr_startup::i0_check_input()`, `mo_mpr_startup::i0_variable_init()`, `mo_multi_param_reg::mpr()`, `mo_mpr_read_config::mpr_read_config()`, `mo_mpr_soilmoist::mpr_sm()`, `mo_read_wrapper::read_data()`, and `mo_soil_database::read_soil_lut()`.

### 17.29.2.7 inputformat\_gridded\_lai

`character(256), public mo_mpr_global_variables::inputformat_gridded_lai`

Definition at line 81 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_read_config::mpr_read_config()`, and `mo_prepare_gridded_lai::prepare_gridded_daily_lai_data()`.

### 17.29.2.8 i0\_asp

`real(dp), dimension(:), allocatable, public mo_mpr_global_variables::i0_asp`

Definition at line 106 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::i0_check_input()`, `mo_mpr_eval::mpr_eval()`, and `mo_read_wrapper::read_data()`.

### 17.29.2.9 i0\_geounit

`integer(i4), dimension(:), allocatable, public mo_mpr_global_variables::i0_geounit`

Definition at line 109 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::i0_check_input()`, `mo_mpr_eval::mpr_eval()`, and `mo_read_wrapper::read_data()`.

### 17.29.2.10 i0\_gridded\_lai

`real(dp), dimension(:, :), allocatable, public mo_mpr_global_variables::i0_gridded_lai`

Definition at line 102 of file `mo_mpr_global_variables.f90`.

Referenced by mo\_mpr\_startup::l0\_check\_input(), mo\_mpr\_eval::mpr\_eval(), mo\_prepare\_gridded\_lai::prepare\_↔\_gridded\_daily\_lai\_data(), mo\_prepare\_gridded\_lai::prepare\_gridded\_mean\_monthly\_lai\_data(), and mo\_read\_↔\_wrapper::read\_data().

#### 17.29.2.11 l0\_slope

```
real(dp), dimension(:), allocatable, public mo_mpr_global_variables::l0_slope
```

Definition at line 105 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::l0\_check\_input(), mo\_mpr\_startup::l0\_variable\_init(), mo\_mrm\_net\_startup::l11\_↔\_calc\_celerity(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), and mo\_↔\_read\_wrapper::read\_data().

#### 17.29.2.12 l0\_slope\_emp

```
real(dp), dimension(:), allocatable, public mo_mpr_global_variables::l0_slope_emp
```

Definition at line 100 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::l0\_variable\_init(), and mo\_mpr\_eval::mpr\_eval().

#### 17.29.2.13 l0\_soilid

```
integer(i4), dimension(:, :), allocatable, public mo_mpr_global_variables::l0_soilid
```

Definition at line 108 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::l0\_check\_input(), mo\_mpr\_startup::l0\_variable\_init(), mo\_mpr\_eval::mpr\_eval(), and mo\_read\_wrapper::read\_data().

#### 17.29.2.14 l1\_aeroresist

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_aeroresist
```

Definition at line 137 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_↔\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

#### 17.29.2.15 l1\_alpha

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_alpha
```

Definition at line 124 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_↔\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.16 l1\_degday**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_degday
```

Definition at line 130 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.17 l1\_degdayinc**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_degdayinc
```

Definition at line 125 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.18 l1\_degdaymax**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_degdaymax
```

Definition at line 127 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.19 l1\_degdaynopre**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_degdaynopre
```

Definition at line 128 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.20 l1\_fasp**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_fasp
```

Definition at line 132 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.21 l1\_froots**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_froots
```

Definition at line 139 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.22 l1\_fsealed**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_fsealed
```

Definition at line 122 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.23 l1\_harsamcoeff**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_harsamcoeff
```

Definition at line 135 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.24 l1\_jarvis\_thresh\_c1**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_jarvis_thresh_c1
```

Definition at line 151 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.25 l1\_karstloss**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_karstloss
```

Definition at line 131 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.26 l1\_kbaseflow**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_kbaseflow
```

Definition at line 144 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.27 l1\_kfastflow**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_kfastflow
```

Definition at line 142 of file mo\_mpr\_global\_variables.f90.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_mhm_eval::mhm_eval()`, `mo_mpr_eval::mpr_eval()`, `mo_↔ restart::read_restart_states()`, `mo_init_states::variables_default_init()`, and `mo_mpr_restart::write_eff_params()`.

#### 17.29.2.28 l1\_kperco

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_kperco
```

Definition at line 145 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_mhm_eval::mhm_eval()`, `mo_mpr_eval::mpr_eval()`, `mo_↔ restart::read_restart_states()`, `mo_init_states::variables_default_init()`, and `mo_mpr_restart::write_eff_params()`.

#### 17.29.2.29 l1\_kslowflow

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_kslowflow
```

Definition at line 143 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_mhm_eval::mhm_eval()`, `mo_mpr_eval::mpr_eval()`, `mo_↔ restart::read_restart_states()`, `mo_init_states::variables_default_init()`, and `mo_mpr_restart::write_eff_params()`.

#### 17.29.2.30 l1\_maxinter

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_maxinter
```

Definition at line 140 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_mhm_eval::mhm_eval()`, `mo_mpr_eval::mpr_eval()`, `mo_↔ restart::read_restart_states()`, `mo_init_states::variables_default_init()`, and `mo_mpr_restart::write_eff_params()`.

#### 17.29.2.31 l1\_petlaicorfactor

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_petlaicorfactor
```

Definition at line 133 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_mhm_eval::mhm_eval()`, `mo_mpr_eval::mpr_eval()`, `mo_↔ restart::read_restart_states()`, `mo_init_states::variables_default_init()`, and `mo_mpr_restart::write_eff_params()`.

#### 17.29.2.32 l1\_prietayalpha

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_prietayalpha
```

Definition at line 136 of file `mo_mpr_global_variables.f90`.

Referenced by `mo_mpr_startup::init_eff_params()`, `mo_mhm_eval::mhm_eval()`, `mo_mpr_eval::mpr_eval()`, `mo_↔ restart::read_restart_states()`, `mo_init_states::variables_default_init()`, and `mo_mpr_restart::write_eff_params()`.

### 17.29.2.33 l1\_sealedthresh

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_sealedthresh
```

Definition at line 155 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

### 17.29.2.34 l1\_soilmoistexp

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_soilmoistexp
```

Definition at line 149 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

### 17.29.2.35 l1\_soilmoistfc

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_soilmoistfc
```

Definition at line 146 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

### 17.29.2.36 l1\_soilmoistsat

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_soilmoistsat
```

Definition at line 148 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

### 17.29.2.37 l1\_surfresist

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_surfresist
```

Definition at line 138 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

### 17.29.2.38 l1\_tempthresh

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_tempthresh
```

Definition at line 153 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.39 l1\_unsatthresh**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_unsatthresh
```

Definition at line 154 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.40 l1\_wiltingpoint**

```
real(dp), dimension(:, :, :), allocatable, public mo_mpr_global_variables::l1_wiltingpoint
```

Definition at line 157 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_startup::init\_eff\_params(), mo\_mhm\_eval::mhm\_eval(), mo\_mpr\_eval::mpr\_eval(), mo\_restart::read\_restart\_states(), mo\_init\_states::variables\_default\_init(), and mo\_mpr\_restart::write\_eff\_params().

**17.29.2.41 laiboundaries**

```
real(dp), dimension(:), allocatable, public mo_mpr_global_variables::laiboundaries
```

Definition at line 87 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_prepare\_gridded\_lai::prepare\_gridded\_daily\_lai\_data(), mo\_prepare\_gridded\_lai::prepare\_gridded\_mean\_monthly\_lai\_data(), mo\_read\_wrapper::read\_data(), and mo\_restart::write\_restart\_files().

**17.29.2.42 lailut**

```
real(dp), dimension(:, :), allocatable, public mo_mpr_global_variables::lailut
```

Definition at line 89 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_read\_wrapper::read\_data().

**17.29.2.43 laiper**

```
type(period), dimension(:), allocatable, public mo_mpr_global_variables::laiper
```

Definition at line 91 of file mo\_mpr\_global\_variables.f90.

**17.29.2.44 laiunitlist**

```
integer(i4), dimension(:), allocatable, public mo_mpr_global_variables::laiunitlist
```

Definition at line 88 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_read\_wrapper::read\_data().



#### 17.29.2.45 ngeounits

```
integer(i4), public mo_mpr_global_variables::ngeounits
```

Definition at line 74 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_read\_config::mpr\_read\_config(), and mo\_read\_wrapper::read\_data().

#### 17.29.2.46 nlai

```
integer(i4), public mo_mpr_global_variables::nlai
```

Definition at line 86 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_mpr\_startup::init\_eff\_params(), mo\_prepare\_gridded\_lai::prepare\_gridded\_daily\_lai\_data(), mo\_prepare\_gridded\_lai::prepare\_gridded\_mean\_monthly\_lai\_data(), mo\_read\_wrapper::read\_data(), mo\_restart::read\_restart\_states(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

#### 17.29.2.47 nlaiclass

```
integer(i4), public mo_mpr_global_variables::nlaiclass
```

Definition at line 85 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_read\_wrapper::read\_data().

#### 17.29.2.48 nsoilhorizons\_mhm

```
integer(i4), public mo_mpr_global_variables::nsoilhorizons_mhm
```

Definition at line 30 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_soil\_database::generate\_soil\_database(), mo\_mpr\_startup::init\_eff\_params(), mo\_mpr\_startup::io\_check\_input(), mo\_mpr\_startup::io\_variable\_init(), mo\_mhm\_eval::mhm\_eval(), mo\_mhm\_read\_config::mhm\_read\_config(), mo\_multi\_param\_reg::mpr(), mo\_mpr\_read\_config::mpr\_read\_config(), mo\_write\_fluxes\_states::newoutputdataset(), mo\_read\_wrapper::read\_data(), mo\_restart::read\_restart\_states(), mo\_soil\_database::read\_soil\_lut(), mo\_write\_fluxes\_states::updatedataset(), mo\_init\_states::variables\_alloc(), mo\_init\_states::variables\_default\_init(), mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_restart::write\_restart\_files().

#### 17.29.2.49 nsoiltypes

```
integer(i4), public mo_mpr_global_variables::nsoiltypes
```

Definition at line 28 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_soil\_database::generate\_soil\_database(), mo\_mpr\_startup::io\_variable\_init(), and mo\_soil\_database::read\_soil\_lut().

### 17.29.2.50 soildb

```
type(soiltype), public mo_mpr_global_variables::soildb
```

Definition at line 69 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_soil\_database::generate\_soil\_database(), mo\_mpr\_startup::l0\_variable\_init(), mo\_multi\_↔ param\_reg::mpr(), mo\_read\_wrapper::read\_data(), and mo\_soil\_database::read\_soil\_lut().

### 17.29.2.51 tillagedepth

```
real(dp), public mo_mpr_global_variables::tillagedepth
```

Definition at line 26 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_mpr\_read\_config::mpr\_read\_config(), and mo\_soil\_database::read\_soil\_lut().

### 17.29.2.52 timestep\_lai\_input

```
integer(i4), public mo_mpr_global_variables::timestep_lai_input
```

Definition at line 82 of file mo\_mpr\_global\_variables.f90.

Referenced by mo\_common\_datetime\_type::datetimeinfo\_update\_lai\_timestep(), mo\_mpr\_startup::l0\_check\_↔ input(), mo\_mpr\_read\_config::mpr\_read\_config(), mo\_prepare\_gridded\_lai::prepare\_gridded\_daily\_lai\_data(), and mo\_read\_wrapper::read\_data().

## 17.30 mo\_mpr\_pet Module Reference

TODO: add description.

### Functions/Subroutines

- subroutine, public [pet\\_correctbylai](#) (param, nodata, LCOVER0, LAI0, mask0, cell\_id0, upp\_row\_L1, low\_↔ row\_L1, lef\_col\_L1, rig\_col\_L1, nL0\_in\_L1, L1\_petLAIcorFactor)  
*estimate PET correction factor based on LAI at L1*
- subroutine, public [pet\\_correctbyasp](#) (ld0, latitude\_l0, Asp0, param, nodata, fAsp0)  
*correction of PET*
- subroutine, public [priestley\\_taylor\\_alpha](#) (LAI0, param, mask0, nodata, cell\_id0, nL0\_in\_L1, Upp\_row\_L1, Low\_row\_L1, Lef\_col\_L1, Rig\_col\_L1, priestley\_taylor\_alpha1)  
*Regionalization of priestley taylor alpha.*
- subroutine, public [bulksurface\\_resistance](#) (LAI0, param, mask0, nodata, cell\_id0, nL0\_in\_L1, Upp\_row\_L1, Low\_row\_L1, Lef\_col\_L1, Rig\_col\_L1, bulksurface\_resistance1)  
*Regionalization of bulk surface resistance.*

### 17.30.1 Detailed Description

TODO: add description.

This module sets up pet correction factor at level-1 based on LAI

**Authors**

Mehmet Cuneyd Demirel, Simon Stisen

**Date**

May 2017

**17.30.2 Function/Subroutine Documentation****17.30.2.1 bulksurface\_resistance()**

```
subroutine, public mo_mpr_pet::bulksurface_resistance (
    real(dp), dimension(:, :), intent(in) LAI0,
    real(dp), intent(in) param,
    logical, dimension(:, :), intent(in) mask0,
    real(dp), intent(in) nodata,
    integer(i4), dimension(:), intent(in) cell_id0,
    integer(i4), dimension(:), intent(in) nL0_in_L1,
    integer(i4), dimension(:), intent(in) Upp_row_L1,
    integer(i4), dimension(:), intent(in) Low_row_L1,
    integer(i4), dimension(:), intent(in) Lef_col_L1,
    integer(i4), dimension(:), intent(in) Rig_col_L1,
    real(dp), dimension(:, :), intent(out) bulksurface_resistance1 )
```

Regionalization of bulk surface resistance.

estimation of bulk surface resistance Global parameters needed (see mhm\_parameter.nml):

- param(1) = stomatal\_resistance

**Parameters**

|     |   |                                     |
|-----|---|-------------------------------------|
| in  | <i>real(dp), dimension(:, :) :: LAI0</i>                    | LAI at level-0                      |
| in  | <i>real(dp) :: param</i>                                    | - global parameter                  |
| in  | <i>logical, dimension(:, :) :: mask0</i>                    | mask at level 0                     |
| in  | <i>real(dp) :: nodata</i>                                   | - nodata value                      |
| in  | <i>integer(i4), dimension(:) :: cell_id0</i>                | Cell ids of hi res field            |
| in  | <i>integer(i4), dimension(:) :: nL0_in_L1</i>               | number of l0 cells within a l1 cell |
| in  | <i>integer(i4), dimension(:) :: Upp_row_L1</i>              | upper row of a l1 cell in l0 grid   |
| in  | <i>integer(i4), dimension(:) :: Low_row_L1</i>              | lower row of a l1 cell in l0 grid   |
| in  | <i>integer(i4), dimension(:) :: Lef_col_L1</i>              | left col of a l1 cell in l0 grid    |
| in  | <i>integer(i4), dimension(:) :: Rig_col_L1</i>              | right col of a l1 cell in l0 grid   |
| out | <i>real(dp), dimension(:, :) :: bulksurface_resistance1</i> | bulk surface resistance             |

**Authors**

Matthias Zink

**Date**

Apr 2013

Definition at line 408 of file mo\_mpr\_pet.f90.

References `mo_mpr_constants::lai_factor_surfresi`, `mo_mpr_constants::lai_offset_surfresi`, `mo_mpr_constants::max_surfresist`, and `mo_upscaling_operators::upscale_arithmetic_mean()`.

Referenced by `mo_multi_param_reg::mpr()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.30.2.2 pet\_correctbyasp()

```

subroutine, public mo_mpr_pet::pet_correctbyasp (
    integer(i4), dimension(:), intent(in) Id0,
    real(dp), dimension(:), intent(in) latitude_l0,
    real(dp), dimension(:), intent(in) Asp0,
    real(dp), dimension(3), intent(in) param,
    real(dp), intent(in) nodata,
    real(dp), dimension(:), intent(out) fAsp0 )
  
```

correction of PET

Correction of PET based on L0 aspect data. Global parameters needed (see `mhm_parameter.nml`):

- `param(1)` = `minCorrectionFactorPET`
- `param(2)` = `maxCorrectionFactorPET`
- `param(3)` = `aspectTresholdPET`

#### Parameters

|     |  |                            |
|-----|--|----------------------------|
| in  | <i>integer(i4), dimension(:) :: id0</i>      | Level 0 cell id            |
| in  | <i>real(dp), dimension(:) :: latitude_l0</i> | latitude on l0             |
| in  | <i>real(dp), dimension(:) :: Asp0</i>        | [degree] Aspect at Level 0 |
| in  | <i>real(dp), dimension(3) :: param</i>       | process parameters         |
| in  | <i>real(dp) :: nodata</i>                    | - no data value            |
| out | <i>real(dp), dimension(:) :: fAsp0</i>       | PET correction for Aspect  |

**Authors**

Stephan Thober, Rohini Kumar

**Date**

Dec 2012

Definition at line 215 of file mo\_mpr\_pet.f90.

Referenced by mo\_multi\_param\_reg::mpr().

Here is the caller graph for this function:

**17.30.2.3 pet\_correctbylai()**

```

subroutine, public mo_mpr_pet::pet_correctbylai (
    real(dp), dimension(5), intent(in) param,
    real(dp), intent(in) nodata,
    integer(i4), dimension(:), intent(in) LCOVER0,
    real(dp), dimension(:, :), intent(in) LAI0,
    logical, dimension(:, :), intent(in) mask0,
    integer(i4), dimension(:), intent(in) cell_id0,
    integer(i4), dimension(:), intent(in) upp_row_L1,
    integer(i4), dimension(:), intent(in) low_row_L1,
    integer(i4), dimension(:), intent(in) lef_col_L1,
    integer(i4), dimension(:), intent(in) rig_col_L1,
    integer(i4), dimension(:), intent(in) nLO_in_L1,
    real(dp), dimension(:, :), intent(inout) L1_petLAIcorFactor )
  
```

estimate PET correction factor based on LAI at L1

estimate PET correction factor based on LAI at L1 for a given Leaf Area Index field. Global parameters needed (see mhm\_parameter.nml): Process Case 5:

- param(1) = PET\_a\_forest
- param(2) = PET\_a\_impervious
- param(3) = PET\_a\_pervious
- param(4) = PET\_b
- param(5) = PET\_c Example DSF=PET\_a+PET\_b\*(1-exp(PET\_c\*LAI)) Similar to the crop coefficient concept  $Kc=a+b*(1-\exp(c*LAI))$  by Allen, R. G., L. S. Pereira, D. Raes, and M. Smith (1998), Crop evapotranspiration - Guidelines for computing crop water requirements., FAO Irrigation and drainage paper 56. See Chapter 9, Equation 97 <http://www.fao.org/docrep/X0490E/x0490e0f.htm> Date: 17/5/2017

**Parameters**

|    |                                      |                       |
|----|--------------------------------------|-----------------------|
| in | real(dp), dimension(5) :: param      | parameters            |
| in | real(dp) :: nodata                   | - nodata value        |
| in | integer(i4), dimension(:) :: LCOVER0 | Land cover at level 0 |

## Parameters

|         |  |                                    |
|---------|--|------------------------------------|
| in      | <i>real(dp), dimension(:, :) :: LAI0</i>               | LAI at level-0                     |
| in      | <i>logical, dimension(:, :) :: mask0</i>               | mask at L0                         |
| in      | <i>integer(i4), dimension(:) :: cell_id0</i>           | Cell ids of hi res field           |
| in      | <i>integer(i4), dimension(:) :: upp_row_L1</i>         | Upper row of hi res block          |
| in      | <i>integer(i4), dimension(:) :: low_row_L1</i>         | Lower row of hi res block          |
| in      | <i>integer(i4), dimension(:) :: lef_col_L1</i>         | Left column of hi res block        |
| in      | <i>integer(i4), dimension(:) :: rig_col_L1</i>         | Right column of hi res block       |
| in      | <i>integer(i4), dimension(:) :: nLO_in_L1</i>          | Number of L0 cells within a L1 cel |
| in, out | <i>real(dp), dimension(:, :) :: L1_petLAIcorFactor</i> | pet cor factor at level-1          |

## Authors

M. Cunejd Demirel and Simon Stisen from GEUS.dk

## Date

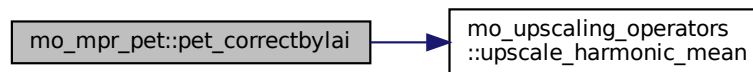
May, 2017

Definition at line 81 of file mo\_mpr\_pet.f90.

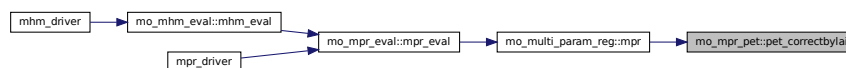
References mo\_upscaling\_operators::upscale\_harmonic\_mean().

Referenced by mo\_multi\_param\_reg::mpr().

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.30.2.4 priestley\_taylor\_alpha()

```

subroutine, public mo_mpr_pet::priestley_taylor_alpha (
  real(dp), dimension(:, :), intent(in) LAI0,
  real(dp), dimension(:), intent(in) param,
  logical, dimension(:, :), intent(in) mask0,
  real(dp), intent(in) nodata,
  integer(i4), dimension(:), intent(in) cell_id0,

```

```

integer(i4), dimension(:), intent(in) nL0_in_L1,
integer(i4), dimension(:), intent(in) Upp_row_L1,
integer(i4), dimension(:), intent(in) Low_row_L1,
integer(i4), dimension(:), intent(in) Lef_col_L1,
integer(i4), dimension(:), intent(in) Rig_col_L1,
real(dp), dimension(:, :), intent(out) priestley_taylor_alpha )

```

Regionalization of priestley taylor alpha.

estimation of priestley taylor alpha Global parameters needed (see mhm\_parameter.nml):

- param(1) = PriestleyTaylorCoeff
- param(2) = PriestleyTaylorLAlcorr

#### Parameters

|     |   |                                     |
|-----|---|-------------------------------------|
| in  | <i>real(dp), dimension(:, :) :: LAI0</i>                    | LAI at level-0                      |
| in  | <i>real(dp), dimension(:) :: param</i>                      | input parameter                     |
| in  | <i>logical, dimension(:, :) :: mask0</i>                    | mask at level 0                     |
| in  | <i>real(dp) :: nodata</i>                                   | - nodata value                      |
| in  | <i>integer(i4), dimension(:) :: cell_id0</i>                | Cell ids of hi res field            |
| in  | <i>integer(i4), dimension(:) :: nL0_in_L1</i>               | number of l0 cells within a l1 cell |
| in  | <i>integer(i4), dimension(:) :: Upp_row_L1</i>              | upper row of a l1 cell in l0 grid   |
| in  | <i>integer(i4), dimension(:) :: Low_row_L1</i>              | lower row of a l1 cell in l0 grid   |
| in  | <i>integer(i4), dimension(:) :: Lef_col_L1</i>              | left col of a l1 cell in l0 grid    |
| in  | <i>integer(i4), dimension(:) :: Rig_col_L1</i>              | right col of a l1 cell in l0 grid   |
| out | <i>real(dp), dimension(:, :) :: priestley_taylor_alpha1</i> | bulk surface resistance             |

#### Authors

Matthias Zink

#### Date

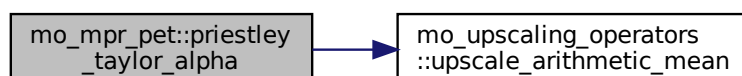
Apr 2013

Definition at line 312 of file mo\_mpr\_pet.f90.

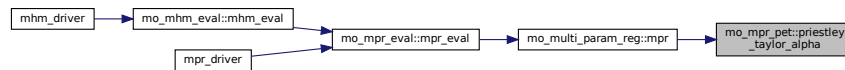
References mo\_upscaling\_operators::upscale\_arithmetic\_mean().

Referenced by mo\_multi\_param\_reg::mpr().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.31 mo\_mpr\_read\_config Module Reference

read mpr config

### Functions/Subroutines

- subroutine, public [mpr\\_read\\_config](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param)  
*Read the general config of mpr.*

#### 17.31.1 Detailed Description

read mpr config

This module contains all mpr subroutines related to reading the mpr configuration from file.

#### Authors

Stephan Thober

#### Date

Aug 2015

#### 17.31.2 Function/Subroutine Documentation

##### 17.31.2.1 mpr\_read\_config()

```

subroutine, public mo_mpr_read_config::mpr_read_config (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist,
    character(*), intent(in) file_namelist_param,
    integer, intent(in) unamelist_param )
  
```

Read the general config of mpr.

Depending on the variable `mrm_coupling_config`, the mRM config is either read from `mrm.nml` and parameters from `mrm_parameter.nml` or copied from mHM.

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>character(*) :: file_namelist</i>       |  |
| in | <i>integer :: unamelist</i>                |  |
| in | <i>character(*) :: file_namelist_param</i> |  |
| in | <i>integer :: unamelist_param</i>          |  |



**Authors**

Stephan Thober

**Date**

Aug 2015

Definition at line 54 of file mo\_mpr\_read\_config.f90.

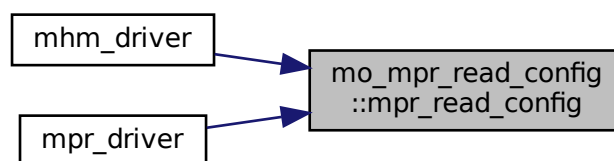
References mo\_mpr\_global\_variables::dirgridded\_lai, mo\_common\_variables::domainmeta, mo\_common\_↵ constants::eps\_dp, mo\_mpr\_global\_variables::fracsealed\_cityarea, mo\_common\_variables::global\_parameters, mo\_common\_variables::global\_parameters\_name, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_mpr\_↵ \_global\_variables::iflag\_soildb, mo\_common\_functions::in\_bound(), mo\_mpr\_global\_variables::inputformat\_↵ gridded\_lai, mo\_mpr\_constants::maxgeounit, mo\_common\_constants::maxnodomains, mo\_mpr\_constants\_↵ ::maxnosoilhorizons, mo\_common\_constants::ncolpars, mo\_mpr\_global\_variables::ngeounits, mo\_common\_↵ constants::nodata\_dp, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_common\_variables::processmatrix, mo\_mpr\_global\_variables::tillagedepth, and mo\_mpr\_global\_variables::timestep\_lai\_input.

Referenced by mhm\_driver(), and mpr\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.32 mo\_mpr\_restart Module Reference

reading and writing states, fluxes and configuration for restart of mHM.

**Data Types**

- interface [unpack\\_field\\_and\\_write](#)

*TODO: add description.*

## Functions/Subroutines

- subroutine, public [write\\_mpr\\_restart\\_files](#) (OutFile)  
*write restart files for each domain*
- subroutine, public [write\\_eff\\_params](#) (mask1, s1, e1, rows1, cols1, soil1, lcscenes, lais, nc)  
*TODO: add description.*
- subroutine [unpack\\_field\\_and\\_write\\_1d\\_i4](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_1d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_2d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_3d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)

### 17.32.1 Detailed Description

reading and writing states, fluxes and configuration for restart of mHM.

routines are seperated for reading and writing variables for:

- states and fluxes, and
- configuration. Reading of L11 configuration is also seperated from the rest, since it is only required when routing is activated.

#### Authors

Stephan Thober

#### Date

Jul 2013

### 17.32.2 Function/Subroutine Documentation

#### 17.32.2.1 [unpack\\_field\\_and\\_write\\_1d\\_dp\(\)](#)

```
subroutine mo_mpr_restart::unpack_field_and_write_1d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name )
```

Definition at line 413 of file mo\_mpr\_restart.f90.

#### 17.32.2.2 [unpack\\_field\\_and\\_write\\_1d\\_i4\(\)](#)

```
subroutine mo_mpr_restart::unpack_field_and_write_1d_i4 (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    integer(i4), intent(in) fill_value,
    integer(i4), dimension(:), intent(in) data,
```

```

logical, dimension(:, :), intent(in) mask,
character(*), intent(in), optional var_long_name )

```

Definition at line 368 of file mo\_mpr\_restart.f90.

### 17.32.2.3 unpack\_field\_and\_write\_2d\_dp()

```

subroutine mo_mpr_restart::unpack_field_and_write_2d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:, :), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name )

```

Definition at line 458 of file mo\_mpr\_restart.f90.

### 17.32.2.4 unpack\_field\_and\_write\_3d\_dp()

```

subroutine mo_mpr_restart::unpack_field_and_write_3d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:, :, :), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name )

```

Definition at line 515 of file mo\_mpr\_restart.f90.

### 17.32.2.5 write\_eff\_params()

```

subroutine, public mo_mpr_restart::write_eff_params (
    logical, dimension(:, :), intent(in), allocatable mask1,
    integer(i4), intent(in) s1,
    integer(i4), intent(in) e1,
    type(ncdimension), intent(in) rows1,
    type(ncdimension), intent(in) cols1,
    type(ncdimension), intent(in) soil1,
    type(ncdimension), intent(in) lcscenes,
    type(ncdimension), intent(in) lais,
    type(ncdataset), intent(inout) nc )

```

TODO: add description.

TODO: add description

#### Parameters

|    |   |                        |
|----|---|------------------------|
| in | <i>logical, dimension(:, :) :: mask1</i>                        | mask at level 1        |
| in | <i>integer(i4) :: s1</i>  | start index at level 1 |
| in | <i>integer(i4) :: e1</i>  | end index at level 1   |
| in | <i>type(NcDimension) :: rows1, cols1, soil1, lcscenes, lais</i> |                        |

## Parameters

|         |   |  |
|---------|---|--|
| in      | <code>type(NcDimension) :: rows1, cols1, soil1, lcscenes, lais</code> |  |
| in      | <code>type(NcDimension) :: rows1, cols1, soil1, lcscenes, lais</code> |  |
| in      | <code>type(NcDimension) :: rows1, cols1, soil1, lcscenes, lais</code> |  |
| in      | <code>type(NcDimension) :: rows1, cols1, soil1, lcscenes, lais</code> |  |
| in, out | <code>type(NcDataset) :: nc</code>                                    |  |

## Authors

Robert Schweppe

## Date

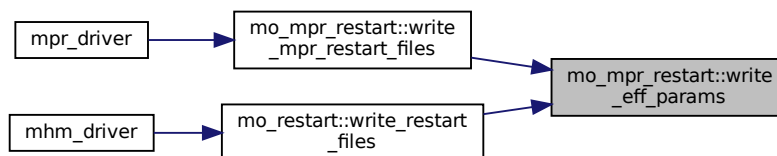
Jun 2018

Definition at line 209 of file `mo_mpr_restart.f90`.

References `mo_mpr_global_variables::l1_aeroresist`, `mo_mpr_global_variables::l1_alpha`, `mo_mpr_global_variables::l1_degday`, `mo_mpr_global_variables::l1_degdayinc`, `mo_mpr_global_variables::l1_degdaymax`, `mo_mpr_global_variables::l1_degdaynopre`, `mo_mpr_global_variables::l1_fasp`, `mo_mpr_global_variables::l1_froots`, `mo_mpr_global_variables::l1_fsealed`, `mo_mpr_global_variables::l1_harsamcoeff`, `mo_mpr_global_variables::l1_jarvis_thresh_c1`, `mo_mpr_global_variables::l1_karstloss`, `mo_mpr_global_variables::l1_kbaseflow`, `mo_mpr_global_variables::l1_kfastflow`, `mo_mpr_global_variables::l1_kperco`, `mo_mpr_global_variables::l1_kslowflow`, `mo_mpr_global_variables::l1_maxinter`, `mo_mpr_global_variables::l1_petlaicofactor`, `mo_mpr_global_variables::l1_prietayalpha`, `mo_mpr_global_variables::l1_sealedthresh`, `mo_mpr_global_variables::l1_soilmoistexp`, `mo_mpr_global_variables::l1_soilmoistfc`, `mo_mpr_global_variables::l1_soilmoistsat`, `mo_mpr_global_variables::l1_surfresist`, `mo_mpr_global_variables::l1_tempthresh`, `mo_mpr_global_variables::l1_unsatthresh`, `mo_mpr_global_variables::l1_wiltingpoint`, `mo_common_variables::lc_year_end`, `mo_common_variables::lc_year_start`, `mo_common_constants::nodata_dp`, `mo_common_constants::nodata_i4`, and `mo_common_variables::processmatrix`.

Referenced by `write_mpr_restart_files()`, and `mo_restart::write_restart_files()`.

Here is the caller graph for this function:



### 17.32.2.6 write\_mpr\_restart\_files()

```

subroutine, public mo_mpr_restart::write_mpr_restart_files (
    character(256), dimension(:), intent(in) OutFile )
  
```

write restart files for each domain

write restart files for each domain. For each domain three restart files are written. These are `xxx_states.nc`, `xxx_L11_config.nc`, and `xxx_config.nc` (`xxx` being the three digit domain index). If a variable is added here, it should also be added in the read restart routines below. **ADDITIONAL INFORMATION** `write_restart`

## Parameters

|    |  |                             |
|----|--|-----------------------------|
| in | <i>character(256), dimension(:) :: OutFile</i> | Output Path for each domain |
|----|--|-----------------------------|

## Authors

Stephan Thober

## Date

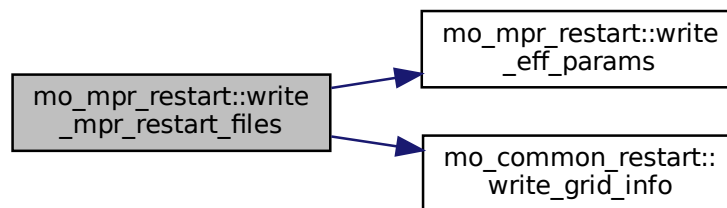
Jun 2014

Definition at line 98 of file mo\_mpr\_restart.f90.

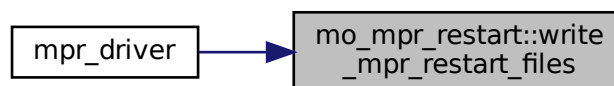
References mo\_common\_variables::domainmeta, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_common\_↵  
 \_constants::laivarname, mo\_common\_constants::landcoverperiodsvarname, mo\_common\_variables::lc\_year↵  
 \_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::level1, mo\_mpr\_global\_variables::nlai,  
 mo\_common\_variables::nlcoverscene, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_common\_constants↵  
 ::soilhorizonsvarname, write\_eff\_params(), and mo\_common\_restart::write\_grid\_info().

Referenced by mpr\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.33 mo\_mpr\_runoff Module Reference

multiscale parameter regionalization for runoff generation

## Functions/Subroutines

- subroutine, public `mpr_runoff` (LCOVER0, mask0, SMs\_FC0, slope\_emp0, KsVar\_H0, param, cell\_id0, upp\_row\_L1, low\_row\_L1, lef\_col\_L1, rig\_col\_L1, nL0\_in\_L1, L1\_HL1, L1\_K0, L1\_K1, L1\_alpha)  
*multiscale parameter regionalization for runoff parameters*

### 17.33.1 Detailed Description

multiscale parameter regionalization for runoff generation

This contains the routine for multiscale parameter regionalization of the runoff parametrization.

#### Authors

Stephan Thober, Rohini Kumar

#### Date

Dec 2012

### 17.33.2 Function/Subroutine Documentation

#### 17.33.2.1 `mpr_runoff()`

```
subroutine, public mo_mpr_runoff::mpr_runoff (
    integer(i4), dimension(:), intent(in) LCOVER0,
    logical, dimension(:, :), intent(in) mask0,
    real(dp), dimension(:), intent(in) SMs_FC0,
    real(dp), dimension(:), intent(in) slope_emp0,
    real(dp), dimension(:), intent(in) KsVar_H0,
    real(dp), dimension(5), intent(in) param,
    integer(i4), dimension(:), intent(in) cell_id0,
    integer(i4), dimension(:), intent(in) upp_row_L1,
    integer(i4), dimension(:), intent(in) low_row_L1,
    integer(i4), dimension(:), intent(in) lef_col_L1,
    integer(i4), dimension(:), intent(in) rig_col_L1,
    integer(i4), dimension(:), intent(in) nL0_in_L1,
    real(dp), dimension(:), intent(out) L1_HL1,
    real(dp), dimension(:), intent(out) L1_K0,
    real(dp), dimension(:), intent(out) L1_K1,
    real(dp), dimension(:), intent(out) L1_alpha )
```

multiscale parameter regionalization for runoff parameters

Perform the multiscale parameter regionalization for runoff global parameters (see `mhm_parameter.nml`). These are the following five parameters:

- param(1) = interflowStorageCapacityFactor
- param(2) = interflowRecession\_slope
- param(3) = fastInterflowRecession\_forest
- param(4) = slowInterflowRecession\_Ks
- param(5) = exponentSlowInterflow

## Parameters

|     |  |  |
|-----|--|--|
| in  | <i>integer(i4), dimension(:) :: LCOVER0</i>    | land cover at level 0                      |
| in  | <i>logical, dimension(:, :) :: mask0</i>       | mask at Level 0                            |
| in  | <i>real(dp), dimension(:) :: SMs_FC0</i>       | [-] soil mositure deficit from field       |
| in  | <i>real(dp), dimension(:) :: slope_emp0</i>    | empirical quantile values F(slope)         |
| in  | <i>real(dp), dimension(:) :: KsVar_H0</i>      | [-] relative variability of saturated      |
| in  | <i>real(dp), dimension(5) :: param</i>         | global parameters                          |
| in  | <i>integer(i4), dimension(:) :: cell_id0</i>   | Cell ids of hi res field                   |
| in  | <i>integer(i4), dimension(:) :: upp_row_L1</i> | Upper row of hi res block                  |
| in  | <i>integer(i4), dimension(:) :: low_row_L1</i> | Lower row of hi res block                  |
| in  | <i>integer(i4), dimension(:) :: lef_col_L1</i> | Left column of hi res block                |
| in  | <i>integer(i4), dimension(:) :: rig_col_L1</i> | Right column of hi res block               |
| in  | <i>integer(i4), dimension(:) :: nL0_in_L1</i>  | Number of L0 cells within a L1 cell        |
| in  | <i>real(dp) :: c2TSTu</i>                      | unit transformations                       |
| out | <i>real(dp), dimension(:) :: L1_HL1</i>        | [10 <sup>-3</sup> m] Threshold water depth |
| out | <i>real(dp), dimension(:) :: L1_K0</i>         | [10 <sup>-3</sup> m] Recession coefficient |
| out | <i>real(dp), dimension(:) :: L1_K1</i>         | [10 <sup>-3</sup> m] Recession coefficient |
| out | <i>real(dp), dimension(:) :: L1_alpha</i>      | [1] Exponent for the upper reservoir       |

## Authors

Stephan Thober, Rohini Kumar

## Date

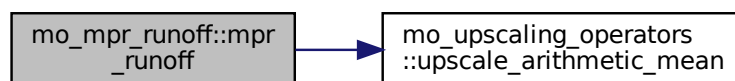
Dec 2012

Definition at line 75 of file mo\_mpr\_runoff.f90.

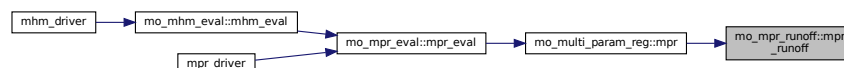
References `mo_common_constants::nodata_dp`, `mo_common_constants::nodata_i4`, and `mo_upscaling_operators::upscale_arithmetic_mean()`.

Referenced by `mo_multi_param_reg::mpr()`.

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.34 mo\_mpr\_smhorizons Module Reference

setting up the soil moisture horizons

### Functions/Subroutines

- subroutine, public `mpr_smhorizons` (`param`, `processMatrix`, `iFlag_soil`, `nHorizons_mHM`, `HorizonDepth`, `LCOVER0`, `soilID0`, `nHorizons`, `nTillHorizons`, `thetaS_till`, `thetaFC_till`, `thetaPW_till`, `thetaS`, `thetaFC`, `thetaPW`, `Wd`, `Db`, `DbM`, `RZdepth`, `mask0`, `cell_id0`, `upp_row_L1`, `low_row_L1`, `lef_col_L1`, `rig_col_L1`, `nL0_in_L1`, `L1_beta`, `L1_SMs`, `L1_FC`, `L1_PW`, `L1_fRoots`)

*upscale soil moisture horizons*

### 17.34.1 Detailed Description

setting up the soil moisture horizons

This module sets up the soil moisture horizons

#### Authors

Stephan Thober, Rohini Kumar

#### Date

Dec 2012

### 17.34.2 Function/Subroutine Documentation

#### 17.34.2.1 mpr\_smhorizons()

```
subroutine, public mo_mpr_smhorizons::mpr_smhorizons (
    real(dp), dimension(:), intent(in) param,
    integer(i4), dimension(:, :), intent(in) processMatrix,
    integer(i4), intent(in) iFlag_soil,
    integer(i4), intent(in) nHorizons_mHM,
    real(dp), dimension(:), intent(in) HorizonDepth,
    integer(i4), dimension(:), intent(in) LCOVER0,
    integer(i4), dimension(:, :), intent(in) soilID0,
    integer(i4), dimension(:), intent(in) nHorizons,
    integer(i4), dimension(:), intent(in) nTillHorizons,
    real(dp), dimension(:, :, :), intent(in) thetaS_till,
    real(dp), dimension(:, :, :), intent(in) thetaFC_till,
    real(dp), dimension(:, :, :), intent(in) thetaPW_till,
    real(dp), dimension(:, :), intent(in) thetaS,
    real(dp), dimension(:, :), intent(in) thetaFC,
    real(dp), dimension(:, :), intent(in) thetaPW,
    real(dp), dimension(:, :, :), intent(in) Wd,
    real(dp), dimension(:, :, :), intent(in) Db,
    real(dp), dimension(:, :), intent(in) DbM,
    real(dp), dimension(:), intent(in) RZdepth,
    logical, dimension(:, :), intent(in) mask0,
    integer(i4), dimension(:), intent(in) cell_id0,
    integer(i4), dimension(:), intent(in) upp_row_L1,
```



```

integer(i4), dimension(:), intent(in) low_row_L1,
integer(i4), dimension(:), intent(in) lef_col_L1,
integer(i4), dimension(:), intent(in) rig_col_L1,
integer(i4), dimension(:), intent(in) nLO_in_L1,
real(dp), dimension(:, :), intent(inout) L1_beta,
real(dp), dimension(:, :), intent(inout) L1_SMs,
real(dp), dimension(:, :), intent(inout) L1_FC,
real(dp), dimension(:, :), intent(inout) L1_PW,
real(dp), dimension(:, :), intent(inout) L1_fRoots )

```

upscale soil moisture horizons

calculate soil properties at the level 1. Global parameters needed (see mhm\_parameter.nml):

- param(1) = rootFractionCoefficient\_forest
- param(2) = rootFractionCoefficient\_impervious
- param(3) = rootFractionCoefficient\_pervious
- param(4) = infiltrationShapeFactor

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: param</i>                | parameters   |
| in | <i>integer(i4), dimension(:, : ) :: processMatrix</i> | - matrix specifying user defined processes                                     |
| in | <i>integer(i4) :: iFlag_soil</i>                      | - flags for handling multiple soil databases                                   |
| in | <i>integer(i4) :: nHorizons_mHM</i>                   | - number of horizons to model  |
| in | <i>real(dp), dimension(:) :: HorizonDepth</i>         | [10 <sup>-3</sup> m] horizon depth from surface, positive downwards            |
| in | <i>integer(i4), dimension(:) :: LCOVER0</i>           | Land cover at level 0  |
| in | <i>integer(i4), dimension(:, : ) :: soilID0</i>       | soil ID at level 0   |
| in | <i>integer(i4), dimension(:) :: nHorizons</i>         | horizons per soil type   |
| in | <i>integer(i4), dimension(:) :: nTillHorizons</i>     | Number of Tillage horizons   |
| in | <i>real(dp), dimension(:, :, : ) :: thetaS_till</i>   | saturated water content of soil horizons upto tillage depth, f(OM, management) |
| in | <i>real(dp), dimension(:, :, : ) :: thetaFC_till</i>  | Field capacity of tillage layers; LUC dependent, f(OM, management)             |
| in | <i>real(dp), dimension(:, :, : ) :: thetaPW_till</i>  | Permanent wilting point of tillage layers; LUC dependent, f(OM, management)    |
| in | <i>real(dp), dimension(:, : ) :: thetaS</i>           | saturated water content of soil horizons after tillage depth                   |
| in | <i>real(dp), dimension(:, : ) :: thetaFC</i>          | Field capacity of deeper layers  |
| in | <i>real(dp), dimension(:, : ) :: thetaPW</i>          | Permanent wilting point of deeper layers                                       |
| in | <i>real(dp), dimension(:, :, : ) :: Wd</i>            | weights of mHM Horizons according to horizons provided in soil database        |
| in | <i>real(dp), dimension(:, :, : ) :: Db</i>            | Bulk density   |
| in | <i>real(dp), dimension(:, : ) :: DbM</i>              | mineral Bulk density   |
| in | <i>real(dp), dimension(:) :: RZdepth</i>              | [mm] Total soil depth  |
| in | <i>logical, dimension(:, : ) :: mask0</i>             | mask at L0   |
| in | <i>integer(i4), dimension(:) :: cell_id0</i>          | Cell ids of hi res field   |
| in | <i>integer(i4), dimension(:) :: upp_row_L1</i>        | Upper row of hi res block  |
| in | <i>integer(i4), dimension(:) :: low_row_L1</i>        | Lower row of hi res block  |
| in | <i>integer(i4), dimension(:) :: lef_col_L1</i>        | Left column of hi res block  |
| in | <i>integer(i4), dimension(:) :: rig_col_L1</i>        | Right column of hi res block   |

## Parameters

|         |   |  |
|---------|---|--|
| in      | <i>integer(i4), dimension(:) :: nL0_in_L1</i> | Number of L0 cells within a L1 cel   |
| in, out | <i>real(dp), dimension(:, :) :: L1_beta</i>   | Parameter that determines the relative contribution to SM, upscaled Bulk density |
| in, out | <i>real(dp), dimension(:, :) :: L1_SMs</i>    | [10 <sup>-3</sup> m] depth of saturated SM cont                                  |
| in, out | <i>real(dp), dimension(:, :) :: L1_FC</i>     | [10 <sup>-3</sup> m] field capacity  |
| in, out | <i>real(dp), dimension(:, :) :: L1_PW</i>     | [10 <sup>-3</sup> m] permanent wilting point                                     |
| in, out | <i>real(dp), dimension(:, :) :: L1_fRoots</i> | fraction of roots in soil horizons   |

## Authors

Luis Samaniego, Rohini Kumar, Stephan Thober

## Date

Dec 2012

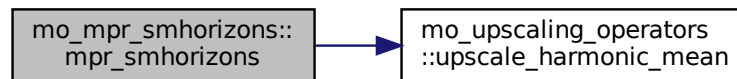
in this case the second dimension of soilld0 = 1

Definition at line 121 of file mo\_mpr\_smhorizons.f90.

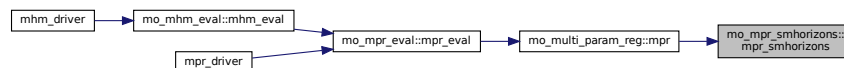
References mo\_common\_constants::nodata\_dp, and mo\_upscaling\_operators::upscale\_harmonic\_mean().

Referenced by mo\_multi\_param\_reg::mpr().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.35 mo\_mpr\_soilmoist Module Reference

Multiscale parameter regionalization (MPR) for soil moisture.

### Functions/Subroutines

- subroutine, public [mpr\\_sm](#) (param, processMatrix, is\_present, nHorizons, nTillHorizons, sand, clay, DbM, ID0, soilld0, LCover0, thetaS\_till, thetaFC\_till, thetaPW\_till, thetaS, thetaFC, thetaPW, Ks, Db, KsVar\_H0, KsVar\_V0, SMs\_FC0)

- multiscale parameter regionalization for soil moisture*
- elemental pure subroutine `pwp` (Genu\_Mual\_n, Genu\_Mual\_alpha, thetaS, thetaPWP)  
*Permanent Wilting point.*
- elemental pure subroutine `field_cap` (thetaFC, Ks, thetaS, Genu\_Mual\_n)  
*calculates the field capacity*
- subroutine `genuchten` (thetaS, Genu\_Mual\_n, Genu\_Mual\_alpha, param, sand, clay, Db)  
*calculates the Genuchten shape parameter*
- subroutine `hydro_cond` (KS, param, sand, clay)  
*calculates the hydraulic conductivity Ks*

### 17.35.1 Detailed Description

Multiscale parameter regionalization (MPR) for soil moisture.

This module contains all routines required for parametrizing soil moisture processes.

#### Authors

Stephan Thober, Rohini Kumar

#### Date

Dec 2012

### 17.35.2 Function/Subroutine Documentation

#### 17.35.2.1 field\_cap()

```
elemental pure subroutine mo_mpr_soilmoist::field_cap (
    real(dp), intent(out) thetaFC,
    real(dp), intent(in) Ks,
    real(dp), intent(in) thetaS,
    real(dp), intent(in) Genu_Mual_n )
```

calculates the field capacity

estimate Field capacity; FC – Flux based approach (Twarakavi, et. al. 2009, WRR) According to the above reference FC is defined as the soil water content at which the drainage from a profile ceases under natural conditions. Since drainage from a soil profile in a simulation never becomes zero, we assume that drainage ceases when the bottom flux from the soil reaches a value that is equivalent to the minimum amount of precipitation that could be recorded (i.e. 0.01 cm/d == 1 mm/d). It is assumed that ThetaR = 0.0\_dp ADDITIONAL INFORMATION Twarakavi, et. al. 2009, WRR

#### Parameters

|     |                                |                                    |
|-----|--------------------------------|------------------------------------|
| out | <i>real(dp) :: thetaFC</i>     | - Field capacity                   |
| in  | <i>real(dp) :: Ks</i>          | - saturated hydraulic conductivity |
| in  | <i>real(dp) :: thetaS</i>      | - saturated water content          |
| in  | <i>real(dp) :: Genu_Mual_n</i> | - Genuchten shape parameter        |

**Authors**

Stephan Thober, Rohini Kumar

**Date**

Dec 2012

Definition at line 564 of file mo\_mpr\_soilmoist.f90.

References mo\_mpr\_constants::field\_cap\_c1, and mo\_mpr\_constants::field\_cap\_c2.

Referenced by mpr\_sm().

Here is the caller graph for this function:

**17.35.2.2 genuchten()**

```

subroutine mo_mpr_soilmoist::genuchten (
    real(dp), intent(out) thetaS,
    real(dp), intent(out) Genu_Mual_n,
    real(dp), intent(out) Genu_Mual_alpha,
    real(dp), dimension(6), intent(in) param,
    real(dp), intent(in) sand,
    real(dp), intent(in) clay,
    real(dp), intent(in) Db )
  
```

calculates the Genuchten shape parameter

estimate SMS\_till &amp; van Genuchten's shape parameter (n) (Zacharias et al, 2007, soil Phy.) Global parameters needed (see mhm\_parameter.nml):

- param( 1) = PTF\_lower66\_5\_constant
- param( 2) = PTF\_lower66\_5\_clay
- param( 3) = PTF\_lower66\_5\_Db
- param( 4) = PTF\_higher66\_5\_constant
- param( 5) = PTF\_higher66\_5\_clay
- param( 6) = PTF\_higher66\_5\_Db ADDITIONAL INFORMATION Zacharias et al, 2007, soil Phy.

**Parameters**

|     |  |   |
|-----|--|---|
| out | <i>real(dp) :: thetaS</i>              | - saturated water content                           |
| out | <i>real(dp) :: Genu_Mual_n</i>         | - van Genuchten shape parameter                     |
| out | <i>real(dp) :: Genu_Mual_alpha</i>     | - van Genuchten shape parameter                     |
| in  | <i>real(dp), dimension(6) :: param</i> | parameters  |
| in  | <i>real(dp) :: sand</i>                | - [%] sand content                                  |
| in  | <i>real(dp) :: clay</i>                | - [%] clay content                                  |
| in  | <i>real(dp) :: Db</i>                  | - [10 <sup>3</sup> kg/m <sup>3</sup> ] bulk density |

**Authors**

Stephan Thober, Rohini Kumar

**Date**

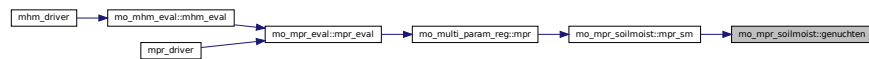
Dec 2012

Definition at line 630 of file mo\_mpr\_soilmoist.f90.

References mo\_mpr\_constants::vgenuchten\_sandtresh, mo\_mpr\_constants::vgenuchtenn\_c1, mo\_mpr\_constants::vgenuchtenn\_c10, mo\_mpr\_constants::vgenuchtenn\_c11, mo\_mpr\_constants::vgenuchtenn\_c12, mo\_mpr\_constants::vgenuchtenn\_c13, mo\_mpr\_constants::vgenuchtenn\_c14, mo\_mpr\_constants::vgenuchtenn\_c15, mo\_mpr\_constants::vgenuchtenn\_c16, mo\_mpr\_constants::vgenuchtenn\_c17, mo\_mpr\_constants::vgenuchtenn\_c18, mo\_mpr\_constants::vgenuchtenn\_c2, mo\_mpr\_constants::vgenuchtenn\_c3, mo\_mpr\_constants::vgenuchtenn\_c4, mo\_mpr\_constants::vgenuchtenn\_c5, mo\_mpr\_constants::vgenuchtenn\_c6, mo\_mpr\_constants::vgenuchtenn\_c7, mo\_mpr\_constants::vgenuchtenn\_c8, and mo\_mpr\_constants::vgenuchtenn\_c9.

Referenced by mpr\_sm().

Here is the caller graph for this function:

**17.35.2.3 hydro\_cond()**

```

subroutine mo_mpr_soilmoist::hydro_cond (
    real(dp), intent(out) KS,
    real(dp), dimension(4), intent(in) param,
    real(dp), intent(in) sand,
    real(dp), intent(in) clay )
  
```

calculates the hydraulic conductivity Ks

By default save this value of Ks, particularly for the deeper layers where OM content plays relatively low or no role  
Global parameters needed (see mhm\_parameter.nml):

- param(1) = PTF\_Ks\_constant
  - param(2) = PTF\_Ks\_sand
  - param(3) = PTF\_Ks\_clay
  - param(4) = PTF\_Ks\_curveSlope
- ADDITIONAL INFORMATION Written, Stephan Thober, Dec 2012

**Parameters**

|     |  |                    |
|-----|--|--------------------|
| out | <i>real(dp) :: KS</i>                  |                    |
| in  | <i>real(dp), dimension(4) :: param</i> |                    |
| in  | <i>real(dp) :: sand</i>                | - [%] sand content |
| in  | <i>real(dp) :: clay</i>                | - [%] clay content |

**Authors**

Stephan Thober, Rohini Kumar

**Date**

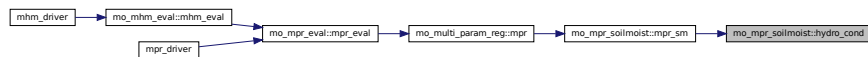
Dec 2012

Definition at line 740 of file mo\_mpr\_soilmoist.f90.

References mo\_mpr\_constants::ks\_c.

Referenced by mpr\_sm().

Here is the caller graph for this function:

**17.35.2.4 mpr\_sm()**

```

subroutine, public mo_mpr_soilmoist::mpr_sm (
    real(dp), dimension(13), intent(in) param,
    integer(i4), dimension(:, :), intent(in) processMatrix,
    integer(i4), dimension(:, :), intent(in) is_present,
    integer(i4), dimension(:, :), intent(in) nHorizons,
    integer(i4), dimension(:, :), intent(in) nTillHorizons,
    real(dp), dimension(:, :), intent(in) sand,
    real(dp), dimension(:, :), intent(in) clay,
    real(dp), dimension(:, :), intent(in) DbM,
    integer(i4), dimension(:, :), intent(in) ID0,
    integer(i4), dimension(:, :), intent(in) soilId0,
    integer(i4), dimension(:, :), intent(in) LCover0,
    real(dp), dimension(:, :, :), intent(out) thetaS_till,
    real(dp), dimension(:, :, :), intent(out) thetaFC_till,
    real(dp), dimension(:, :, :), intent(out) thetaPW_till,
    real(dp), dimension(:, :), intent(out) thetaS,
    real(dp), dimension(:, :), intent(out) thetaFC,
    real(dp), dimension(:, :), intent(out) thetaPW,
    real(dp), dimension(:, :, :), intent(out) Ks,
    real(dp), dimension(:, :, :), intent(out) Db,
    real(dp), dimension(:, :), intent(out) KsVar_H0,
    real(dp), dimension(:, :), intent(out) KsVar_V0,
    real(dp), dimension(:, :), intent(out) SMS_FC0 )
  
```

multiscale parameter regionalization for soil moisture

This subroutine is a wrapper around all soil moisture parameter routines. This subroutine requires 13 parameters. These parameters have to correspond to the parameters in the original parameter array at the following locations: 10-12, 13-18, 27-30. Global parameters needed (see mhm\_parameter.nml):

- param( 1) = orgMatterContent\_forest
- param( 2) = orgMatterContent\_impervious
- param( 3) = orgMatterContent\_pervious

- param( 4) = PTF\_lower66\_5\_constant
- param( 5) = PTF\_lower66\_5\_clay
- param( 6) = PTF\_lower66\_5\_Db
- param( 7) = PTF\_higher66\_5\_constant
- param( 8) = PTF\_higher66\_5\_clay
- param( 9) = PTF\_higher66\_5\_Db
- param(10) = PTF\_Ks\_constant
- param(11) = PTF\_Ks\_sand
- param(12) = PTF\_Ks\_clay
- param(13) = PTF\_Ks\_curveSlope

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>real(dp), dimension(13) :: param</i>             | global parameters   |
| in  | <i>integer(i4), dimension(:) :: is_present</i>      | indicates whether soiltype is present                     |
| in  | <i>integer(i4), dimension(:) :: nHorizons</i>       | Number of Horizons per soiltype                           |
| in  | <i>integer(i4), dimension(:) :: nTillHorizons</i>   | Number of Tillage Horizons                                |
| in  | <i>real(dp), dimension(:, :) :: sand</i>            | sand content  |
| in  | <i>real(dp), dimension(:, :) :: clay</i>            | clay content  |
| in  | <i>real(dp), dimension(:, :) :: DbM</i>             | mineral Bulk density                                      |
| in  | <i>integer(i4), dimension(:) :: ID0</i>             | cell ids at level 0                                       |
| in  | <i>integer(i4), dimension(:, :) :: soilId0</i>      | soil ids at level 0                                       |
| in  | <i>integer(i4), dimension(:) :: LCOVER0</i>         | land cover ids at level 0                                 |
| out | <i>real(dp), dimension(:, :, :) :: thetaS_till</i>  | saturated soil moisture tillage layer                     |
| out | <i>real(dp), dimension(:, :, :) :: thetaFC_till</i> | field capacity tillage layer                              |
| out | <i>real(dp), dimension(:, :, :) :: thetaPW_till</i> | permanent wilting point tillage layer                     |
| out | <i>real(dp), dimension(:, :) :: thetaS</i>          | saturated soil moisture                                   |
| out | <i>real(dp), dimension(:, :) :: thetaFC</i>         | field capacity  |
| out | <i>real(dp), dimension(:, :) :: thetaPW</i>         | permanent wilting point                                   |
| out | <i>real(dp), dimension(:, :, :) :: Ks</i>           | saturated hydraulic conductivity                          |
| out | <i>real(dp), dimension(:, :, :) :: Db</i>           | Bulk density  |
| out | <i>real(dp), dimension(:) :: KsVar_H0</i>           | rel. var. of Ks for horizontal flow                       |
| out | <i>real(dp), dimension(:) :: KsVar_V0</i>           | rel. var. of Ks for vertical flow                         |
| out | <i>real(dp), dimension(:) :: SMS_FC0</i>            | soil moisture deficit from field cap. w.r.t to saturation |

#### Authors

Stephan Thober, Rohini Kumar

#### Date

Dec 2012

here = ncells0

in this case the second dimension of soilId0 = 1

non-till

till layers

HORIZON

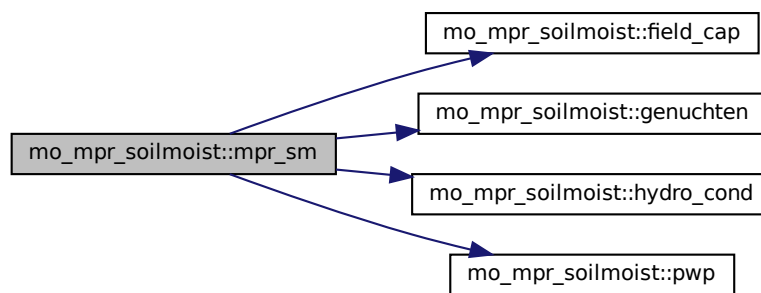
SOIL TYPE

Definition at line 100 of file mo\_mpr\_soilmoist.f90.

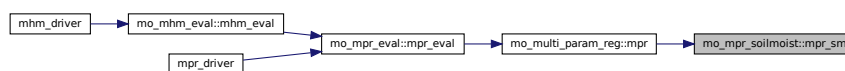
References mo\_mpr\_constants::bulkdens\_orgmatter, field\_cap(), genuchten(), hydro\_cond(), mo\_mpr\_global\_variables::iflag\_soildb, mo\_common\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, and pwp().

Referenced by mo\_multi\_param\_reg::mpr().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.35.2.5 pwp()

```

elemental pure subroutine mo_mpr_soilmoist::pwp (
    real(dp), intent(in) Genu_Mual_n,
    real(dp), intent(in) Genu_Mual_alpha,
    real(dp), intent(in) thetaS,
    real(dp), intent(out) thetaPWP )
  
```

Permanent Wilting point.

This subroutine calculates the permanent wilting point according to Zacharias et al. (2007, Soil Phy.) and using van Genuchten 1980's equation. For the water retention curve at a matrix potential of -1500 kPa, it is assumed that  $\theta_{R} = 0$ . ADDITIONAL INFORMATION Zacharias et al. 2007, Soil Phy.

#### Parameters

|    |                         |                             |
|----|-------------------------|-----------------------------|
| in | real(dp) :: Genu_Mual_n | - Genuchten shape parameter |
|----|-------------------------|-----------------------------|



## Parameters

|     |                                    |                             |
|-----|------------------------------------|-----------------------------|
| in  | <i>real(dp) :: Genu_Mual_alpha</i> | - Genuchten shape parameter |
| in  | <i>real(dp) :: thetaS</i>          | - saturated water content   |
| out | <i>real(dp) :: thetaPWP</i>        | - Permanent Wilting point   |

## Authors

Stephan Thober, Rohini Kumar

## Date

Dec, 2012

Definition at line 493 of file mo\_mpr\_soilmoist.f90.

References mo\_mpr\_constants::pwp\_c, and mo\_mpr\_constants::pwp\_matpot\_thetar.

Referenced by mpr\_sm().

Here is the caller graph for this function:



## 17.36 mo\_mpr\_startup Module Reference

Startup procedures for mHM.

### Functions/Subroutines

- subroutine, public [mpr\\_initialize](#)  
*Initialize main mHM variables.*
- subroutine [l0\\_check\\_input](#) (iDomain)  
*Check for errors in L0 input data.*
- subroutine [l0\\_variable\\_init](#) (iDomain)  
*level 0 variable initialization*
- subroutine, public [init\\_eff\\_params](#) (ncells1)  
*Allocation of space for mHM related L1 and L11 variables.*

### 17.36.1 Detailed Description

Startup procedures for mHM.

This module initializes all variables required to run mHM. This module needs to be run only one time at the beginning of a simulation if re-starting files do not exist.

## Authors

Luis Samaniego, Rohini Kumar

## Date

Dec 2012

## 17.36.2 Function/Subroutine Documentation

### 17.36.2.1 `init_eff_params()`

```
subroutine, public mo_mpr_startup::init_eff_params (
    integer(i4), intent(in) ncells1 )
```

Allocation of space for mHM related L1 and L11 variables.

Allocation of space for mHM related L1 and L11 variables (e.g., states, fluxes, and parameters) for a given domain. Variables allocated here is defined in them [mo\\_global\\_variables.f90](#) file. After allocating any variable in this routine, initialize them in the following `variables_default_init` subroutine:

#### Parameters

|                 |                                     |
|-----------------|-------------------------------------|
| <code>in</code> | <code>integer(i4) :: ncells1</code> |
|-----------------|-------------------------------------|

#### Authors

Rohini Kumar

#### Date

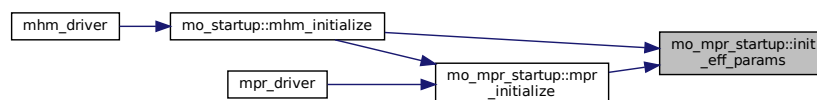
Jan 2013

Definition at line 401 of file `mo_mpr_startup.f90`.

References `mo_mpr_global_variables::l1_aeroresist`, `mo_mpr_global_variables::l1_alpha`, `mo_mpr_global_variables::l1_degday`, `mo_mpr_global_variables::l1_degdayinc`, `mo_mpr_global_variables::l1_degdaymax`, `mo_mpr_global_variables::l1_degdaynopre`, `mo_mpr_global_variables::l1_fasp`, `mo_mpr_global_variables::l1_froots`, `mo_mpr_global_variables::l1_fsealed`, `mo_mpr_global_variables::l1_harsamcoeff`, `mo_mpr_global_variables::l1_jarvis_thresh_c1`, `mo_mpr_global_variables::l1_karstloss`, `mo_mpr_global_variables::l1_kbaseflow`, `mo_mpr_global_variables::l1_kfastflow`, `mo_mpr_global_variables::l1_kperco`, `mo_mpr_global_variables::l1_kslowflow`, `mo_mpr_global_variables::l1_maxinter`, `mo_mpr_global_variables::l1_petlaicorfactor`, `mo_mpr_global_variables::l1_prietayalpha`, `mo_mpr_global_variables::l1_sealedthresh`, `mo_mpr_global_variables::l1_soilmoistexp`, `mo_mpr_global_variables::l1_soilmoistfc`, `mo_mpr_global_variables::l1_soilmoistsat`, `mo_mpr_global_variables::l1_surfresist`, `mo_mpr_global_variables::l1_tempthresh`, `mo_mpr_global_variables::l1_unsatthresh`, `mo_mpr_global_variables::l1_wiltingpoint`, `mo_mpr_global_variables::nlai`, `mo_common_variables::nlcoverscene`, `mo_mpr_global_variables::nsoilhorizons_mhm`, and `mo_common_constants::p1_initstatefluxes`.

Referenced by `mo_startup::mhm_initialize()`, and `mpr_initialize()`.

Here is the caller graph for this function:



### 17.36.2.2 I0\_check\_input()

```
subroutine mo_mpr_startup::l0_check_input (
    integer(i4), intent(in) iDomain )
```

Check for errors in L0 input data.

Check for possible errors in input data (morphological and land cover) at level-0

#### Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | domain id |
|----|-------------------------------|-----------|

#### Authors

Rohini Kumar

#### Date

Jan 2013

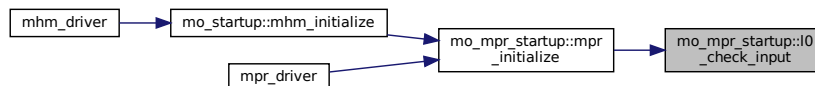
by default; when iFlag\_soilDB = 0

Definition at line 143 of file mo\_mpr\_startup.f90.

References mo\_common\_constants::eps\_dp, mo\_mpr\_global\_variables::iflag\_soildb, mo\_mpr\_global\_variables::l0\_esp, mo\_common\_variables::l0\_elev, mo\_mpr\_global\_variables::l0\_geounit, mo\_mpr\_global\_variables::l0\_gridded\_lai, mo\_common\_variables::l0\_lcover, mo\_mpr\_global\_variables::l0\_slope, mo\_mpr\_global\_variables::l0\_soilid, mo\_common\_variables::level0, mo\_common\_variables::nlcoverscene, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, and mo\_mpr\_global\_variables::timestep\_lai\_input.

Referenced by mpr\_initialize().

Here is the caller graph for this function:



### 17.36.2.3 I0\_variable\_init()

```
subroutine mo_mpr_startup::l0_variable_init (
    integer(i4), intent(in) iDomain )
```

level 0 variable initialization

following tasks are performed for L0 data sets

- cell id & numbering
- storage of cell coordinates (row and column id)
- empirical dist. of terrain slope
- flag to determine the presence of a particular soil id in this configuration of the model run If a variable is added or removed here, then it also has to be added or removed in the subroutine config\_variables\_set in module [mo\\_restart](#) and in the subroutine set\_config in module mo\_set\_netcdf\_restart

## Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | domain id |
|----|-------------------------------|-----------|

## Authors

Rohini Kumar

## Date

Jan 2013

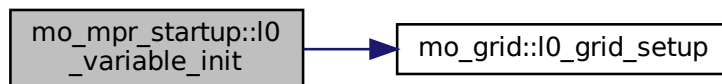
by default; when `iFlag_soilDB = 0`

Definition at line 270 of file `mo_mpr_startup.f90`.

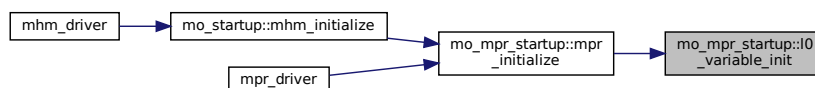
References `mo_mpr_global_variables::iflag_soildb`, `mo_grid::l0_grid_setup()`, `mo_mpr_global_variables::l0_slope`, `mo_mpr_global_variables::l0_slope_emp`, `mo_mpr_global_variables::l0_soilid`, `mo_common_variables::level0`, `mo_mpr_global_variables::nsoilhorizons_mhm`, `mo_mpr_global_variables::nsoiltypes`, and `mo_mpr_global_variables::soildb`.

Referenced by `mpr_initialize()`.

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.36.2.4 mpr\_initialize()

subroutine, public `mo_mpr_startup::mpr_initialize`

Initialize main mHM variables.

Initialize main mHM variables for a given domain. Calls the following procedures in this order:

- Constant initialization.
- Generate soil database.
- Checking inconsistencies input fields.

- Variable initialization at level-0.
- Variable initialization at level-1.
- Variable initialization at level-11.
- Space allocation of remaining variable/parameters. Global variables will be used at this stage.

**Authors**

Luis Samaniego, Rohini Kumar

**Date**

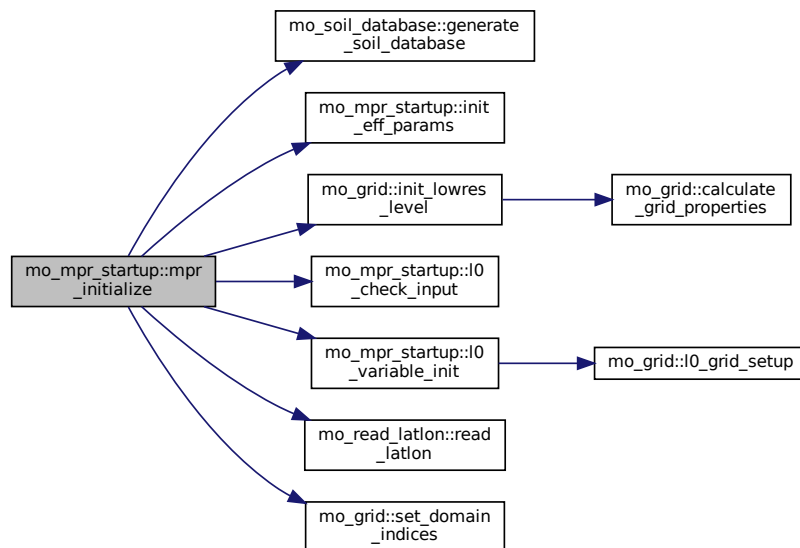
Dec 2012

Definition at line 71 of file mo\_mpr\_startup.f90.

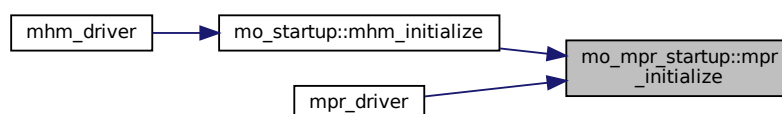
References mo\_common\_variables::domainmeta, mo\_soil\_database::generate\_soil\_database(), init\_eff\_params(), mo\_grid::init\_lowres\_level(), l0\_check\_input(), mo\_common\_variables::l0\_l1\_remap, l0\_variable\_init(), mo\_common\_variables::level0, mo\_common\_variables::level1, mo\_read\_latlon::read\_latlon(), mo\_common\_variables::resolutionhydrology, and mo\_grid::set\_domain\_indices().

Referenced by mo\_startup::mhm\_initialize(), and mpr\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.37 mo\_mrm\_constants Module Reference

Provides mRM specific constants.

### Variables

- integer(i4), parameter, public `noutfluxstate` = 2\_i4
- integer(i4), parameter, public `nroutingstates` = 2
- integer(i4), parameter, public `maxnogauges` = 50\_i4
- real(dp), parameter, public `rout_space_weight` = 0.\_dp
- real(dp), parameter, public `deltah` = 5.000\_dp
- real(dp), dimension(19), parameter `given_ts` = (/ 60.\_dp, 120.\_dp, 180.\_dp, 240.\_dp, 300.\_dp, 360.\_dp, 600.\_dp, 720.\_dp, 900.\_dp, 1200.\_dp, 1800.\_dp, 3600.\_dp, 7200.\_dp, 10800.\_dp, 14400.\_dp, 21600.\_dp, 28800.\_dp, 43200.\_dp, 86400.\_dp/)

### 17.37.1 Detailed Description

Provides mRM specific constants.

Provides mRM specific constants such as flood plain elevation.

#### Authors

Stephan Thober

#### Date

Aug 2015

### 17.37.2 Variable Documentation

#### 17.37.2.1 `deltah`

```
real(dp), parameter, public mo_mrm_constants::deltah = 5.000_dp
```

Definition at line 37 of file `mo_mrm_constants.f90`.

Referenced by `mo_mrm_net_startup::moveup()`.

#### 17.37.2.2 `given_ts`

```
real(dp), dimension(19), parameter mo_mrm_constants::given_ts = (/ 60._dp, 120._dp, 180._dp, 240._dp, 300._dp, 360._dp, 600._dp, 720._dp, 900._dp, 1200._dp, 1800._dp, 3600._dp, 7200._dp, 10800._dp, 14400._dp, 21600._dp, 28800._dp, 43200._dp, 86400._dp/)
```

Definition at line 41 of file `mo_mrm_constants.f90`.

Referenced by `mo_mrm_mpr::mrm_init_param()`, and `mo_mrm_mpr::mrm_update_param()`.

### 17.37.2.3 maxnogauges

```
integer(i4), parameter, public mo_mrm_constants::maxnogauges = 50_i4
```

Definition at line 27 of file mo\_mrm\_constants.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config().

### 17.37.2.4 noutflxstate

```
integer(i4), parameter, public mo_mrm_constants::noutflxstate = 2_i4
```

Definition at line 18 of file mo\_mrm\_constants.f90.

### 17.37.2.5 nroutingstates

```
integer(i4), parameter, public mo_mrm_constants::nroutingstates = 2
```

Definition at line 20 of file mo\_mrm\_constants.f90.

Referenced by mo\_mrm\_riv\_temp\_class::init(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), and mo\_mrm\_init::variables\_alloc\_routing().

### 17.37.2.6 rout\_space\_weight

```
real(dp), parameter, public mo_mrm_constants::rout_space_weight = 0._dp
```

Definition at line 32 of file mo\_mrm\_constants.f90.

Referenced by mo\_mrm\_mpr::mrm\_update\_param().

## 17.38 mo\_mrm\_file Module Reference

Provides file names and units for mRM.

### Variables

- character(len=\*), parameter `version` = '1.0'  
*Current mHM model version.*
- character(len=\*), parameter `version_date` = 'May 2019'  
*Time of current mHM model version release.*
- character(len=\*), parameter `file_main` = 'mrm\_driver.f90'  
*Driver file.*
- character(len=\*), parameter `file_namelist_mrm` = 'mrm.nml'  
*Namelist file name.*
- integer, parameter `unamelist_mrm` = 40  
*Unit for namelist.*
- character(len=\*), parameter `file_namelist_param_mrm` = 'mrm\_parameter.nml'  
*Parameter namelists file name.*
- integer, parameter `unamelist_param_mrm` = 41  
*Unit for namelist.*

- character(len=\*), parameter `file_facc` = 'facc.asc'
- integer, parameter `ufacc` = 56  
*Unit for flow accumulation input data file.*
- character(len=\*), parameter `file_fdir` = 'fdir.asc'  
*flow direction input data file*
- integer, parameter `ufdir` = 57  
*Unit for flow direction input data file.*
- character(len= \*), parameter `file_slope` = 'slope.asc'  
*flow direction input data file*
- integer, parameter `uslope` = 59  
*Unit for flow direction input data file.*
- character(len=\*), parameter `file_gaugeloc` = 'idgauges.asc'  
*gauge location input data file*
- integer, parameter `ugaugeloc` = 62  
*Unit for gauge location input data file.*
- integer, parameter `udischarge` = 66  
*unit for discharge time series*
- character(len=\*), parameter `file_defoutput` = 'mrm\_outputs.nml'  
*file defining mRM's outputs*
- integer, parameter `udefoutput` = 67  
*Unit for file defining mRM's outputs.*
- character(len=\*), parameter `file_config` = 'ConfigFile.log'  
*file defining mHM's outputs*
- integer, parameter `uconfig` = 68  
*Unit for file defining mHM's outputs.*
- character(len=\*), parameter `file_daily_discharge` = 'daily\_discharge.out'  
*file defining optimazation outputs*
- integer, parameter `udaily_discharge` = 74  
*Unit for file optimazation outputs.*
- character(len=\*), parameter `ncfile_discharge` = 'discharge.nc'  
*file defining optimazation outputs*
- character(len=\*), parameter `file_mrm_output` = 'mRM\_Fluxes\_States.nc'  
*file containing mrm output*
- character(len=\*), parameter `file_gw_output` = 'mRM\_gw\_Fluxes\_States.nc'  
*file containing mrm output for groundwater coupling*

### 17.38.1 Detailed Description

Provides file names and units for mRM.

Provides all filenames as well as all units used for the multiscale Routing Model mRM.

#### Authors

Matthias Cuntz, Stephan Thober

#### Date

Aug 2015

### 17.38.2 Variable Documentation



### 17.38.2.1 file\_config

```
character(len = *), parameter mo_mrm_file::file_config = 'ConfigFile.log'
```

file defining mHM's outputs

Definition at line 58 of file mo\_mrm\_file.f90.

### 17.38.2.2 file\_daily\_discharge

```
character(len = *), parameter mo_mrm_file::file_daily_discharge = 'daily_discharge.out'
```

file defining optimization outputs

Definition at line 63 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_write::write\_daily\_obs\_sim\_discharge().

### 17.38.2.3 file\_defoutput

```
character(len = *), parameter mo_mrm_file::file_defoutput = 'mrm_outputs.nml'
```

file defining mRM's outputs

Definition at line 53 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_init::config\_output(), mo\_mrm\_read\_config::mrm\_read\_config(), and mo\_mrm\_init::print\_startup\_message().

### 17.38.2.4 file\_facc

```
character(len = *), parameter mo_mrm_file::file_facc = 'facc.asc'
```

Definition at line 32 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_I0\_data().

### 17.38.2.5 file\_fdir

```
character(len = *), parameter mo_mrm_file::file_fdir = 'fdir.asc'
```

flow direction input data file

Definition at line 36 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_I0\_data().

### 17.38.2.6 file\_gaugeloc

```
character(len = *), parameter mo_mrm_file::file_gaugeloc = 'idgauges.asc'
```

gauge location input data file

Definition at line 45 of file mo\_mrm\_file.f90.

Referenced by `mo_mrm_read_data::mrm_read_l0_data()`.

### 17.38.2.7 file\_gw\_output

```
character(len = *), parameter mo_mrm_file::file_gw_output = 'mRM_gw_Fluxes_States.nc'
```

file containing mrm output for groundwater coupling

Definition at line 73 of file `mo_mrm_file.f90`.

### 17.38.2.8 file\_main

```
character(len = *), parameter mo_mrm_file::file_main = 'mrm_driver.f90'
```

Driver file.

Definition at line 22 of file `mo_mrm_file.f90`.

Referenced by `mo_mrm_init::print_startup_message()`.

### 17.38.2.9 file\_mrm\_output

```
character(len = *), parameter mo_mrm_file::file_mrm_output = 'mRM_Fluxes_States.nc'
```

file containing mrm output

Definition at line 70 of file `mo_mrm_file.f90`.

Referenced by `mo_mrm_write_fluxes_states::createoutputfile()`.

### 17.38.2.10 file\_namelist\_mrm

```
character(len = *), parameter mo_mrm_file::file_namelist_mrm = 'mrm.nml'
```

Namelist file name.

Definition at line 24 of file `mo_mrm_file.f90`.

Referenced by `mo_mrm_init::config_output()`.

### 17.38.2.11 file\_namelist\_param\_mrm

```
character(len = *), parameter mo_mrm_file::file_namelist_param_mrm = 'mrm_parameter.nml'
```

Parameter namelists file name.

Definition at line 28 of file `mo_mrm_file.f90`.

Referenced by `mo_mrm_init::config_output()`.

**17.38.2.12 file\_slope**

```
character(len=*), parameter mo_mrm_file::file_slope = 'slope.asc'
```

flow direction input data file

Definition at line 40 of file mo\_mrm\_file.f90.

**17.38.2.13 ncfile\_discharge**

```
character(len = *), parameter mo_mrm_file::ncfile_discharge = 'discharge.nc'
```

file defining optimization outputs

Definition at line 67 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_write::write\_daily\_obs\_sim\_discharge().

**17.38.2.14 uconfig**

```
integer, parameter mo_mrm_file::uconfig = 68
```

Unit for file defining mHM's outputs.

Definition at line 60 of file mo\_mrm\_file.f90.

**17.38.2.15 udaily\_discharge**

```
integer, parameter mo_mrm_file::udaily_discharge = 74
```

Unit for file optimization outputs.

Definition at line 65 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_write::write\_daily\_obs\_sim\_discharge().

**17.38.2.16 undefoutput**

```
integer, parameter mo_mrm_file::undefoutput = 67
```

Unit for file defining mRM's outputs.

Definition at line 55 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config().

**17.38.2.17 udischarge**

```
integer, parameter mo_mrm_file::udischarge = 66
```

unit for discharge time series

Definition at line 50 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_discharge().

**17.38.2.18 ufact**

```
integer, parameter mo_mrm_file::ufact = 56
```

Unit for flow accumulation input data file.

Definition at line 34 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_I0\_data().

**17.38.2.19 ufdir**

```
integer, parameter mo_mrm_file::ufdir = 57
```

Unit for flow direction input data file.

Definition at line 38 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_I0\_data().

**17.38.2.20 ugaugeloc**

```
integer, parameter mo_mrm_file::ugaugeloc = 62
```

Unit for gauge location input data file.

Definition at line 47 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_I0\_data().

**17.38.2.21 unamelist\_mrm**

```
integer, parameter mo_mrm_file::unamelist_mrm = 40
```

Unit for namelist.

Definition at line 26 of file mo\_mrm\_file.f90.

**17.38.2.22 unamelist\_param\_mrm**

```
integer, parameter mo_mrm_file::unamelist_param_mrm = 41
```

Unit for namelist.

Definition at line 30 of file mo\_mrm\_file.f90.

**17.38.2.23 uslope**

```
integer, parameter mo_mrm_file::uslope = 59
```

Unit for flow direction input data file.

Definition at line 42 of file mo\_mrm\_file.f90.

#### 17.38.2.24 version

```
character(len = *), parameter mo_mrm_file::version = '1.0'
```

Current mHM model version.

Definition at line 18 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_write\_fluxes\_states::createoutputfile(), mo\_mrm\_init::print\_startup\_message(), and mo\_mrm\_write::write\_configfile().

#### 17.38.2.25 version\_date

```
character(len = *), parameter mo_mrm_file::version_date = 'May 2019'
```

Time of current mHM model version release.

Definition at line 20 of file mo\_mrm\_file.f90.

Referenced by mo\_mrm\_init::print\_startup\_message().

## 17.39 mo\_mrm\_global\_variables Module Reference

Global variables for mRM only.

### Data Types

- type [gaugingstation](#)
- type [domaininfo\\_mrm](#)

### Variables

- logical [is\\_start](#)
- integer(i4) [output\\_deflate\\_level\\_mrm](#)
- logical [output\\_double\\_precision\\_mrm](#)
- integer(i4) [timestep\\_model\\_outputs\\_mrm](#)
- logical, dimension(noutfluxstate) [outputfluxstate\\_mrm](#)
- character(256), dimension(:), allocatable, public [dirgauges](#)
- character(256), dimension(:), allocatable, public [dirtotalrunoff](#)
- character(256), public [filenametotalrunoff](#)
- character(256), public [varnametotalrunoff](#)
- character(256), dimension(:), allocatable, public [dirbankfullrunoff](#)
- integer(i4), public [ntstepday](#)
- type([grid](#)), dimension(:), allocatable, target, public [level11](#)
- type([gridremapper](#)), dimension(:), allocatable, public [l0\\_l11\\_remap](#)
- type([gridremapper](#)), dimension(:), allocatable, public [l1\\_l11\\_remap](#)
- real(dp), dimension(:, :), allocatable, public [mrm\\_runoff](#)
- integer(i4), public [ngaugestotal](#)
- integer(i4), public [ngaugeslocal](#)
- integer(i4), public [ninflowgaugestotal](#)

- integer(i4), public [nmeasperday](#)
- type([gaugingstation](#)), public [gauge](#)
- type([gaugingstation](#)), public [inflowgauge](#)
- type([domaininfo\\_mrm](#)), dimension(:), allocatable, target, public [domain\\_mrm](#)
- integer(i4), dimension(:), allocatable, public [l0\\_gaugeloc](#)
- integer(i4), dimension(:), allocatable, public [l0\\_inflowgaugeloc](#)
- integer(i4), dimension(:), allocatable, public [l0\\_facc](#)
- integer(i4), dimension(:), allocatable, public [l0\\_fdir](#)
- integer(i4), dimension(:), allocatable, public [l0\\_drasc](#)
- integer(i4), dimension(:), allocatable, public [l0\\_dracell](#)
- integer(i4), dimension(:), allocatable, public [l0\\_streamnet](#)
- integer(i4), dimension(:), allocatable, public [l0\\_floodplain](#)
- integer(i4), dimension(:), allocatable, public [l0\\_noutlet](#)
- real(dp), dimension(:), allocatable, public [l0\\_celerity](#)
- integer(i4), dimension(:), allocatable, public [l11\\_l1\\_id](#)
- real(dp), dimension(:, :), allocatable, public [l1\\_total\\_runoff\\_in](#)
- integer(i4), dimension(:, :), allocatable, public [l11\\_cellcoor](#)
- integer(i4), dimension(:), allocatable, public [l1\\_l11\\_id](#)
- real(dp), dimension(:), allocatable, public [l11\\_areacell](#)
- real(dp), dimension(:), allocatable, public [l11\\_facc](#)
- integer(i4), dimension(:), allocatable, public [l11\\_fdir](#)
- integer(i4), dimension(:), allocatable, public [l11\\_noutlets](#)
- real(dp), dimension(:), allocatable, public [l11\\_celerity](#)
- real(dp), dimension(:), allocatable, public [l11\\_meandering](#)
- real(dp), dimension(:), allocatable, public [l11\\_linkin\\_facc](#)
- integer(i4), dimension(:), allocatable, public [l11\\_rowout](#)
- integer(i4), dimension(:), allocatable, public [l11\\_colout](#)
- real(dp), dimension(:), allocatable, public [l11\\_qmod](#)
- real(dp), dimension(:), allocatable, public [l11\\_qout](#)
- real(dp), dimension(:, :), allocatable, public [l11\\_qtin](#)
- real(dp), dimension(:, :), allocatable, public [l11\\_qtr](#)
- integer(i4), dimension(:), allocatable, public [l11\\_fromn](#)
- integer(i4), dimension(:), allocatable, public [l11\\_ton](#)
- integer(i4), dimension(:), allocatable, public [l11\\_netperm](#)
- integer(i4), dimension(:), allocatable, public [l11\\_frow](#)
- integer(i4), dimension(:), allocatable, public [l11\\_fcol](#)
- integer(i4), dimension(:), allocatable, public [l11\\_trow](#)
- integer(i4), dimension(:), allocatable, public [l11\\_tcol](#)
- integer(i4), dimension(:), allocatable, public [l11\\_rorder](#)
- integer(i4), dimension(:), allocatable, public [l11\\_label](#)
- logical, dimension(:), allocatable, public [l11\\_sink](#)
- real(dp), dimension(:), allocatable, public [l11\\_length](#)
- real(dp), dimension(:), allocatable, target, public [l11\\_afloodplain](#)
- real(dp), dimension(:), allocatable, public [l11\\_slope](#)
- real(dp), dimension(:, :), allocatable, public [l11\\_nlinkfracfpimp](#)
- real(dp), dimension(:), allocatable, public [l11\\_k](#)
- real(dp), dimension(:), allocatable, public [l11\\_xi](#)
- real(dp), dimension(:), allocatable, public [l11\\_tsrout](#)
- real(dp), dimension(:), allocatable, public [l11\\_c1](#)
- real(dp), dimension(:), allocatable, public [l11\\_c2](#)
- logical [gw\\_coupling](#)
- real(dp), dimension(:), allocatable, public [l11\\_bankfull\\_runoff\\_in](#)
- real(dp), dimension(:), allocatable, public [l0\\_channel\\_depth](#)
- real(dp), dimension(:), allocatable, public [l0\\_channel\\_elevation](#)
- real(dp), dimension(:), allocatable, public [l0\\_river\\_head\\_mon\\_sum](#)
- real(dp), dimension(:), allocatable, public [l0\\_slope](#)
- type([riv\\_temp\\_type](#)), public [riv\\_temp\\_pcs](#)

*This is a container for the river temperature routing process (pcs)*

## 17.39.1 Detailed Description

Global variables for mRM only.

TODO: add description

### Authors

Luis Samaniego, Stephan Thober

### Date

Aug 2015

## 17.39.2 Variable Documentation

### 17.39.2.1 dirbankfullrunoff

```
character(256), dimension(:), allocatable, public mo_mrm_global_variables::dirbankfullrunoff
```

Definition at line 50 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_bankfull\_runoff(), and mo\_mrm\_read\_config::mrm\_read\_config().

### 17.39.2.2 dirgauges

```
character(256), dimension(:), allocatable, public mo_mrm_global_variables::dirgauges
```

Definition at line 46 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_init::config\_output(), mo\_mrm\_read\_config::mrm\_read\_config(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.39.2.3 dirtotalrunoff

```
character(256), dimension(:), allocatable, public mo_mrm_global_variables::dirtotalrunoff
```

Definition at line 47 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_total\_runoff(), and mo\_mrm\_write::write\_configfile().

### 17.39.2.4 domain\_mrm

```
type(domaininfo_mrm), dimension(:), allocatable, target, public mo_mrm_global_variables::domain_mrm
```

Definition at line 116 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_init::config\_output(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_I0\_data(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_write::mrm\_write(),

`mo_mrm_restart::mrm_write_restart()`, `mo_mrm_write::write_configfile()`, and `mo_mrm_write::write_daily_obs_↔  
sim_discharge()`.

#### 17.39.2.5 `filenametotalrunoff`

```
character(256), public mo_mrm_global_variables::filenametotalrunoff
```

Definition at line 48 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_read_config::mrm_read_config()`, and `mo_mrm_read_data::mrm_read_total_runoff()`.

#### 17.39.2.6 `gauge`

```
type(gaugingstation), public mo_mrm_global_variables::gauge
```

Definition at line 87 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_objective_function_runoff::extract_runoff()`, `mo_mrm_read_config::mrm_read_config()`, `mo_mrm_read_data::mrm_read_discharge()`, `mo_mrm_write::mrm_write()`, `mo_mrm_objective_function_runoff_↔  
::multi_objective_ae_fdc_lsv_nse_djf()`, `mo_write_ascii::write_configfile()`, `mo_mrm_write::write_configfile()`, and `mo_mrm_write::write_daily_obs_sim_discharge()`.

#### 17.39.2.7 `gw_coupling`

```
logical mo_mrm_global_variables::gw_coupling
```

Definition at line 215 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_mrm_init::mrm_init()`, and `mo_mrm_read_config::mrm_read_↔  
config()`.

#### 17.39.2.8 `inflowgauge`

```
type(gaugingstation), public mo_mrm_global_variables::inflowgauge
```

Definition at line 88 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mhm_eval::mhm_eval()`, `mo_mrm_read_config::mrm_read_config()`, `mo_mrm_read_data_↔  
::mrm_read_discharge()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

#### 17.39.2.9 `is_start`

```
logical mo_mrm_global_variables::is_start
```

Definition at line 30 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_read_config::mrm_read_config()`, and `mo_mrm_routing::mrm_routing()`.



### 17.39.2.10 IO\_celerity

`real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l0_celerity`

Definition at line 136 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_init::variables_alloc_routing()`.

### 17.39.2.11 IO\_channel\_depth

`real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l0_channel_depth`

Definition at line 219 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_river_head::calc_channel_elevation()`.

### 17.39.2.12 IO\_channel\_elevation

`real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l0_channel_elevation`

Definition at line 220 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_river_head::calc_river_head()`.

### 17.39.2.13 IO\_dracell

`integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_dracell`

Definition at line 132 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_net_startup::l11_set_drain_outlet_gauges()`.

### 17.39.2.14 IO\_drasc

`integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_drasc`

Definition at line 130 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_net_startup::l11_flow_direction()`, `mo_mrm_net_startup::l11_link_location()`, and `mo_mrm_net_startup::l11_set_drain_outlet_gauges()`.

### 17.39.2.15 IO\_facc

`integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_facc`

Definition at line 125 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_river_head::calc_channel_elevation()`, `mo_mrm_init::l0_check_input_routing()`, `mo_mrm_net_startup::l11_flow_direction()`, `mo_mrm_read_data::mrm_read_IO_data()`, `mo_mrm_restart::mrm_read_restart_config()`, and `mo_mrm_restart::mrm_write_restart()`.

### 17.39.2.16 IO\_fdir

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_fdir
```

Definition at line 126 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_river\_head::calc\_channel\_elevation(), mo\_mrm\_init::l0\_check\_input\_routing(), mo\_mrm\_net\_startup::l11\_calc\_celerity(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_read\_data::mrm\_read\_l0\_data(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.17 IO\_floodplain

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_floodplain
```

Definition at line 134 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), and mo\_mrm\_net\_startup::l11\_stream\_features().

### 17.39.2.18 IO\_gaugeloc

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_gaugeloc
```

Definition at line 123 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), and mo\_mrm\_read\_data::mrm\_read\_l0\_data().

### 17.39.2.19 IO\_inflowgaugeloc

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_inflowgaugeloc
```

Definition at line 124 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), and mo\_mrm\_read\_data::mrm\_read\_l0\_data().

### 17.39.2.20 IO\_l11\_remap

```
type(gridremapper), dimension(:), allocatable, public mo_mrm_global_variables::l0_l11_remap
```

Definition at line 62 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_river\_head::calc\_river\_head(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_set\_drain\_outlet\_gauges(), and mo\_mrm\_init::mrm\_init().

### 17.39.2.21 IO\_noutlet

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_noutlet
```

Definition at line 135 of file mo\_mrm\_global\_variables.f90.

**17.39.2.22 l0\_river\_head\_mon\_sum**

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l0_river_head_mon_sum
```

Definition at line 222 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mrm\_init::mrm\_init().

**17.39.2.23 l0\_slope**

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l0_slope
```

Definition at line 223 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_river\_head::calc\_river\_head().

**17.39.2.24 l0\_streamnet**

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l0_streamnet
```

Definition at line 133 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

**17.39.2.25 l11\_afloodplain**

```
real(dp), dimension(:), allocatable, target, public mo_mrm_global_variables::l11_afloodplain
```

Definition at line 195 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

**17.39.2.26 l11\_areacell**

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_areacell
```

Definition at line 158 of file mo\_mrm\_global\_variables.f90.

**17.39.2.27 l11\_bankfull\_runoff\_in**

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_bankfull_runoff_in
```

Definition at line 217 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_river\_head::calc\_river\_head().

### 17.39.2.28 I11\_c1

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_c1
```

Definition at line 208 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_mpr::mrm\_↔ update\_param(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_↔ init::variables\_default\_init\_routing().

### 17.39.2.29 I11\_c2

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_c2
```

Definition at line 209 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_mpr::mrm\_↔ update\_param(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_↔ init::variables\_default\_init\_routing().

### 17.39.2.30 I11\_celerity

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_celerity
```

Definition at line 162 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_mpr::mrm\_update\_param(), mo\_mrm\_restart::mrm\_↔ write\_restart(), and mo\_mrm\_init::variables\_alloc\_routing().

### 17.39.2.31 I11\_cellcoor

```
integer(i4), dimension(:, :), allocatable, public mo_mrm_global_variables::l11_cellcoor
```

Definition at line 155 of file mo\_mrm\_global\_variables.f90.

### 17.39.2.32 I11\_colout

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_colout
```

Definition at line 171 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_↔ restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.33 I11\_facc

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_facc
```

Definition at line 159 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

**17.39.2.34 l11\_fcol**

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_fcol
```

Definition at line 188 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

**17.39.2.35 l11\_fdir**

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_fdir
```

Definition at line 160 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_flow\_accumulation(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_net\_startup::l11\_set\_network\_topology(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

**17.39.2.36 l11\_fromn**

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_fromn
```

Definition at line 184 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_net\_startup::l11\_set\_network\_topology(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.39.2.37 l11\_frow**

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_frow
```

Definition at line 187 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

**17.39.2.38 l11\_k**

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_k
```

Definition at line 204 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_init::variables\_default\_init\_routing().

