

Sample Name: 396-U1574A-1R-CC-HS

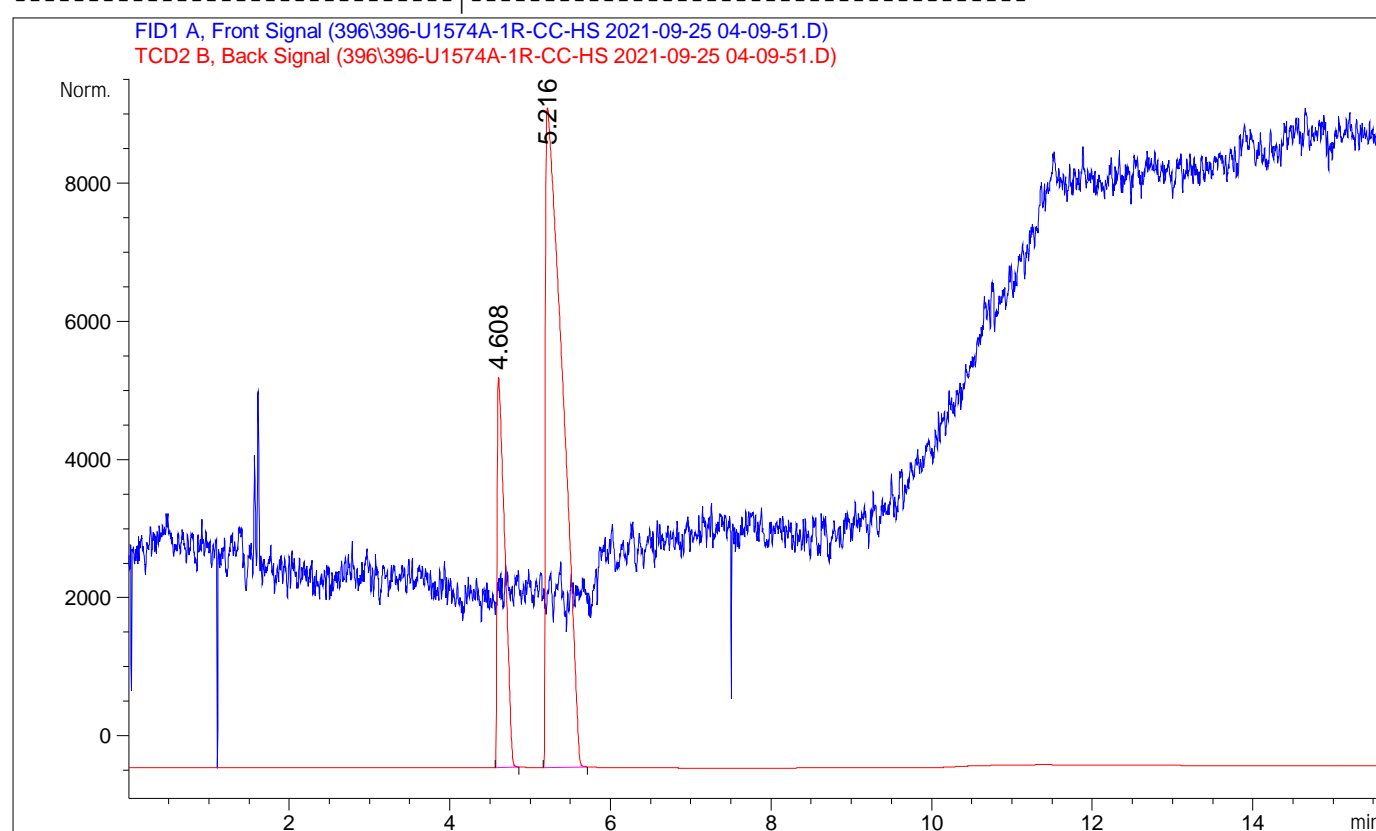
```
=====
Acq. Operator   : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : NGA-2                      Location :      - (F)
Injection Date  : 9/25/2021 4:09:51 AM
                                           Inj Volume : Manual ly

Method          : C:\CHEM32\2\METHODS\396_NGA2.M
Last changed    : 9/24/2021 11:33:14 PM by SYSTEM
                  (modified after loading)
Method Info     : manual headspace injections

Sample Info     : CYL11216811
=====
```

Sample-related custom fields:

Name	Value
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External Standard Report
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Sorted By      :      Signal
Calib. Data Modified : 9/24/2021 11:33:10 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
1.611	-	-	-	-	-	methane
1.963	-	-	-	-	-	ethane
2.570	-	-	-	-	-	ethene
3.794	-	-	-	-	-	propane
6.153	-	-	-	-	-	propene
6.563	-	-	-	-	-	i so-butane
6.803	-	-	-	-	-	n-butane
8.394	-	-	-	-	-	i so-pentane
8.551	-	-	-	-	-	n-pentane
9.742	-	-	-	-	-	i so-hexane
9.918	-	-	-	-	-	n-hexane

Totals : 0.00000

Signal 2: TCD2 B, Back Signal

RetTime [min]	Type	Area [25 μ V*s]	Amt/Area	Amount [ppm]	Grp	Name
10.540	-	-	-	-	-	carbon di oxide

Totals : 0.00000

2 Warnings or Errors :

Warning : Calibration warnings (see calibration table listing)

Warning : Calibrated compound(s) not found

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Summed Peaks Report

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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

Sample Name: 396-U1574A-1R-CC-HS

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
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 μ V*s]	Amount [ppm]
carbon di oxide	0.00000	0.0000

Totals : 0.0000

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

Sorted By : Signal
 Calib. Data Modified : 9/24/2021 11:33:10 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	1.611		0.0000	0.00000	0.00000	methane
2	1.963		0.0000	0.00000	0.00000	ethane
3	2.570		0.0000	0.00000	0.00000	ethene
4	3.794		0.0000	0.00000	0.00000	propane
5	6.153		0.0000	0.00000	0.00000	propene
6	6.563		0.0000	0.00000	0.00000	i so-butane
7	6.803		0.0000	0.00000	0.00000	n-butane
8	8.394		0.0000	0.00000	0.00000	i so-pentane
9	8.551		0.0000	0.00000	0.00000	n-pentane
10	9.742		0.0000	0.00000	0.00000	i so-hexane
11	9.918		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 µV*s]	Area %	Name
1	10.540		0.0000	0.00000	0.00000	carbon di oxide

Totals : 0.00000 0.0000

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

Signal 1: FID1 A, Front Signal

Name	Total Area [25 µV*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 µV*s]	Area %
carbon di oxide	0.00000	0.0000

Totals : 0.0000

Compound-related custom fields:

*** End of Report ***