LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.372	3	10	29	rVB	1971520	1985124	99.61%	11.409%
2	1.507	41	52	74	rVB2	140635	190070	9.54%	1.092%
3	1.610	75	84	95	rBV3	839005	1470552	73.79%	8.452%
4	1.678	101	105	109	rBV	29193	26864	1.35%	0.154%
5	1.707	109	114	121	rVB	52224	56760	2.85%	0.326%
6	1.752	121	128	135	rVB	124349	131142	6.58%	0.754%
7	1.932	175	184	199	rVB2	155586	236958	11.89%	1.362%
8	2.015	202	210	220	rVB	127116	170397	8.55%	0.979%
9	2.179	252	261	269	rBV	190671	260798	13.09%	1.499%
10	2.231	269	277	301	rVB3	200686	422633	21.21%	2.429%
11	2.343	306	312	325	rVB	22610	33171	1.66%	0.191%
12	2.443	332	343	352	rBV2	72955	122986	6.17%	0.707%
13	2.494	353	359	370	rVB2	23093	36499	1.83%	0.210%
14	2.597	378	391	404	rBV	226117	389606	19.55%	2.239%
15	2.668	405	413	429	rVB3	43734	81738	4.10%	0.470%
16	2.890	474	482	492	rVB2	33319	53075	2.66%	0.305%
17	3.009	507	519	536	rBV6	20372	53610	2.69%	0.308%
18	3.115	544	552	561	rVV	17595	31195	1.57%	0.179%
19	3.170	562	569	582	rVB5	11328	23824	1.20%	0.137%
20	3.427	618	649	673	rBV2	84197	276161	13.86%	1.587%
21	3.629	696	712	729	rVB5	58861	152495	7.65%	0.876%
22	3.739	734	746	765	rVB2	33613	73257	3.68%	0.421%
23	3.871	770	787	799	rBV4	13888	35212	1.77%	0.202%
24	3.951	802	812	826	rVB6	8750	22842	1.15%	0.131%
25	4.243	886	903	915	rBV4	18839	48170	2.42%	0.277%
26	4.514	970	987	1000	rBV3	17190	43279	2.17%	0.249%
27	4.716	1025	1050	1067	rBV2	333212	920387	46.18%	5.290%
28	5.115	1154	1174	1197	rBV	184189	445696	22.36%	2.562%
29	5.237	1198	1212	1223	rBV3	9487	22326	1.12%	0.128%
30	5.488	1277	1290	1300	rBV10	9055	22244	1.12%	0.128%
31	5.629	1321	1334	1351	rVB4	12109	25225	1.27%	0.145%
32	5.771	1351	1378	1407	rBV2	379681	1010380	50.70%	5.807%
33	6.009	1441	1452	1460	rBV4	11803	23742	1.19%	0.136%
34	6.067	1461	1470	1489	rVB4	20110	49803	2.50%	0.286%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Title : TRACE VOA SOM01.0

35	6.173	1489	1503	1513	rBV4	13673	29370	1.47%	0.169%
36	6.292	1526	1540	1563	rBV	344670	691083	34.68%	3.972%
37	6.584	1615	1631	1653	rBV	1032634	1992894	100.00%	11.454%
38	6.735	1665	1678	1701	rVB	235925	466146	23.39%	2.679%
39	6.931	1731	1739	1754	rVB3	20363	35202	1.77%	0.202%
40	7.607	1931	1949	1967	rBV	117470	226825	11.38%	1.304%
41	7.935	2034	2051	2064	rBV	442621	784227	39.35%	4.507%
42	8.005	2065	2073	2082	rVB	33752	58010	2.91%	0.333%
43	8.218	2127	2139	2158	rBV	67620	129448	6.50%	0.744%
44	8.584	2241	2253	2269	rVB2	61069	104153	5.23%	0.599%
45	8.690	2270	2286	2319	rBV	342189	1023556	51.36%	5.883%
46	8.832	2321	2330	2362	rVB3	71257	185188	9.29%	1.064%
47	9.449	2508	2522	2559	rBV	457642	852676	42.79%	4.901%
48	9.722	2593	2607	2618	rBV3	11623	23202	1.16%	0.133%
49	10.391	2800	2815	2836	rBV5	27453	62244	3.12%	0.358%
50	10.787	2926	2938	2965	rVB	175753	311651	15.64%	1.791%
51	11.729	3217	3231	3243	rBV2	50558	90033	4.52%	0.517%
52	11.844	3254	3267	3286	rVB	350683	655386	32.89%	3.767%
53	12.224	3369	3385	3405	rVB	332302	607528	30.48%	3.492%
54	12.922	3591	3602	3618	rBV2	56637	92274	4.63%	0.530%
55	14.092	3957	3966	3980	rBV3	16795	30196	1.52%	0.174%

Sum of corrected areas: 17399513

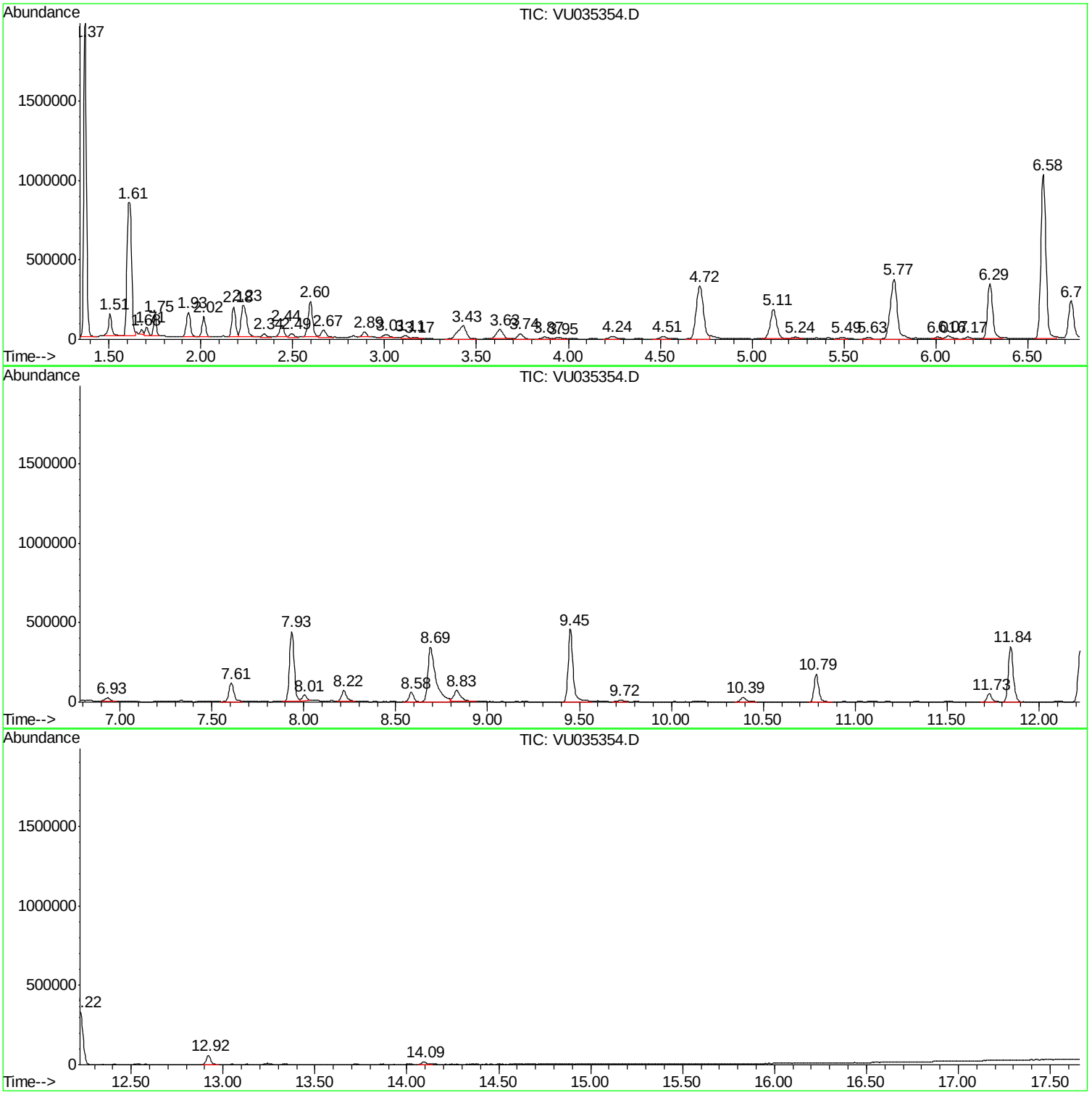
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

