

Sample Name: U1553D 18R7

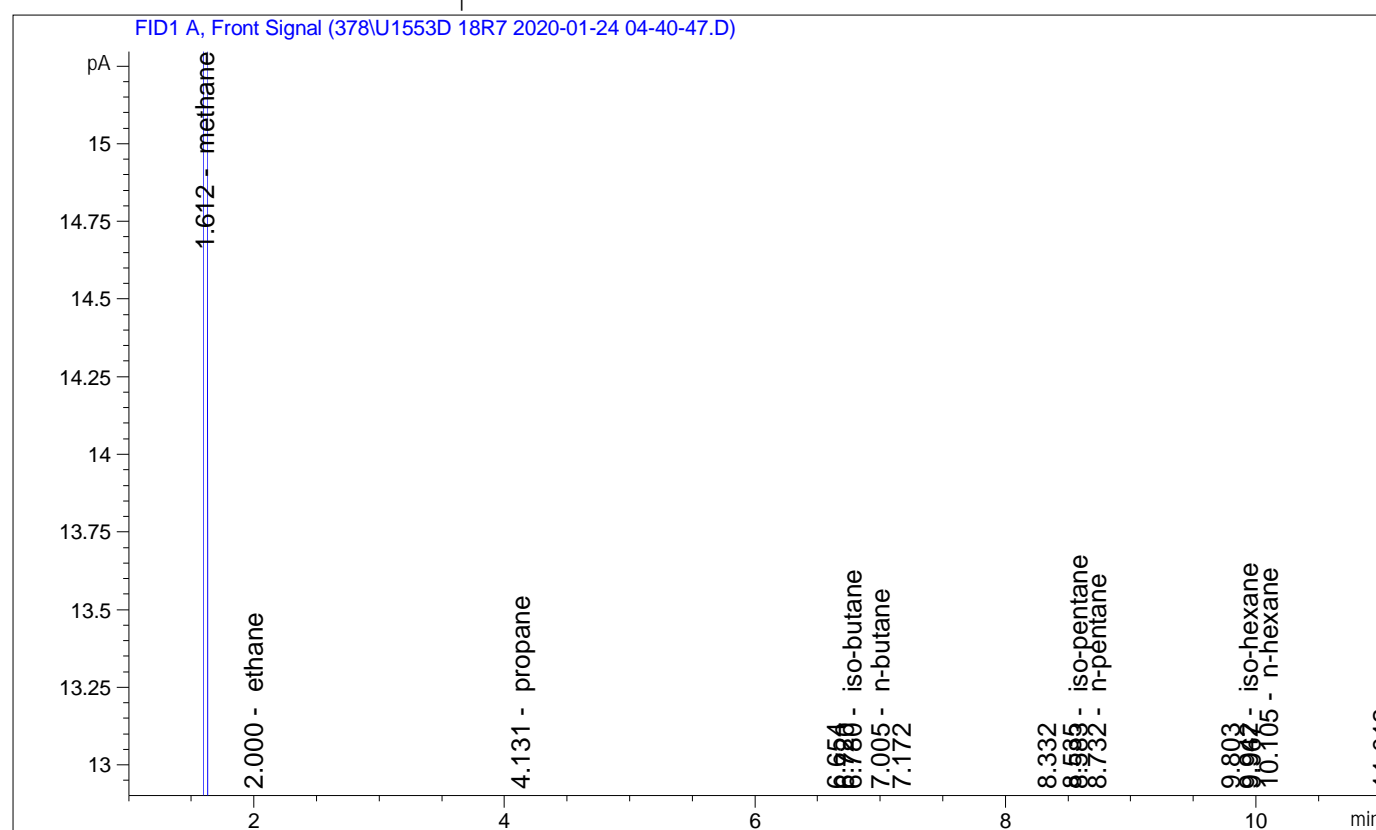
```

=====
Acq. Operator   : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : NGA-2                      Location :    1    (F)
Injection Date  : 1/24/2020 4:40:47 AM
                                           Inj Volume : Manual ly
Method          : C:\Chem32\2\Methods\378_NGA2_TCD0FF_012320.M
Last changed    : 1/24/2020 2:03:08 AM by SYSTEM
Method Info     : manual headspace injections

Sample Info     : othr10789141
  
```

Sample-related custom fields:

Name	Value
------	-------



External Standard Report

```

=====
Sorted By           :      Signal
Calib. Data Modified :      1/24/2020 2:01:48 AM
Multiplier          :      1.0000
Dilution            :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs
  
```

Signal 1: FID1 A, Front Signal

Sample Name: U1553D 18R7

RetTime [min]	Type	Area [pA*s]	Amt/Area	Amount [ppm]	Grp	Name
1.612	BB	61.09828	23.76552	1452.03273		methane
2.000	BB	1.73107	10.19543	17.64898		ethane
2.750		-	-	-		ethene
4.131	BB	2.70952	6.69894	18.15089		propane
6.381		-	-	-		propene
6.780	BV	5.26279e-1	1.96239	1.03277		i so-butane
7.005	BV	2.14193	4.60644	9.86670		n-butane
8.583	BB	1.70451	2.90168	4.94592		i so-pentane
8.732	BB	2.28428	2.63749	6.02478		n-pentane
9.942	BV	1.62231	0.00000	0.00000		i so-hexane
10.105	BB	1.53497	0.00000	0.00000		n-hexane

Totals : 1509.70277

4 Warnings or Errors :

Warning : Calibration warnings (see calibration table listing)

Warning : Calibrated compound(s) not found

Warning : Negative results set to zero (cal. curve intercept), (i so-hexane)

Warning : Negative results set to zero (cal. curve intercept), (n-hexane)

=====

Summed Peaks Report

=====

Signal 1: FID1 A, Front Signal

Empty table.

=====

Final Summed Peaks Report

=====

Signal 1: FID1 A, Front Signal

Name	Total Area [pA*s]	Amount [ppm]
------	----------------------	-----------------

methane	61.09828	1.452e3
ethane	1.73107	17.6490
ethene	0.00000	0.0000
propane	2.70952	18.1509
propene	0.00000	0.0000
i so-butane	5.26279e-1	1.0328
n-butane	2.14193	9.8667
i so-pentane	1.70451	4.9459
n-pentane	2.28428	6.0248
i so-hexane	1.62231	0.0000
n-hexane	1.53497	0.0000

Totals : 1509.7028

Compound-related custom fields:

*** End of Report ***