LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

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\SOMUTR083120WMA.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	1.604	72	82	91	rBV2	48981	73451	1.43%	0.259%
2	1.922	175	181	196	rVB2	42696	84016	1.63%	0.297%
3	2.227	266	276	292	rVB4	47327	85280	1.66%	0.301%
4	2.388	315	326	347	rVB	893366	1368863	26.64%	4.831%
5	2.581	373	386	400	rBV	81748	137840	2.68%	0.486%
6	3.099	538	547	559	rVB2	64952	127174	2.47%	0.449%
7	3.378	616	634	654	rVB2	140958	321507	6.26%	1.135%
8	3.719	724	740	761	rVB2	85308	192711	3.75%	0.680%
9	4.012	810	831	854	rBV	1971537	5139163	100.00%	18.137%
10	4.552	979	999	1017	rBV2	415230	1025525	19.96%	3.619%
11	4.668	1018	1035	1075	rVB2	67122	292493	5.69%	1.032%
12	5.086	1147	1165	1221	rBV2	314272	1013475	19.72%	3.577%
13	5.404	1249	1264	1279	rBV2	53455	120918	2.35%	0.427%
14	5.742	1350	1369	1372	rBV3	165639	360922	7.02%	1.274%
15	5.790	1373	1384	1401	rVB	2083833	4194529	81.62%	14.803%
16	6.263	1517	1531	1550	rVB	178511	338985	6.60%	1.196%
17	6.703	1655	1668	1679	rBV	112453	218142	4.24%	0.770%
18	6.774	1680	1690	1702	rVB2	52030	101560	1.98%	0.358%
19	7.574	1926	1939	1953	rVB	58162	100196	1.95%	0.354%
20	7.722	1972	1985	2003	rVB2	70940	135824	2.64%	0.479%
21	7.906	2018	2042	2053	rBV	187377	333490	6.49%	1.177%
22	7.976	2053	2064	2077	rVB	229904	387629	7.54%	1.368%
23	8.185	2117	2129	2141	rBV2	40114	69796	1.36%	0.246%
24	8.639	2257	2270	2284	rBV	354574	653400	12.71%	2.306%
25	8.800	2308	2320	2338	rBV6	25518	63305	1.23%	0.223%
26	9.449	2499	2522	2551	rVV	2685882	4907079	95.48%	17.318%
27	9.571	2551	2560	2583	rVB4	34489	74677	1.45%	0.264%
28	9.693	2583	2598	2607	rBV	175959	302603	5.89%	1.068%
29	9.761	2607	2619	2633	rVB2	81453	161700	3.15%	0.571%
30	10.102	2714	2725	2739	rVB3	103343	192220	3.74%	0.678%
31	10.256	2757	2773	2793	rBV3	31744	79567	1.55%	0.281%
32	10.481	2831	2843	2854	rBV	278177	438472	8.53%	1.547%
33	10.639	2876	2892	2907	rBV5	36042	108132	2.10%	0.382%
34	10.754	2919	2928	2938	rVB	118920	197770	3.85%	0.698%

LSC Area Percent Report

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Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

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\SOMUTR083120WMA.M
Title : TRACE VOA SOM01.0

35	10.899	2952	2973	2993	rVB2	213483	467390	9.09%	1.650%
36	11.028	3005	3013	3024	rVB3	55753	94486	1.84%	0.333%
37	11.288	3081	3094	3102	rBV	130608	215363	4.19%	0.760%
38	11.462	3139	3148	3166	rVB	74212	124613	2.42%	0.440%
39	11.632	3181	3201	3220	rVB	99940	262907	5.12%	0.928%
40	11.809	3243	3256	3258	rBV	310087	441885	8.60%	1.560%
41	11.886	3272	3280	3299	rVB	164454	270564	5.26%	0.955%
42	12.008	3310	3318	3328	rVB2	48031	74112	1.44%	0.262%
43	12.089	3328	3343	3358	rVV2	465448	758289	14.76%	2.676%
44	12.185	3361	3373	3393	rVB2	269989	547941	10.66%	1.934%
45	12.465	3445	3460	3476	rBV7	37766	94000	1.83%	0.332%
46	12.561	3480	3490	3498	rBV	111685	168613	3.28%	0.595%
47	12.674	3518	3525	3541	rVB3	36332	75207	1.46%	0.265%
48	12.851	3561	3580	3602	rVV	276049	488539	9.51%	1.724%
49	12.963	3605	3615	3622	rVV	61840	104518	2.03%	0.369%
50	13.439	3754	3763	3777	rBV3	38886	72760	1.42%	0.257%
51	13.934	3908	3917	3924	rBV3	50318	83483	1.62%	0.295%
52	13.983	3924	3932	3944	rVB2	104214	171206	3.33%	0.604%
53	14.073	3951	3960	3972	rBV	63453	95872	1.87%	0.338%
54	14.272	4011	4022	4035	rBV9	30301	63604	1.24%	0.224%
55	15.063	4252	4268	4279	rBV2	149892	257113	5.00%	0.907%