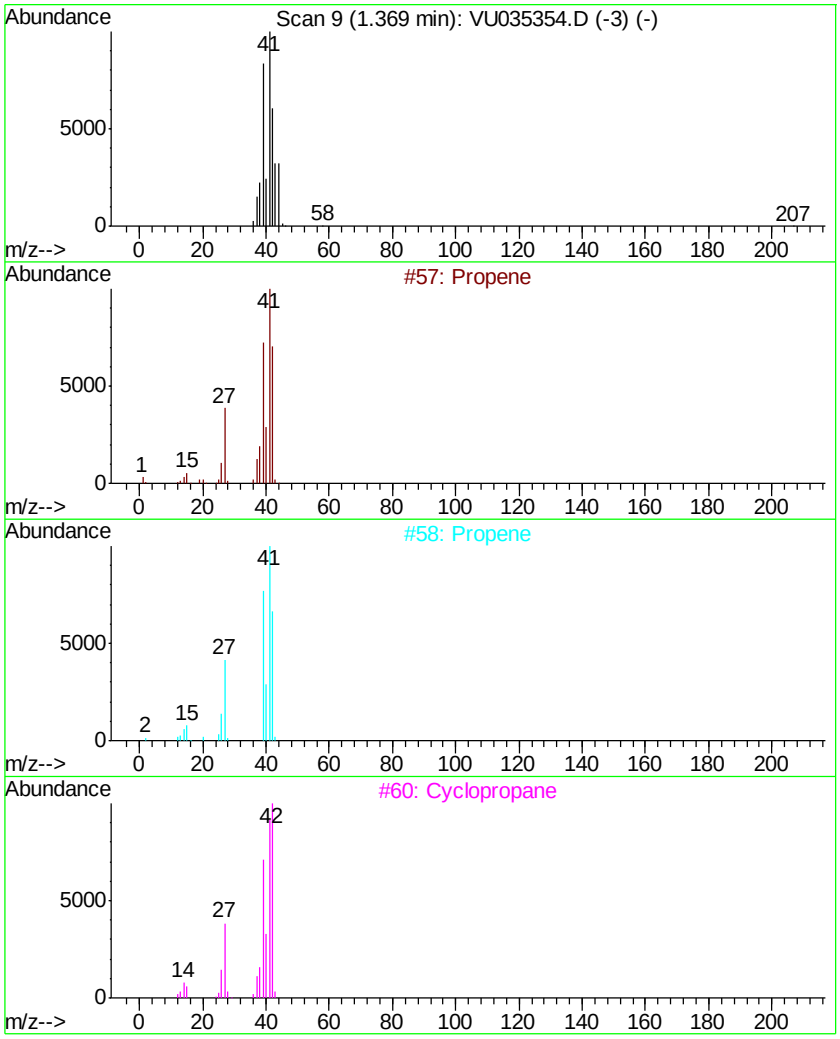
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

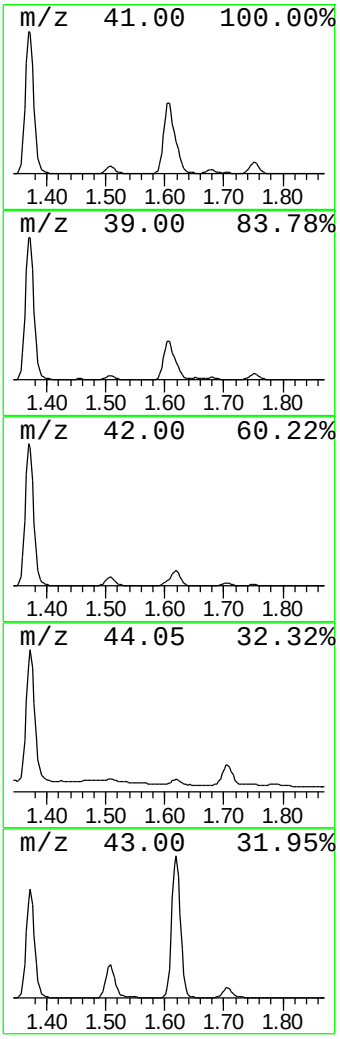
****

Peak Number 1 (DEL) Alkane: Cyclic1.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.37	14.36 ug/L	1985120	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Propene	42	C3H6	000115-07-1	42
3	Cyclopropane	42	C3H6	000075-19-4	7
4	Cyclopropane	42	C3H6	000075-19-4	7
5	Cyclopropene	40	C3H4	002781-85-3	5





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

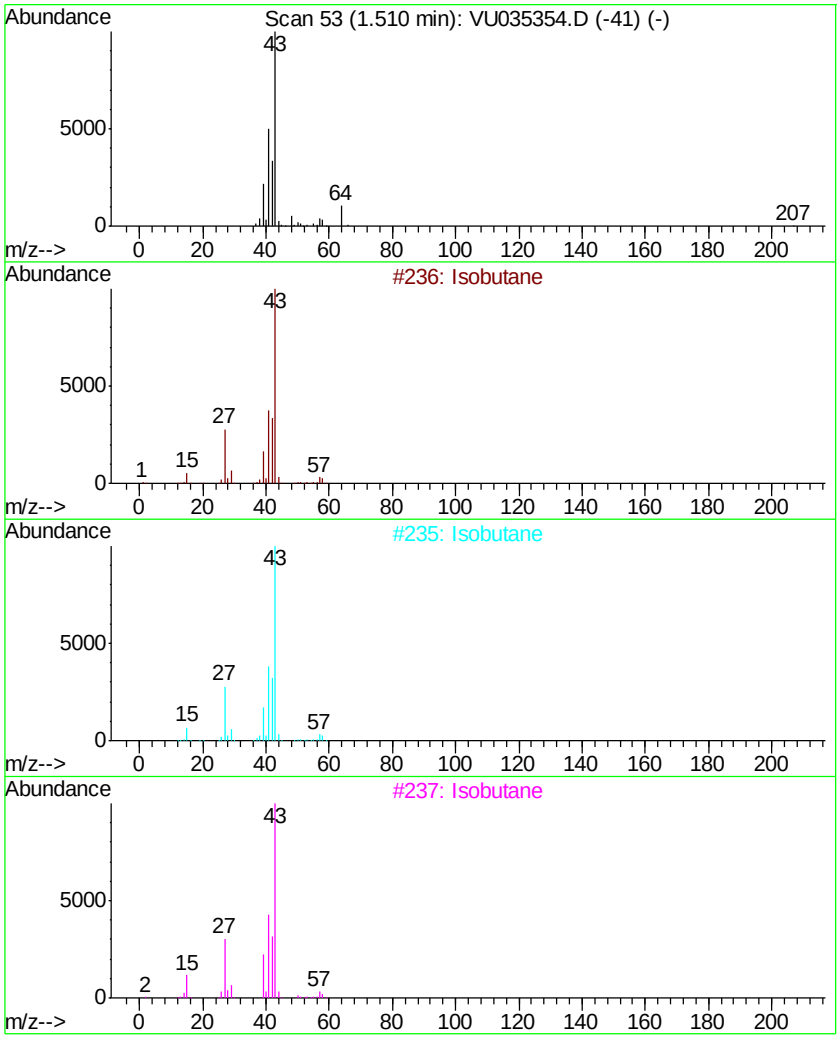
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