**17.39.2.39 l11\_l1\_id**

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_l1_id
```

Definition at line 142 of file mo\_mrm\_global\_variables.f90.

Referenced by `mo_mrm_net_startup::l11_l1_mapping()`, `mo_mhm_eval::mhm_eval()`, `mo_mrm_restart::mrm_read_restart_config()`, `mo_mrm_restart::mrm_write_restart()`, and `mo_mrm_write::write_configfile()`.

#### 17.39.2.40 l11\_label

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_label
```

Definition at line 192 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_net_startup::l11_routing_order()`, `mo_mrm_restart::mrm_read_restart_config()`, `mo_mrm_restart::mrm_write_restart()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

#### 17.39.2.41 l11\_length

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_length
```

Definition at line 194 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_net_startup::l11_stream_features()`, `mo_mhm_eval::mhm_eval()`, `mo_mrm_init::mrm_init()`, `mo_mrm_restart::mrm_read_restart_config()`, `mo_mrm_mpr::mrm_update_param()`, `mo_mrm_restart::mrm_write_restart()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

#### 17.39.2.42 l11\_linkin\_facc

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_linkin_facc
```

Definition at line 166 of file `mo_mrm_global_variables.f90`.

#### 17.39.2.43 l11\_meandering

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_meandering
```

Definition at line 164 of file `mo_mrm_global_variables.f90`.

#### 17.39.2.44 l11\_netperm

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_netperm
```

Definition at line 186 of file `mo_mrm_global_variables.f90`.

Referenced by `mo_mrm_net_startup::l11_link_location()`, `mo_mrm_net_startup::l11_routing_order()`, `mo_mrm_net_startup::l11_stream_features()`, `mo_mhm_eval::mhm_eval()`, `mo_mrm_init::mrm_init()`, `mo_mrm_restart::mrm_read_restart_config()`, `mo_mrm_restart::mrm_write_restart()`, `mo_write_ascii::write_configfile()`, and `mo_mrm_write::write_configfile()`.

#### 17.39.2.45 l11\_nlinkfracpimp

```
real(dp), dimension(:, :), allocatable, public mo_mrm_global_variables::l11_nlinkfracpimp
```

Definition at line 201 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), and mo\_mrm\_restart::mrm\_write\_restart().

#### 17.39.2.46 l11\_noutlets

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_noutlets
```

Definition at line 161 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_mpr::mrm\_update\_param(), and mo\_write\_ascii::write\_configfile().

#### 17.39.2.47 l11\_qmod

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_qmod
```

Definition at line 179 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_init::variables\_default\_init\_routing().

#### 17.39.2.48 l11\_qout

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_qout
```

Definition at line 180 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_init::variables\_default\_init\_routing().

#### 17.39.2.49 l11\_qtin

```
real(dp), dimension(:, :), allocatable, public mo_mrm_global_variables::l11_qtin
```

Definition at line 181 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_init::variables\_default\_init\_routing().

#### 17.39.2.50 l11\_qtr

```
real(dp), dimension(:, :), allocatable, public mo_mrm_global_variables::l11_qtr
```

Definition at line 182 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_init::variables\_default\_init\_routing().

### 17.39.2.51 l11\_rorder

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_rorder
```

Definition at line 191 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.39.2.52 l11\_rowout

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_rowout
```

Definition at line 170 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.53 l11\_sink

```
logical, dimension(:), allocatable, public mo_mrm_global_variables::l11_sink
```

Definition at line 193 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.54 l11\_slope

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_slope
```

Definition at line 197 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.39.2.55 l11\_tcol

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_tcol
```

Definition at line 190 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.56 l11\_ton

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_ton
```

Definition at line 185 of file mo\_mrm\_global\_variables.f90.



Referenced by mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_net\_startup::l11\_set\_network\_topology(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.39.2.57 l11\_trow

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l11_trow
```

Definition at line 189 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mrm\_restart::mrm\_read\_restart\_config(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.58 l11\_tsrouit

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_tsrouit
```

Definition at line 207 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_mpr::mrm\_update\_param(), and mo\_mrm\_restart::mrm\_write\_restart().

### 17.39.2.59 l11\_xi

```
real(dp), dimension(:), allocatable, public mo_mrm_global_variables::l11_xi
```

Definition at line 205 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), and mo\_mrm\_init::variables\_default\_init\_routing().

### 17.39.2.60 l1\_l11\_id

```
integer(i4), dimension(:), allocatable, public mo_mrm_global_variables::l1_l11_id
```

Definition at line 157 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_net\_startup::l11\_l1\_mapping(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

### 17.39.2.61 l1\_l11\_remap

```
type(gridremapper), dimension(:), allocatable, public mo_mrm_global_variables::l1_l11_remap
```

Definition at line 63 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_init::mrm\_init().

**17.39.2.62 l1\_total\_runoff\_in**

```
real(dp), dimension(:, :), allocatable, public mo_mrm_global_variables::l1_total_runoff_in
```

Definition at line 148 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_total\_runoff().

**17.39.2.63 level11**

```
type(grid), dimension(:), allocatable, target, public mo_mrm_global_variables::level11
```

Definition at line 61 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_write\_fluxes\_states::createoutputfile(), mo\_mrm\_net\_startup::l11\_flow\_accumulation(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_floodplain(), mo\_mrm\_net\_startup::l11\_l1\_mapping(), mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_net\_startup::l11\_set\_network\_topology(), mo\_mrm\_net\_startup::l11\_stream\_features(), mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_mpr::mrm\_init\_param(), mo\_mrm\_read\_data::mrm\_read\_bankfull\_runoff(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_mrm\_restart::mrm\_read\_restart\_states(), mo\_mrm\_mpr::mrm\_update\_param(), mo\_mrm\_restart::mrm\_write\_restart(), mo\_mrm\_init::variables\_alloc\_routing(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.39.2.64 mrm\_runoff**

```
real(dp), dimension(:, :), allocatable, public mo_mrm_global_variables::mrm_runoff
```

Definition at line 69 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_read\_data::mrm\_read\_discharge(), and mo\_mrm\_write::mrm\_write().

**17.39.2.65 ngaugeslocal**

```
integer(i4), public mo_mrm_global_variables::ngaugeslocal
```

Definition at line 75 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_discharge(), and mo\_write\_ascii::write\_configfile().

**17.39.2.66 ngaugestotal**

```
integer(i4), public mo_mrm_global_variables::ngaugestotal
```

Definition at line 74 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_write::mrm\_write(), mo\_mrm\_objective\_function\_runoff::single\_objective\_runoff\_master(), mo\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.39.2.67 ninflowgaugestotal**

```
integer(i4), public mo_mrm_global_variables::ninflowgaugestotal
```

Definition at line 76 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_read\_data::mrm\_read\_discharge(), mo\_mrm\_write\_ascii::write\_configfile(), and mo\_mrm\_write::write\_configfile().

**17.39.2.68 nmeasperday**

```
integer(i4), public mo_mrm_global_variables::nmeasperday
```

Definition at line 77 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_objective\_function\_runoff::extract\_runoff(), mo\_mrm\_read\_data::mrm\_read\_discharge(), and mo\_mrm\_objective\_function\_runoff::multi\_objective\_ae\_fdc\_lsv\_nse\_djf().

**17.39.2.69 ntstepday**

```
integer(i4), public mo_mrm_global_variables::ntstepday
```

Definition at line 55 of file mo\_mrm\_global\_variables.f90.

**17.39.2.70 output\_deflate\_level\_mrm**

```
integer(i4) mo_mrm_global_variables::output_deflate_level_mrm
```

Definition at line 36 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), and mo\_mrm\_write\_fluxes\_states::newoutputvariable().

**17.39.2.71 output\_double\_precision\_mrm**

```
logical mo_mrm_global_variables::output_double_precision_mrm
```

Definition at line 37 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config().

**17.39.2.72 outputflxstate\_mrm**

```
logical, dimension(noutflxstate) mo_mrm_global_variables::outputflxstate_mrm
```

Definition at line 39 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_read\_config::mrm\_read\_config(), mo\_mrm\_write\_fluxes\_states::newoutputdataset(), and mo\_mrm\_write\_fluxes\_states::updatedataset().

### 17.39.2.73 riv\_temp\_pcs

```
type(riv_temp_type), public mo_mrm_global_variables::riv_temp_pcs
```

This is a container for the river temperature routing process (pcs)

Definition at line 229 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), mo\_mrm\_init::mrm\_configuration(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_read\_data::mrm\_read\_discharge(), mo\_mrm\_routing::mrm\_routing(), mo\_mrm\_write::mrm\_write\_output\_fluxes(), and mo\_mrm\_write\_fluxes\_states::newoutputdataset().

### 17.39.2.74 timestep\_model\_outputs\_mrm

```
integer(i4) mo_mrm_global_variables::timestep_model_outputs_mrm
```

Definition at line 38 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mhm\_eval::mhm\_eval(), and mo\_mrm\_read\_config::mrm\_read\_config().

### 17.39.2.75 varnametotalrunoff

```
character(256), public mo_mrm_global_variables::varnametotalrunoff
```

Definition at line 49 of file mo\_mrm\_global\_variables.f90.

Referenced by mo\_mrm\_read\_config::mrm\_read\_config(), and mo\_mrm\_read\_data::mrm\_read\_total\_runoff().

## 17.40 mo\_mrm\_init Module Reference

Wrapper for initializing Routing.

### Functions/Subroutines

- subroutine, public [mrm\\_configuration](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param, ReadLatLon)
- subroutine, public [mrm\\_init](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param, ReadLatLon)
  - Initialize all mRM variables at all levels (i.e., L0, L1, and L11).*
- subroutine [print\\_startup\\_message](#) (file\_namelist, file\_namelist\_param)
  - TODO: add description.*
- subroutine [config\\_output](#)
  - TODO: add description.*
- subroutine, public [variables\\_default\\_init\\_routing](#)
  - Default initialization mRM related L11 variables.*
- subroutine [l0\\_check\\_input\\_routing](#) (L0Domain\_iDomain)
  - TODO: add description.*
- subroutine [variables\\_alloc\\_routing](#) (iDomain)
  - TODO: add description.*

### 17.40.1 Detailed Description

Wrapper for initializing Routing.

Calling all routines to initialize all mRM variables

#### Authors

Luis Samaniego, Rohini Kumar and Stephan Thober

#### Date

Aug 2015

### 17.40.2 Function/Subroutine Documentation

#### 17.40.2.1 config\_output()

```
subroutine mo_mrm_init::config_output
```

TODO: add description.

TODO: add description

#### Authors

Robert Schweppe

#### Date

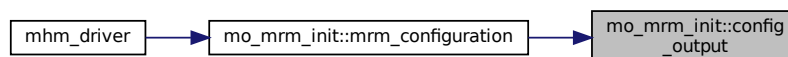
Jun 2018

Definition at line 412 of file mo\_mrm\_init.f90.

References mo\_mrm\_global\_variables::dirgauges, mo\_common\_variables::dircover, mo\_common\_variables::dirmorpho, mo\_common\_variables::dirout, mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_file::file\_defoutput, mo\_mrm\_file::file\_namelist\_mrm, and mo\_mrm\_file::file\_namelist\_param\_mrm.

Referenced by mrm\_configuration().

Here is the caller graph for this function:



#### 17.40.2.2 IO\_check\_input\_routing()

```
subroutine mo_mrm_init::io_check_input_routing (
    integer(i4), intent(in) LODomain_iDomain )
```

TODO: add description.

TODO: add description

#### Parameters

|    |  |
|----|--|
| in | <i>integer(i4) :: L0Domain_iDomain</i> |
|----|--|

#### Authors

Robert Schweppe

#### Date

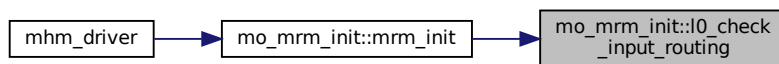
Jun 2018

Definition at line 548 of file mo\_mrm\_init.f90.

References mo\_mrm\_global\_variables::l0\_facc, mo\_mrm\_global\_variables::l0\_fdir, mo\_common\_variables::level0, and mo\_common\_constants::nodata\_i4.

Referenced by mrm\_init().

Here is the caller graph for this function:



### 17.40.2.3 mrm\_configuration()

```

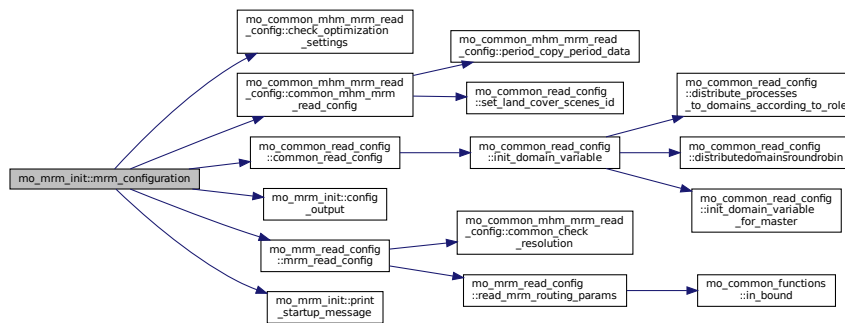
subroutine, public mo_mrm_init::mrm_configuration (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist,
    character(*), intent(in) file_namelist_param,
    integer, intent(in) unamelist_param,
    logical, intent(out) ReadLatLon )
  
```

Definition at line 30 of file mo\_mrm\_init.f90.

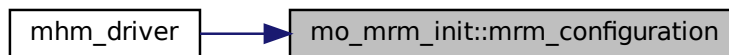
References mo\_common\_mhm\_mrm\_read\_config::check\_optimization\_settings(), mo\_common\_mhm\_mrm\_read\_config::common\_mhm\_mrm\_read\_config(), mo\_common\_read\_config::common\_read\_config(), config\_output(), mo\_common\_mhm\_mrm\_variables::mrm\_coupling\_mode, mo\_mrm\_read\_config::mrm\_read\_config(), print\_startup\_message(), mo\_common\_variables::processmatrix, and mo\_mrm\_global\_variables::riv\_temp\_pcs.

Referenced by mhm\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.40.2.4 mrm\_init()

```
subroutine, public mo_mrm_init::mrm_init (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist,
    character(*), intent(in) file_namelist_param,
    integer, intent(in) unamelist_param,
    logical, intent(inout), optional ReadLatLon )
```

Initialize all mRM variables at all levels (i.e., L0, L1, and L11).

Initialize all mRM variables at all levels (i.e., L0, L1, and L11) either with default values or with values from restart file. The L0 mask (LO\_mask), L0 elevation (LO\_elev), and L0 land cover (LO\_LCover) can be provided as optional variables to save memory because these variable will then not be read in again.

#### Parameters

|    |  |  |
|----|--|--|
| in | character(*) :: file_namelist, file_namelist_param |  |
| in | integer :: unamelist, unamelist_param              |  |
| in | character(*) :: file_namelist, file_namelist_param |  |
| in | integer :: unamelist, unamelist_param              |  |

#### Authors

Stephan Thober

## Date

Aug 2015

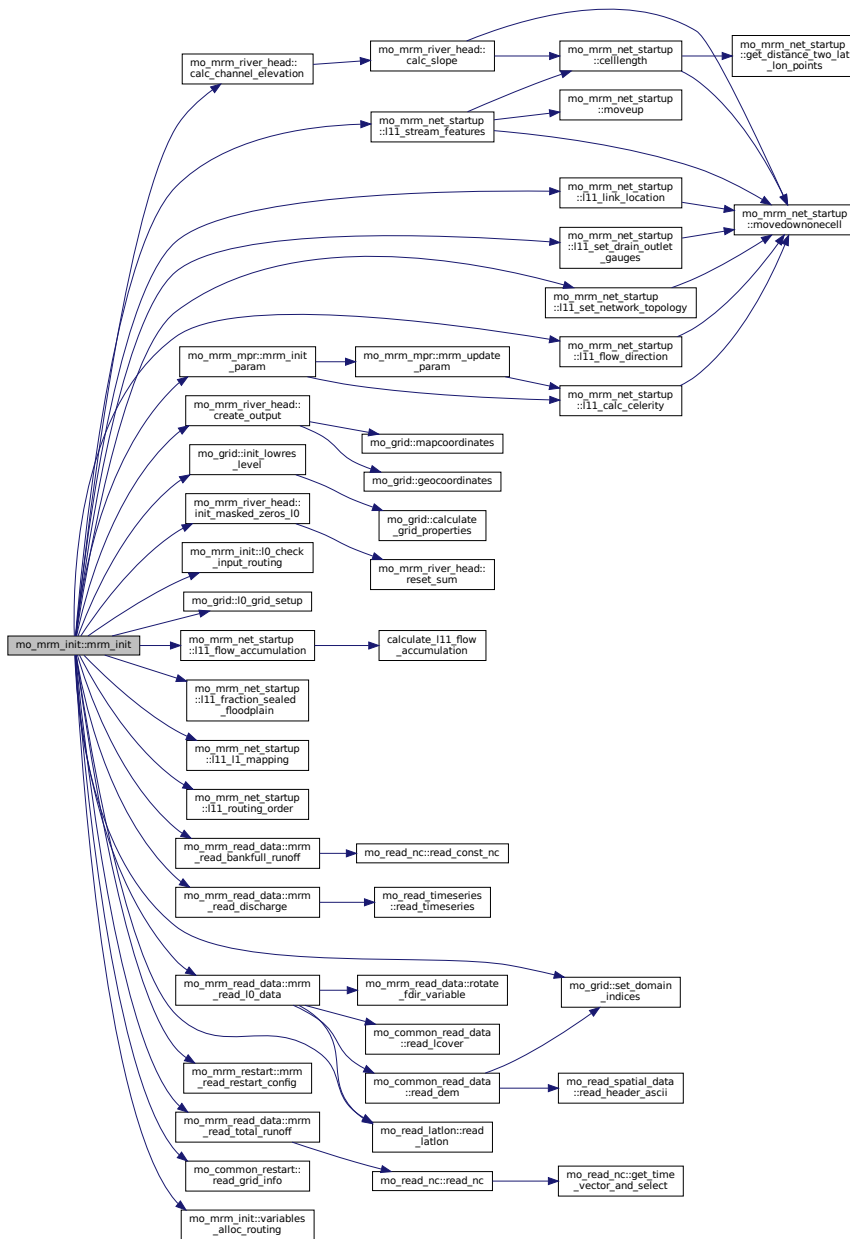
Definition at line 113 of file mo\_mrm\_init.f90.

References mo\_mrm\_river\_head::calc\_channel\_elevation(), mo\_mrm\_river\_head::create\_output(), mo\_mrm\_↔  
 global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_common\_variables::global\_parameters,  
 mo\_mrm\_global\_variables::gw\_coupling, mo\_grid::init\_lowres\_level(), mo\_mrm\_river\_head::init\_masked\_zeros\_↔  
 l0(), l0\_check\_input\_routing(), mo\_grid::l0\_grid\_setup(), mo\_mrm\_global\_variables::l0\_l11\_remap, mo\_common\_↔  
 \_variables::l0\_l1\_remap, mo\_mrm\_global\_variables::l0\_river\_head\_mon\_sum, mo\_mrm\_net\_startup::l11\_↔  
 flow\_accumulation(), mo\_mrm\_net\_startup::l11\_flow\_direction(), mo\_mrm\_net\_startup::l11\_fraction\_sealed\_↔  
 floodplain(), mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_net\_startup::l11\_l1\_mapping(), mo\_mrm\_global\_↔  
 variables::l11\_length, mo\_mrm\_net\_startup::l11\_link\_location(), mo\_mrm\_global\_variables::l11\_netperm, mo\_↔  
 mrm\_global\_variables::l11\_noutlets, mo\_mrm\_net\_startup::l11\_routing\_order(), mo\_mrm\_net\_startup::l11\_set\_↔  
 \_drain\_outlet\_gauges(), mo\_mrm\_net\_startup::l11\_set\_network\_topology(), mo\_mrm\_net\_startup::l11\_stream\_↔  
 \_features(), mo\_mrm\_global\_variables::l1\_l11\_remap, mo\_common\_variables::level0, mo\_common\_variables\_↔  
 ::level1, mo\_mrm\_global\_variables::level11, mo\_common\_mhm\_mrm\_variables::mrm\_coupling\_mode, mo\_↔  
 mrm\_mpr::mrm\_init\_param(), mo\_mrm\_read\_data::mrm\_read\_bankfull\_runoff(), mo\_mrm\_read\_data::mrm\_↔  
 read\_discharge(), mo\_mrm\_read\_data::mrm\_read\_l0\_data(), mo\_mrm\_restart::mrm\_read\_restart\_config(), mo\_↔  
 \_common\_mhm\_mrm\_variables::mrm\_read\_river\_network, mo\_mrm\_read\_data::mrm\_read\_total\_runoff(), mo\_↔  
 \_common\_mhm\_mrm\_variables::mrmfilerestartin, mo\_common\_constants::nodata\_dp, mo\_common\_constants\_↔  
 ::nodata\_i4, mo\_common\_variables::processmatrix, mo\_common\_restart::read\_grid\_info(), mo\_read\_latlon\_↔  
 ::read\_latlon(), mo\_common\_variables::resolutionhydrology, mo\_common\_mhm\_mrm\_variables::resolutionrouting,  
 mo\_mrm\_global\_variables::riv\_temp\_pcs, mo\_grid::set\_domain\_indices(), and variables\_alloc\_routing().

Referenced by mhm\_driver().



Here is the call graph for this function:



Here is the caller graph for this function:



### 17.40.2.5 print\_startup\_message()

```
subroutine mo_mrm_init::print_startup_message (
    character(*), intent(in) file_namelist,
    character(*), intent(in) file_namelist_param )
```

TODO: add description.

TODO: add description

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>character(*) :: file_namelist, file_namelist_param</i> |  |
| in | <i>character(*) :: file_namelist, file_namelist_param</i> |  |

#### Authors

Robert Schweppe

#### Date

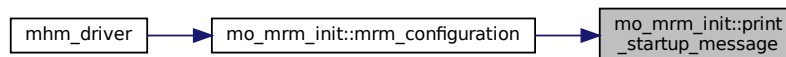
Jun 2018

Definition at line 351 of file mo\_mrm\_init.f90.

References mo\_mrm\_file::file\_defoutput, mo\_mrm\_file::file\_main, mo\_mrm\_file::version, and mo\_mrm\_file::version\_date.

Referenced by mrm\_configuration().

Here is the caller graph for this function:



### 17.40.2.6 variables\_alloc\_routing()

```
subroutine mo_mrm_init::variables_alloc_routing (
    integer(i4), intent(in) iDomain )
```

TODO: add description.

TODO: add description

#### Parameters

|    |                               |  |
|----|-------------------------------|--|
| in | <i>integer(i4) :: iDomain</i> |  |
|----|-------------------------------|--|

**Authors**

Robert Schweppe

**Date**

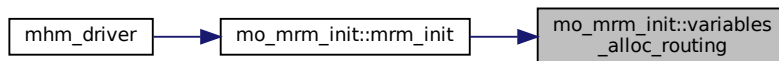
Jun 2018

Definition at line 606 of file mo\_mrm\_init.f90.

References `mo_common_variables::domainmeta`, `mo_mrm_global_variables::l0_celerity`, `mo_mrm_global_variables::l11_c1`, `mo_mrm_global_variables::l11_c2`, `mo_mrm_global_variables::l11_celerity`, `mo_mrm_global_variables::l11_k`, `mo_mrm_global_variables::l11_qmod`, `mo_mrm_global_variables::l11_qout`, `mo_mrm_global_variables::l11_qtin`, `mo_mrm_global_variables::l11_qtr`, `mo_mrm_global_variables::l11_xi`, `mo_common_variables::level0`, `mo_mrm_global_variables::level11`, and `mo_mrm_constants::nroutingstates`.

Referenced by `mrm_init()`.

Here is the caller graph for this function:

**17.40.2.7 variables\_default\_init\_routing()**

```
subroutine, public mo_mrm_init::variables_default_init_routing
```

Default initialization mRM related L11 variables.

Default initialization of mHM related L11 variables (e.g., states, fluxes, and parameters) as per given constant values given in [mo\\_mhm\\_constants](#). Variables initialized here is defined in the [mo\\_global\\_variables.f90](#) file. Only Variables that are defined in the `variables_alloc` subroutine are initialized here. If a variable is added or removed here, then it also has to be added or removed in the subroutine `state_variables_set` in the module [mo\\_restart](#) and in the subroutine `set_state` in the module `mo_set_netcdf_restart`. ADDITIONAL INFORMATION [variables\\_default\\_init\\_routing](#) call [variables\\_default\\_init\(\)](#)

**Authors**

Stephan Thober, Rohini Kumar, and Juliane Mai

**Date**

Aug 2015

**Authors**

Robert Schweppe

**Date**

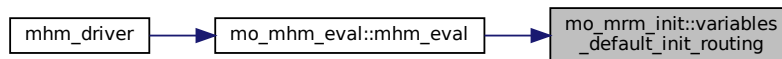
Jun 2018

Definition at line 495 of file mo\_mrm\_init.f90.

References mo\_mrm\_global\_variables::l11\_c1, mo\_mrm\_global\_variables::l11\_c2, mo\_mrm\_global\_variables::l11\_k, mo\_mrm\_global\_variables::l11\_qmod, mo\_mrm\_global\_variables::l11\_qout, mo\_mrm\_global\_variables::l11\_qtin, mo\_mrm\_global\_variables::l11\_qtr, mo\_mrm\_global\_variables::l11\_xi, and mo\_common\_constants::p1\_initstatefluxes.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the caller graph for this function:



## 17.41 mo\_mrm\_mpr Module Reference

Perform Multiscale Parameter Regionalization on Routing Parameters.

### Functions/Subroutines

- subroutine, public [reg\\_rout](#) (param, length, slope, fFPimp, TS, C1, C2)  
*Regionalized routing.*
- subroutine, public [mrm\\_init\\_param](#) (iDomain, param)  
*TODO: add description.*
- subroutine, public [mrm\\_update\\_param](#) (iDomain, param)  
*TODO: add description.*

#### 17.41.1 Detailed Description

Perform Multiscale Parameter Regionalization on Routing Parameters.

This module contains the subroutine for calculating the regionalized routing parameters (beta-parameters) given the five global routing parameters (gamma) at the level 0 scale.

**Authors**

Luis Samaniego, Stephan Thober

**Date**

Aug 2015

#### 17.41.2 Function/Subroutine Documentation

### 17.41.2.1 mrm\_init\_param()

```
subroutine, public mo_mrm_mpr::mrm_init_param (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(:), intent(in) param )
```

TODO: add description.

TODO: add description

#### Parameters

|    |  |   |
|----|--|---|
| in | <i>integer(i4) :: iDomain</i>          | domain number                                 |
| in | <i>real(dp), dimension(:) :: param</i> | input parameter (param(1) is celerity in m/s) |

#### Authors

Stephan Thober, Matthias Kelbling

#### Date

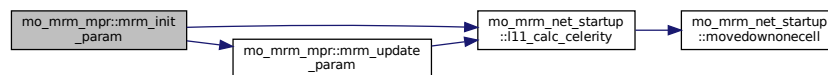
Jun 2018

Definition at line 144 of file mo\_mrm\_mpr.f90.

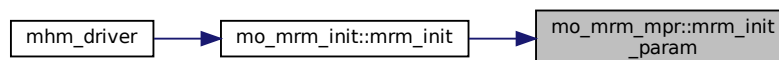
References mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_constants↔  
::given\_ts, mo\_common\_variables::iflag\_cordinate\_sys, mo\_mrm\_net\_startup::l11\_calc\_celerity(), mo\_mrm↔  
global\_variables::l11\_celerity, mo\_mrm\_global\_variables::l11\_tsrou, mo\_mrm\_global\_variables::level11, mrm↔  
\_update\_param(), mo\_common\_mhm\_mrm\_variables::optimize, mo\_common\_variables::processmatrix, mo↔  
common\_mhm\_mrm\_variables::resolutionrouting, and mo\_common\_mhm\_mrm\_variables::timestep.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.41.2.2 mrm\_update\_param()

```
subroutine, public mo_mrm_mpr::mrm_update_param (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(1), intent(in) param )
```

TODO: add description.

TODO: add description

#### Parameters

|    |  |                            |
|----|--|----------------------------|
| in | <i>integer(i4) :: iDomain</i>          | domain number              |
| in | <i>real(dp), dimension(1) :: param</i> | celerity parameter [m s-1] |

#### Authors

Stephan Thober, Matthias Kelbling

#### Date

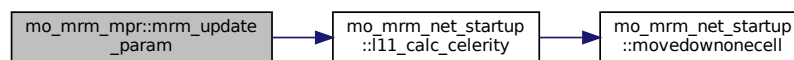
Jun 2018

Definition at line 241 of file mo\_mrm\_mpr.f90.

References `mo_mrm_constants::given_ts`, `mo_common_variables::iflag_cordinate_sys`, `mo_mrm_global_variables::l11_c1`, `mo_mrm_global_variables::l11_c2`, `mo_mrm_net_startup::l11_calc_celerity()`, `mo_mrm_global_variables::l11_celerity`, `mo_mrm_global_variables::l11_length`, `mo_mrm_global_variables::l11_noutlets`, `mo_mrm_global_variables::l11_tsrount`, `mo_mrm_global_variables::level11`, `mo_common_mhm_mrm_variables::optimize`, `mo_common_variables::processmatrix`, `mo_common_mhm_mrm_variables::resolutionrouting`, `mo_mrm_constants::rount_space_weight`, and `mo_common_mhm_mrm_variables::timestep`.

Referenced by `mo_mhm_eval::mhm_eval()`, and `mrm_init_param()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.41.2.3 reg\_rout()

```

subroutine, public mo_mrm_mpr::reg_rout (
    real(dp), dimension(5), intent(in) param,
    real(dp), dimension(:), intent(in) length,
    real(dp), dimension(:), intent(in) slope,
    real(dp), dimension(:), intent(in) fFPimp,
  
```

```

real(dp), intent(in) TS,
real(dp), dimension(:), intent(out) C1,
real(dp), dimension(:), intent(out) C2 )

```

Regionalized routing.

sets up the Regionalized Routing parameters Global parameters needed (see mhm\_parameter.nml):

- param(1) = muskingumTravelTime\_constant
- param(2) = muskingumTravelTime\_riverLength
- param(3) = muskingumTravelTime\_riverSlope
- param(4) = muskingumTravelTime\_impervious
- param(5) = muskingumAttenuation\_riverSlope

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>real(dp), dimension(5) :: param</i>  | input parameter                                   |
| in  | <i>real(dp), dimension(:) :: length</i> | [m] total length                                  |
| in  | <i>real(dp), dimension(:) :: slope</i>  | average slope                                     |
| in  | <i>real(dp), dimension(:) :: fFPimp</i> | fraction of the flood plain with impervious layer |
| in  | <i>real(dp) :: TS</i>                   | - [h] time step in                                |
| out | <i>real(dp), dimension(:) :: C1</i>     | routing parameter C1 (Chow, 25-41)                |
| out | <i>real(dp), dimension(:) :: C2</i>     | routing parameter C2 ("                           |

#### Authors

Stephan Thober, Rohini Kumar

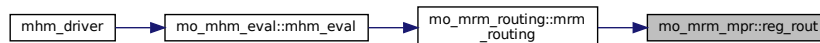
#### Date

Dec 2012

Definition at line 60 of file mo\_mrm\_mpr.f90.

Referenced by mo\_mrm\_routing::mrm\_routing().

Here is the caller graph for this function:



## 17.42 mo\_mrm\_net\_startup Module Reference

Startup drainage network for mHM.

### Functions/Subroutines

- subroutine, public [l11\\_l1\\_mapping](#) (iDomain)  
*TODO: add description.*
- subroutine, public [l11\\_flow\\_direction](#) (iDomain)

- Determine the flow direction of the upscaled river network at level L11.*

  - subroutine, public [l11\\_set\\_network\\_topology](#) (iDomain)
    - Set network topology.*
  - subroutine, public [l11\\_routing\\_order](#) (iDomain)
    - Find routing order, headwater cells and sink.*
  - subroutine, public [l11\\_link\\_location](#) (iDomain)
    - Estimate the LO (row,col) location for each routing link at level L11.*
  - subroutine, public [l11\\_set\\_drain\\_outlet\\_gauges](#) (iDomain)
    - Draining cell identification and Set gauging node.*
  - subroutine, public [l11\\_stream\\_features](#) (iDomain)
    - Stream features (stream network and floodplain)*
  - subroutine, public [l11\\_fraction\\_sealed\\_floodplain](#) (LCClassImp, do\_init)
    - Fraction of the flood plain with impervious cover.*
  - subroutine [moveup](#) (elev0, fDir0, fi, fj, ss, nn)
    - TODO: add description.*
  - subroutine [movedownonecell](#) (fDir, iRow, jCol)
    - TODO: add description.*
  - subroutine [celllength](#) (iDomain, fDir, iRow, jCol, iCoorSystem, length)
    - TODO: add description.*
  - subroutine, public [get\\_distance\\_two\\_lat\\_lon\\_points](#) (lat1, long1, lat2, long2, distance\_out)
    - estimate distance in [m] between two points in a lat-lon*
  - subroutine, public [l11\\_flow\\_accumulation](#) (iDomain)
    - Calculates L11 flow accumulation per grid cell.*
  - subroutine, public [l11\\_calc\\_celerity](#) (iDomain, param)
    - L11 celerity based on L0 elevation and L0 fAcc.*

### 17.42.1 Detailed Description

Startup drainage network for mHM.

This module initializes the drainage network at L11 in mHM.

- Delineation of drainage network at level 11.
- Setting network topology (i.e. nodes and link).
- Determining routing order.
- Determining cell locations for network links.
- Find drainage outlet.
- Determine stream (links) features.

#### Authors

Luis Samaniego

#### Date

Dec 2012

### 17.42.2 Function/Subroutine Documentation



## 17.42.2.1 celllength()

```

subroutine mo_mrm_net_startup::celllength (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) fDir,
    integer(i4), intent(in) iRow,
    integer(i4), intent(in) jCol,
    integer(i4), intent(in) iCoorSystem,
    real(dp), intent(out) length )

```

TODO: add description.

TODO: add description

## Parameters

|     |                                   |  |
|-----|-----------------------------------|--|
| in  | <i>integer(i4) :: iDomain</i>     |  |
| in  | <i>integer(i4) :: fDir</i>        |  |
| in  | <i>integer(i4) :: iRow</i>        |  |
| in  | <i>integer(i4) :: jCol</i>        |  |
| in  | <i>integer(i4) :: iCoorSystem</i> |  |
| out | <i>real(dp) :: length</i>         |  |

## Authors

Robert Schweppe

## Date

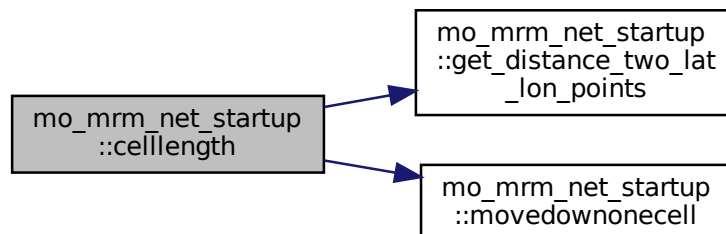
Jun 2018

Definition at line 1823 of file mo\_mrm\_net\_startup.f90.

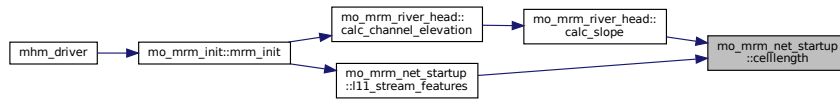
References mo\_common\_variables::domainmeta, get\_distance\_two\_lat\_lon\_points(), mo\_common\_variables↵  
::level0, and movedownonecell().

Referenced by mo\_mrm\_river\_head::calc\_slope(), and l11\_stream\_features().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.42.2.2 get\_distance\_two\_lat\_lon\_points()

```

subroutine, public mo_mrm_net_startup::get_distance_two_lat_lon_points (
    real(dp), intent(in) lat1,
    real(dp), intent(in) long1,
    real(dp), intent(in) lat2,
    real(dp), intent(in) long2,
    real(dp), intent(out) distance_out )
  
```

estimate distance in [m] between two points in a lat-lon

estimate distance in [m] between two points in a lat-lon Code is based on one that is implemented in the VIC-3L model

#### Parameters

|     |   |                                 |
|-----|---|---------------------------------|
| in  | <i>real(dp) :: lat1, long1, lat2, long2</i> | latitude of point-1             |
| in  | <i>real(dp) :: lat1, long1, lat2, long2</i> | longitude of point-1            |
| in  | <i>real(dp) :: lat1, long1, lat2, long2</i> | latitude of point-2             |
| in  | <i>real(dp) :: lat1, long1, lat2, long2</i> | longitude of point-2            |
| out | <i>real(dp) :: distance_out</i>             | distance between two points [m] |

#### Authors

Rohini Kumar

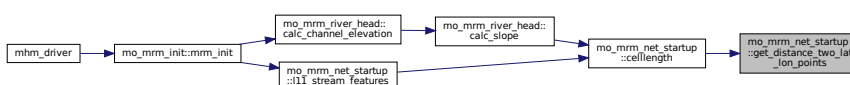
#### Date

May 2014

Definition at line 1917 of file mo\_mrm\_net\_startup.f90.

Referenced by celllength().

Here is the caller graph for this function:



### 17.42.2.3 l11\_calc\_celerity()

```
subroutine, public mo_mrm_net_startup::l11_calc_celerity (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(:), intent(in) param )
```

L11 celerity based on L0 elevation and L0 fAcc.

L11 celerity based on L0 elevation and L0 fAcc

#### Parameters

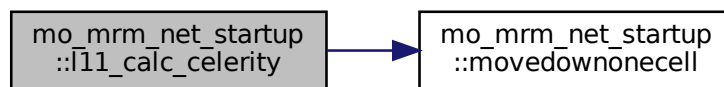
|    |  |  |
|----|--|--|
| in |  |  |
|----|--|--|

Definition at line 2203 of file mo\_mrm\_net\_startup.f90.

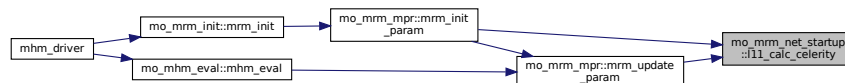
References mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_fdir, mo\_mpr\_global\_variables::l0\_slope, movedownonecell(), mo\_common\_constants::nodata\_dp, and mo\_common\_constants::nodata\_i4.

Referenced by mo\_mrm\_mpr::mrm\_init\_param(), and mo\_mrm\_mpr::mrm\_update\_param().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.42.2.4 l11\_flow\_accumulation()

```
subroutine, public mo_mrm_net_startup::l11_flow_accumulation (
    integer(i4), intent(in) iDomain )
```

Calculates L11 flow accumulation per grid cell.

Calculates L11 flow accumulation per grid cell using L11\_fDir and L11\_cellarea. L11\_flow\_accumulation contains the recursive subroutine `calculate_L11_flow_accumulation` iDomain

#### Author

Matthias Kelbling

## Date

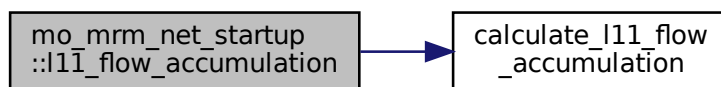
Aug 2017

Definition at line 2013 of file `mo_mrm_net_startup.f90`.

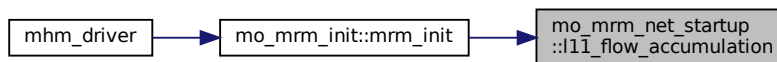
References `calculate_l11_flow_accumulation()`, `mo_mrm_global_variables::l11_fdir`, `mo_mrm_global_variables::level11`, `mo_common_constants::nodata_dp`, and `mo_common_constants::nodata_i4`.

Referenced by `mo_mrm_init::mrm_init()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.42.2.5 l11\_flow\_direction()

```

subroutine, public mo_mrm_net_startup::l11_flow_direction (
    integer(i4), intent(in) iDomain )
  
```

Determine the flow direction of the upscaled river network at level L11.

The hydrographs generated at each cell are routed through the drainage network at level-11 towards their outlets. The drainage network at level-11 is conceptualized as a graph whose nodes are hypothetically located at the center of each grid cell connected by links that represent the river reaches. The flow direction of a link correspond to the direction towards a neighboring cell in which the net flow accumulation (outflows minus inflows) attains its maximum value. The net flow accumulation across a cell's boundary at level-11 is estimated based on flow direction and flow accumulation obtained at level-0 ([Routing Network](#)). Note: level-1 denotes the modeling level, whereas level-L11 is at least as coarse as level-1. Experience has shown that routing can be done at a coarser resolution as level-1, hence the level-11 was introduced.

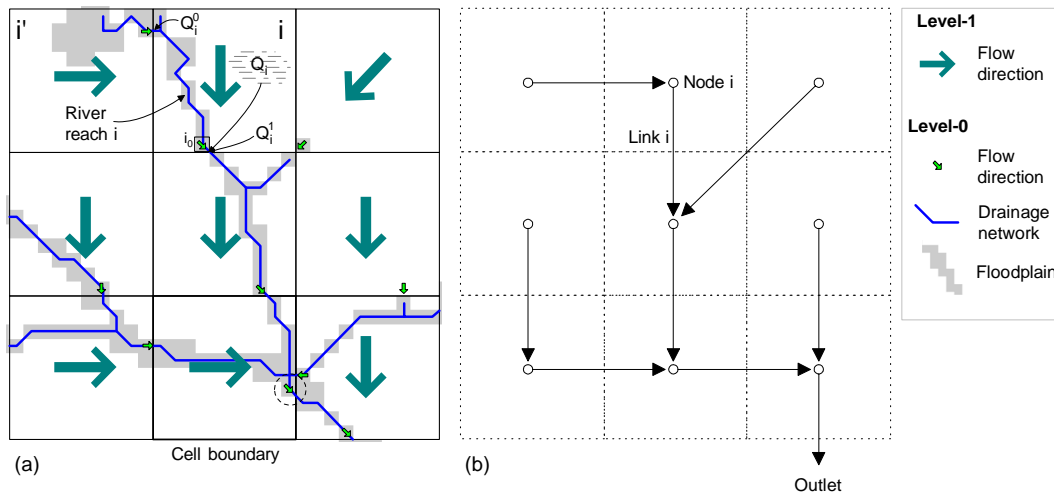


Figure 17.1: Upscaling routing network from L0 to L1 (or L11)

The left panel depicts a schematic derivation of a drainage network at the level-11 based on level-0 flow direction and flow accumulation. The dotted line circle denotes the point with the highest flow accumulation within a grid cell. The topology of a typical drainage routing network at level-11 is shown in the right panel. Gray color areas denote the flood plains estimated in mo\_init\_mrm, where the network upscaling is also carried out. For the sake of simplicity, it is assumed that all runoff leaving a given cell would exit through a major direction. Note that multiple outlets can exist within the modelling domain. If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module mo\_restart and in the subroutine set\_L11\_config in module mo\_set\_netcdf\_restart ADDITIONAL INFORMATION L11\_flow\_direction

Parameters

|    |                        |           |
|----|------------------------|-----------|
| in | integer(i4) :: iDomain | Domain Id |
|----|------------------------|-----------|

Authors

Luis Samaniego

Date

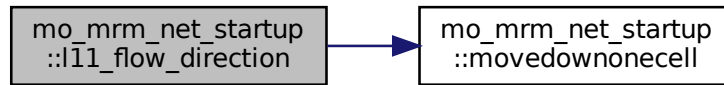
Dec 2005

Definition at line 221 of file mo\_mrm\_net\_startup.f90.

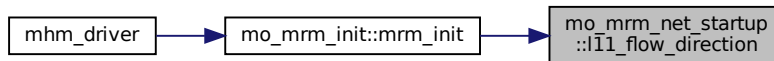
References mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_drasc, mo\_mrm\_global\_variables::l0\_facc, mo\_mrm\_global\_variables::l0\_fdir, mo\_mrm\_global\_variables::l0\_l11\_remap, mo\_mrm\_global\_variables::l11\_colout, mo\_mrm\_global\_variables::l11\_fdir, mo\_mrm\_global\_variables::l11\_noutlets, mo\_mrm\_global\_variables::l11\_rowout, mo\_common\_variables::level0, mo\_mrm\_global\_variables::level11, movedownonecell(), and mo\_common\_constants::nodata\_i4.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.42.2.6 l11\_fraction\_sealed\_floodplain()

```

subroutine, public mo_mrm_net_startup::l11_fraction_sealed_floodplain (
    integer(i4), intent(in) LCClassImp,
    logical, intent(in) do_init )
  
```

Fraction of the flood plain with impervious cover.

Fraction of the flood plain with impervious cover ([Routing Network](#)). This proportion is used to regionalize the Muskingum parameters. Samaniego et al. [15] found out that this fraction is one of the statistically significant predictor variables of peak discharge in mesoscale Domains. If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module [mo\\_restart](#) and in the subroutine set\_↔ L11\_config in module mo\_set\_netcdf\_restart

#### Parameters

|    |                                  |   |
|----|----------------------------------|---|
| in | <i>integer(i4) :: LCClassImp</i> | Impervious land cover class Id, e.g. = 2 (old code) |
| in | <i>logical :: do_init</i>        |   |

#### Authors

Luis Samaniego

#### Date

Dec 2005

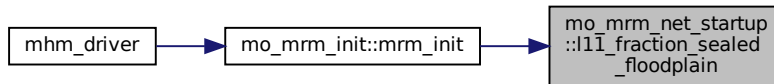
Definition at line 1503 of file mo\_mrm\_net\_startup.f90.

References mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_floodplain, mo\_common\_↔ variables::l0\_lcover, mo\_mrm\_global\_variables::l11\_afloodplain, mo\_mrm\_global\_variables::l11\_nlinkfracpimp,

mo\_mrm\_global\_variables::l11\_noutlets, mo\_common\_variables::level0, mo\_mrm\_global\_variables::level11, mo\_↔  
\_common\_variables::nlcoverscene, and mo\_common\_constants::nodata\_dp.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the caller graph for this function:



### 17.42.2.7 l11\_l1\_mapping()

```

subroutine, public mo_mrm_net_startup::l11_l1_mapping (
    integer(i4), intent(in) iDomain )
  
```

TODO: add description.

TODO: add description

#### Parameters

|    |                               |        |
|----|-------------------------------|--------|
| in | <i>integer(i4) :: iDomain</i> | domain |
|----|-------------------------------|--------|

#### Authors

Robert Schweppe

#### Date

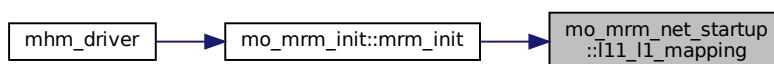
Jun 2018

Definition at line 55 of file mo\_mrm\_net\_startup.f90.

References mo\_mrm\_global\_variables::l11\_l1\_id, mo\_mrm\_global\_variables::l1\_l11\_id, mo\_common\_variables\_↔  
::level1, mo\_mrm\_global\_variables::level11, and mo\_common\_constants::nodata\_i4.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the caller graph for this function:



### 17.42.2.8 l11\_link\_location()

```
subroutine, public mo_mrm_net_startup::l11_link_location (
    integer(i4), intent(in) iDomain )
```

Estimate the LO (row,col) location for each routing link at level L11.

If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module [mo\\_restart](#) and in the subroutine set\_L11\_config in module [mo\\_set\\_netcdf\\_restart](#)

#### Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | Domain Id |
|----|-------------------------------|-----------|

#### Authors

Luis Samaniego

#### Date

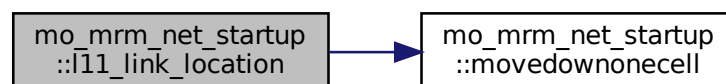
Dec 2005

Definition at line 881 of file [mo\\_mrm\\_net\\_startup.f90](#).

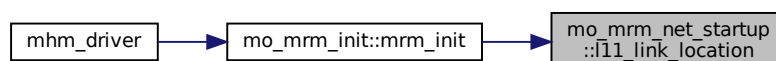
References [mo\\_mrm\\_global\\_variables::domain\\_mrm](#), [mo\\_common\\_variables::domainmeta](#), [mo\\_mrm\\_global\\_variables::l0\\_drasc](#), [mo\\_mrm\\_global\\_variables::l0\\_fdir](#), [mo\\_mrm\\_global\\_variables::l11\\_colout](#), [mo\\_mrm\\_global\\_variables::l11\\_fcol](#), [mo\\_mrm\\_global\\_variables::l11\\_fromn](#), [mo\\_mrm\\_global\\_variables::l11\\_frow](#), [mo\\_mrm\\_global\\_variables::l11\\_netperm](#), [mo\\_mrm\\_global\\_variables::l11\\_noutlets](#), [mo\\_mrm\\_global\\_variables::l11\\_rowout](#), [mo\\_mrm\\_global\\_variables::l11\\_tcol](#), [mo\\_mrm\\_global\\_variables::l11\\_trow](#), [mo\\_common\\_variables::level0](#), [mo\\_mrm\\_global\\_variables::level11](#), [movedownonecell\(\)](#), and [mo\\_common\\_constants::nodata\\_i4](#).

Referenced by [mo\\_mrm\\_init::mrm\\_init\(\)](#).

Here is the call graph for this function:



Here is the caller graph for this function:





### 17.42.2.9 l11\_routing\_order()

```
subroutine, public mo_mrm_net_startup::l11_routing_order (
    integer(i4), intent(in) iDomain )
```

Find routing order, headwater cells and sink.

Find routing order, headwater cells and sink. If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module [mo\\_restart](#) and in the subroutine set\_L11\_config in module [mo\\_set\\_netcdf\\_restart](#)

#### Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | Domain Id |
|----|-------------------------------|-----------|

#### Authors

Luis Samaniego

#### Date

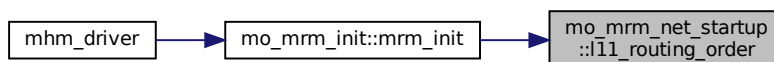
Dec 2005

Definition at line 722 of file [mo\\_mrm\\_net\\_startup.f90](#).

References [mo\\_mrm\\_global\\_variables::l11\\_fdir](#), [mo\\_mrm\\_global\\_variables::l11\\_fromn](#), [mo\\_mrm\\_global\\_variables::l11\\_label](#), [mo\\_mrm\\_global\\_variables::l11\\_netperm](#), [mo\\_mrm\\_global\\_variables::l11\\_noutlets](#), [mo\\_mrm\\_global\\_variables::l11\\_rorder](#), [mo\\_mrm\\_global\\_variables::l11\\_sink](#), [mo\\_mrm\\_global\\_variables::l11\\_ton](#), [mo\\_mrm\\_global\\_variables::level11](#), and [mo\\_common\\_constants::nodata\\_i4](#).

Referenced by [mo\\_mrm\\_init::mrm\\_init\(\)](#).

Here is the caller graph for this function:



### 17.42.2.10 l11\_set\_drain\_outlet\_gauges()

```
subroutine, public mo_mrm_net_startup::l11_set_drain_outlet_gauges (
    integer(i4), intent(in) iDomain )
```

Draining cell identification and Set gauging node.

Perform the following tasks:

- Draining cell identification (cell at L0 to draining cell outlet at L11).
- Set gauging nodes If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module [mo\\_restart](#) and in the subroutine set\_L11\_config in module [mo\\_set\\_netcdf\\_restart](#)

## Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | Domain Id |
|----|-------------------------------|-----------|

## Authors

Luis Samaniego

## Date

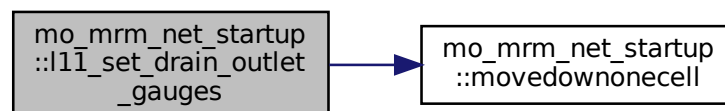
Dec 2005

Definition at line 1083 of file mo\_mrm\_net\_startup.f90.

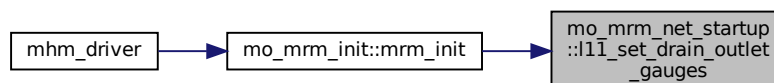
References mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_dracell, mo\_mrm\_global\_variables::l0\_drasc, mo\_mrm\_global\_variables::l0\_fdir, mo\_mrm\_global\_variables::l0\_gaugeloc, mo\_mrm\_global\_variables::l0\_inflowgaugeloc, mo\_mrm\_global\_variables::l0\_l11\_remap, mo\_common\_variables::level0, movedownonecell(), and mo\_common\_constants::nodata\_i4.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.42.2.11 l11\_set\_network\_topology()

```

subroutine, public mo_mrm_net_startup::l11_set_network_topology (
    integer(i4), intent(in) iDomain )
  
```

Set network topology.

Set network topology from and to node for all links at level-11 ([Routing Network](#)). If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module [mo\\_restart](#) and in the subroutine set\_L11\_config in module mo\_set\_netcdf\_restart.

## Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | Domain Id |
|----|-------------------------------|-----------|

## Authors

Luis Samaniego

## Date

Dec 2005

Definition at line 624 of file mo\_mrm\_net\_startup.f90.

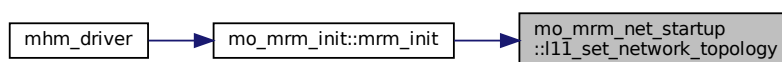
References mo\_mrm\_global\_variables::l11\_fdir, mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_global\_variables::l11\_ton, mo\_mrm\_global\_variables::level11, movedownonecell(), and mo\_common\_constants::nodata\_i4.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.42.2.12 l11\_stream\_features()

```

subroutine, public mo_mrm_net_startup::l11_stream_features (
    integer(i4), intent(in) iDomain )
  
```

Stream features (stream network and floodplain)

Stream features (stream network and floodplain) If a variable is added or removed here, then it also has to be added or removed in the subroutine L11\_config\_set in module [mo\\_restart](#) and in the subroutine set\_L11\_config in module [mo\\_set\\_netcdf\\_restart](#)

## Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | Domain Id |
|----|-------------------------------|-----------|

## Authors

Luis Samaniego

## Date

Dec 2005

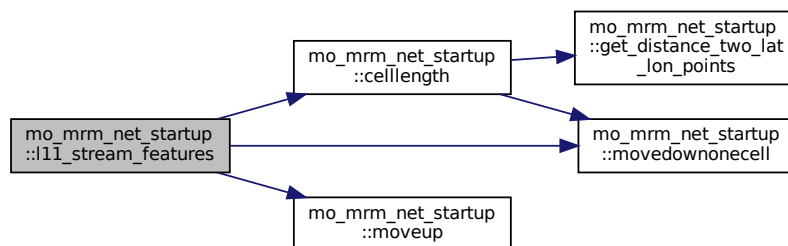
stack(nNodes, 2)

Definition at line 1227 of file mo\_mrm\_net\_startup.f90.

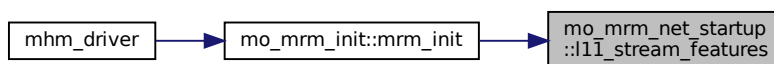
References celllength(), mo\_common\_variables::domainmeta, mo\_common\_variables::iflag\_cordinate\_sys, mo\_common\_variables::l0\_elev, mo\_mrm\_global\_variables::l0\_fdir, mo\_mrm\_global\_variables::l0\_floodplain, mo\_mrm\_global\_variables::l0\_streamnet, mo\_mrm\_global\_variables::l11\_afloodplain, mo\_mrm\_global\_variables::l11\_fcol, mo\_mrm\_global\_variables::l11\_frow, mo\_mrm\_global\_variables::l11\_length, mo\_mrm\_global\_variables::l11\_netperm, mo\_mrm\_global\_variables::l11\_noutlets, mo\_mrm\_global\_variables::l11\_slope, mo\_mrm\_global\_variables::l11\_tcol, mo\_mrm\_global\_variables::l11\_trow, mo\_common\_variables::level0, mo\_mrm\_global\_variables::level11, movedownonecell(), moveup(), mo\_common\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, and mo\_common\_variables::processmatrix.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.42.2.13 movedownonecell()**

```

subroutine mo_mrm_net_startup::movedownonecell (
    integer(i4), intent(in) fDir,
    integer(i4), intent(inout) iRow,
    integer(i4), intent(inout) jCol )
  
```

TODO: add description.

TODO: add description

Parameters

|         |                                  |  |
|---------|----------------------------------|--|
| in      | <i>integer(i4) :: fDir</i>       |  |
| in, out | <i>integer(i4) :: iRow, jCol</i> |  |
| in, out | <i>integer(i4) :: iRow, jCol</i> |  |

Authors

Robert Schweppe

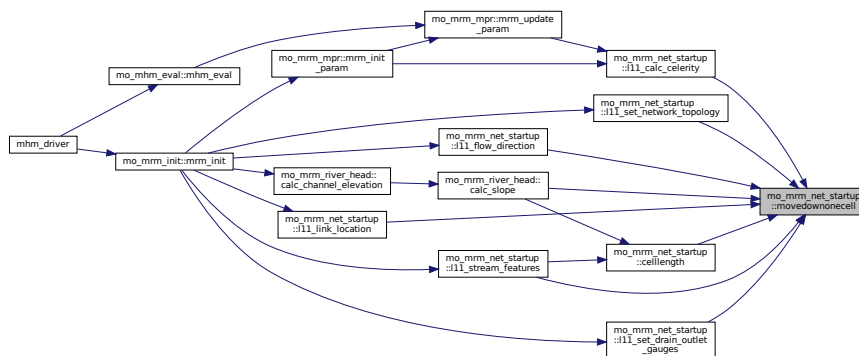
Date

Jun 2018

Definition at line 1760 of file mo\_mrm\_net\_startup.f90.

Referenced by mo\_mrm\_river\_head::calc\_slope(), celloength(), l11\_calc\_celerity(), l11\_flow\_direction(), l11\_link\_location(), l11\_set\_drain\_outlet\_gauges(), l11\_set\_network\_topology(), and l11\_stream\_features().

Here is the caller graph for this function:



17.42.2.14 moveup()

```

subroutine mo_mrm_net_startup::moveup (
    real(dp), dimension(:, :), intent(in), allocatable elev0,
    integer(i4), dimension(:, :), intent(in), allocatable fDir0,
    integer(i4), intent(in) fi,
    integer(i4), intent(in) fj,
    integer(i4), dimension(:, :), intent(inout) ss,
    integer(i4), intent(inout) nn )

```

TODO: add description.

TODO: add description

## Parameters

|         |  |                               |
|---------|--|-------------------------------|
| in      | <i>real(dp), dimension(:, :) :: elev0</i>    |                               |
| in      | <i>integer(i4), dimension(:, :) :: fDir0</i> |                               |
| in      | <i>integer(i4) :: fi, fj</i>                 | co-ordinate of the stream bed |
| in      | <i>integer(i4) :: fi, fj</i>                 | co-ordinate of the stream bed |
| in, out | <i>integer(i4), dimension(:, :) :: ss</i>    |                               |
| in, out | <i>integer(i4) :: nn</i>                     |                               |

## Authors

Robert Schweppe

## Date

Jun 2018

Definition at line 1587 of file mo\_mrm\_net\_startup.f90.

References mo\_mrm\_constants::deltah.

Referenced by l11\_stream\_features().

Here is the caller graph for this function:



## 17.43 mo\_mrm\_objective\_function\_runoff Module Reference

Objective Functions for Optimization of mHM/mRM against runoff.

### Functions/Subroutines

- real(dp) function, public [single\\_objective\\_runoff](#) (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions optimizing against runoff.*
- real(dp) function, public [single\\_objective\\_runoff\\_master](#) (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions optimizing against runoff.*
- subroutine, public [single\\_objective\\_runoff\\_subprocess](#) (eval, arg1, arg2, arg3)  
*Wrapper for objective functions optimizing against runoff.*
- subroutine, public [multi\\_objective\\_runoff](#) (parameterset, eval, multi\_objectives)  
*Wrapper for multi-objective functions where at least one is regarding runoff.*
- real(dp) function [loglikelihood\\_stddev](#) (parameterset, eval, stddev, stddev\_new, likeli\_new)  
*Logarithmic likelihood function with removed linear trend and Lag(1)-autocorrelation.*
- real(dp) function [loglikelihood\\_evin2013\\_2](#) (parameterset, eval, regularize)  
*Logarithmised likelihood with linear error model and lag(1)-autocorrelation of the relative errors.*
- real(dp) function [parameter\\_regularization](#) (paraset, prior, bounds, mask)  
*TODO: add description.*
- real(dp) function [loglikelihood\\_trend\\_no\\_autocorr](#) (parameterset, eval, stddev\_old, stddev\_new, likeli\_new)

- Logarithmic likelihood function with linear trend removed.*

  - real(dp) function [objective\\_lnnse](#) (parameterset, eval)

*Objective function of logarithmic NSE.*
- real(dp) function [objective\\_sse](#) (parameterset, eval)

*Objective function of SSE.*
- real(dp) function [objective\\_nse](#) (parameterset, eval)

*Objective function of NSE.*
- real(dp) function [objective\\_equal\\_nse\\_lnnse](#) (parameterset, eval)

*Objective function equally weighting NSE and lnNSE.*
- real(dp) function, dimension(2) [multi\\_objective\\_nse\\_lnnse](#) (parameterset, eval)

*Multi-objective function with NSE and lnNSE.*
- real(dp) function, dimension(2) [multi\\_objective\\_lnnse\\_highflow\\_lnnse\\_lowflow](#) (parameterset, eval)

*Multi-objective function with NSE and lnNSE.*
- real(dp) function, dimension(2) [multi\\_objective\\_lnnse\\_highflow\\_lnnse\\_lowflow\\_2](#) (parameterset, eval)

*Multi-objective function with NSE and lnNSE.*
- real(dp) function, dimension(2) [multi\\_objective\\_ae\\_fdc\\_lsv\\_nse\\_djf](#) (parameterset, eval)

*Multi-objective function with absolute error of Flow Duration Curves low-segment volume and nse of DJF's discharge.*
- real(dp) function [objective\\_power6\\_nse\\_lnnse](#) (parameterset, eval)

*Objective function of combined NSE and lnNSE with power of 5 i.e. the p-norm with p=5.*
- real(dp) function [objective\\_kge](#) (parameterset, eval)

*Objective function of KGE.*
- real(dp) function [objective\\_multiple\\_gauges\\_kge\\_power6](#) (parameterset, eval)

*combined objective function based on KGE raised to the power 6*
- real(dp) function [objective\\_weighted\\_nse](#) (parameterset, eval)

*Objective function of weighted NSE.*
- real(dp) function [objective\\_sse\\_boxcox](#) (parameterset, eval)

*Objective function of sum of squared errors of transformed streamflow.*
- subroutine, public [extract\\_runoff](#) (gaugeld, runoff, runoff\_agg, runoff\_obs, runoff\_obs\_mask)

*extracts runoff data from global variables*

### 17.43.1 Detailed Description

Objective Functions for Optimization of mHM/mRM against runoff.

This module provides a wrapper for several objective functions used to optimize mRM/mHM against runoff. If the objective contains besides runoff another variable like TWS move it to [mHM/mo\\_objective\\_function.f90](#). If it is only regarding runoff implement it here. All the objective functions are supposed to be minimized! (1) SO: Q: 1.0 - NSE (2) SO: Q: 1.0 - lnNSE (3) SO: Q: 1.0 - 0.5\*(NSE+lnNSE) (4) SO: Q: -1.0 \* loglikelihood with trend removed from absolute errors and then lag(1)-autocorrelation removed (5) SO: Q: ((1-NSE)\*\*6+(1-lnNSE)\*\*6)\*\*(1/6) (6) SO: Q: SSE (7) SO: Q: -1.0 \* loglikelihood with trend removed from absolute errors (8) SO: Q: -1.0 \* loglikelihood with trend removed from the relative errors and then lag(1)-autocorrelation removed (9) SO: Q: 1.0 - KGE (Kling-Gupta efficiency measure) (14) SO: Q: sum[(((1.0-KGE\_i)/ nGauges)\*\*6]\*\*(1/6) > combination of KGE of every gauging station based on a power-6 norm (16) MO: Q: 1st objective: 1.0 - NSE Q: 2nd objective: 1.0 - lnNSE (18) MO: Q: 1st objective: 1.0 - lnNSE(Q\_highflow) (95% percentile) Q: 2nd objective: 1.0 - lnNSE(Q\_lowflow) (5% of data range) (19) MO: Q: 1st objective: 1.0 - lnNSE(Q\_highflow) (non-low flow) Q: 2nd objective: 1.0 - lnNSE(Q\_lowflow) (5% of data range)eshold for Q (20) MO: Q: 1st objective: absolute difference in FDC's low-segment volume Q: 2nd objective: 1.0 - NSE of discharge of months DJF (31) SO: Q: 1.0 - wNSE - weighted NSE (32) SO: Q: SSE of boxcox-transformed streamflow

#### Authors

Juliane Mai

#### Date

Dec 2012

## 17.43.2 Function/Subroutine Documentation

### 17.43.2.1 `extract_runoff()`

```
subroutine, public mo_mrm_objective_function_runoff::extract_runoff (
    integer(i4), intent(in) gaugeId,
    real(dp), dimension(:, :), intent(in) runoff,
    real(dp), dimension(:), intent(out), allocatable runoff_agg,
    real(dp), dimension(:), intent(out), allocatable runoff_obs,
    logical, dimension(:), intent(out), allocatable runoff_obs_mask )
```

extracts runoff data from global variables

extracts simulated and measured runoff from global variables, such that they overlay exactly. For measured runoff, only the runoff during the evaluation period are cut, not succeeding nodata values. For simulated runoff, warming days as well as succeeding nodata values are neglected and the simulated runoff is aggregated to the resolution of the observed runoff. see use in this module above

#### Parameters

|     |   |                        |
|-----|---|------------------------|
| in  | <i>integer(i4) :: gaugeId</i>                   | current gauge Id       |
| in  | <i>real(dp), dimension(:, :) :: runoff</i>      | simulated runoff       |
| out | <i>real(dp), dimension(:) :: runoff_agg</i>     | aggregated simulated   |
| out | <i>real(dp), dimension(:) :: runoff_obs</i>     | extracted measured     |
| out | <i>logical, dimension(:) :: runoff_obs_mask</i> | mask of no data values |

#### Authors

Stephan Thober

#### Date

Jan 2015

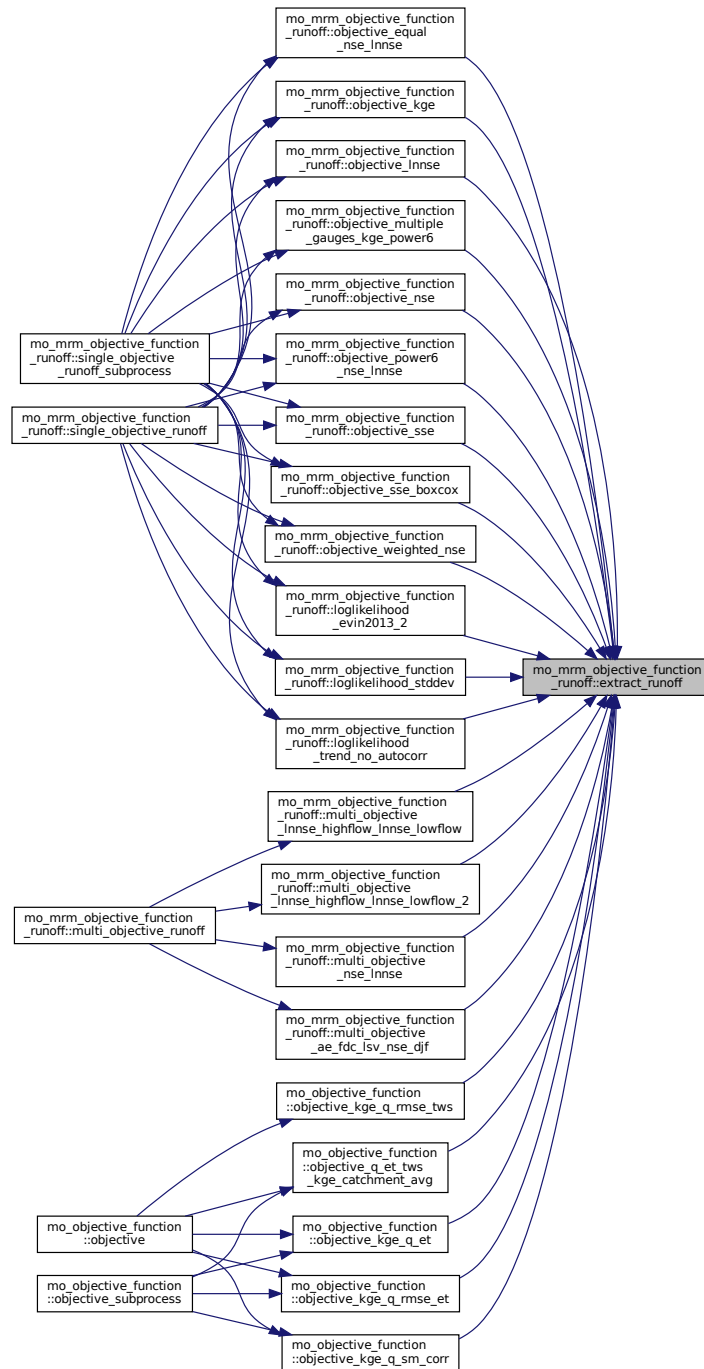
Definition at line 2475 of file `mo_mrm_objective_function_runoff.f90`.

References `mo_common_mhm_mrm_variables::evalper`, `mo_mrm_global_variables::gauge`, `mo_mrm_global_variables::nmeasperday`, `mo_common_mhm_mrm_variables::ntstepday`, and `mo_common_mhm_mrm_variables::warmingdays`.

Referenced by `loglikelihood_evin2013_2()`, `loglikelihood_stddev()`, `loglikelihood_trend_no_autocorr()`, `multi_objective_ae_fdc_lsv_nse_djf()`, `multi_objective_lnnse_highflow_lnnse_lowflow()`, `multi_objective_lnnse_highflow_lnnse_lowflow_2()`, `multi_objective_nse_lnnse()`, `objective_equal_nse_lnnse()`, `objective_kge()`, `mo_objective_function::objective_kge_q_et()`, `mo_objective_function::objective_kge_q_rmse_et()`, `mo_objective_function::objective_kge_q_rmse_tws()`, `mo_objective_function::objective_kge_q_sm_corr()`, `objective_lnnse()`, `objective_multiple_gauges_kge_power6()`, `objective_nse()`, `objective_power6_nse_lnnse()`, `mo_objective_function::objective_q_et_tws_kge_catchment_avg()`, `objective_sse()`, `objective_sse_boxcox()`, and `objective_weighted_nse()`.



Here is the caller graph for this function:



### 17.43.2.2 loglikelihood\_evin2013\_2()

```
real(dp) function mo_mrm_objective_function_runoff::loglikelihood_evin2013_2 (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
```

```
logical, intent(in), optional regularize )
```

Logarithmised likelihood with linear error model and lag(1)-autocorrelation of the relative errors.

This loglikelihood uses a linear error model and a lag(1)-autocorrelation on the relative errors. This is approach 2 of the paper Evin et al. (WRR, 2013). This is `opti_function = 8`. mHM then adds two extra (local) parameters for the error model in `mhm_driver`, which get optimised together with the other, global parameters. ADDITIONAL INFORMATION Evin et al., WRR 49, 4518-4524, 2013

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |
| in | <i>logical, optional :: regularize</i>        |  |

#### Returns

`real(dp) :: loglikelihood_evin2013_2` — logarithmic likelihood using given `stddev` but remove optimal trend and lag(1)-autocorrelation in errors (absolute between running model with `parameterset` and observation)

#### Authors

Juliane Mai and Matthias Cuntz

#### Date

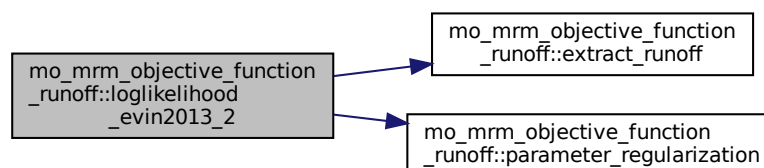
Mar 2014

Definition at line 732 of file `mo_mrm_objective_function_runoff.f90`.

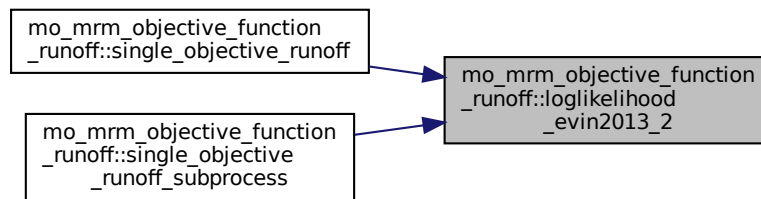
References `extract_runoff()`, `mo_common_variables::global_parameters`, and `parameter_regularization()`.

Referenced by `single_objective_runoff()`, and `single_objective_runoff_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.3 loglikelihood\_stddev()

```

real(dp) function mo_mrm_objective_function_runoff::loglikelihood_stddev (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in) stddev,
    real(dp), intent(out), optional stddev_new,
    real(dp), intent(out), optional likeli_new )
  
```

Logarithmic likelihood function with removed linear trend and Lag(1)-autocorrelation.

The logarithmic likelihood function is used when mHM runs in MCMC mode. It can also be used for optimization when selecting the likelihood in the namelist as *opti\_function*.

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>real(dp), dimension(:) :: parameterset</i> |   |
| in  | <i>procedure(eval_interface) :: eval</i>      |   |
| in  | <i>real(dp) :: stddev</i>                     | standard deviation of data  |
| out | <i>real(dp), optional :: stddev_new</i>       | standard deviation of errors with removed trend and correlation between model run using parameter set and observation |
| out | <i>real(dp), optional :: likeli_new</i>       | logarithmic likelihood determined with <i>stddev_new</i> instead of <i>stddev</i>                                     |

#### Returns

*real(dp) :: loglikelihood\_stddev* — logarithmic likelihood using given *stddev* but remove optimal trend and lag(1)-autocorrelation in errors (absolute between running model with *parameterset* and observation)

#### Authors

Juliane Mai

#### Date

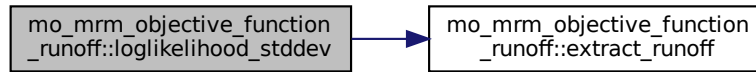
Dec 2012

Definition at line 590 of file *mo\_mrm\_objective\_function\_runoff.f90*.

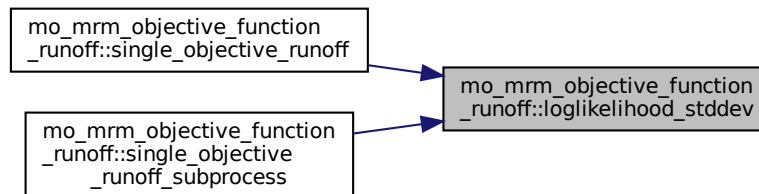
References `extract_runoff()`.

Referenced by `single_objective_runoff()`, and `single_objective_runoff_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.43.2.4 loglikelihood\_trend\_no\_autocorr()

```

real(dp) function mo_mrm_objective_function_runoff::loglikelihood_trend_no_autocorr (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in) stddev_old,
    real(dp), intent(out), optional stddev_new,
    real(dp), intent(out), optional likeli_new )
  
```

Logarithmic likelihood function with linear trend removed.

The logarithmic likelihood function is used when mHM runs in MCMC mode. It can also be used for optimization when selecting the likelihood in the namelist as *opti\_function*.

##### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>real(dp), dimension(:) :: parameterset</i> |   |
| in  | <i>procedure(eval_interface) :: eval</i>      |   |
| in  | <i>real(dp) :: stddev_old</i>                 | standard deviation of data  |
| out | <i>real(dp), optional :: stddev_new</i>       | standard deviation of errors with removed trend between model run using parameter set and observation |
| out | <i>real(dp), optional :: likeli_new</i>       | logarithmic likelihood determined with <i>stddev_new</i> instead of <i>stddev</i>                     |

**Returns**

real(dp) :: loglikelihood\_trend\_no\_autocorr — logarithmic likelihood using given stddev but remove optimal trend in errors (absolute between running model with parameterset and observation)

**Authors**

Juliane Mai and Matthias Cuntz

**Date**

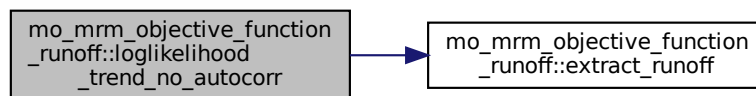
Mar 2014

Definition at line 951 of file mo\_mrm\_objective\_function\_runoff.f90.

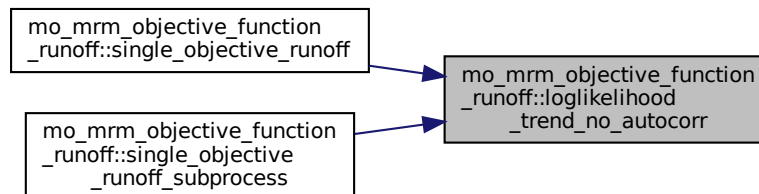
References extract\_runoff().

Referenced by single\_objective\_runoff(), and single\_objective\_runoff\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.43.2.5 multi\_objective\_ae\_fdc\_lsv\_nse\_djf()**

```

real(dp) function, dimension(2) mo_mrm_objective_function_runoff::multi_objective_ae_fdc_lsv_nse_djf (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Multi-objective function with absolute error of Flow Duration Curves low-segment volume and nse of DJF's discharge.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. The first objective is using the routine "`FlowDurationCurves`" from "`mo_signatures`" to determine the low-segment volume of the FDC. The objective is the absolute difference between the observed volume and the simulated volume. For the second objective the discharge of the winter months December, January and February are extracted from the time series. The objective is then the Nash-Sutcliffe efficiency NSE of the observed winter discharge against the simulated winter discharge. The observed data  $Q_{obs}$  are global in this module. To calibrate this objective you need a multi-objective optimizer like PA-DDS.

#### Parameters

|    |   |  |
|----|---|--|
| in | <code>real(dp), dimension(:) :: parameterset</code> |  |
| in | <code>procedure(eval_interface) :: eval</code>      |  |

#### Returns

`real(dp), dimension(2) :: multi_objective_ae_fdc_lsv_nse_djf` — objective function value (which will be e.g. minimized by an optimization routine like PA-DDS)

#### Authors

Juliane Mai

#### Date

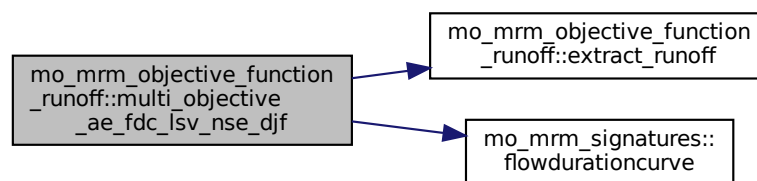
Feb 2016

Definition at line 1834 of file `mo_mrm_objective_function_runoff.f90`.

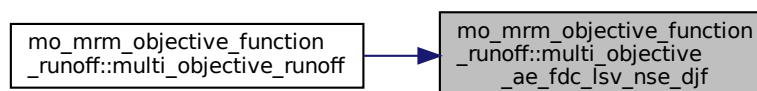
References `mo_common_mhm_mrm_variables::evalper`, `extract_runoff()`, `mo_mrm_signatures::flowdurationcurve()`, `mo_mrm_global_variables::gauge`, and `mo_mrm_global_variables::nmeasperday`.

Referenced by `multi_objective_runoff()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.6 multi\_objective\_Innse\_highflow\_Innse\_lowflow()

```
real(dp) function, dimension(2) mo_mrm_objective_function_runoff::multi_objective_innse_↵
highflow_innse_lowflow (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
```

Multi-objective function with NSE and lnNSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. A timepoint  $t$  of the observed data is marked as a lowflow timepoint  $t_{low}$  if

$$Q_{obs}(t) < \min(Q_{obs}) + 0.05 * (\max(Q_{obs}) - \min(Q_{obs}))$$

and a timepoint  $t$  of the observed data is marked as a highflow timepoint  $t_{high}$  if

$$t_{high} \text{ if } Q_{obs}(i) > \text{percentile}(Q_{obs}, 95.)$$

This timepoint identification is only performed for the observed data. The first objective is the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE_{high}$  of discharge values at high-flow timepoints

$$lnNSE_{high} = 1 - \frac{\sum_{i=1}^{N_{high}} (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

. The second objective is the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE_{low}$  of discharge values at low-flow timepoints

$$lnNSE_{low} = 1 - \frac{\sum_{i=1}^{N_{low}} (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

. Both objectives are returned. The observed data  $Q_{obs}$  are global in this module. To calibrate this objective you need a multi-objective optimizer like PA-DDS.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

real(dp), dimension(2) :: multi\_objective\_Innse\_highflow\_Innse\_lowflow — objective function value (which will be e.g. minimized by an optimization routine like PA-DDS)

#### Authors

Juliane Mai

#### Date

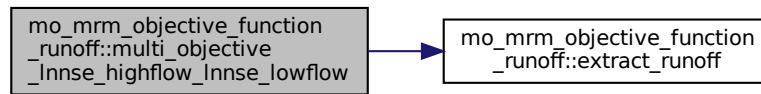
Oct 2015

Definition at line 1558 of file mo\_mrm\_objective\_function\_runoff.f90.

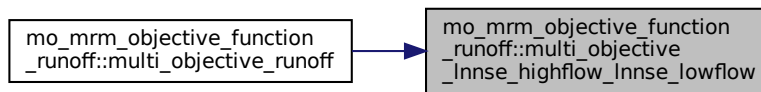
References extract\_runoff().

Referenced by multi\_objective\_runoff().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.7 multi\_objective\_lnnse\_highflow\_lnnse\_lowflow\_2()

```

real(dp) function, dimension(2) mo_mrm_objective_function_runoff::multi_objective_lnnse_↵
highflow_lnnse_lowflow_2 (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Multi-objective function with NSE and lnNSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. A timepoint  $t$  of the observed data is marked as a lowflow timepoint  $t_{low}$  if

$$Q_{obs}(t) < \min(Q_{obs}) + 0.05 * (\max(Q_{obs}) - \min(Q_{obs}))$$

and all other timepoints are marked as a highflow timepoints  $t_{high}$ . This timepoint identification is only performed for the observed data. The first objective is the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE_{high}$  of discharge values at high-flow timepoints

$$lnNSE_{high} = 1 - \frac{\sum_{i=1}^{N_{high}} (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

. The second objective is the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE_{low}$  of discharge values at low-flow timepoints

$$lnNSE_{low} = 1 - \frac{\sum_{i=1}^{N_{low}} (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

. Both objectives are returned. The observed data  $Q_{obs}$  are global in this module. To calibrate this objective you need a multi-objective optimizer like PA-DDS.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |



**Returns**

real(dp), dimension(2) :: multi\_objective\_lnnse\_highflow\_lnnse\_lowflow\_2 — objective function value (which will be e.g. minimized by an optimization routine like PA-DDS)

**Authors**

Juliane Mai

**Date**

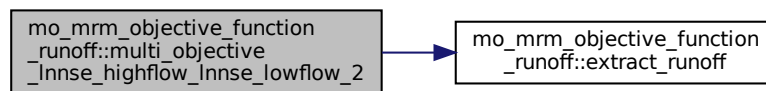
Oct 2015

Definition at line 1702 of file mo\_mrm\_objective\_function\_runoff.f90.

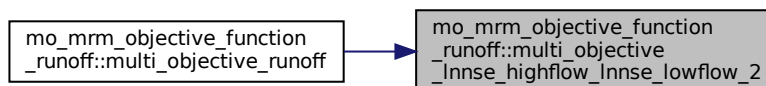
References extract\_runoff().

Referenced by multi\_objective\_runoff().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.43.2.8 multi\_objective\_nse\_lnnse()**

```

real(dp) function, dimension(2) mo_mrm_objective_function_runoff::multi_objective_nse_lnnse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Multi-objective function with NSE and lnNSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Nash-Sutcliffe model efficiency coefficient *NSE*

$$NSE = 1 - \frac{\sum_{i=1}^N (Q_{obs}(i) - Q_{model}(i))^2}{\sum_{i=1}^N (Q_{obs}(i) - \bar{Q}_{obs})^2}$$

and the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE$

$$lnNSE = 1 - \frac{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

are calculated and both returned. The observed data  $Q_{obs}$  are global in this module. To calibrate this objective you need a multi-objective optimizer like PA-DDS.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

*real(dp), dimension(2) :: multi\_objective\_nse\_lnnse* — objective function value (which will be e.g. minimized by an optimization routine like PA-DDS)

#### Authors

Juliane Mai

#### Date

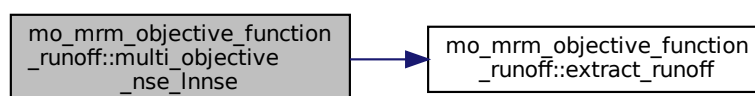
Oct 2015

Definition at line 1453 of file `mo_mrm_objective_function_runoff.f90`.

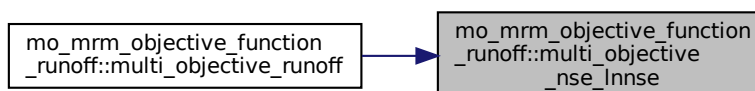
References `extract_runoff()`.

Referenced by `multi_objective_runoff()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.9 multi\_objective\_runoff()

```
subroutine, public mo_mrm_objective_function_runoff::multi_objective_runoff (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), dimension(:), intent(out), allocatable multi_objectives )
```

Wrapper for multi-objective functions where at least one is regarding runoff.

The functions selects the objective function case defined in a namelist, i.e. the global variable *opti\_function*. It return the multiple objective function values for a specific parameter set.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>REAL(dp), DIMENSION(:) :: parameterset</i>     |  |
| in  | <i>procedure(eval_interface) :: eval</i>          |  |
| out | <i>REAL(dp), DIMENSION(:) :: multi_objectives</i> |  |

#### Authors

Juliane Mai

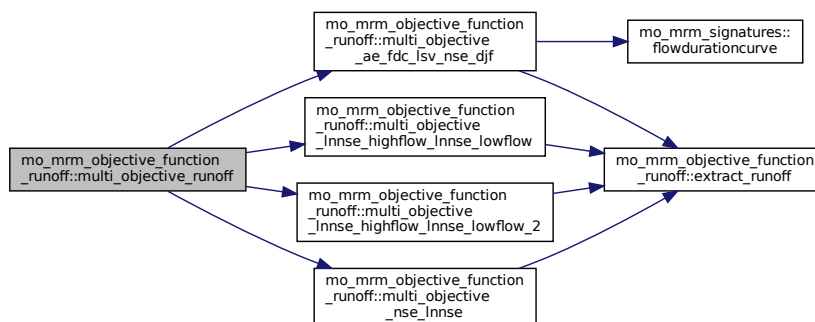
#### Date

Oct 2015

Definition at line 516 of file mo\_mrm\_objective\_function\_runoff.f90.

References `multi_objective_ae_fdc_lsv_nse_djf()`, `multi_objective_lnnse_highflow_lnnse_lowflow()`, `multi_objective_lnnse_highflow_lnnse_lowflow_2()`, `multi_objective_nse_lnnse()`, and `mo_common_mhm_mrm_variables::opti_function`.

Here is the call graph for this function:



### 17.43.2.10 objective\_equal\_nse\_lnnse()

```
real(dp) function mo_mrm_objective_function_runoff::objective_equal_nse_lnnse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
```

Objective function equally weighting NSE and lnNSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Nash-Sutcliffe model efficiency coefficient  $NSE$

$$NSE = 1 - \frac{\sum_{i=1}^N (Q_{obs}(i) - Q_{model}(i))^2}{\sum_{i=1}^N (Q_{obs}(i) - \bar{Q}_{obs})^2}$$

and the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE$

$$lnNSE = 1 - \frac{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

are calculated and added up equally weighted:

$$obj\_value = \frac{1}{2}(NSE + lnNSE)$$

The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <code>real(dp), dimension(:) :: parameterset</code> |  |
| in | <code>procedure(eval_interface) :: eval</code>      |  |

#### Returns

`real(dp) :: objective_equal_nse_lnnse` — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Juliane Mai

#### Date

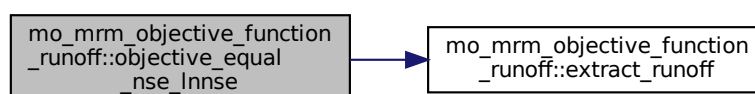
May 2013

Definition at line 1356 of file `mo_mrm_objective_function_runoff.f90`.

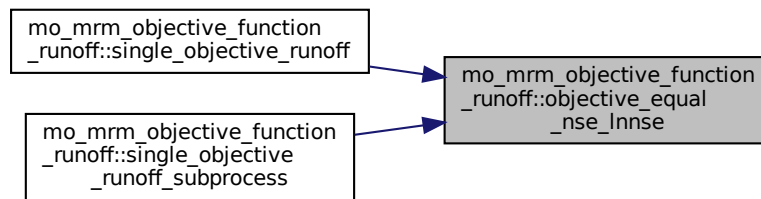
References `extract_runoff()`.

Referenced by `single_objective_runoff()`, and `single_objective_runoff_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.11 objective\_kge()

```

real(dp) function mo_mrm_objective_function_runoff::objective_kge (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of KGE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Kling-Gupta model efficiency coefficient  $KGE$

$$KGE = 1.0 - \sqrt{((1-r)^2 + (1-\alpha)^2 + (1-\beta)^2)}$$

where  $r$  = Pearson product-moment correlation coefficient  $\alpha$  = ratio of simulated mean to observed mean  $\beta$  = ratio of simulated standard deviation to observed standard deviation is calculated and the objective function is

$$obj\_value = 1.0 - KGE$$

$(1 - KGE)$  is the objective since we always apply minimization methods. The minimal value of  $(1 - KGE)$  is 0 for the optimal KGE of 1.0. The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

*real(dp) :: objective\_kge* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Rohini Kumar

## Date

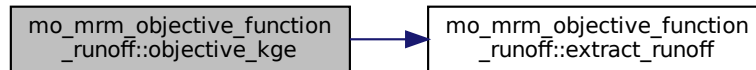
August 2014

Definition at line 2094 of file `mo_mrm_objective_function_runoff.f90`.

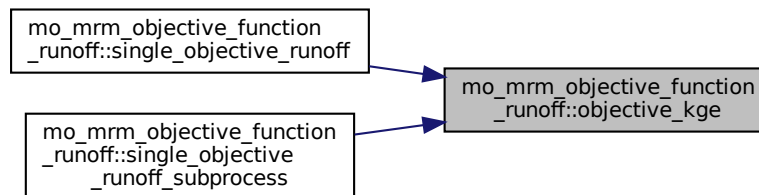
References `extract_runoff()`.

Referenced by `single_objective_runoff()`, and `single_objective_runoff_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.12 objective\_lnnse()

```

real(dp) function mo_mrm_objective_function_runoff::objective_lnnse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of logarithmic NSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the logarithmic Nash-Sutcliffe model efficiency coefficient  $\ln NSE$

$$\ln NSE = 1 - \frac{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

is calculated.

$$obj\_value = \ln NSE$$

The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <code>real(dp), dimension(:) :: parameterset</code> |  |
| in | <code>procedure(eval_interface) :: eval</code>      |  |

**Returns**

real(dp) :: objective\_Innse — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Juliane Mai

**Date**

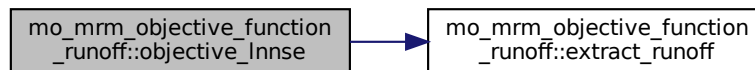
May 2013

Definition at line 1084 of file mo\_mrm\_objective\_function\_runoff.f90.

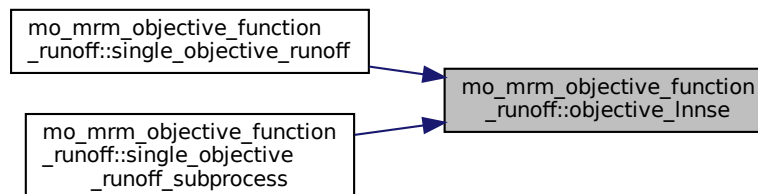
References extract\_runoff().

Referenced by single\_objective\_runoff(), and single\_objective\_runoff\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.43.2.13 objective\_multiple\_gauges\_kge\_power6()**

```

real(dp) function mo_mrm_objective_function_runoff::objective_multiple_gauges_kge_power6 (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

combined objective function based on KGE raised to the power 6

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Kling-Gupta model efficiency coefficient *KGE* for a given gauging station

$$KGE = 1.0 - \sqrt{((1-r)^2 + (1-\alpha)^2 + (1-\beta)^2)}$$

where  $r$  = Pearson product-moment correlation coefficient  $\alpha$  = ratio of simulated mean to observed mean  $\beta$  = ratio of simulated standard deviation to observed standard deviation is calculated and the objective function for a given gauging station ( $i$ ) is

$$\phi_i = 1.0 - KGE_i$$

$\phi_i$  is the objective since we always apply minimization methods. The minimal value of  $\phi_i$  is 0 for the optimal KGE of 1.0. Finally, the overall  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all gauging stations ( $N$ ).

$$OF = \sqrt[6]{\sum ((1.0 - KGE_i)/N)^6}$$

. The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

*real(dp) :: objective\_multiple\_gauges\_kge\_power6* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Rohini Kumar

#### Date

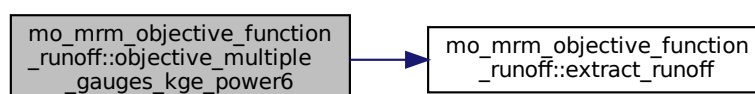
March 2015

Definition at line 2196 of file `mo_mrm_objective_function_runoff.f90`.

References `extract_runoff()`.

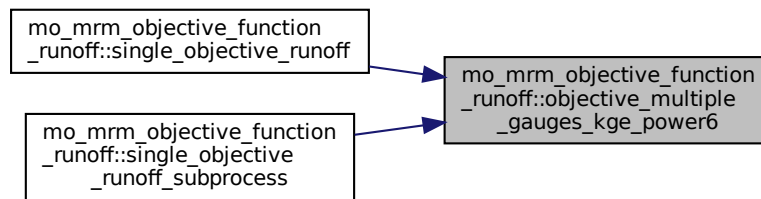
Referenced by `single_objective_runoff()`, and `single_objective_runoff_subprocess()`.

Here is the call graph for this function:





Here is the caller graph for this function:



#### 17.43.2.14 objective\_nse()

```

real(dp) function mo_mrm_objective_function_runoff::objective_nse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of NSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Nash-Sutcliffe model efficiency coefficient  $NSE$

$$NSE = 1 - \frac{\sum_{i=1}^N (Q_{obs}(i) - Q_{model}(i))^2}{\sum_{i=1}^N (Q_{obs}(i) - \bar{Q}_{obs})^2}$$

is calculated and the objective function is

$$obj\_value = 1 - NSE$$

The observed data  $Q_{obs}$  are global in this module.

##### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

##### Returns

real(dp) :: objective\_nse — objective function value (which will be e.g. minimized by an optimization routine like DDS)

##### Authors

Juliane Mai

##### Date

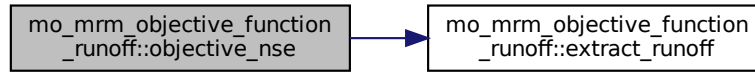
May 2013

Definition at line 1263 of file mo\_mrm\_objective\_function\_runoff.f90.

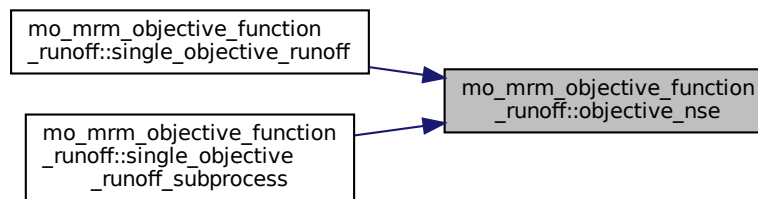
References extract\_runoff().

Referenced by single\_objective\_runoff(), and single\_objective\_runoff\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.15 objective\_power6\_nse\_lnnse()

```

real(dp) function mo_mrm_objective_function_runoff::objective_power6_nse_lnnse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of combined NSE and lnNSE with power of 5 i.e. the p-norm with p=5.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Nash-Sutcliffe model efficiency coefficient  $NSE$

$$NSE = 1 - \frac{\sum_{i=1}^N (Q_{obs}(i) - Q_{model}(i))^2}{\sum_{i=1}^N (Q_{obs}(i) - \bar{Q}_{obs})^2}$$

and the logarithmic Nash-Sutcliffe model efficiency coefficient  $lnNSE$

$$lnNSE = 1 - \frac{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln Q_{model}(i))^2}{\sum_{i=1}^N (\ln Q_{obs}(i) - \ln \bar{Q}_{obs})^2}$$

are calculated and added up equally weighted:

$$obj\_value = \sqrt[6]{(1 - NSE)^6 + (1 - lnNSE)^6}$$

The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

**Returns**

real(dp) :: objective\_power6\_nse\_lnnse — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Juliane Mai and Matthias Cuntz

**Date**

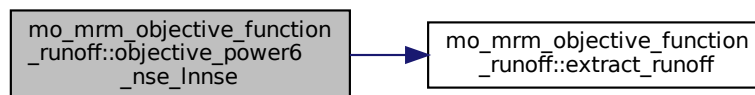
March 2014

Definition at line 1994 of file mo\_mrm\_objective\_function\_runoff.f90.