Sum of corrected areas: 28334879

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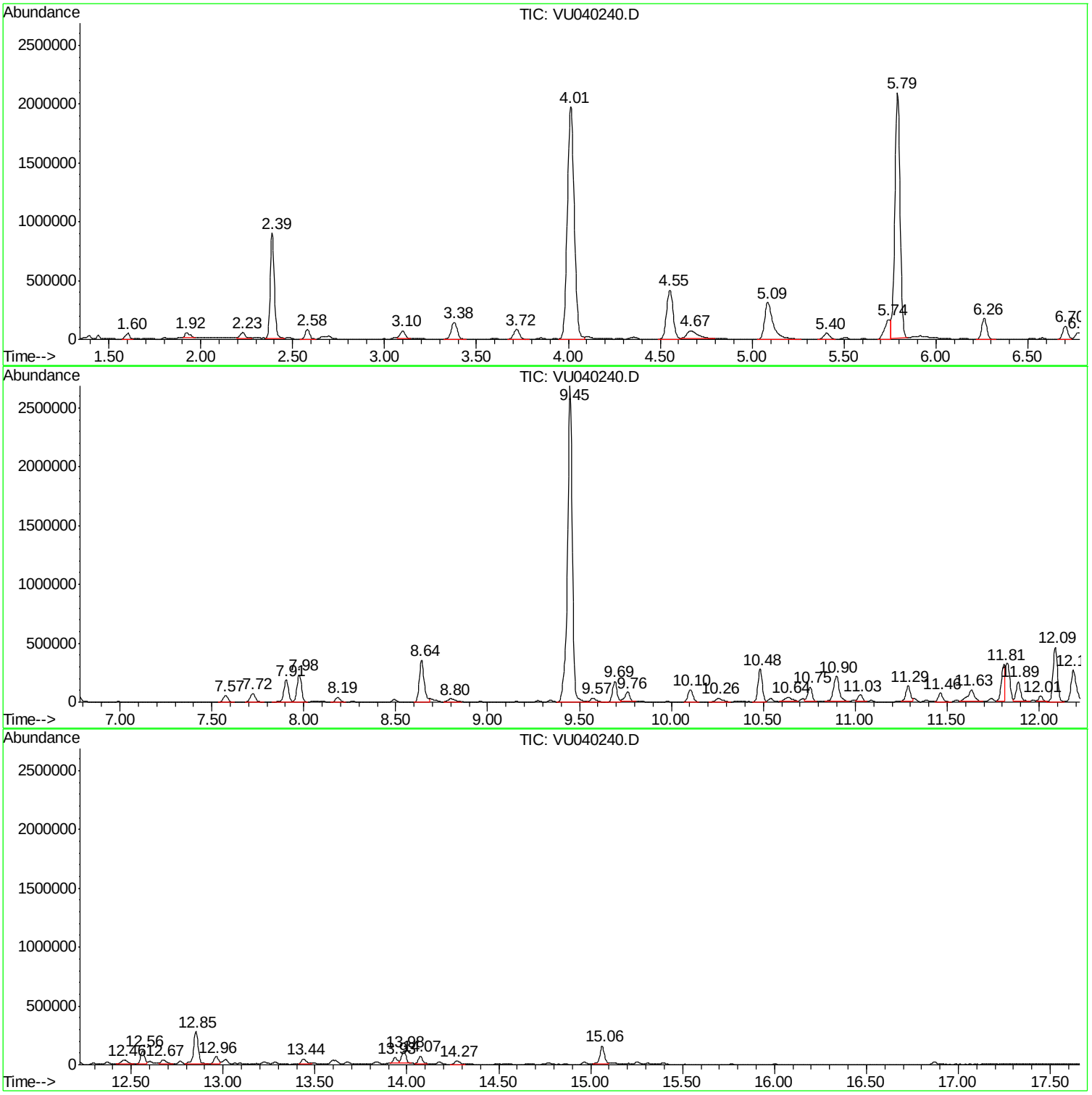
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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



Library Search Compound Report

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Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
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Quant Title : TRACE VOA SOM01.0