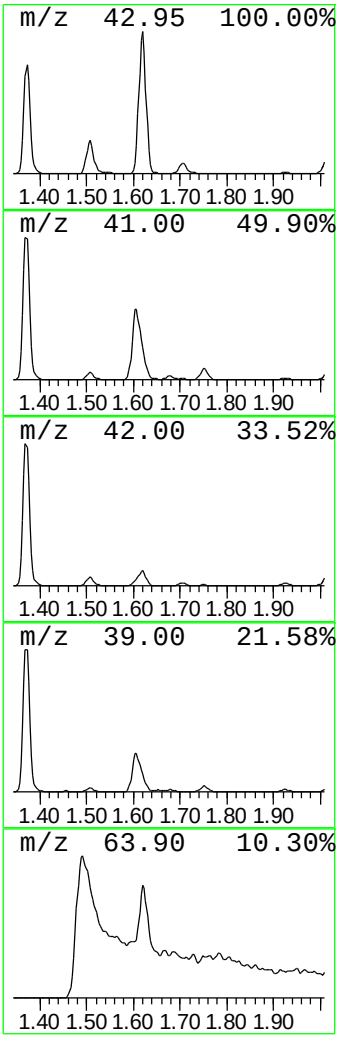
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.51	1.38 ug/L	190070	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isobutane	58	C4H10	000075-28-5	53
2	Isobutane	58	C4H10	000075-28-5	50
3	Isobutane	58	C4H10	000075-28-5	50
4	Isobutane	58	C4H10	000075-28-5	38
5	Propane, 2-nitro-	89	C3H7NO2	000079-46-9	9





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

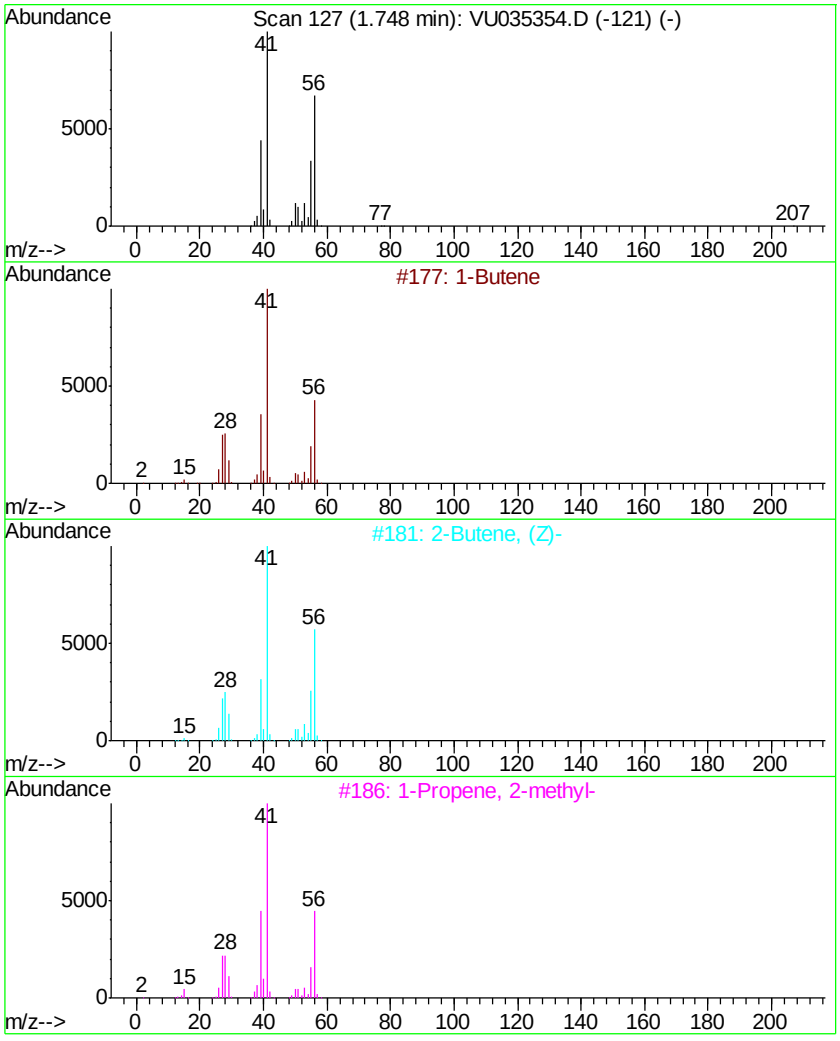
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

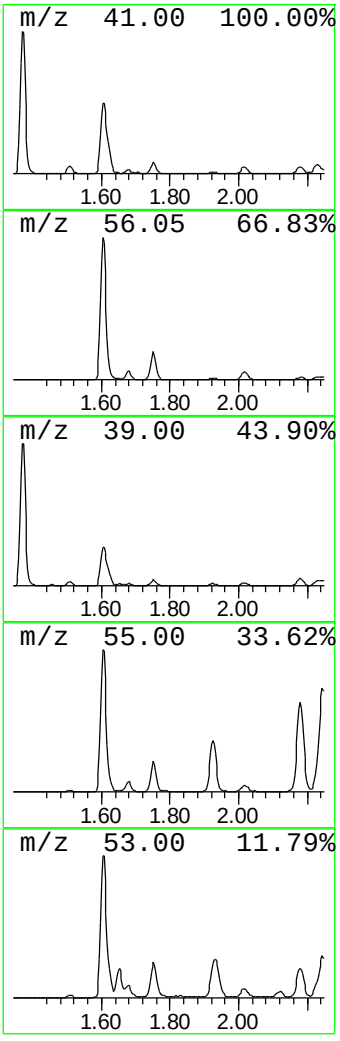
****

Peak Number 3 (DEL) Alkane: Cyclic1.75 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.75	0.95 ug/L	131142	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butene	56	C4H8	000106-98-9	87
2	2-Butene, (Z)-	56	C4H8	000590-18-1	83
3	1-Propene, 2-methyl-	56	C4H8	000115-11-7	83
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	72
5	2-Butene, (Z)-	56	C4H8	000590-18-1	68





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

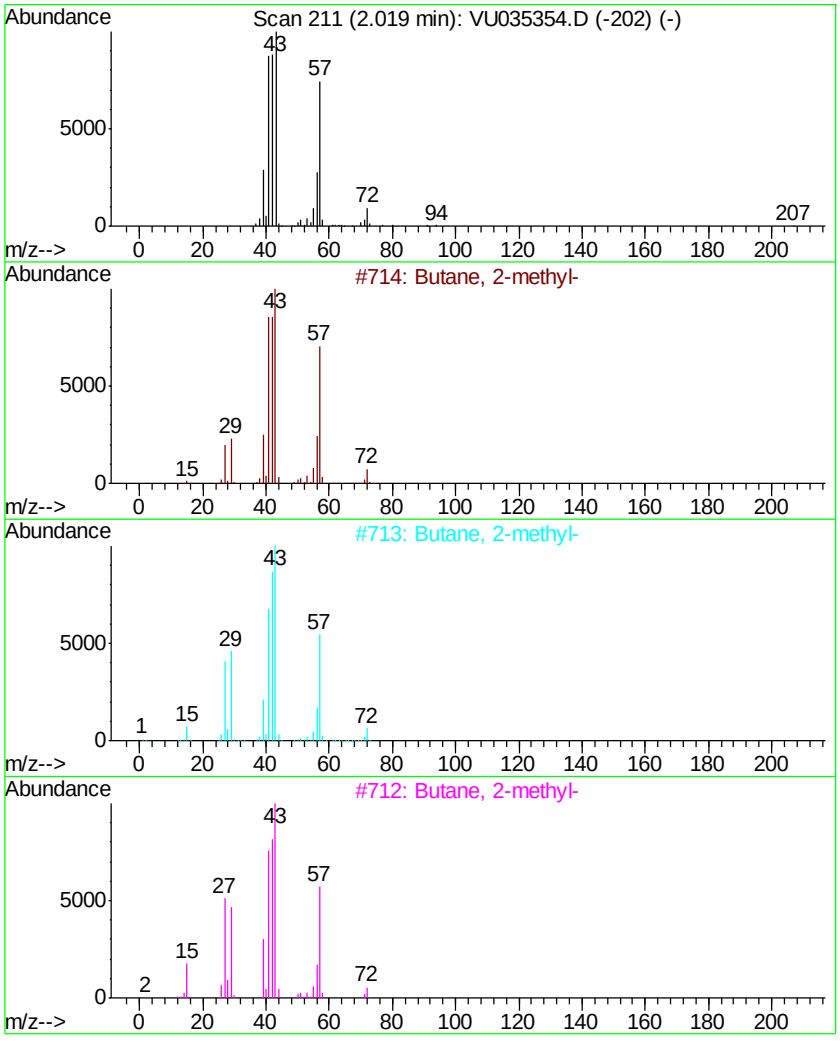
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

