

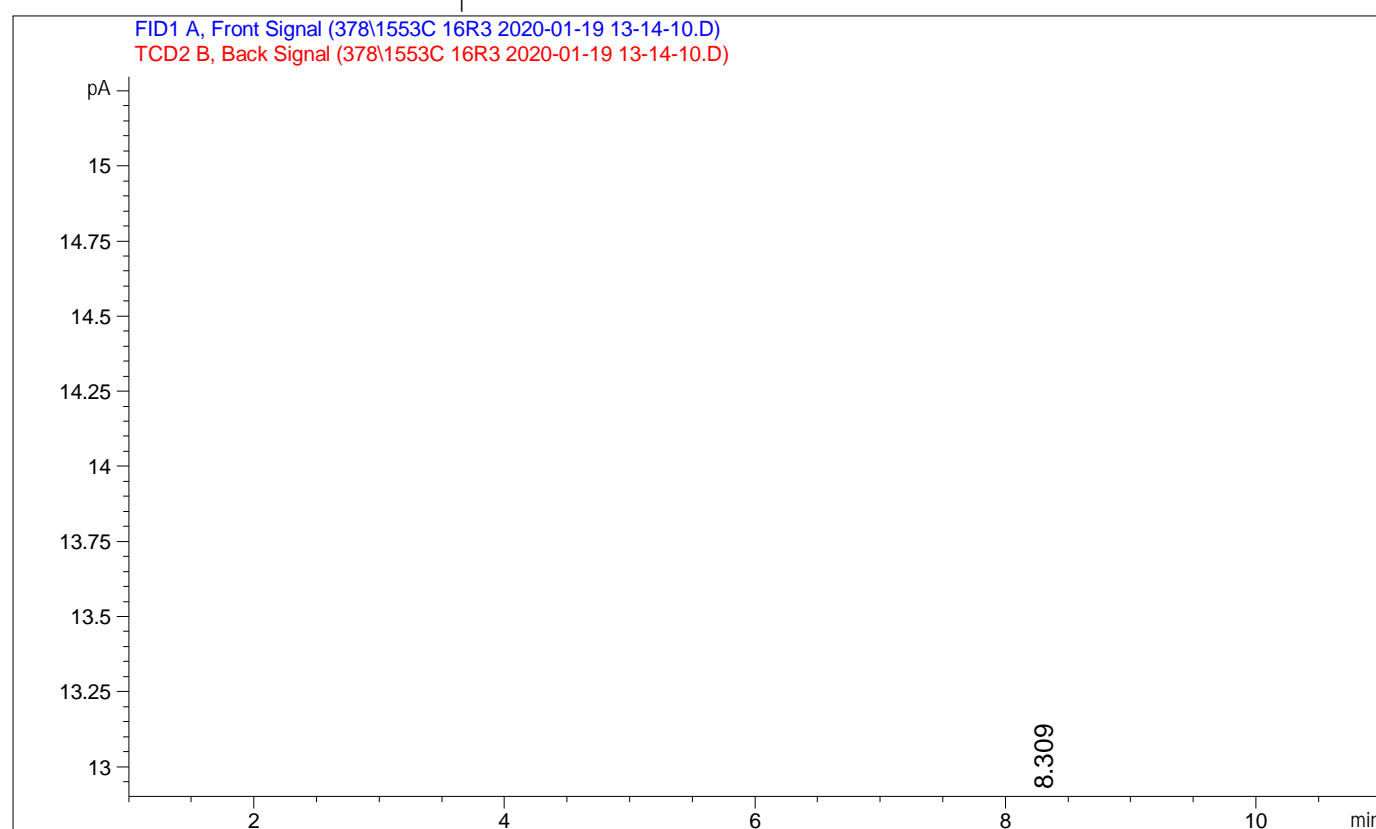
Sample Name: 1553C 16R3

```
=====
Acq. Operator   : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : NGA-2                      Location :    1  (F)
Injection Date  : 1/19/2020 1:14:10 PM
                                           Inj Volume : Manual ly
Method          : C:\Chem32\2\Methods\378T_NGA2_TCD0FF.M
Last changed    : 1/17/2020 4:53:59 AM by SYSTEM
                  (modified after loading)
Method Info     : manual headspace injections

Sample Info     : CYL10775701
=====
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Sample-related custom fields:

Name	Value
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```
=====
External Standard Report
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```