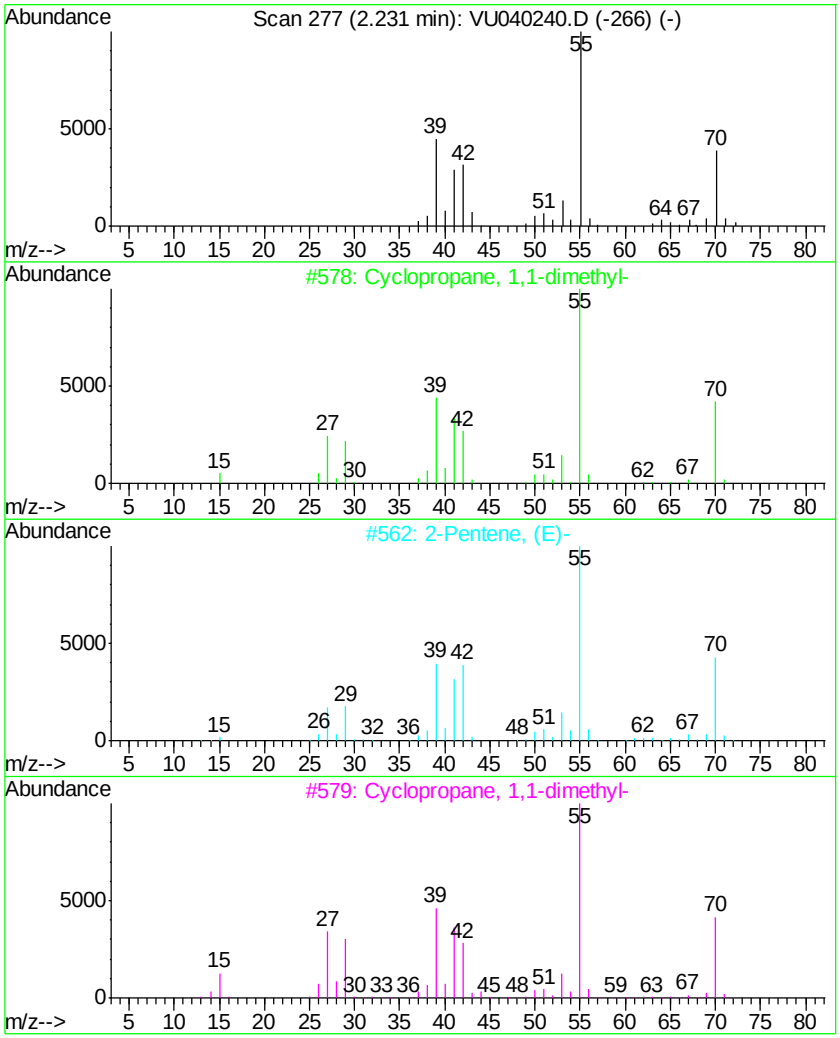
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

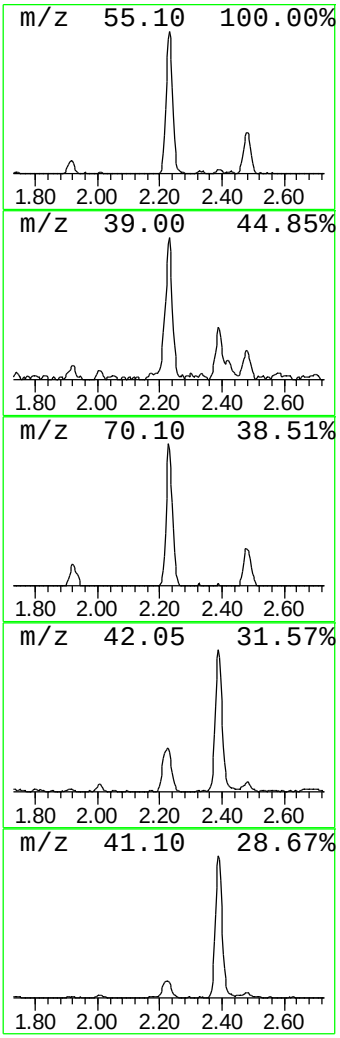
****

Peak Number 1 (DEL) Alkane: Cyclic2.23 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	1.26 ug/L	85280	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	90
2	2-Pentene, (E)-	70	C5H10	000646-04-8	90
3	Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	90
4	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	87
5	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