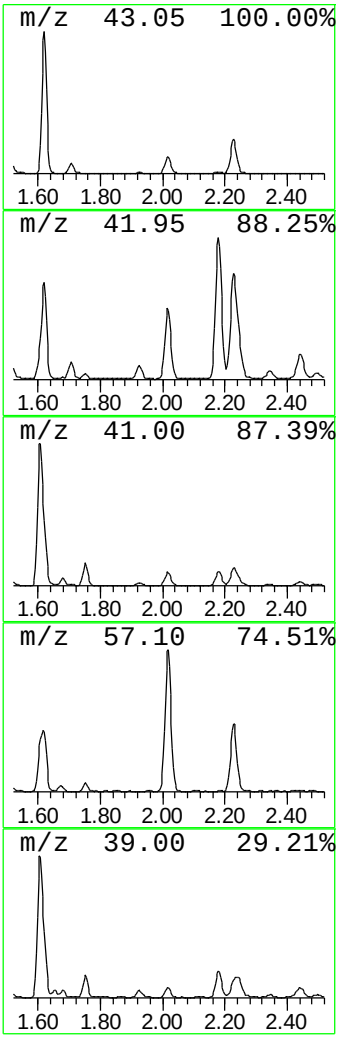
****

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.02	1.23 ug/L	170397	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	91
3	Butane, 2-methyl-	72	C5H12	000078-78-4	80
4	3-Buten-1-ol	72	C4H8O	000627-27-0	47
5	Pentane	72	C5H12	000109-66-0	36





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

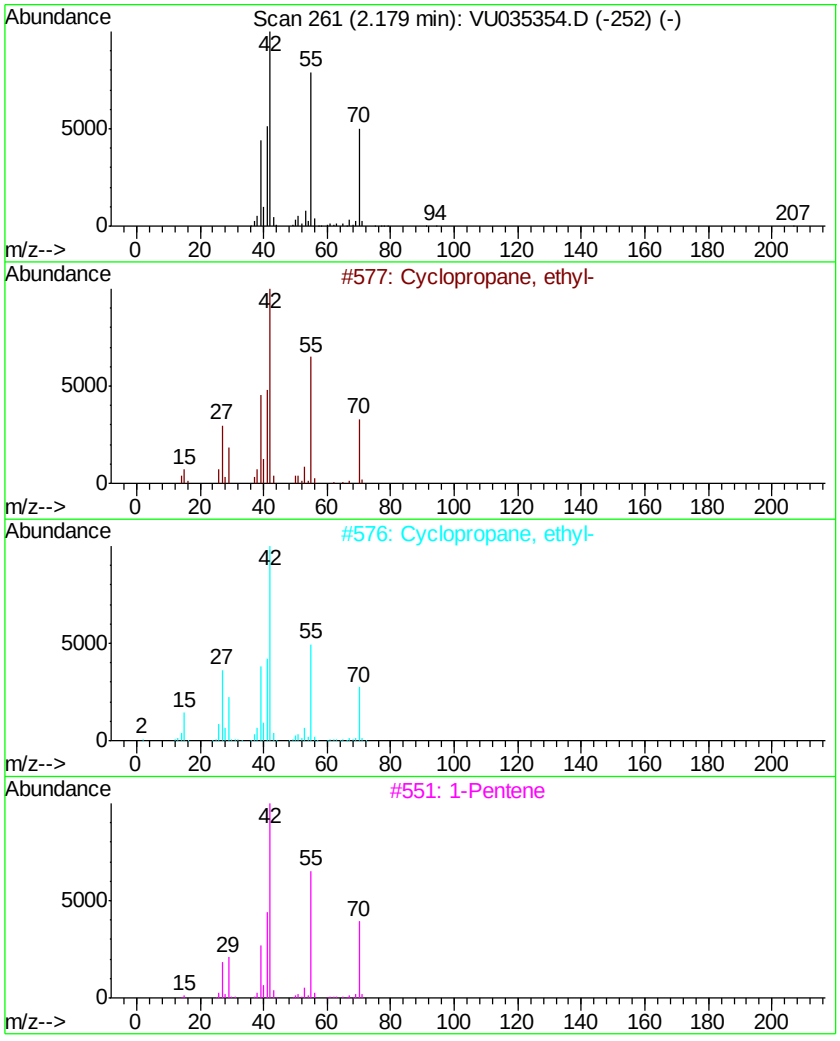
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

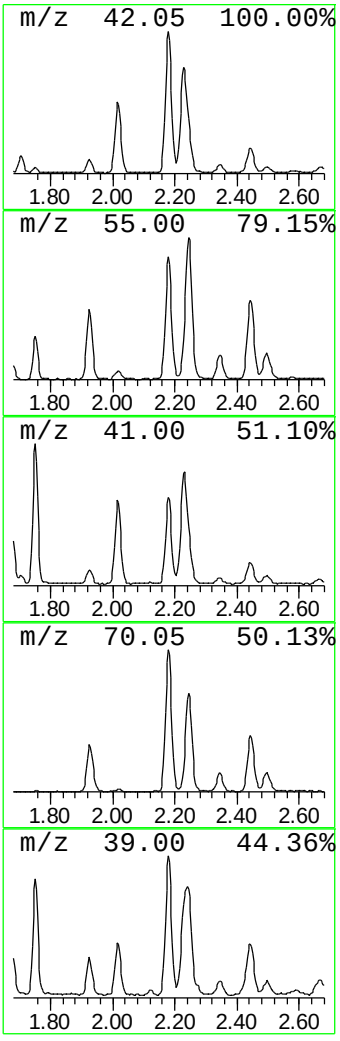
****

Peak Number 5 (DEL) Alkane: Cyclic2.18 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.18	1.89 ug/L	260798	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, ethyl-	70	C5H10	001191-96-4	86
2	Cyclopropane, ethyl-	70	C5H10	001191-96-4	78
3	1-Pentene	70	C5H10	000109-67-1	74
4	1-Pentene	70	C5H10	000109-67-1	68
5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	68





Library Search Compound Report

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Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

