

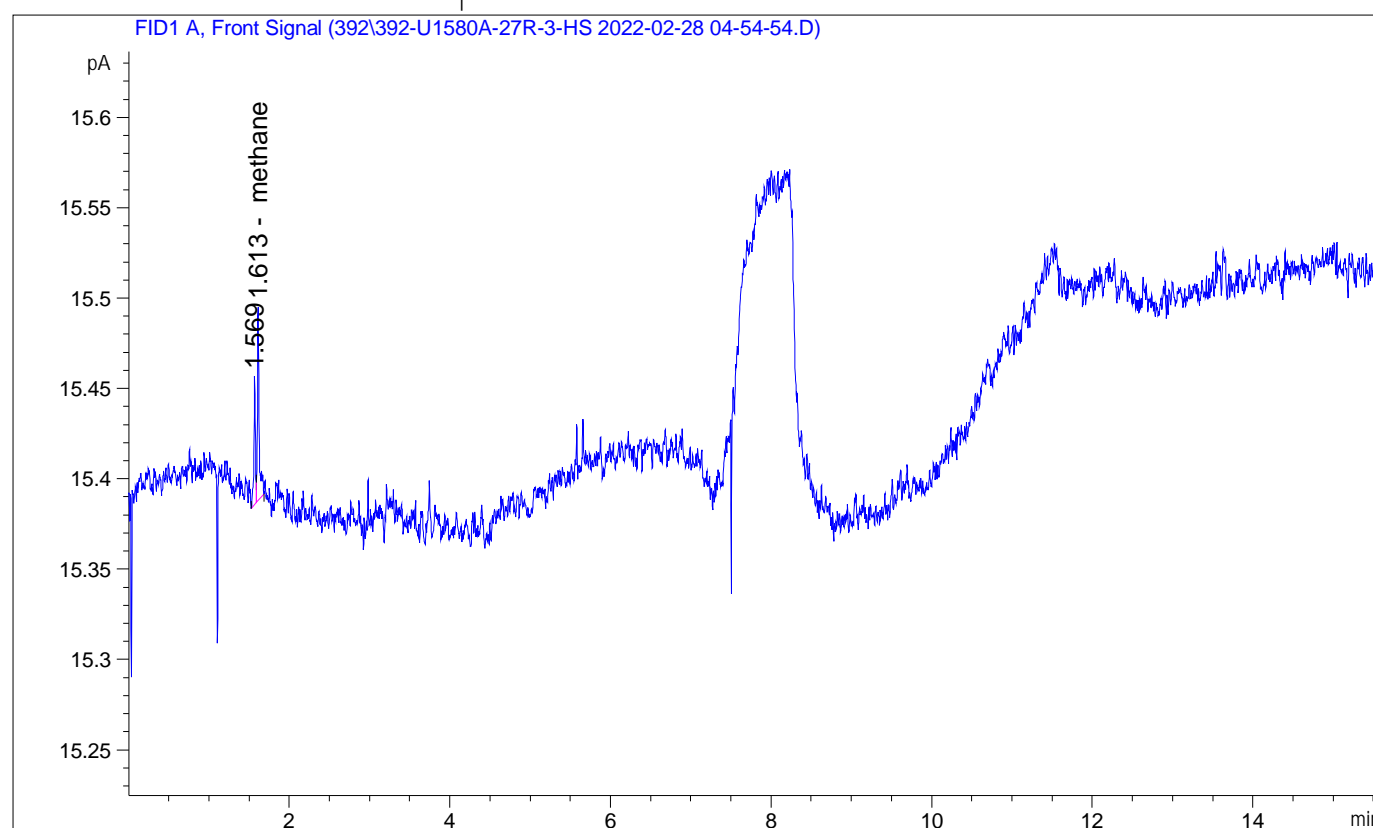
Sample Name: 392-U1580A-27R-3-HS

```
=====
Acq. Operator   : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : NGA-2                      Location :      - (F)
Injection Date  : 2/28/2022 4:54:54 AM
                                           Inj Volume : Manual ly
Method          : C:\Chem32\2\Methods\392_NGA2_TCD_UNCHECKED.M
Last changed    : 2/11/2022 4:53:36 PM by SYSTEM
Method Info     : manual headspace injections
=====
```

Sample Info : CYL11406481

Sample-related custom fields:

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



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

```
Sorted By      :      Signal
Calib. Data Modified :      2/10/2022 4:41:44 AM
Multiplier     :      1.0000
Dilution       :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs
```

Signal 1: FID1 A, Front Signal

RetTime [min]	Type	Area [pA*s]	Amt/Area	Amount [ppm]	Grp	Name
1.613	VB	1.60860e-1	0.00000	0.00000		methane
1.990		-	-	-		ethane
2.715		-	-	-		ethene
4.036		-	-	-		propane
6.361		-	-	-		propene
6.716		-	-	-		i so-butane
6.944		-	-	-		n-butane
8.516		-	-	-		i so-pentane
8.667		-	-	-		n-pentane
9.865		-	-	-		i so-hexane
10.029		-	-	-		n-hexane

Totals : 0.00000

Signal 2: TCD2 B, Back Signal not found

RetTime [min]	Type	Area [25 $\mu$ V*s]	Amt/Area	Amount [ppm]	Grp	Name
10.416		-	-	-		carbon dioxide

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.

Signal 2: TCD2 B, Back Signal not found

Empty table.

=====

#### Final Summed Peaks Report

=====

Signal 1: FID1 A, Front Signal

Name	Total Area [25 $\mu$ V*s]	Amount [ppm]
methane	1.60860e-1	0.0000
ethane	0.00000	0.0000

Name	Total Area [25 $\mu$ V*s]	Amount [ppm]
----- ----- -----		
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

Signal 2: TCD2 B, Back Signal not found

Name	Total Area [25 $\mu$ V*s]	Amount [ppm]
----- ----- -----		
carbon di oxide	0.00000	0.0000

Totals : 0.0000

Compound-related custom fields:

\*\*\* End of Report \*\*\*