

Sample Name: 396-U1565B-1R-3-HS

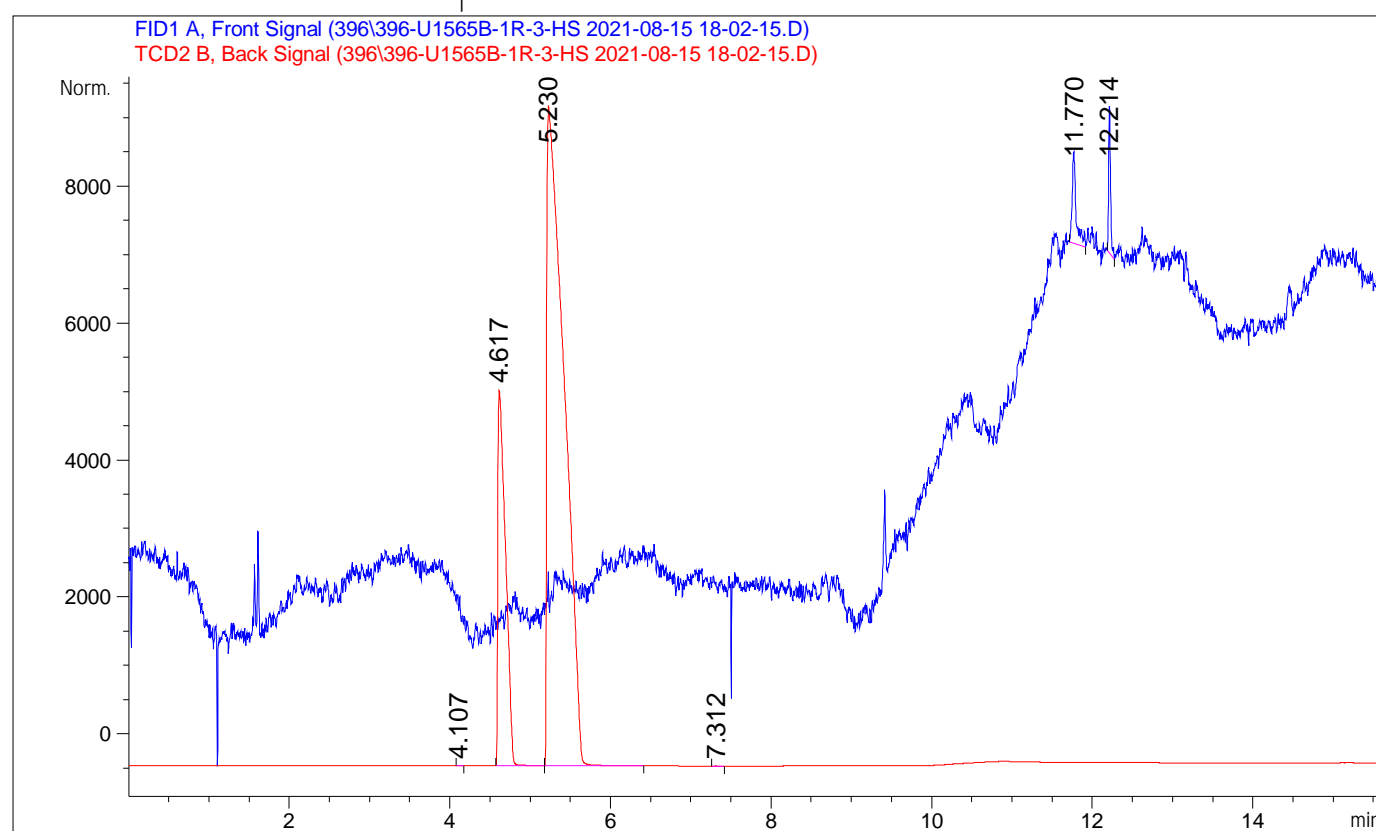
```
=====
Acq. Operator   : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : NGA-2                      Location :      - (F)
Injection Date  : 8/15/2021 6:02:15 PM
                                           Inj Volume : Manual ly

Method          : C:\CHEM32\2\METHODS\396_NGA2.M
Last changed    : 8/15/2021 3:52:08 AM by SYSTEM
                  (modified after loading)
Method Info     : manual headspace injections

Sample Info     : CYL11054471
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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 :      8/11/2021 10:16:43 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
1.614	-	-	-	-	-	methane
1.990	-	-	-	-	-	ethane
2.731	-	-	-	-	-	ethene
3.989	-	-	-	-	-	propane
6.355	-	-	-	-	-	propene
6.717	-	-	-	-	-	i so-butane
6.947	-	-	-	-	-	n-butane
8.519	-	-	-	-	-	i so-pentane
8.672	-	-	-	-	-	n-pentane
9.871	-	-	-	-	-	i so-hexane
10.036	-	-	-	-	-	n-hexane

Totals : 0.00000

Signal 2: TCD2 B, Back Signal

2 Warnings or Errors :

Warning : Calibration warnings (see calibration table listing)

Warning : Calibrated 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
=====
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Signal 1: FID1 A, Front Signal

Name	Total Area [25 μ 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

Name	Total Area [25 μ V*s]	Amount [ppm]
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

Area Percent Report

Sorted By : Signal
Calib. Data Modified : 8/11/2021 10:16:43 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	1.614		0.0000	0.00000	0.00000	methane
2	1.990		0.0000	0.00000	0.00000	ethane
3	2.731		0.0000	0.00000	0.00000	ethene
4	3.989		0.0000	0.00000	0.00000	propane
5	6.355		0.0000	0.00000	0.00000	propene
6	6.717		0.0000	0.00000	0.00000	i so-butane
7	6.947		0.0000	0.00000	0.00000	n-butane
8	8.519		0.0000	0.00000	0.00000	i so-pentane
9	8.672		0.0000	0.00000	0.00000	n-pentane
10	9.871		0.0000	0.00000	0.00000	i so-hexane
11	10.036		0.0000	0.00000	0.00000	n-hexane

Totals : 0.00000 0.0000

Signal 2: TCD2 B, Back Signal

2 Warnings or Errors :

Warning : Calibration warnings (see calibration table listing)
Warning : Calibrated 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

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Signal 1: FID1 A, Front Signal

Name	Total Area [25 μ 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

Compound-related custom fields:

*** End of Report ***