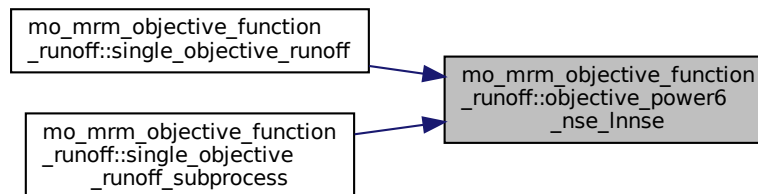
References extract\_runoff().

Referenced by single\_objective\_runoff(), and single\_objective\_runoff\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.43.2.16 objective\_sse()**

```

real(dp) function mo_mrm_objective_function_runoff::objective_sse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of SSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the sum squared errors

$$SSE = \sum_{i=1}^N (Q_{obs}(i) - Q_{model}(i))^2$$

is calculated and the objective function is

$$obj\_value = SSE$$

The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

`real(dp) :: objective_sse` — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Juliane Mai and Matthias Cuntz

#### Date

March 2014

Definition at line 1173 of file `mo_mrm_objective_function_runoff.f90`.

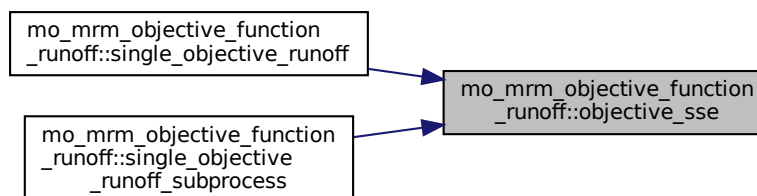
References `extract_runoff()`.

Referenced by `single_objective_runoff()`, and `single_objective_runoff_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



**17.43.2.17 objective\_sse\_boxcox()**

```
real(dp) function mo_mrm_objective_function_runoff::objective_sse_boxcox (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
```

Objective function of sum of squared errors of transformed streamflow.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the sum of squared error  $tSSE$

$$tSSE = \sum_{i=1}^N (z(Q_{obs}(i), \lambda) - z(Q_{model}(i), \lambda))^2$$

is calculated where  $z$  is the transform and given by

$$z(x, \lambda) = \frac{x^\lambda - 1}{\lambda}$$

for  $\lambda$  unequal to zero and

$$z(x, \lambda) = \log x$$

for  $\lambda$  equal to zero. The objective function is

$$obj\_value = tSSE$$

The observed data  $Q_{obs}$  are global in this module.

The boxcox transformation uses a parameter of 0.2, suggested by Woldemeskel et al. Hydrol Earth Syst Sci, 2018 vol. 22 (12) pp. 6257-6278. "Evaluating post-processing approaches for monthly and seasonal streamflow forecasts." <https://www.hydrol-earth-syst-sci.net/22/6257/2018/>

**Parameters**

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

**Returns**

real(dp) :: objective\_sse\_boxcox — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Stephan Thober, Dmitri Kavetski

## Date

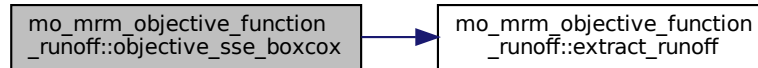
Aug 2019

Definition at line 2377 of file mo\_mrm\_objective\_function\_runoff.f90.

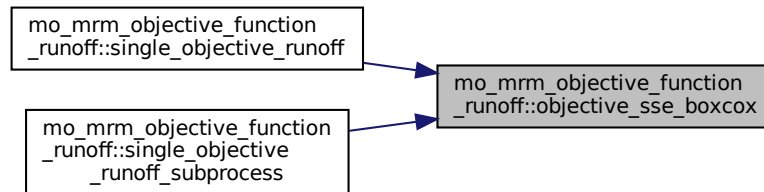
References extract\_runoff().

Referenced by single\_objective\_runoff(), and single\_objective\_runoff\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.43.2.18 objective\_weighted\_nse()

```

real(dp) function mo_mrm_objective_function_runoff::objective_weighted_nse (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of weighted NSE.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the weighted Nash-Sutcliffe model efficiency coefficient  $NSE$

$$wNSE = 1 - \frac{\sum_{i=1}^N Q_{obs}(i) * (Q_{obs}(i) - Q_{model}(i))^2}{\sum_{i=1}^N Q_{obs}(i) * (Q_{obs}(i) - \bar{Q}_{obs})^2}$$

is calculated and the objective function is

$$obj\_value = 1 - wNSE$$

The observed data  $Q_{obs}$  are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

**Returns**

real(dp) :: objective\_weighted\_nse — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Stephan Thober, Bjoern Guse

**Date**

May 2018

Definition at line 2284 of file mo\_mrm\_objective\_function\_runoff.f90.

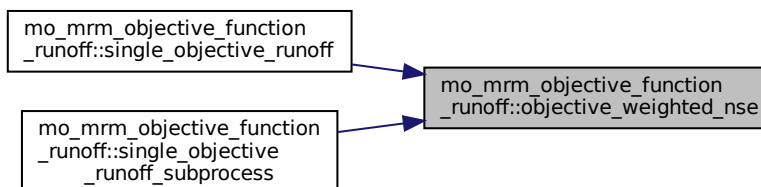
References extract\_runoff().

Referenced by single\_objective\_runoff(), and single\_objective\_runoff\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.43.2.19 parameter\_regularization()**

```

real(dp) function mo_mrm_objective_function_runoff::parameter_regularization (
    real(dp), dimension(:), intent(in) paraset,
    real(dp), dimension(size(paraset)), intent(in) prior,
    real(dp), dimension(size(paraset), 2), intent(in) bounds,
    logical, dimension(size(paraset)), intent(in) mask )
  
```

TODO: add description.

TODO: add description

## Parameters

|    |  |            |
|----|--|------------|
| in | <i>real(dp), dimension(:) :: paraset</i>               |            |
| in | <i>real(dp), dimension(size(paraset)) :: prior</i>     |            |
| in | <i>real(dp), dimension(size(paraset), 2) :: bounds</i> | (min, max) |
| in | <i>logical, dimension(size(paraset)) :: mask</i>       |            |

## Authors

Robert Schweppe

## Date

Jun 2018

Definition at line 867 of file `mo_mrm_objective_function_runoff.f90`.

Referenced by `loglikelihood_evin2013_2()`.

Here is the caller graph for this function:

**17.43.2.20 single\_objective\_runoff()**

```

real(dp) function, public mo_mrm_objective_function_runoff::single_objective_runoff (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in), optional arg1,
    real(dp), intent(out), optional arg2,
    real(dp), intent(out), optional arg3 )
  
```

Wrapper for objective functions optimizing against runoff.

The function selects the objective function case defined in a namelist, i.e. the global variable `opti_function`. It returns the objective function value for a specific parameter set.

## Parameters

|     |   |  |
|-----|---|--|
| in  | <i>REAL(dp), DIMENSION(:) :: parameterset</i> |  |
| in  | <i>procedure(eval_interface) :: eval</i>      |  |
| in  | <i>real(dp), optional :: arg1</i>             |  |
| out | <i>real(dp), optional :: arg2</i>             |  |
| out | <i>real(dp), optional :: arg3</i>             |  |



**Returns**

real(dp) :: objective — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Juliane Mai

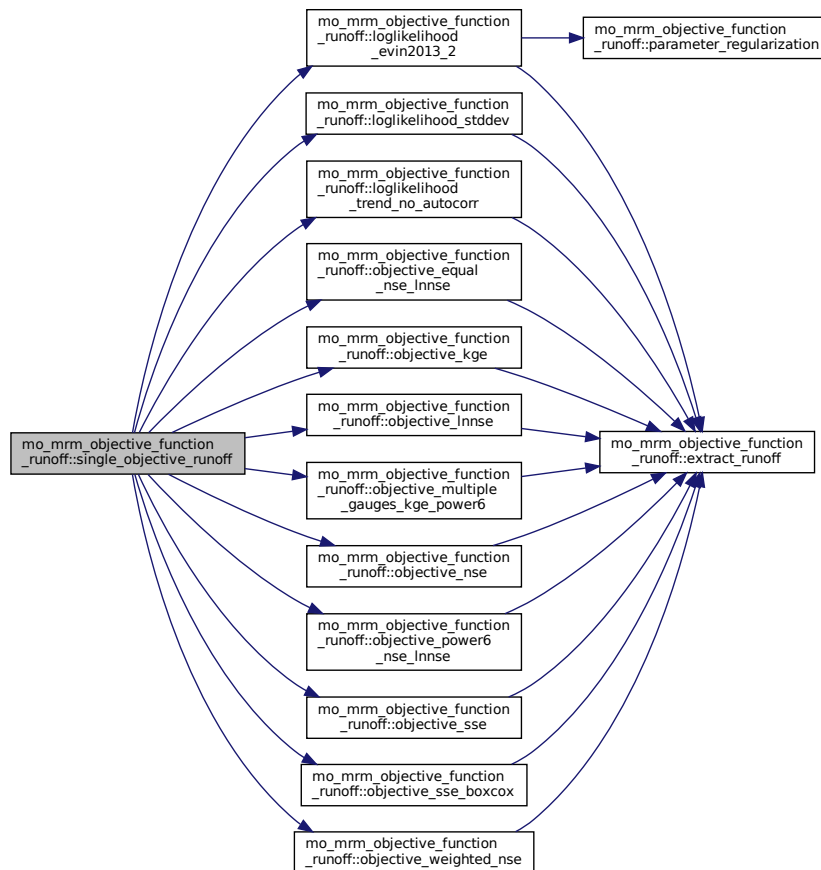
**Date**

Dec 2012

Definition at line 116 of file mo\_mrm\_objective\_function\_runoff.f90.

References `loglikelihood_evin2013_2()`, `loglikelihood_stddev()`, `loglikelihood_trend_no_autocorr()`, `objective_equal_nse_lnnse()`, `objective_kge()`, `objective_lnnse()`, `objective_multiple_gauges_kge_power6()`, `objective_nse()`, `objective_power6_nse_lnnse()`, `objective_sse()`, `objective_sse_boxcox()`, `objective_weighted_nse()`, `mo_common_mhm_mrm_variables::opti_function`, and `mo_common_mhm_mrm_variables::opti_method`.

Here is the call graph for this function:

**17.43.2.21 single\_objective\_runoff\_master()**

```

real(dp) function, public mo_mrm_objective_function_runoff::single_objective_runoff_master (
    real(dp), dimension(:), intent(in) parameterset,

```

```

procedure(eval_interface), intent(in), pointer eval,
real(dp), intent(in), optional arg1,
real(dp), intent(out), optional arg2,
real(dp), intent(out), optional arg3 )

```

Wrapper for objective functions optimizing against runoff.

The functions selects the objective function case defined in a namelist, i.e. the global variable *opti\_function*. It return the objective function value for a specific parameter set.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>REAL(dp), DIMENSION(:) :: parameterset</i> |  |
| in  | <i>procedure(eval_interface) :: eval</i>      |  |
| in  | <i>real(dp), optional :: arg1</i>             |  |
| out | <i>real(dp), optional :: arg2</i>             |  |
| out | <i>real(dp), optional :: arg3</i>             |  |

#### Returns

real(dp) :: objective — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Juliane Mai

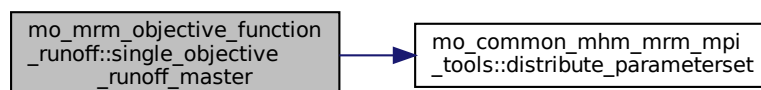
#### Date

Dec 2012

Definition at line 233 of file mo\_mrm\_objective\_function\_runoff.f90.

References mo\_common\_mhm\_mrm\_mpi\_tools::distribute\_parameterset(), mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::ngaugestotal, mo\_common\_mhm\_mrm\_variables::opti\_function, and mo\_common\_mhm\_mrm\_variables::opti\_method.

Here is the call graph for this function:



#### 17.43.2.22 single\_objective\_runoff\_subprocess()

```

subroutine, public mo_mrm_objective_function_runoff::single_objective_runoff_subprocess (
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in), optional arg1,

```

```

real(dp), intent(out), optional arg2,
real(dp), intent(out), optional arg3 )

```

Wrapper for objective functions optimizing against runoff.

The function selects the objective function case defined in a namelist, i.e. the global variable *opti\_function*. It returns the objective function value for a specific parameter set.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>REAL(dp), DIMENSION(:) :: parameterset</i> |  |
| in  | <i>procedure(eval_interface) :: eval</i>      |  |
| in  | <i>real(dp), optional :: arg1</i>             |  |
| out | <i>real(dp), optional :: arg2</i>             |  |
| out | <i>real(dp), optional :: arg3</i>             |  |

#### Returns

*real(dp) :: objective* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Juliane Mai

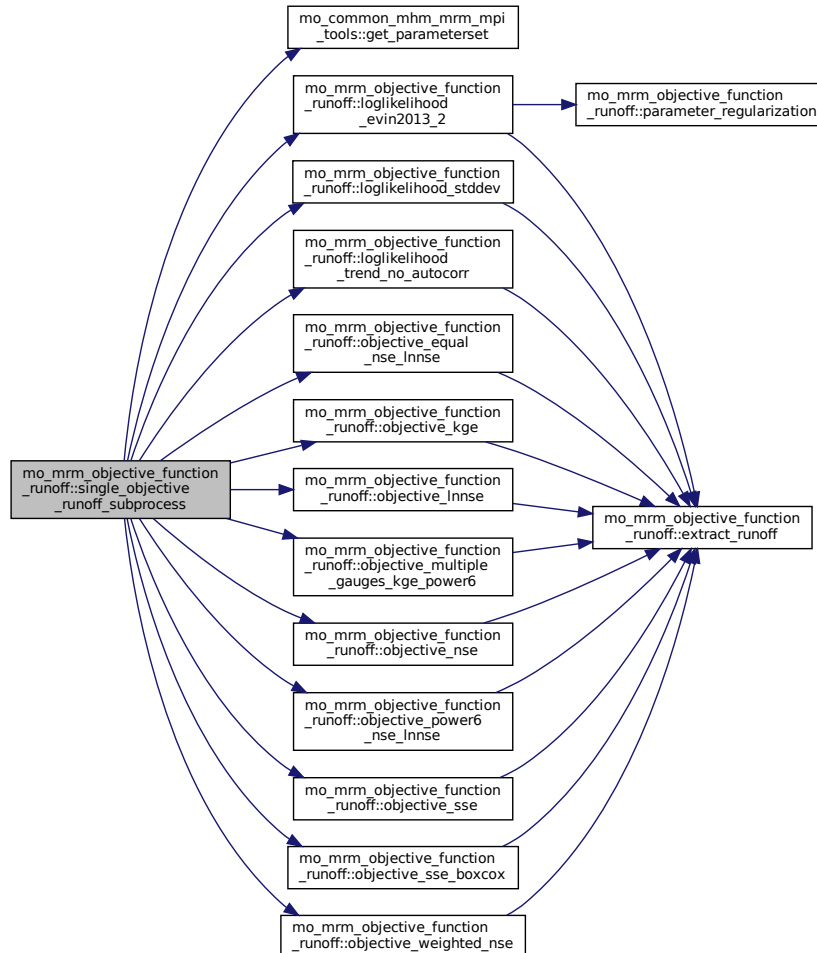
#### Date

Dec 2012

Definition at line 381 of file mo\_mrm\_objective\_function\_runoff.f90.

References mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_mpi\_tools::get\_parameterset(), loglikelihood\_evin2013\_2(), loglikelihood\_stddev(), loglikelihood\_trend\_no\_autocorr(), objective\_equal\_nse\_↔\_lnnse(), objective\_kge(), objective\_lnnse(), objective\_multiple\_gauges\_kge\_power6(), objective\_nse(), objective\_↔\_power6\_nse\_lnnse(), objective\_sse(), objective\_sse\_boxcox(), objective\_weighted\_nse(), mo\_common\_mhm\_↔\_mrm\_variables::opti\_function, and mo\_common\_mhm\_mrm\_variables::opti\_method.

Here is the call graph for this function:



## 17.44 mo\_mrm\_pre\_routing Module Reference

Performs pre-processing for routing for mHM at level L11.

### Functions/Subroutines

- subroutine, public [l11\\_runoff\\_acc](#) (qAll, efecArea, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, TS, map\_flag, qAcc)  
*total runoff accumulation at L11.*
- subroutine, public [add\\_inflow](#) (nInflowGauges, InflowIndexList, InflowHeadwater, InflowNodeList, QInflow, q↔  
Out)  
*Adds inflow discharge to the runoff produced at the cell where the inflow is occurring.*
- subroutine [l11\\_e\\_acc](#) (qAll, efecArea, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, TS, map\_flag, qAcc)  
*temperature energy accumulation at L11.*
- subroutine, public [calc\\_l1\\_runoff\\_e](#) (fSealed\_area\_fraction, fast\_interflow, slow\_interflow, baseflow, direct\_↔  
runoff, temp\_air, mean\_temp\_air, lateral\_E)  
*calculate lateral temperature energy from runoff components.*

- subroutine, public [l11\\_meteo\\_acc](#) (meteo\_all, efecArea, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, map\_flag, meteo\_acc)

*meteo forcing accumulation at L11 for temperature routing.*

### 17.44.1 Detailed Description

Performs pre-processing for routing for mHM at level L11.

This module performs runoff accumulation from L1 to L11 and inflow summation.

#### Authors

Luis Samaniego

#### Date

Dec 2012

### 17.44.2 Function/Subroutine Documentation

#### 17.44.2.1 add\_inflow()

```
subroutine, public mo_mrm_pre_routing::add_inflow (
    integer(i4), intent(in) nInflowGauges,
    integer(i4), dimension(:), intent(in) InflowIndexList,
    logical, dimension(:), intent(in) InflowHeadwater,
    integer(i4), dimension(:), intent(in) InflowNodeList,
    real(dp), dimension(:), intent(in) QInflow,
    real(dp), dimension(:), intent(inout) qOut )
```

Adds inflow discharge to the runoff produced at the cell where the inflow is occurring.

If an inflow gauge is given, then this routine is adding the values to the runoff produced at the grid cell where the inflow is happening. The values are not directly added to the river network. If this cell is not a headwater then the streamflow produced upstream will be neglected.

#### Parameters

|         |   |  |
|---------|---|--|
| in      | <i>integer(i4) :: nInflowGauges</i>                 | [-] number of inflow points                        |
| in      | <i>integer(i4), dimension(:) :: InflowIndexList</i> | [-] index of inflow points                         |
| in      | <i>logical, dimension(:) :: InflowHeadwater</i>     | [-] if to consider headwater cells of inflow gauge |
| in      | <i>integer(i4), dimension(:) :: InflowNodeList</i>  | [-] L11 ID of inflow points                        |
| in      | <i>real(dp), dimension(:) :: QInflow</i>            | [m3 s-1] inflowing water                           |
| in, out | <i>real(dp), dimension(:) :: qOut</i>               | [m3 s-1] Series of attenuated runoff               |

#### Authors

Stephan Thober & Matthias Zink

## Date

Jul 2016

Definition at line 177 of file mo\_mrm\_pre\_routing.f90.

Referenced by mo\_mrm\_routing::mrm\_routing().

Here is the caller graph for this function:



### 17.44.2.2 calc\_l1\_runoff\_e()

```

subroutine, public mo_mrm_pre_routing::calc_l1_runoff_e (
    real(dp), dimension(:), intent(in) fSealed_area_fraction,
    real(dp), dimension(:), intent(in) fast_interflow,
    real(dp), dimension(:), intent(in) slow_interflow,
    real(dp), dimension(:), intent(in) baseflow,
    real(dp), dimension(:), intent(in) direct_runoff,
    real(dp), dimension(:), intent(in) temp_air,
    real(dp), dimension(:), intent(in) mean_temp_air,
    real(dp), dimension(:), intent(inout) lateral_E )
  
```

calculate lateral temperature energy from runoff components.

calculate lateral temperature energy from runoff components.

#### Parameters

|     |  |   |
|-----|--|---|
| in  | <i>REAL(dp) :: fSealed_area_fraction</i> | sealed area fraction [1]                            |
| in  | <i>REAL(dp) :: fast_interflow</i>        | $q_0$ Fast runoff component [mm TS-1]               |
| in  | <i>REAL(dp) :: slow_interflow</i>        | $q_1$ Slow runoff component [mm TS-1]               |
| in  | <i>REAL(dp) :: baseflow</i>              | $q_2$ Baseflow [mm TS-1]                            |
| in  | <i>REAL(dp) :: direct_runoff</i>         | $q_D$ Direct runoff from impervious areas [mm TS-1] |
| in  | <i>REAL(dp) :: temp_air</i>              | air temperature [K]                                 |
| in  | <i>REAL(dp) :: mean_temp_air</i>         | annual mean air temperature [K]                     |
| out | <i>REAL(dp) :: lateral_E</i>             | $E_T$ Generated runoff [K mm TS-1]                  |

#### Authors

Sebastian Mueller

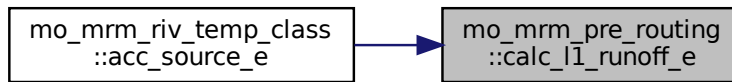
#### Date

Jun 2020

Definition at line 346 of file mo\_mrm\_pre\_routing.f90.

Referenced by mo\_mrm\_riv\_temp\_class::acc\_source\_e().

Here is the caller graph for this function:



### 17.44.2.3 l11\_e\_acc()

```

subroutine mo_mrm_pre_routing::l11_e_acc (
    real(dp), dimension(:), intent(in) qAll,
    real(dp), dimension(:), intent(in) efecArea,
    integer(i4), dimension(:), intent(in) L1_L11_Id,
    real(dp), dimension(:), intent(in) L11_areaCell,
    integer(i4), dimension(:), intent(in) L11_L1_Id,
    integer(i4), intent(in) TS,
    logical, intent(in) map_flag,
    real(dp), dimension(:), intent(out) qAcc )
  
```

temperature energy accumulation at L11.

Upscales energy in space from L1 to L11 if routing resolution is higher than hydrology resolution (`map_flag` equals `.true.`) or downscales runoff from L1 to L11 if routing resolution is lower than hydrology resolution.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>real(dp), dimension(:) :: qall</i>         | total runoff L1 [mm K TS-1]  |
| in  | <i>real(dp), dimension(:) :: efecarea</i>     | effective area in [km2] at Level 1                                       |
| in  | <i>integer(i4), dimension(:) :: L1_L11_Id</i> | L11 lds mapped on L1   |
| in  | <i>real(dp), dimension(:) :: L11_areacell</i> | effective area in [km2] at Level 11                                      |
| in  | <i>integer(i4), dimension(:) :: L11_L1_Id</i> | L1 lds mapped on L11   |
| in  | <i>integer(i4) :: TS</i>                      | time step in [s]   |
| in  | <i>logical :: map_flag</i>                    | Flag indicating whether routing resolution is higher than hydrologic one |
| out | <i>real(dp), dimension(:) :: qAcc</i>         | aggregated runoff at L11 [m3 K s-1]                                      |

#### Authors

Luis Samaniego

#### Date

Jan 2013

Definition at line 253 of file `mo_mrm_pre_routing.f90`.

References `mo_common_constants::nodata_dp`.

### 17.44.2.4 l11\_meteo\_acc()

```
subroutine, public mo_mrm_pre_routing::l11_meteo_acc (
    real(dp), dimension(:), intent(in) meteo_all,
    real(dp), dimension(:), intent(in) efecArea,
    integer(i4), dimension(:), intent(in) L1_L11_Id,
    real(dp), dimension(:), intent(in) L11_areaCell,
    integer(i4), dimension(:), intent(in) L11_L1_Id,
    logical, intent(in) map_flag,
    real(dp), dimension(:), intent(out) meteo_acc )
```

meteo forcing accumulation at L11 for temperature routing.

Upscales meteo forcing in space from L1 to L11 if routing resolution is higher than hydrology resolution (`map_flag` equals `.true.`) or downscales meteo forcing from L1 to L11 if routing resolution is lower than hydrology resolution.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>real(dp), dimension(:) :: meteo_all</i>    | meteo forcing  |
| in  | <i>real(dp), dimension(:) :: efecarea</i>     | effective area in [km2] at Level 1                                       |
| in  | <i>integer(i4), dimension(:) :: L1_L11_Id</i> | L11 lds mapped on L1   |
| in  | <i>real(dp), dimension(:) :: L11_areacell</i> | effective area in [km2] at Level 11                                      |
| in  | <i>integer(i4), dimension(:) :: L11_L1_Id</i> | L1 lds mapped on L11   |
| in  | <i>logical :: map_flag</i>                    | Flag indicating whether routing resolution is higher than hydrologic one |
| out | <i>real(dp), dimension(:) :: meteo_acc</i>    | aggregated meteo forcing   |

#### Authors

Sebastian Mueller

#### Date

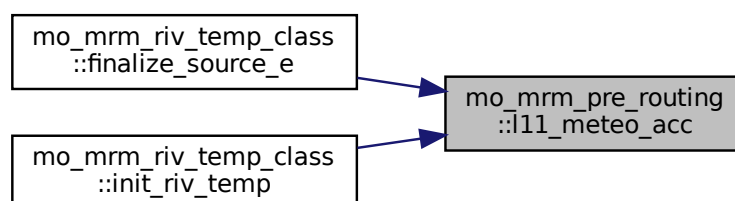
Jul 2020

Definition at line 420 of file `mo_mrm_pre_routing.f90`.

References `mo_common_constants::nodata_dp`.

Referenced by `mo_mrm_riv_temp_class::finalize_source_e()`, and `mo_mrm_riv_temp_class::init_riv_temp()`.

Here is the caller graph for this function:





### 17.44.2.5 l11\_runoff\_acc()

```
subroutine, public mo_mrm_pre_routing::l11_runoff_acc (
    real(dp), dimension(:), intent(in) qAll,
    real(dp), dimension(:), intent(in) efecArea,
    integer(i4), dimension(:), intent(in) L1_L11_Id,
    real(dp), dimension(:), intent(in) L11_areaCell,
    integer(i4), dimension(:), intent(in) L11_L1_Id,
    integer(i4), intent(in) TS,
    logical, intent(in) map_flag,
    real(dp), dimension(:), intent(out) qAcc )
```

total runoff accumulation at L11.

Upscales runoff in space from L1 to L11 if routing resolution is higher than hydrology resolution (map\_flag equals .true.) or downscales runoff from L1 to L11 if routing resolution is lower than hydrology resolution.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>real(dp), dimension(:) :: qall</i>         | total runoff L1 [mm TS-1]  |
| in  | <i>real(dp), dimension(:) :: efecarea</i>     | effective area in [km2] at Level 1                                       |
| in  | <i>integer(i4), dimension(:) :: L1_L11_Id</i> | L11 lds mapped on L1   |
| in  | <i>real(dp), dimension(:) :: L11_areacell</i> | effective area in [km2] at Level 11                                      |
| in  | <i>integer(i4), dimension(:) :: L11_L1_Id</i> | L1 lds mapped on L11   |
| in  | <i>integer(i4) :: TS</i>                      | time step in [s]   |
| in  | <i>logical :: map_flag</i>                    | Flag indicating whether routing resolution is higher than hydrologic one |
| out | <i>real(dp), dimension(:) :: qAcc</i>         | aggregated runoff at L11 [m3 s-1]  |

#### Authors

Luis Samaniego

#### Date

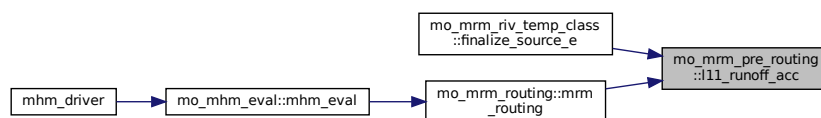
Jan 2013

Definition at line 75 of file mo\_mrm\_pre\_routing.f90.

References mo\_common\_constants::nodata\_dp.

Referenced by mo\_mrm\_riv\_temp\_class::finalize\_source\_e(), and mo\_mrm\_routing::mrm\_routing().

Here is the caller graph for this function:



## 17.45 mo\_mrm\_read\_config Module Reference

read mRM config

## Functions/Subroutines

- subroutine, public `mrm_read_config` (`file_namelist`, `unamelist`, `file_namelist_param`, `unamelist_param`, `do_message`, `readLatLon`)  
*Read the general config of mRM.*
- subroutine `read_mrm_routing_params` (`processCase`, `file_namelist_param`, `unamelist_param`)  
*TODO: add description.*

### 17.45.1 Detailed Description

read mRM config

This module contains all mRM subroutines related to reading the mRM configuration either from file or copy from mHM.

#### Authors

Stephan Thober

#### Date

Aug 2015

### 17.45.2 Function/Subroutine Documentation

#### 17.45.2.1 `mrm_read_config()`

```
subroutine, public mo_mrm_read_config::mrm_read_config (
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist,
    character(*), intent(in) file_namelist_param,
    integer, intent(in) unamelist_param,
    logical, intent(in) do_message,
    logical, intent(out) readLatLon )
```

Read the general config of mRM.

Depending on the variable `mrm_coupling_config`, the mRM config is either read from `mrm.nml` and parameters from `mrm_parameter.nml` or copied from mHM.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <code>character(*) :: file_namelist, file_namelist_param</code> |  |
| in  | <code>integer :: unamelist, unamelist_param</code>              |  |
| in  | <code>character(*) :: file_namelist, file_namelist_param</code> |  |
| in  | <code>integer :: unamelist, unamelist_param</code>              |  |
| in  | <code>logical :: do_message</code>                              | - flag for writing mHM standard messages |
| out | <code>logical :: readLatLon</code>                              | - flag for reading LatLon file           |

#### Authors

Stephan Thober

## Date

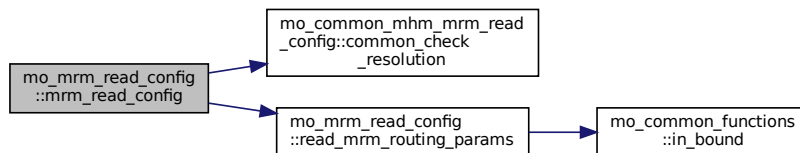
Aug 2015

Definition at line 56 of file mo\_mrm\_read\_config.f90.

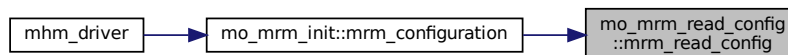
References mo\_common\_variables::alma\_convention, mo\_common\_mhm\_mrm\_read\_config::common\_check\_↔\_resolution(), mo\_mrm\_global\_variables::dirbankfullrunoff, mo\_mrm\_global\_variables::dirgauges, mo\_mrm\_↔\_global\_variables::dirtotalrunoff, mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_file::file\_defoutput, mo\_mrm\_global\_variables::filenametotalrunoff, mo\_mrm\_global\_variables::gauge, mo\_mrm\_global\_variables::gw\_coupling, mo\_mrm\_global\_variables::inflowgauge, mo\_mrm\_global\_variables\_↔::is\_start, mo\_common\_constants::maxnodomains, mo\_mrm\_constants::maxnogauges, mo\_mrm\_global\_↔\_variables::ngaugeslocal, mo\_mrm\_global\_variables::ngaugestotal, mo\_mrm\_global\_variables::ninflowgaugestotal, mo\_common\_constants::nodata\_i4, mo\_mrm\_global\_variables::output\_deflate\_level\_mrm, mo\_mrm\_global\_↔\_variables::output\_double\_precision\_mrm, mo\_mrm\_global\_variables::outputflxstate\_mrm, mo\_common\_↔\_variables::processmatrix, read\_mrm\_routing\_params(), mo\_mrm\_global\_variables::timestep\_model\_outputs\_mrm, mo\_mrm\_file::udefoutput, and mo\_mrm\_global\_variables::varnametotalrunoff.

Referenced by mo\_mrm\_init::mrm\_configuration().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.45.2.2 read\_mrm\_routing\_params()

```

subroutine mo_mrm_read_config::read_mrm_routing_params (
    integer(i4), intent(in) processCase,
    character(*), intent(in) file_namelist_param,
    integer(i4), intent(in) unamelist_param )
  
```

TODO: add description.

TODO: add description

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>integer(i4) :: processCase</i>          | it is the default case, should be one      |
| in | <i>character(*) :: file_namelist_param</i> | file name containing parameter namelist    |
| in | <i>integer(i4) :: unamelist_param</i>      | file name id containing parameter namelist |

**Authors**

Robert Schweppe

**Date**

Jun 2018

Definition at line 429 of file mo\_mrm\_read\_config.f90.

References mo\_common\_variables::global\_parameters, mo\_common\_variables::global\_parameters\_name, mo\_common\_functions::in\_bound(), mo\_common\_constants::ncolpars, and mo\_common\_variables::processmatrix.

Referenced by mrm\_read\_config().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.46 mo\_mrm\_read\_data Module Reference

This module contains all routines to read mRM data from file.

### Functions/Subroutines

- subroutine, public [mrm\\_read\\_l0\\_data](#) (do\_reinit, do\_readlatlon, do\_readlcover)  
*read L0 data from file*
- subroutine, public [mrm\\_read\\_discharge](#)  
*Read discharge timeseries from file.*
- subroutine, public [mrm\\_read\\_total\\_runoff](#) (iDomain)  
*read simulated runoff that is to be routed*
- subroutine, public [mrm\\_read\\_bankfull\\_runoff](#) (iDomain)
- subroutine [rotate\\_fdir\\_variable](#) (x)  
*TODO: add description.*

### 17.46.1 Detailed Description

This module contains all routines to read mRM data from file.

TODO: add description

**Authors**

Stephan Thober

**Date**

Aug 2015

**17.46.2 Function/Subroutine Documentation****17.46.2.1 mrm\_read\_bankfull\_runoff()**

```
subroutine, public mo_mrm_read_data::mrm_read_bankfull_runoff (
    integer(i4), intent(in) iDomain )
```

**Parameters**

|    |                |  |
|----|----------------|--|
| in | <i>idomain</i> |  |
|----|----------------|--|

**Note**

The file read in must contain a double precision float variable with the name "Q\_bkfl".

Definition at line 432 of file mo\_mrm\_read\_data.f90.

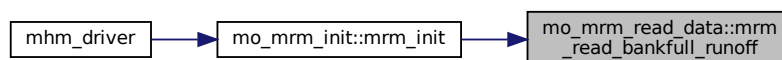
References mo\_mrm\_global\_variables::dirbankfullrunoff, mo\_mrm\_global\_variables::level11, and mo\_read\_nc::read\_const\_nc().

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.46.2.2 mrm\_read\_discharge()

```
subroutine, public mo_mrm_read_data::mrm_read_discharge
```

Read discharge timeseries from file.

Read Observed discharge at the outlet of a catchment and at the inflow of a catchment. Allocate global runoff variable that contains the simulated runoff after the simulation.

#### Authors

Matthias Zink & Stephan Thober

#### Date

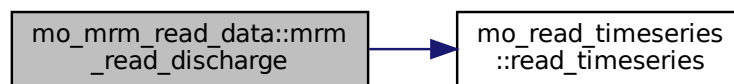
Aug 2015

Definition at line 256 of file mo\_mrm\_read\_data.f90.

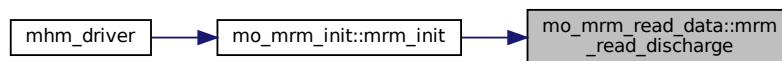
References mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_variables::evalper, mo\_mrm↔  
\_global\_variables::gauge, mo\_mrm\_global\_variables::inflowgauge, mo\_mrm\_global\_variables::mrm\_runoff,  
mo\_mrm\_global\_variables::ngaugeslocal, mo\_mrm\_global\_variables::ninflowgaugestotal, mo\_mrm\_global↔  
variables::nmeasperday, mo\_common\_constants::nodata\_dp, mo\_common\_mhm\_mrm\_variables::ntstepday,  
mo\_common\_mhm\_mrm\_variables::opti\_function, mo\_common\_mhm\_mrm\_variables::optimize, mo\_read↔  
timeseries::read\_timeseries(), mo\_mrm\_global\_variables::riv\_temp\_pcs, mo\_common\_mhm\_mrm\_variables↔  
::simper, and mo\_mrm\_file::udischarge.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.46.2.3 mrm\_read\_l0\_data()

```
subroutine, public mo_mrm_read_data::mrm_read_l0_data (
    logical, intent(in) do_reinit,
    logical, intent(in) do_readlatlon,
    logical, intent(in) do_readlcover )
```

read L0 data from file

With the exception of L0\_mask, L0\_elev, and L0\_LCover, all L0 variables are read from file. The former three are only read if they are not provided as variables.

#### Parameters

|    |                                 |  |
|----|---------------------------------|--|
| in | <i>logical :: do_reinit</i>     |  |
| in | <i>logical :: do_readlatlon</i> |  |
| in | <i>logical :: do_readlcover</i> |  |

#### Authors

Juliane Mai, Matthias Zink, and Stephan Thober

#### Date

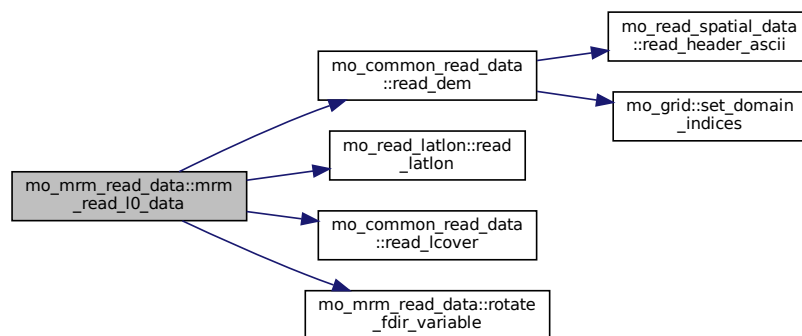
Aug 2015

Definition at line 50 of file mo\_mrm\_read\_data.f90.

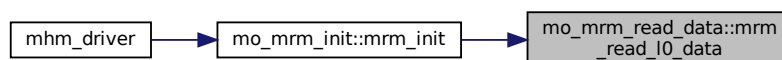
References mo\_common\_variables::dirmorpho, mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_file::file\_facc, mo\_mrm\_file::file\_fdir, mo\_mrm\_file::file\_gaugeloc, mo\_mpr\_file::file\_slope, mo\_mrm\_global\_variables::l0\_facc, mo\_mrm\_global\_variables::l0\_fdir, mo\_mrm\_global\_variables::l0\_gaugeloc, mo\_mrm\_global\_variables::l0\_inflowgaugeloc, mo\_common\_variables::l0\_lcover, mo\_common\_variables::level0, mo\_common\_constants::nodata\_i4, mo\_common\_variables::processmatrix, mo\_common\_read\_data::read\_dem(), mo\_read\_latlon::read\_latlon(), mo\_common\_read\_data::read\_lcover(), rotate\_fdir\_variable(), mo\_mrm\_file::ufacc, mo\_mrm\_file::ufdir, mo\_mrm\_file::ugaugeloc, and mo\_mpr\_file::uslope.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.46.2.4 mrm\_read\_total\_runoff()

```
subroutine, public mo_mrm_read_data::mrm_read_total_runoff (
    integer(i4), intent(in) iDomain )
```

read simulated runoff that is to be routed

read spatio-temporal field of total runoff that has been simulated by a hydrologic model or land surface model. This total runoff will then be aggregated to the level 11 resolution and then routed through the stream network.

#### Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | domain id |
|----|-------------------------------|-----------|

#### Authors

Stephan Thober

#### Date

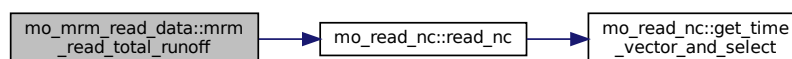
Sep 2015

Definition at line 373 of file mo\_mrm\_read\_data.f90.

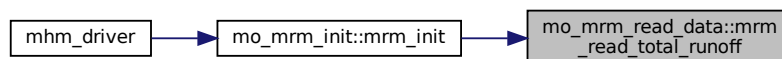
References mo\_common\_variables::alma\_convention, mo\_mrm\_global\_variables::dirtotalrunoff, mo\_mrm\_global\_variables::filenametotalrunoff, mo\_mrm\_global\_variables::l1\_total\_runoff\_in, mo\_common\_variables::level1, mo\_common\_constants::nodata\_dp, mo\_read\_nc::read\_nc(), mo\_common\_mhm\_mrm\_variables::simper, mo\_common\_mhm\_mrm\_variables::timestep, and mo\_mrm\_global\_variables::varnametotalrunoff.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.46.2.5 rotate\_fdir\_variable()

```
subroutine mo_mrm_read_data::rotate_fdir_variable (
    integer(i4), dimension(:, :), intent(inout) x )
```



TODO: add description.

TODO: add description

#### Parameters

|         |  |
|---------|--|
| in, out | <i>integer(i4), dimension(:, :) :: x</i> |
|---------|--|

#### Authors

L. Samaniego & R. Kumar

#### Date

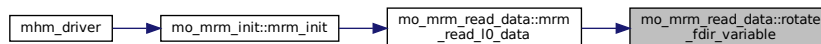
Jun 2018

Definition at line 533 of file mo\_mrm\_read\_data.f90.

References mo\_common\_constants::nodata\_i4.

Referenced by mrm\_read\_I0\_data().

Here is the caller graph for this function:



## 17.47 mo\_mrm\_restart Module Reference

Restart routines.

### Functions/Subroutines

- subroutine, public [mrm\\_write\\_restart](#) (iDomain, domainID, OutFile)  
*write routing states and configuration*
- subroutine, public [mrm\\_read\\_restart\\_states](#) (iDomain, domainID, InFile)  
*read routing states*
- subroutine, public [mrm\\_read\\_restart\\_config](#) (iDomain, domainID, InFile)  
*reads Level 11 configuration from a restart directory*

#### 17.47.1 Detailed Description

Restart routines.

This module contains the subroutines for reading and writing routing related variables to file.

#### Authors

Stephan Thober

#### Date

Aug 2015

## 17.47.2 Function/Subroutine Documentation

### 17.47.2.1 mrm\_read\_restart\_config()

```
subroutine, public mo_mrm_restart::mrm_read_restart_config (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) domainID,
    character(256), intent(in) InFile )
```

reads Level 11 configuration from a restart directory

read Level 11 configuration variables from a given restart directory and initializes all Level 11 configuration variables, that are initialized in L11\_variable\_init, contained in module [mo\\_startup](#).

#### Parameters

|    |                                 |                                     |
|----|---------------------------------|-------------------------------------|
| in | <i>integer(i4) :: iDomain</i>   | number of Domain                    |
| in | <i>character(256) :: InFile</i> | Input Path including trailing slash |

#### Authors

Stephan Thober

#### Date

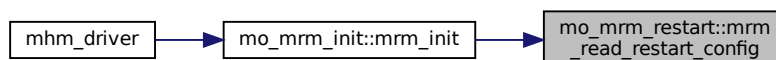
Apr 2013

Definition at line 603 of file mo\_mrm\_restart.f90.

References mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_facc, mo\_mrm\_global\_variables::l0\_fdir, mo\_mpr\_global\_variables::l0\_slope, mo\_mrm\_global\_variables::l0\_streamnet, mo\_mrm\_global\_variables::l11\_afloodplain, mo\_mrm\_global\_variables::l11\_colout, mo\_mrm\_global\_variables::l11\_facc, mo\_mrm\_global\_variables::l11\_fcol, mo\_mrm\_global\_variables::l11\_fdir, mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_global\_variables::l11\_frow, mo\_mrm\_global\_variables::l11\_l1\_id, mo\_mrm\_global\_variables::l11\_label, mo\_mrm\_global\_variables::l11\_length, mo\_mrm\_global\_variables::l11\_netperm, mo\_mrm\_global\_variables::l11\_noutlets, mo\_mrm\_global\_variables::l11\_rorder, mo\_mrm\_global\_variables::l11\_rowout, mo\_mrm\_global\_variables::l11\_sink, mo\_mrm\_global\_variables::l11\_slope, mo\_mrm\_global\_variables::l11\_tcol, mo\_mrm\_global\_variables::l11\_ton, mo\_mrm\_global\_variables::l11\_trow, mo\_mrm\_global\_variables::l11\_tsrou, mo\_mrm\_global\_variables::l1\_l1\_id, mo\_common\_variables::level0, mo\_common\_variables::level1, mo\_mrm\_global\_variables::level11, mo\_common\_constants::nodata\_dp, and mo\_common\_variables::processmatrix.

Referenced by mo\_mrm\_init::mrm\_init().

Here is the caller graph for this function:



### 17.47.2.2 mrm\_read\_restart\_states()

```
subroutine, public mo_mrm_restart::mrm_read_restart_states (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) domainID,
    character(256), intent(in) InFile )
```

read routing states

This subroutine reads the routing states from `mRM_states_<domain_id>.nc` that has to be in the given path directory. This subroutine has to be called directly each time the `mHM_eval` or `mRM_eval` is called such that the the states are always the same at the first simulation time step, crucial for optimization.

#### Parameters

|    |                                 |                                     |
|----|---------------------------------|-------------------------------------|
| in | <i>integer(i4) :: iDomain</i>   | number of domains                   |
| in | <i>character(256) :: InFile</i> | Input Path including trailing slash |

#### Authors

Stephan Thober

#### Date

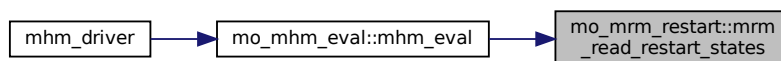
Sep 2015

Definition at line 460 of file `mo_mrm_restart.f90`.

References `mo_mrm_global_variables::l11_c1`, `mo_mrm_global_variables::l11_c2`, `mo_mrm_global_variables::l11_k`, `mo_mrm_global_variables::l11_nlinkfracpimp`, `mo_mrm_global_variables::l11_qmod`, `mo_mrm_global_variables::l11_qout`, `mo_mrm_global_variables::l11_qtin`, `mo_mrm_global_variables::l11_qtr`, `mo_mrm_global_variables::l11_xi`, `mo_mrm_global_variables::level11`, `mo_common_variables::nlcoverscene`, and `mo_mrm_constants::nroutingstates`.

Referenced by `mo_mhm_eval::mhm_eval()`.

Here is the caller graph for this function:



### 17.47.2.3 mrm\_write\_restart()

```
subroutine, public mo_mrm_restart::mrm_write_restart (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) domainID,
    character(256), dimension(:), intent(in) OutFile )
```

write routing states and configuration

write configuration and state variables to a given restart directory.

## Parameters

|    |  |                                 |
|----|--|---------------------------------|
| in | <i>integer(i4) :: iDomain</i>                  | number of domain                |
| in | <i>character(256), dimension(:) :: OutFile</i> | list of Output paths per Domain |

## Authors

Stephan Thober

## Date

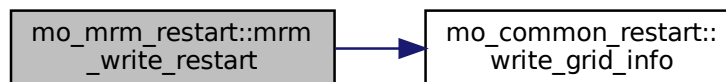
Aug 2015

Definition at line 55 of file mo\_mrm\_restart.f90.

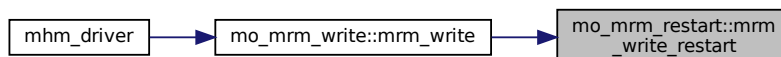
References mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_faccc, mo\_mrm\_global\_variables::l0\_fdir, mo\_mpr\_global\_variables::l0\_slope, mo\_mrm\_global\_variables::l0\_streamnet, mo\_mrm\_global\_variables::l11\_afloodplain, mo\_mrm\_global\_variables::l11\_c1, mo\_mrm\_global\_variables::l11\_c2, mo\_mrm\_global\_variables::l11\_celerity, mo\_mrm\_global\_variables::l11\_colout, mo\_mrm\_global\_variables::l11\_faccc, mo\_mrm\_global\_variables::l11\_fcol, mo\_mrm\_global\_variables::l11\_fdir, mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_global\_variables::l11\_frow, mo\_mrm\_global\_variables::l11\_k, mo\_mrm\_global\_variables::l11\_l1\_id, mo\_mrm\_global\_variables::l11\_label, mo\_mrm\_global\_variables::l11\_length, mo\_mrm\_global\_variables::l11\_netperm, mo\_mrm\_global\_variables::l11\_nlinkfracpimp, mo\_mrm\_global\_variables::l11\_qmod, mo\_mrm\_global\_variables::l11\_qout, mo\_mrm\_global\_variables::l11\_qtin, mo\_mrm\_global\_variables::l11\_qtr, mo\_mrm\_global\_variables::l11\_rorder, mo\_mrm\_global\_variables::l11\_rowout, mo\_mrm\_global\_variables::l11\_sink, mo\_mrm\_global\_variables::l11\_slope, mo\_mrm\_global\_variables::l11\_tcol, mo\_mrm\_global\_variables::l11\_ton, mo\_mrm\_global\_variables::l11\_trow, mo\_mrm\_global\_variables::l11\_tsout, mo\_mrm\_global\_variables::l11\_xi, mo\_mrm\_global\_variables::l1\_l1\_id, mo\_common\_constants::landcoverperiodsvarname, mo\_common\_variables::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::level0, mo\_common\_variables::level1, mo\_mrm\_global\_variables::level11, mo\_common\_variables::nlcoverscene, mo\_common\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, mo\_mrm\_constants::nroutingstates, mo\_common\_variables::processmatrix, and mo\_common\_restart::write\_grid\_info().

Referenced by mo\_mrm\_write::mrm\_write().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.48 mo\_mrm\_riv\_temp\_class Module Reference

Class for the river temperature calculations.

### Data Types

- module [riv\\_temp\\_type](#)

*This is a container to define the river temperature routing in the current time step.*

### Functions/Subroutines

- subroutine [config](#) (self, file\_namelist, unamelist, file\_namelist\_param, unamelist\_param)  
*configure the [riv\\_temp\\_type](#) class from the mhm namelist*
- subroutine [init](#) (self, nCells)  
*initialize the [riv\\_temp\\_type](#) class for the current domain*
- subroutine [init\\_area](#) (self, iDomain, L11\_netPerm, L11\_fromN, L11\_length, nLinks, nCells, nrows, ncols, L11\_mask)  
*initialize the river area of [riv\\_temp\\_type](#) class for the current domain*
- subroutine [init\\_riv\\_temp](#) (self, time, ntimesteps\_day, temp\_air, read\_meteo\_weights, temp\_weights, fday\_temp, fnight\_temp, efecarea, L1\_L11\_Id, L11\_areacell, L11\_L1\_Id, map\_flag)  
*initialize the river temperature of [riv\\_temp\\_type](#) class for the current domain*
- subroutine [reset\\_timestep](#) (self)  
*reset [riv\\_temp\\_type](#) class for next timestep*
- subroutine [alloc\\_lateral](#) (self, nCells)  
*allocate lateral temp components of [riv\\_temp\\_type](#) class for current domain*
- subroutine [dealloc\\_lateral](#) (self)  
*deallocate lateral temp components of [riv\\_temp\\_type](#)*
- subroutine [acc\\_source\\_e](#) (self, time, ntimesteps\_day, fSealed\_area\_fraction, fast\_interflow, slow\_interflow, baseflow, direct\_runoff, temp\_air, mean\_temp\_air, ssrd\_day, strd\_day, read\_meteo\_weights, temp\_weights, fday\_temp, fnight\_temp, fday\_ssrd, fnight\_ssrd, fday\_strd, fnight\_strd)  
*accumulate energy sources of [riv\\_temp\\_type](#)*
- subroutine [finalize\\_source\\_e](#) (self, efecarea, L1\_L11\_Id, L11\_areacell, L11\_L1\_Id, timestep, map\_flag)  
*finalize energy sources of [riv\\_temp\\_type](#)*
- real(dp) function [get\\_lrad\\_out](#) (self, riv\_temp)  
*get outgoing longwave radiation of [riv\\_temp\\_type](#)*
- real(dp) function [get\\_lat\\_heat](#) (self, air\_temp, netrad)  
*latent heat flux of [riv\\_temp\\_type](#)*
- real(dp) function [get\\_sens\\_heat](#) (self, air\_temp, riv\_temp)  
*sensible heat flux of [riv\\_temp\\_type](#)*
- real(dp) function [get\\_e\\_io](#) (self, riv\_temp, cell)  
*get complete energy source of [riv\\_temp\\_type](#) at given cell*
- subroutine [l11\\_routing\\_e](#) (self, nLinks, netPerm, netLink\_fromN, netLink\_toN, netLink\_C1, netLink\_C2, nInflowGauges, InflowHeadwater, InflowNodeList, L11\_qTR, L11\_Qmod)  
*execute the temperature routing of [riv\\_temp\\_type](#)*
- subroutine [init\\_iter](#) (self)  
*initialize iterative solver of [riv\\_temp\\_type](#)*
- subroutine [next\\_iter](#) (self, T\_est, T\_rout)  
*execute next iteration with iterative solver of [riv\\_temp\\_type](#)*

### 17.48.1 Detailed Description

Class for the river temperature calculations.

#### Warning

This feature is still experimental!

#### Version

0.1

#### Authors

Sebastian Mueller

#### Date

Sep 2020

### 17.48.2 Function/Subroutine Documentation

#### 17.48.2.1 `acc_source_e()`

```

subroutine mo_mrm_riv_temp_class::acc_source_e (
    class(riv_temp_type), intent(inout) self,
    real(dp), intent(in) time,
    real(dp), intent(in) ntimesteps_day,
    real(dp), dimension(:), intent(in) fsealed_area_fraction,
    real(dp), dimension(:), intent(in) fast_interflow,
    real(dp), dimension(:), intent(in) slow_interflow,
    real(dp), dimension(:), intent(in) baseflow,
    real(dp), dimension(:), intent(in) direct_runoff,
    real(dp), dimension(:), intent(in) temp_air,
    real(dp), dimension(:), intent(in) mean_temp_air,
    real(dp), dimension(:), intent(in) ssrd_day,
    real(dp), dimension(:), intent(in) strd_day,
    logical, intent(in) read_meteo_weights,
    real(dp), dimension(:, :, :), intent(in) temp_weights,
    real(dp), dimension(:), intent(in) fday_temp,
    real(dp), dimension(:), intent(in) fnight_temp,
    real(dp), dimension(:), intent(in) fday_ssrd,
    real(dp), dimension(:), intent(in) fnight_ssrd,
    real(dp), dimension(:), intent(in) fday_strd,
    real(dp), dimension(:), intent(in) fnight_strd ) [private]

```

accumulate energy sources of [riv\\_temp\\_type](#)

#### Parameters

|    |                              |   |
|----|------------------------------|---|
| in | <i>time</i>                  | current decimal Julian day                          |
| in | <i>ntimesteps_day</i>        | number of time intervals per day, transformed in dp |
| in | <i>fsealed_area_fraction</i> | sealed area fraction [1]                            |
| in | <i>fast_interflow</i>        | $q_0$ Fast runoff component [mm TS-1]               |

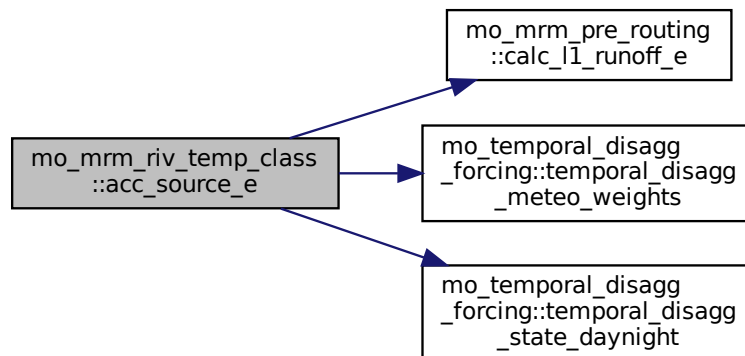
## Parameters

|    |                           |  |
|----|---------------------------|--|
| in | <i>slow_interflow</i>     | $q_1$ Slow runoff component [mm TS-1]                              |
| in | <i>baseflow</i>           | $q_2$ Baseflow [mm TS-1]   |
| in | <i>direct_runoff</i>      | $q_D$ Direct runoff from impervious areas [mm TS-1]                |
| in | <i>temp_air</i>           | air temperature [K]  |
| in | <i>mean_temp_air</i>      | annual mean air temperature [K]                                    |
| in | <i>ssrd_day</i>           | Daily mean short radiation   |
| in | <i>strd_day</i>           | Daily mean longwave radiation                                      |
| in | <i>read_meteo_weights</i> | flag whether weights for tavg and pet have read and should be used |
| in | <i>temp_weights</i>       | multiplicative weights for temperature (deg K)                     |
| in | <i>fday_temp</i>          | [-] day factor mean temp   |
| in | <i>fnight_temp</i>        | [-] night factor mean temp   |
| in | <i>fday_ssrd</i>          | Daytime fraction of ssrd   |
| in | <i>fnight_ssrd</i>        | Nighttime fraction of ssrd   |
| in | <i>fday_strd</i>          | Daytime fraction of strd   |
| in | <i>fnight_strd</i>        | Nighttime fraction of strd   |

Definition at line 441 of file mo\_mrm\_riv\_temp\_class.f90.

References mo\_mrm\_pre\_routing::calc\_l1\_runoff\_e(), mo\_temporal\_disagg\_forcing::temporal\_disagg\_meteo\_weights(), and mo\_temporal\_disagg\_forcing::temporal\_disagg\_state\_daynight().

Here is the call graph for this function:



### 17.48.2.2 alloc\_lateral()

```

subroutine mo_mrm_riv_temp_class::alloc_lateral (
    class(riv_temp_type), intent(inout) self,
    integer(i4), intent(in) nCells ) [private]
  
```

allocate lateral temp components of `riv_temp_type` class for current domain

## Parameters

|    |               |  |
|----|---------------|--|
| in | <i>ncells</i> | number of cells for the current domain |
|----|---------------|--|

Definition at line 407 of file `mo_mrm_riv_temp_class.f90`.

**17.48.2.3 config()**

```
subroutine mo_mrm_riv_temp_class::config (
    class(riv_temp_type), intent(inout) self,
    character(*), intent(in) file_namelist,
    integer, intent(in) unamelist,
    character(*), intent(in) file_namelist_param,
    integer, intent(in) unamelist_param )
```

configure the `riv_temp_type` class from the mhm namelist

## Parameters

|    |                            |                             |
|----|----------------------------|-----------------------------|
| in | <i>file_namelist</i>       | mhm namelist file           |
| in | <i>file_namelist_param</i> | mhm parameter namelist file |

Definition at line 117 of file `mo_mrm_riv_temp_class.f90`.

References `mo_check::check_dir()`, `mo_common_variables::domainmeta`, `mo_common_constants::maxnodomains`, and `mo_common_constants::nodata_i4`.

Here is the call graph for this function:

**17.48.2.4 dealloc\_lateral()**

```
subroutine mo_mrm_riv_temp_class::dealloc_lateral (
    class(riv_temp_type), intent(inout) self ) [private]
```

deallocate lateral temp components of `riv_temp_type`

Definition at line 426 of file `mo_mrm_riv_temp_class.f90`.

**17.48.2.5 finalize\_source\_e()**

```
subroutine mo_mrm_riv_temp_class::finalize_source_e (
```



```

class(riv_temp_type), intent(inout) self,
real(dp), dimension(:), intent(in) efecarea,
integer(i4), dimension(:), intent(in) L1_L11_Id,
real(dp), dimension(:), intent(in) L11_areacell,
integer(i4), dimension(:), intent(in) L11_L1_Id,
integer(i4), intent(in) timestep,
logical, intent(in) map_flag )

```

finalize energy sources of [riv\\_temp\\_type](#)

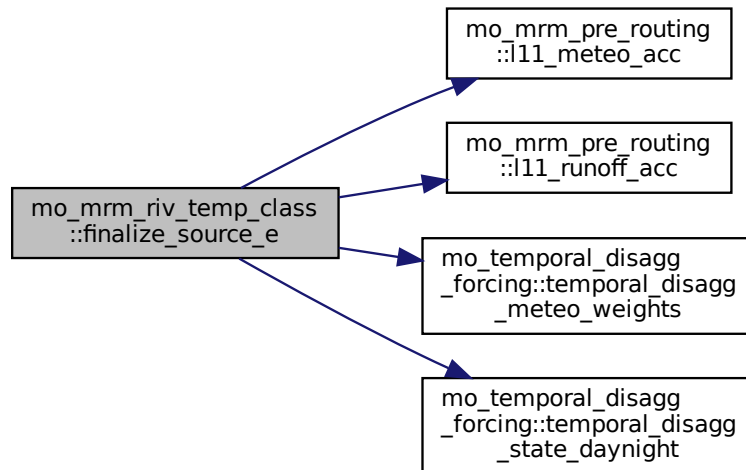
#### Parameters

|    |                     |  |
|----|---------------------|--|
| in | <i>efecarea</i>     | effective area in [km2] at Level 1                                       |
| in | <i>l1_l11_id</i>    | L11 lds mapped on L1   |
| in | <i>l11_areacell</i> | effective area in [km2] at Level 11                                      |
| in | <i>l11_l1_id</i>    | L1 lds mapped on L11   |
| in | <i>timestep</i>     | simulation timestep in [h]   |
| in | <i>map_flag</i>     | Flag indicating whether routing resolution is higher than hydrologic one |

Definition at line 543 of file mo\_mrm\_riv\_temp\_class.f90.

References [mo\\_mhm\\_constants::h2odens](#), [mo\\_mrm\\_pre\\_routing::l11\\_meteo\\_acc\(\)](#), [mo\\_mrm\\_pre\\_routing::l11\\_runoff\\_acc\(\)](#), [mo\\_temporal\\_disagg\\_forcing::temporal\\_disagg\\_meteo\\_weights\(\)](#), and [mo\\_temporal\\_disagg\\_forcing::temporal\\_disagg\\_state\\_daynight\(\)](#).

Here is the call graph for this function:



#### 17.48.2.6 get\_e\_io()

```

real(dp) function mo_mrm_riv_temp_class::get_e_io (
    class(riv_temp_type), intent(in) self,
    real(dp), intent(in) riv_temp,
    integer(i4), intent(in) cell ) [private]

```

get complete energy source of [riv\\_temp\\_type](#) at given cell

#### Returns

energy IO

#### Parameters

|    |                 |   |
|----|-----------------|---|
| in | <i>riv_temp</i> | given river temperature in K to calculate heat fluxes |
| in | <i>cell</i>     | cell index in the current domain                      |

net radiation calc from short and longwave radiation in/out

Definition at line 667 of file `mo_mrm_riv_temp_class.f90`.

References `mo_mhm_constants::h2odens`.

#### 17.48.2.7 `get_lat_heat()`

```
real(dp) function mo_mrm_riv_temp_class::get_lat_heat (
    class(riv\_temp\_type), intent(in) self,
    real(dp), intent(in) air_temp,
    real(dp), intent(in) netrad )
```

latent heat flux of [riv\\_temp\\_type](#)

#### Returns

latent heat flux

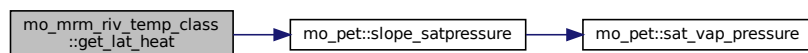
#### Parameters

|    |                 |                          |
|----|-----------------|--------------------------|
| in | <i>air_temp</i> | air temperature in deg C |
| in | <i>netrad</i>   | net radiation in W * m-2 |

Definition at line 628 of file `mo_mrm_riv_temp_class.f90`.

References `mo_pet::slope_satpressure()`.

Here is the call graph for this function:



#### 17.48.2.8 `get_lrad_out()`

```
real(dp) function mo_mrm_riv_temp_class::get_lrad_out (
    class(riv\_temp\_type), intent(in) self,
    real(dp), intent(in) riv_temp )
```

get outgoing longwave radiation of [riv\\_temp\\_type](#)

#### Returns

outgoing longwave radiation

#### Parameters

|    |                 |                        |
|----|-----------------|------------------------|
| in | <i>riv_temp</i> | river temperature in K |
|----|-----------------|------------------------|

Definition at line 611 of file mo\_mrm\_riv\_temp\_class.f90.

#### 17.48.2.9 get\_sens\_heat()

```
real(dp) function mo_mrm_riv_temp_class::get_sens_heat (
    class(riv\_temp\_type), intent(in) self,
    real(dp), intent(in) air_temp,
    real(dp), intent(in) riv_temp )
```

sensible heat flux of [riv\\_temp\\_type](#)

#### Returns

sensible heat flux

#### Parameters

|    |                 |                              |
|----|-----------------|------------------------------|
| in | <i>air_temp</i> | air temperature in [deg C]   |
| in | <i>riv_temp</i> | river temperature in [deg C] |

Definition at line 652 of file mo\_mrm\_riv\_temp\_class.f90.

#### 17.48.2.10 init()

```
subroutine mo_mrm_riv_temp_class::init (
    class(riv\_temp\_type), intent(inout) self,
    integer(i4), intent(in) nCells )
```

initialize the [riv\\_temp\\_type](#) class for the current domain

#### Parameters

|    |               |  |
|----|---------------|--|
| in | <i>ncells</i> | number of cells for the current domain |
|----|---------------|--|

Definition at line 215 of file mo\_mrm\_riv\_temp\_class.f90.

References mo\_mrm\_constants::nroutingstates.

### 17.48.2.11 `init_area()`

```

subroutine mo_mrm_riv_temp_class::init_area (
    class(riv_temp_type), intent(inout) self,
    integer(i4), intent(in) iDomain,
    integer(i4), dimension(:), intent(in) L11_netPerm,
    integer(i4), dimension(:), intent(in) L11_fromN,
    real(dp), dimension(:), intent(in) L11_length,
    integer(i4), intent(in) nLinks,
    integer(i4), intent(in) nCells,
    integer(i4), intent(in) nrows,
    integer(i4), intent(in) ncols,
    logical, dimension(:, :), intent(in) L11_mask )

```

initialize the river area of `riv_temp_type` class for the current domain

#### Parameters

|    |                          |  |
|----|--------------------------|--|
| in | <code>idomain</code>     | Domain ID  |
| in | <code>l11_netperm</code> | L11 routing order  |
| in | <code>l11_fromn</code>   | L11 source grid cell order                                   |
| in | <code>l11_length</code>  | L11 link length  |
| in | <code>nlinks</code>      | number of L11 links in the current domain                    |
| in | <code>ncells</code>      | number of L11 cells of the current domain                    |
| in | <code>ncols</code>       | Number of columns  |
| in | <code>nrows</code>       | Number of rows   |
| in | <code>l11_mask</code>    | the mask for valid cells in the original grid (nrows, ncols) |

Definition at line 258 of file `mo_mrm_riv_temp_class.f90`.

References `mo_read_nc::read_const_nc()`.

Here is the call graph for this function:



### 17.48.2.12 `init_iter()`

```

subroutine mo_mrm_riv_temp_class::init_iter (
    class(riv_temp_type), intent(inout) self )

```

initialize iterative solver of `riv_temp_type`

Definition at line 812 of file `mo_mrm_riv_temp_class.f90`.

**17.48.2.13 init\_riv\_temp()**

```

subroutine mo_mrm_riv_temp_class::init_riv_temp (
    class(riv_temp_type), intent(inout) self,
    real(dp), intent(in) time,
    real(dp), intent(in) ntimesteps_day,
    real(dp), dimension(:), intent(in) temp_air,
    logical, intent(in) read_meteo_weights,
    real(dp), dimension(:, :, :), intent(in) temp_weights,
    real(dp), dimension(:), intent(in) fday_temp,
    real(dp), dimension(:), intent(in) fnight_temp,
    real(dp), dimension(:), intent(in) efecarea,
    integer(i4), dimension(:), intent(in) L1_L11_Id,
    real(dp), dimension(:), intent(in) L11_areacell,
    integer(i4), dimension(:), intent(in) L11_L1_Id,
    logical, intent(in) map_flag )

```

initialize the river temperature of `riv_temp_type` class for the current domain

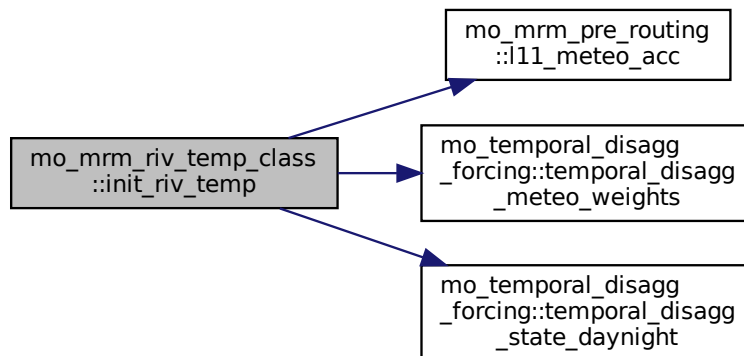
**Parameters**

|    |                           |  |
|----|---------------------------|--|
| in | <i>time</i>               | current decimal Julian day   |
| in | <i>ntimesteps_day</i>     | number of time intervals per day, transformed in dp                      |
| in | <i>temp_air</i>           | air temperature [K]  |
| in | <i>read_meteo_weights</i> | flag whether weights for tavg and pet have read and should be used       |
| in | <i>temp_weights</i>       | multiplicative weights for temperature (deg K)                           |
| in | <i>fday_temp</i>          | [-] day factor mean temp   |
| in | <i>fnight_temp</i>        | [-] night factor mean temp   |
| in | <i>efecarea</i>           | effective area in [km2] at Level 1                                       |
| in | <i>l1_l11_id</i>          | L11 lds mapped on L1   |
| in | <i>l11_areacell</i>       | effective area in [km2] at Level 11                                      |
| in | <i>l11_l1_id</i>          | L1 lds mapped on L11   |
| in | <i>map_flag</i>           | Flag indicating whether routing resolution is higher than hydrologic one |

Definition at line 322 of file mo\_mrm\_riv\_temp\_class.f90.

References `mo_mrm_pre_routing::l11_meteo_acc()`, `mo_temporal_disagg_forcing::temporal_disagg_meteo_↔weights()`, and `mo_temporal_disagg_forcing::temporal_disagg_state_daynight()`.

Here is the call graph for this function:



#### 17.48.2.14 l11\_routing\_e()

```

subroutine mo_mrm_riv_temp_class::l11_routing_e (
    class(riv_temp_type), intent(inout) self,
    integer(i4), intent(in) nLinks,
    integer(i4), dimension(:), intent(in) netPerm,
    integer(i4), dimension(:), intent(in) netLink_fromN,
    integer(i4), dimension(:), intent(in) netLink_toN,
    real(dp), dimension(:), intent(in) netLink_C1,
    real(dp), dimension(:), intent(in) netLink_C2,
    integer(i4), intent(in) nInflowGauges,
    logical, dimension(:), intent(in) InflowHeadwater,
    integer(i4), dimension(:), intent(in) InflowNodeList,
    real(dp), dimension(:), intent(in) L11_qTR,
    real(dp), dimension(:), intent(in) L11_Qmod )
  
```

execute the temperature routing of [riv\\_temp\\_type](#)

#### Parameters

|    |                        |  |
|----|------------------------|--|
| in | <i>nlinks</i>          | number of stream segment (reaches)   |
| in | <i>netperm</i>         | routing order of a given domain (permutation)                              |
| in | <i>netlink_fromn</i>   | from node  |
| in | <i>netlink_ton</i>     | to node  |
| in | <i>netlink_c1</i>      | routing parameter C1 ([7] p. 25-41)  |
| in | <i>netlink_c2</i>      | routing parameters C2 (id)   |
| in | <i>ninflowgauges</i>   | [-] number of inflow points  |
| in | <i>inflowheadwater</i> | [-] if to consider headwater cells of inflow gauge                         |
| in | <i>inflowodelist</i>   | [-] L11 ID of inflow points  |
| in | <i>l11_qtr</i>         | [m3 s-1] Transformed outflow leaving node I at current timestep(Muskingum) |
| in | <i>l11_qmod</i>        | [m3 s-1] Simulated routed discharge  |

Definition at line 696 of file mo\_mrm\_riv\_temp\_class.f90.

#### 17.48.2.15 next\_iter()

```
subroutine mo_mrm_riv_temp_class::next_iter (
    class(riv_temp_type), intent(inout) self,
    real(dp), intent(inout) T_est,
    real(dp), intent(in) T_rout ) [private]
```

execute next iteration with iterative solver of [riv\\_temp\\_type](#)

##### Parameters

|         |                        |                                       |
|---------|------------------------|---------------------------------------|
| in, out | <a href="#">t_est</a>  | estimated river temperature           |
| in      | <a href="#">t_rout</a> | calculated (routed) river temperature |

Definition at line 826 of file mo\_mrm\_riv\_temp\_class.f90.

#### 17.48.2.16 reset\_timestep()

```
subroutine mo_mrm_riv_temp_class::reset_timestep (
    class(riv_temp_type), intent(inout) self )
```

reset [riv\\_temp\\_type](#) class for next timestep

Definition at line 391 of file mo\_mrm\_riv\_temp\_class.f90.

## 17.49 mo\_mrm\_river\_head Module Reference

### Functions/Subroutines

- subroutine, public [init\\_masked\\_zeros\\_I0](#) (iDomain, data)
- subroutine, private [reset\\_sum](#) (iDomain, data)
- subroutine, public [calc\\_channel\\_elevation](#) ()
- subroutine, public [calc\\_river\\_head](#) (iDomain, L11\_Qmod, river\_head)
- real(dp) function [calc\\_slope](#) (iDomain, elev0, fDir0, i, j)
- subroutine, public [avg\\_and\\_write\\_timestep](#) (iDomain, timestep, data)
- subroutine, public [create\\_output](#) (iDomain, OutPath)

### Variables

- type(ncvariable), dimension(:), allocatable [nc\\_time](#)
- type(ncvariable), dimension(:), allocatable [nc\\_riverhead](#)
- integer(i4), dimension(:), allocatable [time\\_counter](#)
- integer(i4), dimension(:), allocatable [sum\\_counter](#)

### 17.49.1 Function/Subroutine Documentation

### 17.49.1.1 avg\_and\_write\_timestep()

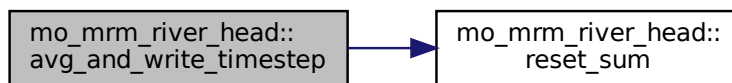
```
subroutine, public mo_mrm_river_head::avg_and_write_timestep (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) timestep,
    real(dp), dimension(:), intent(inout) data )
```

Definition at line 188 of file mo\_mrm\_river\_head.f90.

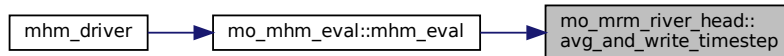
References mo\_common\_variables::level0, nc\_riverhead, nc\_time, mo\_common\_constants::nodata\_dp, reset\_  
sum(), sum\_counter, and time\_counter.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.49.1.2 calc\_channel\_elevation()

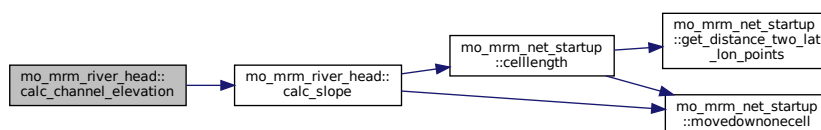
```
subroutine, public mo_mrm_river_head::calc_channel_elevation
```

Definition at line 67 of file mo\_mrm\_river\_head.f90.

References calc\_slope(), mo\_common\_variables::domainmeta, mo\_mrm\_global\_variables::l0\_channel\_depth, mo\_common\_variables::l0\_elev, mo\_mrm\_global\_variables::l0\_facc, mo\_mrm\_global\_variables::l0\_fdir, and mo\_  
\_common\_constants::nodata\_i4.

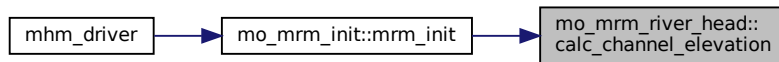
Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:





Here is the caller graph for this function:



### 17.49.1.3 calc\_river\_head()

```

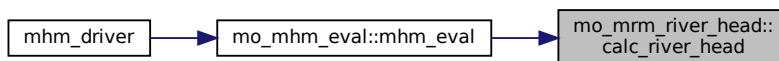
subroutine, public mo_mrm_river_head::calc_river_head (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(:), intent(in) L11_Qmod,
    real(dp), dimension(:), intent(inout), allocatable river_head )
  
```

Definition at line 133 of file mo\_mrm\_river\_head.f90.

References mo\_mrm\_global\_variables::l0\_channel\_elevation, mo\_mrm\_global\_variables::l0\_l11\_remap, mo\_mrm\_global\_variables::l0\_slope, mo\_mrm\_global\_variables::l11\_bankfull\_runoff\_in, mo\_common\_variables::level0, and sum\_counter.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the caller graph for this function:



### 17.49.1.4 calc\_slope()

```

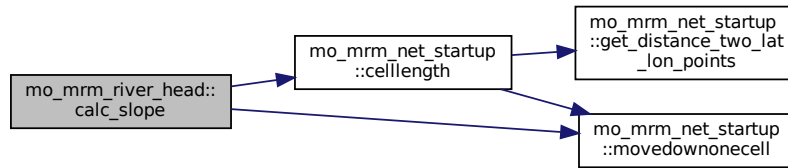
real(dp) function mo_mrm_river_head::calc_slope (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(:, :), intent(in), allocatable elev0,
    integer(i4), dimension(:, :), intent(in), allocatable fDir0,
    integer(i4), intent(in) i,
    integer(i4), intent(in) j ) [private]
  
```

Definition at line 162 of file mo\_mrm\_river\_head.f90.

References mo\_mrm\_net\_startup::celllength(), mo\_common\_variables::iflag\_cordinate\_sys, and mo\_mrm\_net\_startup::movedownonecell().

Referenced by calc\_channel\_elevation().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.49.1.5 create\_output()

```

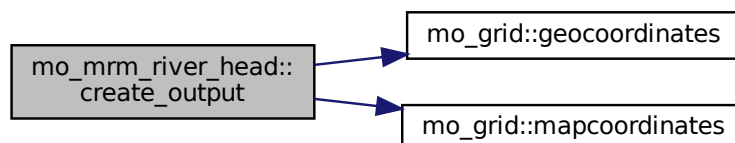
subroutine, public mo_mrm_river_head::create_output (
    integer(i4), intent(in) iDomain,
    character(256), intent(in) OutPath )
  
```

Definition at line 203 of file `mo_mrm_river_head.f90`.

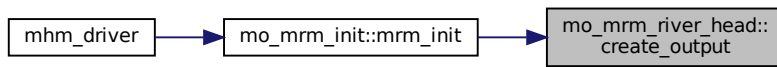
References `mo_common_variables::contact`, `mo_common_variables::conventions`, `mo_common_mhm_mrm_variables::evalper`, `mo_grid::geocoordinates()`, `mo_common_variables::history`, `mo_grid::mapcoordinates()`, `mo_common_variables::mhm_details`, `nc_riverhead`, `nc_time`, `mo_common_constants::nodata_dp`, `mo_common_variables::project_details`, `mo_common_variables::setup_description`, `mo_common_variables::simulation_type`, `sum_counter`, `time_counter`, and `mo_file::version`.

Referenced by `mo_mrm_init::mrm_init()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.49.1.6 init\_masked\_zeros\_l0()

```

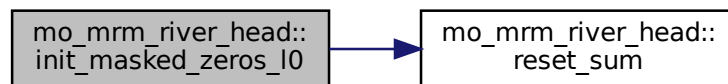
subroutine, public mo_mrm_river_head::init_masked_zeros_l0 (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(:), intent(inout), allocatable data )
  
```

Definition at line 38 of file mo\_mrm\_river\_head.f90.

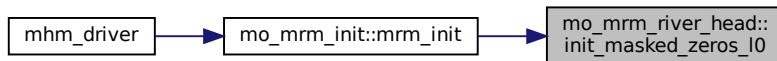
References mo\_common\_variables::level0, mo\_common\_constants::nodata\_dp, and reset\_sum().

Referenced by mo\_mrm\_init::mrm\_init().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.49.1.7 reset\_sum()

```

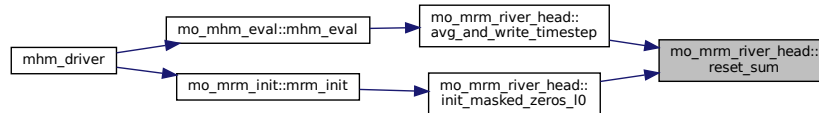
subroutine, private mo_mrm_river_head::reset_sum (
    integer(i4), intent(in) iDomain,
    real(dp), dimension(:), intent(inout) data ) [private]
  
```

Definition at line 51 of file mo\_mrm\_river\_head.f90.

References `mo_common_variables::level0`.

Referenced by `avg_and_write_timestep()`, and `init_masked_zeros_l0()`.

Here is the caller graph for this function:



## 17.49.2 Variable Documentation

### 17.49.2.1 nc\_riverhead

```
type(ncvariable), dimension(:), allocatable mo_mrm_river_head::nc_riverhead [private]
```

Definition at line 28 of file `mo_mrm_river_head.f90`.

Referenced by `avg_and_write_timestep()`, and `create_output()`.

### 17.49.2.2 nc\_time

```
type(ncvariable), dimension(:), allocatable mo_mrm_river_head::nc_time [private]
```

Definition at line 28 of file `mo_mrm_river_head.f90`.

Referenced by `avg_and_write_timestep()`, and `create_output()`.

### 17.49.2.3 sum\_counter

```
integer(i4), dimension(:), allocatable mo_mrm_river_head::sum_counter [private]
```

Definition at line 32 of file `mo_mrm_river_head.f90`.

Referenced by `avg_and_write_timestep()`, `calc_river_head()`, and `create_output()`.

### 17.49.2.4 time\_counter

```
integer(i4), dimension(:), allocatable mo_mrm_river_head::time_counter [private]
```

Definition at line 30 of file `mo_mrm_river_head.f90`.

Referenced by `avg_and_write_timestep()`, and `create_output()`.

## 17.50 mo\_mrm\_routing Module Reference

Performs runoff routing for mHM at level L11.

## Functions/Subroutines

- subroutine, public [mrm\\_routing](#) (read\_states, processCase, global\_routing\_param, L1\_total\_runoff, L1\_areaCell, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, L11\_netPerm, L11\_fromN, L11\_toN, L11\_nOutlets, timestep, tsRoutFactor, nNodes, nInflowGauges, InflowGaugeIndexList, InflowGaugeHeadwater, InflowGaugeNodeList, InflowDischarge, nGauges, gaugeIndexList, gaugeNodeList, map\_flag, L11\_length, L11\_slope, L11\_FracFPimp, L11\_C1, L11\_C2, L11\_qOut, L11\_qTIN, L11\_qTR, L11\_qMod, GaugeDischarge)  
*route water given runoff*
- subroutine [l11\\_routing](#) (nNodes, nLinks, netPerm, netLink\_fromN, netLink\_toN, netLink\_C1, netLink\_C2, netNode\_qOUT, nInflowGauges, InflowHeadwater, InflowNodeList, netNode\_qTIN, netNode\_qTR, netNode\_qMod)  
*Performs runoff routing for mHM at L11 upscaled network ([Routing Network](#)).*

### 17.50.1 Detailed Description

Performs runoff routing for mHM at level L11.

This module performs flood routing at a given time step through the stream network at level L11 to the sink cell. The Muskingum flood routing algorithm is used.

#### Authors

Luis Samaniego

#### Date

Dec 2012

### 17.50.2 Function/Subroutine Documentation

#### 17.50.2.1 l11\_routing()

```
subroutine mo_mrm_routing::l11_routing (
    integer(i4), intent(in) nNodes,
    integer(i4), intent(in) nLinks,
    integer(i4), dimension(:), intent(in) netPerm,
    integer(i4), dimension(:), intent(in) netLink_fromN,
    integer(i4), dimension(:), intent(in) netLink_toN,
    real(dp), dimension(:), intent(in) netLink_C1,
    real(dp), dimension(:), intent(in) netLink_C2,
    real(dp), dimension(:), intent(in) netNode_qOUT,
    integer(i4), intent(in) nInflowGauges,
    logical, dimension(:), intent(in) InflowHeadwater,
    integer(i4), dimension(:), intent(in) InflowNodeList,
    real(dp), dimension(:, :), intent(inout) netNode_qTIN,
    real(dp), dimension(:, :), intent(inout) netNode_qTR,
    real(dp), dimension(nnodes), intent(out) netNode_qmod )
```

Performs runoff routing for mHM at L11 upscaled network ([Routing Network](#)).

Hydrograph routing is carried out with the Muskingum algorithm [7]. This simplification of the St. Venant equations is justified in mHM because the potential areas of application of this model would hardly exhibit abruptly changing hydrographs with supercritical flows. The discharge leaving the river reach located on cell  $i$   $Q_i^1(t)$  at time step  $t$  can be determined by

$$Q_i^1(t) = Q_i^1(t-1) + c_1 (Q_i^0(t-1) - Q_i^1(t-1)) + c_2 (Q_i^0(t) - Q_i^0(t-1))$$

with

$$Q_i^0(t) = Q_i^i(t) + Q_i^1(t)$$

$$c_1 = \frac{\Delta t}{\kappa(1 - \xi) + \frac{\Delta t}{2}}$$

$$c_2 = \frac{\frac{\Delta t}{2} - \kappa\xi}{\kappa(1 - \xi) + \frac{\Delta t}{2}}$$

where  $Q_i^0$  and  $Q_i^1$  denote the discharge entering and leaving the river reach located on cell  $i$  respectively.  $Q_i^i$  is the contribution from the upstream cell  $i$ .  $\kappa$  Muskingum travel time parameter.  $\xi$  Muskingum attenuation parameter.  $\Delta t$  time interval in hours.  $t$  Time index for each  $\Delta t$  interval. To improve performance, a routing sequence "netPerm" is required. This permutation is determined in the mo\_init\_mrm routine.

TODO: add description

#### Parameters

|         |  |  |
|---------|--|--|
| in      | <i>integer(i4) :: nNodes</i>                       | number of network nodes = nCells1                                  |
| in      | <i>integer(i4) :: nLinks</i>                       | number of stream segment (reaches)                                 |
| in      | <i>integer(i4), dimension(:) :: netPerm</i>        | routing order of a given domain (permutation)                      |
| in      | <i>integer(i4), dimension(:) :: netLink_fromN</i>  | from node  |
| in      | <i>integer(i4), dimension(:) :: netLink_toN</i>    | to node  |
| in      | <i>real(dp), dimension(:) :: netLink_C1</i>        | routing parameter C1 ([7] p. 25-41)                                |
| in      | <i>real(dp), dimension(:) :: netLink_C2</i>        | routing parameters C2 (id)   |
| in      | <i>real(dp), dimension(:) :: netNode_qOUT</i>      | Total outflow from cells (given domain) L11 at time tt in [m3 s-1] |
| in      | <i>integer(i4) :: nInflowGauges</i>                | [-] number of inflow points  |
| in      | <i>logical, dimension(:) :: InflowHeadwater</i>    | [-] if to consider headwater cells of inflow gauge                 |
| in      | <i>integer(i4), dimension(:) :: InflowNodeList</i> | [-] L11 ID of inflow points  |
| in, out | <i>real(dp), dimension(:, :) :: netNode_qTIN</i>   | [m3 s-1] Total inputs at t-1 and t                                 |
| in, out | <i>real(dp), dimension(:, :) :: netNode_qTR</i>    | [m3 s-1] Transformed outflow leaving node l (Muskingum)            |
| out     | <i>real(dp), dimension(nNodes) :: netNode_Qmod</i> | [m3 s-1] Simulated routed discharge                                |

#### Authors

Luis Samaniego

#### Date

Dec 2005

Definition at line 379 of file mo\_mrm\_routing.f90.

Referenced by mrm\_routing().

Here is the caller graph for this function:



## 17.50.2.2 mrm\_routing()

```

subroutine, public mo_mrm_routing::mrm_routing (
    logical, intent(in) read_states,
    integer(i4), intent(in) processCase,
    real(dp), dimension(:), intent(in) global_routing_param,
    real(dp), dimension(:), intent(in) L1_total_runoff,
    real(dp), dimension(:), intent(in) L1_areaCell,
    integer(i4), dimension(:), intent(in) L1_L11_Id,
    real(dp), dimension(:), intent(in) L11_areaCell,
    integer(i4), dimension(:), intent(in) L11_L1_Id,
    integer(i4), dimension(:), intent(in) L11_netPerm,
    integer(i4), dimension(:), intent(in) L11_fromN,
    integer(i4), dimension(:), intent(in) L11_toN,
    integer(i4), intent(in) L11_nOutlets,
    integer(i4), intent(in) timestep,
    real(dp), intent(in) tsRoutFactor,
    integer(i4), intent(in) nNodes,
    integer(i4), intent(in) nInflowGauges,
    integer(i4), dimension(:), intent(in) InflowGaugeIndexList,
    logical, dimension(:), intent(in) InflowGaugeHeadwater,
    integer(i4), dimension(:), intent(in) InflowGaugeNodeList,
    real(dp), dimension(:), intent(in) InflowDischarge,
    integer(i4), intent(in) nGauges,
    integer(i4), dimension(:), intent(in) gaugeIndexList,
    integer(i4), dimension(:), intent(in) gaugeNodeList,
    logical, intent(in) map_flag,
    real(dp), dimension(:), intent(in) L11_length,
    real(dp), dimension(:), intent(in) L11_slope,
    real(dp), dimension(:), intent(in) L11_FracFPimp,
    real(dp), dimension(:), intent(inout) L11_C1,
    real(dp), dimension(:), intent(inout) L11_C2,
    real(dp), dimension(:), intent(inout) L11_qOut,
    real(dp), dimension(:, :), intent(inout) L11_qTIN,
    real(dp), dimension(:, :), intent(inout) L11_qTR,
    real(dp), dimension(:), intent(inout) L11_qMod,
    real(dp), dimension(:), intent(inout) GaugeDischarge )

```

route water given runoff

This routine first performs mpr for the routing variables if required, then accumulates the runoff to the routing resolution and eventually routes the water in a third step. The last step is repeated multiple times if the routing timestep is smaller than the timestep of the hydrological timestep

## Parameters

|    |  |  |
|----|--|--|
| in | <i>logical</i> :: <i>read_states</i>                         | whether states are derived from restart file |
| in | <i>integer(i4)</i> :: <i>processCase</i>                     | Process switch for routing                   |
| in | <i>real(dp), dimension(:)</i> :: <i>global_routing_param</i> | routing parameters                           |
| in | <i>real(dp), dimension(:)</i> :: <i>L1_total_runoff</i>      | total runoff from L1 grid cells              |
| in | <i>real(dp), dimension(:)</i> :: <i>L1_areaCell</i>          | L1 cell area                                 |
| in | <i>integer(i4), dimension(:)</i> :: <i>L1_L11_Id</i>         | L1 cell ids on L11                           |
| in | <i>real(dp), dimension(:)</i> :: <i>L11_areaCell</i>         | L11 cell area                                |
| in | <i>integer(i4), dimension(:)</i> :: <i>L11_L1_Id</i>         | L11 cell ids on L1                           |
| in | <i>integer(i4), dimension(:)</i> :: <i>L11_netPerm</i>       | L11 routing order                            |
| in | <i>integer(i4), dimension(:)</i> :: <i>L11_fromN</i>         | L11 source grid cell order                   |
| in | <i>integer(i4), dimension(:)</i> :: <i>L11_toN</i>           | L11 target grid cell order                   |

## Parameters

|         |  |  |
|---------|--|--|
| in      | <i>integer(i4) :: L11_nOutlets</i>                       | L11 number of outlets/sinks  |
| in      | <i>integer(i4) :: timestep</i>                           | simulation timestep in [h]   |
| in      | <i>real(dp) :: tsRoutFactor</i>                          | factor between routing timestep and hydrological timestep                        |
| in      | <i>integer(i4) :: nNodes</i>                             | number of nodes  |
| in      | <i>integer(i4) :: nInflowGauges</i>                      | number of inflow gauges  |
| in      | <i>integer(i4), dimension(:) :: InflowGaugeIndexList</i> | index list of inflow gauges  |
| in      | <i>logical, dimension(:) :: InflowGaugeHeadwater</i>     | flag for headwater cell of inflow gauge  |
| in      | <i>integer(i4), dimension(:) :: InflowGaugeNodeList</i>  | gauge node list at L11   |
| in      | <i>real(dp), dimension(:) :: InflowDischarge</i>         | inflowing discharge at discharge gauge at current day                            |
| in      | <i>integer(i4) :: nGauges</i>                            | number of recording gauges   |
| in      | <i>integer(i4), dimension(:) :: gaugeIndexList</i>       | index list for outflow gauges  |
| in      | <i>integer(i4), dimension(:) :: gaugeNodeList</i>        | gauge node list at L11   |
| in      | <i>logical :: map_flag</i>                               | flag indicating whether routing resolution is coarser than hydrologic resolution |
| in      | <i>real(dp), dimension(:) :: L11_length</i>              | L11 link length  |
| in      | <i>real(dp), dimension(:) :: L11_slope</i>               | L11 slope  |
| in      | <i>real(dp), dimension(:) :: L11_FracFPimp</i>           | L11 fraction of flood plain with impervious cover                                |
| in, out | <i>real(dp), dimension(:) :: L11_C1</i>                  | L11 muskingum parameter 1  |
| in, out | <i>real(dp), dimension(:) :: L11_C2</i>                  | L11 muskingum parameter 2  |
| in, out | <i>real(dp), dimension(:) :: L11_qOut</i>                | total runoff from L11 grid cells   |
| in, out | <i>real(dp), dimension(:, :) :: L11_qTIN</i>             | L11 inflow to the reach  |
| in, out | <i>real(dp), dimension(:, :) :: L11_qTR</i>              | L11 routed outflow   |
| in, out | <i>real(dp), dimension(:) :: L11_qMod</i>                | modelled discharge at each grid cell   |
| in, out | <i>real(dp), dimension(:) :: GaugeDischarge</i>          | modelled discharge at each gauge   |

## Authors

Stephan Thober



## Date

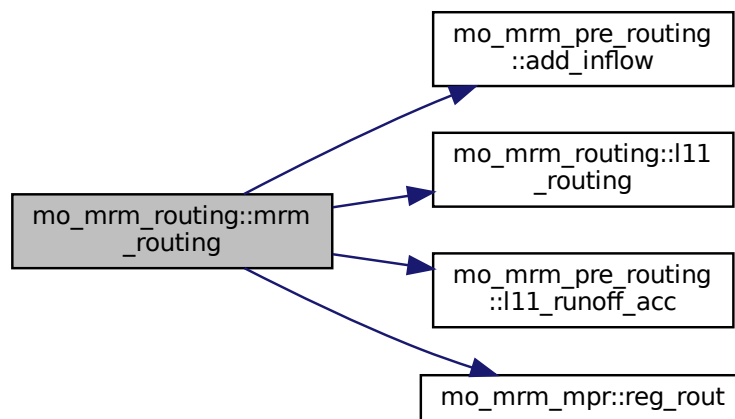
Aug 2015

Definition at line 103 of file mo\_mrm\_routing.f90.

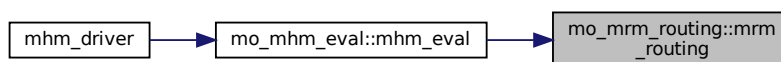
References mo\_mrm\_pre\_routing::add\_inflow(), mo\_mrm\_global\_variables::is\_start, l11\_routing(), mo\_mrm\_pre\_routing::l11\_runoff\_acc(), mo\_mrm\_mpr::reg\_rout(), and mo\_mrm\_global\_variables::riv\_temp\_pcs.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.51 mo\_mrm\_signatures Module Reference

Module with calculations for several hydrological signatures.

### Functions/Subroutines

- real(dp) function, dimension(size(lags, 1)), public [autocorrelation](#) (data, lags, mask)  
*Autocorrelation of a given data series.*
- real(dp) function, dimension(size(quantiles, 1)), public [flowdurationcurve](#) (data, quantiles, mask, concavity\_index, mid\_segment\_slope, mhigh\_segment\_volume, high\_segment\_volume, low\_segment\_volume)  
*Flow duration curves.*
- subroutine, public [limb\\_densities](#) (data, mask, RLD, DLD)

*Calculates limb densities.*

- real(dp) function [maximummonthlyflow](#) (data, mask, yr\_start, mo\_start, dy\_start)

*Maximum of average flows per months.*

- subroutine, public [moments](#) (data, mask, mean\_data, stddev\_data, median\_data, max\_data, mean\_log, stddev\_log, median\_log, max\_log)

*Moments of data and log-transformed data, e.g. mean and standard deviation.*

- real(dp) function, dimension(size(quantiles, 1)), public [peakdistribution](#) (data, quantiles, mask, slope\_peak←\_distribution)

*Calculates the peak distribution.*

- real(dp) function, public [runoffratio](#) (data, domain\_area, mask, precip\_series, precip\_sum, log\_data)

*Runoff ratio (accumulated daily discharge [mm/d] / accumulated daily precipitation [mm/d]).*

- real(dp) function, public [zeroflowratio](#) (data, mask)

*Ratio of zero values to total number of data points.*

### 17.51.1 Detailed Description

Module with calculations for several hydrological signatures.

This module contains calculations for hydrological signatures. It contains:

- Autocorrelation
- Rising and declining limb densities
- Flow duration curves
- Peak distribution

#### Authors

Remko Nijzink,

#### Date

March 2014

### 17.51.2 Function/Subroutine Documentation

#### 17.51.2.1 autocorrelation()

```
real(dp) function, dimension(size(lags, 1)), public mo_mrm_signatures::autocorrelation (
    real(dp), dimension(:), intent(in) data,
    integer(i4), dimension(:), intent(in) lags,
    logical, dimension(size(data, 1)), intent(in), optional mask )
```

Autocorrelation of a given data series.