```
Sorted By      :      Signal
Calib. Data Modified :      1/15/2020 5:12:59 PM
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
1.610		-	-	-		methane
1.990		-	-	-		ethane
2.714		-	-	-		ethene
4.024		-	-	-		propane
6.340		-	-	-		propene
6.736		-	-	-		i so-butane
6.967		-	-	-		n-butane
8.544		-	-	-		i so-pentane
8.697		-	-	-		n-pentane
9.902		-	-	-		i so-hexane
10.068		-	-	-		n-hexane

Totals : 0.00000

Signal 2: TCD2 B, Back Signal

RetTime [min]	Type	Area [25 µV*s]	Amt/Area	Amount [ppm]	Grp	Name
4.598		-	-	-		O
5.414		-	-	-		N

Totals : 0.00000

2 Warnings or Errors :

Warning : Calibration warnings (see calibration table listing)  
Warning : Time reference compound(s) not found

```
=====
Summed Peaks Report
=====
```

Signal 1: FID1 A, Front Signal  
Empty table.

Signal 2: TCD2 B, Back Signal  
Empty table.

```
=====
Final Summed Peaks Report
=====
```

Signal 1: FID1 A, Front Signal

Sample Name: 1553C 16R3

Name	Total Area [25 $\mu$ V*s]	Amount [ppm]
methane	0.00000	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

Signal 2: TCD2 B, Back Signal

Name	Total Area [25 $\mu$ V*s]	Amount [ppm]
O	0.00000	0.0000
N	0.00000	0.0000

Totals : 0.0000

=====  
Area Percent Report  
=====

Sorted By : Signal  
 Calib. Data Modified : 1/15/2020 5:12:59 PM  
 Multiplier : 1.0000  
 Dilution : 1.0000  
 Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: FID1 A, Front Signal

Peak #	RetTime [min]	Type	Width [min]	Area [pA*s]	Area %	Name
1	1.610	+	0.0000	0.00000	0.00000	methane
2	1.990	+	0.0000	0.00000	0.00000	ethane
3	2.714	+	0.0000	0.00000	0.00000	ethene
4	4.024	+	0.0000	0.00000	0.00000	propane
5	6.340	+	0.0000	0.00000	0.00000	propene
6	6.736	+	0.0000	0.00000	0.00000	i so-butane
7	6.967	+	0.0000	0.00000	0.00000	n-butane
8	8.544	+	0.0000	0.00000	0.00000	i so-pentane
9	8.697	+	0.0000	0.00000	0.00000	n-pentane
10	9.902	+	0.0000	0.00000	0.00000	i so-hexane
11	10.068	+	0.0000	0.00000	0.00000	n-hexane

Totals : 0.00000 0.0000

Signal 2: TCD2 B, Back Signal

Peak #	RetTime [min]	Type	Width [min]	Area [25 $\mu$ V*s]	Area %	Name
1	4.598		0.0000	0.00000	0.00000	O
2	5.414		0.0000	0.00000	0.00000	N

Totals : 0.00000 0.0000

2 Warnings or Errors :

Warning : Calibration warnings (see calibration table listing)

Warning : Time reference compound(s) not found

=====  
Summed Peaks Report  
=====

Signal 1: FID1 A, Front Signal  
Empty table.

Signal 2: TCD2 B, Back Signal  
Empty table.

=====  
Final Summed Peaks Report  
=====

Signal 1: FID1 A, Front Signal

Name	Total Area [25 $\mu$ V*s]	Area %
methane	0.00000	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

Signal 2: TCD2 B, Back Signal

Name	Total Area [25 $\mu$ V*s]	Area %
O	0.00000	0.0000
N	0.00000	0.0000

Name	Total Area [25 $\mu$ V*s]	Area %
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Totals :		0.0000

Compound-related custom fields:

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