



PART III: Beyond secondary creep: anisotropic flow laws and the theory of continuous diversity

- 1. Fabric and its evolution: Available anisotropic flow laws
- 2. Continuous diversity of polycrystalline ice masses: the most comprehensive theory to model induced anisotropy







1. Fabric and its evolution: Available anisotropic laws

microscopic	macroscopic	phemomenological
v. d. Veen & Whillans 1994 I dyn. anisotropic viscous power law for indiv. grains Azuma & Goto-Azuma 1996 H static anisotropic flow law no fabric evolution	Lliboutry 1993 static anisotropic flow law Meyssonnier & Philipp 1996 dynamic anisotropic flow law (transv. isotropic) based on VPSC and ODF (Orientation Distribution Function) implemented (simplified version) by Gagliardini & Meyssonnier 1999/2000	Morland & Staroszczyk 1998~2001 obtain evolving anisotropy from instantaneous states of deformation without explicit reference to fabric or grain size (! reversibility of anisotropy)
Duval, Castelnau et al. 1983 ~ 2005 VPSC (Visco-Plastic Self Contained) dyn. anisotr. linear flow law for individual grains	Svendsen/Gödert & Hutter 1996, 1998 S/SI dyn. anisotr. (transv. isotr./orthotr.) flow law based on ODF/ODF + indiv. grains Gillet-Chaulet et al. 2005 S stat. anisotr. (orthotr.) flow law based on ODF and parameters from physical µ-M models, designed for l.s. num. modeling	







1. Fabric and its evolution: Available anisotropic laws

H - (homogenization) models: based on averages of individual grains

Schmidt tensors: $\dot{S}_{ij} = 1/N \sum_{g=1}^{n} m_i^{(g)} c_j^{(g)}$

m_i(g): unit vector parallel to resolved shear stress in the basal plane

 $c_{_{_{i}}}^{\;\;(g)}$: unit vector parallel to the c-axis orientation

n: Number of grains g

S - (statistical) models: based on an Orientation Distribution Function

ODF: orientation density $f=f(\mathbf{x},t,\mathbf{n})$, \mathbf{n} vector of unit length in S^2

$$\int_{S^2} f(\mathbf{x}, t, \mathbf{n}) d^2 \mathbf{n} = 1$$

alignment/structure/anisotropy tensors: $\mathbf{A} := \int_{S^2} f(\mathbf{x}, t, \mathbf{n}) \ \mathbf{n} \otimes \mathbf{n} \ d^2\mathbf{n}$







The theory of mixtures with continuous diversity (MCD)

- has been developped by S. H. Faria from ~ 2001
- conforms to the principles of Rational Mechanics Modeling of Materials
- · is a thermodynamic theory
- is the most comprehensive theory to model heterogeneity, in particluar induced anisotropy
- · has many other applications

In the context of ice sheet modeling, MCD allows for the simultaneous modeling of

- texture evolution (rotation of c-axis)
- recrystallization, polygonization, recovery









Background to the MCD:

Single constituent continua:

5 scalar balance laws for independent primary fields $\rho(x,t)$, v(x,t), T(x,t)

general balance law:

$$\partial$$
 (*)/ ∂t + div [(*) \mathbf{v} + ϕ)] - $s = p$

(*): additive quantity, **v**: velocity, φ, s, p: flux, supply and production of (*)

div: divergence operator in Euclidean space E³

conservation equation: p=0







Multiconstituent continua (chemically reacting mixtures, granular media,...):

N discrete constituents, indexed by α , typically N \leq 3

5N balance laws for primary fields: $\rho_{x}(\mathbf{x},t)$ $\mathbf{v}_{x}(\mathbf{x},t)$, $T_{x}(\mathbf{x},t)$

$$\rho_{\alpha}(\mathbf{x},t) \mathbf{v}_{\alpha}(\mathbf{x},t), T_{\alpha}(\mathbf{x},t)$$