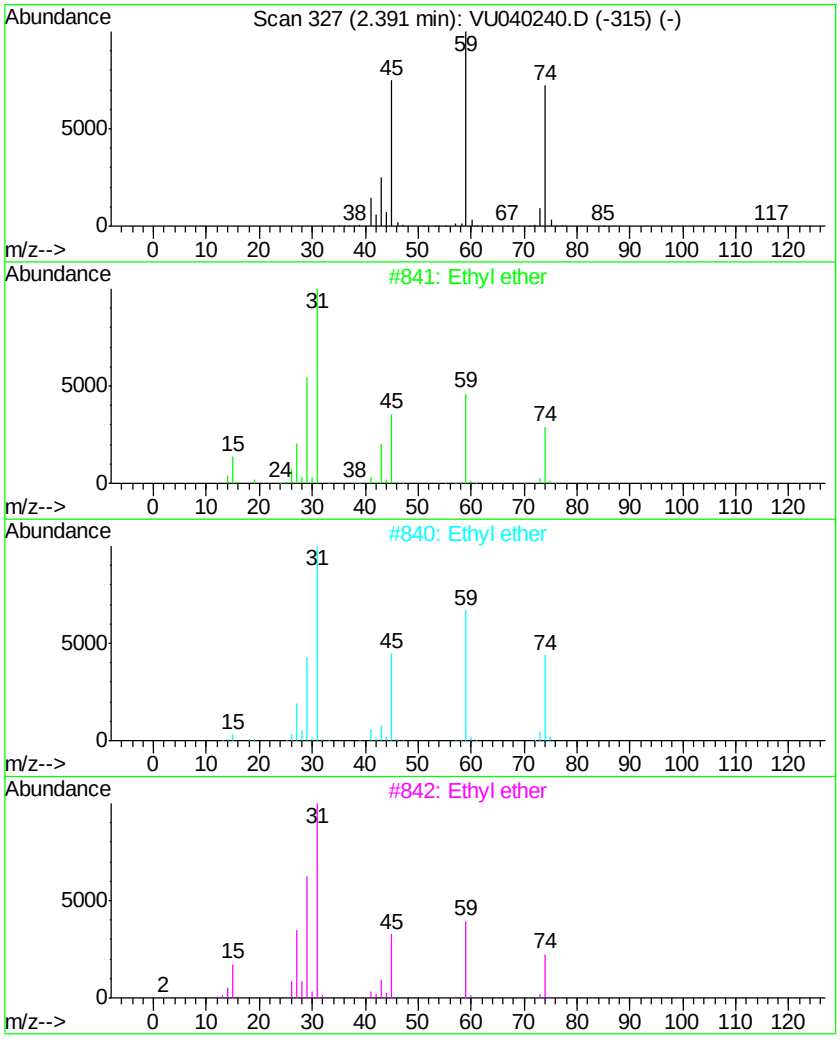
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

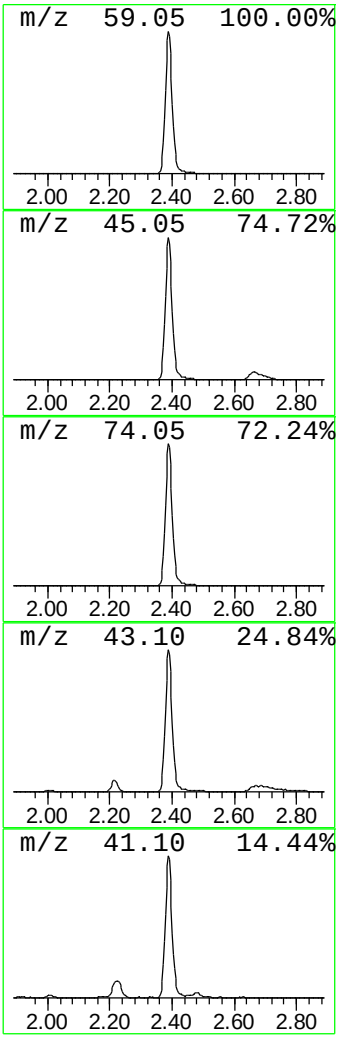
****

Peak Number 2 Ethyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.39	20.19 ug/L	1368860	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl ether	74	C4H10O	000060-29-7	90
2	Ethyl ether	74	C4H10O	000060-29-7	90
3	Ethyl ether	74	C4H10O	000060-29-7	90
4	Ethyl ether	74	C4H10O	000060-29-7	78
5	Ethyl ether	74	C4H10O	000060-29-7	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