Calculates the autocorrelation of a data series at given lags. An optional argument for masking data points can be given. The function is basically a wrapper of the function autocorr from the module mo\_corr. An optional mask of data points can be specified. ADDITIONAL INFORMATION Used as hydrologic signature with lag 1 in Euser, T., Winsemius, H. C., Hrachowitz, M., Fenicia, F., Uhlenbrook, S., & Savenije, H. H. G. (2013). A framework to assess the realism of model structures using hydrological signatures. Hydrology and Earth System Sciences, 17(5), 1893-1912. doi:10.5194/hess-17-1893-2013

#### Parameters

|    |                                       |               |
|----|---------------------------------------|---------------|
| in | <i>real(dp), dimension(:) :: data</i> | Array of data |
|----|---------------------------------------|---------------|

## Parameters

|    |  |   |
|----|--|---|
| in | <i>integer(i4), dimension(:) :: lags</i>                   | Array of lags where autocorrelation is requested                          |
| in | <i>logical, dimension(size(data, 1)), optional :: mask</i> | Mask for data points givenWorks only with 1d double precision input data. |

## Authors

Juliane Mai

## Date

Jun 2015

Definition at line 71 of file mo\_mrm\_signatures.f90.

## 17.51.2.2 flowdurationcurve()

```

real(dp) function, dimension(size(quantiles, 1)), public mo_mrm_signatures::flowdurationcurve
(
    real(dp), dimension(:), intent(in) data,
    real(dp), dimension(:), intent(in) quantiles,
    logical, dimension(:), intent(in), optional mask,
    real(dp), intent(out), optional concavity_index,
    real(dp), intent(out), optional mid_segment_slope,
    real(dp), intent(out), optional mhigh_segment_volume,
    real(dp), intent(out), optional high_segment_volume,
    real(dp), intent(out), optional low_segment_volume )

```

Flow duration curves.

Calculates the flow duration curves for a given data vector. The Flow duration curve at a certain quantile  $x$  is the data point  $p$  where  $x\%$  of the data points are above the value  $p$ . Hence the function percentile of the module `mo_mrm_signatures::percentile` is used. But percentile is determining the point  $p$  where  $x\%$  of the data points are below that value. Therefore, the given quantiles are transformed by  $(1.0 - \text{quantile})$  to get the percentiles of exceedance probabilities. Optionally, the concavity index  $CI$  can be calculated [Zhang2014].  $CI$  is defined by

$$CI = \frac{q_{10\%} - q_{99\%}}{q_{1\%} - q_{99\%}}$$

where  $q_x$  is the data point where  $x\%$  of the data points are above that value. Hence, exceedance probabilities are used. Optionally, the FDC mid-segment slope  $FDC_{MSS}$  as used by Shafii et. al (2014) can be returned. The  $FDC_{MSS}$  is defined as

$$FDC_{MSS} = \log(q_{m_1}) - \log(q_{m_2})$$

where  $m_1$  and  $m_2$  are the lowest and highest flow exceedance probabilities within the midsegment of FDC. The settings  $m_1 = 0.2$  and  $0.7$  are used by Shafii et. al (2014) and are implemented like that. Optionally, the FDC medium high-segment volume  $FDC_{MHSV}$  as used by Shafii et. al (2014) can be returned. The  $FDC_{MHSV}$  is defined as

$$FDC_{MHSV} = \sum_{h=1}^H q_h$$

where  $h = 1, 2, \dots, H$  are flow indices located within the high-flow segment (exceedance probabilities lower than  $m_1$ ).  $H$  is the index of the maximum flow. The settings  $m_1 = 0.2$  is used here to be consistent with the definitions of the low-segment (0.7-1.0) and the mid-segment (0.2-0.7). Optionally, the FDC high-segment volume  $FDC_{HSV}$  as used by Shafii et. al (2014) can be returned. The  $FDC_{HSV}$  is defined as

$$FDC_{HSV} = \sum_{h=1}^H q_h$$

where  $h = 1, 2, \dots, H$  are flow indices located within the high-flow segment (exceedance probabilities lower than  $m_1$ ).  $H$  is the index of the maximum flow. The settings  $m_1 = 0.02$  is used by Shafii et. al (2014) and is implemented like that. Optionally, the FDC low-segment volume  $FDC_{LSV}$  as used by Shafii et. al (2014) can be returned. The  $FDC_{LSV}$  is defined as

$$FDC_{LSV} = - \sum_{l=1}^L (\log(q_l) - \log(q_L))$$

where  $l = 1, 2, \dots, L$  are flow indices located within the low-flow segment (exceedance probabilities larger than  $m_1$ ).  $L$  is the index of the minimum flow. The settings  $m_1 = 0.7$  is used by Shafii et. al (2014) and is implemented like that. An optional mask of data points can be specified. ADDITIONAL INFORMATION Thresholds in `mid_segment_slope`, `mhigh_segment_volume`, `high_segment_volume`, `low_segment_volume` are hard coded. FDC is used as hydrologic signature (quantiles not specified) in Euser, T., Winsemius, H. C., Hrachowitz, M., Fenicia, F., Uhlenbrook, S., & Savenije, H. H. G. (2013). A framework to assess the realism of model structures using hydrological signatures. *Hydrology and Earth System Sciences*, 17(5), 1893-1912. doi:10.5194/hess-17-1893-2013 Concavity Index used as hydrologic signature in Zhang, Y., Vaze, J., Chiew, F. H. S., Teng, J., & Li, M. (2014). Predicting hydrological signatures in ungauged catchments using spatial interpolation, index model, and rainfall-runoff modelling. *Journal of Hydrology*, 517(C), 936-948. doi:10.1016/j.jhydrol.2014.06.032 Concavity index is defined using exceedance probabilities by Sauquet, E., & Catalogne, C. (2011). Comparison of catchment grouping methods for flow duration curve estimation at ungauged sites in France. *Hydrology and Earth System Sciences*, 15(8), 2421-2435. doi:10.5194/hess-15-2421-2011 `mid_segment_slope`, `high_segment_volume`, `low_segment_volume` used as hydrologic signature in Shafii, M., & Tolson, B. A. (2015). Optimizing hydrological consistency by incorporating hydrological signatures into model calibration objectives. *Water Resources Research*, 51(5), 3796-3814. doi:10.1002/2014WR016520

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>real(dp), dimension(:) :: data</i>             | data series  |
| in  | <i>real(dp), dimension(:) :: quantiles</i>        | Percentages of exceedance (x-axis of FDC)              |
| in  | <i>logical, dimension(:), optional :: mask</i>    | mask of data array                                     |
| out | <i>real(dp), optional :: concavity_index</i>      | concavity index as defined by Sauquet et al. (2011)    |
| out | <i>real(dp), optional :: mid_segment_slope</i>    | mid-segment slope as defined by Shafii et al. (2014)   |
| out | <i>real(dp), optional :: mhigh_segment_volume</i> | medium high-segment volume                             |
| out | <i>real(dp), optional :: high_segment_volume</i>  | high-segment volume as defined by Shafii et al. (2014) |
| out | <i>real(dp), optional :: low_segment_volume</i>   | low-segment volume as defined by Shafii et al. (2014)  |

#### Returns

*real(dp), dimension(size(quantiles,1)) :: FlowDurationCurve* — Flow Duration Curve value at resp. quantile

#### Authors

Remko Nijzink, Juliane Mai

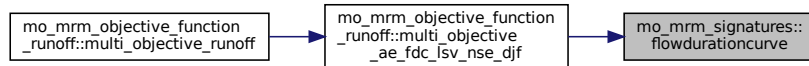
**Date**

March 2014

Definition at line 201 of file mo\_mrm\_signatures.f90.

Referenced by mo\_mrm\_objective\_function\_runoff::multi\_objective\_ae\_fdc\_lsv\_nse\_djf().

Here is the caller graph for this function:

**17.51.2.3 limb\_densities()**

```

subroutine, public mo_mrm_signatures::limb_densities (
    real(dp), dimension(:), intent(in) data,
    logical, dimension(size(data, 1)), intent(in), optional mask,
    real(dp), intent(out), optional RLD,
    real(dp), intent(out), optional DLD )
  
```

Calculates limb densities.

Calculates rising and declining limb densities. The peaks of the given series are first determined by looking for points where preceding and subsequent datapoint are lower. Second, the number of datapoints with rising values (nrise) and declining values (ndecline) are counted basically by comparing neighbors. The duration the data increase (nrise) divided by the number of peaks (npeaks) gives the rising limb density RLD

$$RLD = t_{rise} / n_{peak}$$

whereas the duration the data decrease (ndecline) divided by the number of peaks (npeaks) gives the declining limb density DLD

$$DLD = t_{fall} / n_{peak}.$$

An optional mask of data points can be specified. ADDITIONAL INFORMATION Rising limb density used as hydrologic signature in Euser, T., Winsemius, H. C., Hrachowitz, M., Fenicia, F., Uhlenbrook, S., & Savenije, H. H. G. (2013). A framework to assess the realism of model structures using hydrological signatures. Hydrology and Earth System Sciences, 17(5), 1893-1912. doi:10.5194/hess-17-1893-2013

**Parameters**

|     |   |                        |
|-----|---|------------------------|
| in  | real(dp), dimension(:) :: data                      | data series            |
| in  | logical, dimension(size(data, 1)), optional :: mask | mask for data series   |
| out | real(dp), optional :: RLD                           | rising limb density    |
| out | real(dp), optional :: DLD                           | declining limb density |

**Authors**

Remko Nijzink

**Date**

March 2014

Definition at line 341 of file mo\_mrm\_signatures.f90.

**17.51.2.4 maximummonthlyflow()**

```
real(dp) function mo_mrm_signatures::maximummonthlyflow (
    real(dp), dimension(:), intent(in) data,
    logical, dimension(size(data, 1)), intent(in), optional mask,
    integer(i4), intent(in), optional yr_start,
    integer(i4), intent(in), optional mo_start,
    integer(i4), intent(in), optional dy_start )
```

Maximum of average flows per months.

Maximum of average flow per month is defined as

$$max_{monthlyflow} = Max(F(i), i = 1, ..12)$$

where  $F(i)$  is the average flow of month  $i$ . ADDITIONAL INFORMATION used as hydrologic signature in Shafii, M., & Tolson, B. A. (2015). Optimizing hydrological consistency by incorporating hydrological signatures into model calibration objectives. *Water Resources Research*, 51(5), 3796-3814. doi:10.1002/2014WR016520

**Parameters**

|    |  |  |
|----|--|--|
| in | <i>real(dp), dimension(:) :: data</i>                      | array of data  |
| in | <i>logical, dimension(size(data, 1)), optional :: mask</i> | mask for data points given                           |
| in | <i>integer(i4), optional :: yr_start</i>                   | year of date of first data point given               |
| in | <i>integer(i4), optional :: mo_start</i>                   | month of date of first data point given (default: 1) |
| in | <i>integer(i4), optional :: dy_start</i>                   | month of date of first data point given (default: 1) |

**Returns**

*real(dp) :: MaximumMonthlyFlow* — Maximum of average flow per month Works only with 1d double precision input data. Assumes data are daily values.

**Authors**

Juliane Mai

**Date**

Jun 2015

Definition at line 513 of file mo\_mrm\_signatures.f90.

**17.51.2.5 moments()**

```
subroutine, public mo_mrm_signatures::moments (
    real(dp), dimension(:), intent(in) data,
    logical, dimension(size(data, 1)), intent(in), optional mask,
```

```

real(dp), intent(out), optional mean_data,
real(dp), intent(out), optional stddev_data,
real(dp), intent(out), optional median_data,
real(dp), intent(out), optional max_data,
real(dp), intent(out), optional mean_log,
real(dp), intent(out), optional stddev_log,
real(dp), intent(out), optional median_log,
real(dp), intent(out), optional max_log )

```

Moments of data and log-transformed data, e.g. mean and standard deviation.

Returns several moments of data series given, i.e.

- mean of data
- standard deviation of data
- median of data
- maximum/ peak of data
- mean of log-transformed data
- standard deviation of log-transformed data
- median of log-transformed data
- maximum/ peak of log-transformed data An optional mask of data points can be specified. ADDITIONAL INFORMATION mean\_log and stddev\_log used as hydrologic signature in Zhang, Y., Vaze, J., Chiew, F. H. S., Teng, J., & Li, M. (2014). Predicting hydrological signatures in ungauged catchments using spatial interpolation, index model, and rainfall-runoff modelling. *Journal of Hydrology*, 517(C), 936-948. doi:10.1016/j.jhydrol.2014.06.032 mean\_data, stddev\_data, median\_data, max\_data, mean\_log, and stddev\_log used as hydrologic signature in Shafii, M., & Tolson, B. A. (2015). Optimizing hydrological consistency by incorporating hydrological signatures into model calibration objectives. *Water Resources Research*, 51(5), 3796-3814. doi:10.1002/2014WR016520

#### Parameters

|     |  |  |
|-----|--|--|
| in  | <i>real(dp), dimension(:) :: data</i>                      | array of data  |
| in  | <i>logical, dimension(size(data, 1)), optional :: mask</i> | mask for data points given   |
| out | <i>real(dp), optional :: mean_data</i>                     | mean of data   |
| out | <i>real(dp), optional :: stddev_data</i>                   | standard deviation of data   |
| out | <i>real(dp), optional :: median_data</i>                   | median of data   |
| out | <i>real(dp), optional :: max_data</i>                      | maximum/ peak of data  |
| out | <i>real(dp), optional :: mean_log</i>                      | mean of log-transformed data   |
| out | <i>real(dp), optional :: stddev_log</i>                    | standard deviation of log-transformed data   |
| out | <i>real(dp), optional :: median_log</i>                    | median of log-transformed data   |
| out | <i>real(dp), optional :: max_log</i>                       | maximum/ peak of log-transformed dataWorks only with 1d double precision input data. |

#### Authors

Juliane Mai

#### Date

Jun 2015

Definition at line 660 of file mo\_mrm\_signatures.f90.

### 17.51.2.6 peakdistribution()

```
real(dp) function, dimension(size(quantiles, 1)), public mo_mrm_signatures::peakdistribution (
    real(dp), dimension(:), intent(in) data,
    real(dp), dimension(:), intent(in) quantiles,
    logical, dimension(size(data, 1)), intent(in), optional mask,
    real(dp), intent(out), optional slope_peak_distribution )
```

Calculates the peak distribution.

First, the peaks of the time series given are identified. For the peak distribution only this subset of data points are considered. Second, the peak distribution at the quantiles given is calculated. Calculates the peak distribution at the quantiles given using `mo_percentile`. Since the exceedance probabilities are usually used in hydrology the function percentile is used with (1.0-quantiles). Optionally, the slope of the peak distribution between 10th and 50th percentile, i.e.

$$slope = \frac{peak\_data_{0.1} - peak\_data_{0.5}}{0.9 - 0.5}$$

can be returned. An optional mask for the data points can be given. ADDITIONAL INFORMATION `slope_peak_↔distribution` used as hydrologic signature in Euser, T., Winsemius, H. C., Hrachowitz, M., Fenicia, F., Uhlenbrook, S., & Savenije, H. H. G. (2013). A framework to assess the realism of model structures using hydrological signatures. *Hydrology and Earth System Sciences*, 17(5), 1893-1912. doi:10.5194/hess-17-1893-2013

#### Parameters

|     |  |   |
|-----|--|---|
| in  | <i>real(dp), dimension(:) :: data</i>                      | data array  |
| in  | <i>real(dp), dimension(:) :: quantiles</i>                 | requested quantiles for distribution                            |
| in  | <i>logical, dimension(size(data, 1)), optional :: mask</i> | mask of data array  |
| out | <i>real(dp), optional :: slope_peak_distribution</i>       | slope of the Peak distribution between 10th and 50th percentile |

#### Returns

`real(dp), dimension(size(quantiles,1)) :: PeakDistribution` — Distribution of peak values at resp. quantiles

#### Authors

Remko Nijzink

#### Date

March 2014

Definition at line 788 of file `mo_mrm_signatures.f90`.

### 17.51.2.7 runoffratio()

```
real(dp) function, public mo_mrm_signatures::runoffratio (
    real(dp), dimension(:), intent(in) data,
    real(dp), intent(in) domain_area,
    logical, dimension(size(data, 1)), intent(in), optional mask,
    real(dp), dimension(size(data, 1)), intent(in), optional precip_series,
    real(dp), intent(in), optional precip_sum,
    logical, intent(in), optional log_data )
```

Runoff ratio (accumulated daily discharge [mm/d] / accumulated daily precipitation [mm/d]).



The runoff ratio is defined as

$$runoff\_ratio = \frac{\sum_{t=1}^N q_t}{\sum_{t=1}^N p_t}$$

where  $p_t$  and  $q_t$  are precipitation and discharge, respectively. Therefore, precipitation over the entire domain is required and both discharge and precipitation have to be converted to the same units [mm/d]. Input discharge is given in [m\*\*3/s] as this is mHM default while precipitation has to be given in [mm/km\*\*2 / day]. Either "precip\_sum" or "precip\_series" has to be specified. If "precip\_series" is used the optional mask is also applied to precipitation values. The "precip\_sum" is the accumulated "precip\_series". Optionally, a mask for the data (=discharge) can be given. If optional "log\_data" is set to .true. the runoff ratio will be calculated as

$$runoff\_ratio = \frac{\sum_{t=1}^N \log(q_t)}{\sum_{t=1}^N p_t}$$

where  $p_t$  and  $q_t$  are precipitation and discharge, respectively. ADDITIONAL INFORMATION

### Returns

real(dp), dimension(size(lags,1)) :: RunoffRatio — Ratio of discharge and precipitation Used as hydrologic signature in Shafii, M., & Tolson, B. A. (2015). Optimizing hydrological consistency by incorporating hydrological signatures into model calibration objectives. Water Resources Research, 51(5), 3796-3814. doi:10.1002/2014WR016520

### Parameters

|    |  |  |
|----|--|--|
| in | <i>real(dp), dimension(:) :: data</i>                                | array of data [m**3/s]   |
| in | <i>real(dp) :: domain_area</i>                                       | area of domain [km**2]   |
| in | <i>logical, dimension(size(data, 1)), optional :: mask</i>           | mask for data points given   |
| in | <i>real(dp), dimension(size(data, 1)), optional :: precip_series</i> | daily precipitation values [mm/km**2 / day]                                  |
| in | <i>real(dp), optional :: precip_sum</i>                              | sum of daily precip. values of whole period [mm/km**2 / day]                 |
| in | <i>logical, optional :: log_data</i>                                 | ratio using logarithmic data Works only with 1d double precision input data. |

### Authors

Juliane Mai

### Date

Jun 2015

Definition at line 925 of file mo\_mrm\_signatures.f90.

#### 17.51.2.8 zeroflowratio()

```
real(dp) function, public mo_mrm_signatures::zeroflowratio (
    real(dp), dimension(:), intent(in) data,
    logical, dimension(size(data, 1)), intent(in), optional mask )
```

Ratio of zero values to total number of data points.

An optional mask of data points can be specified. ADDITIONAL INFORMATION

**Returns**

`real(dp), dimension(size(lags,1)) :: ZeroFlowRatio` — Ratio of zero values to total number of data points Used as hydrologic signature in Zhang, Y., Vaze, J., Chiew, F. H. S., Teng, J., & Li, M. (2014). Predicting hydrological signatures in ungauged catchments using spatial interpolation, index model, and rainfall-runoff modelling. *Journal of Hydrology*, 517(C), 936-948. doi:10.1016/j.jhydrol.2014.06.032

**Parameters**

|                 |  |   |
|-----------------|--|---|
| <code>in</code> | <code>real(dp), dimension(:) :: data</code>                      | array of data   |
| <code>in</code> | <code>logical, dimension(size(data, 1)), optional :: mask</code> | mask for data points givenWorks only with 1d double precision input data. |

**Authors**

Juliane Mai

**Date**

Jun 2015

Definition at line 1045 of file `mo_mrm_signatures.f90`.

## 17.52 mo\_mrm\_write Module Reference

write of discharge and restart files

**Functions/Subroutines**

- subroutine, public `mrm_write`  
*write discharge and restart files*
- subroutine `write_configfile`  
*This modules writes the results of the configuration into an ASCII-file.*
- subroutine `write_daily_obs_sim_discharge` (Qobs, Qsim)  
*Write a file for the daily observed and simulated discharge timeseries during the evaluation period for each gauging station.*
- subroutine, public `mrm_write_output_fluxes` (iDomain, nCells, timeStep\_model\_outputs, domainDateTime, tt, timestep, mask11, L11\_qmod)  
*write fluxes to netcdf output files*
- subroutine, public `mrm_write_optifile` (best\_OF, best\_paramSet, param\_names)  
*Write briefly final optimization results.*
- subroutine, public `mrm_write_optinamelist` (parameters, maskpara, parameters\_name)  
*Write final, optimized parameter set in a namelist format.*

**Variables**

- type(`outputdataset`) `nc`

## 17.52.1 Detailed Description

write of discharge and restart files

This module contains the subroutines for writing the discharge files and optionally the restart files.

### Authors

Stephan Thober

### Date

Aug 2015

## 17.52.2 Function/Subroutine Documentation

### 17.52.2.1 mrm\_write()

```
subroutine, public mo_mrm_write::mrm_write
```

write discharge and restart files

First, this subroutine calls the writing of restart files that only succeeds if it happens after the write of mHM restart files because mHM restart files must exist. Second, simulated discharge is aggregated to the daily scale and then written to file jointly with observed discharge

### Authors

Juliane Mai, Rohini Kumar & Stephan Thober

### Date

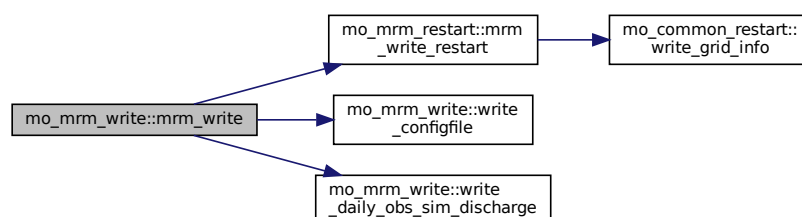
Aug 2015

Definition at line 54 of file mo\_mrm\_write.f90.

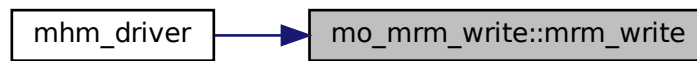
References mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_variables::evalper, mo\_mrm\_global\_variables::gauge, mo\_common\_mhm\_mrm\_variables::mrm\_coupling\_mode, mo\_mrm\_global\_variables::mrm\_runoff, mo\_mrm\_restart::mrm\_write\_restart(), mo\_common\_variables::mrmfilerestartout, mo\_mrm\_global\_variables::ngaugestotal, mo\_common\_mhm\_mrm\_variables::ntstepday, mo\_common\_mhm\_mrm\_variables::simper, mo\_common\_mhm\_mrm\_variables::warmingdays, write\_configfile(), write\_daily\_obs\_sim\_discharge(), and mo\_common\_variables::write\_restart.

Referenced by mhm\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.52.2.2 mrm\_write\_optifile()

```

subroutine, public mo_mrm_write::mrm_write_optifile (
    real(dp), intent(in) best_OF,
    real(dp), dimension(:), intent(in) best_paramSet,
    character(len = *), dimension(:), intent(in) param_names )
  
```

Write briefly final optimization results.

Write overall best objective function and the best optimized parameter set to a file\_opti.

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>real(dp) :: best_OF</i>                             | best objective function value as returned by the optimization routine    |
| in | <i>real(dp), dimension(:) :: best_paramSet</i>         | best associated global parameter set Called only when optimize is .TRUE. |
| in | <i>character(len = *), dimension(:) :: param_names</i> |  |

#### Authors

David Schaefer

#### Date

July 2013

Definition at line 753 of file mo\_mrm\_write.f90.

References mo\_common\_variables::dirconfigout, mo\_common\_mhm\_mrm\_file::file\_opti, and mo\_common\_mhm\_mrm\_file::uopti.

### 17.52.2.3 mrm\_write\_optinamelist()

```

subroutine, public mo_mrm_write::mrm_write_optinamelist (
    real(dp), dimension(:, :), intent(in) parameters,
    logical, dimension(size(parameters, 1)), intent(in) maskpara,
    character(len = *), dimension(size(parameters, 1)), intent(in) parameters_name )
  
```

Write final, optimized parameter set in a namelist format.

Write final, optimized parameter set in a namelist format.

## Parameters

|    |  |                                    |
|----|--|------------------------------------|
| in | <i>real(dp), dimension(:, :) :: parameters</i>                               | (min, max, opti)                   |
| in | <i>logical, dimension(size(parameters, 1)) :: maskpara</i>                   | .true. if parameter was calibrated |
| in | <i>character(len = *), dimension(size(parameters, 1)) :: parameters_name</i> | clear names of parameters          |

## Authors

Juliane Mai

## Date

Dec 2013

Definition at line 832 of file mo\_mrm\_write.f90.

References mo\_common\_variables::dirconfigout, mo\_common\_mhm\_mrm\_file::file\_opti\_nml, mo\_common\_variables::processmatrix, and mo\_common\_mhm\_mrm\_file::uopti\_nml.

## 17.52.2.4 mrm\_write\_output\_fluxes()

```
subroutine, public mo_mrm_write::mrm_write_output_fluxes (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) nCells,
    integer(i4), intent(in) timeStep_model_outputs,
    type(datetimeinfo), intent(in) domainDateTime,
    integer(i4), intent(in) tt,
    integer(i4), intent(in) timestep,
    logical, dimension(:, :), intent(in), pointer mask11,
    real(dp), dimension(:), intent(in) L11_qmod )
```

write fluxes to netcdf output files

This subroutine creates a netcdf data set for writing L11\_QTIN for different time averages.

## Parameters

|    |  |                               |
|----|--|-------------------------------|
| in | <i>integer(i4) :: iDomain</i>                |                               |
| in | <i>integer(i4) :: nCells</i>                 |                               |
| in | <i>integer(i4) :: timeStep_model_outputs</i> | timestep of model outputs     |
| in | <i>integer(i4) :: warmingDays</i>            | number of warming days        |
| in | <i>real(dp) :: newTime</i>                   | julian date of next time step |
| in | <i>integer(i4) :: nTimeSteps</i>             | number of total timesteps     |
| in | <i>integer(i4) :: nTStepDay</i>              | number of timesteps per day   |
| in | <i>integer(i4) :: tt</i>                     | current model timestep        |
| in | <i>integer(i4) :: day</i>                    | current day of the year       |
| in | <i>integer(i4) :: month</i>                  | current month of the year     |
| in | <i>integer(i4) :: year</i>                   | current year                  |
| in | <i>integer(i4) :: timestep</i>               | current model time resolution |
| in | <i>logical, dimension(:, :) :: mask11</i>    | mask at level 11              |
| in | <i>real(dp), dimension(:) :: L11_qMod</i>    | current routed streamflow     |

**Authors**

Stephan Thober

**Date**

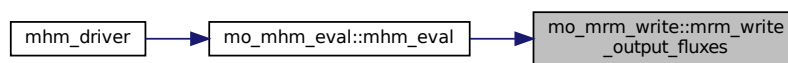
Aug 2015

Definition at line 666 of file mo\_mrm\_write.f90.

References nc, and mo\_mrm\_global\_variables::riv\_temp\_pcs.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the caller graph for this function:

**17.52.2.5 write\_configfile()**

```
subroutine mo_mrm_write::write_configfile
```

This modules writes the results of the configuration into an ASCII-file.

TODO: add description

**Authors**

Christoph Schneider

**Date**

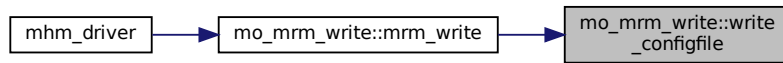
May 2013

Definition at line 156 of file mo\_mrm\_write.f90.

References mo\_common\_variables::dirconfigout, mo\_mrm\_global\_variables::dirgauges, mo\_common\_variables::dircover, mo\_common\_variables::dirmorpho, mo\_common\_variables::dirout, mo\_mrm\_global\_variables::dirtotalrunoff, mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_variables::evalper, mo\_common\_file::file\_config, mo\_mrm\_global\_variables::gauge, mo\_common\_variables::global\_parameters, mo\_common\_variables::global\_parameters\_name, mo\_mrm\_global\_variables::inflowgauge, mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_global\_variables::l11\_l1\_id, mo\_mrm\_global\_variables::l11\_label, mo\_mrm\_global\_variables::l11\_length, mo\_mrm\_global\_variables::l11\_netperm, mo\_mrm\_global\_variables::l11\_rorder, mo\_mrm\_global\_variables::l11\_slope, mo\_mrm\_global\_variables::l11\_ton, mo\_mrm\_global\_variables::l1\_l1\_id, mo\_common\_variables::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::lcfilename, mo\_common\_mhm\_mrm\_variables::lcyyearid, mo\_common\_variables::level0, mo\_common\_variables::level1, mo\_mrm\_global\_variables::level11, mo\_common\_mhm\_mrm\_variables::mrm\_coupling\_mode, mo\_common\_variables::mrmfilerestartout, mo\_mrm\_global\_variables::ngaugestotal, mo\_mrm\_global\_variables::ninflowgaugestotal, mo\_common\_variables::nlcoverscene, mo\_common\_constants::nodata\_dp, mo\_common\_variables::processmatrix, mo\_common\_mhm\_mrm\_variables::read\_restart, mo\_common\_variables::resolutionhydrology, mo\_common\_mhm\_mrm\_variables::resolutionrouting, mo\_common\_mhm\_mrm\_variables::simper, mo\_common\_mhm\_mrm\_variables::timestep, mo\_common\_file::uconfig, mo\_mrm\_file::version, mo\_common\_mhm\_mrm\_variables::warmpcr, and mo\_common\_variables::write\_restart.

Referenced by mrm\_write().

Here is the caller graph for this function:



### 17.52.2.6 write\_daily\_obs\_sim\_discharge()

```

subroutine mo_mrm_write::write_daily_obs_sim_discharge (
    real(dp), dimension(:, :), intent(in) Qobs,
    real(dp), dimension(:, :), intent(in) Qsim )
  
```

Write a file for the daily observed and simulated discharge timeseries during the evaluation period for each gauging station.

Write a file for the daily observed and simulated discharge timeseries during the evaluation period for each gauging station

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>real(dp), dimension(:, :) :: Qobs</i> | daily time series of observed discharge<br>dims = (nModeling_days , nGauges_total) |
| in | <i>real(dp), dimension(:, :) :: Qsim</i> | daily time series of modeled discharge<br>dims = (nModeling_days , nGauges_total)  |

#### Authors

Rohini Kumar

#### Date

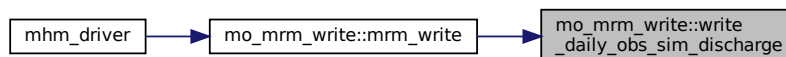
August 2013

Definition at line 476 of file mo\_mrm\_write.f90.

References mo\_common\_variables::dirout, mo\_mrm\_global\_variables::domain\_mrm, mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_variables::evalper, mo\_mrm\_file::file\_daily\_discharge, mo\_mrm\_global\_variables::gauge, mo\_mrm\_file::ncfile\_discharge, mo\_common\_constants::nodata\_dp, and mo\_mrm\_file::udaily\_discharge.

Referenced by mrm\_write().

Here is the caller graph for this function:



### 17.52.3 Variable Documentation

#### 17.52.3.1 nc

```
type(outputdataset) mo_mrm_write::nc [private]
```

Definition at line 29 of file mo\_mrm\_write.f90.

Referenced by mrm\_write\_output\_fluxes().

## 17.53 mo\_mrm\_write\_fluxes\_states Module Reference

Creates NetCDF output for different fluxes and state variables of mHM.

### Data Types

- interface [outputvariable](#)
- interface [outputdataset](#)

### Functions/Subroutines

- type([outputvariable](#)) function [newoutputvariable](#) (nc, name, dtype, dims, ncells, mask, avg)  
*Initialize OutputVariable.*
- subroutine [updatevariable](#) (self, data)  
*Update OutputVariable.*
- subroutine [writevariabletimestep](#) (self, timestep)  
*Write timestep to file.*
- type([outputdataset](#)) function [newoutputdataset](#) (iDomain, mask, nCells)  
*Initialize OutputDataset.*
- subroutine [updatedataset](#) (self, sidx, eidx, L11\_Qmod, L11\_riv\_temp)  
*Update all variables.*
- subroutine [writetimestep](#) (self, timestep)  
*Write all accumulated data.*
- subroutine [close](#) (self)  
*Close the file.*
- type(ncdataset) function [createoutputfile](#) (iDomain)  
*Create and initialize output file for X & Y coordinate system.*
- subroutine [writevariableattributes](#) (var, long\_name, unit)  
*Write output variable attributes.*

#### 17.53.1 Detailed Description

Creates NetCDF output for different fluxes and state variables of mHM.

NetCDF is first initialized and later on variables are put to the NetCDF.

#### Authors

Matthias Zink



**Date**

Apr 2013

**17.53.2 Function/Subroutine Documentation****17.53.2.1 close()**

```
subroutine mo_mrm_write_fluxes_states::close (
    class(outputdataset) self ) [private]
```

Close the file.

Close the file associated with variable of type(OutputDataset)

**Authors**

Rohini Kumar &amp; Stephan Thober

**Date**

August 2013

Definition at line 452 of file mo\_mrm\_write\_fluxes\_states.f90.

References mo\_common\_variables::dirout.

**17.53.2.2 createoutputfile()**

```
type(ncdataset) function mo_mrm_write_fluxes_states::createoutputfile (
    integer(i4), intent(in) iDomain )
```

Create and initialize output file for X &amp; Y coordinate system.

Create output file, write all non-dynamic variables and global attributes for the given domain for X &amp; Y coordinate system

**Returns**

type(NcDataset)

**Parameters**

|    |                               |              |
|----|-------------------------------|--------------|
| in | <i>integer(i4) :: iDomain</i> | -> domain id |
|----|-------------------------------|--------------|

**Authors**

David Schaefer

**Date**

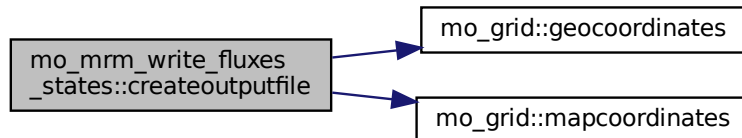
June 2015

Definition at line 494 of file mo\_mrm\_write\_fluxes\_states.f90.

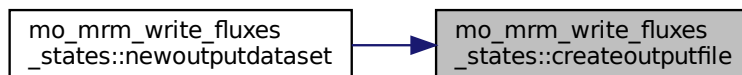
References `mo_common_variables::dirout`, `mo_common_mhm_mrm_variables::evalper`, `mo_mrm_file::file_mrm_↵  
_output`, `mo_grid::geocoordinates()`, `mo_common_variables::iflag_cordinate_sys`, `mo_mrm_global_variables_↵  
::level11`, `mo_grid::mapcoordinates()`, `mo_common_constants::nodata_dp`, and `mo_mrm_file::version`.

Referenced by `newoutputdataset()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.53.2.3 newoutputdataset()

```

type(OutputDataset) function mo_mrm_write_fluxes_states::newoutputdataset (
    integer(i4), intent(in) iDomain,
    logical, dimension(:, :), intent(in), pointer mask,
    integer(i4), intent(in) nCells ) [private]
  
```

Initialize OutputDataset.

Create and initialize the output file. If new a new output variable needs to be written, this is the first of two procedures to change (second: `updateDataset`)

#### Returns

`type(OutputDataset)`

#### Parameters

|    |   |              |
|----|---|--------------|
| in | <code>integer(i4) :: iDomain</code>           | -> domain id |
| in | <code>logical, dimension(:, :) :: mask</code> |              |
| in | <code>integer(i4) :: nCells</code>            |              |

**Authors**

Matthias Zink

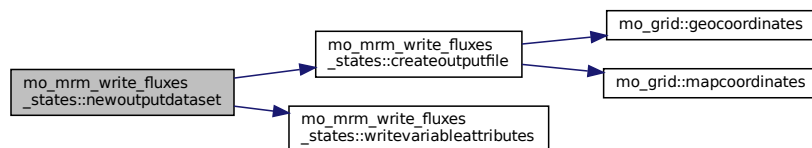
**Date**

Apr 2013

Definition at line 240 of file mo\_mrm\_write\_fluxes\_states.f90.

References `createoutputfile()`, `mo_common_variables::iflag_cordinate_sys`, `mo_mrm_global_variables::outputflxstate`, `mo_mrm`, `mo_mrm_global_variables::riv_temp_pcs`, and `writevariableattributes()`.

Here is the call graph for this function:

**17.53.2.4 newoutputvariable()**

```

type(outputvariable) function mo_mrm_write_fluxes_states::newoutputvariable (
    type(ncdataset), intent(in) nc,
    character(*), intent(in) name,
    character(*), intent(in) dtype,
    character(16), dimension(3), intent(in) dims,
    integer(i4), intent(in) ncells,
    logical, dimension(:, :), intent(in), target mask,
    logical, intent(in), optional avg ) [private]
  
```

Initialize OutputVariable.

TODO: add description

**Returns**

type(OutputVariable)

**Parameters**

|    |  |  |
|----|--|--|
| in | <i>type(NcDataset) :: nc</i>               | -> NcDataset which contains the variable |
| in | <i>character(*) :: name</i>                |  |
| in | <i>character(*) :: dtype</i>               |  |
| in | <i>character(16), dimension(3) :: dims</i> |  |
| in | <i>integer(i4) :: ncells</i>               | -> number of cells in domain             |
| in | <i>logical, dimension(:, :) :: mask</i>    |  |
| in | <i>logical, optional :: avg</i>            | -> average the data before writing       |

**Authors**

David Schaefer

**Date**

June 2015

Definition at line 100 of file mo\_mrm\_write\_fluxes\_states.f90.

References mo\_mrm\_global\_variables::output\_deflate\_level\_mrm.

**17.53.2.5 updatedataset()**

```

subroutine mo_mrm_write_fluxes_states::updatedataset (
    class(outputdataset), intent(inout), target self,
    integer(i4), intent(in) sidx,
    integer(i4), intent(in) eidx,
    real(dp), dimension(:), intent(in) L11_Qmod,
    real(dp), dimension(:), intent(in), optional L11_riv_temp )

```

Update all variables.

Call the type bound procedure updateVariable for all output variables. If a new output variable needs to be written, this is the second of two procedures to change (first: newOutputDataset)

**Parameters**

|         |   |  |
|---------|---|--|
| in, out | <i>class(OutputDataset) :: self</i>                     |  |
| in      | <i>integer(i4) :: sidx, eidx</i>                        | - start index of the domain related data in L1_* arguments |
| in      | <i>integer(i4) :: sidx, eidx</i>                        | - end index of the domain related data in L1_* arguments   |
| in      | <i>real(dp), dimension(:) :: L11_Qmod</i>               |  |
| in      | <i>real(dp), dimension(:), optional :: L11_riv_temp</i> |  |

**Authors**

Matthias Zink

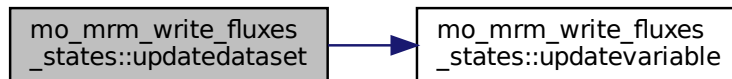
**Date**

Apr 2013

Definition at line 341 of file mo\_mrm\_write\_fluxes\_states.f90.

References mo\_mrm\_global\_variables::outputfluxstate\_mrm, and updatevariable().

Here is the call graph for this function:

**17.53.2.6 updatevariable()**

```

subroutine mo_mrm_write_fluxes_states::updatevariable (
    class(outputvariable), intent(inout) self,
    real(dp), dimension(:), intent(in) data ) [private]
  
```

Update OutputVariable.

Add the array given as actual argument to the derived type's component 'data'

**Returns**

type(OutputVariable)

**Parameters**

|         |                                       |  |
|---------|---------------------------------------|--|
| in, out | <i>class(OutputVariable) :: self</i>  |  |
| in      | <i>real(dp), dimension(:) :: data</i> |  |

**Authors**

David Schaefer

**Date**

June 2015

Definition at line 156 of file mo\_mrm\_write\_fluxes\_states.f90.

Referenced by updatedataset().

Here is the caller graph for this function:

**17.53.2.7 writetimestep()**

```

subroutine mo_mrm_write_fluxes_states::writetimestep (
    class(outputdataset), intent(inout), target self,
    integer(i4), intent(in) timestep )
  
```

Write all accumulated data.

Write all accumulated and potentially averaged data to disk.

**Returns**

type(OutputVariable)

**Parameters**

|         |                                     |                             |
|---------|-------------------------------------|-----------------------------|
| in, out | <i>class(OutputDataset) :: self</i> |                             |
| in      | <i>integer(i4) :: timestep</i>      | The model timestep to write |

**Authors**

David Schaefer

**Date**

June 2015

Definition at line 408 of file mo\_mrm\_write\_fluxes\_states.f90.

**17.53.2.8 writevariableattributes()**

```

subroutine mo_mrm_write_fluxes_states::writevariableattributes (
    type(outputvariable), intent(in) var,
  
```

```

character(*), intent(in) long_name,
character(*), intent(in) unit )

```

Write output variable attributes.

TODO: add description

#### Parameters

|    |  |                  |
|----|--|------------------|
| in | <i>type(OutputVariable) :: var</i>     |                  |
| in | <i>character(*) :: long_name, unit</i> | -> variable name |
| in | <i>character(*) :: long_name, unit</i> | -> physical unit |

#### Authors

David Schaefer

#### Date

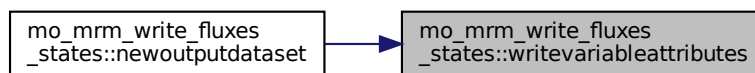
June 2015

Definition at line 649 of file mo\_mrm\_write\_fluxes\_states.f90.

References mo\_common\_constants::nodata\_dp.

Referenced by newoutputdataset().

Here is the caller graph for this function:



#### 17.53.2.9 writevariabletimestep()

```

subroutine mo_mrm_write_fluxes_states::writevariabletimestep (
    class(outputvariable), intent(inout) self,
    integer(i4), intent(in) timestep ) [private]

```

Write timestep to file.

Write the content of the derived types's component 'data' to file, average if necessary

#### Parameters

|         |                                      |  |
|---------|--------------------------------------|--|
| in, out | <i>class(OutputVariable) :: self</i> |  |
| in      | <i>integer(i4) :: timestep</i>       | -> index along the time dimension of the netcdf variable |

**Authors**

David Schafer

**Date**

June 2015

Definition at line 193 of file mo\_mrm\_write\_fluxes\_states.f90.

References mo\_common\_constants::nodata\_dp.

## 17.54 mo\_multi\_param\_reg Module Reference

Multiscale parameter regionalization (MPR).

### Functions/Subroutines

- subroutine, public [mpr](#) (mask0, geoUnit0, soilld0, Asp0, gridded\_LAI0, LCover0, slope\_emp0, y0, ld0, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, n\_subcells1, fSealed1, alpha1, degDayInc1, degDayMax1, degDayNoPre1, fAsp1, HarSamCoeff1, PrieTayAlpha1, aeroResist1, surfResist1, fRoots1, kFastFlow1, kSlowFlow1, kBaseFlow1, kPerco1, karstLoss1, soilMoistFC1, soilMoistSat1, soilMoistExp1, jarvis\_thresh\_c1, tempThresh1, unsatThresh1, sealedThresh1, wiltingPoint1, maxInter1, petLAIcorFactor, parameterset)  
*Regionalizing and Upscaling process parameters.*
- subroutine [baseflow\\_param](#) (param, geoUnit0, k2\_0)  
*baseflow recession parameter*
- subroutine [snow\\_acc\\_melt\\_param](#) (param, fForest1, flperm1, fPerm1, tempThresh1, degDayNoPre1, degDayInc1, degDayMax1)  
*Calculates the snow parameters.*
- subroutine [iper\\_thres\\_runoff](#) (param, sealedThresh1)  
*sets the impervious layer threshold parameter for runoff generation*
- subroutine [karstic\\_layer](#) (param, geoUnit0, mask0, SMs\_FC0, KsVar\_V0, ld0, n\_subcells1, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, karstLoss1, L1\_Kp)  
*calculates the Karstic percolation loss*
- subroutine, public [canopy\\_intercept\\_param](#) (processMatrix, param, LAI0, n\_subcells1, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, ld0, mask0, nodata, max\_intercept1)  
*estimate effective maximum interception capacity at L1*
- subroutine [aerodynamical\\_resistance](#) (LAI0, LCover0, param, mask0, ld0, n\_subcells1, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, aerodyn\_resistance1)  
*Regionalization of aerodynamic resistance.*

### 17.54.1 Detailed Description

Multiscale parameter regionalization (MPR).

This module provides the routines for multiscale parameter regionalization (MPR).

**Authors**

Stephan Thober, Rohini Kumar

**Date**

Dec 2012



## 17.54.2 Function/Subroutine Documentation

### 17.54.2.1 aerodynamical\_resistance()

```

subroutine mo_multi_param_reg::aerodynamical_resistance (
    real(dp), dimension(:, :), intent(in) LAI0,
    integer(i4), dimension(:), intent(in) LCover0,
    real(dp), dimension(6), intent(in) param,
    logical, dimension(:, :), intent(in) mask0,
    integer(i4), dimension(:), intent(in) Id0,
    integer(i4), dimension(:), intent(in) n_subcells1,
    integer(i4), dimension(:), intent(in) upper_bound1,
    integer(i4), dimension(:), intent(in) lower_bound1,
    integer(i4), dimension(:), intent(in) left_bound1,
    integer(i4), dimension(:), intent(in) right_bound1,
    real(dp), dimension(:, :), intent(out) aerodyn_resistance1 )

```

Regionalization of aerodynamic resistance.

estimation of aerodynamical resistance Global parameters needed (see mhm\_parameter.nml):

- param(1) = canopyheight\_forest
- param(2) = canopyheight\_impervious
- param(3) = canopyheight\_pervious
- param(4) = displacementheight\_coeff
- param(5) = roughnesslength\_momentum\_coeff
- param(6) = roughnesslength\_heat\_coeff

#### Parameters

|     |   |                                     |
|-----|---|-------------------------------------|
| in  | <i>real(dp), dimension(:, :) :: LAI0</i>                | LAI at level-0                      |
| in  | <i>integer(i4), dimension(:) :: LCover0</i>             | land cover field                    |
| in  | <i>real(dp), dimension(6) :: param</i>                  | input parameter                     |
| in  | <i>logical, dimension(:, :) :: mask0</i>                | mask at level 0                     |
| in  | <i>integer(i4), dimension(:) :: Id0</i>                 | Cell ids of hi res field            |
| in  | <i>integer(i4), dimension(:) :: n_subcells1</i>         | number of I0 cells within a I1 cell |
| in  | <i>integer(i4), dimension(:) :: upper_bound1</i>        | upper row of a I1 cell in I0 grid   |
| in  | <i>integer(i4), dimension(:) :: lower_bound1</i>        | lower row of a I1 cell in I0 grid   |
| in  | <i>integer(i4), dimension(:) :: left_bound1</i>         | left col of a I1 cell in I0 grid    |
| in  | <i>integer(i4), dimension(:) :: right_bound1</i>        | right col of a I1 cell in I0 grid   |
| out | <i>real(dp), dimension(:, :) :: aerodyn_resistance1</i> | aerodynamical resistance            |

#### Authors

Matthias Zink

#### Date

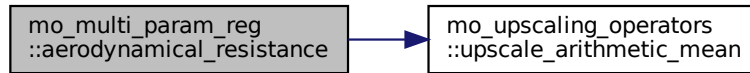
Apr 2013

Definition at line 1232 of file mo\_multi\_param\_reg.f90.

References `mo_common_constants::eps_dp`, `mo_mpr_constants::karman`, `mo_upscaling_operators::upscale_↔ arithmetic_mean()`, and `mo_mpr_constants::windmeasheight`.

Referenced by `mpr()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.54.2.2 baseflow\_param()

```

subroutine mo_multi_param_reg::baseflow_param (
    real(dp), dimension(:), intent(in) param,
    integer(i4), dimension(:), intent(in) geoUnit0,
    real(dp), dimension(:), intent(out) k2_0 )
  
```

baseflow recession parameter

This subroutine calculates the baseflow recession parameter based on the geological units at the Level 0 scale. For each level 0 cell, it assigns the value specified in the parameter array `param` for the geological unit in this cell. Global parameters needed (see `mhm_parameter.nml`):

- `param(1) = GeoParam(1,:)`
- `param(2) = GeoParam(2,:)`
- ...

#### Parameters

|     |  |   |
|-----|--|---|
| in  | <code>real(dp), dimension(:) :: param</code>       | list of required parameters               |
| in  | <code>integer(i4), dimension(:) :: geoUnit0</code> | ids of geological units at L0             |
| out | <code>real(dp), dimension(:) :: k2_0</code>        | - baseflow recession parameter at Level 0 |

#### Authors

Stephan Thober, Rohini Kumar

## Date

Dec 2012

Definition at line 716 of file mo\_multi\_param\_reg.f90.

References mo\_mpr\_global\_variables::geounitlist, and mo\_common\_constants::nodata\_dp.

Referenced by mpr().

Here is the caller graph for this function:



### 17.54.2.3 canopy\_intercept\_param()

```

subroutine, public mo_multi_param_reg::canopy_intercept_param(
  integer(i4), dimension(:, :) :: processMatrix,
  real(dp), dimension(:) :: param,
  real(dp), dimension(:, :) :: LAI0,
  integer(i4), dimension(:) :: n_subcells1,
  integer(i4), dimension(:) :: upper_bound1,
  integer(i4), dimension(:) :: lower_bound1,
  integer(i4), dimension(:) :: left_bound1,
  integer(i4), dimension(:) :: right_bound1,
  integer(i4), dimension(:) :: Id0,
  logical, dimension(:, :) :: mask0,
  real(dp), intent(in) nodata,
  real(dp), dimension(:, :) :: max_intercept1 )
  
```

estimate effective maximum interception capacity at L1

estimate effective maximum interception capacity at L1 for a given Leaf Area Index field. Global parameters needed (see mhm\_parameter.nml): Process Case 1:

- param(1) = canopyInterceptionFactor

#### Parameters

|     |  |  |
|-----|--|--|
| in  | <i>integer(i4), dimension(:, :) :: processMatrix</i> | indicate processes                         |
| in  | <i>real(dp), dimension(:) :: param</i>               | array of global parameters                 |
| in  | <i>real(dp), dimension(:, :) :: LAI0</i>             | LAI at level-0(nCells0, time)              |
| in  | <i>integer(i4), dimension(:) :: n_subcells1</i>      | Number of L0 cells within a L1 cell        |
| in  | <i>integer(i4), dimension(:) :: upper_bound1</i>     | Upper row of high resolution block         |
| in  | <i>integer(i4), dimension(:) :: lower_bound1</i>     | Lower row of high resolution block         |
| in  | <i>integer(i4), dimension(:) :: left_bound1</i>      | Left column of high resolution block       |
| in  | <i>integer(i4), dimension(:) :: right_bound1</i>     | Right column of high resolution block      |
| in  | <i>integer(i4), dimension(:) :: Id0</i>              | Cell ids at level 0                        |
| in  | <i>logical, dimension(:, :) :: mask0</i>             | mask at level 0 field                      |
| in  | <i>real(dp) :: nodata</i>                            | - nodata value                             |
| out | <i>real(dp), dimension(:, :) :: max_intercept1</i>   | max interception at level-1(nCells1, time) |

**Authors**

Rohini Kumar

**Date**

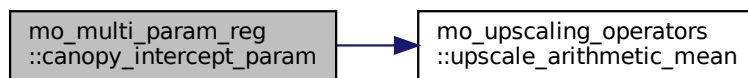
Aug. 2013

Definition at line 1103 of file mo\_multi\_param\_reg.f90.

References mo\_upscaling\_operators::upscale\_arithmetic\_mean().

Referenced by mpr().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.54.2.4 iper\_thres\_runoff()**

```

subroutine mo_multi_param_reg::iper_thres_runoff (
    real(dp), dimension(1), intent(in) param,
    real(dp), dimension(:, :, :), intent(out) sealedThresh1 ) [private]
  
```

sets the impervious layer threshold parameter for runoff generation

to be done by Kumar .... Global parameters needed (see mhm\_parameter.nml):

- param(1) = imperviousStorageCapacity

**Parameters**

|     |   |                             |
|-----|---|-----------------------------|
| in  | <i>real(dp), dimension(1) :: param</i>              | - given threshold parameter |
| out | <i>real(dp), dimension(:, :, : :: sealedThresh1</i> |                             |

**Authors**

Stephan Thober, Rohini Kumar

## Date

Dec 2012

Definition at line 910 of file mo\_multi\_param\_reg.f90.

Referenced by mpr().

Here is the caller graph for this function:



### 17.54.2.5 karstic\_layer()

```

subroutine mo_multi_param_reg::karstic_layer (
    real(dp), dimension(3), intent(in) param,
    integer(i4), dimension(:), intent(in) geoUnit0,
    logical, dimension(:, :) , intent(in) mask0,
    real(dp), dimension(:), intent(in) SMs_FC0,
    real(dp), dimension(:), intent(in) KsVar_V0,
    integer(i4), dimension(:), intent(in) Id0,
    integer(i4), dimension(:), intent(in) n_subcells1,
    integer(i4), dimension(:), intent(in) upper_bound1,
    integer(i4), dimension(:), intent(in) lower_bound1,
    integer(i4), dimension(:), intent(in) left_bound1,
    integer(i4), dimension(:), intent(in) right_bound1,
    real(dp), dimension(:), intent(out) karstLoss1,
    real(dp), dimension(:), intent(out) Ll_Kp ) [private]
  
```

calculates the Karstic percolation loss

This subroutine calls first the karstic\_fraction upscaling routine for determine the karstic fraction area for every Level 1 cell. Then, the karstic percolation loss is estimated given two shape parameters by

$$karstLoss1 = 1 + (fKarArea * param(1)) * ((-1) ** INT(param(2), i4))$$

where *karstLoss1* is the karstic percolation loss and *fKarArea* is the fraction of karstic area at level 1 Global parameters needed (see mhm\_parameter.nml):

- param(1) = rechargeCoefficient
- param(2) = rechargeFactor\_karstic
- param(3) = gain\_loss\_GWreservoir\_karstic

#### Parameters

|    |  |   |
|----|--|---|
| in | <i>real(dp), dimension(3) :: param</i>       | parameters  |
| in | <i>integer(i4), dimension(:) :: geoUnit0</i> | id of the Karstic formation                                       |
| in | <i>logical, dimension(:, :) :: mask0</i>     | mask at level 0   |
| in | <i>real(dp), dimension(:) :: SMs_FC0</i>     | [-] soil moisture deficit from field capacity w.r.t to saturation |
| in | <i>real(dp), dimension(:) :: KsVar_V0</i>    | [-] relative variability of saturated                             |
| in | <i>integer(i4), dimension(:) :: Id0</i>      | Cell ids of hi res field  |

## Parameters

|     |  |                                     |
|-----|--|-------------------------------------|
| in  | <i>integer(i4), dimension(:) :: n_subcells1</i>  | number of I0 cells within a I1 cell |
| in  | <i>integer(i4), dimension(:) :: upper_bound1</i> | upper row of a I1 cell in I0 grid   |
| in  | <i>integer(i4), dimension(:) :: lower_bound1</i> | lower row of a I1 cell in I0 grid   |
| in  | <i>integer(i4), dimension(:) :: left_bound1</i>  | left col of a I1 cell in I0 grid    |
| in  | <i>integer(i4), dimension(:) :: right_bound1</i> | right col of a I1 cell in I0 grid   |
| out | <i>real(dp), dimension(:) :: karstLoss1</i>      | [-] Karstic percolation loss        |
| out | <i>real(dp), dimension(:) :: L1_Kp</i>           | [d-1] percolation coefficient       |

## Authors

Rohini Kumar, Stephan Thober

## Date

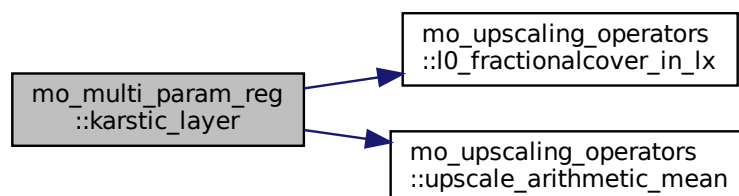
Feb 2013

Definition at line 971 of file mo\_multi\_param\_reg.f90.

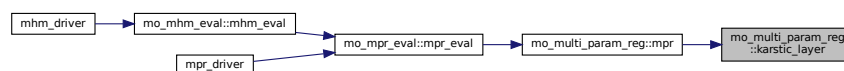
References mo\_mpr\_global\_variables::geounitkar, mo\_mpr\_global\_variables::geounitlist, mo\_upscaling\_operators::I0\_fractionalcover\_in\_Ix(), mo\_common\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, and mo\_upscaling\_operators::upscale\_arithmetic\_mean().

Referenced by mpr().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.54.2.6 mpr()

```

subroutine, public mo_multi_param_reg::mpr (
    logical, dimension(:, :), intent(in) mask0,
  
```

```

integer(i4), dimension(:), intent(in) geoUnit0,
integer(i4), dimension(:, :), intent(in) soilId0,
real(dp), dimension(:), intent(in) Asp0,
real(dp), dimension(:, :), intent(in) gridded_LAI0,
integer(i4), dimension(:, :), intent(in) LCover0,
real(dp), dimension(:), intent(in) slope_emp0,
real(dp), dimension(:), intent(in) y0,
integer(i4), dimension(:), intent(in) Id0,
integer(i4), dimension(:), intent(in) upper_bound1,
integer(i4), dimension(:), intent(in) lower_bound1,
integer(i4), dimension(:), intent(in) left_bound1,
integer(i4), dimension(:), intent(in) right_bound1,
integer(i4), dimension(:), intent(in) n_subcells1,
real(dp), dimension(:, :, :), intent(inout) fSealed1,
real(dp), dimension(:, :, :), intent(inout) alpha1,
real(dp), dimension(:, :, :), intent(inout) degDayIncl,
real(dp), dimension(:, :, :), intent(inout) degDayMax1,
real(dp), dimension(:, :, :), intent(inout) degDayNoPre1,
real(dp), dimension(:, :, :), intent(inout) fAsp1,
real(dp), dimension(:, :, :), intent(inout) HarSamCoeff1,
real(dp), dimension(:, :, :), intent(inout) PrieTayAlpha1,
real(dp), dimension(:, :, :), intent(inout) aeroResist1,
real(dp), dimension(:, :, :), intent(inout) surfResist1,
real(dp), dimension(:, :, :), intent(inout) fRoots1,
real(dp), dimension(:, :, :), intent(inout) kFastFlow1,
real(dp), dimension(:, :, :), intent(inout) kSlowFlow1,
real(dp), dimension(:, :, :), intent(inout) kBaseFlow1,
real(dp), dimension(:, :, :), intent(inout) kPercol,
real(dp), dimension(:, :, :), intent(inout) karstLoss1,
real(dp), dimension(:, :, :), intent(inout) soilMoistFC1,
real(dp), dimension(:, :, :), intent(inout) soilMoistSat1,
real(dp), dimension(:, :, :), intent(inout) soilMoistExpl,
real(dp), dimension(:, :, :), intent(inout) jarvis_thresh_cl,
real(dp), dimension(:, :, :), intent(inout) tempThresh1,
real(dp), dimension(:, :, :), intent(inout) unsatThresh1,
real(dp), dimension(:, :, :), intent(inout) sealedThresh1,
real(dp), dimension(:, :, :), intent(inout) wiltingPoint1,
real(dp), dimension(:, :, :), intent(inout) maxInter1,
real(dp), dimension(:, :, :), intent(inout) petLAIcorFactor,
real(dp), dimension(:), intent(in), optional, target parameterset )

```

Regionalizing and Upscaling process parameters.

calculating process parameters at L0 scale (Regionalization), like:

- Baseflow recession parameter
- Soil moisture parameters
- PET correction for aspect and upscale these parameters to retrieve effective parameters at scale L1. Further parameter regionalizations are done for:
- snow accumulation and melting parameters
- threshold parameter for runoff generation on impervious layer
- karstic percolation loss
- setting up the Regionalized Routing Parameters

## Parameters

|         |   |   |
|---------|---|---|
| in      | <i>logical, dimension(:, :) :: mask0</i>                | mask at level 0 field   |
| in      | <i>integer(i4), dimension(:, :) :: geoUnit0</i>         | L0 geological units   |
| in      | <i>integer(i4), dimension(:, :) :: soilId0</i>          | soil Ids at level 0   |
| in      | <i>real(dp), dimension(:, :) :: Asp0</i>                | [degree] Aspect at Level 0  |
| in      | <i>real(dp), dimension(:, :) :: gridded_LAI0</i>        | LAI grid at level 0, with dim2 = time   |
| in      | <i>integer(i4), dimension(:, :) :: LCOVER0</i>          | land cover at level 0   |
| in      | <i>real(dp), dimension(:, :) :: slope_emp0</i>          | Empirical quantiles of slope  |
| in      | <i>real(dp), dimension(:, :) :: y0</i>                  | y0 at level 0   |
| in      | <i>integer(i4), dimension(:, :) :: Id0</i>              | Cell ids at level 0   |
| in      | <i>integer(i4), dimension(:, :) :: upper_bound1</i>     | Upper row of hi res block   |
| in      | <i>integer(i4), dimension(:, :) :: lower_bound1</i>     | Lower row of hi res block   |
| in      | <i>integer(i4), dimension(:, :) :: left_bound1</i>      | Left column of hi res block   |
| in      | <i>integer(i4), dimension(:, :) :: right_bound1</i>     | Right column of hi res block  |
| in      | <i>integer(i4), dimension(:, :) :: n_subcells1</i>      | Number of L0 cells within a L1 cell   |
| in, out | <i>real(dp), dimension(:, :, :) :: fSealed1</i>         | [1] fraction of sealed area   |
| in, out | <i>real(dp), dimension(:, :, :) :: alpha1</i>           | [1] Exponent for the upper reservoir  |
| in, out | <i>real(dp), dimension(:, :, :) :: degDayInc1</i>       | [d-1 degreeC-1] Increase of the Degree-day factor per mm of increase in precipitation   |
| in, out | <i>real(dp), dimension(:, :, :) :: degDayMax1</i>       | [mm-1 degreeC-1] Maximum Degree-day factor  |
| in, out | <i>real(dp), dimension(:, :, :) :: degDayNoPre1</i>     | [mm-1 degreeC-1] Degree-day factor with no precipitation                                |
| in, out | <i>real(dp), dimension(:, :, :) :: fAsp1</i>            | [1] PET correction for Aspect at level 1  |
| in, out | <i>real(dp), dimension(:, :, :) :: HarSamCoeff1</i>     | [1] PET Hargreaves Samani coeff. at level 1   |
| in, out | <i>real(dp), dimension(:, :, :) :: PrieTayAlpha1</i>    | [1] PET Priestley Taylor coeff. at level 1  |
| in, out | <i>real(dp), dimension(:, :, :) :: aeroResist1</i>      | [s m-1] PET aerodynamical resistance at level 1   |
| in, out | <i>real(dp), dimension(:, :, :) :: surfResist1</i>      | [s m-1] PET bulk surface resistance at level 1  |
| in, out | <i>real(dp), dimension(:, :, :) :: fRoots1</i>          | fraction of roots in soil horizon   |
| in, out | <i>real(dp), dimension(:, :, :) :: kFastFlow1</i>       | [10 <sup>-3</sup> m] Recession coefficient of the upper reservoir, upper outlet         |
| in, out | <i>real(dp), dimension(:, :, :) :: kSlowFlow1</i>       | [10 <sup>-3</sup> m] Recession coefficient of the upper reservoir, lower outlet         |
| in, out | <i>real(dp), dimension(:, :, :) :: kBaseFlow1</i>       | Level 1 baseflow recession  |
| in, out | <i>real(dp), dimension(:, :, :) :: kPerco1</i>          | [d-1] percolation coefficient   |
| in, out | <i>real(dp), dimension(:, :, :) :: karstLoss1</i>       |   |
| in, out | <i>real(dp), dimension(:, :, :) :: soilMoistFC1</i>     | [10 <sup>-3</sup> m] field capacity   |
| in, out | <i>real(dp), dimension(:, :, :) :: soilMoistSat1</i>    | [10 <sup>-3</sup> m] depth of saturated SM  |
| in, out | <i>real(dp), dimension(:, :, :) :: soilMoistExp1</i>    | Parameter that determines the rel. contribution to SM, upscal. Bulk den.                |
| in, out | <i>real(dp), dimension(:, :, :) :: jarvis_thresh_c1</i> | [1] jarvis critical value for norm SWC  |
| in, out | <i>real(dp), dimension(:, :, :) :: tempThresh1</i>      | [degreeC] threshold temperature for snow rain   |
| in, out | <i>real(dp), dimension(:, :, :) :: unsatThresh1</i>     | [10 <sup>-3</sup> m] Threshold water depth in upper reservoir (for Runoff contribution) |
| in, out | <i>real(dp), dimension(:, :, :) :: sealedThresh1</i>    | threshold parameter   |
| in, out | <i>real(dp), dimension(:, :, :) :: wiltingPoint1</i>    | [10 <sup>-3</sup> m] permanent wilting point  |



## Parameters

|         |   |  |
|---------|---|--|
| in, out | <i>real(dp), dimension(:, :, :) :: maxInter1</i>        |  |
| in, out | <i>real(dp), dimension(:, :, :) :: petLAlcorFactor</i>  |  |
| in      | <i>real(dp), dimension(:), optional :: parameterset</i> |  |

## Authors

Stephan Thober, Rohini Kumar

## Date

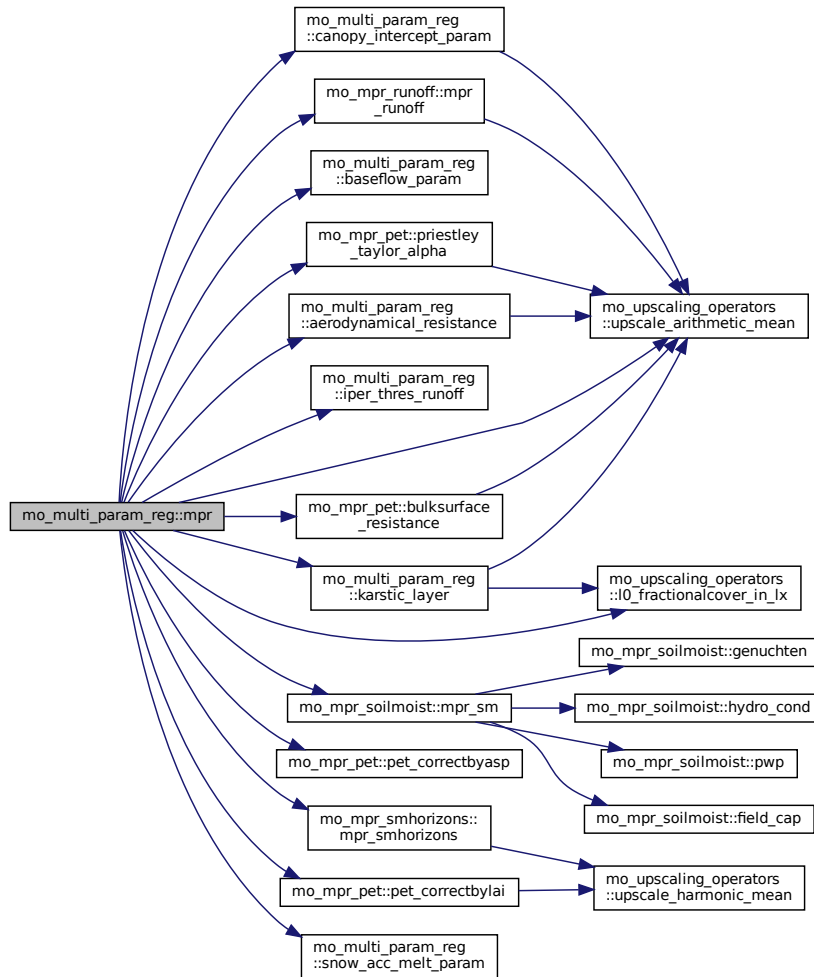
Dec 2012

Definition at line 124 of file mo\_multi\_param\_reg.f90.

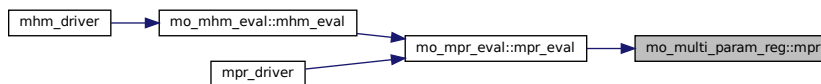
References aerodynamical\_resistance(), baseflow\_param(), mo\_mpr\_pet::bulksurface\_resistance(), canopy\_↔ intercept\_param(), mo\_mpr\_global\_variables::fracsealed\_cityarea, mo\_common\_variables::global\_parameters, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_mpr\_global\_variables::iflag\_soildb, iper\_thres\_runoff(), karstic\_layer(), mo\_upscaling\_operators::l0\_fractionalcover\_in\_lx(), mo\_mpr\_runoff::mpr\_runoff(), mo\_mpr\_↔ \_soilmoist::mpr\_sm(), mo\_mpr\_smhorizons::mpr\_smhorizons(), mo\_common\_constants::nodata\_dp, mo\_↔ common\_constants::nodata\_i4, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_mpr\_pet::pet\_correctbyasp(), mo\_mpr\_pet::pet\_correctbylai(), mo\_mpr\_pet::priestley\_taylor\_alpha(), mo\_common\_variables::processmatrix, snow\_acc\_melt\_param(), mo\_mpr\_global\_variables::soildb, and mo\_upscaling\_operators::upscale\_arithmetic\_↔ mean().

Referenced by mo\_mpr\_eval::mpr\_eval().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.54.2.7 snow\_acc\_melt\_param()

```

subroutine mo_multi_param_reg::snow_acc_melt_param (
    real(dp), dimension(8), intent(in) param,
    real(dp), dimension(:), intent(in) fForest1,
    real(dp), dimension(:), intent(in) fIperml,
    real(dp), dimension(:), intent(in) fPerml,
  
```

```

real(dp), dimension(:), intent(out) tempThresh1,
real(dp), dimension(:), intent(out) degDayNoPre1,
real(dp), dimension(:), intent(out) degDayInc1,
real(dp), dimension(:), intent(out) degDayMax1 )

```

Calculates the snow parameters.

This subroutine calculates the snow parameters threshold temperature (TT), degree-day factor without precipitation (DD) and maximum degree-day factor (DDmax) as well as increase of degree-day factor per mm of increase in precipitation (IDDP). Global parameters needed (see mhm\_parameter.nml):

- param(1) = snowTreshholdTemperature
- param(2) = degreeDayFactor\_forest
- param(3) = degreeDayFactor\_impervious
- param(4) = degreeDayFactor\_pervious
- param(5) = increaseDegreeDayFactorByPrecip
- param(6) = maxDegreeDayFactor\_forest
- param(7) = maxDegreeDayFactor\_impervious
- param(8) = maxDegreeDayFactor\_pervious INTENT(IN)

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>real(dp), dimension(8) :: param</i>        | eight global parameters                       |
| in  | <i>real(dp), dimension(:) :: fForest1</i>     | [1] fraction of forest cover                  |
| in  | <i>real(dp), dimension(:) :: flperm1</i>      | [1] fraction of sealed area                   |
| in  | <i>real(dp), dimension(:) :: fPerm1</i>       | [1] fraction of permeable area                |
| out | <i>real(dp), dimension(:) :: tempThresh1</i>  | [degreeC] threshold temperature for snow rain |
| out | <i>real(dp), dimension(:) :: degDayNoPre1</i> | [mm-1 degreeC-1] Degree-day factor with       |
| out | <i>real(dp), dimension(:) :: degDayInc1</i>   | [d-1 degreeC-1] Increase of the Degree-day    |
| out | <i>real(dp), dimension(:) :: degDayMax1</i>   | [mm-1 degreeC-1] Maximum Degree-day factor    |

#### Authors

Stephan Thober, Rohini Kumar

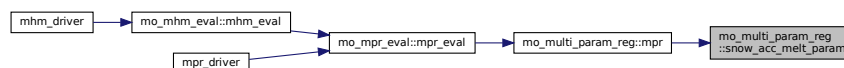
#### Date

Dec 2012

Definition at line 827 of file mo\_multi\_param\_reg.f90.

Referenced by mpr().

Here is the caller graph for this function:



## 17.55 mo\_neutrons Module Reference

Models to predict neutron intensities above soils.

### Functions/Subroutines

- subroutine, public `desiletsn0` (SoilMoisture, Horizons, N0, neutrons)  
*Calculate neutrons from soil moisture in the first layer.*
- subroutine, public `cosmic` (SoilMoisture, Horizons, params, neutron\_integral\_AFast, neutrons)  
*Calculate neutrons from soil moisture in all layers.*
- subroutine `oldintegration` (res, c)  
*TODO: add description.*
- subroutine, public `tabularintegralafast` (integral, maxC)  
*Save approximation data for A\_fast.*
- subroutine `approx_mon_int` (res, f, c, xmin, xmax, eps, steps, fxmin, fxmax)  
*TODO: add description.*
- recursive subroutine `approx_mon_int_steps` (res, f, c, xmin, xmax, eps, steps, fxmin, fxmax)  
*TODO: add description.*
- recursive subroutine `approx_mon_int_eps` (res, f, c, xmin, xmax, eps, fxmin, fxmax)  
*TODO: add description.*
- subroutine `lookupintegral` (res, integral, c)  
*TODO: add description.*
- real(dp) function `intgrandfast` (c, phi)  
*TODO: add description.*

### 17.55.1 Detailed Description

Models to predict neutron intensities above soils.

The number of neutrons above the ground is directly related to the number soil water content in the ground, air, vegetation and/or snow. This module forward-models neutron abundance as a state variable for each cell.

#### Authors

Martin Schroen

#### Date

Mar 2015 THIS MODULE IS WORK IN PROGRESS, DO NOT USE FOR RESEARCH.

### 17.55.2 Function/Subroutine Documentation

#### 17.55.2.1 approx\_mon\_int()

```
subroutine mo_neutrons::approx_mon_int (
    real(dp) res,
    real(dp), external f,
    real(dp), intent(in) c,
    real(dp), intent(in) xmin,
    real(dp), intent(in) xmax,
    real(dp), intent(in), optional eps,
```

```
integer(i4), intent(in), optional steps,
real(dp), intent(in), optional fxmin,
real(dp), intent(in), optional fxmax )
```

TODO: add description.

TODO: add description

#### Parameters

|    |                                       |  |
|----|---------------------------------------|--|
| in | <i>real(dp) :: c</i>                  |  |
| in | <i>real(dp) :: xmin</i>               |  |
| in | <i>real(dp) :: xmax</i>               |  |
| in | <i>real(dp), optional :: eps</i>      |  |
| in | <i>integer(i4), optional :: steps</i> |  |
| in | <i>real(dp), optional :: fxmin</i>    |  |
| in | <i>real(dp), optional :: fxmax</i>    |  |

#### Authors

Robert Schweppe

#### Date

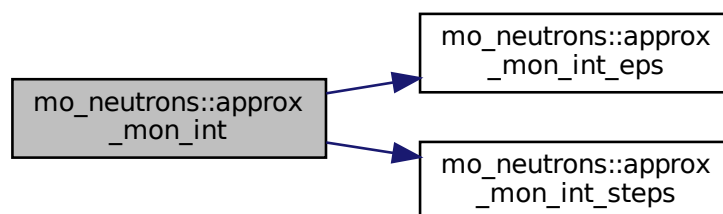
Jun 2018

Definition at line 436 of file mo\_neutrons.f90.

References approx\_mon\_int\_eps(), and approx\_mon\_int\_steps().

Referenced by tabularintegralafast().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.55.2.2 approx\_mon\_int\_eps()

```
recursive subroutine mo_neutrons::approx_mon_int_eps (
    real(dp) res,
    real(dp), external f,
    real(dp), intent(in) c,
    real(dp), intent(in) xmin,
    real(dp), intent(in) xmax,
    real(dp), intent(in) eps,
    real(dp), intent(in) fxmin,
    real(dp), intent(in) fxmax ) [private]
```

TODO: add description.

TODO: add description

#### Parameters

|    |                          |  |
|----|--------------------------|--|
| in | <i>real(dp) :: c</i>     |  |
| in | <i>real(dp) :: xmin</i>  |  |
| in | <i>real(dp) :: xmax</i>  |  |
| in | <i>real(dp) :: eps</i>   |  |
| in | <i>real(dp) :: fxmin</i> |  |
| in | <i>real(dp) :: fxmax</i> |  |

#### Authors

Robert Schweppe

#### Date

Jun 2018

Definition at line 596 of file mo\_neutrons.f90.

Referenced by approx\_mon\_int().

Here is the caller graph for this function:



### 17.55.2.3 approx\_mon\_int\_steps()

```
recursive subroutine mo_neutrons::approx_mon_int_steps (
    real(dp) res,
    real(dp), external f,
    real(dp), intent(in) c,
    real(dp), intent(in) xmin,
    real(dp), intent(in) xmax,
    real(dp), intent(in) eps,
    integer(i4), intent(in) steps,
```

```

real(dp), intent(in) fxmin,
real(dp), intent(in) fxmax ) [private]

```

TODO: add description.

TODO: add description

#### Parameters

|    |                             |  |
|----|-----------------------------|--|
| in | <i>real(dp) :: c</i>        |  |
| in | <i>real(dp) :: xmin</i>     |  |
| in | <i>real(dp) :: xmax</i>     |  |
| in | <i>real(dp) :: eps</i>      |  |
| in | <i>integer(i4) :: steps</i> |  |
| in | <i>real(dp) :: fxmin</i>    |  |
| in | <i>real(dp) :: fxmax</i>    |  |

#### Authors

Robert Schwappe

#### Date

Jun 2018

Definition at line 526 of file mo\_neutrons.f90.

Referenced by approx\_mon\_int().

Here is the caller graph for this function:



#### 17.55.2.4 cosmic()

```

subroutine, public mo_neutrons::cosmic (
  real(dp), dimension(:), intent(in) SoilMoisture,
  real(dp), dimension(:), intent(in) Horizons,
  real(dp), dimension(:), intent(in) params,
  real(dp), dimension(:), intent(in) neutron_integral_AFast,
  real(dp), intent(inout) neutrons )

```

Calculate neutrons from soil moisture in all layers.

Neutron counts above the ground (one value per cell in mHM) can be derived by a simplified physical neutron transport simulation. Fast cosmic-Ray neutrons are generated in the soil and attenuated differently in water and soil. The remaining neutrons that reached the surface relate to the profile of soil water content below. Variables like N, alpha and L3 are site-specific and need to be calibrated. ADDITIONAL INFORMATION COSMIC model based on Shuttleworth et al. 2013 Horizons(:) must not be zero. see supplementaries in literature J. Shuttleworth, R. Rosolem, M. Zreda, and T. Franz, The COsmic-ray Soil Moisture Interaction Code (COSMIC) for use in data assimilation, HESS, 17, 3205-3217, 2013, doi:10.5194/hess-17-3205-2013 Support and Code: <http://cosmos.hwr.arizona.edu/Software/cosmic.html>

## Parameters

|         |   |  |
|---------|---|--|
| in      | <i>real(dp), dimension(:) :: SoilMoisture</i>           | Soil Moisture                          |
| in      | <i>real(dp), dimension(:) :: Horizons</i>               | Horizon depths                         |
| in      | <i>real(dp), dimension(:) :: params</i>                 | ! N0, N1, N2, alpha0, alpha1, L30, L31 |
| in      | <i>real(dp), dimension(:) :: neutron_integral_AFast</i> | Tabular for Int Approx                 |
| in, out | <i>real(dp) :: neutrons</i>                             | Neutron counts                         |

## Authors

Martin Schroen, originally written by Rafael Rosolem

## Date

Mar 2015

Definition at line 145 of file mo\_neutrons.f90.

References mo\_mhm\_constants::cosmic\_alpha, mo\_mhm\_constants::cosmic\_bd, mo\_mhm\_constants::cosmic\_l1, mo\_mhm\_constants::cosmic\_l2, mo\_mhm\_constants::cosmic\_l3, mo\_mhm\_constants::cosmic\_l4, mo\_mhm\_constants::cosmic\_n, mo\_mhm\_constants::cosmic\_vwclat, mo\_mhm\_constants::h2odens, and lookupintegral().

Referenced by mo\_mhm::mhm().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.55.2.5 desiletsn0()

```

subroutine, public mo_neutrons::desiletsn0 (
    real(dp), dimension(:), intent(in) SoilMoisture,
    real(dp), dimension(:), intent(in) Horizons,
    real(dp), intent(in) N0,
    real(dp), intent(inout) neutrons )
  
```

Calculate neutrons from soil moisture in the first layer.

Using the N0-relation derived by Desilets, neutron counts above the ground (one value per cell in mHM) can be derived by a semi-empirical, semi-physical relation. The result depends on N0, the neutron counts for 0% soil



moisture. This variable is site-specific and is a global parameter in mHM. N0 formula based on Desilets et al. 2010 Horizons(1) must not be zero.  $N0=1500\text{cph}$ ,  $\text{SoilMoisture}(1,1)=700\text{mm}$ ,  $\text{Horizons}(1)=200\text{mm}$   $1500*(0.372+0.0808/(70\text{mm}/200\text{mm} + 0.115))$  DesiletsN0 = 819cph Desilets, D., M. Zreda, and T. P. A. Ferre (2010), Nature's neutron probe: Land surface hydrology at an elusive scale with cosmic rays, WRR, 46, W11505, doi:10.1029/2009WR008726.

#### Parameters

|         |   |                    |
|---------|---|--------------------|
| in      | <i>real(dp), dimension(:) :: SoilMoisture</i> | Soil Moisture      |
| in      | <i>real(dp), dimension(:) :: Horizons</i>     | Horizon depths     |
| in      | <i>real(dp) :: N0</i>                         | dry neutron counts |
| in, out | <i>real(dp) :: neutrons</i>                   | Neutron counts     |

#### Authors

Martin Schroen

#### Date

Mar 2015

Definition at line 78 of file mo\_neutrons.f90.

References mo\_mhm\_constants::desilets\_a0, mo\_mhm\_constants::desilets\_a1, and mo\_mhm\_constants::desilets\_a2.

Referenced by mo\_mhm::mhm().

Here is the caller graph for this function:



#### 17.55.2.6 intgrandfast()

```

real(dp) function mo_neutrons::intgrandfast (
    real(dp), intent(in) c,
    real(dp), intent(in) phi )
  
```

TODO: add description.

TODO: add description

#### Parameters

|    |                        |  |
|----|------------------------|--|
| in | <i>real(dp) :: c</i>   |  |
| in | <i>real(dp) :: phi</i> |  |

#### Authors

Robert Schweppe

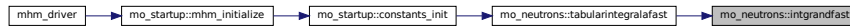
**Date**

Jun 2018

Definition at line 718 of file mo\_neutrons.f90.

Referenced by tabularintegralafast().

Here is the caller graph for this function:

**17.55.2.7 lookupintegral()**

```

subroutine mo_neutrons::lookupintegral (
    real(dp) res,
    real(dp), dimension(:), intent(in) integral,
    real(dp), intent(in) c ) [private]
  
```

TODO: add description.

TODO: add description

**Parameters**

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: integral</i> |  |
| in | <i>real(dp) :: c</i>                      |  |

**Authors**

Robert Schweppe

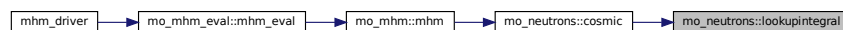
**Date**

Jun 2018

Definition at line 662 of file mo\_neutrons.f90.

Referenced by cosmic().

Here is the caller graph for this function:

**17.55.2.8 oldintegration()**

```

subroutine mo_neutrons::oldintegration (
    real(dp) res,
    real(dp), intent(in) c )
  
```

TODO: add description.

TODO: add description

#### Parameters

|    |                      |  |
|----|----------------------|--|
| in | <i>real(dp) :: c</i> |  |
|----|----------------------|--|

#### Authors

Robert Schweppe

#### Date

Jun 2018

Definition at line 287 of file mo\_neutrons.f90.

#### 17.55.2.9 tabularintegralafast()

```
subroutine, public mo_neutrons::tabularintegralafast (
    real(dp), dimension(:)  integral,
    real(dp), intent(in) maxC )
```

Save approximation data for *A\_fast*.

The COSMIC subroutine needs *A\_fast* to be calculated.  $A\_fast = \int_0^{\pi/2} \exp(-\text{Lambda\_fast}(z)/\cos(\phi)) \, d\phi$ . This subroutine stores data for *intsize* values for  $c = \text{Lambda\_fast}(z)$  between 0 and *maxC*, and will be written into the global array variable *neutron\_integral\_AFast*. The calculation of the values is done with a very precise recursive approximation subroutine. That recursive subroutine should not be used inside the time, cells and layer loops, because it is slow. Inside the loops in the module COSMIC the tabular is used to estimate *A\_fast*, if  $0 < c < \text{maxC}$ , otherwise the recursive approximation is used. *TabularIntegralAFast*: a tabular for calculations with splines *intsize* and *maxC* must be positive *intsize*=8000, *maxC*=20.0\_dp see splines for example

#### Parameters

|    |                         |                                   |
|----|-------------------------|-----------------------------------|
| in | <i>real(dp) :: maxC</i> | ! maximum value for <i>A_fast</i> |
|----|-------------------------|-----------------------------------|

#### Authors

Maren Kaluza

## Date

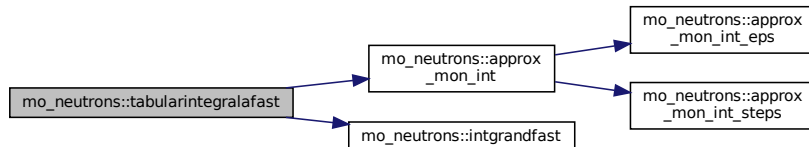
Nov 2017

Definition at line 368 of file mo\_neutrons.f90.

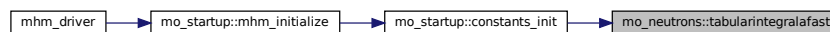
References `approx_mon_int()`, and `intgrandfast()`.

Referenced by `mo_startup::constants_init()`.

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.56 mo\_objective\_function Module Reference

Objective Functions for Optimization of mHM.

### Functions/Subroutines

- `real(dp)` function, public [objective](#) (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions.*
- `real(dp)` function, public [objective\\_master](#) (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions.*
- subroutine, public [objective\\_subprocess](#) (eval, arg1, arg2, arg3)  
*Wrapper for objective functions.*
- `real(dp)` function [objective\\_sm\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for soil moisture.*
- `real(dp)` function, dimension(6) [objective\\_q\\_et\\_tws\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for et, tws and q.*
- subroutine [init\\_indexarray\\_for\\_opti\\_data](#) (domainMeta, optidataOption, nOptiDomains, opti\_domain\_indices)  
*creates an index array of the inidices of the domains eval should MPI process.*
- `real(dp)` function [objective\\_sm\\_corr](#) (parameterset, eval)  
*Objective function for soil moisture.*
- `real(dp)` function [objective\\_sm\\_pd](#) (parameterset, eval)  
*Objective function for soil moisture.*
- `real(dp)` function [objective\\_sm\\_sse\\_standard\\_score](#) (parameterset, eval)  
*Objective function for soil moisture.*

- real(dp) function [objective\\_kge\\_q\\_rmse\\_tws](#) (parameterset, eval)  
*Objective function of KGE for runoff and RMSE for domain\_avg TWS (standardized scores)*
- real(dp) function [objective\\_neutrons\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for neutrons.*
- real(dp) function [objective\\_et\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for evapotranspiration (et).*
- real(dp) function [objective\\_kge\\_q\\_sm\\_corr](#) (parameterset, eval)  
*Objective function of KGE for runoff and correlation for SM.*
- real(dp) function [objective\\_kge\\_q\\_et](#) (parameterset, eval)  
*Objective function of KGE for runoff and KGE for ET.*
- real(dp) function [objective\\_kge\\_q\\_rmse\\_et](#) (parameterset, eval)  
*Objective function of KGE for runoff and RMSE for domain\_avg ET (standardized scores)*
- subroutine [create\\_domain\\_avg\\_tws](#) (iDomain, twsOptiSim, tws\_catch\_avg\_domain, tws\_opti\_catch\_avg, ← domain, mask\_times)
- subroutine [create\\_domain\\_avg\\_et](#) (iDomain, etOptiSim, et\_catch\_avg\_domain, et\_opti\_catch\_avg\_domain, mask\_times)
- subroutine [convert\\_tws\\_to\\_twsa](#) (twsOptiSim, L1\_twsaObs, twsaOptiSim)

### 17.56.1 Detailed Description

Objective Functions for Optimization of mHM.

This module provides a wrapper for several objective functions used to optimize mHM against various variables. If the objective is only regarding runoff move it to [mRRM/mo\\_mrm\\_objective\\_function\\_runoff.f90](#). If it contains besides runoff another variable like TWS implement it here. All the objective functions are supposed to be minimized! (10) SO: SM: 1.0 - KGE of catchment average soilmoisture (11) SO: SM: 1.0 - Pattern dissimilarity (PD) of spatially distributed soil moisture (12) SO: SM: Sum of squared errors (SSE) of spatially distributed standard score (normalization) of soil moisture (13) SO: SM: 1.0 - average temporal correlation of spatially distributed soil moisture (15) SO: Q + TWS: [1.0-KGE(Q)]\*RMSE(domain\_avg\_TWS) - objective function using Q and domain average !ToDo: change comment (standard score) TWS (17) SO: N: 1.0 - KGE of spatio-temporal neutron data, catchment-average (27) SO: ET: 1.0 - KGE of catchment average evapotranspiration

#### Authors

Juliane Mai

#### Date

Dec 2012

### 17.56.2 Function/Subroutine Documentation

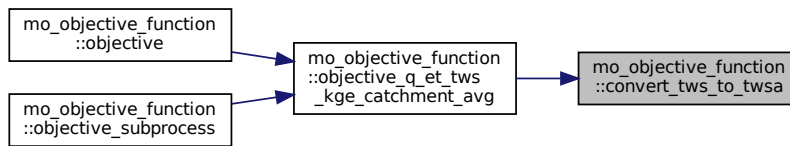
#### 17.56.2.1 convert\_tws\_to\_twsa()

```
subroutine mo_objective_function::convert_tws_to_twsa (
    type(optidata_sim), intent(in) twsOptiSim,
    type(optidata), intent(in) L1_twsaObs,
    type(optidata_sim), intent(inout) twsaOptiSim )
```

Definition at line 2601 of file mo\_objective\_function.f90.

Referenced by [objective\\_q\\_et\\_tws\\_kge\\_catchment\\_avg\(\)](#).

Here is the caller graph for this function:



### 17.56.2.2 create\_domain\_avg\_et()

```

subroutine mo_objective_function::create_domain_avg_et (
  integer(i4), intent(in) iDomain,
  type(optidata_sim), dimension(:), intent(in) etOptiSim,
  real(dp), dimension(:), intent(out), allocatable et_catch_avg_domain,
  real(dp), dimension(:), intent(out), allocatable et_opti_catch_avg_domain,
  logical, dimension(:), intent(out), allocatable mask_times )

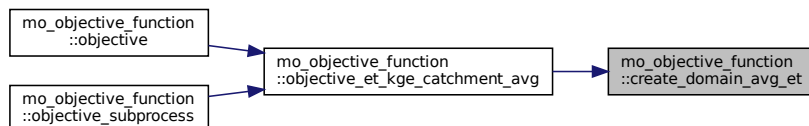
```

Definition at line 2546 of file `mo_objective_function.f90`.

References `mo_global_variables::l1_etobs`, and `mo_common_constants::nodata_dp`.

Referenced by `objective_et_kge_catchment_avg()`.

Here is the caller graph for this function:



### 17.56.2.3 create\_domain\_avg\_tws()

```

subroutine mo_objective_function::create_domain_avg_tws (
  integer(i4), intent(in) iDomain,
  type(optidata_sim), dimension(:), intent(in) twsOptiSim,
  real(dp), dimension(:), intent(out), allocatable tws_catch_avg_domain,
  real(dp), dimension(:), intent(out), allocatable tws_opti_catch_avg_domain,
  logical, dimension(:), intent(out), allocatable mask_times )

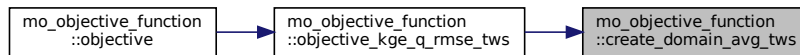
```

Definition at line 2491 of file `mo_objective_function.f90`.

References `mo_global_variables::l1_twsaobs`, and `mo_common_constants::nodata_dp`.

Referenced by `objective_kge_q_rmse_tws()`.

Here is the caller graph for this function:



### 17.56.2.4 init\_indexarray\_for\_opti\_data()

```

subroutine mo_objective_function::init_indexarray_for_opti_data (
    type(domain_meta), intent(in) domainMeta,
    integer(i4), intent(in) optidataOption,
    integer(i4), intent(out) nOptiDomains,
    integer(i4), dimension(:), intent(out), allocatable opti_domain_indices )
  
```

creates an index array of the indices of the domains eval should MPI process.

The data type domainMeta contains an array optidata of size domainMetaDomains, telling us, which domains should be optimized with which opti\_data. This subroutine splits all domains assigned to a process and returns an index list corresponding to the value of domainMetaoptidata.

The index array opti\_domain\_indices can then be passed as an optional argument to the eval subroutine. The eval then instead of using loops over all domains only uses the passed indices.

This subroutine also returns the size of that array since it helps with the calculations of the optimization in the end.

#### Authors

Maren Kaluza

#### Date

July 2019

#### Parameters

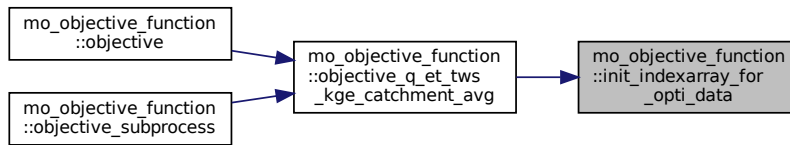
|     |                            |   |
|-----|----------------------------|---|
| in  | <i>domainmeta</i>          | meta data for all domains assigned to that process                              |
| in  | <i>optidataoption</i>      | which opti data should be used in the eval called after calling this subroutine |
| out | <i>noptidomains</i>        | number of domains that will be optimized in the following eval call             |
| out | <i>opti_domain_indices</i> | the indices of the domains that are to be processed in the following eval call  |

domain loop counter

Definition at line 908 of file mo\_objective\_function.f90.

Referenced by objective\_q\_et\_tws\_kge\_catchment\_avg().

Here is the caller graph for this function:



### 17.56.2.5 objective()

```

real(dp) function, public mo_objective_function::objective (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in), optional arg1,
    real(dp), intent(out), optional arg2,
    real(dp), intent(out), optional arg3 )
  
```

Wrapper for objective functions.

The functions selects the objective function case defined in a namelist, i.e. the global variable *opti\_function*. It return the objective function value for a specific parameter set.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>REAL(dp), DIMENSION(:) :: parameterset</i> |  |
| in  | <i>procedure(eval_interface) :: eval</i>      |  |
| in  | <i>real(dp), optional :: arg1</i>             |  |
| out | <i>real(dp), optional :: arg2</i>             |  |
| out | <i>real(dp), optional :: arg3</i>             |  |

#### Returns

*real(dp) :: objective* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Juliane Mai

#### Date

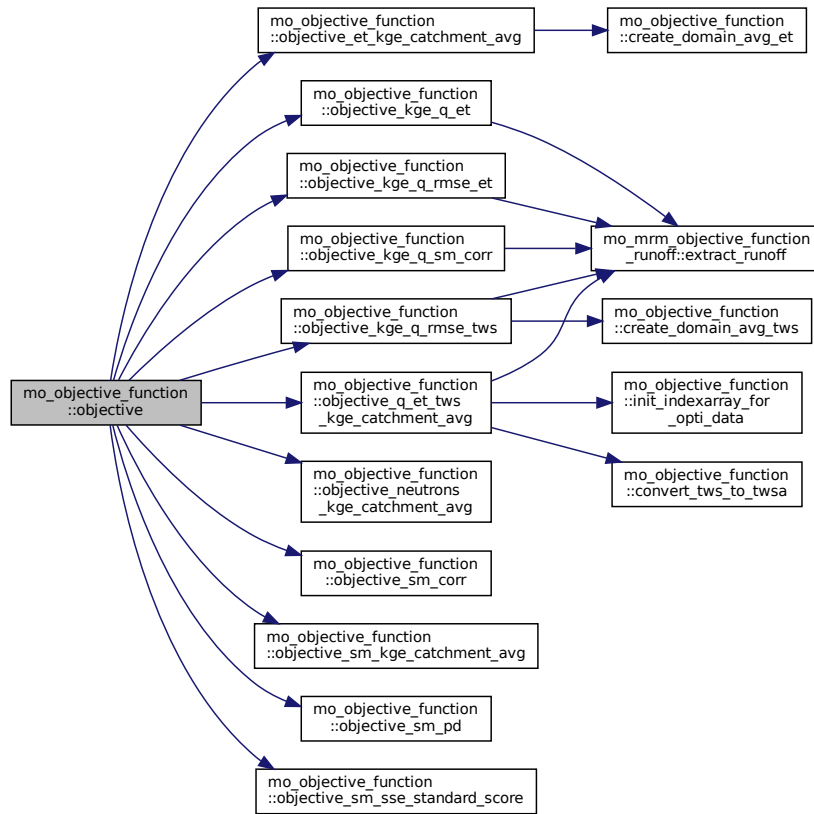
Dec 2012

Definition at line 88 of file *mo\_objective\_function.f90*.

References *mo\_common\_constants::nodata\_dp*, *objective\_et\_kge\_catchment\_avg()*, *objective\_kge\_q\_et()*, *objective\_kge\_q\_rmse\_et()*, *objective\_kge\_q\_rmse\_tws()*, *objective\_kge\_q\_sm\_corr()*, *objective\_neutrons\_kge\_catchment\_avg()*, *objective\_q\_et\_tws\_kge\_catchment\_avg()*, *objective\_sm\_corr()*, *objective\_sm\_kge\_catchment\_avg()*, *objective\_sm\_pd()*, *objective\_sm\_sse\_standard\_score()*, and *mo\_common\_mhm\_mrm\_variables::opti\_function*.