 $\alpha = 1, ...N$

General balance law:

$$\partial \binom{*}{\alpha} / \partial t + \text{div} \left[\binom{*}{\alpha} \mathbf{v}_{\alpha} + \phi_{\alpha} \right] - \mathbf{s}_{\alpha} = \mathbf{p}_{\alpha}$$

Non-conservation equations on constituent level: $p_{\alpha} \neq 0$

Mixture balance laws are derived from the constituent balance laws according to the Rational Mechanics Modeling of Materials approach (Truesdell's third metaphysical principle) and provide homogenization rules:

$$\sum_{\alpha=1}^{N} \rho_{\alpha} = \rho$$
 $\sum_{\alpha=1}^{N} \mathbf{T}_{\alpha} - \rho_{\alpha} \mathbf{u}_{\alpha} \otimes \mathbf{u}_{\alpha} = \mathbf{T}$ $\mathbf{u}_{\alpha} = \mathbf{v}_{\alpha} - \mathbf{v}$ diffusion velocity







Classical continuous mixtures

Countable set of constituents with individual primary variables,

e.g.
$$\rho_{\alpha}(\mathbf{x}, t)$$

Each constituent has countably many properties distinguishing it from other constituents

Nature shows us often the reverse situation:

Mixtures with continuous diversity

Infinitely many constituents with primary variables depending on the continuously varying species label

$$\rho^*(\mathbf{x}, t, \alpha)$$

$$\alpha$$
 in $A = [\alpha_{min}, \alpha_{max}]$ species assemblage

Constituents differ from each other only in very few properties (size, orientation, age, ...)



Polycrystalline ice







Mixtures with continuous diversity

Balance equations for primary fields (note: # does not increase with α) depend on position in i/ Euclidean space and ii/ Species space

$$\rho(\mathbf{x},t,\alpha), \mathbf{v}(\mathbf{x},t,\alpha), T(\mathbf{x},t,\alpha)$$

$$\alpha$$
 in $A = [\alpha_{\min}, \alpha_{\max}]$







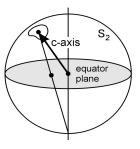
Mixtures with continuous diversity

Balance equations for primary fields (note: # does not increase with α) depend on position in i/ Euclidean space and ii/ Species space

$$\rho(\mathbf{x}, t, \alpha), \mathbf{v}(\mathbf{x}, t, \alpha), T(\mathbf{x}, t, \alpha)$$
 $\alpha \text{ in } A = [\alpha_{\min}, \alpha_{\max}]$

$$\alpha$$
 in $A = [\alpha_{\min}, \alpha_{\max}]$

Polycrystalline ice: species (single crystals) are identified by their orientation, represented by a unit normal vector ${\bf n}$ in S^2









Mixtures with continuous diversity

Balance equations for primary fields (note: # does not increase with α) depend on position in i/ Euclidean space and ii/ Species space

$$\rho(\mathbf{x},t,\alpha), \mathbf{v}(\mathbf{x},t,\alpha), \mathcal{T}(\mathbf{x},t,\alpha)$$

$$\alpha$$
 in $A = [\alpha_{\min}, \alpha_{\max}]$

Primary fields amended by dislocation density ρ_D and c-axis spin velocity **s**: $\rho(\mathbf{x},t,\mathbf{n}), \rho_D(\mathbf{x},t,\mathbf{n}), \mathbf{s}(\mathbf{x},t,\mathbf{n}), \mathbf{v}(\mathbf{x},t,\mathbf{n}), T(\mathbf{x},t,\mathbf{n})$

General Balance Equation

$$\partial$$
 (*)/ ∂t + div_{E³} [(*) \mathbf{v} + ϕ)] + div_{S²} [(*) \mathbf{w} + ψ] - s = p

w interspecies transition rate

ψ interspecies flux





Species balance equation for polycrystals modeled as mixtures with continuous diversity (Faria, 2006, Proc. R. Soc. Lond. A)

Balance of mass:

Balance of dislocation density:

Balance of linear momentum:

Balance of lattice spin velocity:

Balance of internal energy:

includes recrystallization includes interspecies flux density of dislocations and production rate of dislocations includes interspecies stress and high-angle interaction force includes polygonization tensor (interspecies couple stress) and high-angle interaction couple includes dissipative contributions associated with all new interspecies quantities







Homogenization of species balance equations:

explores Rational Mechanics Modeling of Materials approach, is of type

$$\rho = \int_{S^2} \rho(\mathbf{x}, t, \mathbf{n}) \, \mathrm{d}^2 \mathbf{n}$$

$$T = \int_{S^2} (T(x,t, n) - \rho(x,t, n) [v(x,t, n) - v(x,t)] \otimes [v(x,t, n) - v(x,t)]) d^2n$$

continuous mixtures:

$$\sum_{\alpha=1}^{N} \rho_{\alpha} = \rho \qquad \qquad \sum_{\alpha=1}^{N} \mathbf{T}_{\alpha} - \rho_{\alpha} \mathbf{u}_{\alpha} \otimes \mathbf{u}_{\alpha} = \mathbf{T}$$





Constitutive theory:

Work in progress

In Part III: Simplified reduced model

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Creep and recrystallization of large polycrystalline masses. II. Constitutive theory for crystalline media with transversely isotropic grains

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Creep and recrystallization of large polycrystalline masses. I. General continuum theory

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Creep and recryptallization of large polycrystalline masses. III. Continuum theory of ice sheets

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This work sets forth the first thermodynamically consistent constitutive theory for ice sheets undergoing strain-induced anisotropy, polygonization and recrystallization effects.







The simplified model presented in Faria's Part III is still more general than all other anisotropic flow laws.

T= -p 1 +
$$\mu^{(4)}$$
D^E with $\mu^{(4)} = \mu^{(4)}(\rho_D, \mathbf{n}, ...)$

It encompasses the previously suggested models by

- Svendsen/Gödert/Hutter
- Azuma/Goto-Azuma
- and the CAFFE model [Continuum mechancial Anisotropic Flow model based on an anisotropic Flow Enhancement factor] (cf. Placidi & Hutter, 2005, Seddik et al. 2008, Greve et al. (in print), Faria 2008)







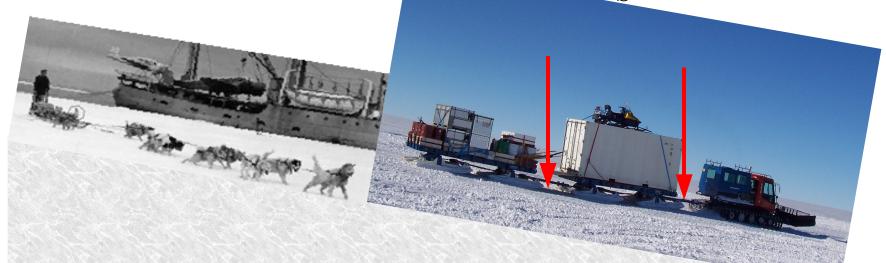
The flow law in the CAFFE model:

 $D = E(T^{D}, A^{(2)}, B^{(4)}) A(T) \sigma^{n-1} T^{D}$

$$\mathbf{A}^{(2)} = \int_{S^2} f(\mathbf{x}, t, \mathbf{n}) \, \mathbf{n} \times \mathbf{n} \, d^2 \mathbf{n}$$

$$\mathbf{B}^{(4)} = \int_{S^2} \mathbf{n} \times \mathbf{n} \, f(\mathbf{x}, t, \mathbf{n}) \, \mathbf{n} \times \mathbf{n} \, d^2 \mathbf{n}$$

$$\mathbf{B}^{(4)} = \int_{\mathbf{S}^2} \mathbf{n} \, \mathbf{r} \, \mathbf{n} \, f(\mathbf{x}, t, \mathbf{n}) \, \mathbf{n} \, \mathbf{x} \, \mathbf{n} \, \mathrm{d}^2 \mathbf{r}$$



The CAFFE model is implemented in Elmer/Ice @ CSC Finland (Th. Zwinger)