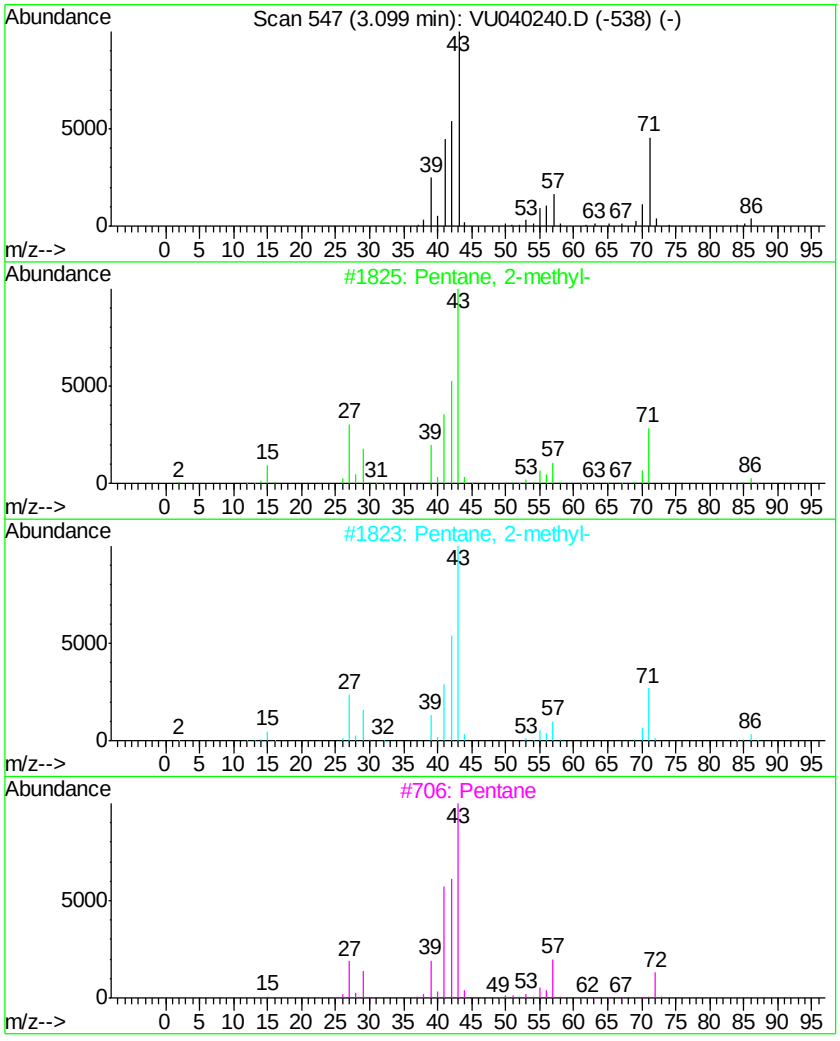
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

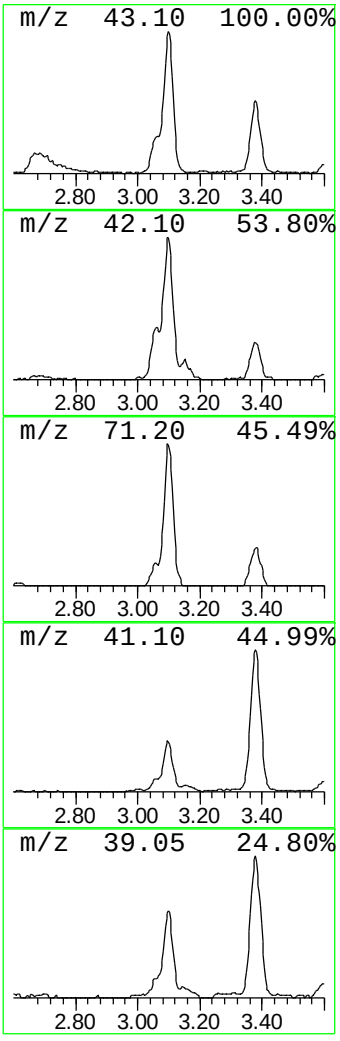
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	1.88 ug/L	127174	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	83
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	78
3	Pentane	72	C5H12	000109-66-0	47
4	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	38
5	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	36





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

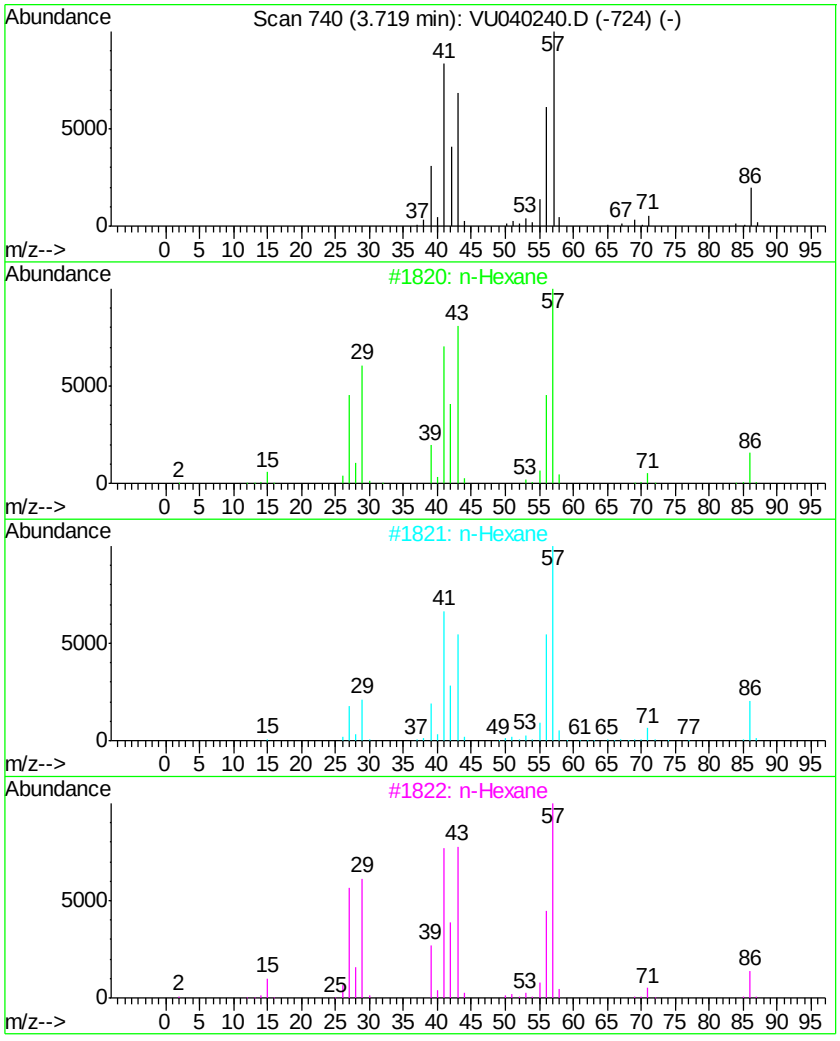
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