Here is the call graph for this function:



### 17.56.2.6 objective\_et\_kge\_catchment\_avg()

```

real(dp) function mo_objective_function::objective_et_kge_catchment_avg (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function for evapotranspiration (et).

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Kling-Gupta model efficiency  $KGE$  of the catchment average evapotranspiration (et) is calculated

$$KGE = 1.0 - \sqrt{((1-r)^2 + (1-\alpha)^2 + (1-\beta)^2)}$$

where  $r$  = Pearson product-moment correlation coefficient  $\alpha$  = ratio of simulated mean to observed mean  $\beta$  = ratio of simulated standard deviation to observed standard deviation is calculated and the objective function for a given domain  $i$  is

$$\phi_i = 1.0 - KGE_i$$

$\phi_i$  is the objective since we always apply minimization methods. The minimal value of  $\phi_i$  is 0 for the optimal  $KGE$  of 1.0. Finally, the overall objective function value  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all domains  $N$ .

$$OF = \sqrt[6]{\sum ((1.0 - KGE_i)/N)^6}$$

The observed data `L1_et`, `L1_et_mask` are global in this module.

## Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

## Returns

*real(dp) :: objective\_et\_kge\_catchment\_avg* — objective function value (which will be e.g. minimized by an optimization routine)

## Authors

Johannes Brenner

## Date

Feb 2017

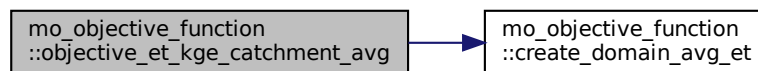
simulated et

Definition at line 1780 of file mo\_objective\_function.f90.

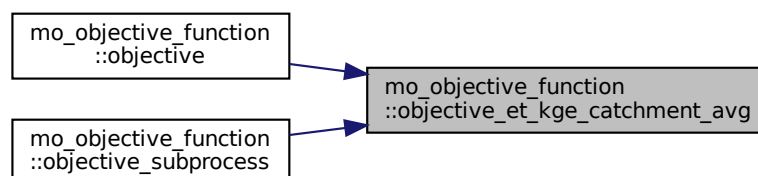
References `create_domain_avg_et()`, `mo_common_variables::domainmeta`, and `mo_common_constants::nodata←_dp`.

Referenced by `objective()`, and `objective_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.56.2.7 objective\_kge\_q\_et()

```
real(dp) function mo_objective_function::objective_kge_q_et (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
```

Objective function of KGE for runoff and KGE for ET.

Objective function of KGE for runoff and KGE for ET. Further details can be found in the documentation of objective functions '14 - objective\_multiple\_gauges\_kge\_power6'.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

real(dp) :: objective\_kge\_q\_et — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Johannes Brenner

#### Date

July 2017

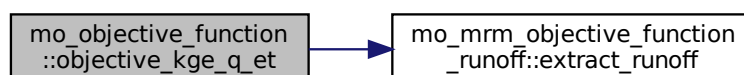
simulated et

Definition at line 2062 of file mo\_objective\_function.f90.

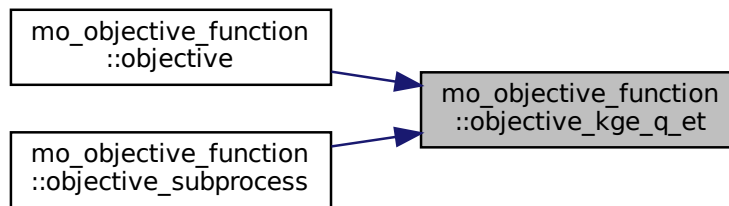
References mo\_common\_variables::domainmeta, mo\_mrm\_objective\_function\_runoff::extract\_runoff(), mo\_↔ global\_variables::l1\_etobs, and mo\_common\_variables::level1.

Referenced by objective(), and objective\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.56.2.8 objective\_kge\_q\_rmse\_et()

```

real(dp) function mo_objective_function::objective_kge_q_rmse_et (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of KGE for runoff and RMSE for domain\_avg ET (standarized scores)

Objective function of KGE for runoff and RMSE for domain\_avg ET (standarized scores)

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

real(dp) :: objective\_kge\_q\_rmse\_et — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Johannes Brenner

#### Date

July 2017

simulated et

Definition at line 2239 of file mo\_objective\_function.f90.

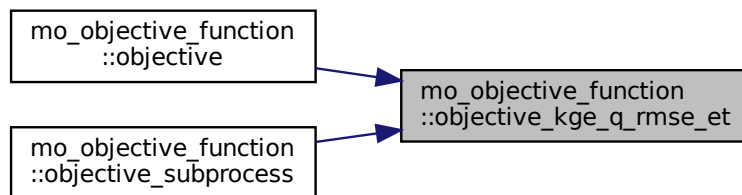
References mo\_common\_variables::domainmeta, mo\_common\_constants::eps\_dp, mo\_common\_mhm\_mrm\_variables::evalper, mo\_mrm\_objective\_function\_runoff::extract\_runoff(), mo\_global\_variables::l1\_etobs, and mo\_common\_constants::nodata\_dp.

Referenced by objective(), and objective\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.56.2.9 objective\_kge\_q\_rmse\_tws()

```

real(dp) function mo_objective_function::objective_kge_q_rmse_tws (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of KGE for runoff and RMSE for domain\_avg TWS (standarized scores)

Objective function of KGE for runoff and RMSE for domain\_avg TWS (standarized scores)

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

`real(dp) :: objective_kge_q_rmse_tws` — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Oldrich Rakovec, Rohini Kumar

## Date

Oct. 2015

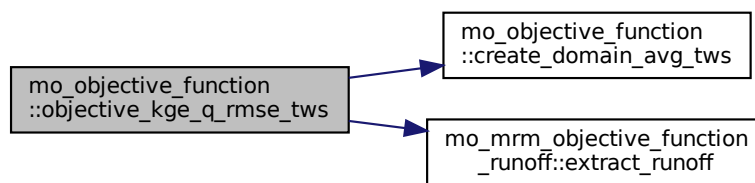
simulated tws

Definition at line 1396 of file mo\_objective\_function.f90.

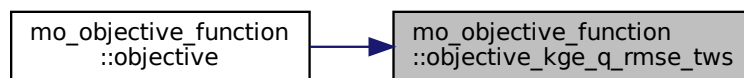
References `create_domain_avg_tws()`, `mo_common_variables::domainmeta`, `mo_common_constants::eps_dp`, `mo_common_mhm_mrm_variables::evalper`, `mo_mrm_objective_function_runoff::extract_runoff()`, `mo_global_variables::l1_twsaobs`, and `mo_common_constants::nodata_dp`.

Referenced by `objective()`.

Here is the call graph for this function:



Here is the caller graph for this function:

**17.56.2.10 objective\_kge\_q\_sm\_corr()**

```

real(dp) function mo_objective_function::objective_kge_q_sm_corr (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function of KGE for runoff and correlation for SM.

Objective function of KGE for runoff and SSE for soil moisture (standarized scores). Further details can be found in the documentation of objective functions '14 - objective\_multiple\_gauges\_kge\_power6' and '13 - objective\_sm\_corr'.

**Parameters**

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

**Returns**

real(dp) :: objective\_kge\_q\_sse\_sm — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Matthias Zink

**Date**

Mar. 2017

Definition at line 1882 of file mo\_objective\_function.f90.

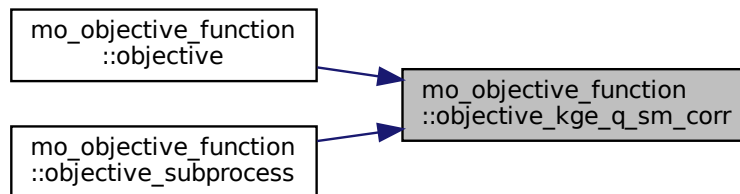
References mo\_common\_variables::domainmeta, mo\_mrm\_objective\_function\_runoff::extract\_runoff(), mo\_global\_variables::l1\_smobs, and mo\_common\_variables::level1.

Referenced by objective(), and objective\_subprocess().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.56.2.11 objective\_master()**

```

real(dp) function, public mo_objective_function::objective_master (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in), optional arg1,
    real(dp), intent(out), optional arg2,
    real(dp), intent(out), optional arg3 )
  
```



Wrapper for objective functions.

The functions sends the parameterset to the subprocess, receives the partial objective and calculates the final objective

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>REAL(dp), DIMENSION(:) :: parameterset</i> |  |
| in  | <i>procedure(eval_interface) :: eval</i>      |  |
| in  | <i>real(dp), optional :: arg1</i>             |  |
| out | <i>real(dp), optional :: arg2</i>             |  |
| out | <i>real(dp), optional :: arg3</i>             |  |

#### Returns

*real(dp) :: objective* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Juliane Mai

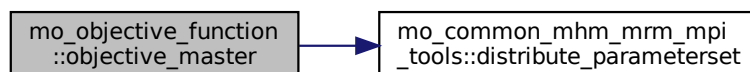
#### Date

Dec 2012

Definition at line 202 of file mo\_objective\_function.f90.

References mo\_common\_mhm\_mrm\_mpi\_tools::distribute\_parameterset(), mo\_common\_variables::domainmeta, mo\_common\_constants::nodata\_dp, and mo\_common\_mhm\_mrm\_variables::opti\_function.

Here is the call graph for this function:



#### 17.56.2.12 objective\_neutrons\_kge\_catchment\_avg()

```

real(dp) function mo_objective_function::objective_neutrons_kge_catchment_avg (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function for neutrons.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Kling-Gupta model efficiency *KGE* of the catchment average neutrons (*N*) is calculated

$$KGE = 1.0 - \sqrt{((1-r)^2 + (1-\alpha)^2 + (1-\beta)^2)}$$

where  $r$  = Pearson product-moment CORRELATION coefficient  $\alpha$  = ratio of simulated mean to observed mean SM  $\beta$  = ratio of simulated standard deviation to observed standard deviation is calculated and the objective function for a given domain  $i$  is

$$\phi_i = 1.0 - KGE_i$$

$\phi_i$  is the objective since we always apply minimization methods. The minimal value of  $\phi_i$  is 0 for the optimal KGE of 1.0. Finally, the overall objective function value  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all domains  $N$ .

$$OF = \sqrt[6]{\sum((1.0 - KGE_i)/N)^6}$$

The observed data L1\_neutronsdata, L1\_neutronsdata\_mask are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

*real(dp) :: objective\_neutrons\_kge\_catchment\_avg* — objective function value (which will be e.g. minimized by an optimization routine)

#### Authors

Martin Schroen

#### Date

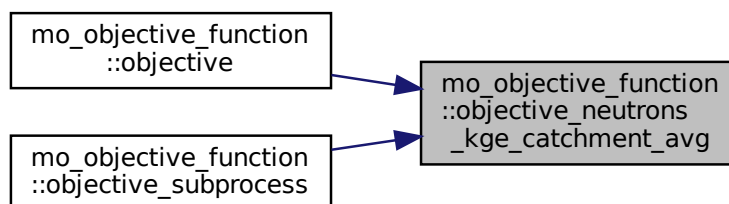
Jun 2015

Definition at line 1633 of file mo\_objective\_function.f90.

References mo\_common\_variables::domainmeta, mo\_global\_variables::l1\_neutronsobs, and mo\_common\_↔ constants::nodata\_dp.

Referenced by objective(), and objective\_subprocess().

Here is the caller graph for this function:



#### 17.56.2.13 objective\_q\_et\_tws\_kge\_catchment\_avg()

```

real(dp) function, dimension(6) mo_objective_function::objective_q_et_tws_kge_catchment_avg (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function for et, tws and q.

The feature of this objective function is the separation of the eval call into four calls, each with another index list. The subroutine eval then only uses the indices from that index list internally instead of having loops over all domains. The integer array domainMetaoptidata decides which indices to use. Therefore the array is split into disjunct subsets, and, if chosen wisely in the namelist, also covers all domains.

With this the eval calls sum up in a way that for each domain eval was called at most once, but for different opti\_data.

#### Authors

Maren Kaluza

#### Date

July 2019

#### Parameters

|    |                     |   |
|----|---------------------|---|
| in | <i>parameterset</i> | the parameterset passed to the eval subroutine        |
| in | <i>eval</i>         | the eval subroutine called by this objective function |

#### Returns

the return value of the objective function. In this case it is an array to provide the possibility to weight the outcome accordingly

domain loop counter

counter for short loops

for sixth root

modelled runoff for a given parameter set

number of all gauges, aquired via runoff

aggregated simulated runoff

measured runoff

mask for measured runoff

kge\_q(nGaugesTotal)

gauges counter

number of q domains

number of et domains

number of tws domains

number of TWS and ET domains (providing both)

index array of ET domains

index array of TWS domains

index array of TWS and ET domains (providing both)

index array of ET domains

simulated et

simulated tws

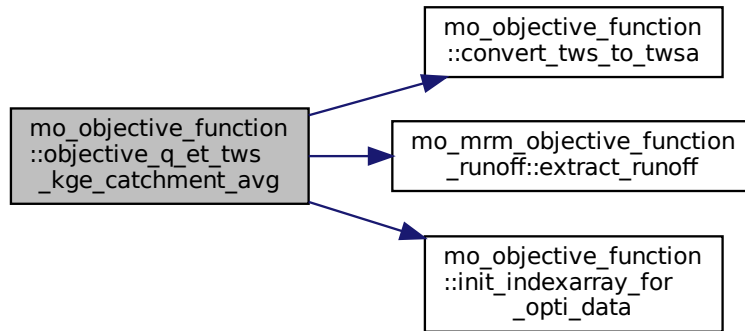
simulated twsa (anomaly)

Definition at line 649 of file mo\_objective\_function.f90.

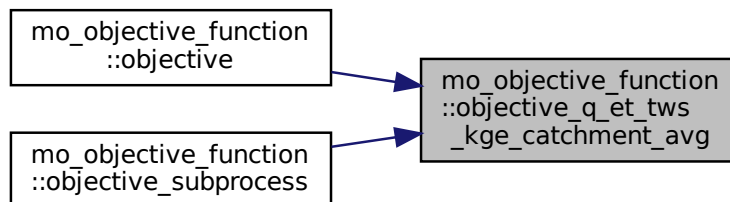
References `convert_tws_to_twsa()`, `mo_common_variables::domainmeta`, `mo_mrm_objective_function_runoff::extract_runoff()`, `init_indexarray_for_opti_data()`, `mo_global_variables::l1_etobs`, `mo_global_variables::l1_twsaobs`, and `mo_common_constants::nodata_dp`.

Referenced by `objective()`, and `objective_subprocess()`.

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.56.2.14 objective\_sm\_corr()

```

real(dp) function mo_objective_function::objective_sm_corr (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function for soil moisture.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore the Pearson correlation between observed and modeled soil moisture on each grid cell  $j$  is compared

$$r_j = r^2(SM_{obs}^j, SM_{sim}^j)$$

where  $r^2$  = Pearson correlation coefficient,  $SM_{obs}$  = observed soil moisture,  $SM_{sim}$  = simulated soil moisture. The observed data  $SM_{obs}$  are global in this module. The the correlation is spatially averaged as

$$\phi_i = \frac{1}{K} \cdot \sum_{j=1}^K r_j$$

where  $K$  denotes the number of valid cells in the study domain. Finally, the overall objective function value  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all domains  $N$ .

$$OF = \sqrt[6]{\sum ((1.0 - \phi_i)/N)^6}$$

The observed data L1\_sm, L1\_sm\_mask are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

#### Returns

real(dp) :: objective\_sm\_corr — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Matthias Zink

#### Date

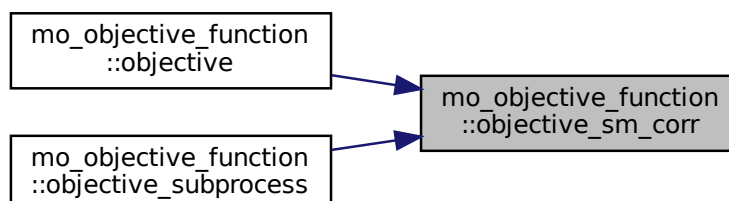
March 2015

Definition at line 986 of file mo\_objective\_function.f90.

References mo\_common\_variables::domainmeta, mo\_global\_variables::l1\_smobs, and mo\_common\_variables↔::level1.

Referenced by objective(), and objective\_subprocess().

Here is the caller graph for this function:



### 17.56.2.15 objective\_sm\_kge\_catchment\_avg()

```
real(dp) function mo_objective_function::objective_sm_kge_catchment_avg (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
```

Objective function for soil moisture.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore, the Kling-Gupta model efficiency  $KGE$  of the catchment average soil moisture (SM) is calculated

$$KGE = 1.0 - \sqrt{((1-r)^2 + (1-\alpha)^2 + (1-\beta)^2)}$$

where  $r$  = Pearson product-moment correlation coefficient  $\alpha$  = ratio of simulated mean to observed mean SM  $\beta$  = ratio of simulated standard deviation to observed standard deviation is calculated and the objective function for a given domain  $i$  is

$$\phi_i = 1.0 - KGE_i$$

$\phi_i$  is the objective since we always apply minimization methods. The minimal value of  $\phi_i$  is 0 for the optimal  $KGE$  of 1.0. Finally, the overall objective function value  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all domains  $N$ .

$$OF = \sqrt[6]{\sum((1.0 - KGE_i)/N)^6}$$

The observed data `L1_sm`, `L1_sm_mask` are global in this module.

#### Parameters

|    |   |  |
|----|---|--|
| in | <code>real(dp), dimension(:) :: parameterset</code> |  |
| in | <code>procedure(eval_interface) :: eval</code>      |  |

#### Returns

`real(dp) :: objective_sm_kge_catchment_avg` — objective function value (which will be e.g. minimized by an optimization routine like DDS)

#### Authors

Matthias Zink

## Date

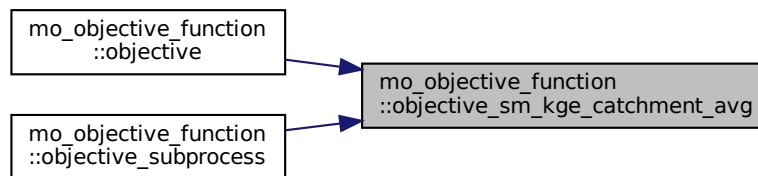
May 2015

Definition at line 504 of file mo\_objective\_function.f90.

References mo\_common\_variables::domainmeta, mo\_global\_variables::l1\_smobs, mo\_common\_variables::level1, and mo\_common\_constants::nodata\_dp.

Referenced by objective(), and objective\_subprocess().

Here is the caller graph for this function:



### 17.56.2.16 objective\_sm\_pd()

```

real(dp) function mo_objective_function::objective_sm_pd (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function for soil moisture.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore the Pattern Dissimilarity (PD) of observed and modeled soil moisture fields is calculated - aim: matching spatial patterns

$$E(t) = PD(SM_{obs}(t), SM_{sim}(t))$$

where  $PD$  = pattern dissimilarity function,  $SM_{obs}$  = observed soil moisture,  $SM_{sim}$  = simulated soil moisture.  $E(t)$  = pattern dissimilarity at timestep  $t$ . The the pattern dissimilarity ( $E$ ) is spatially averaged as

$$\phi_i = \frac{1}{T} \cdot \sum_{t=1}^T E_t$$

where  $T$  denotes the number of time steps. Finally, the overall objective function value  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all domains  $N$ .

$$OF = \sqrt[6]{\sum ((1.0 - \phi_i)/N)^6}$$

The observed data L1\_sm, L1\_sm\_mask are global in this module. The observed data L1\_sm, L1\_sm\_mask are global in this module.

#### Parameters

|    |  |  |
|----|--|--|
| in | real(dp), dimension(:) :: parameterset |  |
| in | procedure(eval_interface) :: eval      |  |

**Returns**

real(dp) :: objective\_sm\_pd — objective function value (which will be e.g. minimized by an optimization routine like DDS)

**Authors**

Matthias Zink

**Date**

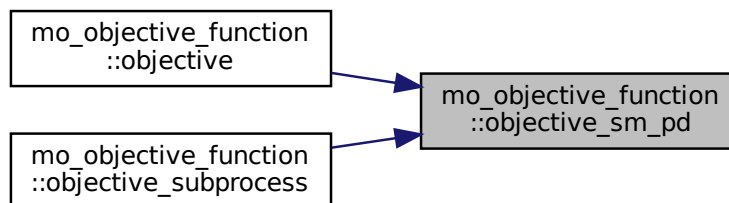
May 2015

Definition at line 1120 of file mo\_objective\_function.f90.

References mo\_common\_variables::domainmeta, mo\_global\_variables::l1\_smobs, mo\_common\_variables::level1, and mo\_common\_constants::nodata\_dp.

Referenced by objective(), and objective\_subprocess().

Here is the caller graph for this function:

**17.56.2.17 objective\_sm\_sse\_standard\_score()**

```

real(dp) function mo_objective_function::objective_sm_sse_standard_score (
    real(dp), dimension(:), intent(in) parameterset,
    procedure(eval_interface), intent(in), pointer eval )
  
```

Objective function for soil moisture.

The objective function only depends on a parameter vector. The model will be called with that parameter vector and the model output is subsequently compared to observed data. Therefore the sum of squared errors (SSE) of the standard score of observed and modeled soil moisture is calculated. The standard score or normalization (anomaly) make the objective function bias insensitive and basically the dynamics of the soil moisture is tried to capture by this objective function.

$$\phi_i = \sum_{j=1}^K \{standard\_score(SM_{obs}(j)) - standard\_score(SM_{sim}(j))\}^2$$

where *standard\_score* = standard score function,  $SM_{obs}$  = observed soil moisture,  $SM_{sim}$  = simulated soil moisture.  $K$  = valid elements in study domain. Finally, the overall objective function value  $OF$  is estimated based on the power-6 norm to combine the  $\phi_i$  from all domains  $N$ .

$$OF = \sqrt[6]{\sum(\phi_i/N)^6}.$$

The observed data L1\_sm, L1\_sm\_mask are global in this module.



## Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: parameterset</i> |  |
| in | <i>procedure(eval_interface) :: eval</i>      |  |

## Returns

*real(dp) :: objective\_sm\_sse\_standard\_score* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

## Authors

Matthias Zink

## Date

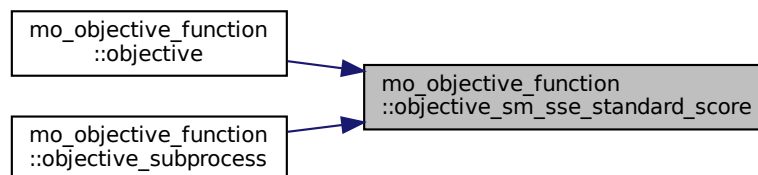
March 2015

Definition at line 1276 of file mo\_objective\_function.f90.

References mo\_common\_variables::domainmeta, mo\_global\_variables::l1\_smobs, and mo\_common\_variables::level1.

Referenced by objective(), and objective\_subprocess().

Here is the caller graph for this function:



## 17.56.2.18 objective\_subprocess()

```

subroutine, public mo_objective_function::objective_subprocess (
    procedure(eval_interface), intent(in), pointer eval,
    real(dp), intent(in), optional arg1,
    real(dp), intent(out), optional arg2,
    real(dp), intent(out), optional arg3 )
  
```

Wrapper for objective functions.

The function receives the parameterset from the master process, selects the objective function case defined in a namelist, i.e. the global variable *opti\_function*. It returns the partial objective function value for a specific parameter set.

## Parameters

|    |   |  |
|----|---|--|
| in | <i>REAL(dp), DIMENSION(:) :: parameterset</i> |  |
|----|---|--|

## Parameters

|     |  |  |
|-----|--|--|
| in  | <i>procedure(eval_interface) :: eval</i> |  |
| in  | <i>real(dp), optional :: arg1</i>        |  |
| out | <i>real(dp), optional :: arg2</i>        |  |
| out | <i>real(dp), optional :: arg3</i>        |  |

## Returns

*real(dp) :: objective* — objective function value (which will be e.g. minimized by an optimization routine like DDS)

## Authors

Juliane Mai

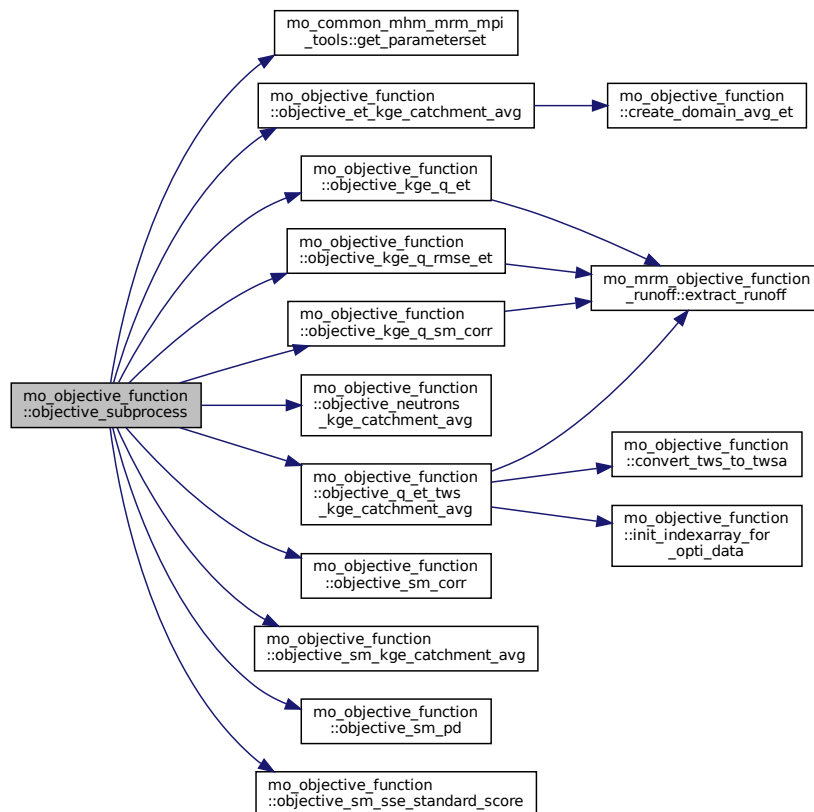
## Date

Dec 2012

Definition at line 354 of file `mo_objective_function.f90`.

References `mo_common_variables::domainmeta`, `mo_common_mhm_mrm_mpi_tools::get_parameterset()`, `mo_common_constants::nodata_dp`, `objective_et_kge_catchment_avg()`, `objective_kge_q_et()`, `objective_kge_q_rmse_et()`, `objective_kge_q_sm_corr()`, `objective_neutrons_kge_catchment_avg()`, `objective_q_et_tws_kge_catchment_avg()`, `objective_sm_corr()`, `objective_sm_kge_catchment_avg()`, `objective_sm_pd()`, `objective_sm_sse_standard_score()`, and `mo_common_mhm_mrm_variables::opti_function`.

Here is the call graph for this function:



## 17.57 mo\_optimization Module Reference

Wrapper subroutine for optimization against runoff and sm.

### Functions/Subroutines

- subroutine, public [optimization](#) (eval, objective, dirConfigOut, funcBest, maskpara)

*Wrapper for optimization.*

#### 17.57.1 Detailed Description

Wrapper subroutine for optimization against runoff and sm.

This module provides a wrapper subroutine for optimization of mRM/mHM against runoff or soil moisture.

#### Authors

Stephan Thober

#### Date

Oct 2015

#### 17.57.2 Function/Subroutine Documentation

##### 17.57.2.1 optimization()

```
subroutine, public mo_optimization::optimization (
    procedure(eval_interface), intent(in), pointer eval,
    procedure(objective_interface), intent(in), pointer objective,
    character(len = *) , intent(in) dirConfigOut,
    real(dp), intent(out) funcBest,
    logical, dimension(:), intent(out), allocatable maskpara )
```

Wrapper for optimization.

This subroutine selects the optimization defined in a namelist, i.e. the global variable *opti\_method*. It return the objective function value for a specific parameter set.

#### Parameters

|     |  |   |
|-----|--|---|
| in  | <i>procedure(eval_interface) :: eval</i>           |   |
| in  | <i>procedure(objective_interface) :: objective</i> | - objective function used in the optimization   |
| in  | <i>character(len = *) :: dirConfigOut</i>          | - directory where to write ascii output   |
| out | <i>real(dp) :: funcbest</i>                        | - best objective function value obtained during optimization  |
| out | <i>logical, dimension(:) :: maskpara</i>           | true = parameter will be optimized = parameter(i,4) = 1<br>false = parameter will not be optimized = parameter(i,4) = 0 |

**Authors**

Matthias Cuntz, Luis Samaniego, Juliane Mai, Matthias Zink and Stephan Thober

**Date**

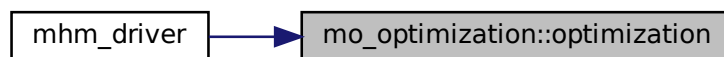
Oct 2015

Definition at line 51 of file mo\_optimization.f90.

References `mo_common_mhm_mrm_variables::dds_r`, `mo_common_variables::domainmeta`, `mo_common_variables::global_parameters`, `mo_common_mhm_mrm_variables::mcmc_error_params`, `mo_common_mhm_mrm_variables::mcmc_opti`, `mo_common_mhm_mrm_variables::niterations`, `mo_common_mhm_mrm_variables::opti_function`, `mo_common_mhm_mrm_variables::opti_method`, `mo_common_mhm_mrm_variables::optimize_restart`, `mo_common_mhm_mrm_variables::sa_temp`, `mo_common_mhm_mrm_variables::sce_ngs`, `mo_common_mhm_mrm_variables::sce_npg`, `mo_common_mhm_mrm_variables::sce_nps`, and `mo_common_mhm_mrm_variables::seed`.

Referenced by `mhm_driver()`.

Here is the caller graph for this function:



## 17.58 mo\_pet Module Reference

Module for calculating reference/potential evapotranspiration [mm d-1].

### Functions/Subroutines

- elemental pure real(dp) function, public [pet\\_hargreaves](#) (HarSamCoeff, HarSamConst, tavg, tmax, tmin, latitude, doy)  
*Reference Evapotranspiration after Hargreaves.*
- elemental pure real(dp) function, public [pet\\_priestly](#) (PriTayParam, Rn, tavg)  
*Reference Evapotranspiration after Priestly-Taylor.*
- elemental pure real(dp) function, public [pet\\_penman](#) (net\_rad, tavg, act\_vap\_pressure, aerodyn\_resistance, bulksurface\_resistance, a\_s, a\_sh)  
*Reference Evapotranspiration after Penman-Monteith.*
- elemental pure real(dp) function, public [extraterr\\_rad\\_approx](#) (doy, latitude)  
*Approximation of extraterrestrial radiation.*
- elemental pure real(dp) function, public [slope\\_satpressure](#) (tavg)  
*slope of saturation vapour pressure curve*
- elemental pure real(dp) function, public [sat\\_vap\\_pressure](#) (tavg)  
*calculation of the saturation vapour pressure*

### 17.58.1 Detailed Description

Module for calculating reference/potential evapotranspiration [mm d-1].

This module calculates PET [mm/d] based on one of the methods

- Hargreaves-Samani (1982)
- Priestly-Taylor (1972)
- Penman-Monteith FAO (1998)

#### Authors

Matthias Zink, Christoph Schneider, Matthias Cuntz

#### Date

Apr 2014

### 17.58.2 Function/Subroutine Documentation

#### 17.58.2.1 extraterr\_rad\_approx()

```
elemental pure real(dp) function, public mo_pet::extraterr_rad_approx (
    integer(i4), intent(in) doy,
    real(dp), intent(in) latitude )
```

Approximation of extraterrestrial radiation.

Approximation of extraterrestrial radiation at the top of the atmosphere  $R_a$  after Duffie and Beckman (1980).  $R_a$  is converted from [ $J m^{-2} d^{-1}$ ] in [ $mm d^{-1}$ ].

$$R_a = \frac{86400}{\pi \cdot \lambda} \cdot E_0 \cdot d_r \cdot (\omega \cdot \sin(\text{latitude}) \cdot \sin(\delta) + \cos(\text{latitude}) \cdot \cos(\delta) \cdot \sin(\omega))$$

where  $E_0 = 1367 J m^{-2} s^{-1}$  is the solar constant and It is dependent on the following sub equations: The relative distance Earth-Sun:

$$d_r = 1 + 0.033 \cdot \cos\left(\frac{2 \cdot \pi \cdot \text{doy}}{365}\right)$$

in which doy is the day of the year. The solar declination [radians] defined by

$$\delta = 0.4093 \cdot \sin\left(\frac{2 \cdot \pi \cdot \text{doy}}{365} - 1.405\right)$$

The sunset hour angle [radians]:

$$\omega = \arccos(-\tan(\text{latitude}) * \tan(\delta))$$

$\lambda = 2.45 \cdot 10^6 J m^{-2} mm^{-1}$  is the latent heat of vaporization.

#### Parameters

|    |                             |                 |
|----|-----------------------------|-----------------|
| in | <i>integer(i4) :: doy</i>   | day of year [-] |
| in | <i>real(dp) :: latitude</i> | latitude [rad]  |

#### Returns

*real(dp) :: extraterr\_rad\_approx* — extraterrestrial radiation approximation [ $W m^{-2}$ ]

**Authors**

Matthias Zink

**Date**

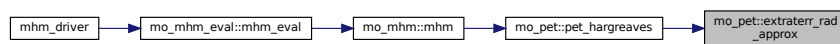
Apr 2014

Definition at line 315 of file mo\_pet.f90.

References mo\_mhm\_constants::duffiedelta1, mo\_mhm\_constants::duffiedelta2, and mo\_mhm\_constants::duffiedr.

Referenced by pet\_hargreaves().

Here is the caller graph for this function:

**17.58.2.2 pet\_hargreaves()**

```

elemental pure real(dp) function, public mo_pet::pet_hargreaves (
    real(dp), intent(in) HarSamCoeff,
    real(dp), intent(in) HarSamConst,
    real(dp), intent(in) tavg,
    real(dp), intent(in) tmax,
    real(dp), intent(in) tmin,
    real(dp), intent(in) latitude,
    integer(i4), intent(in) doy )
  
```

Reference Evapotranspiration after Hargreaves.

Calculates the Reference Evapotranspiration [ $mm\ d^{-1}$ ] based on the Hargreaves-Samani (1982) model for a given cell by applying the equation

$$PET = HarSamCoeff * R_a * (T_{avg} + HarSamConst) * \sqrt{T_{max} - T_{min}}$$

where  $R_a$  [ $W\ m^{-2}$ ] is the incoming solar radiation and  $T_{avg}$ ,  $T_{max}$  and  $T_{min}$  [ $^{\circ}C$ ] are the mean, maximum, and minimum daily temperatures at the given day, respectively.**Note**

Hargreaves, G.H., and Samani, Z.A. (1982). "Estimating potential evapotranspiration." Tech. Note, J. Irrig. and drain. Engrg., ASCE, 108(3):225-230.

**Parameters**

|    |                                |  |
|----|--------------------------------|--|
| in | <i>real(dp) :: HarSamCoeff</i> | coefficient of Hargreaves-Samani equation [-]                |
| in | <i>real(dp) :: HarSamConst</i> | constant of Hargreaves-Samani equation [-]                   |
| in | <i>real(dp) :: tavg</i>        | daily mean temperature [ $^{\circ}C$ ]                       |
| in | <i>real(dp) :: tmax</i>        | daily maximum of temp [ $^{\circ}C$ ]                        |
| in | <i>real(dp) :: tmin</i>        | daily minimum of temp [ $^{\circ}C$ ]                        |
| in | <i>real(dp) :: latitude</i>    | latitude of the cell for $R_a$ estimation [ <i>radians</i> ] |
| in | <i>integer(i4) :: doy</i>      | day of year for $R_a$ estimation                             |

**Returns**

real(dp) :: pet\_hargreaves — Hargreaves-Samani pot. evapotranspiration [mm d-1]

**Authors**

Matthias Zink

**Date**

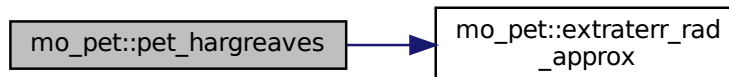
Dec 2012

Definition at line 74 of file mo\_pet.f90.

References extraterr\_rad\_approx().

Referenced by mo\_mhm::mhm().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.58.2.3 pet\_penman()**

```

elemental pure real(dp) function, public mo_pet::pet_penman (
    real(dp), intent(in) net_rad,
    real(dp), intent(in) tavg,
    real(dp), intent(in) act_vap_pressure,
    real(dp), intent(in) aerodyn_resistance,
    real(dp), intent(in) bulksurface_resistance,
    real(dp), intent(in) a_s,
    real(dp), intent(in) a_sh )
  
```

Reference Evapotranspiration after Penman-Monteith.

Calculates the reference evapotranspiration [ $mm\ d^{-1}$ ] based on the Penman-Monteith model for every given cell by applying the equation

$$PET = \frac{1}{\lambda} \cdot \frac{\Delta \cdot R_n + \rho \cdot c_p \cdot (e_s - e) \cdot \frac{a_s h}{r_a}}{\Delta + \gamma \cdot \frac{a_s h}{a_s} \cdot \left(1 + \frac{r_s}{r_a}\right)}$$

where  $R_n$  [ $W m^{-2}$ ] is the net solar radiation,  $\Delta$  [ $kPa K^{-1}$ ] is the slope of the saturation-vapour pressure curve,  $\lambda$  [ $MJ kg^{-1}$ ] is the latent heat of vaporization,  $(e_s - e)$  [ $kPa$ ] is the vapour pressure deficit of the air,  $\rho$  [ $kg m^{-3}$ ] is the mean atmospheric density,  $c_p = 1005.0 J kg^{-1} K^{-1}$  is the specific heat of the air,  $\gamma$  [ $kPa K^{-1}$ ] is the psychrometric constant,  $r_s$  [ $sm^{-1}$ ] is the bulk canopy resistance,  $r_a$  [ $sm^{-1}$ ] is the aerodynamic resistance,  $a_s$  [1] is the fraction of one-sided leaf area covered by stomata (1 if stomata are on one side only, 2 if they are on both sides) and  $a_{sh}$  [-] is the fraction of projected area exchanging sensible heat with the air (2) Implementation refers to the so-called Penman-Montheith equation for transpiration. Adjusting the arguments  $a_{sh}$  and  $a_s$  we obtain the corrected MU equation (for details see Schymanski and Or, 2017). If  $a_{sh} = 1 = a_s$  Penman-Montheith equation for transpiration is preserved. For reproducing characteristics of symmetrical amphistomatous leaves use  $a_{sh} = 2 = a_s$ , in which case the classic PM equation is only missing a factor of 2 in the nominator, as pointed out by Jarvis and McNaughton (1986, Eq. A9). These analytical solutions eliminated the non-linearity problem of the saturation vapour pressure curve, but they do not consider the dependency of the long-wave component of the soil surface or leaf energy balance ( $R_l$ ) on soil or leaf temperature ( $T_l$ ). We assume that net radiation equals the absorbed short-wave radiation, i.e.  $R_N = R_S$  (p.79 in Monteith and Unsworth, 2013).

#### Parameters

|    |   |  |
|----|---|--|
| in | <code>real(dp) :: net_rad</code>                | net radiation [ $W m^{-2}$ ]                                       |
| in | <code>real(dp) :: tavg</code>                   | average daily temperature [ $^{\circ}C$ ]                          |
| in | <code>real(dp) :: act_vap_pressure</code>       | actual vapur pressure [ $kPa$ ]                                    |
| in | <code>real(dp) :: aerodyn_resistance</code>     | aerodynmaical resistance $s m^{-1}$                                |
| in | <code>real(dp) :: bulksurface_resistance</code> | bulk surface resistance $s m^{-1}$                                 |
| in | <code>real(dp) :: a_s</code>                    | fraction of one-sided leaf area covered by stomata 1               |
| in | <code>real(dp) :: a_sh</code>                   | fraction of projected area exchanging sensible heat with the air 1 |

#### Returns

`real(dp) :: pet_penman` — Reference Evapotranspiration [mm d-1]

#### Authors

Matthias Zink

#### Date

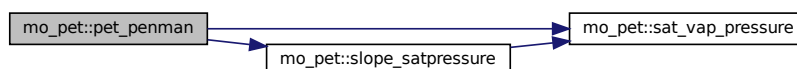
Apr 2014

Definition at line 236 of file `mo_pet.f90`.

References `sat_vap_pressure()`, and `slope_satpressure()`.

Referenced by `mo_mhm::mhm()`.

Here is the call graph for this function:





Here is the caller graph for this function:



#### 17.58.2.4 pet\_priestly()

```

elemental pure real(dp) function, public mo_pet::pet_priestly (
    real(dp), intent(in) PriTayParam,
    real(dp), intent(in) Rn,
    real(dp), intent(in) tavg )
  
```

Reference Evapotranspiration after Priestly-Taylor.

Calculates the Reference Evapotranspiration [ $mm\ d^{-1}$ ] based on the Priestly-Taylor (1972) model for every given cell by applying the equation

$$PET = \alpha * \frac{\Delta}{(\gamma + \Delta)} * R_n$$

where  $R_n$  [ $W\ m^{-2}$ ] is the net solar radiation  $\Delta = f(T_{avg})$  is the slope of the saturation-vapour pressure curve and  $\alpha$  is a emperical coefficient.

##### Parameters

|    |                                |   |
|----|--------------------------------|---|
| in | <i>real(dp) :: PriTayParam</i> | Priestley-Taylor coefficient $\alpha$ [—] |
| in | <i>real(dp) :: Rn</i>          | net solar radiation [ $W\ m^{-2}$ ]       |
| in | <i>real(dp) :: Tavg</i>        |   |

##### Returns

*real(dp) :: pet\_priestly* — Priestley-Taylor pot. evapotranspiration [mm d-1]

##### Authors

Matthias Zink

##### Date

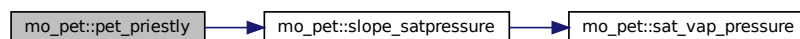
Apr 2014

Definition at line 150 of file mo\_pet.f90.

References slope\_satpressure().

Referenced by mo\_mhm::mhm().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.58.2.5 sat\_vap\_pressure()

```

elemental pure real(dp) function, public mo_pet::sat_vap_pressure (
    real(dp), intent(in) tavg )
  
```

calculation of the saturation vapour pressure

Calculation of the saturation vapour pressure

$$e_s(T_a) = 0.6108 \cdot \exp\left(\frac{17.27 \cdot T_a}{T_a + 237.3}\right)$$

#### Parameters

|    |                         |                    |
|----|-------------------------|--------------------|
| in | <i>real(dp) :: tavg</i> | temperature [degC] |
|----|-------------------------|--------------------|

#### Returns

*real(dp) :: sat\_vap\_pressure* — saturation vapour pressure [kPa]

#### Authors

Matthias Zink

#### Date

Apr 2014

Definition at line 427 of file mo\_pet.f90.

References mo\_mhm\_constants::tetens\_c1, mo\_mhm\_constants::tetens\_c2, and mo\_mhm\_constants::tetens\_c3.

Referenced by pet\_penman(), and slope\_satpressure().

Here is the caller graph for this function:



### 17.58.2.6 slope\_satpressure()

elemental pure real(dp) function, public mo\_pet::slope\_satpressure (  
 real(dp), intent(in) tavg )

slope of saturation vapour pressure curve

slope of saturation vapour pressure curve after Tetens

$$\Delta = \frac{0.6108 * e_s(T_a)}{e^{(2 \cdot \log(T_a + 237.3))}}$$

#### Parameters

|    |                  |                                |
|----|------------------|--------------------------------|
| in | real(dp) :: tavg | average daily temperature [°C] |
|----|------------------|--------------------------------|

#### Returns

real(dp) :: slope\_satpressure — slope of saturation vapour pressure curve [*kPa K*−1]

#### Authors

Matthias Zink

#### Date

Apr 2014

Definition at line 385 of file mo\_pet.f90.

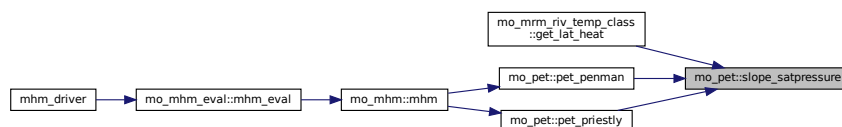
References sat\_vap\_pressure(), mo\_mhm\_constants::satpressureslope1, and mo\_mhm\_constants::tetens\_c3.

Referenced by mo\_mrm\_riv\_temp\_class::get\_lat\_heat(), pet\_penman(), and pet\_priestly().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.59 mo\_prepare\_gridded\_lai Module Reference

Prepare daily LAI fields (e.g., MODIS data) for mHM.

## Functions/Subroutines

- subroutine, public [prepare\\_gridded\\_daily\\_lai\\_data](#) (iDomain, nrows, ncols, mask, LAIPer\_iDomain)  
*Prepare gridded daily LAI data.*
- subroutine, public [prepare\\_gridded\\_mean\\_monthly\\_lai\\_data](#) (iDomain, nrows, ncols, mask)  
*prepare\_gridded\_mean\_monthly\_LAI\_data*

### 17.59.1 Detailed Description

Prepare daily LAI fields (e.g., MODIS data) for mHM.

Prepare daily LAI fields(e.g., MODIS data) for mHM

#### Authors

John Craven & Rohini Kumar

#### Date

Aug 2013

### 17.59.2 Function/Subroutine Documentation

#### 17.59.2.1 [prepare\\_gridded\\_daily\\_lai\\_data\(\)](#)

```
subroutine, public mo_prepare_gridded_lai::prepare_gridded_daily_lai_data (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) nrows,
    integer(i4), intent(in) ncols,
    logical, dimension(:, :) , intent(in) mask,
    type(period), intent(in), optional LAIPer_iDomain )
```

Prepare gridded daily LAI data.

Prepare gridded daily LAI data at Level-0 (e.g., using MODIS datasets)

#### Parameters

|    |   |           |
|----|---|-----------|
| in | <i>integer(i4) :: iDomain, nrows, ncols</i>     | domain Id |
| in | <i>integer(i4) :: iDomain, nrows, ncols</i>     | domain Id |
| in | <i>integer(i4) :: iDomain, nrows, ncols</i>     | domain Id |
| in | <i>logical, dimension(:, :) :: mask</i>         |           |
| in | <i>type(period), optional :: LAIPer_iDomain</i> |           |

#### Authors

John Craven & Rohini Kumar

#### Date

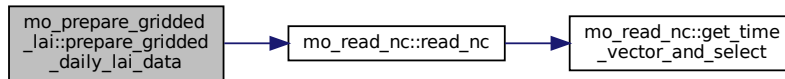
Aug 2013

Definition at line 61 of file mo\_prepare\_gridded\_lai.f90.

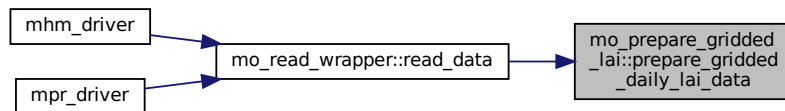
References `mo_mpr_global_variables::dirgridded_lai`, `mo_mpr_global_variables::inputformat_gridded_lai`, `mo_mpr_global_variables::l0_gridded_lai`, `mo_mpr_global_variables::lai_boundaries`, `mo_mpr_global_variables::nlai`, `mo_read_nc::read_nc()`, and `mo_mpr_global_variables::timestep_lai_input`.

Referenced by `mo_read_wrapper::read_data()`.

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.59.2.2 prepare\_gridded\_mean\_monthly\_lai\_data()

```

subroutine, public mo_prepare_gridded_lai::prepare_gridded_mean_monthly_lai_data (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) nrows,
    integer(i4), intent(in) ncols,
    logical, dimension(:, :), intent(in) mask )
  
```

`prepare_gridded_mean_monthly_LAI_data`

Long term mean monthly gridded LAI data at Level-0 (e.g., using MODIS datasets) The netcdf file should contain 12 (calendar months) gridded fields of climatological LAI data at the input L0 data resolution.

#### Parameters

|    |   |           |
|----|---|-----------|
| in | <i>integer(i4) :: iDomain, nrows, ncols</i> | domain Id |
| in | <i>integer(i4) :: iDomain, nrows, ncols</i> | domain Id |
| in | <i>integer(i4) :: iDomain, nrows, ncols</i> | domain Id |
| in | <i>logical, dimension(:, :) :: mask</i>     |           |

#### Authors

Rohini Kumar

## Date

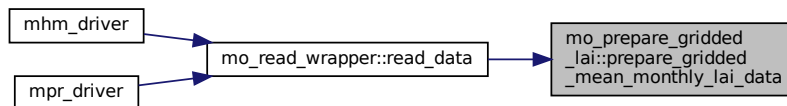
Dec 2016

Definition at line 151 of file mo\_prepare\_gridded\_lai.f90.

References mo\_mpr\_global\_variables::dirgridded\_lai, mo\_mpr\_global\_variables::l0\_gridded\_lai, mo\_mpr\_global\_variables::laiboundaries, and mo\_mpr\_global\_variables::nlai.

Referenced by mo\_read\_wrapper::read\_data().

Here is the caller graph for this function:



## 17.60 mo\_read\_latlon Module Reference

reading latitude and longitude coordinates for each domain

### Functions/Subroutines

- subroutine, public [read\\_latlon](#) (ii, lon\_var\_name, lat\_var\_name, level\_name, level)  
*reads latitude and longitude coordinates*

#### 17.60.1 Detailed Description

reading latitude and longitude coordinates for each domain

TODO: add description

#### Authors

Stephan Thober

#### Date

Nov 2013

#### 17.60.2 Function/Subroutine Documentation

##### 17.60.2.1 read\_latlon()

```

subroutine, public mo_read_latlon::read_latlon (
    integer(i4), intent(in) ii,
    character(*), intent(in) lon_var_name,
    character(*), intent(in) lat_var_name,

```

```
character(*), intent(in) level_name,
type(grid), intent(inout) level )
```

reads latitude and longitude coordinates

reads latitude and longitude coordinates from netcdf file for each domain and appends it to the global variables latitude and longitude.

**Parameters**

|         |                                     |  |
|---------|-------------------------------------|--|
| in      | <i>integer(i4) :: ii</i>            | domain indexFile name of the domains must be xxx_latlon.nc, where xxx is the domain id. Variable names in the netcdf file have to be 'lat' for latitude and 'lon' for longitude. |
| in      | <i>character(*) :: lon_var_name</i> |  |
| in      | <i>character(*) :: lat_var_name</i> |  |
| in      | <i>character(*) :: level_name</i>   |  |
| in, out | <i>type(Grid) :: level</i>          |  |

**Authors**

Stephan Thober

**Date**

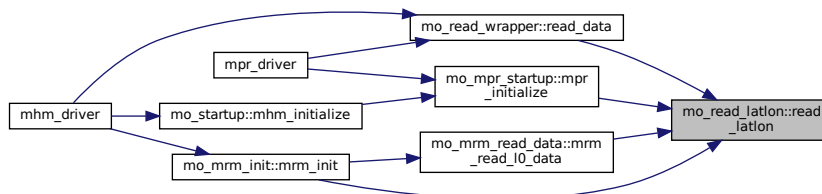
Nov 2013

Definition at line 65 of file mo\_read\_latlon.f90.

References mo\_common\_variables::filelatlon.

Referenced by mo\_mpr\_startup::mpr\_initialize(), mo\_mrm\_init::mrm\_init(), mo\_mrm\_read\_data::mrm\_read\_I0\_data(), and mo\_read\_wrapper::read\_data().

Here is the caller graph for this function:



## 17.61 mo\_read\_lut Module Reference

Routines reading lookup tables (lut).

**Functions/Subroutines**

- subroutine, public [read\\_geof ormation\\_lut](#) (filename, fileunit, nGeo, geo\_unit, geo\_karstic)  
*Reads LUT containing geological formation information.*
- subroutine, public [read\\_lai\\_lut](#) (filename, fileunit, nLAI, LAIIDlist, LAI)  
*Reads LUT containing LAI information.*

### 17.61.1 Detailed Description

Routines reading lookup tables (lut).

This module contains routines reading various lookup tables (lut). (1) LUT containing gauge information. (2) LUT containing geological formation information. (3) LUT containing LAI class information.

#### Authors

Juliane Mai, Matthias Zink

#### Date

Jan 2013

### 17.61.2 Function/Subroutine Documentation

#### 17.61.2.1 read\_geoformation\_lut()

```
subroutine, public mo_read_lut::read_geoformation_lut (
    character(len = *), intent(in) filename,
    integer(i4), intent(in) fileunit,
    integer(i4), intent(out) nGeo,
    integer(i4), dimension(:), intent(out), allocatable geo_unit,
    integer(i4), dimension(:), intent(out), allocatable geo_karstic )
```

Reads LUT containing geological formation information.

The LUT needs to have the following header:

```
nGeo_Formations < Number of lines containing data >
GeoParam(i)   ClassUnit   Karstic   Description
```

The subsequent lines contains the geological formation information:

```
<GeoParam(i)> <ClassUnit_i4> <Karstic_i4> <Description_char>
```

All following lines will be discarded while reading. GeoParam is a running index while ClassUnit is the unit of the map containing the geological formations such that it does not necessarily contains subsequent numbers. The parametrization of this unit is part of the namelist mhm\_parameter.nml under <geoparameter>.

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>character(len = *) :: filename</i>           | File name of LUT                                  |
| in  | <i>integer(i4) :: fileunit</i>                  | Unit to open file                                 |
| out | <i>integer(i4) :: nGeo</i>                      | Number of geological formations                   |
| out | <i>integer(i4), dimension(:) :: geo_unit</i>    | List of id numbers of each geological formations  |
| out | <i>integer(i4), dimension(:) :: geo_karstic</i> | ID of the Karstic formation (0 == does not exist) |

#### Authors

Juliane Mai



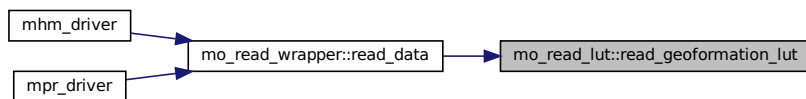
## Date

Jan 2013

Definition at line 71 of file mo\_read\_lut.f90.

Referenced by mo\_read\_wrapper::read\_data().

Here is the caller graph for this function:



## 17.61.2.2 read\_lai\_lut()

```

subroutine, public mo_read_lut::read_lai_lut (
    character(len = *), intent(in) filename,
    integer(i4), intent(in) fileunit,
    integer(i4), intent(out) nLAI,
    integer(i4), dimension(:), intent(out), allocatable LAIIDlist,
    real(dp), dimension(:, :), intent(out), allocatable LAI )
  
```

Reads LUT containing LAI information.

The LUT needs to have the following header:

```

NoLAIclasses <Number of lines containing data>
Id land-use Jan. Feb. Mar. Apr. May Jun. Jul. Aug. Sep. Oct. Nov. Dec.
  
```

The subsequent lines contains the lai class information:

```

<ID_i4> <landuse_char> <val_1_dp> <val_2_dp> <val_3_dp> <val_4_dp> ... <val_12_dp>
  
```

All following lines will be discarded while reading.

## Parameters

|     |   |                                     |
|-----|---|-------------------------------------|
| in  | <i>character(len = *) :: filename</i>         | File name of LUT                    |
| in  | <i>integer(i4) :: fileunit</i>                | Unit to open file                   |
| out | <i>integer(i4) :: nLAI</i>                    | Number of LAI classes               |
| out | <i>integer(i4), dimension(:) :: LAIIDlist</i> | List of ids of LAI classes          |
| out | <i>real(dp), dimension(:, :) :: LAI</i>       | LAI per class (row) and month (col) |

## Authors

Juliane Mai

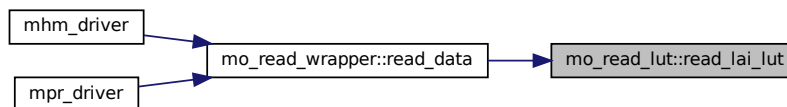
## Date

Jan 2013

Definition at line 151 of file mo\_read\_lut.f90.

Referenced by mo\_read\_wrapper::read\_data().

Here is the caller graph for this function:



## 17.62 mo\_read\_nc Module Reference

Reads forcing input data.

### Functions/Subroutines

- subroutine, public [read\\_nc](#) (folder, nRows, nCols, varName, mask, data, target\_period, lower, upper, nc-timestep, fileName, nocheck, maskout)

*Reads forcing input in NetCDF file format.*

- subroutine, public [read\\_const\\_nc](#) (folder, nRows, nCols, varName, data, fileName)

*Reads time independent forcing input in NetCDF file format.*

- subroutine, public [read\\_weights\\_nc](#) (folder, nRows, nCols, varName, data, mask, lower, upper, nocheck, maskout, fileName)

*Reads weights for meteo forcings input in NetCDF file format.*

- subroutine [get\\_time\\_vector\\_and\\_select](#) (var, fname, inctimestep, time\_start, time\_cnt, target\_period)

*TODO: add description.*

### 17.62.1 Detailed Description

Reads forcing input data.

This module is to read forcing input data contained in netcdf files, e.g. temperature, precipitation, total\_runoff, lai. Timesteps can be hourly, daily, monthly, and annual. The module provides a subroutine for NetCDF files only. First, the dimensions given are cross-checked with header.txt information. Second, the data of the specified period are read from the specified directory. If the optional lower and/or upper bound for the data values is given, the read data are checked for validity. The program is stopped if any value lies out of range.

#### Authors

Juliane Mai

#### Date

Dec 2012

## 17.62.2 Function/Subroutine Documentation

### 17.62.2.1 get\_time\_vector\_and\_select()

```
subroutine mo_read_nc::get_time_vector_and_select (
    type(ncvariable), intent(in) var,
    character(256), intent(in) fname,
    integer(i4), intent(in) inctimestep,
    integer(i4), intent(out) time_start,
    integer(i4), intent(out) time_cnt,
    type(period), intent(in), optional target_period )
```

TODO: add description.

TODO: add description ADDITIONAL INFORMATION get\_time\_vector\_and\_select Extract time vector in unit julian hours and get supposed time step in hours

#### Parameters

|     |  |   |
|-----|--|---|
| in  | <i>type(NcVariable) :: var</i>                 | variable of interest                    |
| in  | <i>character(256) :: fname</i>                 | fname of ncfile for error message       |
| in  | <i>integer(i4) :: inctimestep</i>              | flag for requested time step            |
| out | <i>integer(i4) :: time_start</i>               | time_start index of time selection      |
| out | <i>integer(i4) :: time_cnt</i>                 | time_count of indexes of time selection |
| in  | <i>type(period), optional :: target_period</i> | reference period                        |

#### Authors

Matthias Zink

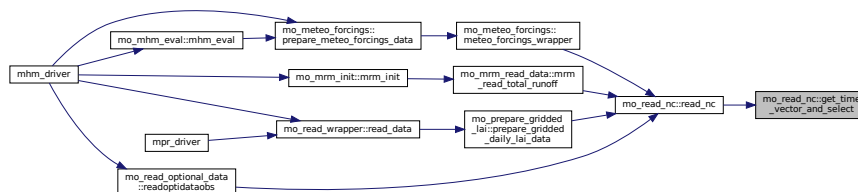
#### Date

Oct 2012

Definition at line 602 of file mo\_read\_nc.f90.

Referenced by read\_nc().

Here is the caller graph for this function:



### 17.62.2.2 read\_const\_nc()

```
subroutine, public mo_read_nc::read_const_nc (
    character(len=*), intent(in) folder,
```

```

integer(i4), intent(in) nRows,
integer(i4), intent(in) nCols,
character(len=*), intent(in) varName,
real(dp), dimension(:, :), intent(out), allocatable data,
character(256), intent(in), optional fileName )

```

Reads time independent forcing input in NetCDF file format.

Reads time independent netCDF forcing files.

First, the dimensions given are cross-checked with header.txt information. Second, the data of the specified period are read from the specified directory. If the optional lower and/or upper bound for the data values is given, the read data are checked for validity. The program is stopped if any value lies out of range.

If the optional argument nocheck is true, the data are not checked for coverage with the input mask. Additionally in this case an mask of vild data points can be received from the routine in maskout.

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>character(len=*) :: folder</i>           | Name of the folder where data are stored        |
| in  | <i>integer(i4) :: nRows</i>                 | Number of datapoints in longitudinal direction  |
| in  | <i>integer(i4) :: nCols</i>                 | Number of datapoints in latitudinal direction   |
| in  | <i>character(len=*) :: varName</i>          | Name of variable name to read                   |
| in  | <i>logical, dimension(:, :) :: mask</i>     | mask of valid data fields                       |
| out | <i>real(dp), dimension(:, :,) :: data</i>   | Data matrix dim_1 = longitude, dim_2 = latitude |
| in  | <i>character(256), optional :: fileName</i> | name of file, defaults to varName data points   |

#### Note

Files have to be called like defined in mo\_files. Furthermore the variable names have to be called like they are defined in the declaration of this subroutine. The NetCDF file has to have 2 dimensions:

1. x, 2. y, It is expected that the variables (especially) within the NetCDF files contain an unit attribute. The timestep has to be equidistant.

#### Author

Lennart Schueler, heavily influenced by read\_nc

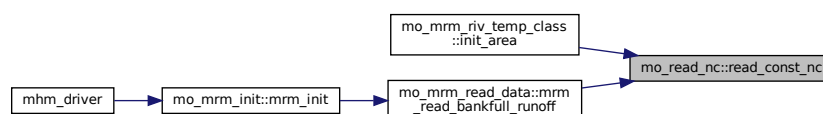
#### Date

May 2018

Definition at line 318 of file mo\_read\_nc.f90.

Referenced by mo\_mrm\_riv\_temp\_class::init\_area(), and mo\_mrm\_read\_data::mrm\_read\_bankfull\_runoff().

Here is the caller graph for this function:



### 17.62.2.3 read\_nc()

```
subroutine, public mo_read_nc::read_nc (
    character(len = *), intent(in) folder,
    integer(i4), intent(in) nRows,
    integer(i4), intent(in) nCols,
    character(len = *), intent(in) varName,
    logical, dimension(:, :), intent(in) mask,
    real(dp), dimension(:, :, :), intent(out), allocatable data,
    type(period), intent(in), optional target_period,
    real(dp), intent(in), optional lower,
    real(dp), intent(in), optional upper,
    integer(i4), intent(in), optional nctimestep,
    character(256), intent(in), optional fileName,
    logical, intent(in), optional nocheck,
    logical, dimension(:, :, :), intent(out), optional, allocatable maskout )
```

Reads forcing input in NetCDF file format.

Reads netCDF forcing files. First, the dimensions given are cross-checked with header.txt information. Second, the data of the specified period are read from the specified directory. If the optional lower and/or upper bound for the data values is given, the read data are checked for validity. The program is stopped if any value lies out of range. If the optional argument nocheck is true, the data are not checked for coverage with the input mask. Additionally in this case an mask of vild data points can be received from the routine in maskout.

#### Parameters

|     |   |  |
|-----|---|--|
| in  | <i>character(len = *) :: folder</i>                     | Name of the folder where data are stored                                       |
| in  | <i>integer(i4) :: nRows</i>                             | Number of datapoints in longitudinal direction                                 |
| in  | <i>integer(i4) :: nCols</i>                             | Number of datapoints in latitudinal direction                                  |
| in  | <i>character(len = *) :: varName</i>                    | Name of variable name to read  |
| in  | <i>logical, dimension(:, :) :: mask</i>                 | mask of valid data fields  |
| out | <i>real(dp), dimension(:, :, :) :: data</i>             | Data matrixdim_1 = longitude, dim_2 = latitude, dim_3 = time                   |
| in  | <i>type(period), optional :: target_period</i>          | Period the data are needed for   |
| in  | <i>real(dp), optional :: lower</i>                      | Lower bound for check of validity of data values                               |
| in  | <i>real(dp), optional :: upper</i>                      | Upper bound for check of validity of data values                               |
| in  | <i>integer(i4), optional :: nctimestep</i>              | timestep in netcdf file  |
| in  | <i>character(256), optional :: fileName</i>             | name of file, defaults to varName  |
| in  | <i>logical, optional :: nocheck</i>                     | .TRUE. if check for nodata values deactivateddefault = .FALSE. - check is done |
| out | <i>logical, dimension(:, :, :), optional :: maskout</i> | ! mask of validdata points   |

#### Authors

Matthias Zink

#### Date

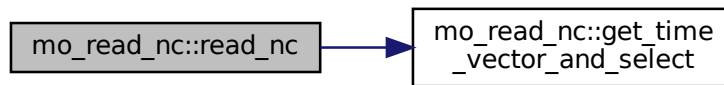
May 2013

Definition at line 85 of file mo\_read\_nc.f90.

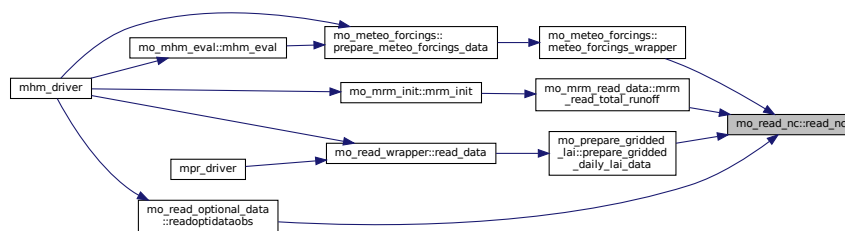
References `get_time_vector_and_select()`.

Referenced by `mo_meteo_forcings::meteo_forcings_wrapper()`, `mo_mrm_read_data::mrm_read_total_runoff()`, `mo_prepare_gridded_lai::prepare_gridded_daily_lai_data()`, and `mo_read_optional_data::readoptidataobs()`.

Here is the call graph for this function:



Here is the caller graph for this function:



#### 17.62.2.4 read\_weights\_nc()

```

subroutine, public mo_read_nc::read_weights_nc (
  character(len = *), intent(in) folder,
  integer(i4), intent(in) nRows,
  integer(i4), intent(in) nCols,
  character(len = *), intent(in) varName,
  real(dp), dimension(:, :, :, :), intent(out), allocatable data,
  logical, dimension(:, :), intent(in) mask,
  real(dp), intent(in), optional lower,
  real(dp), intent(in), optional upper,
  logical, intent(in), optional nocheck,
  logical, dimension(:, :, :, :), intent(out), optional, allocatable maskout,
  character(256), intent(in), optional fileName )
  
```

Reads weights for meteo forcings input in NetCDF file format.

Reads netCDF weight files. First, the dimensions given are cross-checked with header.txt information. If the optional lower and/or upper bound for the data values is given, the read data are checked for validity. The program is stopped if any value lies out of range. If the optional argument nocheck is true, the data are not checked for coverage with the input mask. Additionally in this case an mask of vild data points can be received from the routine in maskout.

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>character(len = *) :: folder</i>            | Name of the folder where data are stored       |
| in | <i>integer(i4) :: nRows</i>                    | Number of datapoints in longitudinal direction |
| in | <i>integer(i4) :: nCols</i>                    | Number of datapoints in latitudinal direction  |
| in | <i>character(len = *) :: varName</i>           | Name of variable name to read                  |
| in | <i>logical, dimension(:, :)</i> :: <i>mask</i> | mask of valid data fields                      |

## Parameters

|     |   |   |
|-----|---|---|
| out | <i>real(dp), dimension(:, :, :, :) :: data</i>            | Data matrix dim_1 = longitude, dim_2 = latitude, dim_3 = months, dim_4 = hours  |
| in  | <i>real(dp), optional :: lower</i>                        | Lower bound for check of validity of data values                                |
| in  | <i>real(dp), optional :: upper</i>                        | Upper bound for check of validity of data values                                |
| in  | <i>logical, optional :: nocheck</i>                       | .TRUE. if check for nodata values deactivated default = .FALSE. - check is done |
| in  | <i>character(256), optional :: fileName</i>               | name of variable, defaults to fileName  |
| out | <i>logical, dimension(:, :, :, ), optional :: maskout</i> | ! mask of valid data points   |

## Authors

Stephan Thober & Matthias Zink

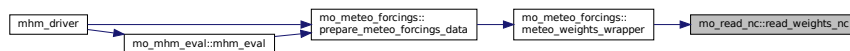
## Date

Jan 2017

Definition at line 422 of file mo\_read\_nc.f90.

Referenced by mo\_meteo\_forcings::meteo\_weights\_wrapper().

Here is the caller graph for this function:



## 17.63 mo\_read\_optional\_data Module Reference

Read optional data for mHM calibration.

### Functions/Subroutines

- subroutine, public [readoptidataobs](#) (iDomain, domainID, L1\_optiObs)  
*Read evapotranspiration data from NetCDF file for calibration.*

#### 17.63.1 Detailed Description

Read optional data for mHM calibration.

Data have to be provided in resolution of the hydrology.

## Authors

Matthias Zink

## Date

Mar 2015

## 17.63.2 Function/Subroutine Documentation

### 17.63.2.1 readoptidataobs()

```
subroutine, public mo_read_optional_data::readoptidataobs (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) domainID,
    type(optidata), intent(inout) Ll_optiObs )
```

Read evapotranspiration data from NetCDF file for calibration.

This routine reads observed evapotranspiration fields which are used for model calibration. The evapotranspiration file is expected to be called "et.nc" with a variable "et" inside. The data are read only for the evaluation period they are intended to be used for calibration. Evapotranspiration data are only read if one of the corresponding objective functions is chosen.

#### Parameters

|    |                               |           |
|----|-------------------------------|-----------|
| in | <i>integer(i4) :: iDomain</i> | domain Id |
|----|-------------------------------|-----------|

#### Authors

Johannes Brenner

#### Date

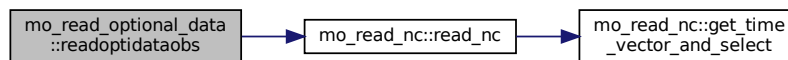
Feb 2017

Definition at line 54 of file mo\_read\_optional\_data.f90.

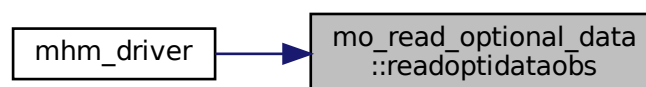
References `mo_common_mhm_mrm_variables::evalper`, `mo_common_variables::level1`, `mo_common_`  
`constants::nodata_dp`, and `mo_read_nc::read_nc()`.

Referenced by `mhm_driver()`.

Here is the call graph for this function:



Here is the caller graph for this function:





## 17.64 mo\_read\_spatial\_data Module Reference

Reads spatial input data.

### Data Types

- interface [read\\_spatial\\_data\\_ascii](#)  
*Reads spatial data files of ASCII format.*

### Functions/Subroutines

- subroutine [read\\_spatial\\_data\\_ascii\\_dp](#) (filename, fileunit, header\_ncols, header\_nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, data, mask)  
*TODO: add description.*
- subroutine [read\\_spatial\\_data\\_ascii\\_i4](#) (filename, fileunit, header\_ncols, header\_nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, data, mask)  
*TODO: add description.*
- subroutine, public [read\\_header\\_ascii](#) (filename, fileunit, header\_ncols, header\_nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, header\_nodata)  
*Reads header lines of ASCII files.*

#### 17.64.1 Detailed Description

Reads spatial input data.

This module is to read spatial input data, e.g. dem, aspect, flow direction. The module provides a subroutine for ASCII files. (Subroutine for NetCDF files will come with release 5.1). The data are read from the specified directory.

#### Authors

Juliane Mai

#### Date

Dec 2012

#### 17.64.2 Function/Subroutine Documentation

##### 17.64.2.1 read\_header\_ascii()

```
subroutine, public mo_read_spatial_data::read_header_ascii (
    character(len = *), intent(in) filename,
    integer(i4), intent(in) fileunit,
    integer(i4), intent(out) header_ncols,
    integer(i4), intent(out) header_nrows,
    real(dp), intent(out) header_xllcorner,
    real(dp), intent(out) header_yllcorner,
    real(dp), intent(out) header_cellsize,
    real(dp), intent(out) header_nodata )
```

Reads header lines of ASCII files.

Reads header lines of ASCII files, e.g. dem, aspect, flow direction.

## Parameters

|     |                                       |                                 |
|-----|---------------------------------------|---------------------------------|
| in  | <i>character(len = *) :: filename</i> | Name of file and its location   |
| in  | <i>integer(i4) :: fileunit</i>        | File unit for open file         |
| out | <i>integer(i4) :: header_nCols</i>    | Reference number of columns     |
| out | <i>integer(i4) :: header_nRows</i>    | Reference number of rows        |
| out | <i>real(dp) :: header_xllcorner</i>   | Reference lower left corner (x) |
| out | <i>real(dp) :: header_yllcorner</i>   | Reference lower left corner (y) |
| out | <i>real(dp) :: header_cellsize</i>    | Reference cell size [m]         |
| out | <i>real(dp) :: header_nodata</i>      | Reference nodata value          |

## Authors

Juliane Mai

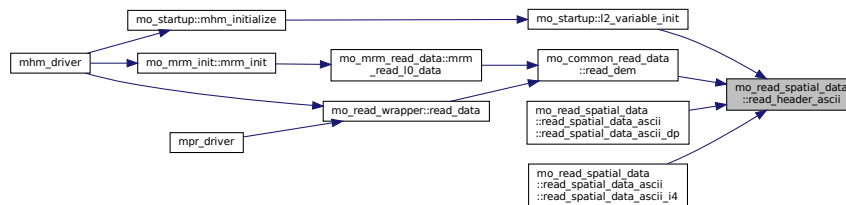
## Date

Jan 2013

Definition at line 375 of file `mo_read_spatial_data.f90`.

Referenced by `mo_startup::l2_variable_init()`, `mo_common_read_data::read_dem()`, `mo_read_spatial_data::read_spatial_data_ascii::read_spatial_data_ascii_dp()`, and `mo_read_spatial_data::read_spatial_data_ascii::read_spatial_data_ascii_i4()`.

Here is the caller graph for this function:



### 17.64.2.2 read\_spatial\_data\_ascii\_dp()

```

subroutine mo_read_spatial_data::read_spatial_data_ascii_dp (
  character(len = *), intent(in) filename,
  integer(i4), intent(in) fileunit,
  integer(i4), intent(in) header_ncols,
  integer(i4), intent(in) header_nrows,
  real(dp), intent(in) header_xllcorner,
  real(dp), intent(in) header_yllcorner,
  real(dp), intent(in) header_cellsize,
  real(dp), dimension(:, :), intent(out), allocatable data,
  logical, dimension(:, :), intent(out), allocatable mask ) [private]

```

TODO: add description.

TODO: add description

## Parameters

|     |  |                                   |
|-----|--|-----------------------------------|
| in  | <i>character(len = *) :: filename</i>    | filename with location            |
| in  | <i>integer(i4) :: fileunit</i>           | unit for opening the file         |
| in  | <i>integer(i4) :: header_nCols</i>       | number of columns of data fields: |
| in  | <i>integer(i4) :: header_nRows</i>       | number of rows of data fields:    |
| in  | <i>real(dp) :: header_xllcorner</i>      | header read in lower left corner  |
| in  | <i>real(dp) :: header_yllcorner</i>      | header read in lower left corner  |
| in  | <i>real(dp) :: header_cellsize</i>       | header read in cellsize           |
| out | <i>real(dp), dimension(:, :) :: data</i> | data                              |
| out | <i>logical, dimension(:, :) :: mask</i>  | mask                              |

## Authors

Robert Schweppe

## Date

Jun 2018

Definition at line 94 of file mo\_read\_spatial\_data.f90.

## 17.64.2.3 read\_spatial\_data\_ascii\_i4()

```
subroutine mo_read_spatial_data::read_spatial_data_ascii_i4 (
    character(len = *), intent(in) filename,
    integer(i4), intent(in) fileunit,
    integer(i4), intent(in) header_ncols,
    integer(i4), intent(in) header_nrows,
    real(dp), intent(in) header_xllcorner,
    real(dp), intent(in) header_yllcorner,
    real(dp), intent(in) header_cellsize,
    integer(i4), dimension(:, :), intent(out), allocatable data,
    logical, dimension(:, :), intent(out), allocatable mask ) [private]
```

TODO: add description.

TODO: add description

## Parameters

|     |   |                                   |
|-----|---|-----------------------------------|
| in  | <i>character(len = *) :: filename</i>       | filename with location            |
| in  | <i>integer(i4) :: fileunit</i>              | unit for opening the file         |
| in  | <i>integer(i4) :: header_nCols</i>          | number of columns of data fields: |
| in  | <i>integer(i4) :: header_nRows</i>          | number of rows of data fields:    |
| in  | <i>real(dp) :: header_xllcorner</i>         | header read in lower left corner  |
| in  | <i>real(dp) :: header_yllcorner</i>         | header read in lower left corner  |
| in  | <i>real(dp) :: header_cellsize</i>          | header read in cellsize           |
| out | <i>integer(i4), dimension(:, :) :: data</i> | data                              |
| out | <i>logical, dimension(:, :) :: mask</i>     | mask                              |

**Authors**

Robert Schweppe

**Date**

Jun 2018

Definition at line 234 of file mo\_read\_spatial\_data.f90.

**17.65 mo\_read\_timeseries Module Reference**

Routines to read files containing timeseries data.

**Functions/Subroutines**

- subroutine, public [read\\_timeseries](#) (filename, fileunit, periodStart, periodEnd, optimize, opti\_function, data, mask, nMeasPerDay)

*Reads time series in ASCII format.*

**17.65.1 Detailed Description**

Routines to read files containing timeseries data.

This routine is reading time series input data for a particular time period. The files need to have a specific header specified in the different routines.

**Authors**

Matthias Zink, Juliane Mai

**Date**

Jan 2013

**17.65.2 Function/Subroutine Documentation****17.65.2.1 read\_timeseries()**

```
subroutine, public mo_read_timeseries::read_timeseries (
    character(len = *), intent(in) filename,
    integer(i4), intent(in) fileunit,
    integer(i4), dimension(3), intent(in) periodStart,
    integer(i4), dimension(3), intent(in) periodEnd,
    logical, intent(in) optimize,
    integer(i4), intent(in) opti_function,
    real(dp), dimension(:), intent(out), allocatable data,
    logical, dimension(:), intent(out), optional, allocatable mask,
    integer(i4), intent(out), optional nMeasPerDay )
```

Reads time series in ASCII format.

Reads time series in ASCII format. Needs specific header lines:

```

<description>
nodata <nodata value>
n <number of measurements per day> measurements per day [1, 1440]
start <YYYY_i4> <MM_i4> <DD_i4> <HH_i4> <MM_i4> (YYYY MM DD HH MM)
end <YYYY_i4> <MM_i4> <DD_i4> <HH_i4> <MM_i4> (YYYY MM DD HH MM)

```

Line 6 is the first line with data in the following format:

```
<YYYY_i4> <MM_i4> <DD_i4> <HH_i4> <MM_i4> <data_dp>
```

The routine checks for missing data points and if data points are equal distanced. The first data point at each day has to be at HH:MM = 00:00.

#### Parameters

|     |   |                                   |
|-----|---|-----------------------------------|
| in  | <i>character(len = *) :: filename</i>           | File name                         |
| in  | <i>integer(i4) :: fileunit</i>                  | Unit to open file                 |
| in  | <i>integer(i4), dimension(3) :: periodStart</i> | Start day of reading (YYYY,MM,DD) |
| in  | <i>integer(i4), dimension(3) :: periodEnd</i>   | End day of reading (YYYY,MM,DD)   |
| in  | <i>logical :: optimize</i>                      | optimization flag                 |
| in  | <i>integer(i4) :: opti_function</i>             |                                   |
| out | <i>real(dp), dimension(:) :: data</i>           | Data vector                       |
| out | <i>logical, dimension(:), optional :: mask</i>  | Mask for nodata values in data    |
| out | <i>integer(i4), optional :: nMeasPerDay</i>     | Number of data points per day     |

#### Authors

Matthias Zink, Juliane Mai

#### Date

Jan 2013

Definition at line 73 of file mo\_read\_timeseries.f90.

Referenced by mo\_mrm\_read\_data::mrm\_read\_discharge().

Here is the caller graph for this function:



## 17.66 mo\_read\_wrapper Module Reference

Wrapper for all reading routines.

### Functions/Subroutines

- subroutine, public [read\\_data](#) (LAIPer)  
*Reads data.*
- subroutine [check\\_consistency\\_lut\\_map](#) (data, lookuptable, filename, unique\_values)  
*Checks if classes in input maps appear in look up tables.*

### 17.66.1 Detailed Description

Wrapper for all reading routines.

This module is to wrap up all reading routines. The general written reading routines are used to store now the read data into global variables.

#### Authors

Juliane Mai, Matthias Zink

#### Date

Jan 2013

### 17.66.2 Function/Subroutine Documentation

#### 17.66.2.1 check\_consistency\_lut\_map()

```
subroutine mo_read_wrapper::check_consistency_lut_map (
    integer(i4), dimension(:), intent(in) data,
    integer(i4), dimension(:), intent(in) lookuptable,
    character(*), intent(in) filename,
    integer(i4), dimension(:), intent(out), optional, allocatable unique_values )
```

Checks if classes in input maps appear in look up tables.

Determines whether a class appearing in the morphological input is occurring in the respective look up table. mHM breaks if inconsistencies are discovered.

#### Parameters

|     |   |                                    |
|-----|---|------------------------------------|
| in  | <i>integer(i4), dimension(:) :: data</i>                    | map of study domain                |
| in  | <i>integer(i4), dimension(:) :: lookuptable</i>             | look up table corresponding to map |
| in  | <i>character(*) :: filename</i>                             | name of the lut file - ERORR warn  |
| out | <i>integer(i4), dimension(:), optional :: unique_values</i> | array of unique values in dataone  |

#### Authors

Matthias Zink

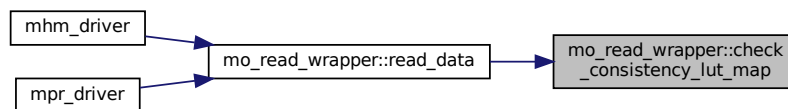
**Date**

Nov 2016

Definition at line 368 of file mo\_read\_wrapper.f90.

Referenced by read\_data().

Here is the caller graph for this function:

**17.66.2.2 read\_data()**

```

subroutine, public mo_read_wrapper::read_data (
    type(period), dimension(:), intent(in), optional LAIPer )
  
```

Reads data.

The namelists are already read by read\_config call. All LUTs are read from their respective directory and information within those files are shared across all domains to be modeled.

**Parameters**

|    |   |
|----|---|
| in | <i>type(period), dimension(:), optional :: LAIPer</i> |
|----|---|

**Authors**

Juliane Mai &amp; Matthias Zink

**Date**

Feb 2013

by default; when iFlag\_soilDB = 0

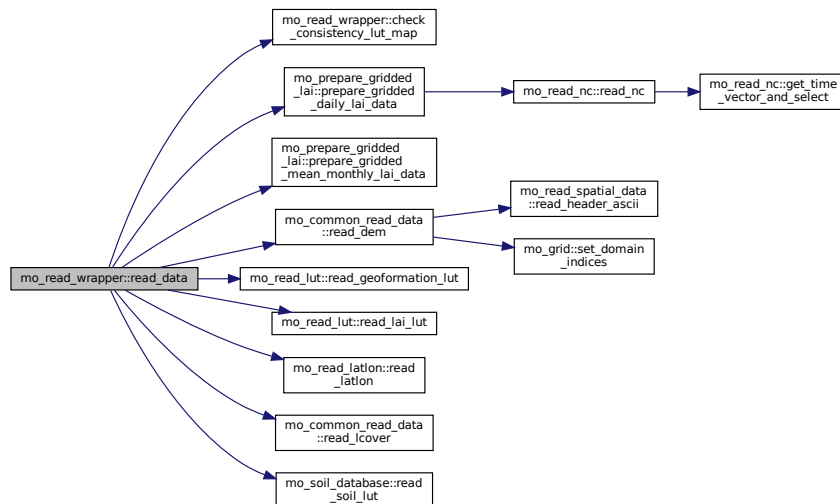
Definition at line 63 of file mo\_read\_wrapper.f90.

References check\_consistency\_lut\_map(), mo\_common\_variables::dircommonfiles, mo\_common\_variables::dirmorpho, mo\_common\_variables::domainmeta, mo\_mpr\_file::file\_aspect, mo\_mpr\_file::file\_geolut, mo\_mpr\_file::file\_hydrogeoclass, mo\_mpr\_file::file\_laiclass, mo\_mpr\_file::file\_lailut, mo\_mpr\_file::file\_slope, mo\_mpr\_file::file\_soil\_database, mo\_mpr\_file::file\_soil\_database\_1, mo\_mpr\_file::file\_soilclass, mo\_mpr\_global\_variables::geounitkar, mo\_mpr\_global\_variables::geounitlist, mo\_common\_variables::global\_parameters, mo\_mpr\_global\_variables::iflag\_soildb, mo\_mpr\_global\_variables::l0\_asp, mo\_mpr\_global\_variables::l0\_geounit, mo\_mpr\_global\_variables::l0\_gridded\_lai, mo\_mpr\_global\_variables::l0\_slope, mo\_mpr\_global\_variables::l0\_soilid, mo\_mpr\_global\_variables::laiboundaries, mo\_mpr\_global\_variables::lailut, mo\_mpr\_global\_variables::laiunitlist, mo\_common\_variables::level0, mo\_mpr\_global\_variables::ngeounits, mo\_mpr\_global\_variables::nlai, mo\_mpr\_global\_variables::nlaiclass, mo\_common\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_prepare\_gridded\_lai::prepare\_gridded\_daily\_lai\_data(), mo\_prepare\_gridded\_lai::prepare\_gridded\_mean\_monthly\_lai\_data(), mo\_common\_variables::processmatrix, mo\_

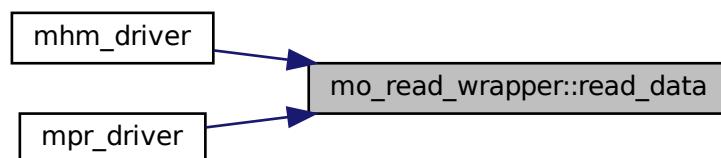
common\_read\_data::read\_dem(), mo\_read\_lut::read\_geoformation\_lut(), mo\_read\_lut::read\_lai\_lut(), mo\_read\_latlon::read\_latlon(), mo\_common\_read\_data::read\_lcover(), mo\_soil\_database::read\_soil\_lut(), mo\_mpr\_global\_variables::soildb, mo\_mpr\_global\_variables::timestep\_lai\_input, mo\_mpr\_file::uaspect, mo\_mpr\_file::ugeolut, mo\_mpr\_file::uhydrogeoclass, mo\_mpr\_file::ulaiclass, mo\_mpr\_file::ulailut, mo\_mpr\_file::uslope, and mo\_mpr\_file::usoilclass.

Referenced by mhm\_driver(), and mpr\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.67 mo\_restart Module Reference

reading and writing states, fluxes and configuration for restart of mHM.

### Data Types

- interface [unpack\\_field\\_and\\_write](#)

*TODO: add description.*



## Functions/Subroutines

- subroutine, public [write\\_restart\\_files](#) (OutFile)  
*write restart files for each domain*
- subroutine, public [read\\_restart\\_states](#) (iDomain, domainID, InFile)  
*reads fluxes and state variables from file*
- subroutine [unpack\\_field\\_and\\_write\\_1d\\_i4](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_1d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_2d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_3d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)

### 17.67.1 Detailed Description

reading and writing states, fluxes and configuration for restart of mHM.

routines are seperated for reading and writing variables for:

- states and fluxes, and
- configuration. Reading of L11 configuration is also seperated from the rest, since it is only required when routing is activated.

#### Authors

Stephan Thober

#### Date

Jul 2013

### 17.67.2 Function/Subroutine Documentation

#### 17.67.2.1 read\_restart\_states()

```
subroutine, public mo_restart::read_restart_states (
    integer(i4), intent(in) iDomain,
    integer(i4), intent(in) domainID,
    character(256), intent(in) InFile )
```

reads fluxes and state variables from file

read fluxes and state variables from given restart directory and initialises all state variables that are initialized in the subroutine initialise, contained in module [mo\\_startup](#).

#### Parameters

|    |                                 |                                     |
|----|---------------------------------|-------------------------------------|
| in | <i>integer(i4) :: iDomain</i>   | number of domains                   |
| in | <i>character(256) :: InFile</i> | Input Path including trailing slash |

#### Authors

Stephan Thober

## Date

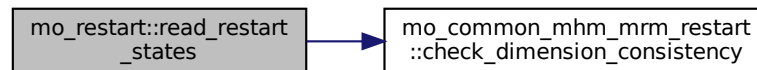
Apr 2013

Definition at line 344 of file mo\_restart.f90.

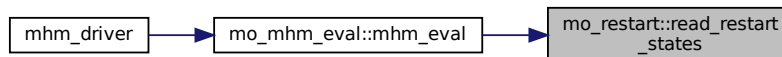
References mo\_common\_mhm\_mrm\_restart::check\_dimension\_consistency(), mo\_mpr\_global\_variables::l1\_aeroresist, mo\_global\_variables::l1\_aetcanopy, mo\_global\_variables::l1\_aetsealed, mo\_global\_variables::l1\_aetsoil, mo\_mpr\_global\_variables::l1\_alpha, mo\_global\_variables::l1\_baseflow, mo\_mpr\_global\_variables::l1\_degday, mo\_mpr\_global\_variables::l1\_degdayinc, mo\_mpr\_global\_variables::l1\_degdaymax, mo\_mpr\_global\_variables::l1\_degdaynopre, mo\_mpr\_global\_variables::l1\_fasp, mo\_global\_variables::l1\_fastrunoff, mo\_mpr\_global\_variables::l1\_froots, mo\_mpr\_global\_variables::l1\_fsealed, mo\_mpr\_global\_variables::l1\_harsamcoeff, mo\_global\_variables::l1\_infilsoil, mo\_global\_variables::l1\_inter, mo\_mpr\_global\_variables::l1\_jarvis\_thresh\_c1, mo\_mpr\_global\_variables::l1\_karstloss, mo\_mpr\_global\_variables::l1\_kbaseflow, mo\_mpr\_global\_variables::l1\_kfastflow, mo\_mpr\_global\_variables::l1\_kperco, mo\_mpr\_global\_variables::l1\_kslowflow, mo\_mpr\_global\_variables::l1\_maxinter, mo\_global\_variables::l1\_melt, mo\_global\_variables::l1\_percol, mo\_mpr\_global\_variables::l1\_petlaicorfactor, mo\_global\_variables::l1\_preeffect, mo\_mpr\_global\_variables::l1\_prietayalpha, mo\_global\_variables::l1\_rain, mo\_global\_variables::l1\_runoffseal, mo\_global\_variables::l1\_satstw, mo\_mpr\_global\_variables::l1\_sealedthresh, mo\_global\_variables::l1\_sealstw, mo\_global\_variables::l1\_slowrunoff, mo\_global\_variables::l1\_snow, mo\_global\_variables::l1\_snowpack, mo\_global\_variables::l1\_soilmoist, mo\_mpr\_global\_variables::l1\_soilmoistexp, mo\_mpr\_global\_variables::l1\_soilmoistfc, mo\_mpr\_global\_variables::l1\_soilmoistsat, mo\_mpr\_global\_variables::l1\_surfresist, mo\_mpr\_global\_variables::l1\_tempthresh, mo\_global\_variables::l1\_throughfall, mo\_global\_variables::l1\_total\_runoff, mo\_global\_variables::l1\_unsatstw, mo\_mpr\_global\_variables::l1\_unsatthresh, mo\_mpr\_global\_variables::l1\_wiltingpoint, mo\_common\_constants::laivarname, mo\_common\_constants::landcoverperiodsvarname, mo\_common\_variables::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::level1, mo\_mpr\_global\_variables::nlai, mo\_common\_variables::nlcoverscene, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_common\_variables::processmatrix, and mo\_common\_constants::soilhorizonsvarname.

Referenced by mo\_mhm\_eval::mhm\_eval().

Here is the call graph for this function:



Here is the caller graph for this function:

**17.67.2.2 unpack\_field\_and\_write\_1d\_dp()**

```

subroutine mo_restart::unpack_field_and_write_1d_dp (
    type(ncdataset), intent(inout) nc,
  
```

```

character(*), intent(in) var_name,
type(ncdimension), dimension(:), intent(in) var_dims,
real(dp), intent(in) fill_value,
real(dp), dimension(:), intent(in) data,
logical, dimension(:, :), intent(in) mask,
character(*), intent(in), optional var_long_name )

```

Definition at line 830 of file mo\_restart.f90.

### 17.67.2.3 unpack\_field\_and\_write\_1d\_i4()

```

subroutine mo_restart::unpack_field_and_write_1d_i4 (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    integer(i4), intent(in) fill_value,
    integer(i4), dimension(:), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name )

```

Definition at line 785 of file mo\_restart.f90.

### 17.67.2.4 unpack\_field\_and\_write\_2d\_dp()

```

subroutine mo_restart::unpack_field_and_write_2d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:, :), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name )

```

Definition at line 875 of file mo\_restart.f90.

### 17.67.2.5 unpack\_field\_and\_write\_3d\_dp()

```

subroutine mo_restart::unpack_field_and_write_3d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:, :, :), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name )

```

Definition at line 932 of file mo\_restart.f90.

### 17.67.2.6 write\_restart\_files()

```

subroutine, public mo_restart::write_restart_files (
    character(256), dimension(:), intent(in) OutFile )

```

write restart files for each domain

write restart files for each domain. For each domain three restart files are written. These are xxx\_states.nc, xxx\_↵ L11\_config.nc, and xxx\_config.nc (xxx being the three digit domain index). If a variable is added here, it should also be added in the read restart routines below.

#### Parameters

|    |  |                             |
|----|--|-----------------------------|
| in | <i>character(256), dimension(:) :: OutFile</i> | Output Path for each domain |
|----|--|-----------------------------|

#### Authors

Stephan Thober

#### Date

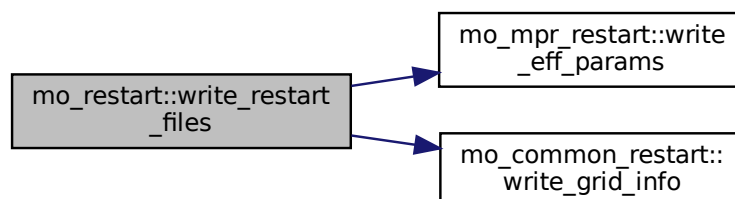
Jun 2014

Definition at line 99 of file mo\_restart.f90.

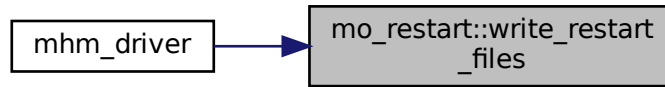
References mo\_common\_variables::domainmeta, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_global\_↵ variables::l1\_aetcanopy, mo\_global\_variables::l1\_aetsealed, mo\_global\_variables::l1\_aetsoil, mo\_global\_↵ variables::l1\_baseflow, mo\_global\_variables::l1\_fastrunoff, mo\_global\_variables::l1\_infilsoil, mo\_global\_variables\_↵ ::l1\_inter, mo\_global\_variables::l1\_melt, mo\_global\_variables::l1\_percol, mo\_global\_variables::l1\_preeffect, mo\_↵ global\_variables::l1\_rain, mo\_global\_variables::l1\_runoffseal, mo\_global\_variables::l1\_satstw, mo\_global\_↵ variables::l1\_sealstw, mo\_global\_variables::l1\_slowrunoff, mo\_global\_variables::l1\_snow, mo\_global\_variables\_↵ ::l1\_snowpack, mo\_global\_variables::l1\_soilmoist, mo\_global\_variables::l1\_throughfall, mo\_global\_variables::l1\_↵ total\_runoff, mo\_global\_variables::l1\_unsatstw, mo\_mpr\_global\_variables::laiboundaries, mo\_common\_variables\_↵ ::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::level1, mo\_mpr\_global\_variables\_↵ ::nlai, mo\_common\_variables::nlcoverscene, mo\_common\_constants::nodata\_dp, mo\_mpr\_global\_variables\_↵ ::nsoilhorizons\_mhm, mo\_mpr\_restart::write\_eff\_params(), and mo\_common\_restart::write\_grid\_info().

Referenced by mhm\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.68 mo\_runoff Module Reference

Runoff generation for the unsaturated zone, saturated zone (or groundwater zone), and runoff accumulation.

### Functions/Subroutines

- subroutine, public [runoff\\_unsat\\_zone](#) (k1, kp, k0, alpha, karst\_loss, pefec\_soil, unsat\_thresh, sat\_storage, unsat\_storage, slow\_interflow, fast\_interflow, perc)  
*Runoff generation for the saturated zone.*
- subroutine, public [runoff\\_sat\\_zone](#) (k2, sat\_storage, baseflow)  
*Runoff generation for the saturated zone.*
- subroutine, public [l1\\_total\\_runoff](#) (fSealed\_area\_fraction, fast\_interflow, slow\_interflow, baseflow, direct\_runoff, total\_runoff)  
*total runoff accumulation at level 1*

### 17.68.1 Detailed Description

Runoff generation for the unsaturated zone, saturated zone (or groundwater zone), and runoff accumulation.

This module generates the runoff for the unsaturated and saturated zones and provides runoff accumulation.