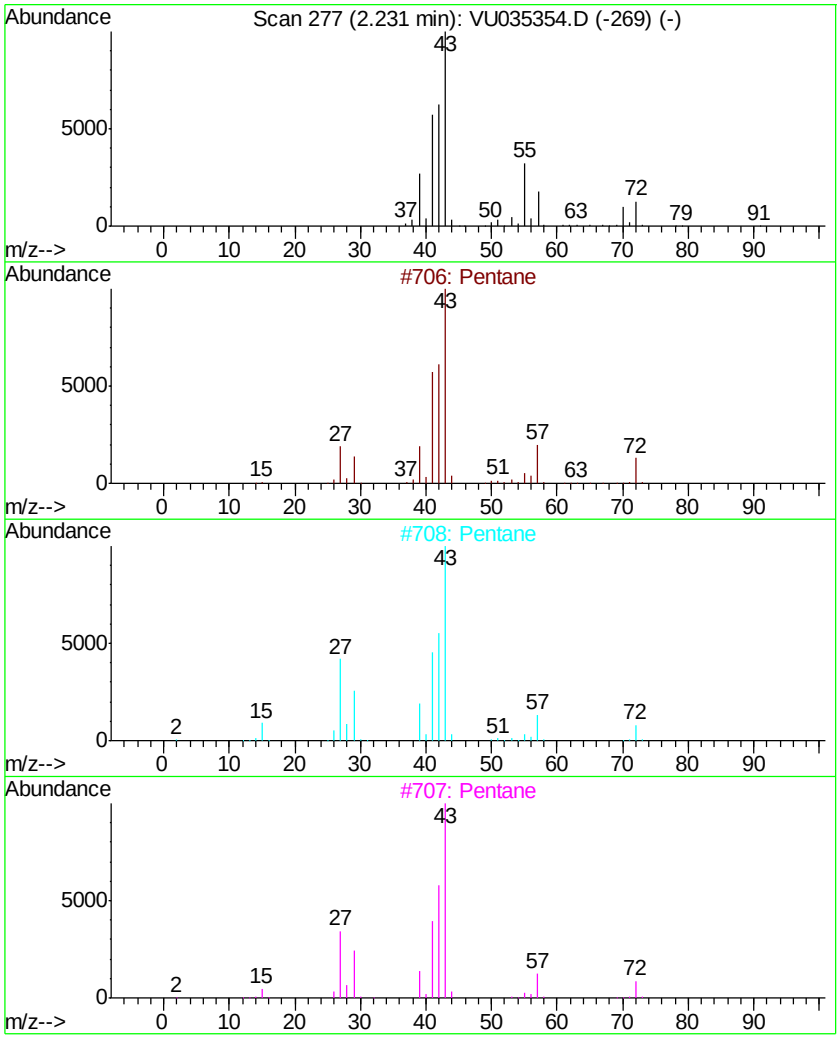
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

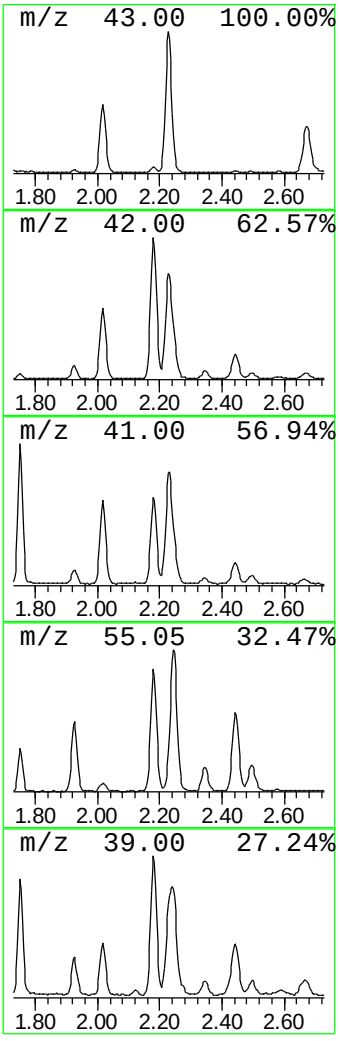
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	3.06 ug/L	422633	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	80
2	Pentane	72	C5H12	000109-66-0	80
3	Pentane	72	C5H12	000109-66-0	72
4	Pentane	72	C5H12	000109-66-0	64
5	Butane, 2-methyl-	72	C5H12	000078-78-4	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

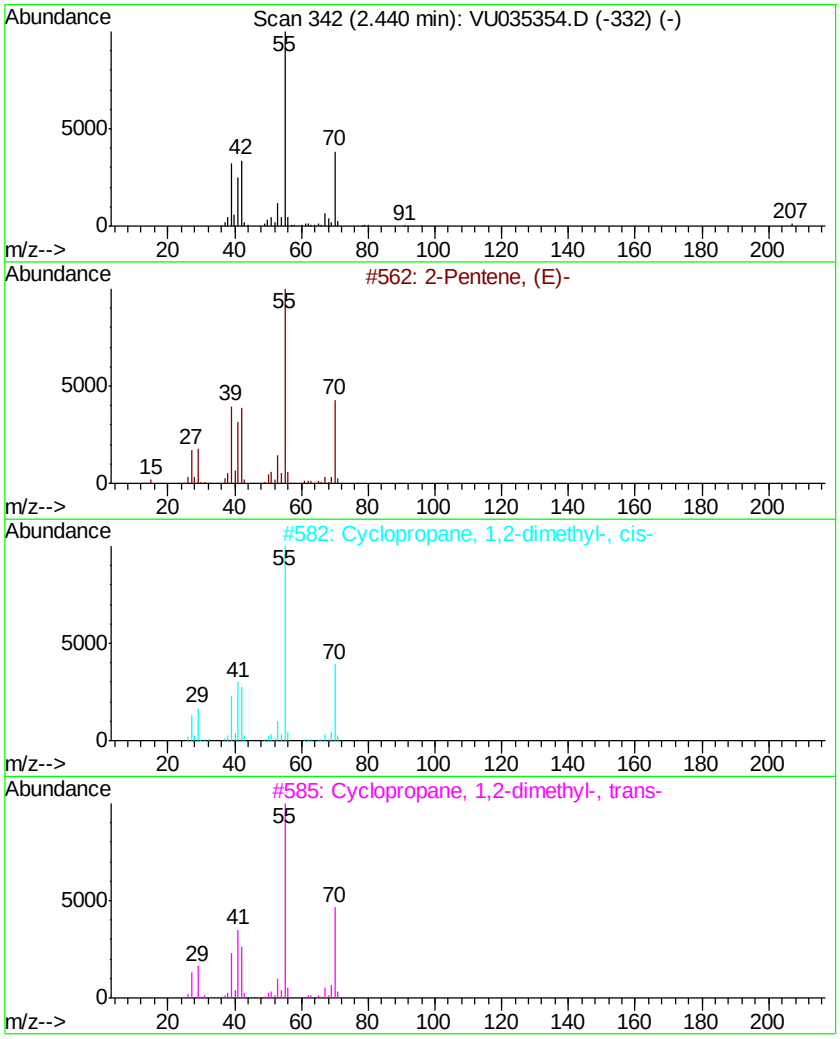
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

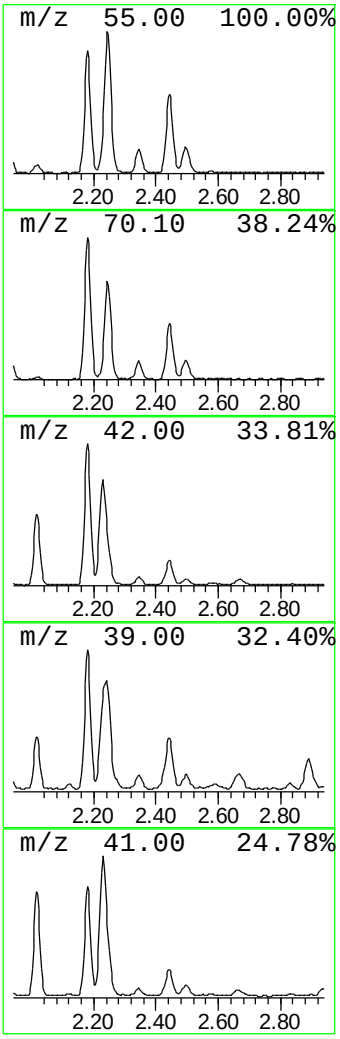
****

Peak Number 7 (DEL) Alkane: Cyclic2.44 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.44	0.89 ug/L	122986	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (E)-	70	C5H10	000646-04-8	93
2	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
3	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	91
4	2-Pentene, (E)-	70	C5H10	000646-04-8	87
5	2-Pentene, (Z)-	70	C5H10	000627-20-3	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