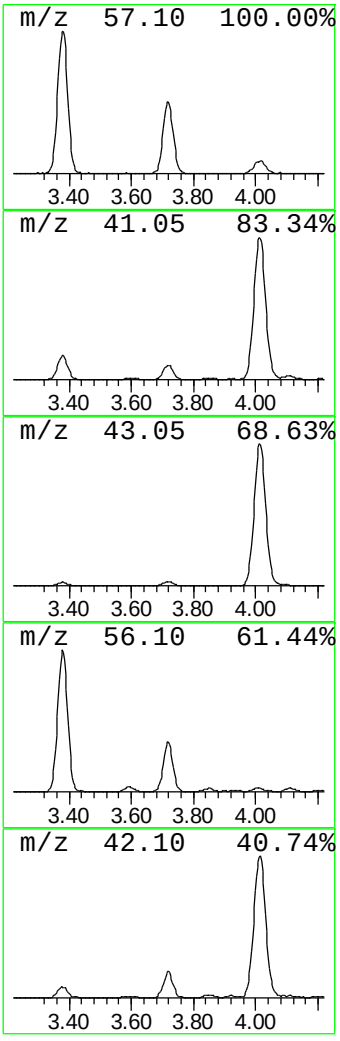
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.72	2.84 ug/L	192711	1,4-Difluorobenzene	6.26

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	87
3	n-Hexane	86	C6H14	000110-54-3	64
4	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	40
5	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	40





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Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

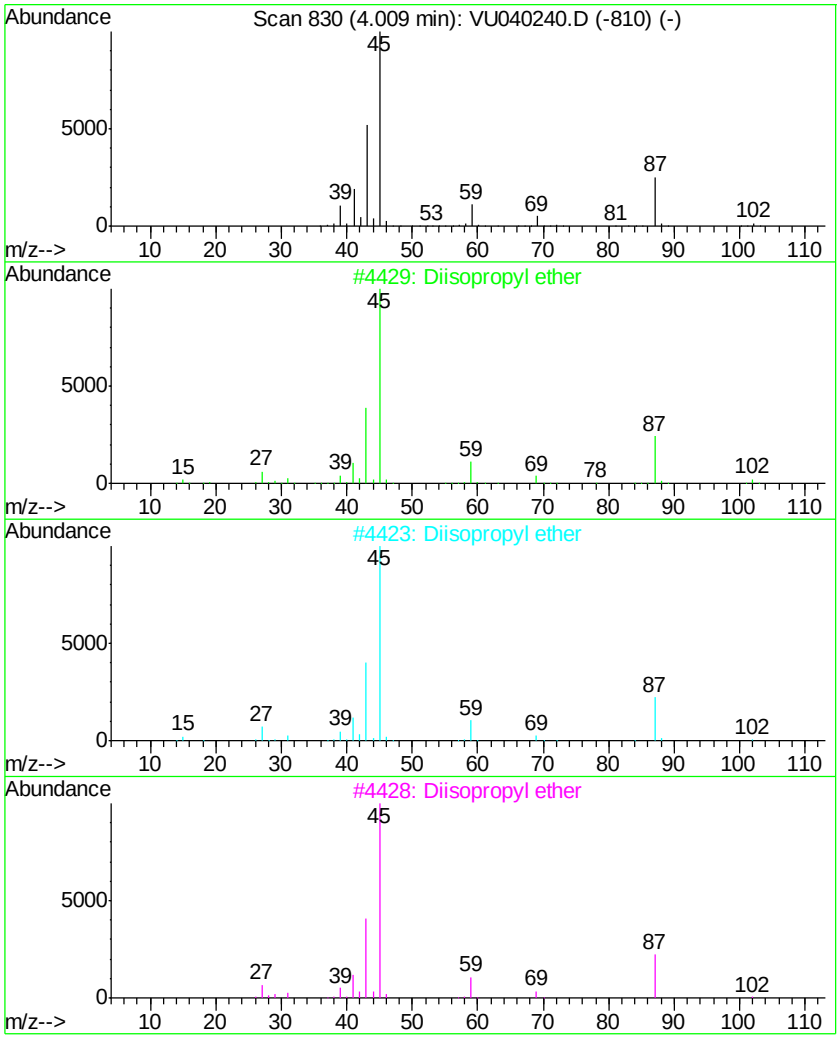
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