#### Authors

Vladyslav Prykhodko

#### Date

Dec 2012

### 17.68.2 Function/Subroutine Documentation

#### 17.68.2.1 l1\_total\_runoff()

```

subroutine, public mo_runoff::l1_total_runoff (
    real(dp), intent(in) fSealed_area_fraction,
    real(dp), intent(in) fast_interflow,
    real(dp), intent(in) slow_interflow,
    real(dp), intent(in) baseflow,

```

```

real(dp), intent(in) direct_runoff,
real(dp), intent(out) total_runoff )

```

total runoff accumulation at level 1

Accumulates runoff.

$$q_T = (q_0 + q_1 + q_2) * (1 - f_{Sealed}) + q_D * f_{Sealed}$$

, where  $f_{Sealed}$  is the fraction of sealed area.

#### Parameters

|     |  |   |
|-----|--|---|
| in  | <i>REAL(dp) :: fSealed_area_fraction</i> | sealed area fraction [1]                            |
| in  | <i>REAL(dp) :: fast_interflow</i>        | $q_0$ Fast runoff component [mm TS-1]               |
| in  | <i>REAL(dp) :: slow_interflow</i>        | $q_1$ Slow runoff component [mm TS-1]               |
| in  | <i>REAL(dp) :: baseflow</i>              | $q_2$ Baseflow [mm TS-1]                            |
| in  | <i>REAL(dp) :: direct_runoff</i>         | $q_D$ Direct runoff from impervious areas [mm TS-1] |
| out | <i>REAL(dp) :: total_runoff</i>          | $q_T$ Generated runoff [mm TS-1]                    |

#### Authors

Vladyslav Prykhodko

#### Date

Dec 2012

Definition at line 251 of file mo\_runoff.f90.

Referenced by mo\_mhm::mhm().

Here is the caller graph for this function:



#### 17.68.2.2 runoff\_sat\_zone()

```

subroutine, public mo_runoff::runoff_sat_zone (
real(dp), intent(in) k2,
real(dp), intent(inout) sat_storage,
real(dp), intent(out) baseflow )

```

Runoff generation for the saturated zone.

Calculates the runoff generation for the saturated zone. If the level of the ground water reservoir is zero, then the baseflow is also zero. If the level of the ground water reservoir is greater than zero, then the baseflow is equal to baseflow recession coefficient times the level of the ground water reservoir, which will be then reduced by the value of baseflow.

#### Parameters

|    |                       |                                       |
|----|-----------------------|---------------------------------------|
| in | <i>REAL(dp) :: k2</i> | Baseflow recession coefficient [TS-1] |
|----|-----------------------|---------------------------------------|

## Parameters

|         |                                |                          |
|---------|--------------------------------|--------------------------|
| in, out | <i>REAL(dp) :: sat_storage</i> | Groundwater storage [mm] |
| out     | <i>REAL(dp) :: baseflow</i>    | Baseflow [mm TS-1]       |

## Authors

Vladyslav Prykhodko

## Date

Dec 2012

Definition at line 194 of file mo\_runoff.f90.

Referenced by mo\_mhm::mhm().

Here is the caller graph for this function:



## 17.68.2.3 runoff\_unsat\_zone()

```

subroutine, public mo_runoff::runoff_unsat_zone (
    real(dp), intent(in) k1,
    real(dp), intent(in) kp,
    real(dp), intent(in) k0,
    real(dp), intent(in) alpha,
    real(dp), intent(in) karst_loss,
    real(dp), intent(in) pefec_soil,
    real(dp), intent(in) unsat_thresh,
    real(dp), intent(inout) sat_storage,
    real(dp), intent(inout) unsat_storage,
    real(dp), intent(out) slow_interflow,
    real(dp), intent(out) fast_interflow,
    real(dp), intent(out) perc )
  
```

Runoff generation for the saturated zone.

Calculates the runoff generation for the unsaturated zone. Calculates percolation, interflow and baseflow. Updates upper soil and groundwater storages.

## Parameters

|    |                               |   |
|----|-------------------------------|---|
| in | <i>REAL(dp) :: k1</i>         | Recession coefficient of the upper reservoir, lower outlet [TS-1] |
| in | <i>REAL(dp) :: kp</i>         | Percolation coefficient [TS-1]                                    |
| in | <i>REAL(dp) :: k0</i>         | Recession coefficient of the upperreservoir, upper outlet [TS-1]  |
| in | <i>REAL(dp) :: alpha</i>      | Exponent for the upper reservoir [-]                              |
| in | <i>REAL(dp) :: karst_loss</i> | Karstic percolation loss [-]                                      |
| in | <i>REAL(dp) :: pefec_soil</i> | Input to the soil layer [mm]                                      |

## Parameters

|         |                                   |  |
|---------|-----------------------------------|--|
| in      | <i>REAL(dp) :: unsat_thresh</i>   | Threshold water depth in upper reservoir(for Runoff contribution) [mm] |
| in, out | <i>REAL(dp) :: sat_storage</i>    | Groundwater storage [mm]   |
| in, out | <i>REAL(dp) :: unsat_storage</i>  | Upper soil storage [mm]  |
| out     | <i>REAL(dp) :: slow_interflow</i> | Slow runoff component [mm TS-1]  |
| out     | <i>REAL(dp) :: fast_interflow</i> | Fast runoff component [mm TS-1]  |
| out     | <i>REAL(dp) :: perc</i>           | Percolation [mm TS-1]  |

## Authors

Vladyslav Prykhodko

## Date

Dec 2012

Definition at line 77 of file mo\_runoff.f90.

References mo\_common\_constants::eps\_dp.

Referenced by mo\_mhm::mhm().

Here is the caller graph for this function:



## 17.69 mo\_snow\_accum\_melt Module Reference

Snow melting and accumulation.

### Functions/Subroutines

- subroutine, public [snow\\_accum\\_melt](#) (deg\_day\_incr, deg\_day\_max, deg\_day\_noprec, prec, temperature, temperature\_thresh, thrfall, snow\_pack, deg\_day, melt, prec\_effect, rain, snow)

*Snow melting and accumulation.*

#### 17.69.1 Detailed Description

Snow melting and accumulation.

This module calculates snow melting and accumulation.

## Authors

Vladyslav Prykhodko

## Date

Dec 2012



## 17.69.2 Function/Subroutine Documentation

### 17.69.2.1 snow\_accum\_melt()

```

subroutine, public mo_snow_accum_melt::snow_accum_melt (
    real(dp), intent(in) deg_day_incr,
    real(dp), intent(in) deg_day_max,
    real(dp), intent(in) deg_day_noprec,
    real(dp), intent(in) prec,
    real(dp), intent(in) temperature,
    real(dp), intent(in) temperature_thresh,
    real(dp), intent(in) thrfall,
    real(dp), intent(inout) snow_pack,
    real(dp), intent(out) deg_day,
    real(dp), intent(out) melt,
    real(dp), intent(out) prec_effect,
    real(dp), intent(out) rain,
    real(dp), intent(out) snow )

```

Snow melting and accumulation.

Separates throughfall into rain and snow by comparing the temperature with the treshhold. by comparing the temperature with the treshhold. Calculates degree daily factor. Calculates snow melting rates. Calculates snow, rain and effective precipitation depth and snow pack.

#### Parameters

|         |                                       |   |
|---------|---------------------------------------|---|
| in      | <i>REAL(dp) :: deg_day_incr</i>       | Increase of the Degree-day factor per mm of increase in precipitation [s-1 degreeC-1] |
| in      | <i>REAL(dp) :: deg_day_max</i>        | Maximum Degree-day factor [m-1 degreeC-1]   |
| in      | <i>REAL(dp) :: deg_day_noprec</i>     | Degree-day factor with no precipitation [m-1 degreeC-1]                               |
| in      | <i>REAL(dp) :: prec</i>               | Daily mean precipitation [m]  |
| in      | <i>REAL(dp) :: temperature</i>        | Daily mean temperature [degreeC]  |
| in      | <i>REAL(dp) :: temperature_thresh</i> | Threshold temperature for snow/rain [degreeC]   |
| in      | <i>REAL(dp) :: thrfall</i>            | Throughfall [m TS-1]  |
| in, out | <i>REAL(dp) :: snow_pack</i>          | Snow pack [m]   |
| out     | <i>REAL(dp) :: deg_day</i>            | Degree-day factor [m s-1 degreeC-1]   |
| out     | <i>REAL(dp) :: melt</i>               | Melting snow depth [m TS-1]   |
| out     | <i>REAL(dp) :: prec_effect</i>        | Effective precipitation depth (snow melt + rain) [m]                                  |
| out     | <i>REAL(dp) :: rain</i>               | Rain precipitation depth [m]  |
| out     | <i>REAL(dp) :: snow</i>               | Snow precipitation depth [m]  |

#### Authors

Vladyslav Prykhodko

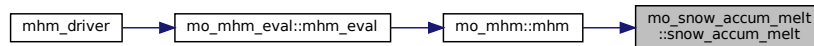
#### Date

Dec 2012

Definition at line 71 of file mo\_snow\_accum\_melt.f90.

Referenced by mo\_mhm::mhm().

Here is the caller graph for this function:



## 17.70 mo\_soil\_database Module Reference

Generating soil database from input file.

### Functions/Subroutines

- subroutine, public [read\\_soil\\_lut](#) (filename)  
*Reads the soil LUT file.*
- subroutine, public [generate\\_soil\\_database](#)  
*Generates soil database.*

#### 17.70.1 Detailed Description

Generating soil database from input file.

This module provides the routines for generating the soil database for mHM from an ASCII input file. One routine *read\_soil\_LUT* reads a soil LookUpTable, performs some consistency checks and returns an initial soil database. The second routine *generate\_soil\_database* calculates based on the initial one the proper soil database.

#### Authors

Juliane Mai

#### Date

Dec 2012

#### 17.70.2 Function/Subroutine Documentation

##### 17.70.2.1 generate\_soil\_database()

```
subroutine, public mo_soil_database::generate_soil_database
```

Generates soil database.

Calculates the proper soil database using the initialized soil database from *read\_soil\_LUT*.

#### Authors

Juliane Mai

**Date**

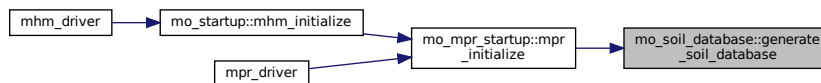
Dec 2012

Definition at line 380 of file mo\_soil\_database.f90.

References mo\_mpr\_global\_variables::horizonddepth\_mhm, mo\_mpr\_global\_variables::iflag\_soildb, mo\_common\_↔  
\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_↔  
\_mpr\_global\_variables::nsoiltypes, and mo\_mpr\_global\_variables::soildb.

Referenced by mo\_mpr\_startup::mpr\_initialize().

Here is the caller graph for this function:

**17.70.2.2 read\_soil\_lut()**

```

subroutine, public mo_soil_database::read_soil_lut (
    character(len = *), intent(in) filename )
  
```

Reads the soil LUT file.

Reads the soil LookUpTable file and checks for consistency.

**Parameters**

|    |                                       |                          |
|----|---------------------------------------|--------------------------|
| in | <i>character(len = *) :: filename</i> | filename of the soil LUT |
|----|---------------------------------------|--------------------------|

**Authors**

Juliane Mai

## Date

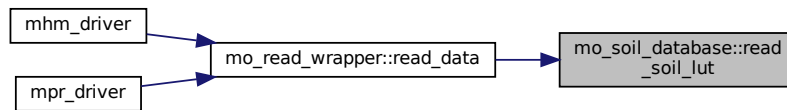
Dec 2012

Definition at line 61 of file mo\_soil\_database.f90.

References mo\_common\_constants::eps\_dp, mo\_mpr\_global\_variables::horizondepth\_mhm, mo\_mpr\_global\_variables::iflag\_soildb, mo\_mpr\_constants::nlcover\_class, mo\_common\_constants::nodata\_dp, mo\_common\_constants::nodata\_i4, mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_mpr\_global\_variables::nsoiltypes, mo\_mpr\_global\_variables::soildb, mo\_mpr\_global\_variables::tillagedepth, and mo\_mpr\_file::usoil\_database.

Referenced by mo\_read\_wrapper::read\_data().

Here is the caller graph for this function:



## 17.71 mo\_soil\_moisture Module Reference

Soil moisture of the different layers.

### Functions/Subroutines

- subroutine, public [soil\\_moisture](#) (processCase, frac\_sealed, water\_thresh\_sealed, pet, evap\_coeff, soil\_moist\_sat, frac\_roots, soil\_moist\_FC, wilting\_point, soil\_moist\_exponen, jarvis\_thresh\_c1, aet\_canopy, prec\_effec, runoff\_sealed, storage\_sealed, infiltration, soil\_moist, aet, aet\_sealed)

*Soil moisture in different soil horizons.*

- elemental pure real(dp) function, public [feddes\\_et\\_reduction](#) (soil\_moist, soil\_moist\_FC, wilting\_point, frac\_roots)

*stress factor for reducing evapotranspiration based on actual soil moisture*

- elemental pure real(dp) function, public [jarvis\\_et\\_reduction](#) (soil\_moist, soil\_moist\_sat, wilting\_point, frac\_roots, jarvis\_thresh\_c1)

*stress factor for reducing evapotranspiration based on actual soil moisture*

### 17.71.1 Detailed Description

Soil moisture of the different layers.

Soil moisture in the different layers is calculated with infiltration as  $(\theta/\theta_{sat})^\beta$ . Then evapotranspiration is calculated from PET with a soil water stress factor  $f_{SM}$  either using the Feddes equation - precessCase(1):  $f_{SM} = \frac{\theta - \theta_{pwp}}{\theta_{fc} - \theta_{pwp}}$  or using the Jarvis equation - precessCase(1):  $f_{SM} = \frac{1}{\theta_{stress-index-C1}} \frac{\theta - \theta_{pwp}}{\theta_{sat} - \theta_{pwp}}$ .

#### Authors

Matthias Cuntz, Luis Samaniego

#### Date

Dec 2012

## 17.71.2 Function/Subroutine Documentation

### 17.71.2.1 feddes\_et\_reduction()

```
elemental pure real(dp) function, public mo_soil_moisture::feddes_et_reduction (
    real(dp), intent(in) soil_moist,
    real(dp), intent(in) soil_moist_FC,
    real(dp), intent(in) wilting_point,
    real(dp), intent(in) frac_roots )
```

stress factor for reducing evapotranspiration based on actual soil moisture

Potential evapotranspiration is reduced to 0 if SM is lower PWP. PET is equal fraction of roots if soil moisture is exceeding field capacity. If soil moisture is in between PWP and FC PET is reduced by fraction of roots times a stress factor. The ET reduction factor  $f$  is estimated as

$$f = \begin{cases} f_{roots} & \text{if } \theta \geq \theta_{fc} \\ f_{roots} \cdot \frac{\theta - \theta_{pwp}}{\theta_{fc} - \theta_{pwp}} & \text{if } \theta < \theta_{fc} \\ 0 & \text{if } \theta < \theta_{pwp} \end{cases}$$

#### Parameters

|    |                                  |   |
|----|----------------------------------|---|
| in | <i>real(dp) :: soil_moist</i>    | Soil moisture of each horizon [mm]                  |
| in | <i>real(dp) :: soil_moist_FC</i> | Soil moisture below which actual ET is reduced [mm] |
| in | <i>real(dp) :: wilting_point</i> | Permanent wilting point                             |
| in | <i>real(dp) :: frac_roots</i>    | Fraction of Roots in soil horizon is reduced [mm]   |

#### Returns

*real(dp) :: feddes\_et\_reduction*; et reduction factor

#### Authors

Matthias Cuntz, Cuenejd Demirel, Matthias Zink

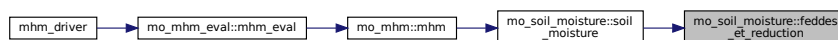
#### Date

March 2017

Definition at line 334 of file mo\_soil\_moisture.f90.

Referenced by soil\_moisture().

Here is the caller graph for this function:



### 17.71.2.2 jarvis\_et\_reduction()

```

elemental pure real(dp) function, public mo_soil_moisture::jarvis_et_reduction (
    real(dp), intent(in) soil_moist,
    real(dp), intent(in) soil_moist_sat,
    real(dp), intent(in) wilting_point,
    real(dp), intent(in) frac_roots,
    real(dp), intent(in) jarvis_thresh_c1 )

```

stress factor for reducing evapotranspiration based on actual soil moisture

The soil moisture stress factor is estimated based on the normalized soil water content. The normalized soil water content  $\theta_{norm}$  is estimated as:

$$\theta_{norm} = \frac{\theta - \theta_{pwp}}{\theta_{sat} - \theta_{pwp}}$$

The ET reduction factor  $f$  is estimated as

$$f = \begin{cases} f_{roots} & \text{if } \theta_{norm} \geq \text{jarvis\_sm\_threshold\_c1} \\ f_{roots} \frac{\theta_{norm}}{\text{jarvis\_sm\_threshold\_c1}} & \text{if } \theta_{norm} < \text{jarvis\_sm\_threshold\_c1} \end{cases}$$

#### Parameters

|    |                                     |   |
|----|-------------------------------------|---|
| in | <i>real(dp) :: soil_moist</i>       | Soil moisture of each horizon [mm]                |
| in | <i>real(dp) :: soil_moist_sat</i>   | saturated Soil moisture content [mm]              |
| in | <i>real(dp) :: wilting_point</i>    | Permanent wilting point                           |
| in | <i>real(dp) :: frac_roots</i>       | Fraction of Roots in soil horizon is reduced [mm] |
| in | <i>real(dp) :: jarvis_thresh_c1</i> | parameter C1 from Jarvis formulation              |

#### Returns

*real(dp) :: jarvis\_et\_reduction*; et reduction factor

#### Authors

Cueneyd Demirel, Matthias Zink

#### Date

March 2017

Definition at line 406 of file mo\_soil\_moisture.f90.

Referenced by soil\_moisture().

Here is the caller graph for this function:



## 17.71.2.3 soil\_moisture()

```

subroutine, public mo_soil_moisture::soil_moisture (
  integer(i4), intent(in) processCase,
  real(dp), intent(in) frac_sealed,
  real(dp), intent(in) water_thresh_sealed,
  real(dp), intent(in) pet,
  real(dp), intent(in) evap_coeff,
  real(dp), dimension(:), intent(in) soil_moist_sat,
  real(dp), dimension(:), intent(in) frac_roots,
  real(dp), dimension(:), intent(in) soil_moist_FC,
  real(dp), dimension(:), intent(in) wilting_point,
  real(dp), dimension(:), intent(in) soil_moist_exponen,
  real(dp), intent(in) jarvis_thresh_c1,
  real(dp), intent(in) aet_canopy,
  real(dp), intent(in) prec_effec,
  real(dp), intent(inout) runoff_sealed,
  real(dp), intent(inout) storage_sealed,
  real(dp), dimension(size(soil_moist_sat, 1)), intent(inout) infiltration,
  real(dp), dimension(size(soil_moist_sat, 1)), intent(inout) soil_moist,
  real(dp), dimension(size(soil_moist_sat, 1)), intent(out) aet,
  real(dp), intent(out) aet_sealed )

```

Soil moisture in different soil horizons.

Infiltration  $I$  from one layer  $k - 1$  to the next  $k$  on pervious areas is calculated as (omit  $t$ )

$$I[k] = I[k - 1](\theta[k]/\theta_{sat}[k])^{\beta[k]}$$

Then soil moisture can be calculated as (omit  $k$ )

$$\theta[t] = \theta[t - 1] + I[t] - ET[t]$$

with  $ET$  (omit  $[k, t]$ ) being

$$ET = f_{roots} \cdot f_{SM} \cdot PET$$

## Parameters

|    |   |  |
|----|---|--|
| in | <i>integer(i4) :: processCase</i>                   | 1 - Feddes equation for PET reduction<br>2 - Jarvis equation for PET reduction<br>3 - Jarvis equation for PET reduction and FC dependency on root fraction coefficient |
| in | <i>real(dp) :: frac_sealed</i>                      | Fraction of sealed area  |
| in | <i>real(dp) :: water_thresh_sealed</i>              | Threshold water depth in impervious areas [mm TS-1]  |
| in | <i>real(dp) :: pet</i>                              | Reference evapotranspiration [mm TS-1]   |
| in | <i>real(dp) :: evap_coeff</i>                       | Evaporation coefficient for free-water surface of that current month   |
| in | <i>real(dp), dimension(:) :: soil_moist_sat</i>     | Saturation soil moisture for each horizon [mm]   |
| in | <i>real(dp), dimension(:) :: frac_roots</i>         | Fraction of Roots in soil horizon  |
| in | <i>real(dp), dimension(:) :: soil_moist_FC</i>      | Soil moisture below which actual ET is reduced [mm]  |
| in | <i>real(dp), dimension(:) :: wilting_point</i>      | Permanent wilting point for each horizon [mm]  |
| in | <i>real(dp), dimension(:) :: soil_moist_exponen</i> | Exponential parameter to how non-linear is the soil water retention  |
| in | <i>real(dp) :: jarvis_thresh_c1</i>                 | Jarvis critical value for normalized soil water content  |
| in | <i>real(dp) :: aet_canopy</i>                       | Actual ET from canopy [mm TS-1]  |

## Parameters

|         |   |  |
|---------|---|--|
| in, out | <i>real(dp) :: prec_effec</i>                                       | Effective precipitation (rain + snow melt) [mm]                                      |
| in, out | <i>real(dp) :: runoff_sealed</i>                                    | Direct runoff from impervious areas  |
| in, out | <i>real(dp) :: storage_sealed</i>                                   | Retention storage of impervious areas  |
| in, out | <i>real(dp), dimension(size(soil_moist_sat, 1)) :: infiltration</i> | Recharge, infiltration intensity oreffective precipitation of each horizon [mm TS-1] |
| in, out | <i>real(dp), dimension(size(soil_moist_sat, 1)) :: soil_moist</i>   | Soil moisture of each horizon [mm]   |
| out     | <i>real(dp), dimension(size(soil_moist_sat, 1)) :: aet</i>          | actual ET [mm TS-1]  |
| out     | <i>real(dp) :: aet_sealed</i>                                       | actual ET from free-water surfaces,i.e impervious cover [mm TS-1]                    |

## Authors

Matthias Cuntz

## Date

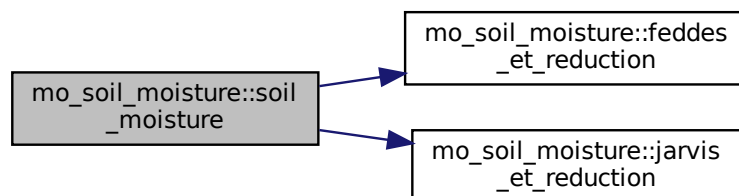
Dec 2012

Definition at line 95 of file mo\_soil\_moisture.f90.

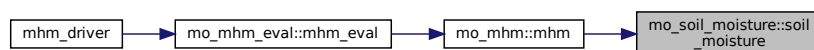
References mo\_common\_constants::eps\_dp, feddes\_et\_reduction(), and jarvis\_et\_reduction().

Referenced by mo\_mhm::mhm().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.72 mo\_spatial\_agg\_disagg\_forcing Module Reference

Spatial aggegation or disaggregation of meteorological input data.



## Data Types

- interface [spatial\\_aggregation](#)  
*Spatial aggregation of meteorological variables.*
- interface [spatial\\_disaggregation](#)  
*Spatial disaggregation of meteorological variables.*

## Functions/Subroutines

- subroutine [spatial\\_aggregation\\_3d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [spatial\\_aggregation\\_4d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [spatial\\_disaggregation\\_3d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [spatial\\_disaggregation\\_4d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)

### 17.72.1 Detailed Description

Spatial aggregation or disaggregation of meteorological input data.

This module contains two subroutines to upscale and downscale, respectively, the level-2 meteorological inputs to a required Level-1 hydrological spatial resolution.

#### Authors

Rohini Kumar

#### Date

Jan 2013

### 17.72.2 Function/Subroutine Documentation

#### 17.72.2.1 [spatial\\_aggregation\\_3d\(\)](#)

```
subroutine mo_spatial_agg_disagg_forcing::spatial_aggregation_3d (
    real(dp), dimension(:, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
    logical, dimension(:, :), intent(in) mask2,
    real(dp), dimension(:, :, :), intent(out), allocatable data1 ) [private]
```

Definition at line 96 of file mo\_spatial\_agg\_disagg\_forcing.f90.

#### 17.72.2.2 [spatial\\_aggregation\\_4d\(\)](#)

```
subroutine mo_spatial_agg_disagg_forcing::spatial_aggregation_4d (
    real(dp), dimension(:, :, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
```

```

logical, dimension(:, :), intent(in) mask2,
real(dp), dimension(:, :, :, :), intent(out), allocatable data1 )

```

Definition at line 204 of file mo\_spatial\_agg\_disagg\_forcing.f90.

### 17.72.2.3 spatial\_disaggregation\_3d()

```

subroutine mo_spatial_agg_disagg_forcing::spatial_disaggregation_3d (
    real(dp), dimension(:, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
    logical, dimension(:, :), intent(in) mask2,
    real(dp), dimension(:, :, :), intent(out), allocatable data1 )

```

Definition at line 315 of file mo\_spatial\_agg\_disagg\_forcing.f90.

### 17.72.2.4 spatial\_disaggregation\_4d()

```

subroutine mo_spatial_agg_disagg_forcing::spatial_disaggregation_4d (
    real(dp), dimension(:, :, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
    logical, dimension(:, :), intent(in) mask2,
    real(dp), dimension(:, :, :, :), intent(out), allocatable data1 )

```

Definition at line 381 of file mo\_spatial\_agg\_disagg\_forcing.f90.

## 17.73 mo\_startup Module Reference

Startup procedures for mHM.

### Functions/Subroutines

- subroutine, public [mhm\\_initialize](#)  
*Initialize main mHM variables.*
- subroutine [constants\\_init](#)  
*Initialize mHM constants.*
- subroutine [l2\\_variable\\_init](#) (iDomain, level0\_iDomain, level2\_iDomain)  
*Initialize Level-2 meteorological forcings data.*

### 17.73.1 Detailed Description

Startup procedures for mHM.

This module initializes all variables required to run mHM. This module needs to be run only one time at the beginning of a simulation if re-starting files do not exist.

#### Authors

Luis Samaniego, Rohini Kumar

## Date

Dec 2012

## 17.73.2 Function/Subroutine Documentation

## 17.73.2.1 constants\_init()

```
subroutine mo_startup::constants_init
```

Initialize mHM constants.

transformation of time units & initialize constants

## Authors

Luis Samaniego

## Date

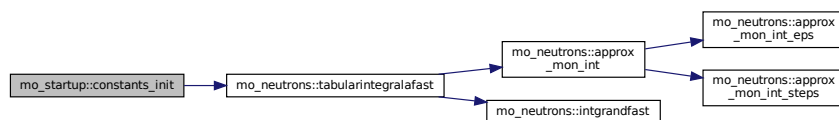
Dec 2012

Definition at line 144 of file mo\_startup.f90.

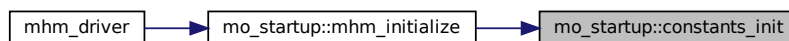
References mo\_common\_mhm\_mrm\_variables::c2tstu, mo\_mpr\_file::file\_hydrogeoclass, mo\_file::file\_namelist\_↔, mhm\_param, mo\_mpr\_global\_variables::geounitlist, mo\_global\_variables::neutron\_integral\_↔, mo\_common\_variables::processmatrix, mo\_neutrons::tabularintegralafast(), and mo\_common\_mhm\_mrm\_variables::timestep.

Referenced by mhm\_initialize().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.73.2.2 I2\_variable\_init()

```
subroutine mo_startup::I2_variable_init (
    integer(i4), intent(in) iDomain,
```

```

type(grid), intent(in) level0_iDomain,
type(grid), intent(inout) level2_iDomain )

```

Initialize Level-2 meteorological forcings data.

following tasks are performed 1) cell id & numbering 2) mask creation 3) append variable of interest to global ones

#### Parameters

|         |                                     |           |
|---------|-------------------------------------|-----------|
| in      | <i>integer(i4) :: iDomain</i>       | domain Id |
| in      | <i>type(Grid) :: level0_iDomain</i> |           |
| in, out | <i>type(Grid) :: level2_iDomain</i> |           |

#### Authors

Rohini Kumar

#### Date

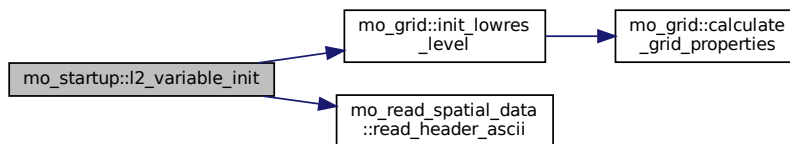
Feb 2013

Definition at line 211 of file mo\_startup.f90.

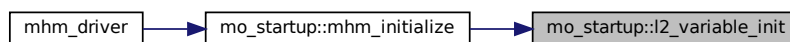
References mo\_global\_variables::dirprecipitation, mo\_mpr\_file::file\_meteo\_header, mo\_grid::init\_lowres\_level(), mo\_read\_spatial\_data::read\_header\_ascii(), and mo\_mpr\_file::umeteo\_header.

Referenced by mhm\_initialize().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.73.2.3 mhm\_initialize()

```

subroutine, public mo_startup::mhm_initialize

```

Initialize main mHM variables.

Initialize main mHM variables for a given domain. Calls the following procedures in this order:

- Constant initialization.
- Generate soil database.
- Checking inconsistencies input fields.
- Variable initialization at level-0.
- Variable initialization at level-1.
- Variable initialization at level-11.
- Space allocation of remaining variable/parameters. Global variables will be used at this stage.

**Authors**

Luis Samaniego, Rohini Kumar

**Date**

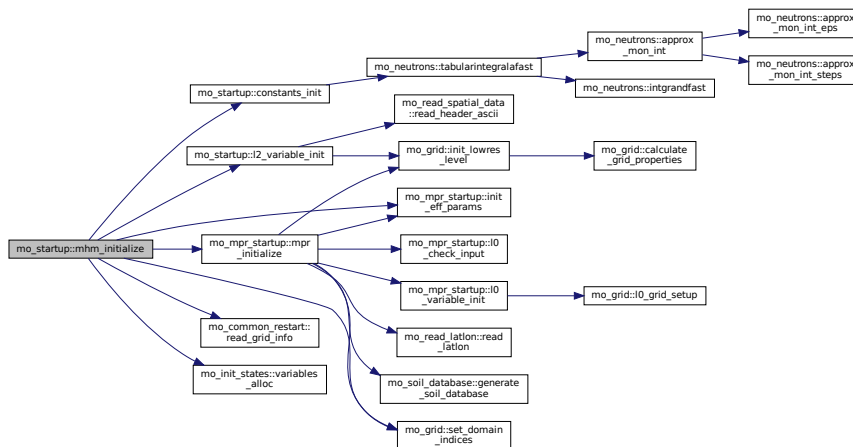
Dec 2012

Definition at line 70 of file mo\_startup.f90.

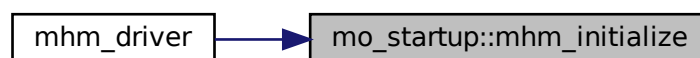
References constants\_init(), mo\_common\_variables::domainmeta, mo\_mpr\_startup::init\_eff\_params(), I2\_variable\_init(), mo\_common\_variables::level0, mo\_common\_variables::level1, mo\_global\_variables::level2, mo\_common\_mhm\_mrm\_variables::mhmfilerestartin, mo\_mpr\_startup::mpr\_initialize(), mo\_common\_restart::read\_grid\_info(), mo\_common\_mhm\_mrm\_variables::read\_restart, mo\_grid::set\_domain\_indices(), and mo\_init\_states::variables\_alloc().

Referenced by mhm\_driver().

Here is the call graph for this function:



Here is the caller graph for this function:



## 17.74 mo\_template Module Reference

Template for future module developments.

### Data Types

- interface [mean](#)  
*The average.*

### Functions/Subroutines

- elemental pure real(dp) function, public [circum](#) (radius)  
*Circumference of a circle.*
- real(dp) function [mean\\_dp](#) (dat, mask)
- real(sp) function [mean\\_sp](#) (dat, mask)

### Variables

- real(dp), parameter, public [pi\\_dp](#) = 3.141592653589793238462643383279502884197\_dp  
*Constant Pi in double precision.*
- real(sp), parameter, public [pi\\_sp](#) = 3.141592653589793238462643383279502884197\_sp  
*Constant Pi in single precision.*
- integer(i4), parameter [itest](#) = 1

### 17.74.1 Detailed Description

Template for future module developments.

This module serves as a template for future model developments. It shows the module structure, the coding style, and documentation. Please read the [Coding and documentation style](#) guide.

#### Authors

Matthias Cuntz, Christoph Schneider

#### Date

Dec 2012

### 17.74.2 Function/Subroutine Documentation

#### 17.74.2.1 [circum\(\)](#)

```
elemental pure real(dp) function, public mo_template::circum (
    real(dp), intent(in) radius )
```

Circumference of a circle.

Calculates the circumference of a circle

$$c = 2\pi r$$

## Parameters

|    |                           |        |
|----|---------------------------|--------|
| in | <i>real(dp) :: radius</i> | Radius |
|----|---------------------------|--------|

## Returns

*real(dp) :: circum* — circumference of circle.

## Authors

Matthias Cuntz

## Date

Dec 2012

Definition at line 122 of file `mo_template.f90`.

References `pi_dp`.

**17.74.2.2 mean\_dp()**

```
real(dp) function mo_template::mean_dp (
    real(dp), dimension(:), intent(in) dat,
    logical, dimension(:), intent(in), optional mask ) [private]
```

Definition at line 137 of file `mo_template.f90`.

Referenced by `mo_template::mean::mean_dp()`.

Here is the caller graph for this function:

**17.74.2.3 mean\_sp()**

```
real(sp) function mo_template::mean_sp (
    real(sp), dimension(:), intent(in) dat,
    logical, dimension(:), intent(in), optional mask ) [private]
```

Definition at line 167 of file `mo_template.f90`.

Referenced by `mo_template::mean::mean_sp()`.

Here is the caller graph for this function:



### 17.74.3 Variable Documentation

#### 17.74.3.1 itest

```
integer(i4), parameter mo_template::itest = 1 [private]
```

Definition at line 92 of file mo\_template.f90.

#### 17.74.3.2 pi\_dp

```
real(dp), parameter, public mo_template::pi_dp = 3.141592653589793238462643383279502884197_dp
```

Constant Pi in double precision.

Definition at line 87 of file mo\_template.f90.

Referenced by circum().

#### 17.74.3.3 pi\_sp

```
real(sp), parameter, public mo_template::pi_sp = 3.141592653589793238462643383279502884197_sp
```

Constant Pi in single precision.

Definition at line 89 of file mo\_template.f90.

## 17.75 mo\_temporal\_disagg\_forcing Module Reference

Temporal disaggregation of daily input values.

### Functions/Subroutines

- elemental pure subroutine, public [temporal\\_disagg\\_forcing](#) (isday, ntimesteps\_day, prec\_day, pet\_day, temp←\_day, fday\_prec, fday\_pet, fday\_temp, fnight\_prec, fnight\_pet, fnight\_temp, temp\_weights, pet\_weights, pre←\_weights, read\_meteo\_weights, prec, pet, temp)
 

*Temporally distribute daily mean forcings onto time step.*
- elemental subroutine, public [temporal\\_disagg\\_meteo\\_weights](#) (meteo\_val\_day, meteo\_val\_weights, meteo←\_val, weights\_correction)



*Temporally distribute daily mean forcings onto time step.*

- elemental subroutine, public [temporal\\_disagg\\_flux\\_daynight](#) (isday, ntimesteps\_day, meteo\_val\_day, fday\_↔meteo\_val, fnight\_meteo\_val, meteo\_val)

*Temporally distribute daily mean forcings onto time step.*

- elemental subroutine, public [temporal\\_disagg\\_state\\_daynight](#) (isday, ntimesteps\_day, meteo\_val\_day, fday\_↔\_meteo\_val, fnight\_meteo\_val, meteo\_val, add\_correction)

*Temporally distribute daily mean state forcings onto time step.*

## 17.75.1 Detailed Description

Temporal disaggregation of daily input values.

Calculate actual values for precipitation, PET and temperature from daily mean inputs ote There is not PET correction for aspect in this routine. Use pet \* fasp before or after the routine.

### Authors

Matthias Cuntz

### Date

Dec 2012

## 17.75.2 Function/Subroutine Documentation

### 17.75.2.1 temporal\_disagg\_flux\_daynight()

```
elemental subroutine, public mo_temporal_disagg_forcing::temporal_disagg_flux_daynight (
    logical, intent(in) isday,
    real(dp), intent(in) ntimesteps_day,
    real(dp), intent(in) meteo_val_day,
    real(dp), intent(in) fday_meteo_val,
    real(dp), intent(in) fnight_meteo_val,
    real(dp), intent(out) meteo_val )
```

Temporally distribute daily mean forcings onto time step.

Calculates actual meteo forcings from daily mean inputs. They are distributed with predefined factors/summands for day and night.

### Parameters

|     |                                     |                                   |
|-----|-------------------------------------|-----------------------------------|
| in  | <i>logical :: isday</i>             | is day or night                   |
| in  | <i>real(dp) :: ntimesteps_day</i>   | # of time steps per day           |
| in  | <i>real(dp) :: meteo_val_day</i>    | Daily mean meteo value            |
| in  | <i>real(dp) :: fday_meteo_val</i>   | Daytime fraction of meteo value   |
| in  | <i>real(dp) :: fnight_meteo_val</i> | Nighttime fraction of meteo value |
| out | <i>real(dp) :: meteo_val</i>        | Actual meteo value                |

### Authors

Sebastian Mueller

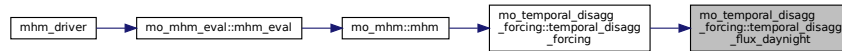
## Date

Jun 2020

Definition at line 222 of file mo\_temporal\_disagg\_forcing.f90.

Referenced by temporal\_disagg\_forcing().

Here is the caller graph for this function:



### 17.75.2.2 temporal\_disagg\_forcing()

```

elemental pure subroutine, public mo_temporal_disagg_forcing::temporal_disagg_forcing (
    logical, intent(in) isday,
    real(dp), intent(in) ntimesteps_day,
    real(dp), intent(in) prec_day,
    real(dp), intent(in) pet_day,
    real(dp), intent(in) temp_day,
    real(dp), intent(in) fday_prec,
    real(dp), intent(in) fday_pet,
    real(dp), intent(in) fday_temp,
    real(dp), intent(in) fnight_prec,
    real(dp), intent(in) fnight_pet,
    real(dp), intent(in) fnight_temp,
    real(dp), intent(in) temp_weights,
    real(dp), intent(in) pet_weights,
    real(dp), intent(in) pre_weights,
    logical, intent(in) read_meteo_weights,
    real(dp), intent(out) prec,
    real(dp), intent(out) pet,
    real(dp), intent(out) temp )
  
```

Temporally distribute daily mean forcings onto time step.

Calculates actual precipitation, PET and temperature from daily mean inputs. Precipitation and PET are distributed with predefined factors onto the day. Temperature gets a predefined amplitude added on day and subtracted at night. Alternatively, weights for each hour and month can be given and disaggregation is using these as factors for PET and temperature. Precipitation is distributed uniformly.

#### Parameters

|    |  |                                   |
|----|--|-----------------------------------|
| in | <i>logical</i> :: <i>isday</i>           | is day or night                   |
| in | <i>real(dp)</i> :: <i>ntimesteps_day</i> | # of time steps per day           |
| in | <i>real(dp)</i> :: <i>prec_day</i>       | Daily mean precipitation [mm/d]   |
| in | <i>real(dp)</i> :: <i>pet_day</i>        | Daily mean ET [mm/d]              |
| in | <i>real(dp)</i> :: <i>temp_day</i>       | Daily mean air temperature [K]    |
| in | <i>real(dp)</i> :: <i>fday_prec</i>      | Daytime fraction of precipitation |
| in | <i>real(dp)</i> :: <i>fday_pet</i>       | Daytime fraction of PET           |
| in | <i>real(dp)</i> :: <i>fday_temp</i>      | Daytime air temperature increase  |
| in | <i>real(dp)</i> :: <i>fnight_prec</i>    | Daytime fraction of precipitation |

## Parameters

|     |                                      |   |
|-----|--------------------------------------|---|
| in  | <i>real(dp) :: fnight_pet</i>        | Daytime fraction of PET                     |
| in  | <i>real(dp) :: fnight_temp</i>       | Daytime air temparture increase             |
| in  | <i>real(dp) :: temp_weights</i>      | weights for average temperature             |
| in  | <i>real(dp) :: pet_weights</i>       | weights for PET                             |
| in  | <i>real(dp) :: pre_weights</i>       | weights for precipitation                   |
| in  | <i>logical :: read_meteo_weights</i> | flag indicating that weights should be used |
| out | <i>real(dp) :: prec</i>              | Actual precipitation [mm/d]                 |
| out | <i>real(dp) :: pet</i>               | Reference ET [mm/d]                         |
| out | <i>real(dp) :: temp</i>              | Air temperature [K]                         |

## Authors

Matthias Cuntz

## Date

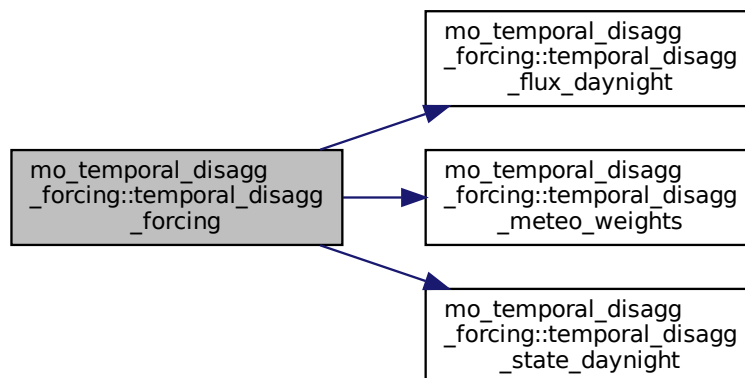
Dec 2012

Definition at line 77 of file mo\_temporal\_disagg\_forcing.f90.

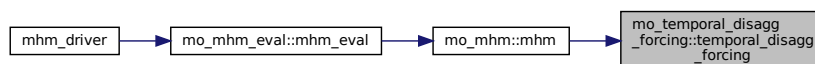
References temporal\_disagg\_flux\_daynight(), temporal\_disagg\_meteo\_weights(), and temporal\_disagg\_state\_↔daynight().

Referenced by mo\_mhm::mhm().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.75.2.3 temporal\_disagg\_meteo\_weights()

```

elemental subroutine, public mo_temporal_disagg_forcing::temporal_disagg_meteo_weights (
    real(dp), intent(in) meteo_val_day,
    real(dp), intent(in) meteo_val_weights,
    real(dp), intent(out) meteo_val,
    real(dp), intent(in), optional weights_correction )

```

Temporally distribute daily mean forcings onto time step.

Calculates actual meteo forcings from daily mean inputs. They are distributed with predefined weights onto the day.

#### Parameters

|     |  |  |
|-----|--|--|
| in  | <i>logical</i> :: <i>isday</i>                                 | is day or night                                      |
| in  | <i>real(dp)</i> :: <i>ntimesteps_day</i>                       | # of time steps per day                              |
| in  | <i>real(dp)</i> :: <i>meteo_val_day</i>                        | Daily mean meteo value                               |
| out | <i>real(dp)</i> :: <i>meteo_val</i>                            | Actual meteo value                                   |
| in  | <i>real(dp)</i> , <i>optional</i> :: <i>weights_correction</i> | Additive correction value before weights are applied |

#### Authors

Sebastian Mueller

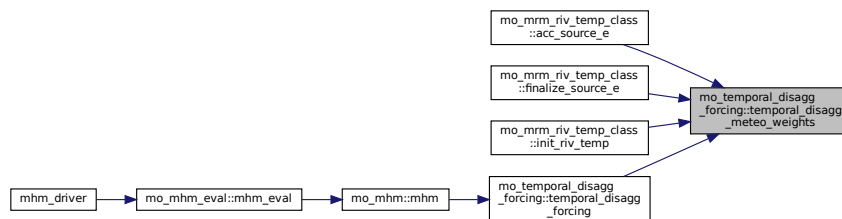
#### Date

Jun 2020

Definition at line 167 of file `mo_temporal_disagg_forcing.f90`.

Referenced by `mo_mrm_riv_temp_class::acc_source_e()`, `mo_mrm_riv_temp_class::finalize_source_e()`, `mo_mrm_riv_temp_class::init_riv_temp()`, and `temporal_disagg_forcing()`.

Here is the caller graph for this function:



### 17.75.2.4 temporal\_disagg\_state\_daynight()

```

elemental subroutine, public mo_temporal_disagg_forcing::temporal_disagg_state_daynight (
    logical, intent(in) isday,
    real(dp), intent(in) ntimesteps_day,
    real(dp), intent(in) meteo_val_day,

```

```

real(dp), intent(in) fday_meteo_val,
real(dp), intent(in) fnight_meteo_val,
real(dp), intent(out) meteo_val,
logical, intent(in), optional add_correction )

```

Temporally distribute daily mean state forcings onto time step.

Calculates meteo forcings from daily mean inputs of a state variable. They are distributed with predefined factors/summands for day and night.

#### Parameters

|     |   |   |
|-----|---|---|
| in  | <i>logical</i> :: <i>isday</i>                    | is day or night   |
| in  | <i>real(dp)</i> :: <i>ntimesteps_day</i>          | # of time steps per day                                   |
| in  | <i>real(dp)</i> :: <i>meteo_val_day</i>           | Daily mean meteo value                                    |
| in  | <i>real(dp)</i> :: <i>fday_meteo_val</i>          | Daytime fraction/addition of meteo value                  |
| in  | <i>real(dp)</i> :: <i>fnight_meteo_val</i>        | Nighttime fraction/addition of meteo value                |
| out | <i>real(dp)</i> :: <i>meteo_val</i>               | Actual meteo value  |
| in  | <i>logical, optional</i> :: <i>add_correction</i> | State if correction should be added instead of multiplied |

#### Authors

Sebastian Mueller

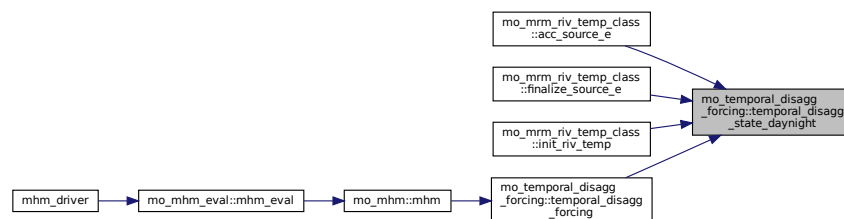
#### Date

Jun 2020

Definition at line 288 of file mo\_temporal\_disagg\_forcing.f90.

Referenced by mo\_mrm\_riv\_temp\_class::acc\_source\_e(), mo\_mrm\_riv\_temp\_class::finalize\_source\_e(), mo\_mrm\_riv\_temp\_class::init\_riv\_temp(), and temporal\_disagg\_forcing().

Here is the caller graph for this function:



## 17.76 mo\_upscaling\_operators Module Reference

Module containing upscaling operators.

### Functions/Subroutines

- integer(i4) function, dimension(size(l1\_upper\_rowld\_cell, 1)), public [majority\\_statistics](#) (nClass, L1\_upper\_rowld\_cell, L1\_lower\_rowld\_cell, L1\_left\_colonld\_cell, L1\_right\_colonld\_cell, L0\_fineScale\_2D\_data)  
*majority\_statistics*

- real(dp) function, dimension(size(l0upbound\_inLx, 1)), public [l0\\_fractionalcover\\_in\\_Lx](#) (dataIn0, classId, mask0, L0upBound\_inLx, L0downBound\_inLx, L0leftBound\_inLx, L0rightBound\_inLx, nTCells0\_inLx)  
*fractional coverage of a given class of L0 fields in Lx field (Lx = L1 or L11)*
- real(dp) function, dimension(size(nl0\_cells\_in\_l1\_cell, 1)), public [upscale\\_arithmetic\\_mean](#) (nL0\_cells\_in\_L1\_cell, L1\_upper\_rowId\_cell, L1\_lower\_rowId\_cell, L1\_left\_colonId\_cell, L1\_right\_colonId\_cell, L0\_cellId, mask0, nodata\_value, L0\_fineScale\_data)  
*arithmetic mean*
- real(dp) function, dimension(size(nl0\_cells\_in\_l1\_cell, 1)), public [upscale\\_harmonic\\_mean](#) (nL0\_cells\_in\_L1\_cell, L1\_upper\_rowId\_cell, L1\_lower\_rowId\_cell, L1\_left\_colonId\_cell, L1\_right\_colonId\_cell, L0\_cellId, mask0, nodata\_value, L0\_fineScale\_data)  
*harmonic mean*
- real(dp) function, dimension(size(l1\_upper\_rowId\_cell, 1)), public [upscale\\_geometric\\_mean](#) (L1\_upper\_rowId\_cell, L1\_lower\_rowId\_cell, L1\_left\_colonId\_cell, L1\_right\_colonId\_cell, mask0, nodata\_value, L0\_fineScale\_data)  
*geometric mean*
- real(dp) function, dimension(size(nl0\_cells\_in\_l1\_cell, 1)) [upscale\\_p\\_norm](#) (nL0\_cells\_in\_L1\_cell, L1\_upper\_rowId\_cell, L1\_lower\_rowId\_cell, L1\_left\_colonId\_cell, L1\_right\_colonId\_cell, L0\_cellId, mask0, nodata\_value, p\_norm, L0\_fineScale\_data)  
*arithmetic mean*

### 17.76.1 Detailed Description

Module containing upscaling operators.

This module provides the routines for upscaling operators.

#### Authors

Giovanni Dalmasso, Rohini Kumar

#### Date

Dec 2012

### 17.76.2 Function/Subroutine Documentation

#### 17.76.2.1 l0\_fractionalcover\_in\_Lx()

```
real(dp) function, dimension(size(l0upbound_inLx, 1)), public mo_upscaling_operators::l0_fractionalcover_in_Lx (
    integer(i4), dimension(:), intent(in) dataIn0,
    integer(i4), intent(in) classId,
    logical, dimension(:, :), intent(in) mask0,
    integer(i4), dimension(:), intent(in) L0upBound_inLx,
    integer(i4), dimension(:), intent(in) L0downBound_inLx,
    integer(i4), dimension(:), intent(in) L0leftBound_inLx,
    integer(i4), dimension(:), intent(in) L0rightBound_inLx,
    integer(i4), dimension(:), intent(in) nTCells0_inLx )
```

fractional coverage of a given class of L0 fields in Lx field (Lx = L1 or L11)

Fractional coverage of a given class of L0 fields in Lx field (Lx = L1 or L11). For example, this routine can be used for calculating the karstic fraction.

## Parameters

|    |   |   |
|----|---|---|
| in | <i>integer(i4), dimension(:) :: dataIn0</i>           | input fields at finer scale                       |
| in | <i>integer(i4) :: classId</i>                         | class id for which fraction has to be estimated   |
| in | <i>logical, dimension(:, :) :: mask0</i>              | finer scale L0 mask                               |
| in | <i>integer(i4), dimension(:) :: L0upBound_inLx</i>    | row start at finer L0 scale                       |
| in | <i>integer(i4), dimension(:) :: L0downBound_inLx</i>  | row end at finer L0 scale                         |
| in | <i>integer(i4), dimension(:) :: L0leftBound_inLx</i>  | col start at finer L0 scale                       |
| in | <i>integer(i4), dimension(:) :: L0rightBound_inLx</i> | col end at finer L0 scale                         |
| in | <i>integer(i4), dimension(:) :: nTCells0_inLx</i>     | total number of valid L0 cells in a given Lx cell |

## Returns

real(dp) :: L0\_fractionalCover\_in\_Lx(:) — packed 1D fraction coverage (Lx) of given class id

## Authors

Rohini Kumar

## Date

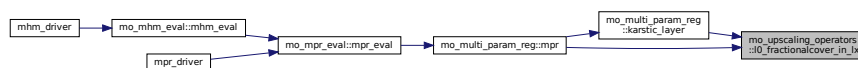
Feb 2013

Definition at line 153 of file mo\_upscaling\_operators.f90.

References mo\_common\_constants::nodata\_i4.

Referenced by mo\_multi\_param\_reg::karstic\_layer(), and mo\_multi\_param\_reg::mpr().

Here is the caller graph for this function:



## 17.76.2.2 majority\_statistics()

```

integer(i4) function, dimension(size(l1_upper_rowId_cell, 1)), public mo_upscaling_operators←
::majority_statistics (
    integer(i4), intent(in) nClass,
    integer(i4), dimension(:), intent(in) L1_upper_rowId_cell,
    integer(i4), dimension(:), intent(in) L1_lower_rowId_cell,
    integer(i4), dimension(:), intent(in) L1_left_colonId_cell,
    integer(i4), dimension(:), intent(in) L1_right_colonId_cell,
    integer(i4), dimension(:, :), intent(in) L0_fineScale_2D_data )

```

majority statistics

upscale grid L0\_fineScale\_2D\_data based on a majority statistics

## Parameters

|    |                              |                   |
|----|------------------------------|-------------------|
| in | <i>integer(i4) :: nClass</i> | number of classes |
|----|------------------------------|-------------------|

## Parameters

|    |   |  |
|----|---|--|
| in | <i>integer(i4), dimension(:) :: L1_upper_rowId_cell</i>     | upper row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: L1_lower_rowId_cell</i>     | lower row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: L1_left_colonId_cell</i>    | left colon boundary (level-0) of a level-1 cell  |
| in | <i>integer(i4), dimension(:) :: L1_right_colonId_cell</i>   | right colon boundary (level-0) of a level-1 cell |
| in | <i>integer(i4), dimension(:, :) :: L0_fineScale_2D_data</i> | high resolution data                             |

## Returns

*integer(i4) :: majority\_statistics(:)* — Upscaled variable based on majority.

## Authors

Giovanni Dalmaso, Rohini Kumar

## Date

Dec 2012

Definition at line 66 of file `mo_upscaling_operators.f90`.

### 17.76.2.3 upscale\_arithmetic\_mean()

```
real(dp) function, dimension(size(nL0_cells_in_L1_cell, 1)), public mo_upscaling_operators←
::upscale_arithmetic_mean (
    integer(i4), dimension(:), intent(in) nL0_cells_in_L1_cell,
    integer(i4), dimension(:), intent(in) L1_upper_rowId_cell,
    integer(i4), dimension(:), intent(in) L1_lower_rowId_cell,
    integer(i4), dimension(:), intent(in) L1_left_colonId_cell,
    integer(i4), dimension(:), intent(in) L1_right_colonId_cell,
    integer(i4), dimension(:), intent(in) L0_cellId,
    logical, dimension(:, :), intent(in) mask0,
    real(dp), intent(in) nodata_value,
    real(dp), dimension(:), intent(in) L0_fineScale_data )
```

arithmetic mean

upsampling of level-0 grid data to level-1 using arithmetic mean

## Parameters

|    |   |  |
|----|---|--|
| in | <i>integer(i4), dimension(:) :: nL0_cells_in_L1_cell</i>  | number of level-0 cells within a level-1 cell    |
| in | <i>integer(i4), dimension(:) :: L1_upper_rowId_cell</i>   | upper row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: L1_lower_rowId_cell</i>   | lower row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: L1_left_colonId_cell</i>  | left colon boundary (level-0) of a level-1 cell  |
| in | <i>integer(i4), dimension(:) :: L1_right_colonId_cell</i> | right colon boundary (level-0) of a level-1 cell |
| in | <i>integer(i4), dimension(:) :: L0_cellId</i>             | cell ID at level-0                               |
| in | <i>logical, dimension(:, :) :: mask0</i>                  | mask at level 0                                  |
| in | <i>real(dp) :: nodata_value</i>                           | no data value                                    |
| in | <i>real(dp), dimension(:) :: L0_fineScale_data</i>        | high resolution data                             |



**Returns**

real(dp) :: upscale\_arithmetic\_mean(:) — Upscaled variable from L0 to L1 using arithmetic mean

**Authors**

Giovanni Dalmaso, Rohini Kumar

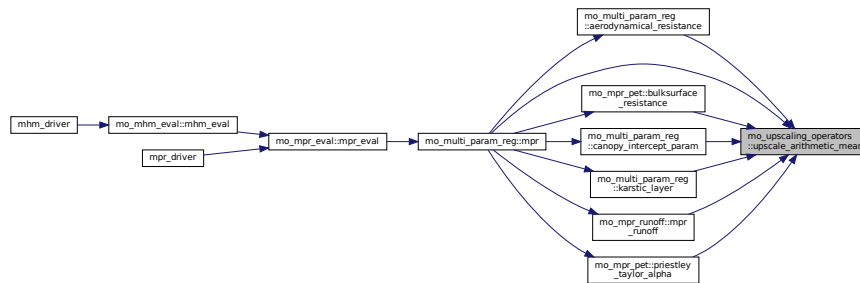
**Date**

Dec 2012

Definition at line 267 of file mo\_upscaling\_operators.f90.

Referenced by mo\_multi\_param\_reg::aerodynamical\_resistance(), mo\_mpr\_pet::bulksurface\_resistance(), mo\_multi\_param\_reg::canopy\_intercept\_param(), mo\_multi\_param\_reg::karstic\_layer(), mo\_multi\_param\_reg::mpr(), mo\_mpr\_runoff::mpr\_runoff(), and mo\_mpr\_pet::priestley\_taylor\_alpha().

Here is the caller graph for this function:

**17.76.2.4 upscale\_geometric\_mean()**

```
real(dp) function, dimension(size(l1_upper_rowId_cell, 1)), public mo_upscaling_operators::upscale_geometric_mean (
    integer(i4), dimension(:), intent(in) L1_upper_rowId_cell,
    integer(i4), dimension(:), intent(in) L1_lower_rowId_cell,
    integer(i4), dimension(:), intent(in) L1_left_colonId_cell,
    integer(i4), dimension(:), intent(in) L1_right_colonId_cell,
    logical, dimension(:, :), intent(in) mask0,
    real(dp), intent(in) nodata_value,
    real(dp), dimension(:), intent(in) L0_fineScale_data )
```

geometric mean

upscaling of level-0 grid data to level-1 using geometric mean

**Parameters**

|    |   |  |
|----|---|--|
| in | <i>integer(i4), dimension(:) :: L1_upper_rowId_cell</i>   | upper row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: L1_lower_rowId_cell</i>   | lower row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: L1_left_colonId_cell</i>  | left colon boundary (level-0) of a level-1 cell  |
| in | <i>integer(i4), dimension(:) :: L1_right_colonId_cell</i> | right colon boundary (level-0) of a level-1 cell |
| in | <i>logical, dimension(:, :) :: mask0</i>                  | mask at level 0                                  |
| in | <i>real(dp) :: nodata_value</i>                           | no data value                                    |
| in | <i>real(dp), dimension(:) :: L0_fineScale_data</i>        | high resolution data                             |

**Returns**

`real(dp) :: upscale_geometric_mean(:)` — Upscaled variable from L0 to L1 using geometric mean

**Authors**

Giovanni Dalmaso, Rohini Kumar

**Date**

Dec 2012

Definition at line 470 of file `mo_upscaling_operators.f90`.

**17.76.2.5 upscale\_harmonic\_mean()**

```
real(dp) function, dimension(size(nl0_cells_in_l1_cell, 1)), public mo_upscaling_operators↵
::upscale_harmonic_mean (
    integer(i4), dimension(:), intent(in) nl0_cells_in_l1_cell,
    integer(i4), dimension(:), intent(in) l1_upper_rowId_cell,
    integer(i4), dimension(:), intent(in) l1_lower_rowId_cell,
    integer(i4), dimension(:), intent(in) l1_left_colonId_cell,
    integer(i4), dimension(:), intent(in) l1_right_colonId_cell,
    integer(i4), dimension(:), intent(in) l0_cellId,
    logical, dimension(:, :), intent(in) mask0,
    real(dp), intent(in) nodata_value,
    real(dp), dimension(:), intent(in) l0_fineScale_data )
```

harmonic mean

upscaling of level-0 grid data to level-1 using harmonic mean

**Parameters**

|    |   |  |
|----|---|--|
| in | <code>integer(i4), dimension(:) :: nl0_cells_in_l1_cell</code>  | number of level-0 cells within a level-1 cell    |
| in | <code>integer(i4), dimension(:) :: l1_upper_rowId_cell</code>   | upper row boundary (level-0) of a level-1 cell   |
| in | <code>integer(i4), dimension(:) :: l1_lower_rowId_cell</code>   | lower row boundary (level-0) of a level-1 cell   |
| in | <code>integer(i4), dimension(:) :: l1_left_colonId_cell</code>  | left colon boundary (level-0) of a level-1 cell  |
| in | <code>integer(i4), dimension(:) :: l1_right_colonId_cell</code> | right colon boundary (level-0) of a level-1 cell |
| in | <code>integer(i4), dimension(:) :: l0_cellId</code>             | cell ID at level-0                               |
| in | <code>logical, dimension(:, :) :: mask0</code>                  | mask at Level 0                                  |
| in | <code>real(dp) :: nodata_value</code>                           | no data value                                    |
| in | <code>real(dp), dimension(:) :: l0_fineScale_data</code>        | high resolution data                             |

**Returns**

`real(dp) :: upscale_harmonic_mean(:)` — Upscaled variable from L0 to L1 using harmonic mean

**Authors**

Giovanni Dalmaso, Rohini Kumar

## Date

Dec 2012

Definition at line 370 of file mo\_upscaling\_operators.f90.

Referenced by mo\_mpr\_smhorizons::mpr\_smhorizons(), and mo\_mpr\_pet::pet\_correctbylai().

Here is the caller graph for this function:



## 17.76.2.6 upscale\_p\_norm()

```

real(dp) function, dimension(size(nl0_cells_in_l1_cell, 1)) mo_upscaling_operators::upscale_p←
_norm (
    integer(i4), dimension(:), intent(in) nl0_cells_in_l1_cell,
    integer(i4), dimension(:), intent(in) l1_upper_rowId_cell,
    integer(i4), dimension(:), intent(in) l1_lower_rowId_cell,
    integer(i4), dimension(:), intent(in) l1_left_colonId_cell,
    integer(i4), dimension(:), intent(in) l1_right_colonId_cell,
    integer(i4), dimension(:), intent(in) l0_cellId,
    logical, dimension(:, :), intent(in) mask0,
    real(dp), intent(in) nodata_value,
    real(dp), intent(in) p_norm,
    real(dp), dimension(:), intent(in) l0_fineScale_data )

```

arithmetic mean

upscaling of level-0 grid data to level-1 using arithmetic mean

## Parameters

|    |   |  |
|----|---|--|
| in | <i>integer(i4), dimension(:) :: nl0_cells_in_l1_cell</i>  | number of level-0 cells within a level-1 cell    |
| in | <i>integer(i4), dimension(:) :: l1_upper_rowId_cell</i>   | upper row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: l1_lower_rowId_cell</i>   | lower row boundary (level-0) of a level-1 cell   |
| in | <i>integer(i4), dimension(:) :: l1_left_colonId_cell</i>  | left colon boundary (level-0) of a level-1 cell  |
| in | <i>integer(i4), dimension(:) :: l1_right_colonId_cell</i> | right colon boundary (level-0) of a level-1 cell |
| in | <i>integer(i4), dimension(:) :: l0_cellId</i>             | cell ID at level-0                               |
| in | <i>logical, dimension(:, : ) :: mask0</i>                 | mask at level 0                                  |
| in | <i>real(dp) :: nodata_value</i>                           | no data value                                    |
| in | <i>real(dp) :: p_norm</i>                                 | p_norm value                                     |
| in | <i>real(dp), dimension(:) :: l0_fineScale_data</i>        | high resolution data                             |

## Returns

real(dp) :: upscale\_arithmetic\_mean() — Upscaled variable from L0 to L1 using arithmetic mean

**Authors**

Giovanni Dalmasso, Rohini Kumar

**Date**

Dec 2012

Definition at line 577 of file mo\_upscaling\_operators.f90.

## 17.77 mo\_write\_ascii Module Reference

Module to write ascii file output.

### Functions/Subroutines

- subroutine, public [write\\_configfile](#)  
*This module writes the results of the configuration into an ASCII-file.*
- subroutine, public [write\\_optifile](#) (best\_OF, best\_paramSet, param\_names)  
*Write briefly final optimization results.*
- subroutine, public [write\\_optinamelist](#) (processMatrix, parameters, maskpara, parameters\_name)  
*Write final, optimized parameter set in a namelist format.*

### 17.77.1 Detailed Description

Module to write ascii file output.

Module to write ascii file output. Writing model output to ASCII should be the exception. Therefore, output is written usually as NetCDF and only: (1) The configuration file of mHM, (2) the final parameter set after optimization, and (3) the simulated vs. observed daily discharge is written in ASCII file format to allow for a quick assurance of proper model runs.

**Authors**

Christoph Schneider, Juliane Mai, Luis Samaniego

**Date**

May 2013

### 17.77.2 Function/Subroutine Documentation

#### 17.77.2.1 write\_configfile()

```
subroutine, public mo_write_ascii::write_configfile
```

This module writes the results of the configuration into an ASCII-file.

TODO: add description

**Authors**

Christoph Schneider

**Date**

May 2013

TODO: add description

TODO: add description

**Authors**

Robert Schwappe

**Date**

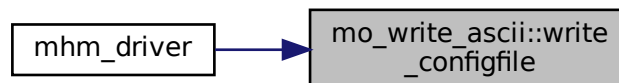
Jun 2018

Definition at line 85 of file mo\_write\_ascii.f90.

References mo\_common\_variables::dirconfigout, mo\_mrm\_global\_variables::dirgauges, mo\_common\_variables::dirlcover, mo\_common\_variables::dirmorpho, mo\_common\_variables::dirout, mo\_global\_variables::dirprecipitation, mo\_global\_variables::dirreferenceet, mo\_global\_variables::dirtemperature, mo\_common\_variables::domainmeta, mo\_common\_mhm\_mrm\_variables::evalper, mo\_common\_file::file\_config, mo\_mrm\_global\_variables::gauge, mo\_common\_variables::global\_parameters, mo\_common\_variables::global\_parameters\_name, mo\_common\_variables::iflag\_cordinate\_sys, mo\_mrm\_global\_variables::inflowgauge, mo\_mrm\_global\_variables::l11\_fromn, mo\_mrm\_global\_variables::l11\_label, mo\_mrm\_global\_variables::l11\_length, mo\_mrm\_global\_variables::l11\_netperm, mo\_mrm\_global\_variables::l11\_noutlets, mo\_mrm\_global\_variables::l11\_rorder, mo\_mrm\_global\_variables::l11\_slope, mo\_mrm\_global\_variables::l11\_ton, mo\_mrm\_global\_variables::l1\_l11\_id, mo\_common\_variables::lc\_year\_end, mo\_common\_variables::lc\_year\_start, mo\_common\_variables::lcfilename, mo\_common\_mhm\_mrm\_variables::lcyetid, mo\_common\_variables::level0, mo\_common\_variables::level1, mo\_mrm\_global\_variables::level11, mo\_common\_variables::mhmfilerestartout, mo\_mrm\_global\_variables::ngaugeslocal, mo\_mrm\_global\_variables::ngaigestotal, mo\_mrm\_global\_variables::ninflowgaigestotal, mo\_common\_variables::nlcoverscene, mo\_common\_constants::nodata\_dp, mo\_common\_variables::processmatrix, mo\_common\_mhm\_mrm\_variables::read\_restart, mo\_common\_variables::resolutionhydrology, mo\_common\_mhm\_mrm\_variables::resolutionrouting, mo\_common\_mhm\_mrm\_variables::simper, mo\_common\_mhm\_mrm\_variables::timestep, mo\_common\_file::uconfig, mo\_file::version, mo\_common\_mhm\_mrm\_variables::warmper, and mo\_common\_variables::write\_restart.

Referenced by mhm\_driver().

Here is the caller graph for this function:

**17.77.2.2 write\_optifile()**

```

subroutine, public mo_write_ascii::write_optifile (
    real(dp), intent(in) best_OF,
    real(dp), dimension(:), intent(in) best_paramSet,
    character(len = *), dimension(:), intent(in) param_names )
  
```

Write briefly final optimization results.

Write overall best objective function and the best optimized parameter set to a file\_opti.

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>real(dp) :: best_OF</i>                             | best objective function value as returned by the optimization routine    |
| in | <i>real(dp), dimension(:) :: best_paramSet</i>         | best associated global parameter set Called only when optimize is .TRUE. |
| in | <i>character(len = *), dimension(:) :: param_names</i> |  |

#### Authors

David Schaefer

#### Date

July 2013

Definition at line 425 of file mo\_write\_ascii.f90.

References mo\_common\_variables::dirconfigout, mo\_common\_mhm\_mrm\_file::file\_opti, and mo\_common\_mhm\_mrm\_file::uopti.

### 17.77.2.3 write\_optinamelist()

```
subroutine, public mo_write_ascii::write_optinamelist (
    integer(i4), dimension(nprocesses, 3), intent(in) processMatrix,
    real(dp), dimension(:, :) :: parameters,
    logical, dimension(size(parameters, 1)) :: maskpara,
    character(len = *), dimension(size(parameters, 1)) :: parameters_name )
```

Write final, optimized parameter set in a namelist format.

Write final, optimized parameter set in a namelist format. Only parameters of processes which were switched on are written to the namelist. All others are discarded.

#### Parameters

|    |  |   |
|----|--|---|
| in | <i>integer(i4), dimension(nProcesses, 3) :: processMatrix</i>                | information about which process case was used |
| in | <i>real(dp), dimension(:, :) :: parameters</i>                               | (min, max, opti)                              |
| in | <i>logical, dimension(size(parameters, 1)) :: maskpara</i>                   | .true. if parameter was calibrated            |
| in | <i>character(len = *), dimension(size(parameters, 1)) :: parameters_name</i> | clear names of parameters                     |

#### Authors

Juliane Mai

**Date**

Dec 2013

Definition at line 512 of file mo\_write\_ascii.f90.

References mo\_common\_variables::dirconfigout, mo\_common\_mhm\_mrm\_file::file\_opti\_nml, mo\_common\_variables::nprocesses, and mo\_common\_mhm\_mrm\_file::uopti\_nml.

## 17.78 mo\_write\_fluxes\_states Module Reference

Creates NetCDF output for different fluxes and state variables of mHM.

### Data Types

- interface [outputvariable](#)
- interface [outputdataset](#)

### Functions/Subroutines

- type([outputvariable](#)) function [newoutputvariable](#) (nc, name, dtype, dims, ncells, mask, avg)  
*Initialize OutputVariable.*
- subroutine [updatevariable](#) (self, data)  
*Update OutputVariable.*
- subroutine [writevariabletimestep](#) (self, timestep)  
*Write timestep to file.*
- type([outputdataset](#)) function, public [newoutputdataset](#) (iDomain, mask1, nCells)  
*Initialize OutputDataset.*
- subroutine [updatedataset](#) (self, sidx, eidx, L1\_fSealed, L1\_fNotSealed, L1\_inter, L1\_snowPack, L1\_soilMoist, L1\_soilMoistSat, L1\_sealSTW, L1\_unsatSTW, L1\_satSTW, L1\_neutrons, L1\_pet, L1\_aETSoil, L1\_aETCanopy, L1\_aETSealed, L1\_total\_runoff, L1\_runoffSeal, L1\_fastRunoff, L1\_slowRunoff, L1\_baseflow, L1\_percol, L1\_infilSoil, L1\_preEffect)  
*Update all variables.*
- subroutine [writetimestep](#) (self, timestep)  
*Write all accumulated data.*
- subroutine [close](#) (self)  
*Close the file.*
- type(ncdataset) function [createoutputfile](#) (iDomain)  
*Create and initialize output file for X & Y coordinate system.*
- subroutine [writevariableattributes](#) (var, long\_name, unit)  
*Write output variable attributes.*
- character(16) function [fluxesunit](#) (iDomain)  
*Generate a unit string.*

### 17.78.1 Detailed Description

Creates NetCDF output for different fluxes and state variables of mHM.

NetCDF is first initialized and later on variables are put to the NetCDF.

**Authors**

Matthias Zink

**Date**

Apr 2013

**17.78.2 Function/Subroutine Documentation****17.78.2.1 close()**

```
subroutine mo_write_fluxes_states::close (
    class(outputdataset) self ) [private]
```

Close the file.

Close the file associated with variable of type(OutputDataset)

**Authors**

Rohini Kumar &amp; Stephan Thober

**Date**

August 2013

Definition at line 851 of file mo\_write\_fluxes\_states.f90.

References mo\_common\_variables::dirout.

**17.78.2.2 createoutputfile()**

```
type(ncdataset) function mo_write_fluxes_states::createoutputfile (
    integer(i4), intent(in) iDomain )
```

Create and initialize output file for X & Y coordinate system.

Create output file, write all non-dynamic variables and global attributes for the given domain for X & Y coordinate system

**Returns**

type(NcDataset)

**Parameters**

|    |                               |              |
|----|-------------------------------|--------------|
| in | <i>integer(i4) :: iDomain</i> | -> domain id |
|----|-------------------------------|--------------|

**Authors**

David Schaefer

**Date**

June 2015

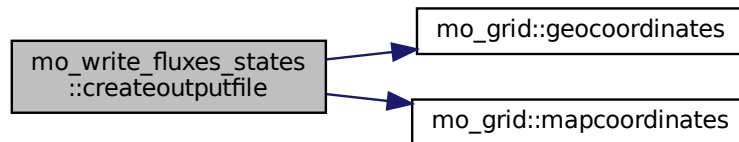
Definition at line 893 of file mo\_write\_fluxes\_states.f90.



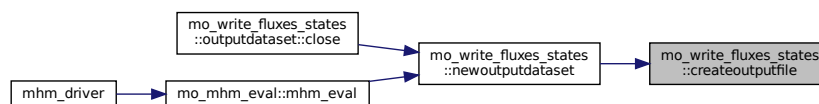
References mo\_common\_variables::dirout, mo\_common\_mhm\_mrm\_variables::evalper, mo\_grid::geocoordinates(), mo\_common\_variables::iflag\_coordinate\_sys, mo\_common\_variables::level1, mo\_grid::mapcoordinates(), mo\_↔common\_constants::nodata\_dp, mo\_global\_variables::output\_double\_precision, and mo\_file::version.

Referenced by newoutputdataset().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.78.2.3 fluxesunit()

```
character(16) function mo_write_fluxes_states::fluxesunit (
    integer(i4), intent(in) iDomain ) [private]
```

Generate a unit string.

Generate the unit string for the output variable netcdf attribute based on modeling timestep

#### Returns

character(16)

#### Parameters

|    |                               |  |
|----|-------------------------------|--|
| in | <i>integer(i4) :: iDomain</i> |  |
|----|-------------------------------|--|

#### Authors

David Schaefer

**Date**

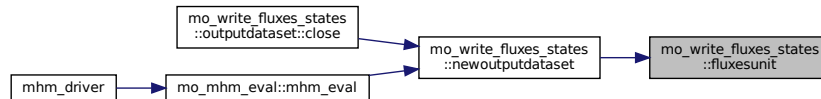
June 2015

Definition at line 1089 of file mo\_write\_fluxes\_states.f90.

References mo\_common\_mhm\_mrm\_variables::ntstepday, mo\_common\_mhm\_mrm\_variables::simper, mo\_↔common\_mhm\_mrm\_variables::timestep, and mo\_global\_variables::timestep\_model\_outputs.

Referenced by newoutputdataset().

Here is the caller graph for this function:

**17.78.2.4 newoutputdataset()**

```

type(outputdataset) function, public mo_write_fluxes_states::newoutputdataset (
    integer(i4), intent(in) iDomain,
    logical, dimension(:, :), intent(in), pointer mask1,
    integer(i4), intent(in) nCells )
  
```

Initialize OutputDataset.

Create and initialize the output file. If new a new output variable needs to be written, this is the first of two procedures to change (second: updateDataset)

**Returns**

type(OutputDataset)

**Parameters**

|    |  |                                    |
|----|--|------------------------------------|
| in | <i>integer(i4) :: iDomain</i>            | -> domain id                       |
| in | <i>logical, dimension(:, :) :: mask1</i> | -> L1 mask to reconstruct the data |
| in | <i>integer(i4) :: nCells</i>             |                                    |

**Authors**

Matthias Zink

**Date**

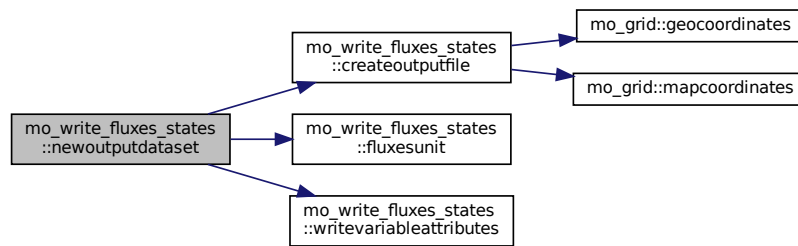
Apr 2013

Definition at line 242 of file mo\_write\_fluxes\_states.f90.

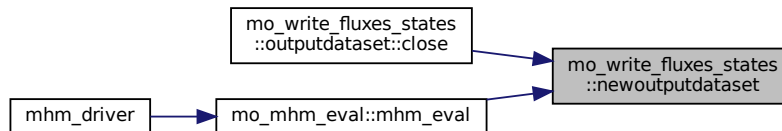
References createoutputfile(), fluxesunit(), mo\_common\_variables::iflag\_cordinate\_sys, mo\_mpr\_global\_↔variables::nsoilhorizons\_mhm, mo\_global\_variables::outputfixstate, and writevariableattributes().

Referenced by mo\_write\_fluxes\_states::outputdataset::close(), and mo\_mhm\_eval::mhm\_eval().

Here is the call graph for this function:



Here is the caller graph for this function:



### 17.78.2.5 newoutputvariable()

```

type(outputvariable) function mo_write_fluxes_states::newoutputvariable (
    type(ncdataset), intent(in) nc,
    character(*), intent(in) name,
    character(*), intent(in) dtype,
    character(16), dimension(3), intent(in) dims,
    integer(i4), intent(in) ncells,
    logical, dimension(:, :), intent(in), target mask,
    logical, intent(in), optional avg ) [private]
  
```

Initialize OutputVariable.

TODO: add description

#### Returns

type(OutputVariable)

#### Parameters

|    |  |  |
|----|--|--|
| in | <i>type(NcDataset) :: nc</i>               | -> NcDataset which contains the variable |
| in | <i>character(*) :: name</i>                |  |
| in | <i>character(*) :: dtype</i>               |  |
| in | <i>character(16), dimension(3) :: dims</i> |  |
| in | <i>integer(i4) :: ncells</i>               | -> number of cells in domain             |

## Parameters

|    |   |                                    |
|----|---|------------------------------------|
| in | <i>logical, dimension(:, :) :: mask</i> |                                    |
| in | <i>logical, optional :: avg</i>         | -> average the data before writing |

## Authors

David Schaefer

## Date

June 2015

Definition at line 103 of file mo\_write\_fluxes\_states.f90.

References mo\_global\_variables::output\_deflate\_level.

### 17.78.2.6 updatedataset()

```

subroutine mo_write_fluxes_states::updatedataset (
    class(outputdataset), intent(inout), target self,
    integer(i4), intent(in) sidx,
    integer(i4), intent(in) eidx,
    real(dp), dimension(:), intent(in) Ll_fSealed,
    real(dp), dimension(:), intent(in) Ll_fNotSealed,
    real(dp), dimension(:), intent(in) Ll_inter,
    real(dp), dimension(:), intent(in) Ll_snowPack,
    real(dp), dimension(:, :), intent(in) Ll_soilMoist,
    real(dp), dimension(:, :), intent(in) Ll_soilMoistSat,
    real(dp), dimension(:), intent(in) Ll_sealSTW,
    real(dp), dimension(:), intent(in) Ll_unsatSTW,
    real(dp), dimension(:), intent(in) Ll_satSTW,
    real(dp), dimension(:), intent(in) Ll_neutrons,
    real(dp), dimension(:), intent(in) Ll_pet,
    real(dp), dimension(:, :), intent(in) Ll_aETSoil,
    real(dp), dimension(:), intent(in) Ll_aETCanopy,
    real(dp), dimension(:), intent(in) Ll_aETSealed,
    real(dp), dimension(:), intent(in) Ll_total_runoff,
    real(dp), dimension(:), intent(in) Ll_runoffSeal,
    real(dp), dimension(:), intent(in) Ll_fastRunoff,
    real(dp), dimension(:), intent(in) Ll_slowRunoff,
    real(dp), dimension(:), intent(in) Ll_baseflow,
    real(dp), dimension(:), intent(in) Ll_percol,
    real(dp), dimension(:, :), intent(in) Ll_infilSoil,
    real(dp), dimension(:), intent(in) Ll_preEffect )

```

Update all variables.

Call the type bound procedure updateVariable for all output variables. If a new output variable needs to be written, this is the second of two procedures to change (first: newOutputDataset)

## Parameters

|         |                                     |  |
|---------|-------------------------------------|--|
| in, out | <i>class(OutputDataset) :: self</i> |  |
| in      | <i>integer(i4) :: sidx, eidx</i>    |  |
| in      | <i>integer(i4) :: sidx, eidx</i>    |  |

## Parameters

|    |   |  |
|----|---|--|
| in | <i>real(dp), dimension(:) :: L1_fSealed</i>         |  |
| in | <i>real(dp), dimension(:) :: L1_fNotSealed</i>      |  |
| in | <i>real(dp), dimension(:) :: L1_inter</i>           |  |
| in | <i>real(dp), dimension(:) :: L1_snowPack</i>        |  |
| in | <i>real(dp), dimension(:, :) :: L1_soilMoist</i>    |  |
| in | <i>real(dp), dimension(:, :) :: L1_soilMoistSat</i> |  |
| in | <i>real(dp), dimension(:) :: L1_sealSTW</i>         |  |
| in | <i>real(dp), dimension(:) :: L1_unsatSTW</i>        |  |
| in | <i>real(dp), dimension(:) :: L1_satSTW</i>          |  |
| in | <i>real(dp), dimension(:) :: L1_neutrons</i>        |  |
| in | <i>real(dp), dimension(:) :: L1_pet</i>             |  |
| in | <i>real(dp), dimension(:, :) :: L1_aETSoil</i>      |  |
| in | <i>real(dp), dimension(:) :: L1_aETCanopy</i>       |  |
| in | <i>real(dp), dimension(:) :: L1_aETSealed</i>       |  |
| in | <i>real(dp), dimension(:) :: L1_total_runoff</i>    |  |
| in | <i>real(dp), dimension(:) :: L1_runoffSeal</i>      |  |
| in | <i>real(dp), dimension(:) :: L1_fastRunoff</i>      |  |
| in | <i>real(dp), dimension(:) :: L1_slowRunoff</i>      |  |
| in | <i>real(dp), dimension(:) :: L1_baseflow</i>        |  |
| in | <i>real(dp), dimension(:) :: L1_percol</i>          |  |
| in | <i>real(dp), dimension(:, :) :: L1_infilSoil</i>    |  |
| in | <i>real(dp), dimension(:) :: L1_preEffect</i>       |  |

## Authors

Matthias Zink

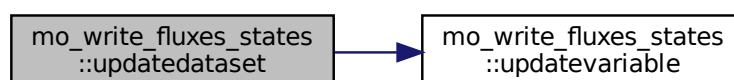
## Date

Apr 2013

Definition at line 520 of file mo\_write\_fluxes\_states.f90.

References mo\_mpr\_global\_variables::nsoilhorizons\_mhm, mo\_global\_variables::outputfluxstate, and updatevariable().

Here is the call graph for this function:



### 17.78.2.7 updatevariable()

```
subroutine mo_write_fluxes_states::updatevariable (
    class(outputvariable), intent(inout) self,
    real(dp), dimension(:), intent(in) data ) [private]
```

Update OutputVariable.

Add the array given as actual argument to the derived type's component 'data'

#### Returns

type(OutputVariable)

#### Parameters

|         |                                       |  |
|---------|---------------------------------------|--|
| in, out | <i>class(OutputVariable) :: self</i>  |  |
| in      | <i>real(dp), dimension(:) :: data</i> |  |

#### Authors

David Schaefer

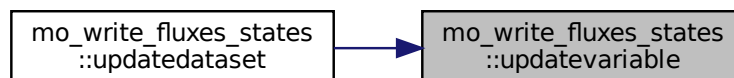
#### Date

June 2015

Definition at line 159 of file mo\_write\_fluxes\_states.f90.

Referenced by updatedataset().

Here is the caller graph for this function:



### 17.78.2.8 writetimestep()

```
subroutine mo_write_fluxes_states::writetimestep (
    class(outputdataset), intent(inout), target self,
    integer(i4), intent(in) timestep )
```

Write all accumulated data.

Write all accumulated and potentially averaged data to disk.

#### Returns

type(OutputVariable)

## Parameters

|         |                                     |                             |
|---------|-------------------------------------|-----------------------------|
| in, out | <i>class(OutputDataset) :: self</i> |                             |
| in      | <i>integer(i4) :: timestep</i>      | The model timestep to write |

## Authors

David Schaefer

## Date

June 2015

Definition at line 808 of file mo\_write\_fluxes\_states.f90.

**17.78.2.9 writevariableattributes()**

```
subroutine mo_write_fluxes_states::writevariableattributes (
    type(outputvariable), intent(in) var,
    character(*), intent(in) long_name,
    character(*), intent(in) unit )
```

Write output variable attributes.

TODO: add description

## Parameters

|    |  |                  |
|----|--|------------------|
| in | <i>type(OutputVariable) :: var</i>     |                  |
| in | <i>character(*) :: long_name, unit</i> | -> variable name |
| in | <i>character(*) :: long_name, unit</i> | -> physical unit |

## Authors

David Schaefer

## Date

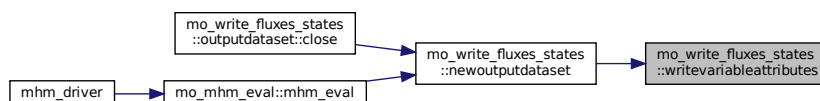
June 2015

Definition at line 1048 of file mo\_write\_fluxes\_states.f90.

References mo\_common\_constants::nodata\_dp.

Referenced by newoutputdataset().

Here is the caller graph for this function:



**17.78.2.10 writevariabletimestep()**

```
subroutine mo_write_fluxes_states::writevariabletimestep (
    class(outputvariable), intent(inout) self,
    integer(i4), intent(in) timestep ) [private]
```

Write timestep to file.

Write the content of the derived types's component 'data' to file, average if necessary

**Parameters**

|         |                                      |  |
|---------|--------------------------------------|--|
| in, out | <i>class(OutputVariable) :: self</i> |  |
| in      | <i>integer(i4) :: timestep</i>       | -> index along the time dimension of the netcdf variable |

**Authors**

David Schafer

**Date**

June 2015

Definition at line 196 of file mo\_write\_fluxes\_states.f90.

References mo\_common\_constants::nodata\_dp.



# Chapter 18

## Data Type Documentation

### 18.1 mo\_common\_mhm\_mrm\_restart::check\_consistency\_element Interface Reference

#### Private Member Functions

- subroutine [check\\_consistency\\_element\\_i4](#) (item1, item2, name, iBasin)
- subroutine [check\\_consistency\\_element\\_dp](#) (item1, item2, name, iBasin)

#### 18.1.1 Detailed Description

Definition at line 21 of file mo\_common\_mHM\_mRM\_restart.f90.

#### 18.1.2 Member Function/Subroutine Documentation

##### 18.1.2.1 check\_consistency\_element\_dp()

```
subroutine mo_common_mhm_mrm_restart::check_consistency_element::check_consistency_element_dp  
(  
    real(dp), intent(in) item1,  
    real(dp), intent(in) item2,  
    character(*), intent(in) name,  
    integer(i4), intent(in) iBasin ) [private]
```

Definition at line 90 of file mo\_common\_mHM\_mRM\_restart.f90.

##### 18.1.2.2 check\_consistency\_element\_i4()

```
subroutine mo_common_mhm_mrm_restart::check_consistency_element::check_consistency_element_i4  
(  
    integer(i4), intent(in) item1,  
    integer(i4), intent(in) item2,  
    character(*), intent(in) name,  
    integer(i4), intent(in) iBasin ) [private]
```

Definition at line 109 of file mo\_common\_mHM\_mRM\_restart.f90.

The documentation for this interface was generated from the following file:

- [src/common\\_mHM\\_mRM/mo\\_common\\_mHM\\_mRM\\_restart.f90](#)

## 18.2 mo\_common\_datetime\_type::datetimeinfo Type Reference

### Private Member Functions

- procedure [init](#) => [datetimeinfo\\_init](#)
- procedure [increment](#) => [datetimeinfo\\_increment](#)
- procedure [update\\_lai\\_timestep](#) => [datetimeinfo\\_update\\_lai\\_timestep](#)
- procedure [writeout](#) => [datetimeinfo\\_writeout](#)

### Private Attributes

- integer(i4) [ntimesteps](#)  
*number of timesteps in simulation period*
- real(dp) [newtime](#)  
*starts with simPer(iDomain)julStart, then increments with julday(...)*
- integer(i4) [day](#)  
*current day*
- integer(i4) [month](#)  
*current month*
- integer(i4) [year](#)  
*current year*
- integer(i4), private [hour](#)
- integer(i4), private [prev\\_day](#)
- integer(i4), private [prev\\_month](#)
- integer(i4), private [prev\\_year](#)
- logical [is\\_new\\_day](#)
- logical [is\\_new\\_month](#)
- logical [is\\_new\\_year](#)
- integer(i4) [ilai](#)
- integer(i4) [yid](#)
- integer(i4) [tindex\\_out](#)

### 18.2.1 Detailed Description

Definition at line 27 of file [mo\\_common\\_datetime\\_type.f90](#).

### 18.2.2 Member Function/Subroutine Documentation

#### 18.2.2.1 increment()

```
procedure mo_common_datetime_type::datetimeinfo::increment [private]
```

Definition at line 62 of file [mo\\_common\\_datetime\\_type.f90](#).

### 18.2.2.2 init()

```
procedure mo_common_datetime_type::datetimeinfo::init [private]
```

Definition at line 61 of file mo\_common\_datetime\_type.f90.

### 18.2.2.3 update\_lai\_timestep()

```
procedure mo_common_datetime_type::datetimeinfo::update_lai_timestep [private]
```

Definition at line 63 of file mo\_common\_datetime\_type.f90.

### 18.2.2.4 writeout()

```
procedure mo_common_datetime_type::datetimeinfo::writeout [private]
```

Definition at line 65 of file mo\_common\_datetime\_type.f90.

## 18.2.3 Member Data Documentation

### 18.2.3.1 day

```
integer(i4) mo_common_datetime_type::datetimeinfo::day [private]
```

current day

Definition at line 35 of file mo\_common\_datetime\_type.f90.

### 18.2.3.2 hour

```
integer(i4), private mo_common_datetime_type::datetimeinfo::hour [private]
```

Definition at line 41 of file mo\_common\_datetime\_type.f90.

### 18.2.3.3 ilai

```
integer(i4) mo_common_datetime_type::datetimeinfo::ilai [private]
```

Definition at line 54 of file mo\_common\_datetime\_type.f90.

### 18.2.3.4 is\_new\_day

```
logical mo_common_datetime_type::datetimeinfo::is_new_day [private]
```

Definition at line 50 of file mo\_common\_datetime\_type.f90.

### 18.2.3.5 is\_new\_month

```
logical mo_common_datetime_type::datetimeinfo::is_new_month [private]
```

Definition at line 51 of file mo\_common\_datetime\_type.f90.

### 18.2.3.6 is\_new\_year

```
logical mo_common_datetime_type::datetimeinfo::is_new_year [private]
```

Definition at line 52 of file mo\_common\_datetime\_type.f90.

### 18.2.3.7 month

```
integer(i4) mo_common_datetime_type::datetimeinfo::month [private]
```

current month

Definition at line 37 of file mo\_common\_datetime\_type.f90.

### 18.2.3.8 newtime

```
real(dp) mo_common_datetime_type::datetimeinfo::newtime [private]
```

starts with simPer(iDomain)julStart, then increments with julday(...)

Definition at line 32 of file mo\_common\_datetime\_type.f90.

### 18.2.3.9 ntimesteps

```
integer(i4) mo_common_datetime_type::datetimeinfo::ntimesteps [private]
```

number of timesteps in simulation period

Definition at line 29 of file mo\_common\_datetime\_type.f90.

### 18.2.3.10 prev\_day

```
integer(i4), private mo_common_datetime_type::datetimeinfo::prev_day [private]
```

Definition at line 45 of file mo\_common\_datetime\_type.f90.

### 18.2.3.11 prev\_month

```
integer(i4), private mo_common_datetime_type::datetimeinfo::prev_month [private]
```

Definition at line 46 of file mo\_common\_datetime\_type.f90.

### 18.2.3.12 prev\_year

```
integer(i4), private mo_common_datetime_type::datetimeinfo::prev_year [private]
```

Definition at line 47 of file mo\_common\_datetime\_type.f90.

### 18.2.3.13 tindex\_out

```
integer(i4) mo_common_datetime_type::datetimeinfo::tindex_out [private]
```

Definition at line 58 of file mo\_common\_datetime\_type.f90.

### 18.2.3.14 year

```
integer(i4) mo_common_datetime_type::datetimeinfo::year [private]
```

current year

Definition at line 39 of file mo\_common\_datetime\_type.f90.

### 18.2.3.15 yid

```
integer(i4) mo_common_datetime_type::datetimeinfo::yid [private]
```

Definition at line 55 of file mo\_common\_datetime\_type.f90.

The documentation for this type was generated from the following file:

- [src/common/mo\\_common\\_datetime\\_type.f90](#)

## 18.3 mo\_common\_variables::domain\_meta Type Reference

### Public Attributes

- integer(i4) [ndomains](#)
- integer(i4) [overallnumberofdomains](#)
- integer(i4), dimension(:), allocatable [indices](#)
- integer(i4), dimension(:), allocatable [I0datafrom](#)
- integer(i4), dimension(:), allocatable [optidata](#)
- logical, dimension(:), allocatable [dorouting](#)
- logical [ismasterincomlocal](#)
- type(mpi\_comm) [commaster](#)
- type(mpi\_comm) [comlocal](#)

### 18.3.1 Detailed Description

Definition at line 138 of file mo\_common\_variables.f90.

### 18.3.2 Member Data Documentation

### 18.3.2.1 comlocal

```
type(mpi_comm) mo_common_variables::domain_meta::comlocal
```

Definition at line 157 of file mo\_common\_variables.f90.

### 18.3.2.2 commaster

```
type(mpi_comm) mo_common_variables::domain_meta::commaster
```

Definition at line 155 of file mo\_common\_variables.f90.

### 18.3.2.3 dorouting

```
logical, dimension(:), allocatable mo_common_variables::domain_meta::dorouting
```

Definition at line 152 of file mo\_common\_variables.f90.

### 18.3.2.4 indices

```
integer(i4), dimension(:), allocatable mo_common_variables::domain_meta::indices
```

Definition at line 141 of file mo\_common\_variables.f90.

### 18.3.2.5 ismasterincomlocal

```
logical mo_common_variables::domain_meta::ismasterincomlocal
```

Definition at line 154 of file mo\_common\_variables.f90.

### 18.3.2.6 l0datafrom

```
integer(i4), dimension(:), allocatable mo_common_variables::domain_meta::l0datafrom
```

Definition at line 142 of file mo\_common\_variables.f90.

### 18.3.2.7 ndomains

```
integer(i4) mo_common_variables::domain_meta::ndomains
```

Definition at line 139 of file mo\_common\_variables.f90.

### 18.3.2.8 optidata

```
integer(i4), dimension(:), allocatable mo_common_variables::domain_meta::optidata
```

Definition at line 151 of file mo\_common\_variables.f90.

### 18.3.2.9 overallnumberofdomains

```
integer(i4) mo_common_variables::domain_meta::overallnumberofdomains
```

Definition at line 140 of file mo\_common\_variables.f90.

The documentation for this type was generated from the following file:

- [src/common/mo\\_common\\_variables.f90](#)

## 18.4 mo\_mrm\_global\_variables::domaininfo\_mrm Type Reference

### Public Attributes

- integer(i4) [ngauges](#)
- integer(i4), dimension(:), allocatable [gaugeidlist](#)
- integer(i4), dimension(:), allocatable [gaugeindexlist](#)
- integer(i4), dimension(:), allocatable [gaugenodelist](#)
- integer(i4) [ninflowgauges](#)
- integer(i4), dimension(:), allocatable [inflowgaugeidlist](#)
- integer(i4), dimension(:), allocatable [inflowgaugeindexlist](#)
- integer(i4), dimension(:), allocatable [inflowgaugenodelist](#)
- logical, dimension(:), allocatable [inflowgaugeheadwater](#)
- integer(i4) [l0\\_outlet](#)
- integer(i4), dimension(:), allocatable [l0\\_rowoutlet](#)
- integer(i4), dimension(:), allocatable [l0\\_coloutlet](#)

### 18.4.1 Detailed Description

Definition at line 93 of file mo\_mrm\_global\_variables.f90.

### 18.4.2 Member Data Documentation

#### 18.4.2.1 gaugeidlist

```
integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::gaugeidlist
```

Definition at line 97 of file mo\_mrm\_global\_variables.f90.

#### 18.4.2.2 gaugeindexlist

```
integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::gaugeindexlist
```

Definition at line 98 of file mo\_mrm\_global\_variables.f90.

### 18.4.2.3 gaugenodelist

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::gaugenodelist`

Definition at line 99 of file `mo_mrm_global_variables.f90`.

### 18.4.2.4 inflowgaugeheadwater

`logical, dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::inflowgaugeheadwater`

Definition at line 106 of file `mo_mrm_global_variables.f90`.