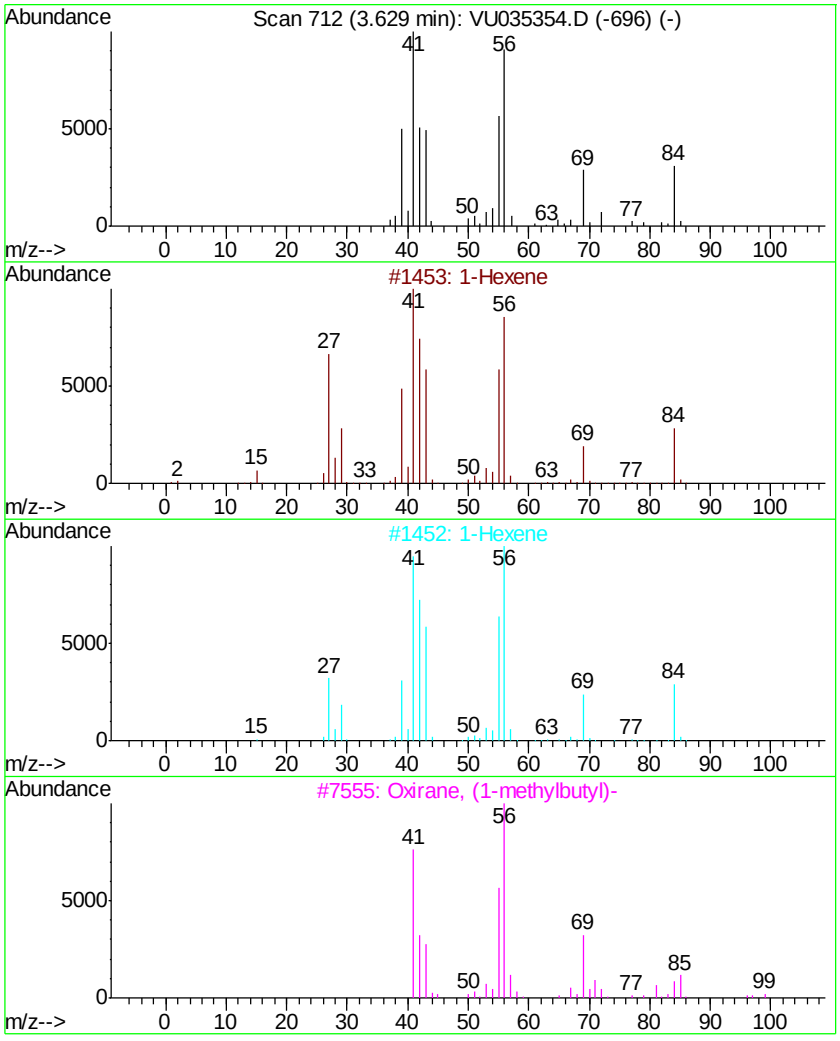
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

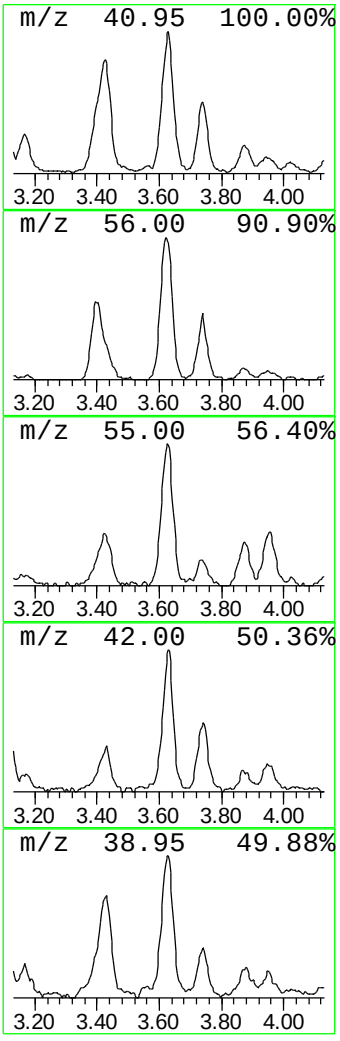
****

Peak Number 8 (DEL) Alkane: Cyclic3.63 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.63	1.10 ug/L	152495	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene	84	C6H12	000592-41-6	81
2	1-Hexene	84	C6H12	000592-41-6	76
3	Oxirane, (1-methylbutyl)-	114	C7H14O	053229-39-3	72
4	2-Hexene, (Z)-	84	C6H12	007688-21-3	52
5	Cyclohexane	84	C6H12	000110-82-7	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

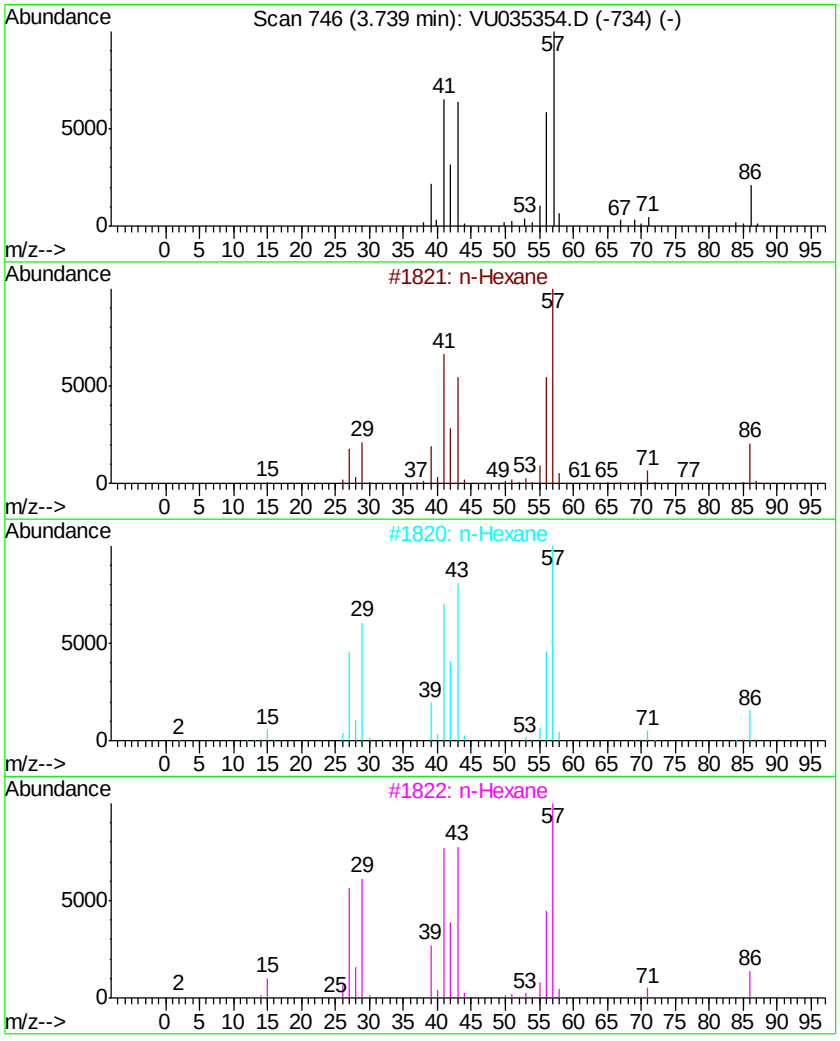
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