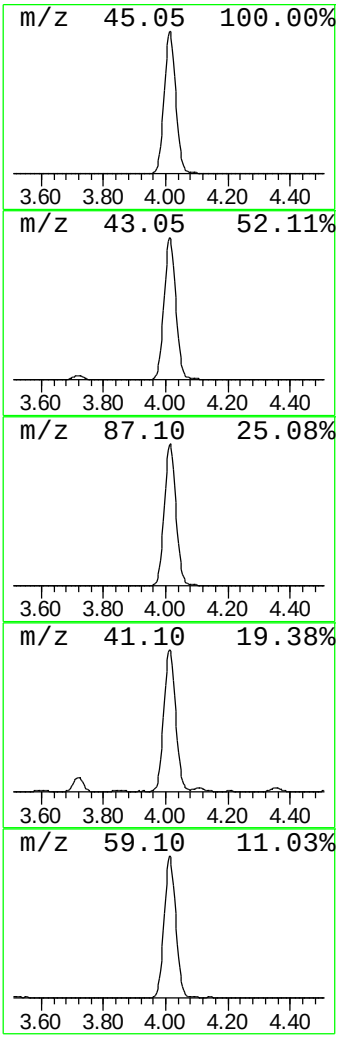
****

Peak Number 5 Diisopropyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.01	75.80 ug/L	5139160	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisopropyl ether	102	C6H14O	000108-20-3	86
2	Diisopropyl ether	102	C6H14O	000108-20-3	83
3	Diisopropyl ether	102	C6H14O	000108-20-3	83
4	1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	72
5	Ether, sec-butyl isopropyl	116	C7H16O	018641-81-1	64





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ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

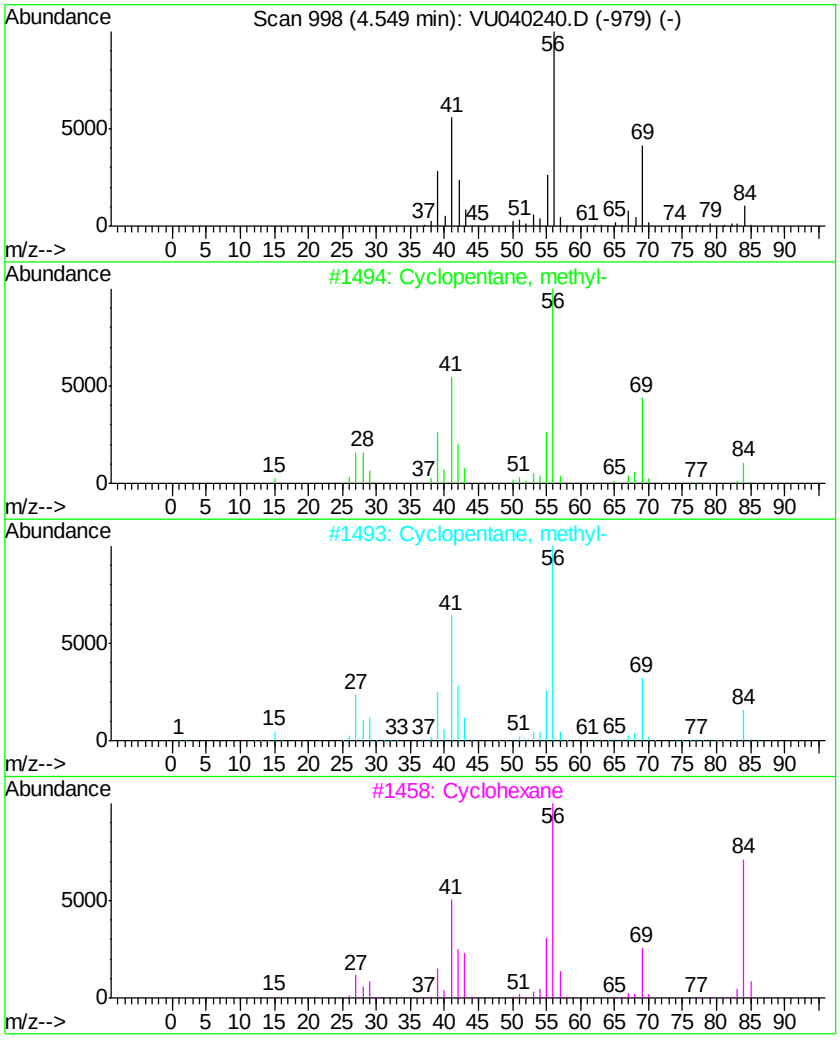
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