### 18.4.2.5 inflowgaugeidlist

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::inflowgaugeidlist`

Definition at line 103 of file `mo_mrm_global_variables.f90`.

### 18.4.2.6 inflowgaugeindexlist

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::inflowgaugeindexlist`

Definition at line 104 of file `mo_mrm_global_variables.f90`.

### 18.4.2.7 inflowgaugenodelist

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::inflowgaugenodelist`

Definition at line 105 of file `mo_mrm_global_variables.f90`.

### 18.4.2.8 l0\_coloutlet

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::l0_coloutlet`

Definition at line 112 of file `mo_mrm_global_variables.f90`.

### 18.4.2.9 l0\_noutlet

`integer(i4) mo_mrm_global_variables::domaininfo_mrm::l0_noutlet`

Definition at line 110 of file `mo_mrm_global_variables.f90`.

### 18.4.2.10 l0\_rowoutlet

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::domaininfo_mrm::l0_rowoutlet`

Definition at line 111 of file `mo_mrm_global_variables.f90`.



### 18.4.2.11 ngauges

```
integer(i4) mo_mrm_global_variables::domaininfo_mrm::ngauges
```

Definition at line 96 of file mo\_mrm\_global\_variables.f90.

### 18.4.2.12 ninflowgauges

```
integer(i4) mo_mrm_global_variables::domaininfo_mrm::ninflowgauges
```

Definition at line 102 of file mo\_mrm\_global\_variables.f90.

The documentation for this type was generated from the following file:

- [src/mRM/mo\\_mrm\\_global\\_variables.f90](#)

## 18.5 mo\_mrm\_global\_variables::gaugingstation Type Reference

### Public Attributes

- integer(i4), dimension(:), allocatable [domainid](#)
- integer(i4), dimension(:), allocatable [gaugeid](#)
- character(256), dimension(:), allocatable [fname](#)
- real(dp), dimension(:, :), allocatable [q](#)
- real(dp), dimension(:, :), allocatable [t](#)

### 18.5.1 Detailed Description

Definition at line 79 of file mo\_mrm\_global\_variables.f90.

### 18.5.2 Member Data Documentation

#### 18.5.2.1 domainid

```
integer(i4), dimension(:), allocatable mo_mrm_global_variables::gaugingstation::domainid
```

Definition at line 80 of file mo\_mrm\_global\_variables.f90.

#### 18.5.2.2 fname

```
character(256), dimension(:), allocatable mo_mrm_global_variables::gaugingstation::fname
```

Definition at line 82 of file mo\_mrm\_global\_variables.f90.

### 18.5.2.3 gaugeid

`integer(i4), dimension(:), allocatable mo_mrm_global_variables::gaugingstation::gaugeid`

Definition at line 81 of file `mo_mrm_global_variables.f90`.

### 18.5.2.4 q

`real(dp), dimension(:, :), allocatable mo_mrm_global_variables::gaugingstation::q`

Definition at line 83 of file `mo_mrm_global_variables.f90`.

### 18.5.2.5 t

`real(dp), dimension(:, :), allocatable mo_mrm_global_variables::gaugingstation::t`

Definition at line 85 of file `mo_mrm_global_variables.f90`.

The documentation for this type was generated from the following file:

- [src/mRM/mo\\_mrm\\_global\\_variables.f90](#)

## 18.6 mo\_common\_variables::grid Type Reference

### Public Attributes

- `integer(i4) ncols`
- `integer(i4) nrows`
- `integer(i4) ncells`
- `real(dp) xllcorner`
- `real(dp) yllcorner`
- `real(dp) cellsize`
- `real(dp) nodata_value`
- `real(dp), dimension(:, :), allocatable x`
- `real(dp), dimension(:, :), allocatable y`
- `logical, dimension(:, :), allocatable mask`
- `integer(i4) istart`
- `integer(i4) iend`
- `integer(i4), dimension(:, :), allocatable cellcoor`
- `real(dp), dimension(:), allocatable cellarea`
- `integer(i4), dimension(:), allocatable id`

### 18.6.1 Detailed Description

Definition at line 76 of file `mo_common_variables.f90`.

### 18.6.2 Member Data Documentation

### 18.6.2.1 cellarea

```
real(dp), dimension(:), allocatable mo_common_variables::grid::cellarea
```

Definition at line 93 of file mo\_common\_variables.f90.

### 18.6.2.2 cellcoor

```
integer(i4), dimension(:, :), allocatable mo_common_variables::grid::cellcoor
```

Definition at line 92 of file mo\_common\_variables.f90.

### 18.6.2.3 cellsize

```
real(dp) mo_common_variables::grid::cellsize
```

Definition at line 83 of file mo\_common\_variables.f90.

### 18.6.2.4 id

```
integer(i4), dimension(:), allocatable mo_common_variables::grid::id
```

Definition at line 94 of file mo\_common\_variables.f90.

### 18.6.2.5 iend

```
integer(i4) mo_common_variables::grid::iend
```

Definition at line 90 of file mo\_common\_variables.f90.

### 18.6.2.6 istart

```
integer(i4) mo_common_variables::grid::istart
```

Definition at line 89 of file mo\_common\_variables.f90.

### 18.6.2.7 mask

```
logical, dimension(:, :), allocatable mo_common_variables::grid::mask
```

Definition at line 87 of file mo\_common\_variables.f90.

### 18.6.2.8 ncells

```
integer(i4) mo_common_variables::grid::ncells
```

Definition at line 80 of file mo\_common\_variables.f90.

### 18.6.2.9 ncols

```
integer(i4) mo_common_variables::grid::ncols
```

Definition at line 78 of file mo\_common\_variables.f90.

### 18.6.2.10 nodata\_value

```
real(dp) mo_common_variables::grid::nodata_value
```

Definition at line 84 of file mo\_common\_variables.f90.

### 18.6.2.11 nrows

```
integer(i4) mo_common_variables::grid::nrows
```

Definition at line 79 of file mo\_common\_variables.f90.

### 18.6.2.12 x

```
real(dp), dimension(:, :), allocatable mo_common_variables::grid::x
```

Definition at line 85 of file mo\_common\_variables.f90.

### 18.6.2.13 xllcorner

```
real(dp) mo_common_variables::grid::xllcorner
```

Definition at line 81 of file mo\_common\_variables.f90.

### 18.6.2.14 y

```
real(dp), dimension(:, :), allocatable mo_common_variables::grid::y
```

Definition at line 86 of file mo\_common\_variables.f90.

### 18.6.2.15 yllcorner

```
real(dp) mo_common_variables::grid::yllcorner
```

Definition at line 82 of file mo\_common\_variables.f90.

The documentation for this type was generated from the following file:

- [src/common/mo\\_common\\_variables.f90](#)

## 18.7 mo\_common\_variables::gridmapper Type Reference

### Public Attributes

- type([grid](#)), pointer [high\\_res\\_grid](#)
- type([grid](#)), pointer [low\\_res\\_grid](#)
- integer(i4), dimension(:), allocatable [lower\\_bound](#)
- integer(i4), dimension(:), allocatable [upper\\_bound](#)
- integer(i4), dimension(:), allocatable [left\\_bound](#)
- integer(i4), dimension(:), allocatable [right\\_bound](#)
- integer(i4), dimension(:), allocatable [n\\_subcells](#)
- integer(i4), dimension(:, :), allocatable [lowres\\_id\\_on\\_highres](#)

### 18.7.1 Detailed Description

Definition at line 101 of file mo\_common\_variables.f90.

### 18.7.2 Member Data Documentation

#### 18.7.2.1 high\_res\_grid

type([grid](#)), pointer mo\_common\_variables::gridmapper::high\_res\_grid

Definition at line 102 of file mo\_common\_variables.f90.

#### 18.7.2.2 left\_bound

integer(i4), dimension(:), allocatable mo\_common\_variables::gridmapper::left\_bound

Definition at line 108 of file mo\_common\_variables.f90.

#### 18.7.2.3 low\_res\_grid

type([grid](#)), pointer mo\_common\_variables::gridmapper::low\_res\_grid

Definition at line 103 of file mo\_common\_variables.f90.

#### 18.7.2.4 lower\_bound

integer(i4), dimension(:), allocatable mo\_common\_variables::gridmapper::lower\_bound

Definition at line 106 of file mo\_common\_variables.f90.

### 18.7.2.5 lowres\_id\_on\_highres

```
integer(i4), dimension(:, :), allocatable mo_common_variables::gridmapper::lowres_id_on_↵
highres
```

Definition at line 111 of file mo\_common\_variables.f90.

### 18.7.2.6 n\_subcells

```
integer(i4), dimension(:), allocatable mo_common_variables::gridmapper::n_subcells
```

Definition at line 110 of file mo\_common\_variables.f90.

### 18.7.2.7 right\_bound

```
integer(i4), dimension(:), allocatable mo_common_variables::gridmapper::right_bound
```

Definition at line 109 of file mo\_common\_variables.f90.

### 18.7.2.8 upper\_bound

```
integer(i4), dimension(:), allocatable mo_common_variables::gridmapper::upper_bound
```

Definition at line 107 of file mo\_common\_variables.f90.

The documentation for this type was generated from the following file:

- [src/common/mo\\_common\\_variables.f90](#)

## 18.8 mo\_template::mean Interface Reference

The average.

### Public Member Functions

- real(sp) function [mean\\_sp](#) (dat, mask)
- real(dp) function [mean\\_dp](#) (dat, mask)

### 18.8.1 Detailed Description

The average.

Calculates the average value of a vector, i.e. the first moment of a series of numbers:

$$\bar{x} = \frac{1}{N} \sum_{i=1}^N x_i$$

If an optional mask is given, the mean is only over those locations that correspond to true values in the mask. x can be single or double precision. The result will have the same numerical precision. ADDITIONAL INFORMATION  
 vec = (/ 1., 2, 3., -999., 5., 6. /) m = mean(vec, mask=(vec >= 0.)) -> see also example in test directory Sokal

RR & Rohlf FJ - Biometry: the principle and practice of statistics in biological research, Freeman & Co., ISBN 0-7167-2411-1 Press WH, Teukolsky SA, Vetterling WT, & Flannery BP - Numerical Recipes in Fortran 90 - The Art of Parallel Scientific Computing, 2nd Edition, Volume 2 of Fortran Numerical Recipes, Cambridge University Press, UK, 1996

#### Returns

real(sp/dp) :: mean —  $\bar{x}$  average of all elements in vec

#### Authors

Matthias Cuntz

#### Date

Nov 2011

Definition at line 79 of file mo\_template.f90.

## 18.8.2 Member Function/Subroutine Documentation

### 18.8.2.1 mean\_dp()

```
real(dp) function mo_template::mean::mean_dp (  
    real(dp), dimension(:), intent(in) dat,  
    logical, dimension(:), intent(in), optional mask )
```

Definition at line 137 of file mo\_template.f90.

References mo\_template::mean\_dp().

Here is the call graph for this function:



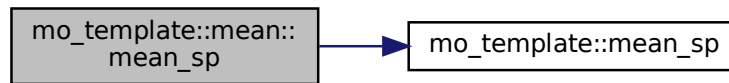
### 18.8.2.2 mean\_sp()

```
real(sp) function mo_template::mean::mean_sp (  
    real(sp), dimension(:), intent(in) dat,  
    logical, dimension(:), intent(in), optional mask )
```

Definition at line 167 of file mo\_template.f90.

References mo\_template::mean\_sp().

Here is the call graph for this function:



The documentation for this interface was generated from the following file:

- [src/common/mo\\_template.f90](#)

## 18.9 mo\_mrm\_write\_fluxes\_states::outputdataset Interface Reference

### Public Member Functions

- procedure, public [updatedataset](#)
- procedure, public [writetimestep](#)
- procedure, public [close](#)

### Public Attributes

- integer(i4) [idomain](#)
- integer(i4) [domain](#)
- integer(i4) [id](#)
- type(ncdataset) [nc](#)
- type(ncdataset) [ncdataset](#)
- type(ncdataset) [to](#)
- type(ncdataset) [write](#)
- type(outputvariable), dimension(:), allocatable [vars](#)
- type(outputvariable), allocatable [store](#)
- type(outputvariable), allocatable [all](#)
- type(outputvariable), dimension(dynamic), allocatable [created](#)
- type(outputvariable), allocatable [variables](#)
- integer(i4) [counter](#) = 0
- integer(i4) [count](#)
- integer(i4) [written](#)
- integer(i4) [time](#)
- integer(i4) [steps](#)

### 18.9.1 Detailed Description

Definition at line 45 of file [mo\\_mrm\\_write\\_fluxes\\_states.f90](#).

### 18.9.2 Member Function/Subroutine Documentation



### 18.9.2.1 close()

```
procedure, public mo_mrm_write_fluxes_states::outputdataset::close
```

Definition at line 54 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.2.2 updatedataset()

```
procedure, public mo_mrm_write_fluxes_states::outputdataset::updatedataset
```

Definition at line 52 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.2.3 writetimestep()

```
procedure, public mo_mrm_write_fluxes_states::outputdataset::writetimestep
```

Definition at line 53 of file mo\_mrm\_write\_fluxes\_states.f90.

## 18.9.3 Member Data Documentation

### 18.9.3.1 all

```
type(outputvariable), allocatable mo_mrm_write_fluxes_states::outputdataset::all
```

Definition at line 48 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.2 count

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::count
```

Definition at line 49 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.3 counter

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::counter = 0
```

Definition at line 49 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.4 created

```
type(outputvariable), dimension (dynamic), allocatable mo_mrm_write_fluxes_states::outputdataset←  
::created
```

Definition at line 48 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.5 domain

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::domain
```

Definition at line 46 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.6 id

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::id
```

Definition at line 46 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.7 idomain

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::idomain
```

Definition at line 46 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.8 nc

```
type(ncdataset) mo_mrm_write_fluxes_states::outputdataset::nc
```

Definition at line 47 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.9 ncdataset

```
type(ncdataset) mo_mrm_write_fluxes_states::outputdataset::ncdataset
```

Definition at line 47 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.10 steps

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::steps
```

Definition at line 49 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.11 store

```
type(outputvariable), allocatable mo_mrm_write_fluxes_states::outputdataset::store
```

Definition at line 48 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.12 time

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::time
```

Definition at line 49 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.13 to

```
type(ncdataset) mo_mrm_write_fluxes_states::outputdataset::to
```

Definition at line 47 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.14 variables

```
type(outputvariable), allocatable mo_mrm_write_fluxes_states::outputdataset::variables
```

Definition at line 48 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.15 vars

```
type(outputvariable), dimension(:), allocatable mo_mrm_write_fluxes_states::outputdataset↔  
::vars
```

Definition at line 48 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.16 write

```
type(ncdataset) mo_mrm_write_fluxes_states::outputdataset::write
```

Definition at line 47 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.9.3.17 written

```
integer(i4) mo_mrm_write_fluxes_states::outputdataset::written
```

Definition at line 49 of file mo\_mrm\_write\_fluxes\_states.f90.

The documentation for this interface was generated from the following file:

- [src/mRM/mo\\_mrm\\_write\\_fluxes\\_states.f90](#)

## 18.10 mo\_write\_fluxes\_states::outputdataset Interface Reference

### Public Member Functions

- procedure, public [updatedataset](#)
- procedure, public [writetimestep](#)
- procedure, public [close](#)

### Public Attributes

- integer(i4) [idomain](#)
- integer(i4) [domain](#)

- integer(i4) `id`
- type(ncdataset) `nc`
- type(ncdataset) `ncdataset`
- type(ncdataset) `to`
- type(ncdataset) `write`
- type(outputvariable), dimension(:), allocatable `vars`
- type(outputvariable), allocatable `store`
- type(outputvariable), allocatable `all`
- type(outputvariable), dimension(dynamic), allocatable `created`
- type(outputvariable), allocatable `variables`
- integer(i4) `counter` = 0
- integer(i4) `count`
- integer(i4) `written`
- integer(i4) `time`
- integer(i4) `steps`

### 18.10.1 Detailed Description

Definition at line 45 of file `mo_write_fluxes_states.f90`.

### 18.10.2 Member Function/Subroutine Documentation

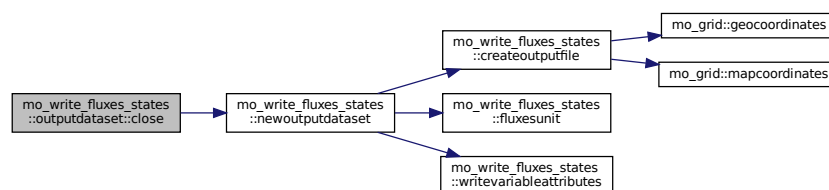
#### 18.10.2.1 `close()`

```
procedure, public mo_write_fluxes_states::outputdataset::close
```

Definition at line 54 of file `mo_write_fluxes_states.f90`.

References `mo_write_fluxes_states::newoutputdataset()`.

Here is the call graph for this function:



#### 18.10.2.2 `updatedataset()`

```
procedure, public mo_write_fluxes_states::outputdataset::updatedataset
```

Definition at line 52 of file `mo_write_fluxes_states.f90`.

### 18.10.2.3 writetimestep()

procedure, public mo\_write\_fluxes\_states::outputdataset::writetimestep

Definition at line 53 of file mo\_write\_fluxes\_states.f90.

## 18.10.3 Member Data Documentation

### 18.10.3.1 all

type(outputvariable), allocatable mo\_write\_fluxes\_states::outputdataset::all

Definition at line 48 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.2 count

integer(i4) mo\_write\_fluxes\_states::outputdataset::count

Definition at line 49 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.3 counter

integer(i4) mo\_write\_fluxes\_states::outputdataset::counter = 0

Definition at line 49 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.4 created

type(outputvariable), dimension (dynamic), allocatable mo\_write\_fluxes\_states::outputdataset←  
::created

Definition at line 48 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.5 domain

integer(i4) mo\_write\_fluxes\_states::outputdataset::domain

Definition at line 46 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.6 id

integer(i4) mo\_write\_fluxes\_states::outputdataset::id

Definition at line 46 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.7 idomain

```
integer(i4) mo_write_fluxes_states::outputdataset::idomain
```

Definition at line 46 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.8 nc

```
type(ncdataset) mo_write_fluxes_states::outputdataset::nc
```

Definition at line 47 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.9 ncdataset

```
type(ncdataset) mo_write_fluxes_states::outputdataset::ncdataset
```

Definition at line 47 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.10 steps

```
integer(i4) mo_write_fluxes_states::outputdataset::steps
```

Definition at line 49 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.11 store

```
type(outputvariable), allocatable mo_write_fluxes_states::outputdataset::store
```

Definition at line 48 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.12 time

```
integer(i4) mo_write_fluxes_states::outputdataset::time
```

Definition at line 49 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.13 to

```
type(ncdataset) mo_write_fluxes_states::outputdataset::to
```

Definition at line 47 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.14 variables

```
type(outputvariable), allocatable mo_write_fluxes_states::outputdataset::variables
```

Definition at line 48 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.15 vars

```
type(outputvariable), dimension(:), allocatable mo_write_fluxes_states::outputdataset::vars
```

Definition at line 48 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.16 write

```
type(ncdataset) mo_write_fluxes_states::outputdataset::write
```

Definition at line 47 of file mo\_write\_fluxes\_states.f90.

### 18.10.3.17 written

```
integer(i4) mo_write_fluxes_states::outputdataset::written
```

Definition at line 49 of file mo\_write\_fluxes\_states.f90.

The documentation for this interface was generated from the following file:

- [src/mHM/mo\\_write\\_fluxes\\_states.f90](#)

## 18.11 mo\_mrm\_write\_fluxes\_states::outputvariable Interface Reference

### Public Member Functions

- procedure, public [updatevariable](#)
- procedure, public [writevariabletimestep](#)

### Public Attributes

- type(ncvariable) [nc](#)
- type(ncvariable) [ncdataset](#)
- type(ncvariable) [which](#)
- type(ncvariable) [contains](#)
- type(ncvariable), allocatable [the](#)
- type(ncvariable) [variable](#)
- logical [avg](#) = .false.
- logical [average](#)
- logical, dimension(:), allocatable, pointer [data](#)
- logical [before](#)
- logical [writing](#)
- logical, dimension(:, :), pointer [mask](#)
- logical, pointer [to](#)
- logical, pointer [reconstruct](#)
- real(dp), dimension(:), allocatable, pointer [data](#)
- real(dp), allocatable [store](#)
- real(dp), allocatable [the](#)
- real(dp), allocatable [between](#)

- `real(dp)`, allocatable [writes](#)
- `integer(i4)` `counter = 0`
- `integer(i4)` `count`
- `integer(i4)`, allocatable [the](#)
- `integer(i4)` `number`
- `integer(i4)` `of`
- `integer(i4)`, public [updatevariable](#)
- `integer(i4)` `calls`

### 18.11.1 Detailed Description

Definition at line 27 of file `mo_mrm_write_fluxes_states.f90`.

### 18.11.2 Member Function/Subroutine Documentation

#### 18.11.2.1 `updatevariable()`

```
procedure, public mo_mrm_write_fluxes_states::outputvariable::updatevariable
```

Definition at line 35 of file `mo_mrm_write_fluxes_states.f90`.

#### 18.11.2.2 `writevariabletimestep()`

```
procedure, public mo_mrm_write_fluxes_states::outputvariable::writevariabletimestep
```

Definition at line 36 of file `mo_mrm_write_fluxes_states.f90`.

### 18.11.3 Member Data Documentation

#### 18.11.3.1 `average`

```
logical mo_mrm_write_fluxes_states::outputvariable::average
```

Definition at line 29 of file `mo_mrm_write_fluxes_states.f90`.

#### 18.11.3.2 `avg`

```
logical mo_mrm_write_fluxes_states::outputvariable::avg = .false.
```

Definition at line 29 of file `mo_mrm_write_fluxes_states.f90`.



### 18.11.3.3 before

```
logical mo_mrm_write_fluxes_states::outputvariable::before
```

Definition at line 29 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.4 between

```
real(dp), allocatable mo_mrm_write_fluxes_states::outputvariable::between
```

Definition at line 31 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.5 calls

```
integer(i4) mo_mrm_write_fluxes_states::outputvariable::calls
```

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.6 contains

```
type(ncvariable) mo_mrm_write_fluxes_states::outputvariable::contains
```

Definition at line 28 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.7 count

```
integer(i4) mo_mrm_write_fluxes_states::outputvariable::count
```

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.8 counter

```
integer(i4) mo_mrm_write_fluxes_states::outputvariable::counter = 0
```

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.9 data [1/2]

```
logical, dimension(:), allocatable, pointer mo_mrm_write_fluxes_states::outputvariable::data
```

Definition at line 29 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.10 data [2/2]

```
real(dp), dimension(:), allocatable, pointer mo_mrm_write_fluxes_states::outputvariable::data
```

Definition at line 31 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.11 mask**

logical, dimension(:, :), pointer mo\_mrm\_write\_fluxes\_states::outputvariable::mask

Definition at line 30 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.12 nc**

type(ncvariable) mo\_mrm\_write\_fluxes\_states::outputvariable::nc

Definition at line 28 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.13 ncdataset**

type(ncvariable) mo\_mrm\_write\_fluxes\_states::outputvariable::ncdataset

Definition at line 28 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.14 number**

integer(i4) mo\_mrm\_write\_fluxes\_states::outputvariable::number

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.15 of**

integer(i4) mo\_mrm\_write\_fluxes\_states::outputvariable::of

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.16 reconstruct**

logical, pointer mo\_mrm\_write\_fluxes\_states::outputvariable::reconstruct

Definition at line 30 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.17 store**

real(dp), allocatable mo\_mrm\_write\_fluxes\_states::outputvariable::store

Definition at line 31 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.18 the** [1/3]

```
type(ncvariable), allocatable mo_mrm_write_fluxes_states::outputvariable::the
```

Definition at line 28 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.19 the** [2/3]

```
real(dp), allocatable mo_mrm_write_fluxes_states::outputvariable::the
```

Definition at line 31 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.20 the** [3/3]

```
integer(i4), allocatable mo_mrm_write_fluxes_states::outputvariable::the
```

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.21 to**

```
logical, pointer mo_mrm_write_fluxes_states::outputvariable::to
```

Definition at line 30 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.22 updatevariable**

```
integer(i4), public mo_mrm_write_fluxes_states::outputvariable::updatevariable
```

Definition at line 32 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.23 variable**

```
type(ncvariable) mo_mrm_write_fluxes_states::outputvariable::variable
```

Definition at line 28 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.24 which**

```
type(ncvariable) mo_mrm_write_fluxes_states::outputvariable::which
```

Definition at line 28 of file mo\_mrm\_write\_fluxes\_states.f90.

**18.11.3.25 writes**

```
real(dp), allocatable mo_mrm_write_fluxes_states::outputvariable::writes
```

Definition at line 31 of file mo\_mrm\_write\_fluxes\_states.f90.

### 18.11.3.26 writing

logical mo\_mrm\_write\_fluxes\_states::outputvariable::writing

Definition at line 29 of file mo\_mrm\_write\_fluxes\_states.f90.

The documentation for this interface was generated from the following file:

- [src/mRM/mo\\_mrm\\_write\\_fluxes\\_states.f90](#)

## 18.12 mo\_write\_fluxes\_states::outputvariable Interface Reference

### Public Member Functions

- procedure, public [updatevariable](#)
- procedure, public [writevariabletimestep](#)

### Public Attributes

- type(ncvariable) [nc](#)
- type(ncvariable) [ncdataset](#)
- type(ncvariable) [which](#)
- type(ncvariable) [contains](#)
- type(ncvariable), allocatable [the](#)
- type(ncvariable) [variable](#)
- logical [avg](#) = .false.
- logical [average](#)
- logical, dimension(:), allocatable, pointer [data](#)
- logical [before](#)
- logical [writing](#)
- logical, dimension(:, :), pointer [mask](#)
- logical, pointer [to](#)
- logical, pointer [reconstruct](#)
- real(dp), dimension(:), allocatable, pointer [data](#)
- real(dp), allocatable [store](#)
- real(dp), allocatable [the](#)
- real(dp), allocatable [between](#)
- real(dp), allocatable [writes](#)
- integer(i4) [counter](#) = 0
- integer(i4) [count](#)
- integer(i4), allocatable [the](#)
- integer(i4) [number](#)
- integer(i4) [of](#)
- integer(i4), public [updatevariable](#)
- integer(i4) [calls](#)

### 18.12.1 Detailed Description

Definition at line 27 of file mo\_write\_fluxes\_states.f90.

## 18.12.2 Member Function/Subroutine Documentation

### 18.12.2.1 updatevariable()

procedure, public mo\_write\_fluxes\_states::outputvariable::updatevariable

Definition at line 35 of file mo\_write\_fluxes\_states.f90.

### 18.12.2.2 writevariabletimestep()

procedure, public mo\_write\_fluxes\_states::outputvariable::writevariabletimestep

Definition at line 36 of file mo\_write\_fluxes\_states.f90.

## 18.12.3 Member Data Documentation

### 18.12.3.1 average

logical mo\_write\_fluxes\_states::outputvariable::average

Definition at line 29 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.2 avg

logical mo\_write\_fluxes\_states::outputvariable::avg = .false.

Definition at line 29 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.3 before

logical mo\_write\_fluxes\_states::outputvariable::before

Definition at line 29 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.4 between

real(dp), allocatable mo\_write\_fluxes\_states::outputvariable::between

Definition at line 31 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.5 calls

integer(i4) mo\_write\_fluxes\_states::outputvariable::calls

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.6 contains

```
type(ncvariable) mo_write_fluxes_states::outputvariable::contains
```

Definition at line 28 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.7 count

```
integer(i4) mo_write_fluxes_states::outputvariable::count
```

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.8 counter

```
integer(i4) mo_write_fluxes_states::outputvariable::counter = 0
```

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.9 data [1/2]

```
logical, dimension(:), allocatable, pointer mo_write_fluxes_states::outputvariable::data
```

Definition at line 29 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.10 data [2/2]

```
real(dp), dimension(:), allocatable, pointer mo_write_fluxes_states::outputvariable::data
```

Definition at line 31 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.11 mask

```
logical, dimension(:, :), pointer mo_write_fluxes_states::outputvariable::mask
```

Definition at line 30 of file mo\_write\_fluxes\_states.f90.

#### 18.12.3.12 nc

```
type(ncvariable) mo_write_fluxes_states::outputvariable::nc
```

Definition at line 28 of file mo\_write\_fluxes\_states.f90.

**18.12.3.13 ncdataset**

```
type(ncvariable) mo_write_fluxes_states::outputvariable::ncdataset
```

Definition at line 28 of file mo\_write\_fluxes\_states.f90.

**18.12.3.14 number**

```
integer(i4) mo_write_fluxes_states::outputvariable::number
```

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

**18.12.3.15 of**

```
integer(i4) mo_write_fluxes_states::outputvariable::of
```

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

**18.12.3.16 reconstruct**

```
logical, pointer mo_write_fluxes_states::outputvariable::reconstruct
```

Definition at line 30 of file mo\_write\_fluxes\_states.f90.

**18.12.3.17 store**

```
real(dp), allocatable mo_write_fluxes_states::outputvariable::store
```

Definition at line 31 of file mo\_write\_fluxes\_states.f90.

**18.12.3.18 the [1/3]**

```
type(ncvariable), allocatable mo_write_fluxes_states::outputvariable::the
```

Definition at line 28 of file mo\_write\_fluxes\_states.f90.

**18.12.3.19 the [2/3]**

```
real(dp), allocatable mo_write_fluxes_states::outputvariable::the
```

Definition at line 31 of file mo\_write\_fluxes\_states.f90.

**18.12.3.20 the [3/3]**

```
integer(i4), allocatable mo_write_fluxes_states::outputvariable::the
```

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.21 to

logical, pointer mo\_write\_fluxes\_states::outputvariable::to

Definition at line 30 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.22 updatevariable

integer(i4), public mo\_write\_fluxes\_states::outputvariable::updatevariable

Definition at line 32 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.23 variable

type(ncvariable) mo\_write\_fluxes\_states::outputvariable::variable

Definition at line 28 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.24 which

type(ncvariable) mo\_write\_fluxes\_states::outputvariable::which

Definition at line 28 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.25 writes

real(dp), allocatable mo\_write\_fluxes\_states::outputvariable::writes

Definition at line 31 of file mo\_write\_fluxes\_states.f90.

### 18.12.3.26 writing

logical mo\_write\_fluxes\_states::outputvariable::writing

Definition at line 29 of file mo\_write\_fluxes\_states.f90.

The documentation for this interface was generated from the following file:

- [src/mHM/mo\\_write\\_fluxes\\_states.f90](#)

## 18.13 mo\_common\_variables::period Type Reference

### Public Attributes

- integer(i4) [dstart](#)
- integer(i4) [mstart](#)



- integer(i4) [ystart](#)
- integer(i4) [dend](#)
- integer(i4) [mend](#)
- integer(i4) [yend](#)
- integer(i4) [julstart](#)
- integer(i4) [julend](#)
- integer(i4) [nobs](#)

### 18.13.1 Detailed Description

Definition at line 61 of file mo\_common\_variables.f90.

### 18.13.2 Member Data Documentation

#### 18.13.2.1 dend

```
integer(i4) mo_common_variables::period::dend
```

Definition at line 65 of file mo\_common\_variables.f90.

#### 18.13.2.2 dstart

```
integer(i4) mo_common_variables::period::dstart
```

Definition at line 62 of file mo\_common\_variables.f90.

#### 18.13.2.3 julend

```
integer(i4) mo_common_variables::period::julend
```

Definition at line 69 of file mo\_common\_variables.f90.

#### 18.13.2.4 julstart

```
integer(i4) mo_common_variables::period::julstart
```

Definition at line 68 of file mo\_common\_variables.f90.

#### 18.13.2.5 mend

```
integer(i4) mo_common_variables::period::mend
```

Definition at line 66 of file mo\_common\_variables.f90.

### 18.13.2.6 mstart

```
integer(i4) mo_common_variables::period::mstart
```

Definition at line 63 of file mo\_common\_variables.f90.

### 18.13.2.7 nobS

```
integer(i4) mo_common_variables::period::nobS
```

Definition at line 70 of file mo\_common\_variables.f90.

### 18.13.2.8 yend

```
integer(i4) mo_common_variables::period::yend
```

Definition at line 67 of file mo\_common\_variables.f90.

### 18.13.2.9 ystart

```
integer(i4) mo_common_variables::period::ystart
```

Definition at line 64 of file mo\_common\_variables.f90.

The documentation for this type was generated from the following file:

- [src/common/mo\\_common\\_variables.f90](#)

## 18.14 mo\_read\_spatial\_data::read\_spatial\_data\_ascii Interface Reference

Reads spatial data files of ASCII format.

### Public Member Functions

- subroutine [read\\_spatial\\_data\\_ascii\\_i4](#) (filename, fileunit, header\_ncols, header\_nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, data, mask)

*TODO: add description.*

- subroutine [read\\_spatial\\_data\\_ascii\\_dp](#) (filename, fileunit, header\_ncols, header\_nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, data, mask)

*TODO: add description.*

### 18.14.1 Detailed Description

Reads spatial data files of ASCII format.

Reads spatial input data, e.g. dem, aspect, flow direction.

**Authors**

Juliane Mai

**Date**

Jan 2013

Definition at line 52 of file mo\_read\_spatial\_data.f90.

**18.14.2 Member Function/Subroutine Documentation****18.14.2.1 read\_spatial\_data\_ascii\_dp()**

```

subroutine mo_read_spatial_data::read_spatial_data_ascii::read_spatial_data_ascii_dp (
    character(len = *), intent(in) filename,
    integer(i4), intent(in) fileunit,
    integer(i4), intent(in) header_ncols,
    integer(i4), intent(in) header_nrows,
    real(dp), intent(in) header_xllcorner,
    real(dp), intent(in) header_yllcorner,
    real(dp), intent(in) header_cellsize,
    real(dp), dimension(:, :), intent(out), allocatable data,
    logical, dimension(:, :), intent(out), allocatable mask )

```

TODO: add description.

TODO: add description

**Parameters**

|     |  |                                   |
|-----|--|-----------------------------------|
| in  | <i>character(len = *) :: filename</i>    | filename with location            |
| in  | <i>integer(i4) :: fileunit</i>           | unit for opening the file         |
| in  | <i>integer(i4) :: header_nCols</i>       | number of columns of data fields: |
| in  | <i>integer(i4) :: header_nRows</i>       | number of rows of data fields:    |
| in  | <i>real(dp) :: header_xllcorner</i>      | header read in lower left corner  |
| in  | <i>real(dp) :: header_yllcorner</i>      | header read in lower left corner  |
| in  | <i>real(dp) :: header_cellsize</i>       | header read in cellsize           |
| out | <i>real(dp), dimension(:, :) :: data</i> | data                              |
| out | <i>logical, dimension(:, :) :: mask</i>  | mask                              |

**Authors**

Robert Schweppe

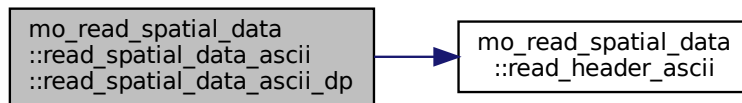
**Date**

Jun 2018

Definition at line 94 of file mo\_read\_spatial\_data.f90.

References mo\_read\_spatial\_data::read\_header\_ascii().

Here is the call graph for this function:



### 18.14.2.2 read\_spatial\_data\_ascii\_i4()

```

subroutine mo_read_spatial_data::read_spatial_data_ascii::read_spatial_data_ascii_i4 (
    character(len = *) , intent(in) filename,
    integer(i4) , intent(in) fileunit,
    integer(i4) , intent(in) header_ncols,
    integer(i4) , intent(in) header_nrows,
    real(dp) , intent(in) header_xllcorner,
    real(dp) , intent(in) header_yllcorner,
    real(dp) , intent(in) header_cellsize,
    integer(i4) , dimension(:, :), intent(out), allocatable data,
    logical, dimension(:, :), intent(out), allocatable mask )
  
```

TODO: add description.

TODO: add description

#### Parameters

|     |   |                                   |
|-----|---|-----------------------------------|
| in  | <i>character(len = *) :: filename</i>       | filename with location            |
| in  | <i>integer(i4) :: fileunit</i>              | unit for opening the file         |
| in  | <i>integer(i4) :: header_nCols</i>          | number of columns of data fields: |
| in  | <i>integer(i4) :: header_nRows</i>          | number of rows of data fields:    |
| in  | <i>real(dp) :: header_xllcorner</i>         | header read in lower left corner  |
| in  | <i>real(dp) :: header_yllcorner</i>         | header read in lower left corner  |
| in  | <i>real(dp) :: header_cellsize</i>          | header read in cellsize           |
| out | <i>integer(i4), dimension(:, :) :: data</i> | data                              |
| out | <i>logical, dimension(:, :) :: mask</i>     | mask                              |

#### Authors

Robert Schweppe

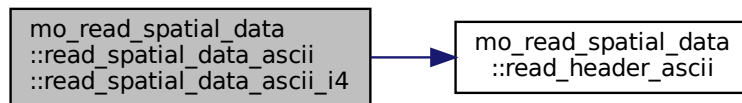
#### Date

Jun 2018

Definition at line 234 of file mo\_read\_spatial\_data.f90.

References mo\_read\_spatial\_data::read\_header\_ascii().

Here is the call graph for this function:



The documentation for this interface was generated from the following file:

- [src/common/mo\\_read\\_spatial\\_data.f90](#)

## 18.15 mo\_mrm\_riv\_temp\_class::riv\_temp\_type Module Reference

This is a container to define the river temperature routing in the current time step.

### Public Member Functions

- procedure [config](#)  
*configure the riv\_temp\_type class from the mhm namelist*
- procedure [init](#)  
*initialize the riv\_temp\_type class for the current domain*
- procedure [init\\_area](#)  
*initialize the river area of riv\_temp\_type class for the current domain*
- procedure [init\\_riv\\_temp](#)  
*initialize the river temperature of riv\_temp\_type class for the current domain*
- procedure [acc\\_source\\_e](#)  
*accumulate energy sources of riv\_temp\_type*
- procedure [finalize\\_source\\_e](#)  
*finalize energy sources of riv\_temp\_type*
- procedure [get\\_lrad\\_out](#)  
*get outgoing longwave radiation of riv\_temp\_type*
- procedure [get\\_lat\\_heat](#)  
*latent heat flux of riv\_temp\_type*
- procedure [get\\_sens\\_heat](#)  
*sensible heat flux of riv\_temp\_type*
- procedure [get\\_e\\_io](#)  
*get complete energy source of riv\_temp\_type at given cell*
- procedure [l11\\_routing\\_e](#)  
*execute the temperature routing of riv\_temp\_type*
- procedure [init\\_iter](#)  
*initialize iterative solver of riv\_temp\_type*
- procedure [next\\_iter](#)  
*execute next iteration with iterative solver of riv\_temp\_type*
- procedure [reset\\_timestep](#)  
*reset riv\_temp\_type class for next timestep*

- procedure [alloc\\_lateral](#)  
*allocate lateral temp components of [riv\\_temp\\_type](#) class for current domain*
- procedure [dealloc\\_lateral](#)  
*deallocate lateral temp components of [riv\\_temp\\_type](#)*

## Public Attributes

- logical [active](#) = .false.  
*state if this process is active*
- integer(i4) [case](#) = 0\_i4  
*the selected process-case option*
- character(256) [nml\\_name](#) = 'config\_riv\_temp'  
*namelist name in mhm namelist*
- character(256), dimension(:), allocatable [dir\\_riv\\_widths](#)  
*Directory where river widths are stored.*
- character(256) [riv\\_widths\\_file](#)  
*file name for river widths*
- character(256) [riv\\_widths\\_name](#)  
*variable name for river widths*
- real(dp), dimension(:), allocatable, public [l11\\_riv\\_widths](#)  
*river widths in L11*
- real(dp), dimension(:), allocatable, public [l11\\_riv\\_areas](#)  
*river area in L11*
- real(dp) [albedo\\_water](#)  
*albedo of open water*
- real(dp) [pt\\_a\\_water](#)  
*priestley taylor alpha parameter for PET on open water*
- real(dp) [emissivity\\_water](#)  
*emissivity of water*
- real(dp) [turb\\_heat\\_ex\\_coeff](#)  
*lateral heat exchange coefficient water <-> air*
- real(dp) [delta\\_t](#) = 0.1\_dp  
*cutoff value for temperature*
- integer(i4) [max\\_iter](#)  
*input: maximal number of iterations done*
- real(dp) [delta\\_iter](#)  
*input: convergence criteria for iterative solver*
- real(dp) [step\\_iter](#)  
*input: step-size for linear search*
- logical [first\\_iter](#)  
*whether it is at the first iteration (to determine search direction)*
- logical [up\\_iter](#)  
*whether the search direction is upwards*
- logical [bisection\\_iter](#)  
*whether to do the bisection search part (after the interval is found)*
- real(dp) [up\\_bnd\\_iter](#)  
*upper bound for the current bisection step*
- real(dp) [low\\_bnd\\_iter](#)  
*lower bound for the current bisection step*
- real(dp), dimension(:), allocatable [l1\\_runoff\\_e](#)

- *runoff energy at L1 level*
- real(dp), dimension(:), allocatable [l1\\_acc\\_ssrd](#)  
*accumulated shortwave radiation at L1 level*
- real(dp), dimension(:), allocatable [l1\\_acc\\_strd](#)  
*accumulated longwave radiation at L1 level*
- real(dp), dimension(:), allocatable [l1\\_acc\\_temp](#)  
*accumulated air temperature radiation at L1 level*
- integer(i4) [ts\\_cnt](#)  
*sub time-step counter for accumulation of meteo*
- integer(i4) [s11](#)  
*starting index for current L11 domain*
- integer(i4) [e11](#)  
*ending index for current L11 domain*
- real(dp), dimension(:, :), allocatable [netnode\\_e\\_in](#)  
*Total energy inputs at t-1 and t.*
- real(dp), dimension(:, :), allocatable [netnode\\_e\\_r](#)  
*energy leaving at t-1 and t*
- real(dp), dimension(:), allocatable [netnode\\_e\\_mod](#)  
*Simulated routed energy.*
- real(dp), dimension(:), allocatable [netnode\\_e\\_out](#)  
*total energy source from cell in L11*
- real(dp), dimension(:), allocatable [l11\\_srad\\_net](#)  
*net short wave radiation at L11*
- real(dp), dimension(:), allocatable [l11\\_lrad\\_in](#)  
*incoming long wave radiation at L11*
- real(dp), dimension(:), allocatable [l11\\_air\\_temp](#)  
*air temp at L11*
- real(dp), dimension(:), allocatable [river\\_temp](#)  
*resulting river temp at L11 in [deg C]*

### 18.15.1 Detailed Description

This is a container to define the river temperature routing in the current time step.

This class provides all procedures to rout river tperature through the river network.

#### Warning

This feature is still experimental! This first version doesn't provide ice covering, which means, that the river temperature can drop below 0 [deg C] in winter.

Definition at line 23 of file mo\_mrm\_riv\_temp\_class.f90.

### 18.15.2 Member Function/Subroutine Documentation

#### 18.15.2.1 acc\_source\_e()

procedure mo\_mrm\_riv\_temp\_class::riv\_temp\_type::acc\_source\_e  
accumulate energy sources of [riv\\_temp\\_type](#)

## Parameters

|    |                              |  |
|----|------------------------------|--|
| in | <i>time</i>                  | current decimal Julian day   |
| in | <i>ntimesteps_day</i>        | number of time intervals per day, transformed in dp                |
| in | <i>fsealed_area_fraction</i> | sealed area fraction [1]   |
| in | <i>fast_interflow</i>        | $q_0$ Fast runoff component [mm TS-1]                              |
| in | <i>slow_interflow</i>        | $q_1$ Slow runoff component [mm TS-1]                              |
| in | <i>baseflow</i>              | $q_2$ Baseflow [mm TS-1]   |
| in | <i>direct_runoff</i>         | $q_D$ Direct runoff from impervious areas [mm TS-1]                |
| in | <i>temp_air</i>              | air temperature [K]  |
| in | <i>mean_temp_air</i>         | annual mean air temperature [K]                                    |
| in | <i>ssrd_day</i>              | Daily mean short radiation   |
| in | <i>strd_day</i>              | Daily mean longwave radiation                                      |
| in | <i>read_meteo_weights</i>    | flag whether weights for tavg and pet have read and should be used |
| in | <i>temp_weights</i>          | multiplicative weights for temperature (deg K)                     |
| in | <i>fday_temp</i>             | [-] day factor mean temp   |
| in | <i>fnight_temp</i>           | [-] night factor mean temp   |
| in | <i>fday_ssrd</i>             | Daytime fraction of ssrd   |
| in | <i>fnight_ssrd</i>           | Nighttime fraction of ssrd   |
| in | <i>fday_strd</i>             | Daytime fraction of strd   |
| in | <i>fnight_strd</i>           | Nighttime fraction of strd   |

## See also

[mo\\_mrm\\_riv\\_temp\\_class::acc\\_source\\_e](#)

Definition at line 82 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.2 alloc\_lateral()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::alloc_lateral
allocate lateral temp components of riv\_temp\_type class for current domain
```

## Parameters

|    |               |  |
|----|---------------|--|
| in | <i>ncells</i> | number of cells for the current domain |
|----|---------------|--|

## See also

[mo\\_mrm\\_riv\\_temp\\_class::alloc\\_lateral](#)

Definition at line 108 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.3 config()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::config
configure the riv\_temp\_type class from the mhm namelist
```



## Parameters

|    |                            |                             |
|----|----------------------------|-----------------------------|
| in | <i>file_namelist</i>       | mhm namelist file           |
| in | <i>file_namelist_param</i> | mhm parameter namelist file |

## See also

[mo\\_mrm\\_riv\\_temp\\_class::config](#)

Definition at line 72 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.4 dealloc\_lateral()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::dealloc_lateral
```

deallocate lateral temp components of [riv\\_temp\\_type](#)

## See also

[mo\\_mrm\\_riv\\_temp\\_class::dealloc\\_lateral](#)

Definition at line 110 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.5 finalize\_source\_e()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::finalize_source_e
```

finalize energy sources of [riv\\_temp\\_type](#)

## Parameters

|    |                     |  |
|----|---------------------|--|
| in | <i>efecarea</i>     | effective area in [km2] at Level 1                                       |
| in | <i>l1_l11_id</i>    | L11 lds mapped on L1   |
| in | <i>l11_areacell</i> | effective area in [km2] at Level 11                                      |
| in | <i>l11_l1_id</i>    | L1 lds mapped on L11   |
| in | <i>timestep</i>     | simulation timestep in [h]   |
| in | <i>map_flag</i>     | Flag indicating whether routing resolution is higher than hydrologic one |

## See also

[mo\\_mrm\\_riv\\_temp\\_class::finalize\\_source\\_e](#)

Definition at line 84 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.6 get\_e\_io()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::get_e_io
```

get complete energy source of [riv\\_temp\\_type](#) at given cell

**Returns**

energy IO

**Parameters**

|    |                 |   |
|----|-----------------|---|
| in | <i>riv_temp</i> | given river temperature in K to calculate heat fluxes |
| in | <i>cell</i>     | cell index in the current domain                      |

**See also**

[mo\\_mrm\\_riv\\_temp\\_class::get\\_e\\_io](#)

Definition at line 94 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.7 get\_lat\_heat()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::get_lat_heat
```

latent heat flux of [riv\\_temp\\_type](#)

**Returns**

latent heat flux

**Parameters**

|    |                 |                          |
|----|-----------------|--------------------------|
| in | <i>air_temp</i> | air temperature in deg C |
| in | <i>netrad</i>   | net radiation in W * m-2 |

**See also**

[mo\\_mrm\\_riv\\_temp\\_class::get\\_lat\\_heat](#)

Definition at line 90 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.8 get\_lrad\_out()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::get_lrad_out
```

get outgoing longwave radiation of [riv\\_temp\\_type](#)

**Returns**

outgoing longwave radiation

**Parameters**

|    |                 |                        |
|----|-----------------|------------------------|
| in | <i>riv_temp</i> | river temperature in K |
|----|-----------------|------------------------|

See also

[mo\\_mrm\\_riv\\_temp\\_class::get\\_lrad\\_out](#)

Definition at line 88 of file mo\_mrm\_riv\_temp\_class.f90.

### 18.15.2.9 get\_sens\_heat()

procedure mo\_mrm\_riv\_temp\_class::riv\_temp\_type::get\_sens\_heat  
sensible heat flux of [riv\\_temp\\_type](#)

Returns

sensible heat flux

Parameters

|    |                 |                              |
|----|-----------------|------------------------------|
| in | <i>air_temp</i> | air temperature in [deg C]   |
| in | <i>riv_temp</i> | river temperature in [deg C] |

See also

[mo\\_mrm\\_riv\\_temp\\_class::get\\_sens\\_heat](#)

Definition at line 92 of file mo\_mrm\_riv\_temp\_class.f90.

### 18.15.2.10 init()

procedure mo\_mrm\_riv\_temp\_class::riv\_temp\_type::init  
initialize the [riv\\_temp\\_type](#) class for the current domain

Parameters

|    |               |  |
|----|---------------|--|
| in | <i>ncells</i> | number of cells for the current domain |
|----|---------------|--|

See also

[mo\\_mrm\\_riv\\_temp\\_class::init](#)

Definition at line 74 of file mo\_mrm\_riv\_temp\_class.f90.

### 18.15.2.11 init\_area()

procedure mo\_mrm\_riv\_temp\_class::riv\_temp\_type::init\_area  
initialize the river area of [riv\\_temp\\_type](#) class for the current domain

Parameters

|    |                |           |
|----|----------------|-----------|
| in | <i>idomain</i> | Domain ID |
|----|----------------|-----------|

## Parameters

|    |                    |  |
|----|--------------------|--|
| in | <i>l11_netperm</i> | L11 routing order  |
| in | <i>l11_fromn</i>   | L11 source grid cell order                                   |
| in | <i>l11_length</i>  | L11 link length  |
| in | <i>nlinks</i>      | number of L11 links in the current domain                    |
| in | <i>ncells</i>      | number of L11 cells of the current domain                    |
| in | <i>ncols</i>       | Number of columns  |
| in | <i>nrows</i>       | Number of rows   |
| in | <i>l11_mask</i>    | the mask for valid cells in the original grid (nrows, ncols) |

## See also

[mo\\_mrm\\_riv\\_temp\\_class::init\\_area](#)

Definition at line 76 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.12 init\_iter()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::init_iter
```

initialize iterative solver of [riv\\_temp\\_type](#)

## See also

[mo\\_mrm\\_riv\\_temp\\_class::init\\_iter](#)

Definition at line 100 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.2.13 init\_riv\_temp()**

```
procedure mo_mrm_riv_temp_class::riv_temp_type::init_riv_temp
```

initialize the river temperature of [riv\\_temp\\_type](#) class for the current domain

## Parameters

|    |                           |  |
|----|---------------------------|--|
| in | <i>time</i>               | current decimal Julian day   |
| in | <i>ntimesteps_day</i>     | number of time intervals per day, transformed in dp                      |
| in | <i>temp_air</i>           | air temperature [K]  |
| in | <i>read_meteo_weights</i> | flag whether weights for tavg and pet have read and should be used       |
| in | <i>temp_weights</i>       | multiplicative weights for temperature (deg K)                           |
| in | <i>fday_temp</i>          | [-] day factor mean temp   |
| in | <i>fnight_temp</i>        | [-] night factor mean temp   |
| in | <i>efecarea</i>           | effective area in [km2] at Level 1                                       |
| in | <i>l1_l11_id</i>          | L11 lds mapped on L1   |
| in | <i>l11_areacell</i>       | effective area in [km2] at Level 11                                      |
| in | <i>l1_l1_id</i>           | L1 lds mapped on L11   |
| in | <i>map_flag</i>           | Flag indicating whether routing resolution is higher than hydrologic one |

See also

[mo\\_mrm\\_riv\\_temp\\_class::init\\_riv\\_temp](#)

Definition at line 78 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.2.14 l11\_routing\_e()

```
procedure mo_mrm_riv_temp_class::riv_temp_type::l11_routing_e
```

execute the temperature routing of [riv\\_temp\\_type](#)

Parameters

|    |                        |  |
|----|------------------------|--|
| in | <i>nlinks</i>          | number of stream segment (reaches)   |
| in | <i>netperm</i>         | routing order of a given domain (permutation)                              |
| in | <i>netlink_fromn</i>   | from node  |
| in | <i>netlink_ton</i>     | to node  |
| in | <i>netlink_c1</i>      | routing parameter C1 ([7] p. 25-41)  |
| in | <i>netlink_c2</i>      | routing parameters C2 (id)   |
| in | <i>ninflowgauges</i>   | [-] number of inflow points  |
| in | <i>inflowheadwater</i> | [-] if to consider headwater cells of inflow gauge                         |
| in | <i>inflowodelist</i>   | [-] L11 ID of inflow points  |
| in | <i>l11_qtr</i>         | [m3 s-1] Transformed outflow leaving node l at current timestep(Muskingum) |
| in | <i>l11_qmod</i>        | [m3 s-1] Simulated routed discharge  |

See also

[mo\\_mrm\\_riv\\_temp\\_class::l11\\_routing\\_e](#)

Definition at line 96 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.2.15 next\_iter()

```
procedure mo_mrm_riv_temp_class::riv_temp_type::next_iter
```

execute next iteration with iterative solver of [riv\\_temp\\_type](#)

Parameters

|         |               |                                       |
|---------|---------------|---------------------------------------|
| in, out | <i>t_est</i>  | estimated river temperature           |
| in      | <i>t_rout</i> | calculated (routed) river temperature |

See also

[mo\\_mrm\\_riv\\_temp\\_class::next\\_iter](#)

Definition at line 102 of file mo\_mrm\_riv\_temp\_class.f90.

### 18.15.2.16 reset\_timestep()

```
procedure mo_mrm_riv_temp_class::riv_temp_type::reset_timestep
```

reset [riv\\_temp\\_type](#) class for next timestep

See also

[mo\\_mrm\\_riv\\_temp\\_class::reset\\_timestep](#)

Definition at line 106 of file `mo_mrm_riv_temp_class.f90`.

## 18.15.3 Member Data Documentation

### 18.15.3.1 active

```
logical mo_mrm_riv_temp_class::riv_temp_type::active = .false.
```

state if this process is active

Definition at line 25 of file `mo_mrm_riv_temp_class.f90`.

### 18.15.3.2 albedo\_water

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::albedo_water
```

albedo of open water

Definition at line 35 of file `mo_mrm_riv_temp_class.f90`.

### 18.15.3.3 bisect\_iter

```
logical mo_mrm_riv_temp_class::riv_temp_type::bisect_iter
```

whether to do the bisection search part (after the interval is found)

Definition at line 48 of file `mo_mrm_riv_temp_class.f90`.

### 18.15.3.4 case

```
integer(i4) mo_mrm_riv_temp_class::riv_temp_type::case = 0_i4
```

the selected process-case option

Definition at line 26 of file `mo_mrm_riv_temp_class.f90`.

### 18.15.3.5 delta\_iter

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::delta_iter
```

input: convergence criteria for iterative solver

Definition at line 44 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.6 delta\_t

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::delta_t = 0.1_dp
```

cutoff value for temperature

Definition at line 41 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.7 dir\_riv\_widths

```
character(256), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::dir_riv_widths
```

Directory where river widths are stored.

Definition at line 29 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.8 e11

```
integer(i4) mo_mrm_riv_temp_class::riv_temp_type::e11
```

ending index for current L11 domain

Definition at line 59 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.9 emissivity\_water

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::emissivity_water
```

emissivity of water

Definition at line 38 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.10 first\_iter

```
logical mo_mrm_riv_temp_class::riv_temp_type::first_iter
```

whether it is at the first iteration (to determine search direction)

Definition at line 46 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.11 l11\_air\_temp

```
real(dp), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::l11_air_temp
```

air temp at L11

Definition at line 66 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.12 l11\_lrad\_in**

```
real(dp), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::l11_lrad_in
```

incoming long wave radiation at L11

Definition at line 65 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.13 l11\_riv\_areas**

```
real(dp), dimension(:), allocatable, public mo_mrm_riv_temp_class::riv_temp_type::l11_riv_↵  
areas
```

river area in L11

Definition at line 33 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.14 l11\_riv\_widths**

```
real(dp), dimension(:), allocatable, public mo_mrm_riv_temp_class::riv_temp_type::l11_riv_↵  
widths
```

river widths in L11

Definition at line 32 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.15 l11\_srad\_net**

```
real(dp), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::l11_srad_net
```

net short wave radiation at L11

Definition at line 64 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.16 l1\_acc\_ssrđ**

```
real(dp), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::l1_acc_ssrđ
```

accumulated shortwave radiation at L1 level

Definition at line 53 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.17 l1\_acc\_strđ**

```
real(dp), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::l1_acc_strđ
```

accumulated longwave radiation at L1 level

Definition at line 54 of file mo\_mrm\_riv\_temp\_class.f90.



**18.15.3.18 l1\_acc\_temp**

real(dp), dimension(:), allocatable mo\_mrm\_riv\_temp\_class::riv\_temp\_type::l1\_acc\_temp

accumulated air temperature radiation at L1 level

Definition at line 55 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.19 l1\_runoff\_e**

real(dp), dimension(:), allocatable mo\_mrm\_riv\_temp\_class::riv\_temp\_type::l1\_runoff\_e

runoff energy at L1 level

Definition at line 52 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.20 low\_bnd\_iter**

real(dp) mo\_mrm\_riv\_temp\_class::riv\_temp\_type::low\_bnd\_iter

lower bound for the current bisection step

Definition at line 50 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.21 max\_iter**

integer(i4) mo\_mrm\_riv\_temp\_class::riv\_temp\_type::max\_iter

input: maximal number of iterations done

Definition at line 43 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.22 netnode\_e\_in**

real(dp), dimension(:, :), allocatable mo\_mrm\_riv\_temp\_class::riv\_temp\_type::netnode\_e\_in

Total energy inputs at t-1 and t.

Definition at line 60 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.23 netnode\_e\_mod**

real(dp), dimension(:), allocatable mo\_mrm\_riv\_temp\_class::riv\_temp\_type::netnode\_e\_mod

Simulated routed energy.

Definition at line 62 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.24 netnode\_e\_out**

real(dp), dimension(:), allocatable mo\_mrm\_riv\_temp\_class::riv\_temp\_type::netnode\_e\_out

total energy source from cell in L11

Definition at line 63 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.25 netnode\_e\_r

```
real(dp), dimension(:, :), allocatable mo_mrm_riv_temp_class::riv_temp_type::netnode_e_r
```

energy leaving at t-1 and t

Definition at line 61 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.26 nml\_name

```
character(256) mo_mrm_riv_temp_class::riv_temp_type::nml_name = 'config_riv_temp'
```

namelist name in mhm namelist

Definition at line 27 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.27 pt\_a\_water

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::pt_a_water
```

priestley taylor alpha parameter for PET on open water

Definition at line 36 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.28 riv\_widths\_file

```
character(256) mo_mrm_riv_temp_class::riv_temp_type::riv_widths_file
```

file name for river widths

Definition at line 30 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.29 riv\_widths\_name

```
character(256) mo_mrm_riv_temp_class::riv_temp_type::riv_widths_name
```

variable name for river widths

Definition at line 31 of file mo\_mrm\_riv\_temp\_class.f90.

#### 18.15.3.30 river\_temp

```
real(dp), dimension(:), allocatable mo_mrm_riv_temp_class::riv_temp_type::river_temp
```

resulting river temp at L11 in [deg C]

Definition at line 68 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.31 s11**

```
integer(i4) mo_mrm_riv_temp_class::riv_temp_type::s11
```

starting index for current L11 domain

Definition at line 58 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.32 step\_iter**

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::step_iter
```

input: step-size for linear search

Definition at line 45 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.33 ts\_cnt**

```
integer(i4) mo_mrm_riv_temp_class::riv_temp_type::ts_cnt
```

sub time-step counter for accumulation of meteo

Definition at line 56 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.34 turb\_heat\_ex\_coeff**

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::turb_heat_ex_coeff
```

lateral heat exchange coefficient water <-> air

Definition at line 39 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.35 up\_bnd\_iter**

```
real(dp) mo_mrm_riv_temp_class::riv_temp_type::up_bnd_iter
```

upper bound for the current bisection step

Definition at line 49 of file mo\_mrm\_riv\_temp\_class.f90.

**18.15.3.36 up\_iter**

```
logical mo_mrm_riv_temp_class::riv_temp_type::up_iter
```

whether the search direction is upwards

Definition at line 47 of file mo\_mrm\_riv\_temp\_class.f90.

The documentation for this module was generated from the following file:

- [src/mRM/mo\\_mrm\\_riv\\_temp\\_class.f90](#)

## 18.16 mo\_mpr\_global\_variables::soiltype Type Reference

### Private Attributes

- integer(i4), dimension(:), allocatable [id](#)
- integer(i4), dimension(:), allocatable [nhorizons](#)
- integer(i4), dimension(:), allocatable [is\\_present](#)
- real(dp), dimension(:, :), allocatable [ud](#)
- real(dp), dimension(:, :), allocatable [ld](#)
- real(dp), dimension(:, :), allocatable [clay](#)
- real(dp), dimension(:, :), allocatable [sand](#)
- real(dp), dimension(:, :), allocatable [dbm](#)
- real(dp), dimension(:, :), allocatable [depth](#)
- real(dp), dimension(:), allocatable [rzdepth](#)
- real(dp), dimension(:, :, :), allocatable [wd](#)
- integer(i4), dimension(:), allocatable [ntillhorizons](#)
- real(dp), dimension(:, :, :), allocatable [thetas\\_till](#)
- real(dp), dimension(:, :), allocatable [thetas](#)
- real(dp), dimension(:, :, :), allocatable [db](#)
- real(dp), dimension(:, :, :), allocatable [thetafc\\_till](#)
- real(dp), dimension(:, :), allocatable [thetafc](#)
- real(dp), dimension(:, :, :), allocatable [thetapw\\_till](#)
- real(dp), dimension(:, :), allocatable [thetapw](#)
- real(dp), dimension(:, :, :), allocatable [ks](#)

### 18.16.1 Detailed Description

Definition at line 34 of file `mo_mpr_global_variables.f90`.

### 18.16.2 Member Data Documentation

#### 18.16.2.1 clay

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::clay [private]
```

Definition at line 45 of file `mo_mpr_global_variables.f90`.

#### 18.16.2.2 db

```
real(dp), dimension(:, :, :), allocatable mo_mpr_global_variables::soiltype::db [private]
```

Definition at line 59 of file `mo_mpr_global_variables.f90`.

#### 18.16.2.3 dbm

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::dbm [private]
```

Definition at line 47 of file `mo_mpr_global_variables.f90`.

#### 18.16.2.4 depth

real(dp), dimension(:, :), allocatable mo\_mpr\_global\_variables::soiltype::depth [private]

Definition at line 48 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.5 id

integer(i4), dimension(:), allocatable mo\_mpr\_global\_variables::soiltype::id [private]

Definition at line 39 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.6 is\_present

integer(i4), dimension(:), allocatable mo\_mpr\_global\_variables::soiltype::is\_present [private]

Definition at line 41 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.7 ks

real(dp), dimension(:, :, :), allocatable mo\_mpr\_global\_variables::soiltype::ks [private]

Definition at line 67 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.8 ld

real(dp), dimension(:, :), allocatable mo\_mpr\_global\_variables::soiltype::ld [private]

Definition at line 44 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.9 nhorizons

integer(i4), dimension(:), allocatable mo\_mpr\_global\_variables::soiltype::nhorizons [private]

Definition at line 40 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.10 ntillhorizons

integer(i4), dimension(:), allocatable mo\_mpr\_global\_variables::soiltype::ntillhorizons [private]

Definition at line 52 of file mo\_mpr\_global\_variables.f90.

#### 18.16.2.11 rzdepth

real(dp), dimension(:), allocatable mo\_mpr\_global\_variables::soiltype::rzdepth [private]

Definition at line 49 of file mo\_mpr\_global\_variables.f90.

**18.16.2.12 sand**

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::sand [private]
```

Definition at line 46 of file mo\_mpr\_global\_variables.f90.

**18.16.2.13 thetafc**

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::thetafc [private]
```

Definition at line 63 of file mo\_mpr\_global\_variables.f90.

**18.16.2.14 thetafc\_till**

```
real(dp), dimension(:, :, :), allocatable mo_mpr_global_variables::soiltype::thetafc_till  
[private]
```

Definition at line 61 of file mo\_mpr\_global\_variables.f90.

**18.16.2.15 thetapw**

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::thetapw [private]
```

Definition at line 66 of file mo\_mpr\_global\_variables.f90.

**18.16.2.16 thetapw\_till**

```
real(dp), dimension(:, :, :), allocatable mo_mpr_global_variables::soiltype::thetapw_till  
[private]
```

Definition at line 64 of file mo\_mpr\_global\_variables.f90.

**18.16.2.17 thetas**

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::thetas [private]
```

Definition at line 57 of file mo\_mpr\_global\_variables.f90.

**18.16.2.18 thetas\_till**

```
real(dp), dimension(:, :, :), allocatable mo_mpr_global_variables::soiltype::thetas_till [private]
```

Definition at line 55 of file mo\_mpr\_global\_variables.f90.

### 18.16.2.19 ud

```
real(dp), dimension(:, :), allocatable mo_mpr_global_variables::soiltype::ud [private]
```

Definition at line 43 of file mo\_mpr\_global\_variables.f90.

### 18.16.2.20 wd

```
real(dp), dimension(:, :, :), allocatable mo_mpr_global_variables::soiltype::wd [private]
```

Definition at line 50 of file mo\_mpr\_global\_variables.f90.

The documentation for this type was generated from the following file:

- [src/MPR/mo\\_mpr\\_global\\_variables.f90](#)

## 18.17 mo\_spatial\_agg\_disagg\_forcing::spatial\_aggregation Interface Reference

Spatial aggregation of meteorological variables.

### Public Member Functions

- subroutine [spatial\\_aggregation\\_3d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [spatial\\_aggregation\\_4d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)

### 18.17.1 Detailed Description

Spatial aggregation of meteorological variables.

Aggregate (or upscale) the given level-2 meteorological data to the required level-1 spatial resolution for the mHM run.

#### Authors

Rohini Kumar

#### Date

Jan 2013

Definition at line 49 of file mo\_spatial\_agg\_disagg\_forcing.f90.

### 18.17.2 Member Function/Subroutine Documentation

#### 18.17.2.1 spatial\_aggregation\_3d()

```
subroutine mo_spatial_agg_disagg_forcing::spatial_aggregation::spatial_aggregation_3d (  
    real(dp), dimension(:, :, :), intent(in) data2,  
    real(dp), intent(in) cellsize2,
```

```

real(dp), intent(in) cellsize1,
logical, dimension(:, :), intent(in) mask1,
logical, dimension(:, :), intent(in) mask2,
real(dp), dimension(:, :, :), intent(out), allocatable data1 )

```

Definition at line 96 of file `mo_spatial_agg_disagg_forcing.f90`.

References `mo_common_constants::nodata_dp`.

### 18.17.2.2 `spatial_aggregation_4d()`

```

subroutine mo_spatial_agg_disagg_forcing::spatial_aggregation::spatial_aggregation_4d (
    real(dp), dimension(:, :, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
    logical, dimension(:, :), intent(in) mask2,
    real(dp), dimension(:, :, :, :), intent(out), allocatable data1 )

```

Definition at line 204 of file `mo_spatial_agg_disagg_forcing.f90`.

References `mo_common_constants::nodata_dp`.

The documentation for this interface was generated from the following file:

- [src/mHM/mo\\_spatial\\_agg\\_disagg\\_forcing.f90](#)

## 18.18 `mo_spatial_agg_disagg_forcing::spatial_disaggregation` Interface Reference

Spatial disaggregation of meteorological variables.

### Public Member Functions

- subroutine [spatial\\_disaggregation\\_3d](#) (`data2`, `cellsize2`, `cellsize1`, `mask1`, `mask2`, `data1`)
- subroutine [spatial\\_disaggregation\\_4d](#) (`data2`, `cellsize2`, `cellsize1`, `mask1`, `mask2`, `data1`)

### 18.18.1 Detailed Description

Spatial disaggregation of meteorological variables.

Disaggregate (or downscale) the given level-2 meteorological data to the required level-1 spatial resolution for the mHM run.

#### Parameters

|     |   |                    |
|-----|---|--------------------|
| in  | <i>real(dp), dimension(:, :, :)</i> :: <i>data2</i> | Level-2 data       |
| in  | <i>real(dp)</i> :: <i>cellsize2</i>                 | Level-2 resolution |
| in  | <i>real(dp)</i> :: <i>cellsize1</i>                 | Level-1 resolution |
| in  | <i>logical, dimension(:, :)</i> :: <i>mask1</i>     | Level-1 mask       |
| in  | <i>logical, dimension(:, :)</i> :: <i>mask2</i>     | Level-2 mask       |
| out | <i>real(dp), dimension(:, :, :)</i> :: <i>data1</i> | Level-1 data       |



**Authors**

Rohini Kumar

**Date**

Jan 2013

Definition at line 84 of file mo\_spatial\_agg\_disagg\_forcing.f90.

**18.18.2 Member Function/Subroutine Documentation****18.18.2.1 spatial\_disaggregation\_3d()**

```

subroutine mo_spatial_agg_disagg_forcing::spatial_disaggregation::spatial_disaggregation_3d (
    real(dp), dimension(:, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
    logical, dimension(:, :), intent(in) mask2,
    real(dp), dimension(:, :, :), intent(out), allocatable data1 )

```

Definition at line 315 of file mo\_spatial\_agg\_disagg\_forcing.f90.

References mo\_common\_constants::nodata\_dp.

**18.18.2.2 spatial\_disaggregation\_4d()**

```

subroutine mo_spatial_agg_disagg_forcing::spatial_disaggregation::spatial_disaggregation_4d (
    real(dp), dimension(:, :, :, :), intent(in) data2,
    real(dp), intent(in) cellsize2,
    real(dp), intent(in) cellsize1,
    logical, dimension(:, :), intent(in) mask1,
    logical, dimension(:, :), intent(in) mask2,
    real(dp), dimension(:, :, :, :), intent(out), allocatable data1 )

```

Definition at line 381 of file mo\_spatial\_agg\_disagg\_forcing.f90.

References mo\_common\_constants::nodata\_dp.

The documentation for this interface was generated from the following file:

- [src/mHM/mo\\_spatial\\_agg\\_disagg\\_forcing.f90](#)

**18.19 mo\_mpr\_restart::unpack\_field\_and\_write Interface Reference**

TODO: add description.

**Private Member Functions**

- subroutine [unpack\\_field\\_and\\_write\\_1d\\_i4](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_1d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_2d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_3d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)

### 18.19.1 Detailed Description

TODO: add description.

TODO: add description

#### Parameters

|         |  |                                    |
|---------|--|------------------------------------|
| in, out | <i>type(NcDataset) :: nc</i>                       | NcDataset to add variable to       |
| in      | <i>character(*) :: var_name</i>                    | variable name                      |
| in      | <i>type(NcDimension), dimension(:) :: var_dims</i> | vector of Variable dimensions      |
| in      | <i>integer(i4) :: fill_value</i>                   | fill value used for missing values |
| in      | <i>integer(i4), dimension(:) :: data</i>           | packed data to be set to variable  |
| in      | <i>logical, dimension(:, :) :: mask</i>            | mask used for unpacking            |
| in      | <i>character(*), optional :: var_long_name</i>     | variable long name attribute       |

#### Authors

Robert Schweppe

#### Date

Jun 2018

Definition at line 59 of file mo\_mpr\_restart.f90.

### 18.19.2 Member Function/Subroutine Documentation

#### 18.19.2.1 `unpack_field_and_write_1d_dp()`

```
subroutine mo_mpr_restart::unpack_field_and_write::unpack_field_and_write_1d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 413 of file mo\_mpr\_restart.f90.

#### 18.19.2.2 `unpack_field_and_write_1d_i4()`

```
subroutine mo_mpr_restart::unpack_field_and_write::unpack_field_and_write_1d_i4 (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    integer(i4), intent(in) fill_value,
    integer(i4), dimension(:), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 368 of file mo\_mpr\_restart.f90.

### 18.19.2.3 unpack\_field\_and\_write\_2d\_dp()

```
subroutine mo_mpr_restart::unpack_field_and_write::unpack_field_and_write_2d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:, :), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 458 of file mo\_mpr\_restart.f90.

### 18.19.2.4 unpack\_field\_and\_write\_3d\_dp()

```
subroutine mo_mpr_restart::unpack_field_and_write::unpack_field_and_write_3d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:, :, :), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 515 of file mo\_mpr\_restart.f90.

The documentation for this interface was generated from the following file:

- [src/MPR/mo\\_mpr\\_restart.f90](#)

## 18.20 mo\_restart::unpack\_field\_and\_write Interface Reference

TODO: add description.

### Private Member Functions

- subroutine [unpack\\_field\\_and\\_write\\_1d\\_i4](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_1d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_2d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [unpack\\_field\\_and\\_write\\_3d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)

### 18.20.1 Detailed Description

TODO: add description.

TODO: add description

#### Parameters

|         |                              |                              |
|---------|------------------------------|------------------------------|
| in, out | <i>type(NcDataset) :: nc</i> | NcDataset to add variable to |
|---------|------------------------------|------------------------------|

## Parameters

|    |  |                                    |
|----|--|------------------------------------|
| in | <i>character(*) :: var_name</i>                    | variable name                      |
| in | <i>type(NcDimension), dimension(:) :: var_dims</i> | vector of Variable dimensions      |
| in | <i>integer(i4) :: fill_value</i>                   | fill value used for missing values |
| in | <i>integer(i4), dimension(:) :: data</i>           | packed data to be set to variable  |
| in | <i>logical, dimension(:, :) :: mask</i>            | mask used for unpacking            |
| in | <i>character(*), optional :: var_long_name</i>     | variable long name attribute       |

## Authors

Robert Schweppe

## Date

Jun 2018

Definition at line 60 of file mo\_restart.f90.

## 18.20.2 Member Function/Subroutine Documentation

### 18.20.2.1 unpack\_field\_and\_write\_1d\_dp()

```
subroutine mo_restart::unpack_field_and_write::unpack_field_and_write_1d_dp (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    real(dp), intent(in) fill_value,
    real(dp), dimension(:), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 830 of file mo\_restart.f90.

### 18.20.2.2 unpack\_field\_and\_write\_1d\_i4()

```
subroutine mo_restart::unpack_field_and_write::unpack_field_and_write_1d_i4 (
    type(ncdataset), intent(inout) nc,
    character(*), intent(in) var_name,
    type(ncdimension), dimension(:), intent(in) var_dims,
    integer(i4), intent(in) fill_value,
    integer(i4), dimension(:), intent(in) data,
    logical, dimension(:, :), intent(in) mask,
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 785 of file mo\_restart.f90.

### 18.20.2.3 unpack\_field\_and\_write\_2d\_dp()

```
subroutine mo_restart::unpack_field_and_write::unpack_field_and_write_2d_dp (  
    type(ncdataset), intent(inout) nc,  
    character(*), intent(in) var_name,  
    type(ncdimension), dimension(:), intent(in) var_dims,  
    real(dp), intent(in) fill_value,  
    real(dp), dimension(:, :), intent(in) data,  
    logical, dimension(:, :), intent(in) mask,  
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 875 of file mo\_restart.f90.

### 18.20.2.4 unpack\_field\_and\_write\_3d\_dp()

```
subroutine mo_restart::unpack_field_and_write::unpack_field_and_write_3d_dp (  
    type(ncdataset), intent(inout) nc,  
    character(*), intent(in) var_name,  
    type(ncdimension), dimension(:), intent(in) var_dims,  
    real(dp), intent(in) fill_value,  
    real(dp), dimension(:, :, :), intent(in) data,  
    logical, dimension(:, :, :), intent(in) mask,  
    character(*), intent(in), optional var_long_name ) [private]
```

Definition at line 932 of file mo\_restart.f90.

The documentation for this interface was generated from the following file:

- [src/mHM/mo\\_restart.f90](#)



## Chapter 19

# File Documentation

- 19.1 doc/1-main.dox File Reference**
- 19.2 doc/2-get\_started.dox File Reference**
- 19.3 doc/3-data\_preparation.dox File Reference**
- 19.4 doc/4-visualise\_out.dox File Reference**
- 19.5 doc/5-calibration.dox File Reference**
- 19.6 doc/6-style\_guide.dox File Reference**
- 19.7 doc/7-test\_basin.dox File Reference**
- 19.8 doc/8-protocols\_for\_setup\_new\_mHM\_basin.dox File Reference**
- 19.9 doc/9-mRM.dox File Reference**
- 19.10 doc/DEPENDENCIES.md File Reference**
- 19.11 doc/INSTALL.md File Reference**
- 19.12 doc/mhm\_papers.md File Reference**
- 19.13 doc/RELEASES.md File Reference**
- 19.14 src/common/mo\_check.f90 File Reference**

Input checking routines.

## Modules

- module [mo\\_check](#)

## Functions/Subroutines

- subroutine, public [mo\\_check::check\\_dir](#) (path, text\_, throwError\_, tab\_, text\_length\_)  
*Check if a given directory exists.*

### 19.14.1 Detailed Description

Input checking routines.

This module provides sanity checks for the input data.

#### Authors

Sebastian Mueller

#### Date

Nov 2020

## 19.15 src/common/mo\_common\_constants.f90 File Reference

### Modules

- module [mo\\_common\\_constants](#)  
*Provides constants commonly used by mHM, mRM and MPR.*

### Variables

- real(dp), parameter, public [mo\\_common\\_constants::eps\\_dp](#) = epsilon(1.0\_dp)  
*epsilon(1.0) in double precision*
- real(sp), parameter, public [mo\\_common\\_constants::eps\\_sp](#) = epsilon(1.0\_sp)  
*epsilon(1.0) in single precision*
- integer(i4), parameter, public [mo\\_common\\_constants::nodata\\_i4](#) = -9999\_i4
- real(dp), parameter, public [mo\\_common\\_constants::nodata\\_dp](#) = -9999.\_dp
- real(dp), parameter, public [mo\\_common\\_constants::p1\\_initstatefluxes](#) = 0.00\_dp
- integer(i4), parameter, public [mo\\_common\\_constants::ncolpars](#) = 5\_i4
- integer(i4), parameter, public [mo\\_common\\_constants::maxnodomains](#) = 50\_i4
- integer(i4), parameter, public [mo\\_common\\_constants::maxnlcovers](#) = 50\_i4
- character(64), parameter, public [mo\\_common\\_constants::soilhorizonsvarname](#) = "L1\_SoilHorizons"
- character(64), parameter, public [mo\\_common\\_constants::landcoverperiodsvarname](#) = "L1\_LandCover←  
Periods"
- character(64), parameter, public [mo\\_common\\_constants::laivarname](#) = "L1\_LAITimesteps"

## 19.16 src/common/mo\_common\_datetime\_type.f90 File Reference

### Data Types

- type [mo\\_common\\_datetime\\_type::datetimeinfo](#)



## Modules

- module [mo\\_common\\_datetime\\_type](#)

## Functions/Subroutines

- subroutine [mo\\_common\\_datetime\\_type::datetimeinfo\\_init](#) (this, iDomain)
- subroutine [mo\\_common\\_datetime\\_type::datetimeinfo\\_increment](#) (this)
- subroutine [mo\\_common\\_datetime\\_type::datetimeinfo\\_update\\_lai\\_timestep](#) (this)
- logical function [mo\\_common\\_datetime\\_type::datetimeinfo\\_writeout](#) (this, timeStep\_model\_outputs, tt)

## 19.17 src/common/mo\_common\_file.f90 File Reference

### Modules

- module [mo\\_common\\_file](#)  
*Provides file names and units for mRM.*

### Variables

- character(len= \*), parameter [mo\\_common\\_file::file\\_dem](#) = 'dem.asc'  
*DEM input data file.*
- integer, parameter [mo\\_common\\_file::udem](#) = 53  
*Unit for DEM input data file.*
- integer, parameter [mo\\_common\\_file::ulcoverclass](#) = 61  
*Unit for LCover input data file.*
- character(len= \*), parameter [mo\\_common\\_file::file\\_config](#) = 'ConfigFile.log'  
*file defining mHM's outputs*
- integer, parameter [mo\\_common\\_file::uconfig](#) = 68  
*Unit for file defining mHM's outputs.*

## 19.18 src/common/mo\_common\_functions.f90 File Reference

### Modules

- module [mo\\_common\\_functions](#)  
*Provides small utility functions used by multiple parts of the code (mHM, mRM, MPR)*

### Functions/Subroutines

- logical function, public [mo\\_common\\_functions::in\\_bound](#) (params)  
*TODO: add description.*

## 19.19 src/common/mo\_common\_read\_config.f90 File Reference

### Modules

- module [mo\\_common\\_read\\_config](#)  
*Reading of main model configurations.*

## Functions/Subroutines

- subroutine, public [mo\\_common\\_read\\_config::common\\_read\\_config](#) (file\_namelist, unamelist)  
*Read main configurations commonly used by mHM, mRM and MPR.*
- subroutine, public [mo\\_common\\_read\\_config::set\\_land\\_cover\\_scenes\\_id](#) (sim\_Per, LCyear\_Id)  
*Read main configurations commonly used by mHM, mRM and MPR.*
- subroutine [mo\\_common\\_read\\_config::init\\_domain\\_variable](#) (nDomains, optiData, domainMeta)
- subroutine [mo\\_common\\_read\\_config::init\\_domain\\_variable\\_for\\_master](#) (domainMeta, colMasters, col←Domain)
- subroutine [mo\\_common\\_read\\_config::distributeddomainsroundrobin](#) (nproc, rank, domainMeta)
- subroutine [mo\\_common\\_read\\_config::distribute\\_processes\\_to\\_domains\\_according\\_to\\_role](#) (optiData, rank, domainMeta, colMasters, colDomain)

## 19.20 src/common/mo\_common\_read\_data.f90 File Reference

### Modules

- module [mo\\_common\\_read\\_data](#)  
*TODO: add description.*

### Functions/Subroutines

- subroutine, public [mo\\_common\\_read\\_data::read\\_dem](#)  
*TODO: add description.*
- subroutine, public [mo\\_common\\_read\\_data::read\\_lcover](#)  
*TODO: add description.*

## 19.21 src/common/mo\_common\_restart.f90 File Reference

### Modules

- module [mo\\_common\\_restart](#)  
*TODO: add description.*

### Functions/Subroutines

- subroutine, public [mo\\_common\\_restart::write\\_grid\\_info](#) (grid\_in, level\_name, nc)  
*write restart files for each domain*
- subroutine, public [mo\\_common\\_restart::read\\_grid\\_info](#) (domainID, InFile, level\_name, new\_grid)  
*reads configuration apart from Level 11 configuration from a restart directory*

## 19.22 src/common/mo\_common\_variables.f90 File Reference

### Data Types

- type [mo\\_common\\_variables::period](#)
- type [mo\\_common\\_variables::grid](#)
- type [mo\\_common\\_variables::gridremapper](#)
- type [mo\\_common\\_variables::domain\\_meta](#)

## Modules

- module [mo\\_common\\_variables](#)  
*Provides structures needed by mHM, mRM and/or mpr.*

## Variables

- character(1024), public [mo\\_common\\_variables::project\\_details](#)
- character(1024), public [mo\\_common\\_variables::setup\\_description](#)
- character(1024), public [mo\\_common\\_variables::simulation\\_type](#)
- character(256), public [mo\\_common\\_variables::conventions](#)
- character(1024), public [mo\\_common\\_variables::contact](#)
- character(1024), public [mo\\_common\\_variables::mhm\\_details](#)
- character(1024), public [mo\\_common\\_variables::history](#)
- integer(i4), public [mo\\_common\\_variables::iflag\\_cordinate\\_sys](#)
- real(dp), dimension(:), allocatable, public [mo\\_common\\_variables::resolutionhydrology](#)
- integer(i4), dimension(:), allocatable, public [mo\\_common\\_variables::i0\\_domain](#)
- logical, public [mo\\_common\\_variables::write\\_restart](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::mhmfilerestartout](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::mrmfilerestartout](#)
- character(256), public [mo\\_common\\_variables::dirconfigout](#)
- character(256), public [mo\\_common\\_variables::dircommonfiles](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::dirmorpho](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::dircover](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::dirout](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::filelatlon](#)
- type(grid), dimension(:), allocatable, target, public [mo\\_common\\_variables::level0](#)
- type(grid), dimension(:), allocatable, target, public [mo\\_common\\_variables::level1](#)
- type(gridremapper), dimension(:), allocatable, public [mo\\_common\\_variables::i0\\_i1\\_remap](#)
- real(dp), dimension(:), allocatable, public [mo\\_common\\_variables::i0\\_elev](#)
- integer(i4), dimension(:, :), allocatable, public [mo\\_common\\_variables::i0\\_cover](#)
- type(mpi\_comm) [mo\\_common\\_variables::comm](#)
- type(domain\_meta), public [mo\\_common\\_variables::domainmeta](#)
- integer(i4), public [mo\\_common\\_variables::nuniquei0domains](#)
- integer(i4), public [mo\\_common\\_variables::nlcoverscene](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::lcfilename](#)
- integer(i4), dimension(:), allocatable, public [mo\\_common\\_variables::lc\\_year\\_start](#)
- integer(i4), dimension(:), allocatable, public [mo\\_common\\_variables::lc\\_year\\_end](#)
- integer(i4), parameter, public [mo\\_common\\_variables::nprocesses](#) = 11
- integer(i4), dimension(nprocesses, 3), public [mo\\_common\\_variables::processmatrix](#)
- real(dp), dimension(:, :), allocatable, target, public [mo\\_common\\_variables::global\\_parameters](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_variables::global\\_parameters\\_name](#)
- logical [mo\\_common\\_variables::alma\\_convention](#)

## 19.23 src/common/mo\_grid.f90 File Reference

### Modules

- module [mo\\_grid](#)  
*TODO: add description.*

## Functions/Subroutines

- subroutine, public [mo\\_grid::init\\_lowres\\_level](#) (highres, target\_resolution, lowres, highres\_lowres\_remap)  
*Level-1 variable initialization.*
- subroutine, public [mo\\_grid::set\\_domain\\_indices](#) (grids, indices)  
*TODO: add description.*
- subroutine, public [mo\\_grid::l0\\_grid\\_setup](#) (new\_grid)  
*level 0 variable initialization*
- subroutine, public [mo\\_grid::mapcoordinates](#) (level, y, x)  
*Generate map coordinates.*
- subroutine, public [mo\\_grid::geocoordinates](#) (level, lat, lon)  
*Generate geographic coordinates.*
- subroutine [mo\\_grid::calculate\\_grid\\_properties](#) (nrowsIn, ncolsIn, xllcornerIn, yllcornerIn, cellsizeIn, aiming←  
Resolution, nrowsOut, ncolsOut, xllcornerOut, yllcornerOut, cellsizeOut)  
*Calculates basic grid properties at a required coarser level using information of a given finer level. Calculates basic grid properties at a required coarser level (e.g., L11) using information of a given finer level (e.g., L0). Basic grid properties such as nrows, ncols, xllcorner, yllcorner cellsize are estimated in this routine.*

## 19.24 src/common/mo\_read\_latlon.f90 File Reference

### Modules

- module [mo\\_read\\_latlon](#)  
*reading latitude and longitude coordinates for each domain*

### Functions/Subroutines

- subroutine, public [mo\\_read\\_latlon::read\\_latlon](#) (ii, lon\_var\_name, lat\_var\_name, level\_name, level)  
*reads latitude and longitude coordinates*

## 19.25 src/common/mo\_read\_nc.f90 File Reference

### Modules

- module [mo\\_read\\_nc](#)  
*Reads forcing input data.*

### Functions/Subroutines

- subroutine, public [mo\\_read\\_nc::read\\_nc](#) (folder, nRows, nCols, varName, mask, data, target\_period, lower, upper, nctimestep, fileName, nocheck, maskout)  
*Reads forcing input in NetCDF file format.*
- subroutine, public [mo\\_read\\_nc::read\\_const\\_nc](#) (folder, nRows, nCols, varName, data, fileName)  
*Reads time independent forcing input in NetCDF file format.*
- subroutine, public [mo\\_read\\_nc::read\\_weights\\_nc](#) (folder, nRows, nCols, varName, data, mask, lower, upper, nocheck, maskout, fileName)  
*Reads weights for meteo forcings input in NetCDF file format.*
- subroutine [mo\\_read\\_nc::get\\_time\\_vector\\_and\\_select](#) (var, fname, inctimestep, time\_start, time\_cnt, target←  
\_period)  
*TODO: add description.*

## 19.26 src/common/mo\_read\_spatial\_data.f90 File Reference

### Data Types

- interface [mo\\_read\\_spatial\\_data::read\\_spatial\\_data\\_ascii](#)  
*Reads spatial data files of ASCII format.*

### Modules

- module [mo\\_read\\_spatial\\_data](#)  
*Reads spatial input data.*

### Functions/Subroutines

- subroutine [mo\\_read\\_spatial\\_data::read\\_spatial\\_data\\_ascii\\_dp](#) (filename, fileunit, header\_ncols, header\_↔nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, data, mask)  
*TODO: add description.*
- subroutine [mo\\_read\\_spatial\\_data::read\\_spatial\\_data\\_ascii\\_i4](#) (filename, fileunit, header\_ncols, header\_↔nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, data, mask)  
*TODO: add description.*
- subroutine, public [mo\\_read\\_spatial\\_data::read\\_header\\_ascii](#) (filename, fileunit, header\_ncols, header\_↔nrows, header\_xllcorner, header\_yllcorner, header\_cellsize, header\_nodata)  
*Reads header lines of ASCII files.*

## 19.27 src/common/mo\_read\_timeseries.f90 File Reference

### Modules

- module [mo\\_read\\_timeseries](#)  
*Routines to read files containing timeseries data.*

### Functions/Subroutines

- subroutine, public [mo\\_read\\_timeseries::read\\_timeseries](#) (filename, fileunit, periodStart, periodEnd, optimize, opti\_function, data, mask, nMeasPerDay)  
*Reads time series in ASCII format.*

## 19.28 src/common/mo\_template.f90 File Reference

### Data Types

- interface [mo\\_template::mean](#)  
*The average.*

### Modules

- module [mo\\_template](#)  
*Template for future module developments.*

## Functions/Subroutines

- elemental pure real(dp) function, public `mo_template::circum` (radius)  
*Circumference of a circle.*
- real(dp) function `mo_template::mean_dp` (dat, mask)
- real(sp) function `mo_template::mean_sp` (dat, mask)

## Variables

- real(dp), parameter, public `mo_template::pi_dp` = 3.141592653589793238462643383279502884197\_dp  
*Constant Pi in double precision.*
- real(sp), parameter, public `mo_template::pi_sp` = 3.141592653589793238462643383279502884197\_sp  
*Constant Pi in single precision.*
- integer(i4), parameter `mo_template::itest` = 1

## 19.29 src/common\_mHM\_mRM/mo\_common\_mHM\_mRM\_file.f90 File Reference

### Modules

- module `mo_common_mhm_mrm_file`  
*Provides file names and units for mHM.*

### Variables

- character(len=\*), parameter `mo_common_mhm_mrm_file::file_opti` = 'FinalParam.out'  
*file defining optimization outputs (objective and parameter set)*
- integer, parameter `mo_common_mhm_mrm_file::uopti` = 72  
*Unit for file optimization outputs (objective and parameter set)*
- character(len=\*), parameter `mo_common_mhm_mrm_file::file_opti_nml` = 'FinalParam.nml'  
*file defining optimization outputs in a namelist format (parameter set)*
- integer, parameter `mo_common_mhm_mrm_file::uopti_nml` = 73  
*Unit for file optimization outputs in a namelist format (parameter set)*

## 19.30 src/common\_mHM\_mRM/mo\_common\_mHM\_mRM\_MPI\_tools.f90 File Reference

### Modules

- module `mo_common_mhm_mrm_mpi_tools`  
*tools for MPI communication that are mHM or mRM specific*

### Functions/Subroutines

- subroutine, public `mo_common_mhm_mrm_mpi_tools::distribute_parameterset` (parameterset)
- subroutine, public `mo_common_mhm_mrm_mpi_tools::get_parameterset` (parameterset)

## 19.31 src/common\_mHM\_mRM/mo\_common\_mHM\_mRM\_read\_config.f90 File Reference

### Modules

- module [mo\\_common\\_mhm\\_mrm\\_read\\_config](#)

*Reading of main model configurations.*

### Functions/Subroutines

- subroutine, public [mo\\_common\\_mhm\\_mrm\\_read\\_config::common\\_mhm\\_mrm\\_read\\_config](#) (file\_namelist, unamelist)  
*Read main configurations for common parts.*
- subroutine, public [mo\\_common\\_mhm\\_mrm\\_read\\_config::check\\_optimization\\_settings](#)  
*TODO: add description.*
- subroutine, public [mo\\_common\\_mhm\\_mrm\\_read\\_config::common\\_check\\_resolution](#) (do\_message, allow\_subgrid\_routing)  
*TODO: add description.*
- subroutine [mo\\_common\\_mhm\\_mrm\\_read\\_config::period\\_copy\\_period\\_data](#) (toPeriod, fromPeriod)

## 19.32 src/common\_mHM\_mRM/mo\_common\_mHM\_mRM\_restart.f90 File Reference

### Data Types

- interface [mo\\_common\\_mhm\\_mrm\\_restart::check\\_consistency\\_element](#)

### Modules

- module [mo\\_common\\_mhm\\_mrm\\_restart](#)

*TODO: add description.*

### Functions/Subroutines

- subroutine, public [mo\\_common\\_mhm\\_mrm\\_restart::check\\_dimension\\_consistency](#) (iBasin, nSoilHorizons\_temp, soilHorizonBoundaries\_temp, nLAIs\_temp, LAIBoundaries\_temp, nLandCoverPeriods\_temp, landCoverPeriodBoundaries\_temp)  
*checks dimension configurations read from restart file*
- subroutine [mo\\_common\\_mhm\\_mrm\\_restart::check\\_consistency\\_element\\_dp](#) (item1, item2, name, iBasin)
- subroutine [mo\\_common\\_mhm\\_mrm\\_restart::check\\_consistency\\_element\\_i4](#) (item1, item2, name, iBasin)

## 19.33 src/common\_mHM\_mRM/mo\_common\_mHM\_mRM\_variables.f90 File Reference

### Modules

- module [mo\\_common\\_mhm\\_mrm\\_variables](#)

*Provides structures needed by mHM, mRM and/or mpr.*

## Variables

- integer(i4) [mo\\_common\\_mhm\\_mrm\\_variables::mrm\\_coupling\\_mode](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::timestep](#)
- real(dp), public [mo\\_common\\_mhm\\_mrm\\_variables::c2tstu](#)
- real(dp), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::resolutionrouting](#)
- logical, public [mo\\_common\\_mhm\\_mrm\\_variables::read\\_restart](#)
- logical, public [mo\\_common\\_mhm\\_mrm\\_variables::mrm\\_read\\_river\\_network](#)
- type(period), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::warmper](#)
- type(period), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::evalper](#)
- type(period), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::simper](#)
- type(period), public [mo\\_common\\_mhm\\_mrm\\_variables::readper](#)
- integer(i4), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::warmingdays](#)
- integer(i4), dimension(:, :), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::lcyyearid](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::ntstepday](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::mhmfilerestartin](#)
- character(256), dimension(:), allocatable, public [mo\\_common\\_mhm\\_mrm\\_variables::mrmfilerestartin](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::opti\\_method](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::opti\\_function](#)
- logical, public [mo\\_common\\_mhm\\_mrm\\_variables::optimize](#)
- logical, public [mo\\_common\\_mhm\\_mrm\\_variables::optimize\\_restart](#)
- integer(i8), public [mo\\_common\\_mhm\\_mrm\\_variables::seed](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::iterations](#)
- real(dp), public [mo\\_common\\_mhm\\_mrm\\_variables::dds\\_r](#)
- real(dp), public [mo\\_common\\_mhm\\_mrm\\_variables::sa\\_temp](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::sce\\_ngs](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::sce\\_npg](#)
- integer(i4), public [mo\\_common\\_mhm\\_mrm\\_variables::sce\\_nps](#)
- logical, public [mo\\_common\\_mhm\\_mrm\\_variables::mcmc\\_opti](#)
- integer(i4), parameter, public [mo\\_common\\_mhm\\_mrm\\_variables::nerror\\_model](#) = 2
- real(dp), dimension(nerror\_model), public [mo\\_common\\_mhm\\_mrm\\_variables::mcmc\\_error\\_params](#)

## 19.34 src/common\_mHM\_mRM/mo\_optimization.f90 File Reference

### Modules

- module [mo\\_optimization](#)  
*Wrapper subroutine for optimization against runoff and sm.*

### Functions/Subroutines

- subroutine, public [mo\\_optimization::optimization](#) (eval, objective, dirConfigOut, funcBest, maskpara)  
*Wrapper for optimization.*

## 19.35 src/mHM/mhm\_driver.f90 File Reference

### Functions/Subroutines

- program [mhm\\_driver](#)  
*Distributed precipitation-runoff model mHM.*



## 19.35.1 Function/Subroutine Documentation

### 19.35.1.1 mhm\_driver()

