

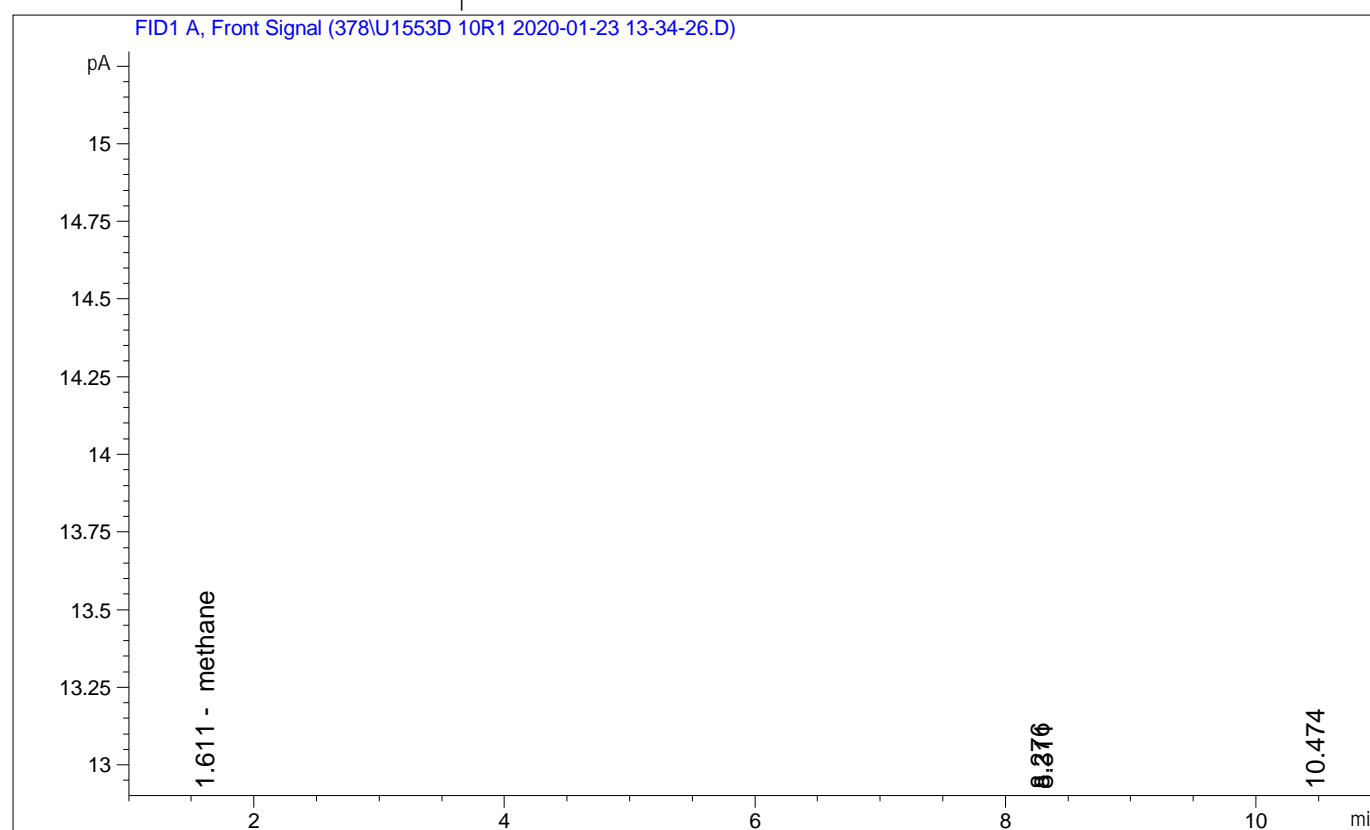
Sample Name: U1553D 10R1

```
=====
Acq. Operator   : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : NGA-2                      Location :    1   (F)
Injection Date  : 1/23/2020 1:34:26 PM
                                           Inj Volume : Manual ly
Method          : C:\Chem32\2\Methods\378_NGA2_TCD0FF_012320.M
Last changed    : 1/23/2020 9:30:40 AM by SYSTEM
Method Info     : manual headspace injections

Sample Info     : CYL10787381
=====
```

Sample-related custom fields:

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



```
=====
External Standard Report
=====
```

```
Sorted By      :      Signal
Calib. Data Modified :      1/22/2020 9:00:53 PM
Multiplier     :      1.0000
Dilution       :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs
```

Signal 1: FID1 A, Front Signal

Sample Name: U1553D 10R1

RetTime [min]	Type	Area [pA*s]	Amt/Area	Amount [ppm]	Grp	Name
1.611	VB	1.73206e-1	0.00000	0.00000		methane
1.999		-	-	-		ethane
2.755		-	-	-		ethene
4.125		-	-	-		propane
6.394		-	-	-		propene
6.779		-	-	-		i so-butane
7.005		-	-	-		n-butane
8.582		-	-	-		i so-pentane
8.731		-	-	-		n-pentane
9.941		-	-	-		i so-hexane
10.104		-	-	-		n-hexane

Totals : 0.00000

3 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), (methane)

=====

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	1.73206e-1	0.0000
ethane	0.00000	0.0000
ethene	0.00000	0.0000
propane	0.00000	0.0000
propene	0.00000	0.0000
i so-butane	0.00000	0.0000
n-butane	0.00000	0.0000
i so-pentane	0.00000	0.0000
n-pentane	0.00000	0.0000
i so-hexane	0.00000	0.0000
n-hexane	0.00000	0.0000

Totals : 0.0000

Compound-related custom fields:

*** End of Report ***