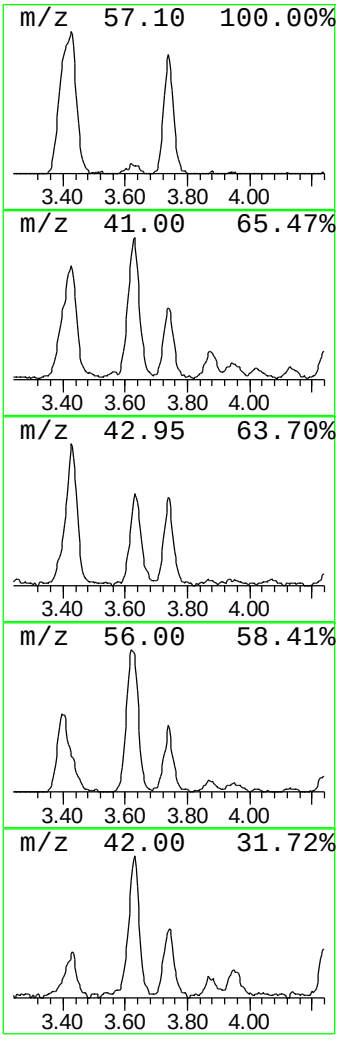
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.74	0.53 ug/L	73257	1,4-Difluorobenzene	6.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	90
2	n-Hexane	86	C6H14	000110-54-3	59
3	n-Hexane	86	C6H14	000110-54-3	53
4	Pentane, 3-methyl-	86	C6H14	000096-14-0	50
5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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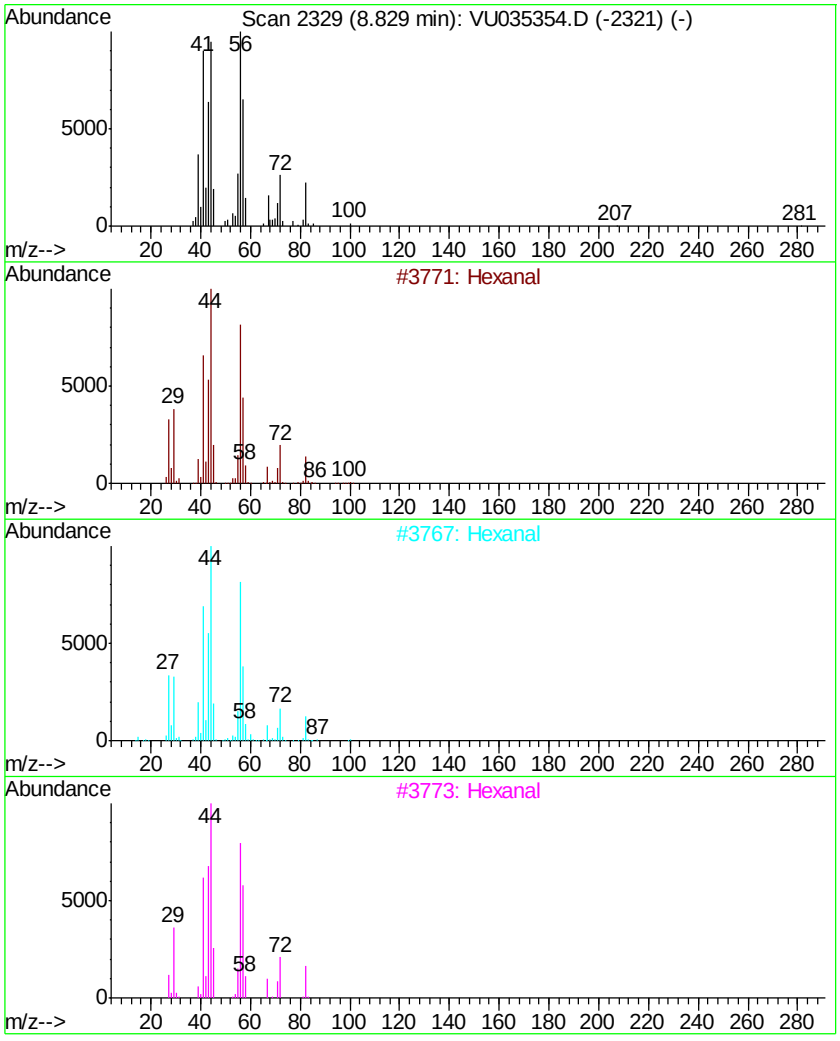
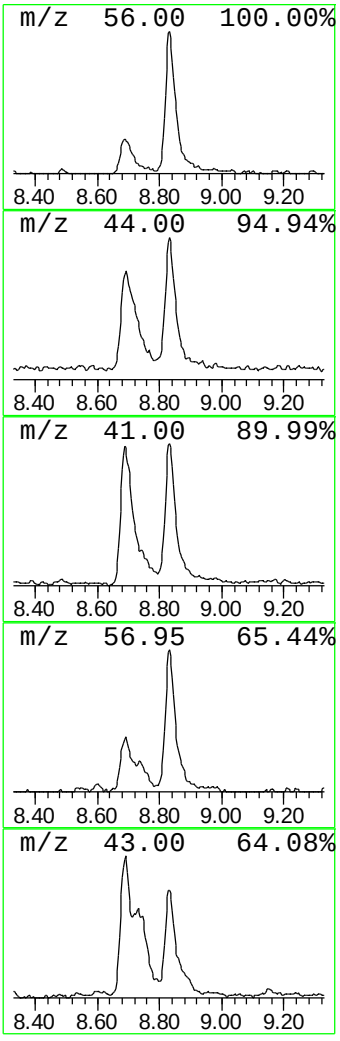
Peak Number 11 Hexanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.83	1.09 ug/L	185188	Chlorobenzene-d5	9.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal	100	C6H12O	000066-25-1	87
2	Hexanal	100	C6H12O	000066-25-1	83
3	Hexanal	100	C6H12O	000066-25-1	83
4	Hexanal	100	C6H12O	000066-25-1	78
5	Pentanal, 3-methyl-	100	C6H12O	015877-57-3	43





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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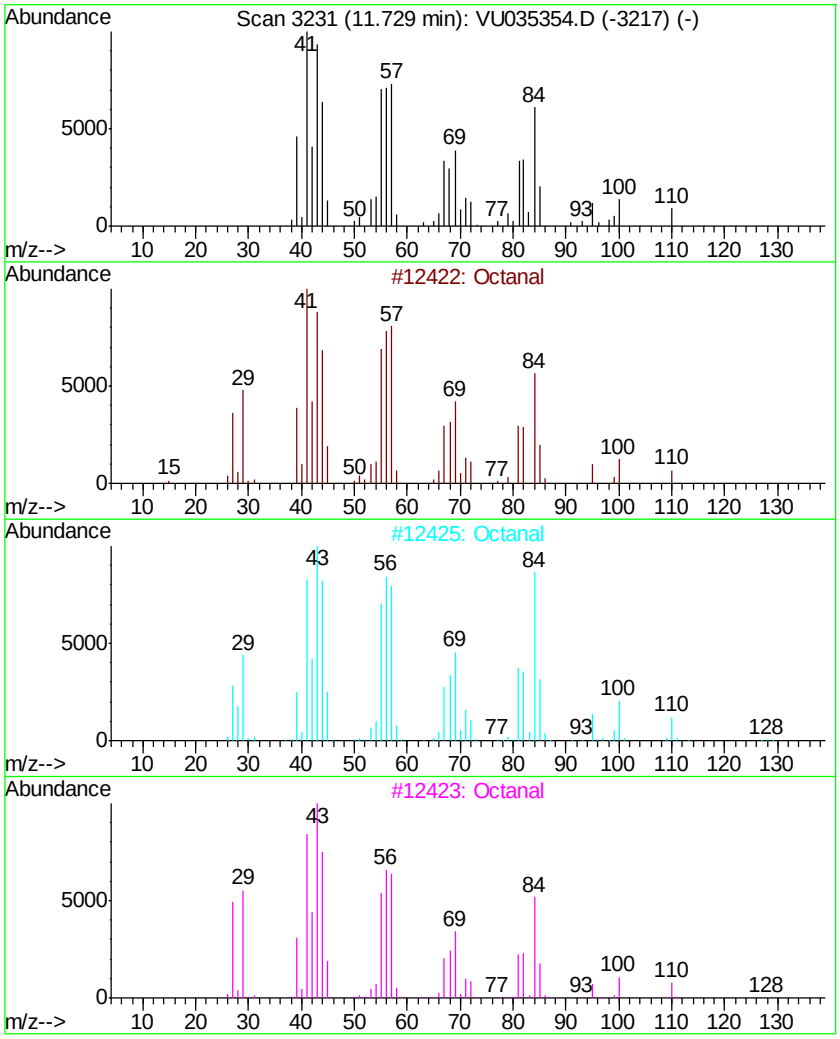
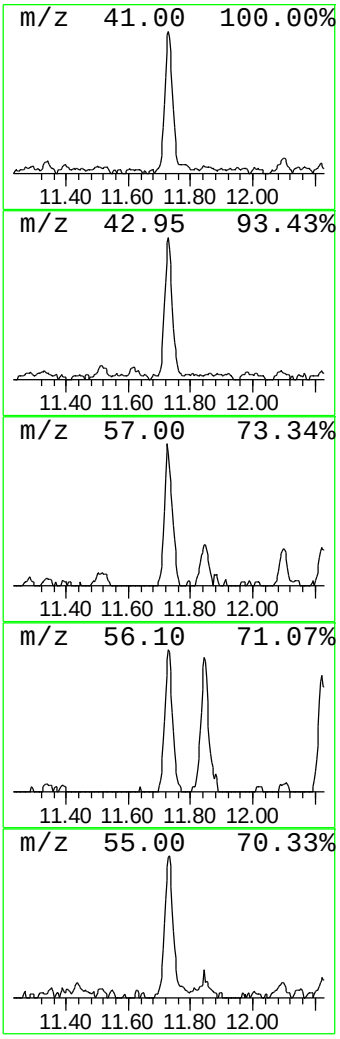
Peak Number 12 Octanal Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	0.69 ug/L	90033	1,4-Dichlorobenzene-d4	11.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanal	128	C8H16O	000124-13-0	91
2	Octanal	128	C8H16O	000124-13-0	87
3	Octanal	128	C8H16O	000124-13-0	72
4	Octanal	128	C8H16O	000124-13-0	72
5	1-Pentene, 3-methyl-	84	C6H12	000760-20-3	43