```
program mhm_driver
```

Distributed precipitation-runoff model mHM.

This is the main driver of mHM, which calls one instance of mHM for a multiple domains and a given period.

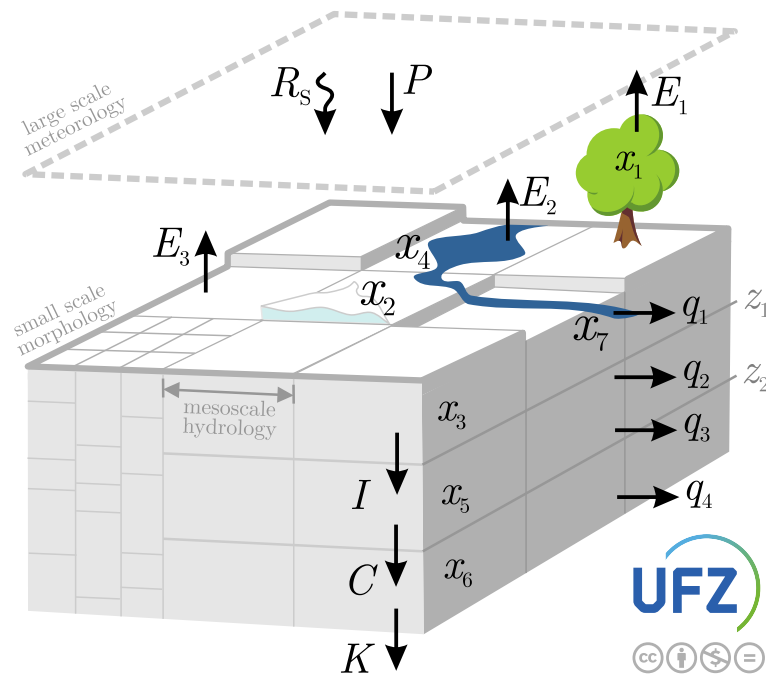


Figure 19.1: Typical mHM cell

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#### Version

5.11.2

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Samaniego L., Kumar R., Attinger S. (2010): Multiscale parameter regionalization of a grid-based hydrologic model at the mesoscale. *Water Resour. Res.*, 46, W05523, doi:10.1029/2008WR007327.

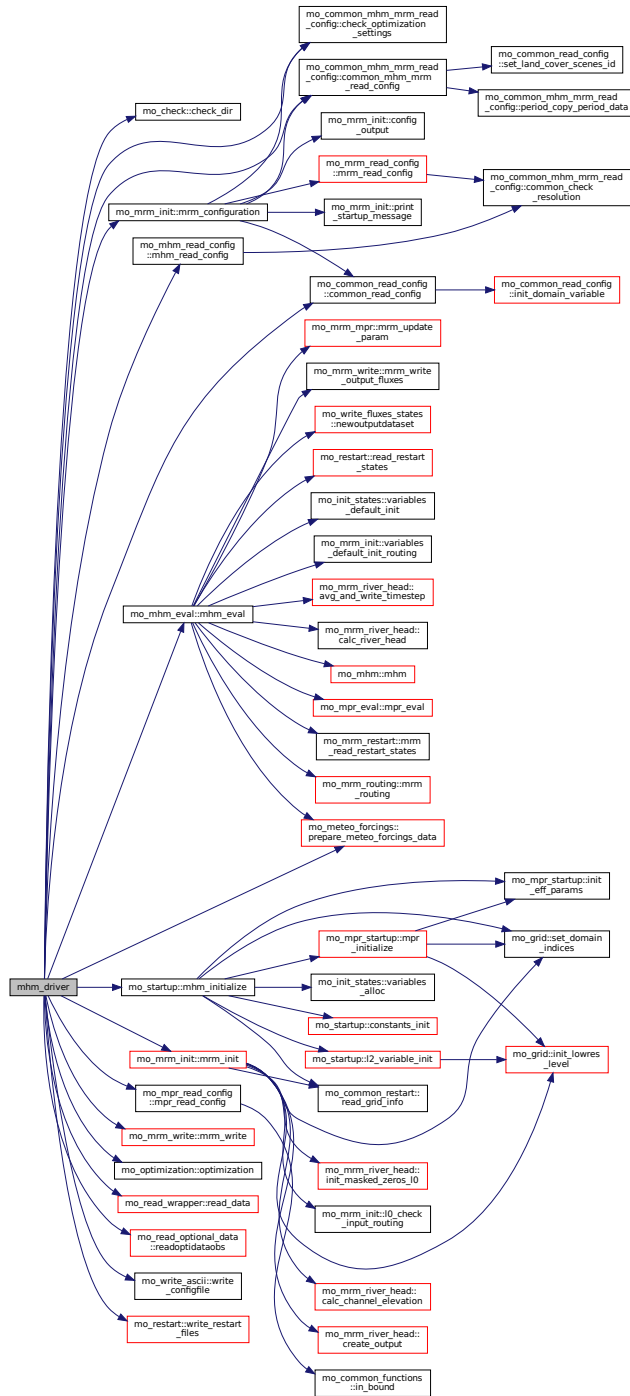
Kumar, R., L. Samaniego, and S. Attinger (2013), Implications of distributed hydrologic model parameterization on water fluxes at multiple scales and locations, *Water Resour. Res.*, 49, doi:10.1029/2012WR012195.

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Definition at line 79 of file `mhm_driver.f90`.

References `mo_check::check_dir()`, `mo_common_mhm_mrm_read_config::check_optimization_settings()`, `mo↔_common_mhm_mrm_read_config::common_mhm_mrm_read_config()`, `mo_common_read_config::common↔_read_config()`, `mo_global_variables::dirprecipitation`, `mo_file::file_main`, `mo_mhm_eval::mhm_eval()`, `mo↔_startup::mhm_initialize()`, `mo_mhm_read_config::mhm_read_config()`, `mo_mpr_read_config::mpr_read_config()`, `mo_mrm_init::mrm_configuration()`, `mo_mrm_init::mrm_init()`, `mo_mrm_write::mrm_write()`, `mo_common_mhm↔_mrm_variables::ntstepday`, `mo_optimization::optimization()`, `mo_meteo_forcings::prepare_meteo_forcings_data()`, `mo_read_wrapper::read_data()`, `mo_read_optional_data::readoptidataobs()`, `mo_file::version`, `mo_file::version↔_date`, `mo_write_ascii::write_configfile()`, `mo_common_variables::write_restart`, and `mo_restart::write_restart_files()`.

Here is the call graph for this function:



## 19.36 src/mHM/mo\_canopy\_interc.f90 File Reference

### Modules

- module [mo\\_canopy\\_interc](#)

*Canopy interception.*

## Functions/Subroutines

- elemental pure subroutine, public `mo_canopy_interc::canopy_interc` (pet, interc\_max, precip, interc, throughfall, evap\_canopy)  
*Canopy interception.*

## 19.37 src/mHM/mo\_file.f90 File Reference

### Modules

- module `mo_file`  
*Provides file names and units for mHM.*

### Variables

- character(len=\*), parameter `mo_file::version` = '5.11.2'  
*Current mHM model version (will be set to 5.11.2)*
- character(len=\*), parameter `mo_file::version_date` = 'Jul 2021'  
*Time of current mHM model version release.*
- character(len=\*), parameter `mo_file::file_main` = 'mhm\_driver.f90'  
*Driver file.*
- character(len=\*), parameter `mo_file::file_namelist_mhm` = 'mhm.nml'  
*Namelist file name.*
- integer, parameter `mo_file::unamelist_mhm` = 30  
*Unit for namelist.*
- character(len=\*), parameter `mo_file::file_namelist_mhm_param` = 'mhm\_parameter.nml'  
*Parameter namelists file name.*
- integer, parameter `mo_file::unamelist_mhm_param` = 31  
*Unit for namelist.*
- character(len=\*), parameter `mo_file::file_defoutput` = 'mhm\_outputs.nml'  
*file defining mHM's outputs*
- integer, parameter `mo_file::udefoutput` = 67  
*Unit for file defining mHM's outputs.*

## 19.38 src/mHM/mo\_global\_variables.f90 File Reference

### Modules

- module `mo_global_variables`  
*Global variables ONLY used in reading, writing and startup.*

### Variables

- integer(i4) `mo_global_variables::output_deflate_level`
- logical `mo_global_variables::output_double_precision`
- integer(i4) `mo_global_variables::timestep_model_outputs`
- logical, dimension(noutflxstate) `mo_global_variables::outputflxstate`
- integer(i4), dimension(:), allocatable, public `mo_global_variables::timestep_model_inputs`
- logical, public `mo_global_variables::read_meteo_weights`
- character(256), public `mo_global_variables::inputformat_meteo_forcings`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirprecipitation`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirtemperature`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirmintemperature`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirmaxtemperature`

- character(256), dimension(:), allocatable, public `mo_global_variables::dirnetradiation`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirabsvappressure`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirwindspeed`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirreferencecet`
- character(256), dimension(:), allocatable, public `mo_global_variables::dirradiation`
- integer(i4), parameter, public `mo_global_variables::routingstates = 2`
- type(grid), dimension(:), allocatable, public `mo_global_variables::level2`
- real(dp), dimension(:, :, :), allocatable, public `mo_global_variables::l1_temp_weights`
- real(dp), dimension(:, :, :), allocatable, public `mo_global_variables::l1_pet_weights`
- real(dp), dimension(:, :, :), allocatable, public `mo_global_variables::l1_pre_weights`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_pre`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_temp`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_pet`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_tmin`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_tmax`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_netrad`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_absvappress`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_windspeed`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_ssrd`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_strd`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_tann`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_sm`
- logical, dimension(:, :), allocatable, public `mo_global_variables::l1_sm_mask`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_neutronsdata`
- logical, dimension(:, :), allocatable, public `mo_global_variables::l1_neutronsdata_mask`
- integer(i4) `mo_global_variables::nsoilhorizons_sm_input`
- type(optidata), dimension(:), allocatable, public `mo_global_variables::l1_smobs`
- type(optidata), dimension(:), allocatable, public `mo_global_variables::l1_neutronsobs`
- type(optidata), dimension(:), allocatable, public `mo_global_variables::l1_etobs`
- type(optidata), dimension(:), allocatable, public `mo_global_variables::l1_twsaobs`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_inter`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_snowpack`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_sealstw`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_soilmoist`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_unsatstw`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_satstw`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_neutrons`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_pet_calc`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_aetsoil`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_aetcanopy`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_aetsealed`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_baseflow`
- real(dp), dimension(:, :), allocatable, public `mo_global_variables::l1_infilsoil`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_fastrunoff`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_melt`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_percol`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_preeffect`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_rain`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_runoffseal`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_slowrunoff`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_snow`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_throughfall`
- real(dp), dimension(:), allocatable, public `mo_global_variables::l1_total_runoff`
- real(dp), dimension(yearmonths\_i4), public `mo_global_variables::evap_coeff`
- real(dp), dimension(yearmonths\_i4), public `mo_global_variables::fday_prec`

- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fnight\\_prec](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fday\\_pet](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fnight\\_pet](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fday\\_temp](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fnight\\_temp](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fday\\_ssrđ](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fnight\\_ssrđ](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fday\\_strđ](#)
- real(dp), dimension(yearmonths\_i4), public [mo\\_global\\_variables::fnight\\_strđ](#)
- real(dp), dimension(:), allocatable, public [mo\\_global\\_variables::neutron\\_integral\\_ąfast](#)

## 19.39 src/mHM/mo\_init\_states.f90 File Reference

### Modules

- module [mo\\_init\\_states](#)  
*Initialization of all state variables of mHM.*

### Functions/Subroutines

- subroutine, public [mo\\_init\\_states::variables\\_alloc](#) (ncells1)  
*Allocation of space for mHM related L1 and L11 variables.*
- subroutine, public [mo\\_init\\_states::variables\\_default\\_init](#)  
*Default initialization mHM related L1 variables.*

## 19.40 src/mHM/mo\_meteo\_forcings.f90 File Reference

### Modules

- module [mo\\_meteo\\_forcings](#)  
*Prepare meteorological forcings data for mHM.*

### Functions/Subroutines

- subroutine, public [mo\\_meteo\\_forcings::prepare\\_meteo\\_forcings\\_data](#) (iDomain, domainID, tt)  
*Prepare meteorological forcings data for a given variable.*
- subroutine [mo\\_meteo\\_forcings::meteo\\_forcings\\_wrapper](#) (iDomain, dataPath, inputFormat, dataOut1, lower, upper, ncvrName)  
*Prepare meteorological forcings data for mHM at Level-1.*
- subroutine [mo\\_meteo\\_forcings::meteo\\_weights\\_wrapper](#) (iDomain, read\_meteo\_weights, dataPath, dataOut1, lower, upper, ncvrName)  
*Prepare weights for meteorological forcings data for mHM at Level-1.*
- subroutine [mo\\_meteo\\_forcings::chunk\\_config](#) (iDomain, tt, read\_flag, readPer)  
*determines the start date, end date, and read\_flag given Domain id and current timestep*
- logical function [mo\\_meteo\\_forcings::is\\_read](#) (iDomain, tt)  
*evaluate whether new chunk should be read at this timestep*
- subroutine [mo\\_meteo\\_forcings::chunk\\_size](#) (iDomain, tt, readPer)  
*calculate beginning and end of read Period, i.e. that is length of current chunk to read*

## 19.41 src/mHM/mo\_mhm.f90 File Reference

### Modules

- module [mo\\_mhm](#)  
*Call all main processes of mHM.*

## Functions/Subroutines

- subroutine, public `mo_mhm::mhm` (`read_states`, `tt`, `time`, `processMatrix`, `horizon_depth`, `nCells1`, `nHorizons`, `mHM`, `ntimesteps_day`, `c2TSTu`, `neutron_integral_AFast`, `global_parameters`, `latitude`, `evap_coeff`, `fday`, `prec`, `fnight_prec`, `fday_pet`, `fnight_pet`, `fday_temp`, `fnight_temp`, `temp_weights`, `pet_weights`, `pre_weights`, `read_meteo_weights`, `pet_in`, `tmin_in`, `tmax_in`, `netrad_in`, `absvappres_in`, `windspeed_in`, `prec_in`, `temp_in`, `fSealed1`, `interc`, `snowpack`, `sealedStorage`, `soilMoisture`, `unsatStorage`, `satStorage`, `neutrons`, `pet_calc`, `aet_soil`, `aet_canopy`, `aet_sealed`, `baseflow`, `infiltration`, `fast_interflow`, `melt`, `perc`, `prec_effect`, `rain`, `runoff_sealed`, `slow_interflow`, `snow`, `throughfall`, `total_runoff`, `alpha`, `deg_day_incr`, `deg_day_max`, `deg_day_noprec`, `deg_day`, `fAsp`, `petLAlcorFactorL1`, `HarSamCoeff`, `PrieTayAlpha`, `aeroResist`, `surfResist`, `frac_roots`, `interc_max`, `karst_loss`, `k0`, `k1`, `k2`, `kp`, `soil_moist_FC`, `soil_moist_sat`, `soil_moist_exponen`, `jarvis_thresh_c1`, `temp_thresh`, `unsat_thresh`, `water_thresh_sealed`, `wilting_point`)

*Pure mHM calculations.*

## 19.42 src/mHM/mo\_mhm\_constants.f90 File Reference

### Modules

- module `mo_mhm_constants`

*Provides mHM specific constants.*

### Variables

- real(dp), parameter, public `mo_mhm_constants::h2odens` = 1000.0\_dp
- real(dp), parameter, public `mo_mhm_constants::p2_initstatefluxes` = 15.00\_dp
- real(dp), parameter, public `mo_mhm_constants::p3_initstatefluxes` = 10.00\_dp
- real(dp), parameter, public `mo_mhm_constants::p4_initstatefluxes` = 75.00\_dp
- real(dp), parameter, public `mo_mhm_constants::p5_initstatefluxes` = 1500.00\_dp
- real(dp), parameter, public `mo_mhm_constants::c1_initstatesm` = 0.25\_dp
- integer(i4), parameter, public `mo_mhm_constants::noutflxstate` = 20\_i4
- real(dp), parameter, public `mo_mhm_constants::harsamconst` = 17.800\_dp

*Hargreaves-Samani ref. ET formula [deg C].*

- real(dp), parameter, public `mo_mhm_constants::duffiedr` = 0.0330\_dp
- real(dp), parameter, public `mo_mhm_constants::duffiedelta1` = 0.4090\_dp
- real(dp), parameter, public `mo_mhm_constants::duffiedelta2` = 1.3900\_dp
- real(dp), parameter, public `mo_mhm_constants::tetens_c1` = 0.6108\_dp

*Tetens's formula to calculate saturated vapour pressure.*

- real(dp), parameter, public `mo_mhm_constants::tetens_c2` = 17.270\_dp
- real(dp), parameter, public `mo_mhm_constants::tetens_c3` = 237.30\_dp
- real(dp), parameter, public `mo_mhm_constants::satpressureslope1` = 4098.0\_dp

*calculation of the slope of the saturation vapour pressure curve following Tetens*

- real(dp), parameter, public `mo_mhm_constants::desilets_a0` = 0.0808\_dp

*Neutrons and moisture: N0 formula, Desilets et al. 2010.*

- real(dp), parameter, public `mo_mhm_constants::desilets_a1` = 0.372\_dp
- real(dp), parameter, public `mo_mhm_constants::desilets_a2` = 0.115\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_bd` = 1.4020\_dp

*Neutrons and moisture: COSMIC, Shuttleworth et al. 2013.*

- real(dp), parameter, public `mo_mhm_constants::cosmic_vwclat` = 0.0753\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_n` = 348.33\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_alpha` = 0.2392421548\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_l1` = 161.98621864\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_l2` = 129.14558985\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_l3` = 107.82204562\_dp
- real(dp), parameter, public `mo_mhm_constants::cosmic_l4` = 3.1627190566\_dp

## 19.43 src/mHM/mo\_mhm\_eval.f90 File Reference

### Modules

- module [mo\\_mhm\\_eval](#)

*Runs mhm with a specific parameter set and returns required variables, e.g. runoff.*

### Functions/Subroutines

- subroutine, public [mo\\_mhm\\_eval::mhm\\_eval](#) (parameterset, opti\_domain\_indices, runoff, smOptiSim, neutronsOptiSim, etOptiSim, twsOptiSim)

*Runs mhm with a specific parameter set and returns required variables, e.g. runoff.*

## 19.44 src/mHM/mo\_mhm\_read\_config.f90 File Reference

### Modules

- module [mo\\_mhm\\_read\\_config](#)

*Reading of main model configurations.*

### Functions/Subroutines

- subroutine, public [mo\\_mhm\\_read\\_config::mhm\\_read\\_config](#) (file\_namelist, unamelist)

*Read main configurations for mHM.*

## 19.45 src/mHM/mo\_neutrons.f90 File Reference

### Modules

- module [mo\\_neutrons](#)

*Models to predict neutron intensities above soils.*

### Functions/Subroutines

- subroutine, public [mo\\_neutrons::desiletsn0](#) (SoilMoisture, Horizons, N0, neutrons)  
*Calculate neutrons from soil moisture in the first layer.*
- subroutine, public [mo\\_neutrons::cosmic](#) (SoilMoisture, Horizons, params, neutron\_integral\_AFast, neutrons)  
*Calculate neutrons from soil moisture in all layers.*
- subroutine [mo\\_neutrons::oldintegration](#) (res, c)  
*TODO: add description.*
- subroutine, public [mo\\_neutrons::tabularintegralafast](#) (integral, maxC)  
*Save approximation data for A\_fast.*
- subroutine [mo\\_neutrons::approx\\_mon\\_int](#) (res, f, c, xmin, xmax, eps, steps, fxmin, fxmax)  
*TODO: add description.*
- recursive subroutine [mo\\_neutrons::approx\\_mon\\_int\\_steps](#) (res, f, c, xmin, xmax, eps, steps, fxmin, fxmax)  
*TODO: add description.*
- recursive subroutine [mo\\_neutrons::approx\\_mon\\_int\\_eps](#) (res, f, c, xmin, xmax, eps, fxmin, fxmax)  
*TODO: add description.*
- subroutine [mo\\_neutrons::lookupintegral](#) (res, integral, c)  
*TODO: add description.*
- real(dp) function [mo\\_neutrons::intgrandfast](#) (c, phi)  
*TODO: add description.*



## 19.46 src/mHM/mo\_objective\_function.f90 File Reference

### Modules

- module [mo\\_objective\\_function](#)  
*Objective Functions for Optimization of mHM.*

### Functions/Subroutines

- real(dp) function, public [mo\\_objective\\_function::objective](#) (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions.*
- real(dp) function, public [mo\\_objective\\_function::objective\\_master](#) (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions.*
- subroutine, public [mo\\_objective\\_function::objective\\_subprocess](#) (eval, arg1, arg2, arg3)  
*Wrapper for objective functions.*
- real(dp) function [mo\\_objective\\_function::objective\\_sm\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for soil moisture.*
- real(dp) function, dimension(6) [mo\\_objective\\_function::objective\\_q\\_et\\_tws\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for et, tws and q.*
- subroutine [mo\\_objective\\_function::init\\_indexarray\\_for\\_opti\\_data](#) (domainMeta, optidataOption, nOptiDomains, opti\_domain\_indices)  
*creates an index array of the inidices of the domains eval should MPI process.*
- real(dp) function [mo\\_objective\\_function::objective\\_sm\\_corr](#) (parameterset, eval)  
*Objective function for soil moisture.*
- real(dp) function [mo\\_objective\\_function::objective\\_sm\\_pd](#) (parameterset, eval)  
*Objective function for soil moisture.*
- real(dp) function [mo\\_objective\\_function::objective\\_sm\\_sse\\_standard\\_score](#) (parameterset, eval)  
*Objective function for soil moisture.*
- real(dp) function [mo\\_objective\\_function::objective\\_kge\\_q\\_rmse\\_tws](#) (parameterset, eval)  
*Objective function of KGE for runoff and RMSE for domain\_avg TWS (standarized scores)*
- real(dp) function [mo\\_objective\\_function::objective\\_neutrons\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for neutrons.*
- real(dp) function [mo\\_objective\\_function::objective\\_et\\_kge\\_catchment\\_avg](#) (parameterset, eval)  
*Objective function for evpotranspirstion (et).*
- real(dp) function [mo\\_objective\\_function::objective\\_kge\\_q\\_sm\\_corr](#) (parameterset, eval)  
*Objective function of KGE for runoff and correlation for SM.*
- real(dp) function [mo\\_objective\\_function::objective\\_kge\\_q\\_et](#) (parameterset, eval)  
*Objective function of KGE for runoff and KGE for ET.*
- real(dp) function [mo\\_objective\\_function::objective\\_kge\\_q\\_rmse\\_et](#) (parameterset, eval)  
*Objective function of KGE for runoff and RMSE for domain\_avg ET (standarized scores)*
- subroutine [mo\\_objective\\_function::create\\_domain\\_avg\\_tws](#) (iDomain, twsOptiSim, tws\_catch\_avg\_domain, tws\_opti\_catch\_avg\_domain, mask\_times)
- subroutine [mo\\_objective\\_function::create\\_domain\\_avg\\_et](#) (iDomain, etOptiSim, et\_catch\_avg\_domain, et\_opti\_catch\_avg\_domain, mask\_times)
- subroutine [mo\\_objective\\_function::convert\\_tws\\_to\\_twsa](#) (twsOptiSim, L1\_twsaObs, twsaOptiSim)

## 19.47 src/mHM/mo\_pet.f90 File Reference

### Modules

- module [mo\\_pet](#)  
*Module for calculating reference/potential evapotranspiration [mm d-1].*

## Functions/Subroutines

- elemental pure real(dp) function, public [mo\\_pet::pet\\_hargreaves](#) (HarSamCoeff, HarSamConst, tavg, tmax, tmin, latitude, doy)  
*Reference Evapotranspiration after Hargreaves.*
- elemental pure real(dp) function, public [mo\\_pet::pet\\_priestly](#) (PriTayParam, Rn, tavg)  
*Reference Evapotranspiration after Priestly-Taylor.*
- elemental pure real(dp) function, public [mo\\_pet::pet\\_penman](#) (net\_rad, tavg, act\_vap\_pressure, aerodyn\_↔resistance, bulksurface\_resistance, a\_s, a\_sh)  
*Reference Evapotranspiration after Penman-Monteith.*
- elemental pure real(dp) function, public [mo\\_pet::extraterr\\_rad\\_approx](#) (doy, latitude)  
*Approximation of extraterrestrial radiation.*
- elemental pure real(dp) function, public [mo\\_pet::slope\\_satpressure](#) (tavg)  
*slope of saturation vapour pressure curve*
- elemental pure real(dp) function, public [mo\\_pet::sat\\_vap\\_pressure](#) (tavg)  
*calculation of the saturation vapour pressure*

## 19.48 src/mHM/mo\_read\_optional\_data.f90 File Reference

### Modules

- module [mo\\_read\\_optional\\_data](#)  
*Read optional data for mHM calibration.*

### Functions/Subroutines

- subroutine, public [mo\\_read\\_optional\\_data::readoptidataobs](#) (iDomain, domainID, L1\_optiObs)  
*Read evapotranspiration data from NetCDF file for calibration.*

## 19.49 src/mHM/mo\_restart.f90 File Reference

### Data Types

- interface [mo\\_restart::unpack\\_field\\_and\\_write](#)  
*TODO: add description.*

### Modules

- module [mo\\_restart](#)  
*reading and writing states, fluxes and configuration for restart of mHM.*

### Functions/Subroutines

- subroutine, public [mo\\_restart::write\\_restart\\_files](#) (OutFile)  
*write restart files for each domain*
- subroutine, public [mo\\_restart::read\\_restart\\_states](#) (iDomain, domainID, InFile)  
*reads fluxes and state variables from file*
- subroutine [mo\\_restart::unpack\\_field\\_and\\_write\\_1d\\_i4](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_↔\_long\_name)
- subroutine [mo\\_restart::unpack\\_field\\_and\\_write\\_1d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_↔\_long\_name)
- subroutine [mo\\_restart::unpack\\_field\\_and\\_write\\_2d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_↔\_long\_name)
- subroutine [mo\\_restart::unpack\\_field\\_and\\_write\\_3d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_↔\_long\_name)

## 19.50 src/mHM/mo\_runoff.f90 File Reference

### Modules

- module [mo\\_runoff](#)

*Runoff generation for the unsaturated zone, saturated zone (or groundwater zone), and runoff accumulation.*

### Functions/Subroutines

- subroutine, public [mo\\_runoff::runoff\\_unsat\\_zone](#) (k1, kp, k0, alpha, karst\_loss, pefec\_soil, unsat\_thresh, sat\_storage, unsat\_storage, slow\_interflow, fast\_interflow, perc)

*Runoff generation for the saturated zone.*

- subroutine, public [mo\\_runoff::runoff\\_sat\\_zone](#) (k2, sat\_storage, baseflow)

*Runoff generation for the saturated zone.*

- subroutine, public [mo\\_runoff::l1\\_total\\_runoff](#) (fSealed\_area\_fraction, fast\_interflow, slow\_interflow, baseflow, direct\_runoff, total\_runoff)

*total runoff accumulation at level 1*

## 19.51 src/mHM/mo\_snow\_accum\_melt.f90 File Reference

### Modules

- module [mo\\_snow\\_accum\\_melt](#)

*Snow melting and accumulation.*

### Functions/Subroutines

- subroutine, public [mo\\_snow\\_accum\\_melt::snow\\_accum\\_melt](#) (deg\_day\_incr, deg\_day\_max, deg\_day\_max\_noprec, prec, temperature, temperature\_thresh, thrfall, snow\_pack, deg\_day, melt, prec\_effect, rain, snow)

*Snow melting and accumulation.*

## 19.52 src/mHM/mo\_soil\_moisture.f90 File Reference

### Modules

- module [mo\\_soil\\_moisture](#)

*Soil moisture of the different layers.*

### Functions/Subroutines

- subroutine, public [mo\\_soil\\_moisture::soil\\_moisture](#) (processCase, frac\_sealed, water\_thresh\_sealed, pet, evap\_coeff, soil\_moist\_sat, frac\_roots, soil\_moist\_FC, wilting\_point, soil\_moist\_exponen, jarvis\_thresh\_c1, aet\_canopy, prec\_effec, runoff\_sealed, storage\_sealed, infiltration, soil\_moist, aet, aet\_sealed)

*Soil moisture in different soil horizons.*

- elemental pure real(dp) function, public [mo\\_soil\\_moisture::feddes\\_et\\_reduction](#) (soil\_moist, soil\_moist\_FC, wilting\_point, frac\_roots)

*stress factor for reducing evapotranspiration based on actual soil moisture*

- elemental pure real(dp) function, public [mo\\_soil\\_moisture::jarvis\\_et\\_reduction](#) (soil\_moist, soil\_moist\_sat, wilting\_point, frac\_roots, jarvis\_thresh\_c1)

*stress factor for reducing evapotranspiration based on actual soil moisture*

## 19.53 src/mHM/mo\_spatial\_agg\_disagg\_forcing.f90 File Reference

### Data Types

- interface [mo\\_spatial\\_agg\\_disagg\\_forcing::spatial\\_aggregation](#)  
*Spatial aggregation of meteorological variables.*
- interface [mo\\_spatial\\_agg\\_disagg\\_forcing::spatial\\_disaggregation](#)  
*Spatial disaggregation of meteorological variables.*

### Modules

- module [mo\\_spatial\\_agg\\_disagg\\_forcing](#)  
*Spatial aggregation or disaggregation of meteorological input data.*

### Functions/Subroutines

- subroutine [mo\\_spatial\\_agg\\_disagg\\_forcing::spatial\\_aggregation\\_3d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [mo\\_spatial\\_agg\\_disagg\\_forcing::spatial\\_aggregation\\_4d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [mo\\_spatial\\_agg\\_disagg\\_forcing::spatial\\_disaggregation\\_3d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)
- subroutine [mo\\_spatial\\_agg\\_disagg\\_forcing::spatial\\_disaggregation\\_4d](#) (data2, cellsize2, cellsize1, mask1, mask2, data1)

## 19.54 src/mHM/mo\_startup.f90 File Reference

### Modules

- module [mo\\_startup](#)  
*Startup procedures for mHM.*

### Functions/Subroutines

- subroutine, public [mo\\_startup::mhm\\_initialize](#)  
*Initialize main mHM variables.*
- subroutine [mo\\_startup::constants\\_init](#)  
*Initialize mHM constants.*
- subroutine [mo\\_startup::l2\\_variable\\_init](#) (iDomain, level0\_iDomain, level2\_iDomain)  
*Initialize Level-2 meteorological forcings data.*

## 19.55 src/mHM/mo\_temporal\_disagg\_forcing.f90 File Reference

### Modules

- module [mo\\_temporal\\_disagg\\_forcing](#)  
*Temporal disaggregation of daily input values.*

### Functions/Subroutines

- elemental pure subroutine, public [mo\\_temporal\\_disagg\\_forcing::temporal\\_disagg\\_forcing](#) (isday, ntimesteps↔\_day, prec\_day, pet\_day, temp\_day, fday\_prec, fday\_pet, fday\_temp, fnight\_prec, fnight\_pet, fnight\_temp, temp\_weights, pet\_weights, pre\_weights, read\_meteo\_weights, prec, pet, temp)  
*Temporally distribute daily mean forcings onto time step.*

- elemental subroutine, public [mo\\_temporal\\_disagg\\_forcing::temporal\\_disagg\\_meteo\\_weights](#) (meteo\_val↔  
day, meteo\_val\_weights, meteo\_val, weights\_correction)  
*Temporally distribute daily mean forcings onto time step.*
- elemental subroutine, public [mo\\_temporal\\_disagg\\_forcing::temporal\\_disagg\\_flux\\_daynight](#) (isday, ntimesteps↔  
\_day, meteo\_val\_day, fday\_meteo\_val, fnight\_meteo\_val, meteo\_val)  
*Temporally distribute daily mean forcings onto time step.*
- elemental subroutine, public [mo\\_temporal\\_disagg\\_forcing::temporal\\_disagg\\_state\\_daynight](#) (isday,  
ntimesteps\_day, meteo\_val\_day, fday\_meteo\_val, fnight\_meteo\_val, meteo\_val, add\_correction)  
*Temporally distribute daily mean state forcings onto time step.*

## 19.56 src/mHM/mo\_write\_ascii.f90 File Reference

### Modules

- module [mo\\_write\\_ascii](#)  
*Module to write ascii file output.*

### Functions/Subroutines

- subroutine, public [mo\\_write\\_ascii::write\\_configfile](#)  
*This modules writes the results of the configuration into an ASCII-file.*
- subroutine, public [mo\\_write\\_ascii::write\\_optifile](#) (best\_OF, best\_paramSet, param\_names)  
*Write briefly final optimization results.*
- subroutine, public [mo\\_write\\_ascii::write\\_optinamelist](#) (processMatrix, parameters, maskpara, parameters↔  
name)  
*Write final, optimized parameter set in a namelist format.*

## 19.57 src/mHM/mo\_write\_fluxes\_states.f90 File Reference

### Data Types

- interface [mo\\_write\\_fluxes\\_states::outputvariable](#)
- interface [mo\\_write\\_fluxes\\_states::outputdataset](#)

### Modules

- module [mo\\_write\\_fluxes\\_states](#)  
*Creates NetCDF output for different fluxes and state variables of mHM.*

### Functions/Subroutines

- type(outputvariable) function [mo\\_write\\_fluxes\\_states::newoutputvariable](#) (nc, name, dtype, dims, ncells,  
mask, avg)  
*Initialize OutputVariable.*
- subroutine [mo\\_write\\_fluxes\\_states::updatevariable](#) (self, data)  
*Update OutputVariable.*
- subroutine [mo\\_write\\_fluxes\\_states::writevariabletimestep](#) (self, timestep)  
*Write timestep to file.*
- type(outputdataset) function, public [mo\\_write\\_fluxes\\_states::newoutputdataset](#) (iDomain, mask1, nCells)  
*Initialize OutputDataset.*
- subroutine [mo\\_write\\_fluxes\\_states::updatedataset](#) (self, sidx, eidx, L1\_fSealed, L1\_fNotSealed, L1\_inter,  
L1\_snowPack, L1\_soilMoist, L1\_soilMoistSat, L1\_sealSTW, L1\_unsatSTW, L1\_satSTW, L1\_neutrons, L1↔  
pet, L1\_aETSoil, L1\_aETCanopy, L1\_aETSealed, L1\_total\_runoff, L1\_runoffSeal, L1\_fastRunoff, L1\_slow↔  
Runoff, L1\_baseflow, L1\_percol, L1\_infilSoil, L1\_preEffect)

- Update all variables.*

  - subroutine `mo_write_fluxes_states::writetimestep` (self, timestep)

*Write all accumulated data.*

  - subroutine `mo_write_fluxes_states::close` (self)

*Close the file.*

  - type(ncdataset) function `mo_write_fluxes_states::createoutputfile` (iDomain)

*Create and initialize output file for X & Y coordinate system.*

  - subroutine `mo_write_fluxes_states::writevariableattributes` (var, long\_name, unit)

*Write output variable attributes.*

  - character(16) function `mo_write_fluxes_states::fluxesunit` (iDomain)

*Generate a unit string.*

## 19.58 src/MPR/mo\_mpr\_constants.f90 File Reference

### Modules

- module `mo_mpr_constants`
- Provides MPR specific constants.*

### Variables

- integer(i4), parameter, public `mo_mpr_constants::nlcover_class` = 3\_i4
- integer(i4), parameter, public `mo_mpr_constants::maxgeounit` = 25\_i4
- integer(i4), parameter, public `mo_mpr_constants::maxnosoilhorizons` = 10\_i4
- real(dp), parameter, public `mo_mpr_constants::p2_initstatefluxes` = 15.00\_dp
- real(dp), parameter, public `mo_mpr_constants::p3_initstatefluxes` = 10.00\_dp
- real(dp), parameter, public `mo_mpr_constants::p4_initstatefluxes` = 75.00\_dp
- real(dp), parameter, public `mo_mpr_constants::p5_initstatefluxes` = 1500.00\_dp
- real(dp), parameter, public `mo_mpr_constants::c1_initstatesm` = 0.25\_dp
- real(dp), parameter, public `mo_mpr_constants::bulkdens_ormatter` = 0.224\_dp
- real(dp), parameter, public `mo_mpr_constants::field_cap_c1` = -0.60\_dp
- real(dp), parameter, public `mo_mpr_constants::field_cap_c2` = 2.0\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchten_sandfresh` = 66.5\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c1` = 1.392\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c2` = 0.418\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c3` = -0.024\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c4` = 1.212\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c5` = -0.704\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c6` = -0.648\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c7` = 0.023\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c8` = 0.044\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c9` = 3.168\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c10` = -2.562\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c11` = 7.0E-9\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c12` = 4.004\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c13` = 3.750\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c14` = -0.016\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c15` = -4.197\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c16` = 0.013\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c17` = 0.076\_dp
- real(dp), parameter, public `mo_mpr_constants::vgenuchtenn_c18` = 0.276\_dp
- real(dp), parameter, public `mo_mpr_constants::ks_c` = 10.0\_dp
- real(dp), parameter, public `mo_mpr_constants::pwp_c` = 1.0\_dp
- real(dp), parameter, public `mo_mpr_constants::pwp_matpot_thetar` = 15000.0\_dp

- real(dp), parameter, public `mo_mpr_constants::windmeasheight` = 10.0\_dp  
*assumed meteorol. measurement hight for estimation of aeroResist and surfResist*
- real(dp), parameter, public `mo_mpr_constants::karman` = 0.41\_dp  
*von karman constant*
- real(dp), parameter, public `mo_mpr_constants::lai_factor_surfresi` = 0.3\_dp  
*LAI factor for bulk surface resistance formulation.*
- real(dp), parameter, public `mo_mpr_constants::lai_offset_surfresi` = 1.2\_dp  
*LAI offset for bulk surface resistance formulation.*
- real(dp), parameter, public `mo_mpr_constants::max_surfresist` = 250.0\_dp  
*maximum bulk surface resistance*

## 19.59 src/MPR/mo\_mpr\_eval.f90 File Reference

### Modules

- module `mo_mpr_eval`  
*Runs MPR and writes to global effective parameters.*

### Functions/Subroutines

- subroutine, public `mo_mpr_eval::mpr_eval` (parameterset, opti\_domain\_indices)  
*Runs MPR and writes to global effective parameters.*

## 19.60 src/MPR/mo\_mpr\_file.f90 File Reference

### Modules

- module `mo_mpr_file`  
*Provides file names and units for mRM.*

### Variables

- character(len=\*), parameter `mo_mpr_file::version` = '0.1'  
*Current mHM model version.*
- character(len=\*), parameter `mo_mpr_file::version_date` = 'Jun 2019'  
*Time of current mHM model version release.*
- character(len=\*), parameter `mo_mpr_file::file_main` = 'mpr\_driver.f90'  
*Driver file.*
- character(len=\*), parameter `mo_mpr_file::file_namelist_mpr` = 'mpr.nml'  
*Namelist file name.*
- integer, parameter `mo_mpr_file::unamelist_mpr` = 80  
*Unit for namelist.*
- character(len=\*), parameter `mo_mpr_file::file_namelist_mpr_param` = 'mpr\_parameter.nml'  
*Parameter namelists file name.*
- integer, parameter `mo_mpr_file::unamelist_mpr_param` = 31  
*Unit for namelist.*
- character(len=\*), parameter `mo_mpr_file::file_soil_database` = 'soil\_classdefinition.txt'  
*Soil database file (iFlag\_soilDB = 0) = classical mHM format.*
- character(len=\*), parameter `mo_mpr_file::file_soil_database_1` = 'soil\_classdefinition\_iFlag\_soilDB\_1.txt'  
*Soil database file (iFlag\_soilDB = 1)*
- integer, parameter `mo_mpr_file::usoil_database` = 52  
*Unit for soil data base.*

- character(len=\*), parameter `mo_mpr_file::file_slope` = 'slope.asc'  
*slope input data file*
- integer, parameter `mo_mpr_file::uslope` = 54  
*Unit for slope input data file.*
- character(len=\*), parameter `mo_mpr_file::file_aspect` = 'aspect.asc'  
*aspect input data file*
- integer, parameter `mo_mpr_file::uaspect` = 55  
*Unit for aspect input data file.*
- character(len=\*), parameter `mo_mpr_file::file_hydrogeoclass` = 'geology\_class.asc'  
*hydrogeological classes input data file*
- integer, parameter `mo_mpr_file::uhydrogeoclass` = 58  
*Unit for hydrogeological classes input data file.*
- character(len=\*), parameter `mo_mpr_file::file_soilclass` = 'soil\_class.asc'  
*soil classes input data file*
- integer, parameter `mo_mpr_file::usoilclass` = 59  
*Unit for soil classes input data file.*
- character(len=\*), parameter `mo_mpr_file::file_laiclass` = 'LAI\_class.asc'  
*LAI classes input data file.*
- integer, parameter `mo_mpr_file::ulaiclass` = 60  
*Unit for LAI input data file.*
- character(len=\*), parameter `mo_mpr_file::file_geolut` = 'geology\_classdefinition.txt'  
*geological formation lookup table file*
- integer, parameter `mo_mpr_file::ugeolut` = 64  
*Unit for geological formation lookup table file.*
- character(len=\*), parameter `mo_mpr_file::file_lailut` = 'LAI\_classdefinition.txt'  
*LAI classes lookup table file.*
- integer, parameter `mo_mpr_file::ulailut` = 65  
*Unit for LAI classes lookup table file.*
- character(len=\*), parameter `mo_mpr_file::file_meteo_header` = 'header.txt'  
*Input nCols and nRows of binary meteo and LAI files are in header file.*
- integer, parameter `mo_mpr_file::umeteo_header` = 50  
*Unit for meteo header file.*
- character(len=\*), parameter `mo_mpr_file::file_meteo_binary_end` = '.bin'  
*File ending of meteo files.*
- integer, parameter `mo_mpr_file::umeteo` = 51  
*Unit for meteo files.*

## 19.61 src/MPR/mo\_mpr\_global\_variables.f90 File Reference

### Data Types

- type `mo_mpr_global_variables::soiltype`

### Modules

- module `mo_mpr_global_variables`  
*Global variables for mpr only.*



## Variables

- real(dp), public `mo_mpr_global_variables::tillagedepth`
- integer(i4), public `mo_mpr_global_variables::nsoiltypes`
- integer(i4), public `mo_mpr_global_variables::iflag_soildb`
- integer(i4), public `mo_mpr_global_variables::nsoilhorizons_mhm`
- real(dp), dimension(:), allocatable, public `mo_mpr_global_variables::horizondepth_mhm`
- type(soiltype), public `mo_mpr_global_variables::soildb`
- integer(i4), public `mo_mpr_global_variables::ngeounits`
- integer(i4), dimension(:), allocatable, public `mo_mpr_global_variables::geounitlist`
- integer(i4), dimension(:), allocatable, public `mo_mpr_global_variables::geounitkar`
- character(256), public `mo_mpr_global_variables::inputformat_gridded_lai`
- integer(i4), public `mo_mpr_global_variables::timestep_lai_input`
- integer(i4), public `mo_mpr_global_variables::nlaiclass`
- integer(i4), public `mo_mpr_global_variables::nlai`
- real(dp), dimension(:), allocatable, public `mo_mpr_global_variables::laiboundaries`
- integer(i4), dimension(:), allocatable, public `mo_mpr_global_variables::laiunitlist`
- real(dp), dimension(:, :), allocatable, public `mo_mpr_global_variables::lailut`
- type(period), dimension(:), allocatable, public `mo_mpr_global_variables::laiper`
- real(dp), public `mo_mpr_global_variables::fracsealed_cityarea`
- real(dp), dimension(:), allocatable, public `mo_mpr_global_variables::l0_slope_emp`
- real(dp), dimension(:, :), allocatable, public `mo_mpr_global_variables::l0_gridded_lai`
- real(dp), dimension(:), allocatable, public `mo_mpr_global_variables::l0_slope`
- real(dp), dimension(:), allocatable, public `mo_mpr_global_variables::l0_asp`
- integer(i4), dimension(:, :), allocatable, public `mo_mpr_global_variables::l0_soilid`
- integer(i4), dimension(:), allocatable, public `mo_mpr_global_variables::l0_geounit`
- character(256), dimension(:), allocatable, public `mo_mpr_global_variables::dirgridded_lai`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_fsealed`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_alpha`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_degdayinc`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_degdaymax`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_degdaynopre`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_degday`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_karstloss`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_fasp`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_petlaicorfactor`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_harsamcoeff`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_prietayalpha`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_aeroresist`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_surfresist`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_frosts`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_maxinter`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_kfastflow`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_kslowflow`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_kbaseflow`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_kperco`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_soilmoistfc`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_soilmoistsat`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_soilmoistexp`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_jarvis_thresh_c1`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_tempthresh`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_unsatthresh`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_sealedthresh`
- real(dp), dimension(:, :, :), allocatable, public `mo_mpr_global_variables::l1_wiltingpoint`

## 19.62 src/MPR/mo\_mpr\_pet.f90 File Reference

### Modules

- module [mo\\_mpr\\_pet](#)  
*TODO: add description.*

### Functions/Subroutines

- subroutine, public [mo\\_mpr\\_pet::pet\\_correctbylai](#) (param, nodata, LCOVER0, LAI0, mask0, cell\_id0, upp\_↔row\_L1, low\_row\_L1, lef\_col\_L1, rig\_col\_L1, nL0\_in\_L1, L1\_petLAIcorFactor)  
*estimate PET correction factor based on LAI at L1*
- subroutine, public [mo\\_mpr\\_pet::pet\\_correctbyasp](#) (ld0, latitude\_l0, Asp0, param, nodata, fAsp0)  
*correction of PET*
- subroutine, public [mo\\_mpr\\_pet::priestley\\_taylor\\_alpha](#) (LAI0, param, mask0, nodata, cell\_id0, nL0\_in\_L1, Upp\_row\_L1, Low\_row\_L1, Lef\_col\_L1, Rig\_col\_L1, priestley\_taylor\_alpha1)  
*Regionalization of priestley taylor alpha.*
- subroutine, public [mo\\_mpr\\_pet::bulksurface\\_resistance](#) (LAI0, param, mask0, nodata, cell\_id0, nL0\_in\_L1, Upp\_row\_L1, Low\_row\_L1, Lef\_col\_L1, Rig\_col\_L1, bulksurface\_resistance1)  
*Regionalization of bulk surface resistance.*

## 19.63 src/MPR/mo\_mpr\_read\_config.f90 File Reference

### Modules

- module [mo\\_mpr\\_read\\_config](#)  
*read mpr config*

### Functions/Subroutines

- subroutine, public [mo\\_mpr\\_read\\_config::mpr\\_read\\_config](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param)  
*Read the general config of mpr.*

## 19.64 src/MPR/mo\_mpr\_restart.f90 File Reference

### Data Types

- interface [mo\\_mpr\\_restart::unpack\\_field\\_and\\_write](#)  
*TODO: add description.*

### Modules

- module [mo\\_mpr\\_restart](#)  
*reading and writing states, fluxes and configuration for restart of mHM.*

### Functions/Subroutines

- subroutine, public [mo\\_mpr\\_restart::write\\_mpr\\_restart\\_files](#) (OutFile)  
*write restart files for each domain*
- subroutine, public [mo\\_mpr\\_restart::write\\_eff\\_params](#) (mask1, s1, e1, rows1, cols1, soil1, lcscenes, lais, nc)  
*TODO: add description.*
- subroutine [mo\\_mpr\\_restart::unpack\\_field\\_and\\_write\\_1d\\_i4](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)

- subroutine [mo\\_mpr\\_restart::unpack\\_field\\_and\\_write\\_1d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [mo\\_mpr\\_restart::unpack\\_field\\_and\\_write\\_2d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)
- subroutine [mo\\_mpr\\_restart::unpack\\_field\\_and\\_write\\_3d\\_dp](#) (nc, var\_name, var\_dims, fill\_value, data, mask, var\_long\_name)

## 19.65 src/MPR/mo\_mpr\_runoff.f90 File Reference

### Modules

- module [mo\\_mpr\\_runoff](#)  
*multiscale parameter regionalization for runoff generation*

### Functions/Subroutines

- subroutine, public [mo\\_mpr\\_runoff::mpr\\_runoff](#) (LCOVER0, mask0, SMs\_FC0, slope\_emp0, KsVar\_H0, param, cell\_id0, upp\_row\_L1, low\_row\_L1, lef\_col\_L1, rig\_col\_L1, nL0\_in\_L1, L1\_HL1, L1\_K0, L1\_K1, L1\_alpha)  
*multiscale parameter regionalization for runoff parameters*

## 19.66 src/MPR/mo\_mpr\_smhorizons.f90 File Reference

### Modules

- module [mo\\_mpr\\_smhorizons](#)  
*setting up the soil moisture horizons*

### Functions/Subroutines

- subroutine, public [mo\\_mpr\\_smhorizons::mpr\\_smhorizons](#) (param, processMatrix, iFlag\_soil, nHorizons\_mHM, HorizonDepth, LCOVER0, soilID0, nHorizons, nTillHorizons, thetaS\_till, thetaFC\_till, thetaPW\_till, thetaS, thetaFC, thetaPW, Wd, Db, DbM, RZdepth, mask0, cell\_id0, upp\_row\_L1, low\_row\_L1, lef\_col\_L1, rig\_col\_L1, nL0\_in\_L1, L1\_beta, L1\_SMs, L1\_FC, L1\_PW, L1\_fRoots)  
*upscale soil moisture horizons*

## 19.67 src/MPR/mo\_mpr\_soilmoist.f90 File Reference

### Modules

- module [mo\\_mpr\\_soilmoist](#)  
*Multiscale parameter regionalization (MPR) for soil moisture.*

### Functions/Subroutines

- subroutine, public [mo\\_mpr\\_soilmoist::mpr\\_sm](#) (param, processMatrix, is\_present, nHorizons, nTillHorizons, sand, clay, DbM, ID0, soilID0, LCover0, thetaS\_till, thetaFC\_till, thetaPW\_till, thetaS, thetaFC, thetaPW, Ks, Db, KsVar\_H0, KsVar\_V0, SMs\_FC0)  
*multiscale parameter regionalization for soil moisture*
- elemental pure subroutine [mo\\_mpr\\_soilmoist::pwp](#) (Genu\_Mual\_n, Genu\_Mual\_alpha, thetaS, thetaPWP)  
*Permanent Wilting point.*
- elemental pure subroutine [mo\\_mpr\\_soilmoist::field\\_cap](#) (thetaFC, Ks, thetaS, Genu\_Mual\_n)  
*calculates the field capacity*
- subroutine [mo\\_mpr\\_soilmoist::genuchten](#) (thetaS, Genu\_Mual\_n, Genu\_Mual\_alpha, param, sand, clay, Db)

- calculates the Genuchten shape parameter*
- subroutine `mo_mpr_soilmoist::hydro_cond` (KS, param, sand, clay)
  - calculates the hydraulic conductivity Ks*

## 19.68 src/MPR/mo\_mpr\_startup.f90 File Reference

### Modules

- module `mo_mpr_startup`
  - Startup procedures for mHM.*

### Functions/Subroutines

- subroutine, public `mo_mpr_startup::mpr_initialize`
  - Initialize main mHM variables.*
- subroutine `mo_mpr_startup::l0_check_input` (iDomain)
  - Check for errors in L0 input data.*
- subroutine `mo_mpr_startup::l0_variable_init` (iDomain)
  - level 0 variable initialization*
- subroutine, public `mo_mpr_startup::init_eff_params` (ncells1)
  - Allocation of space for mHM related L1 and L11 variables.*

## 19.69 src/MPR/mo\_multi\_param\_reg.f90 File Reference

### Modules

- module `mo_multi_param_reg`
  - Multiscale parameter regionalization (MPR).*

### Functions/Subroutines

- subroutine, public `mo_multi_param_reg::mpr` (mask0, geoUnit0, soilld0, Asp0, gridded\_LAI0, LCover0, slope\_emp0, y0, ld0, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, n\_subcells1, fSealed1, alpha1, degDayInc1, degDayMax1, degDayNoPre1, fAsp1, HarSamCoeff1, PrieTayAlpha1, aeroResist1, surf↔Resist1, fRoots1, kFastFlow1, kSlowFlow1, kBaseFlow1, kPerco1, karstLoss1, soilMoistFC1, soilMoist↔Sat1, soilMoistExp1, jarvis\_thresh\_c1, tempThresh1, unsatThresh1, sealedThresh1, wiltingPoint1, max↔Inter1, petLAIcorFactor, parameterset)
  - Regionalizing and Upscaling process parameters.*
- subroutine `mo_multi_param_reg::baseflow_param` (param, geoUnit0, k2\_0)
  - baseflow recession parameter*
- subroutine `mo_multi_param_reg::snow_acc_melt_param` (param, fForest1, flperm1, fPerm1, tempThresh1, degDayNoPre1, degDayInc1, degDayMax1)
  - Calculates the snow parameters.*
- subroutine `mo_multi_param_reg::iper_thres_runoff` (param, sealedThresh1)
  - sets the impervious layer threshold parameter for runoff generation*
- subroutine `mo_multi_param_reg::karstic_layer` (param, geoUnit0, mask0, SMs\_FC0, KsVar\_V0, ld0, n↔subcells1, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, karstLoss1, L1\_Kp)
  - calculates the Karstic percolation loss*
- subroutine, public `mo_multi_param_reg::canopy_intercept_param` (processMatrix, param, LAI0, n\_subcells1, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, ld0, mask0, nodata, max\_intercept1)
  - estimate effective maximum interception capacity at L1*
- subroutine `mo_multi_param_reg::aerodynamical_resistance` (LAI0, LCover0, param, mask0, ld0, n↔subcells1, upper\_bound1, lower\_bound1, left\_bound1, right\_bound1, aerodyn\_resistance1)
  - Regionalization of aerodynamic resistance.*

## 19.70 src/MPR/mo\_prepare\_gridded\_lai.f90 File Reference

### Modules

- module [mo\\_prepare\\_gridded\\_lai](#)  
*Prepare daily LAI fields (e.g., MODIS data) for mHM.*

### Functions/Subroutines

- subroutine, public [mo\\_prepare\\_gridded\\_lai::prepare\\_gridded\\_daily\\_lai\\_data](#) (iDomain, nrows, ncols, mask, LAIPer\_iDomain)  
*Prepare gridded daily LAI data.*
- subroutine, public [mo\\_prepare\\_gridded\\_lai::prepare\\_gridded\\_mean\\_monthly\\_lai\\_data](#) (iDomain, nrows, ncols, mask)  
*prepare\_gridded\_mean\_monthly\_LAI\_data*

## 19.71 src/MPR/mo\_read\_lut.f90 File Reference

### Modules

- module [mo\\_read\\_lut](#)  
*Routines reading lookup tables (lut).*

### Functions/Subroutines

- subroutine, public [mo\\_read\\_lut::read\\_geoformation\\_lut](#) (filename, fileunit, nGeo, geo\_unit, geo\_karstic)  
*Reads LUT containing geological formation information.*
- subroutine, public [mo\\_read\\_lut::read\\_lai\\_lut](#) (filename, fileunit, nLAI, LAIIDlist, LAI)  
*Reads LUT containing LAI information.*

## 19.72 src/MPR/mo\_read\_wrapper.f90 File Reference

### Modules

- module [mo\\_read\\_wrapper](#)  
*Wrapper for all reading routines.*

### Functions/Subroutines

- subroutine, public [mo\\_read\\_wrapper::read\\_data](#) (LAIPer)  
*Reads data.*
- subroutine [mo\\_read\\_wrapper::check\\_consistency\\_lut\\_map](#) (data, lookuptable, filename, unique\_values)  
*Checks if classes in input maps appear in look up tables.*

## 19.73 src/MPR/mo\_soil\_database.f90 File Reference

### Modules

- module [mo\\_soil\\_database](#)  
*Generating soil database from input file.*

## Functions/Subroutines

- subroutine, public [mo\\_soil\\_database::read\\_soil\\_lut](#) (filename)  
*Reads the soil LUT file.*
- subroutine, public [mo\\_soil\\_database::generate\\_soil\\_database](#)  
*Generates soil database.*

## 19.74 src/MPR/mo\_upscaling\_operators.f90 File Reference

### Modules

- module [mo\\_upscaling\\_operators](#)  
*Module containing upscaling operators.*

### Functions/Subroutines

- integer(i4) function, dimension(size(l1\_upper\_rowld\_cell, 1)), public [mo\\_upscaling\\_operators::majority\\_statistics](#) (nClass, L1\_upper\_rowld\_cell, L1\_lower\_rowld\_cell, L1\_left\_colonld\_cell, L1\_right\_colonld\_cell, L0\_fine↔Scale\_2D\_data)  
*majority statistics*
- real(dp) function, dimension(size(l0upbound\_inlx, 1)), public [mo\\_upscaling\\_operators::l0\\_fractionalcover\\_in\\_lx](#) (dataIn0, classId, mask0, L0upBound\_inLx, L0downBound\_inLx, L0leftBound\_inLx, L0rightBound\_inLx, n↔TCells0\_inLx)  
*fractional coverage of a given class of L0 fields in Lx field (Lx = L1 or L11)*
- real(dp) function, dimension(size(nl0\_cells\_in\_l1\_cell, 1)), public [mo\\_upscaling\\_operators::upscale\\_arithmetic\\_mean](#) (nL0\_cells\_in\_L1\_cell, L1\_upper\_rowld\_cell, L1\_lower\_rowld\_cell, L1\_left\_colonld\_cell, L1\_right\_colonld\_↔cell, L0\_cellld, mask0, nodata\_value, L0\_fineScale\_data)  
*arithmetic mean*
- real(dp) function, dimension(size(nl0\_cells\_in\_l1\_cell, 1)), public [mo\\_upscaling\\_operators::upscale\\_harmonic\\_mean](#) (nL0\_cells\_in\_L1\_cell, L1\_upper\_rowld\_cell, L1\_lower\_rowld\_cell, L1\_left\_colonld\_cell, L1\_right\_colonld\_↔cell, L0\_cellld, mask0, nodata\_value, L0\_fineScale\_data)  
*harmonic mean*
- real(dp) function, dimension(size(l1\_upper\_rowld\_cell, 1)), public [mo\\_upscaling\\_operators::upscale\\_geometric\\_mean](#) (L1\_upper\_rowld\_cell, L1\_lower\_rowld\_cell, L1\_left\_colonld\_cell, L1\_right\_colonld\_cell, mask0, nodata\_↔value, L0\_fineScale\_data)  
*geometric mean*
- real(dp) function, dimension(size(nl0\_cells\_in\_l1\_cell, 1)) [mo\\_upscaling\\_operators::upscale\\_p\\_norm](#) (nL0\_↔cells\_in\_L1\_cell, L1\_upper\_rowld\_cell, L1\_lower\_rowld\_cell, L1\_left\_colonld\_cell, L1\_right\_colonld\_cell, L0\_cellld, mask0, nodata\_value, p\_norm, L0\_fineScale\_data)  
*arithmetic mean*

## 19.75 src/MPR/mpr\_driver.f90 File Reference

### Modules

- module [dummy\\_mpr](#)

### Functions/Subroutines

- program [mpr\\_driver](#)  
*Distributed precipitation-runoff model mHM.*

### 19.75.1 Function/Subroutine Documentation

## 19.75.1.1 mpr\_driver()

```
program mpr_driver
```

Distributed precipitation-runoff model mHM.

This is the main driver of mHM, which calls one instance of mHM for a multiple domains and a given period.

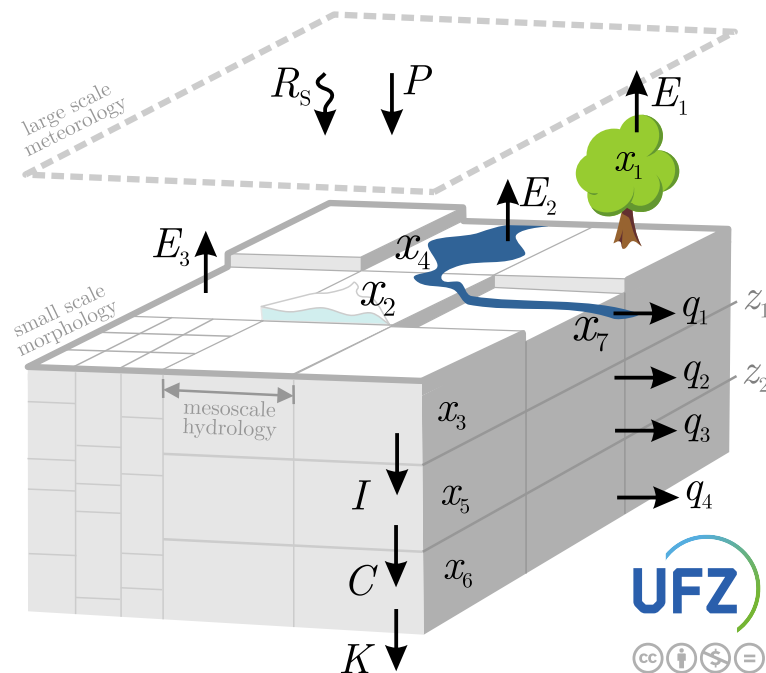


Figure 19.2: Typical mHM cell

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#### Version

0.1

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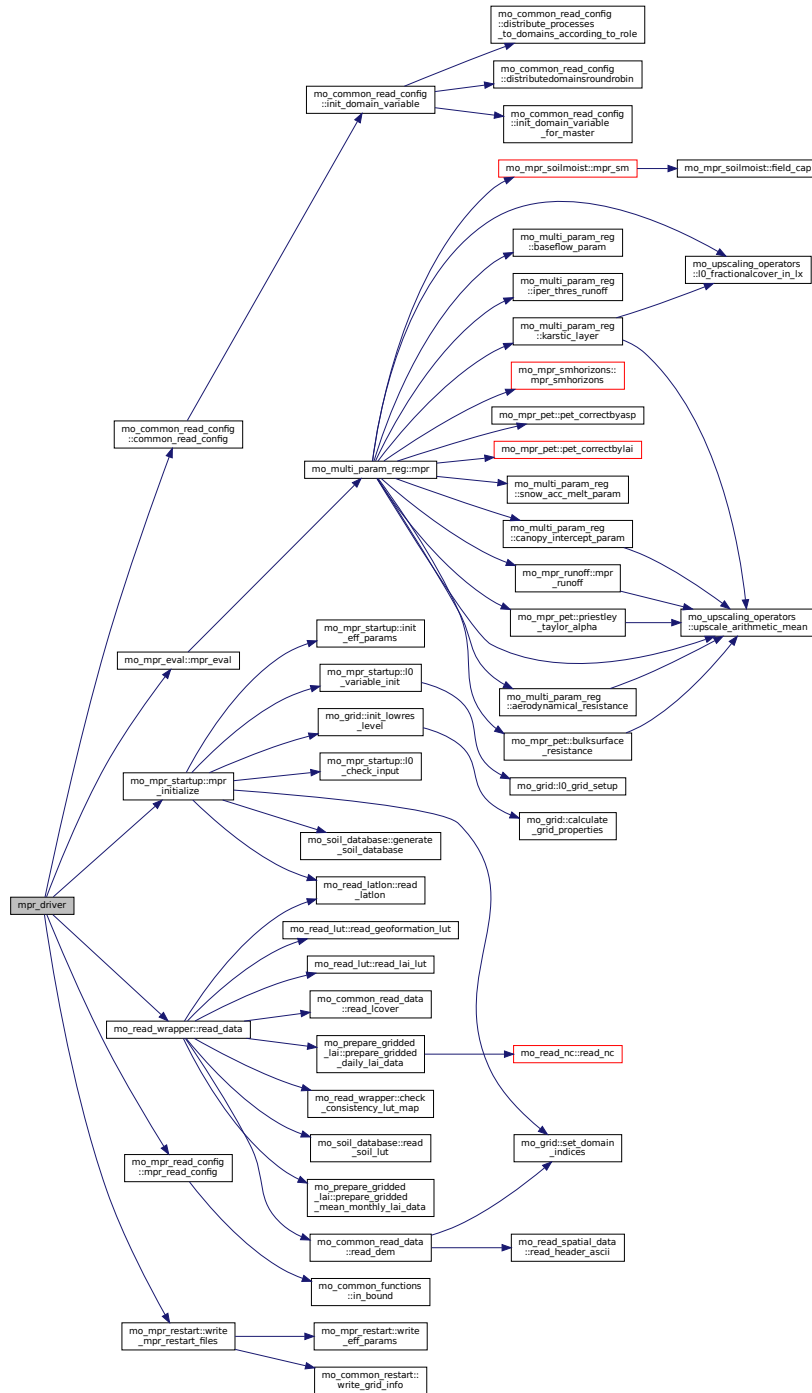
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Definition at line 59 of file mpr\_driver.f90.

References mo\_common\_read\_config::common\_read\_config(), mo\_mpr\_file::file\_namelist\_mpr\_param, mo\_common\_variables::mhmfilerestartout, mo\_mpr\_eval::mpr\_eval(), mo\_mpr\_startup::mpr\_initialize(), mo\_mpr\_read\_config::mpr\_read\_config(), mo\_read\_wrapper::read\_data(), mo\_mpr\_file::unamelist\_mpr\_param, mo\_mpr\_restart::write\_mpr\_restart\_files(), and mo\_common\_variables::write\_restart.

Here is the call graph for this function:





## 19.76 src/mRM/mo\_mrm\_constants.f90 File Reference

### Modules

- module [mo\\_mrm\\_constants](#)  
*Provides mRM specific constants.*

### Variables

- integer(i4), parameter, public [mo\\_mrm\\_constants::noutflxstate](#) = 2\_i4
- integer(i4), parameter, public [mo\\_mrm\\_constants::nroutingstates](#) = 2
- integer(i4), parameter, public [mo\\_mrm\\_constants::maxnogauges](#) = 50\_i4
- real(dp), parameter, public [mo\\_mrm\\_constants::rout\\_space\\_weight](#) = 0.\_dp
- real(dp), parameter, public [mo\\_mrm\\_constants::deltah](#) = 5.000\_dp
- real(dp), dimension(19), parameter [mo\\_mrm\\_constants::given\\_ts](#) = (/ 60.\_dp, 120.\_dp, 180.\_dp, 240.\_dp, 300.\_dp, 360.\_dp, 600.\_dp, 720.\_dp, 900.\_dp, 1200.\_dp, 1800.\_dp, 3600.\_dp, 7200.\_dp, 10800.\_dp, 14400.\_dp, 21600.\_dp, 28800.\_dp, 43200.\_dp, 86400.\_dp/)

## 19.77 src/mRM/mo\_mrm\_file.f90 File Reference

### Modules

- module [mo\\_mrm\\_file](#)  
*Provides file names and units for mRM.*

### Variables

- character(len=\*), parameter [mo\\_mrm\\_file::version](#) = '1.0'  
*Current mHM model version.*
- character(len=\*), parameter [mo\\_mrm\\_file::version\\_date](#) = 'May 2019'  
*Time of current mHM model version release.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_main](#) = 'mrm\_driver.f90'  
*Driver file.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_namelist\\_mrm](#) = 'mrm.nml'  
*Namelist file name.*
- integer, parameter [mo\\_mrm\\_file::unamelist\\_mrm](#) = 40  
*Unit for namelist.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_namelist\\_param\\_mrm](#) = 'mrm\_parameter.nml'  
*Parameter namelists file name.*
- integer, parameter [mo\\_mrm\\_file::unamelist\\_param\\_mrm](#) = 41  
*Unit for namelist.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_facc](#) = 'facc.asc'
- integer, parameter [mo\\_mrm\\_file::ufacc](#) = 56  
*Unit for flow accumulation input data file.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_fdir](#) = 'fdir.asc'  
*flow direction input data file*
- integer, parameter [mo\\_mrm\\_file::ufdir](#) = 57  
*Unit for flow direction input data file.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_slope](#) = 'slope.asc'  
*flow direction input data file*
- integer, parameter [mo\\_mrm\\_file::uslope](#) = 59  
*Unit for flow direction input data file.*
- character(len=\*), parameter [mo\\_mrm\\_file::file\\_gaugeloc](#) = 'idgauges.asc'

- gauge location input data file*

  - integer, parameter `mo_mrm_file::ugaugeloc` = 62  
*Unit for gauge location input data file.*
- integer, parameter `mo_mrm_file::udischarge` = 66  
*unit for discharge time series*
- character(len=\*), parameter `mo_mrm_file::file_defoutput` = 'mrm\_outputs.nml'  
*file defining mRM's outputs*
- integer, parameter `mo_mrm_file::udefoutput` = 67  
*Unit for file defining mRM's outputs.*
- character(len=\*), parameter `mo_mrm_file::file_config` = 'ConfigFile.log'  
*file defining mHM's outputs*
- integer, parameter `mo_mrm_file::uconfig` = 68  
*Unit for file defining mHM's outputs.*
- character(len=\*), parameter `mo_mrm_file::file_daily_discharge` = 'daily\_discharge.out'  
*file defining optimization outputs*
- integer, parameter `mo_mrm_file::udaily_discharge` = 74  
*Unit for file optimization outputs.*
- character(len=\*), parameter `mo_mrm_file::ncfile_discharge` = 'discharge.nc'  
*file defining optimization outputs*
- character(len=\*), parameter `mo_mrm_file::file_mrm_output` = 'mRM\_Fluxes\_States.nc'  
*file containing mrm output*
- character(len=\*), parameter `mo_mrm_file::file_gw_output` = 'mRM\_gw\_Fluxes\_States.nc'  
*file containing mrm output for groundwater coupling*

## 19.78 src/mRM/mo\_mrm\_global\_variables.f90 File Reference

### Data Types

- type `mo_mrm_global_variables::gaugingstation`
- type `mo_mrm_global_variables::domaininfo_mrm`

### Modules

- module `mo_mrm_global_variables`  
*Global variables for mRM only.*

### Variables

- logical `mo_mrm_global_variables::is_start`
- integer(i4) `mo_mrm_global_variables::output_deflate_level_mrm`
- logical `mo_mrm_global_variables::output_double_precision_mrm`
- integer(i4) `mo_mrm_global_variables::timestep_model_outputs_mrm`
- logical, dimension(noutflxstate) `mo_mrm_global_variables::outputflxstate_mrm`
- character(256), dimension(:), allocatable, public `mo_mrm_global_variables::dirgauges`
- character(256), dimension(:), allocatable, public `mo_mrm_global_variables::dirtotalrunoff`
- character(256), public `mo_mrm_global_variables::filenametotalrunoff`
- character(256), public `mo_mrm_global_variables::varnametotalrunoff`
- character(256), dimension(:), allocatable, public `mo_mrm_global_variables::dirbankfullrunoff`
- integer(i4), public `mo_mrm_global_variables::ntstepday`
- type(grid), dimension(:), allocatable, target, public `mo_mrm_global_variables::level11`
- type(gridremapper), dimension(:), allocatable, public `mo_mrm_global_variables::l0_l11_remap`
- type(gridremapper), dimension(:), allocatable, public `mo_mrm_global_variables::l1_l11_remap`
- real(dp), dimension(:, :), allocatable, public `mo_mrm_global_variables::mrm_runoff`

- integer(i4), public `mo_mrm_global_variables::ngaugestotal`
- integer(i4), public `mo_mrm_global_variables::ngaugeslocal`
- integer(i4), public `mo_mrm_global_variables::ninflowgaugestotal`
- integer(i4), public `mo_mrm_global_variables::nmeasperday`
- type(gaugingstation), public `mo_mrm_global_variables::gauge`
- type(gaugingstation), public `mo_mrm_global_variables::inflowgauge`
- type(domaininfo\_mrm), dimension(:), allocatable, target, public `mo_mrm_global_variables::domain_mrm`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_gaugeloc`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_inflowgaugeloc`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_facc`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_fdir`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_drasc`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_dracell`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_streamnet`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_floodplain`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l0_noutlet`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l0_celerity`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_l1_id`
- real(dp), dimension(:, :), allocatable, public `mo_mrm_global_variables::l1_total_runoff_in`
- integer(i4), dimension(:, :), allocatable, public `mo_mrm_global_variables::l11_cellcoor`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l1_l11_id`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_areacell`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_facc`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_fdir`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_noutlets`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_celerity`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_meandering`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_linkin_facc`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_rowout`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_colout`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_qmod`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_qout`
- real(dp), dimension(:, :), allocatable, public `mo_mrm_global_variables::l11_qtin`
- real(dp), dimension(:, :), allocatable, public `mo_mrm_global_variables::l11_qtr`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_fromn`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_ton`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_netperm`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_frow`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_fcol`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_trow`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_tcol`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_rorder`
- integer(i4), dimension(:), allocatable, public `mo_mrm_global_variables::l11_label`
- logical, dimension(:), allocatable, public `mo_mrm_global_variables::l11_sink`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_length`
- real(dp), dimension(:), allocatable, target, public `mo_mrm_global_variables::l11_afloodplain`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_slope`
- real(dp), dimension(:, :), allocatable, public `mo_mrm_global_variables::l11_nlinkfracpimp`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_k`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_xi`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_tsrout`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_c1`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_c2`
- logical `mo_mrm_global_variables::gw_coupling`
- real(dp), dimension(:), allocatable, public `mo_mrm_global_variables::l11_bankfull_runoff_in`

- real(dp), dimension(:), allocatable, public [mo\\_mrm\\_global\\_variables::l0\\_channel\\_depth](#)
- real(dp), dimension(:), allocatable, public [mo\\_mrm\\_global\\_variables::l0\\_channel\\_elevation](#)
- real(dp), dimension(:), allocatable, public [mo\\_mrm\\_global\\_variables::l0\\_river\\_head\\_mon\\_sum](#)
- real(dp), dimension(:), allocatable, public [mo\\_mrm\\_global\\_variables::l0\\_slope](#)
- type(riv\_temp\_type), public [mo\\_mrm\\_global\\_variables::riv\\_temp\\_pcs](#)

*This is a container for the river temperature routing process (pcs)*

## 19.79 src/mRM/mo\_mrm\_init.f90 File Reference

### Modules

- module [mo\\_mrm\\_init](#)  
*Wrapper for initializing Routing.*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_init::mrm\\_configuration](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param, ReadLatLon)
- subroutine, public [mo\\_mrm\\_init::mrm\\_init](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param, ReadLatLon)  
*Initialize all mRM variables at all levels (i.e., L0, L1, and L11).*
- subroutine [mo\\_mrm\\_init::print\\_startup\\_message](#) (file\_namelist, file\_namelist\_param)  
*TODO: add description.*
- subroutine [mo\\_mrm\\_init::config\\_output](#)  
*TODO: add description.*
- subroutine, public [mo\\_mrm\\_init::variables\\_default\\_init\\_routing](#)  
*Default initialization mRM related L11 variables.*
- subroutine [mo\\_mrm\\_init::l0\\_check\\_input\\_routing](#) (L0Domain\_iDomain)  
*TODO: add description.*
- subroutine [mo\\_mrm\\_init::variables\\_alloc\\_routing](#) (iDomain)  
*TODO: add description.*

## 19.80 src/mRM/mo\_mrm\_mpr.f90 File Reference

### Modules

- module [mo\\_mrm\\_mpr](#)  
*Perform Multiscale Parameter Regionalization on Routing Parameters.*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_mpr::reg\\_rout](#) (param, length, slope, fFPimp, TS, C1, C2)  
*Regionalized routing.*
- subroutine, public [mo\\_mrm\\_mpr::mrm\\_init\\_param](#) (iDomain, param)  
*TODO: add description.*
- subroutine, public [mo\\_mrm\\_mpr::mrm\\_update\\_param](#) (iDomain, param)  
*TODO: add description.*

## 19.81 src/mRM/mo\_mrm\_net\_startup.f90 File Reference

### Modules

- module [mo\\_mrm\\_net\\_startup](#)  
*Startup drainage network for mHM.*

## Functions/Subroutines

- subroutine, public `mo_mrm_net_startup::l11_l1_mapping` (iDomain)  
*TODO: add description.*
- subroutine, public `mo_mrm_net_startup::l11_flow_direction` (iDomain)  
*Determine the flow direction of the upscaled river network at level L11.*
- subroutine, public `mo_mrm_net_startup::l11_set_network_topology` (iDomain)  
*Set network topology.*
- subroutine, public `mo_mrm_net_startup::l11_routing_order` (iDomain)  
*Find routing order, headwater cells and sink.*
- subroutine, public `mo_mrm_net_startup::l11_link_location` (iDomain)  
*Estimate the LO (row,col) location for each routing link at level L11.*
- subroutine, public `mo_mrm_net_startup::l11_set_drain_outlet_gauges` (iDomain)  
*Draining cell identification and Set gauging node.*
- subroutine, public `mo_mrm_net_startup::l11_stream_features` (iDomain)  
*Stream features (stream network and floodplain)*
- subroutine, public `mo_mrm_net_startup::l11_fraction_sealed_floodplain` (LCClassImp, do\_init)  
*Fraction of the flood plain with impervious cover.*
- subroutine `mo_mrm_net_startup::moveup` (elev0, fDir0, fi, fj, ss, nn)  
*TODO: add description.*
- subroutine `mo_mrm_net_startup::movedownonecell` (fDir, iRow, jCol)  
*TODO: add description.*
- subroutine `mo_mrm_net_startup::celllength` (iDomain, fDir, iRow, jCol, iCoorSystem, length)  
*TODO: add description.*
- subroutine, public `mo_mrm_net_startup::get_distance_two_lat_lon_points` (lat1, long1, lat2, long2, distance\_out)  
*estimate distance in [m] between two points in a lat-lon*
- subroutine, public `mo_mrm_net_startup::l11_flow_accumulation` (iDomain)  
*Calculates L11 flow accumulation per grid cell.*
- recursive subroutine `calculate_l11_flow_accumulation` (fDir, fAcc, ii, jj, nrow, ncol)
- subroutine, public `mo_mrm_net_startup::l11_calc_celerity` (iDomain, param)  
*L11 celerity based on L0 elevation and L0 fAcc.*

### 19.81.1 Function/Subroutine Documentation

#### 19.81.1.1 calculate\_l11\_flow\_accumulation()

```
recursive subroutine l11_flow_accumulation::calculate_l11_flow_accumulation (
    integer(i4), dimension(:, :), intent(in) fDir,
    real(dp), dimension(:, :), intent(inout) fAcc,
    integer(i4), intent(in) ii,
    integer(i4), intent(in) jj,
    integer(i4), intent(in) nrow,
    integer(i4), intent(in) ncol )
```

Definition at line 2075 of file `mo_mrm_net_startup.f90`.

Referenced by `mo_mrm_net_startup::l11_flow_accumulation()`.

Here is the caller graph for this function:



## 19.82 src/mRM/mo\_mrm\_objective\_function\_runoff.f90 File Reference

### Modules

- module `mo_mrm_objective_function_runoff`  
*Objective Functions for Optimization of mHM/mRM against runoff.*

### Functions/Subroutines

- real(dp) function, public `mo_mrm_objective_function_runoff::single_objective_runoff` (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions optimizing against runoff.*
- real(dp) function, public `mo_mrm_objective_function_runoff::single_objective_runoff_master` (parameterset, eval, arg1, arg2, arg3)  
*Wrapper for objective functions optimizing against runoff.*
- subroutine, public `mo_mrm_objective_function_runoff::single_objective_runoff_subprocess` (eval, arg1, arg2, arg3)  
*Wrapper for objective functions optimizing against runoff.*
- subroutine, public `mo_mrm_objective_function_runoff::multi_objective_runoff` (parameterset, eval, multi\_↔ objectives)  
*Wrapper for multi-objective functions where at least one is regarding runoff.*
- real(dp) function `mo_mrm_objective_function_runoff::loglikelihood_stddev` (parameterset, eval, stddev, stddev\_new, likeli\_new)  
*Logarithmic likelihood function with removed linear trend and Lag(1)-autocorrelation.*
- real(dp) function `mo_mrm_objective_function_runoff::loglikelihood_evin2013_2` (parameterset, eval, regularize)  
*Logarithmised likelihood with linear error model and lag(1)-autocorrelation of the relative errors.*
- real(dp) function `mo_mrm_objective_function_runoff::parameter_regularization` (paraset, prior, bounds, mask)  
*TODO: add description.*
- real(dp) function `mo_mrm_objective_function_runoff::loglikelihood_trend_no_autocorr` (parameterset, eval, stddev\_old, stddev\_new, likeli\_new)  
*Logarithmic likelihood function with linear trend removed.*
- real(dp) function `mo_mrm_objective_function_runoff::objective_lnnse` (parameterset, eval)  
*Objective function of logarithmic NSE.*
- real(dp) function `mo_mrm_objective_function_runoff::objective_sse` (parameterset, eval)  
*Objective function of SSE.*
- real(dp) function `mo_mrm_objective_function_runoff::objective_nse` (parameterset, eval)  
*Objective function of NSE.*
- real(dp) function `mo_mrm_objective_function_runoff::objective_equal_nse_lnnse` (parameterset, eval)  
*Objective function equally weighting NSE and lnNSE.*
- real(dp) function, dimension(2) `mo_mrm_objective_function_runoff::multi_objective_nse_lnnse` (parameterset, eval)  
*Multi-objective function with NSE and lnNSE.*
- real(dp) function, dimension(2) `mo_mrm_objective_function_runoff::multi_objective_lnnse_highflow_lnnse_lowflow` (parameterset, eval)  
*Multi-objective function with NSE and lnNSE.*
- real(dp) function, dimension(2) `mo_mrm_objective_function_runoff::multi_objective_lnnse_highflow_lnnse_lowflow_2` (parameterset, eval)  
*Multi-objective function with NSE and lnNSE.*
- real(dp) function, dimension(2) `mo_mrm_objective_function_runoff::multi_objective_ae_fdc_lsv_nse_djf` (parameterset, eval)  
*Multi-objective function with absolute error of Flow Duration Curves low-segment volume and nse of DJF's discharge.*

- real(dp) function [mo\\_mrm\\_objective\\_function\\_runoff::objective\\_power6\\_nse\\_lnnse](#) (parameterset, eval)  
*Objective function of combined NSE and lnNSE with power of 5 i.e. the p-norm with p=5.*
- real(dp) function [mo\\_mrm\\_objective\\_function\\_runoff::objective\\_kge](#) (parameterset, eval)  
*Objective function of KGE.*
- real(dp) function [mo\\_mrm\\_objective\\_function\\_runoff::objective\\_multiple\\_gauges\\_kge\\_power6](#) (parameter-set, eval)  
*combined objective function based on KGE raised to the power 6*
- real(dp) function [mo\\_mrm\\_objective\\_function\\_runoff::objective\\_weighted\\_nse](#) (parameterset, eval)  
*Objective function of weighted NSE.*
- real(dp) function [mo\\_mrm\\_objective\\_function\\_runoff::objective\\_sse\\_boxcox](#) (parameterset, eval)  
*Objective function of sum of squared errors of transformed streamflow.*
- subroutine, public [mo\\_mrm\\_objective\\_function\\_runoff::extract\\_runoff](#) (gaugeld, runoff, runoff\_agg, runoff\_↔obs, runoff\_obs\_mask)  
*extracts runoff data from global variables*

## 19.83 src/mRM/mo\_mrm\_pre\_routing.f90 File Reference

### Modules

- module [mo\\_mrm\\_pre\\_routing](#)  
*Performs pre-processing for routing for mHM at level L11.*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_pre\\_routing::l11\\_runoff\\_acc](#) (qAll, efecArea, L1\_L11\_Id, L11\_areaCell, L11\_↔L1\_Id, TS, map\_flag, qAcc)  
*total runoff accumulation at L11.*
- subroutine, public [mo\\_mrm\\_pre\\_routing::add\\_inflow](#) (nInflowGauges, InflowIndexList, InflowHeadwater, InflowNodeList, Qinflow, qOut)  
*Adds inflow discharge to the runoff produced at the cell where the inflow is occurring.*
- subroutine [mo\\_mrm\\_pre\\_routing::l11\\_e\\_acc](#) (qAll, efecArea, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, TS, map\_flag, qAcc)  
*temperature energy accumulation at L11.*
- subroutine, public [mo\\_mrm\\_pre\\_routing::calc\\_l1\\_runoff\\_e](#) (fSealed\_area\_fraction, fast\_interflow, slow\_↔interflow, baseflow, direct\_runoff, temp\_air, mean\_temp\_air, lateral\_E)  
*calculate lateral temperature energy from runoff components.*
- subroutine, public [mo\\_mrm\\_pre\\_routing::l11\\_meteo\\_acc](#) (meteo\_all, efecArea, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, map\_flag, meteo\_acc)  
*meteo forcing accumulation at L11 for temperature routing.*

## 19.84 src/mRM/mo\_mrm\_read\_config.f90 File Reference

### Modules

- module [mo\\_mrm\\_read\\_config](#)  
*read mRM config*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_read\\_config::mrm\\_read\\_config](#) (file\_namelist, unamelist, file\_namelist\_param, unamelist\_param, do\_message, readLatLon)  
*Read the general config of mRM.*
- subroutine [mo\\_mrm\\_read\\_config::read\\_mrm\\_routing\\_params](#) (processCase, file\_namelist\_param, unamelist\_param)  
*TODO: add description.*

## 19.85 src/mRM/mo\_mrm\_read\_data.f90 File Reference

### Modules

- module [mo\\_mrm\\_read\\_data](#)  
*This module contains all routines to read mRM data from file.*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_read\\_data::mrm\\_read\\_l0\\_data](#) (do\_reinit, do\_readlatlon, do\_readlcover)  
*read L0 data from file*
- subroutine, public [mo\\_mrm\\_read\\_data::mrm\\_read\\_discharge](#)  
*Read discharge timeseries from file.*
- subroutine, public [mo\\_mrm\\_read\\_data::mrm\\_read\\_total\\_runoff](#) (iDomain)  
*read simulated runoff that is to be routed*
- subroutine, public [mo\\_mrm\\_read\\_data::mrm\\_read\\_bankfull\\_runoff](#) (iDomain)
- subroutine [mo\\_mrm\\_read\\_data::rotate\\_fdir\\_variable](#) (x)  
*TODO: add description.*

## 19.86 src/mRM/mo\_mrm\_restart.f90 File Reference

### Modules

- module [mo\\_mrm\\_restart](#)  
*Restart routines.*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_restart::mrm\\_write\\_restart](#) (iDomain, domainID, OutFile)  
*write routing states and configuration*
- subroutine, public [mo\\_mrm\\_restart::mrm\\_read\\_restart\\_states](#) (iDomain, domainID, InFile)  
*read routing states*
- subroutine, public [mo\\_mrm\\_restart::mrm\\_read\\_restart\\_config](#) (iDomain, domainID, InFile)  
*reads Level 11 configuration from a restart directory*

## 19.87 src/mRM/mo\_mrm\_riv\_temp\_class.f90 File Reference

### Data Types

- module [mo\\_mrm\\_riv\\_temp\\_class::riv\\_temp\\_type](#)  
*This is a container to define the river temperature routing in the current time step.*

### Modules

- module [mo\\_mrm\\_riv\\_temp\\_class](#)  
*Class for the river temperature calculations.*

### Functions/Subroutines

- subroutine [mo\\_mrm\\_riv\\_temp\\_class::config](#) (self, file\_namelist, unamelist, file\_namelist\_param, unamelist\_param)  
*configure the [riv\\_temp\\_type](#) class from the mhm namelist*
- subroutine [mo\\_mrm\\_riv\\_temp\\_class::init](#) (self, nCells)  
*initialize the [riv\\_temp\\_type](#) class for the current domain*



- subroutine `mo_mrm_riv_temp_class::init_area` (self, iDomain, L11\_netPerm, L11\_fromN, L11\_length, nLinks, nCells, nrows, ncols, L11\_mask)
  - initialize the river area of riv\_temp\_type class for the current domain*
- subroutine `mo_mrm_riv_temp_class::init_riv_temp` (self, time, ntimesteps\_day, temp\_air, read\_meteo↵ weights, temp\_weights, fday\_temp, fnight\_temp, efecarea, L1\_L11\_Id, L11\_areacell, L11\_L1\_Id, map\_flag)
  - initialize the river temperature of riv\_temp\_type class for the current domain*
- subroutine `mo_mrm_riv_temp_class::reset_timestep` (self)
  - reset riv\_temp\_type class for next timestep*
- subroutine `mo_mrm_riv_temp_class::alloc_lateral` (self, nCells)
  - allocate lateral temp components of riv\_temp\_type class for current domain*
- subroutine `mo_mrm_riv_temp_class::dealloc_lateral` (self)
  - deallocate lateral temp components of riv\_temp\_type*
- subroutine `mo_mrm_riv_temp_class::acc_source_e` (self, time, ntimesteps\_day, fSealed\_area\_fraction, fast\_interflow, slow\_interflow, baseflow, direct\_runoff, temp\_air, mean\_temp\_air, ssrd\_day, strd\_day, read\_meteo↵ weights, temp\_weights, fday\_temp, fnight\_temp, fday\_ssrd, fnight\_ssrd, fday\_strd, fnight\_strd)
  - accumulate energy sources of riv\_temp\_type*
- subroutine `mo_mrm_riv_temp_class::finalize_source_e` (self, efecarea, L1\_L11\_Id, L11\_areacell, L11\_L1\_↵ Id, timestep, map\_flag)
  - finalize energy sources of riv\_temp\_type*
- real(dp) function `mo_mrm_riv_temp_class::get_lrad_out` (self, riv\_temp)
  - get outgoing longwave radiation of riv\_temp\_type*
- real(dp) function `mo_mrm_riv_temp_class::get_lat_heat` (self, air\_temp, netrad)
  - latent heat flux of riv\_temp\_type*
- real(dp) function `mo_mrm_riv_temp_class::get_sens_heat` (self, air\_temp, riv\_temp)
  - sensible heat flux of riv\_temp\_type*
- real(dp) function `mo_mrm_riv_temp_class::get_e_io` (self, riv\_temp, cell)
  - get complete energy source of riv\_temp\_type at given cell*
- subroutine `mo_mrm_riv_temp_class::l11_routing_e` (self, nLinks, netPerm, netLink\_fromN, netLink\_toN, netLink\_C1, netLink\_C2, nInflowGauges, InflowHeadwater, InflowNodeList, L11\_qTR, L11\_Qmod)
  - execute the temperature routing of riv\_temp\_type*
- subroutine `mo_mrm_riv_temp_class::init_iter` (self)
  - initialize iterative solver of riv\_temp\_type*
- subroutine `mo_mrm_riv_temp_class::next_iter` (self, T\_est, T\_rout)
  - execute next iteration with iterative solver of riv\_temp\_type*

## 19.88 src/mRM/mo\_mrm\_river\_head.f90 File Reference

### Modules

- module `mo_mrm_river_head`

### Functions/Subroutines

- subroutine, public `mo_mrm_river_head::init_masked_zeros_I0` (iDomain, data)
- subroutine, private `mo_mrm_river_head::reset_sum` (iDomain, data)
- subroutine, public `mo_mrm_river_head::calc_channel_elevation` ()
- subroutine, public `mo_mrm_river_head::calc_river_head` (iDomain, L11\_Qmod, river\_head)
- real(dp) function `mo_mrm_river_head::calc_slope` (iDomain, elev0, fDir0, i, j)
- subroutine, public `mo_mrm_river_head::avg_and_write_timestep` (iDomain, timestep, data)
- subroutine, public `mo_mrm_river_head::create_output` (iDomain, OutPath)

## Variables

- type(ncvariable), dimension(:), allocatable [mo\\_mrm\\_river\\_head::nc\\_time](#)
- type(ncvariable), dimension(:), allocatable [mo\\_mrm\\_river\\_head::nc\\_riverhead](#)
- integer(i4), dimension(:), allocatable [mo\\_mrm\\_river\\_head::time\\_counter](#)
- integer(i4), dimension(:), allocatable [mo\\_mrm\\_river\\_head::sum\\_counter](#)

## 19.89 src/mRM/mo\_mrm\_routing.f90 File Reference

### Modules

- module [mo\\_mrm\\_routing](#)  
*Performs runoff routing for mHM at level L11.*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_routing::mrm\\_routing](#) (read\_states, processCase, global\_routing\_param, L1↔\_total\_runoff, L1\_areaCell, L1\_L11\_Id, L11\_areaCell, L11\_L1\_Id, L11\_netPerm, L11\_fromN, L11\_toN, L11\_nOutlets, timestep, tsRoutFactor, nNodes, nInflowGauges, InflowGaugeIndexList, InflowGauge↔Headwater, InflowGaugeNodeList, InflowDischarge, nGauges, gaugeIndexList, gaugeNodeList, map\_flag, L11\_length, L11\_slope, L11\_FracFPimp, L11\_C1, L11\_C2, L11\_qOut, L11\_qTIN, L11\_qTR, L11\_qMod, GaugeDischarge)  
*route water given runoff*
- subroutine [mo\\_mrm\\_routing::l11\\_routing](#) (nNodes, nLinks, netPerm, netLink\_fromN, netLink\_toN, netLink↔\_C1, netLink\_C2, netNode\_qOUT, nInflowGauges, InflowHeadwater, InflowNodeList, netNode\_qTIN, net↔Node\_qTR, netNode\_Qmod)  
*Performs runoff routing for mHM at L11 upscaled network ([Routing Network](#)).*

## 19.90 src/mRM/mo\_mrm\_signatures.f90 File Reference

### Modules

- module [mo\\_mrm\\_signatures](#)  
*Module with calculations for several hydrological signatures.*

### Functions/Subroutines

- real(dp) function, dimension(size(lags, 1)), public [mo\\_mrm\\_signatures::autocorrelation](#) (data, lags, mask)  
*Autocorrelation of a given data series.*
- real(dp) function, dimension(size(quantiles, 1)), public [mo\\_mrm\\_signatures::flowdurationcurve](#) (data, quantiles, mask, concavity\_index, mid\_segment\_slope, mhigh\_segment\_volume, high\_segment\_volume, low\_↔segment\_volume)  
*Flow duration curves.*
- subroutine, public [mo\\_mrm\\_signatures::limb\\_densities](#) (data, mask, RLD, DLD)  
*Calculates limb densities.*
- real(dp) function [mo\\_mrm\\_signatures::maximummonthlyflow](#) (data, mask, yr\_start, mo\_start, dy\_start)  
*Maximum of average flows per months.*
- subroutine, public [mo\\_mrm\\_signatures::moments](#) (data, mask, mean\_data, stddev\_data, median\_data, max\_data, mean\_log, stddev\_log, median\_log, max\_log)  
*Moments of data and log-transformed data, e.g. mean and standard deviation.*
- real(dp) function, dimension(size(quantiles, 1)), public [mo\\_mrm\\_signatures::peakdistribution](#) (data, quantiles, mask, slope\_peak\_distribution)  
*Calculates the peak distribution.*
- real(dp) function, public [mo\\_mrm\\_signatures::runoffratio](#) (data, domain\_area, mask, precip\_series, precip↔\_sum, log\_data)

*Runoff ratio (accumulated daily discharge [mm/d] / accumulated daily precipitation [mm/d]).*

- real(dp) function, public [mo\\_mrm\\_signatures::zeroflowratio](#) (data, mask)

*Ratio of zero values to total number of data points.*

## 19.91 src/mRM/mo\_mrm\_write.f90 File Reference

### Modules

- module [mo\\_mrm\\_write](#)  
*write of discharge and restart files*

### Functions/Subroutines

- subroutine, public [mo\\_mrm\\_write::mrm\\_write](#)  
*write discharge and restart files*
- subroutine [mo\\_mrm\\_write::write\\_configfile](#)  
*This modules writes the results of the configuration into an ASCII-file.*
- subroutine [mo\\_mrm\\_write::write\\_daily\\_obs\\_sim\\_discharge](#) (Qobs, Qsim)  
*Write a file for the daily observed and simulated discharge timeseries during the evaluation period for each gauging station.*
- subroutine, public [mo\\_mrm\\_write::mrm\\_write\\_output\\_fluxes](#) (iDomain, nCells, timeStep\_model\_outputs, domainDateTime, tt, timestep, mask11, L11\_qmod)  
*write fluxes to netcdf output files*
- subroutine, public [mo\\_mrm\\_write::mrm\\_write\\_optifile](#) (best\_OF, best\_paramSet, param\_names)  
*Write briefly final optimization results.*
- subroutine, public [mo\\_mrm\\_write::mrm\\_write\\_optinamelist](#) (parameters, maskpara, parameters\_name)  
*Write final, optimized parameter set in a namelist format.*

### Variables

- type(outputdataset) [mo\\_mrm\\_write::nc](#)

## 19.92 src/mRM/mo\_mrm\_write\_fluxes\_states.f90 File Reference

### Data Types

- interface [mo\\_mrm\\_write\\_fluxes\\_states::outputvariable](#)
- interface [mo\\_mrm\\_write\\_fluxes\\_states::outputdataset](#)

### Modules

- module [mo\\_mrm\\_write\\_fluxes\\_states](#)  
*Creates NetCDF output for different fluxes and state variables of mHM.*

### Functions/Subroutines

- type(outputvariable) function [mo\\_mrm\\_write\\_fluxes\\_states::newoutputvariable](#) (nc, name, dtype, dims, ncells, mask, avg)  
*Initialize OutputVariable.*
- subroutine [mo\\_mrm\\_write\\_fluxes\\_states::updatevariable](#) (self, data)  
*Update OutputVariable.*
- subroutine [mo\\_mrm\\_write\\_fluxes\\_states::writevariabletimestep](#) (self, timestep)  
*Write timestep to file.*
- type(outputdataset) function [mo\\_mrm\\_write\\_fluxes\\_states::newoutputdataset](#) (iDomain, mask, nCells)

*Initialize OutputDataset.*

- subroutine `mo_mrm_write_fluxes_states::updatedataset` (self, sidx, eidx, L11\_Qmod, L11\_riv\_temp)

*Update all variables.*

- subroutine `mo_mrm_write_fluxes_states::writetimestep` (self, timestep)

*Write all accumulated data.*

- subroutine `mo_mrm_write_fluxes_states::close` (self)

*Close the file.*

- type(ncdataset) function `mo_mrm_write_fluxes_states::createoutputfile` (iDomain)

*Create and initialize output file for X & Y coordinate system.*

- subroutine `mo_mrm_write_fluxes_states::writevariableattributes` (var, long\_name, unit)

*Write output variable attributes.*

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