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

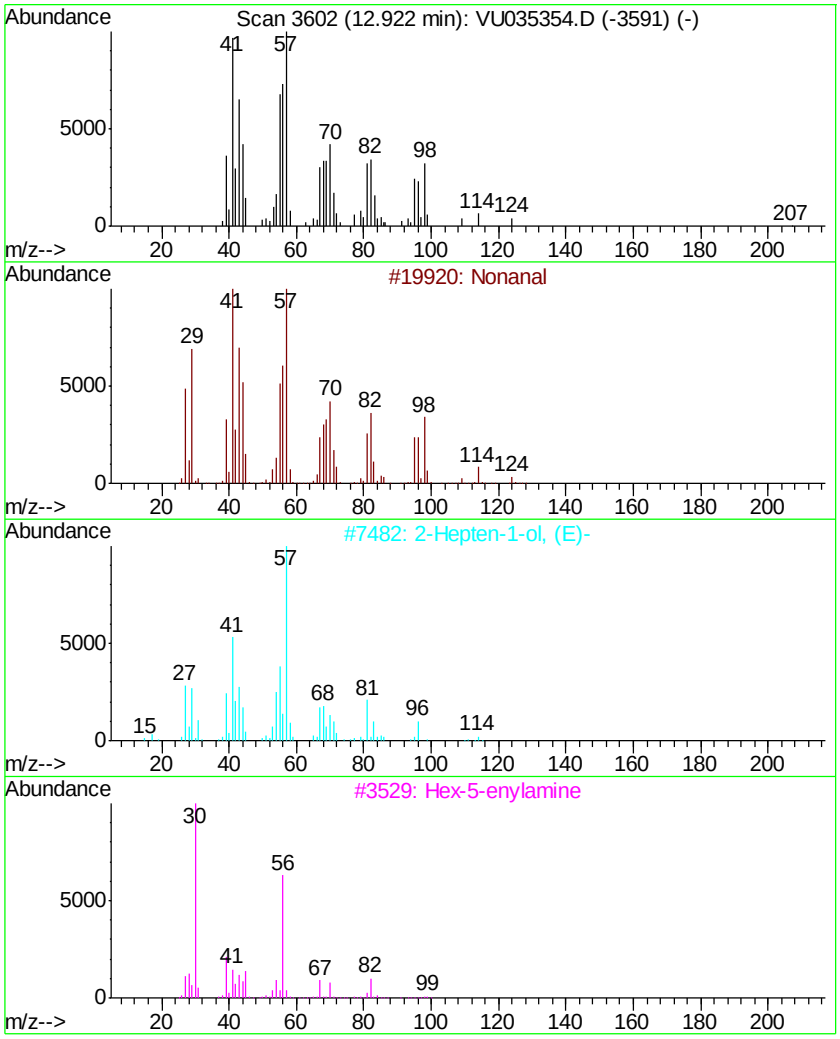
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

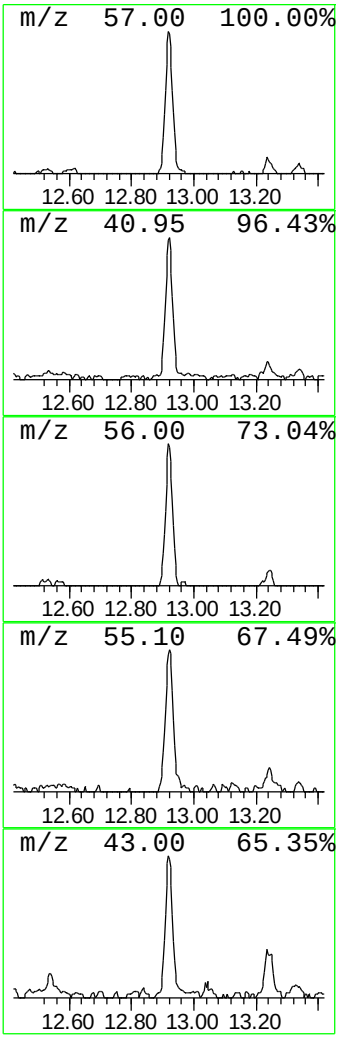
****

Peak Number 13 Nonanal Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	0.70 ug/L	92274	1,4-Dichlorobenzene-d4	11.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanal	142	C9H18O	000124-19-6	91
2	2-Hepten-1-ol, (E)-	114	C7H14O	033467-76-4	38
3	Hex-5-enylamine	99	C6H13N	034825-70-2	25
4	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	22
5	2,3-Dimethyl-aziridine	71	C4H9N	002549-68-0	22





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU102219\
Data File : VU035354.D
Acq On : 22 Oct 2019 10:42
Operator : JC/SP
Sample : K5461-13
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : H0G66

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR101719WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Cyc...	1.37	14.4	ug/L	1985120	1	6.29	691083	5.0
(DEL) Alkane: Str...	1.51	1.4	ug/L	190070	1	6.29	691083	5.0
(DEL) Alkane: Cyc...	1.75	0.9	ug/L	131142	1	6.29	691083	5.0
(DEL) Alkane: Str...	2.02	1.2	ug/L	170397	1	6.29	691083	5.0
(DEL) Alkane: Cyc...	2.18	1.9	ug/L	260798	1	6.29	691083	5.0
(DEL) Alkane: Str...	2.23	3.1	ug/L	422633	1	6.29	691083	5.0
(DEL) Alkane: Cyc...	2.44	0.9	ug/L	122986	1	6.29	691083	5.0
(DEL) Alkane: Cyc...	3.63	1.1	ug/L	152495	1	6.29	691083	5.0
(DEL) Alkane: Str...	3.74	0.5	ug/L	73257	1	6.29	691083	5.0
Hexanal	8.83	1.1	ug/L	185188	2	9.45	852676	5.0
Octanal	11.73	0.7	ug/L	90033	3	11.84	655386	5.0
Nonanal	12.92	0.7	ug/L	92274	3	11.84	655386	5.0