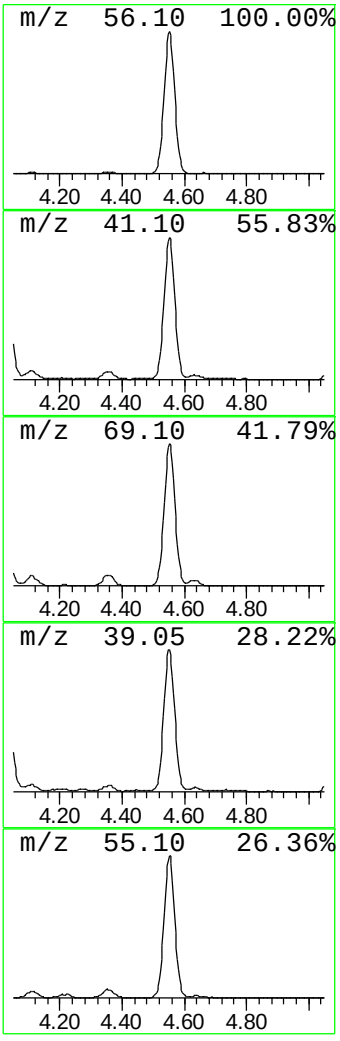
****

Peak Number 6 (DEL) Alkane: Cyclic4.55 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.55	15.13 ug/L	1025530	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
3	Cyclohexane	84	C6H12	000110-82-7	78
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	78
5	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72





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ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

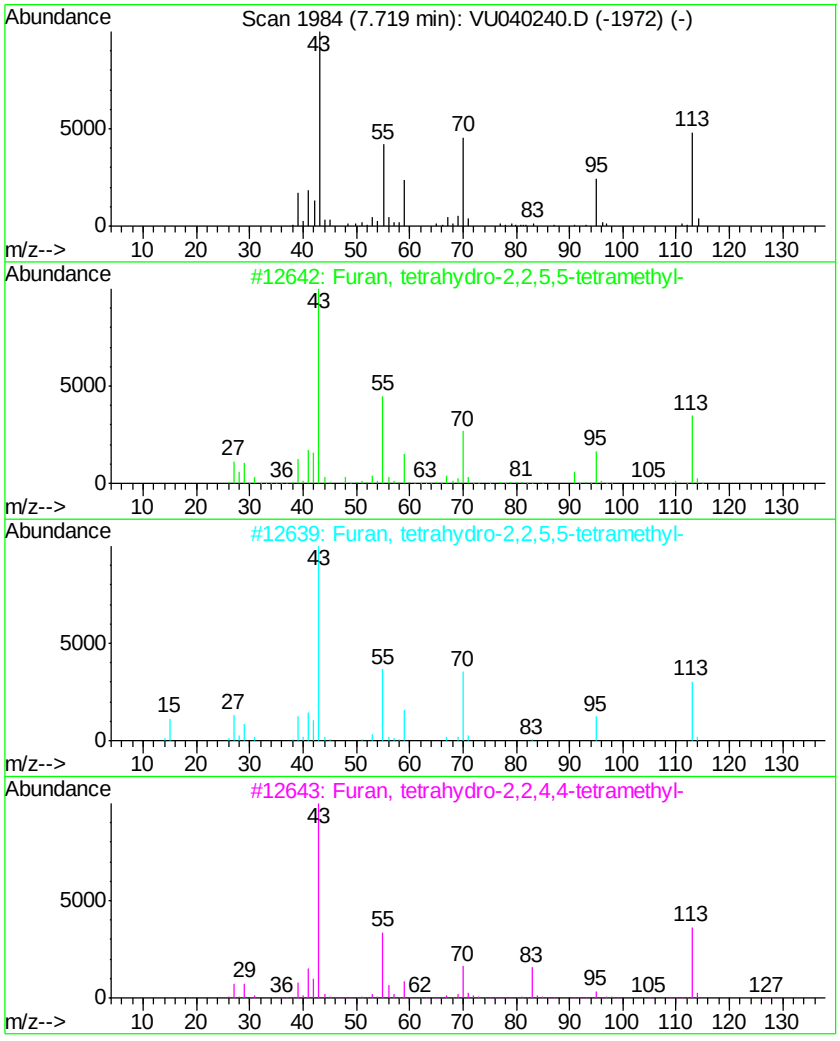
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

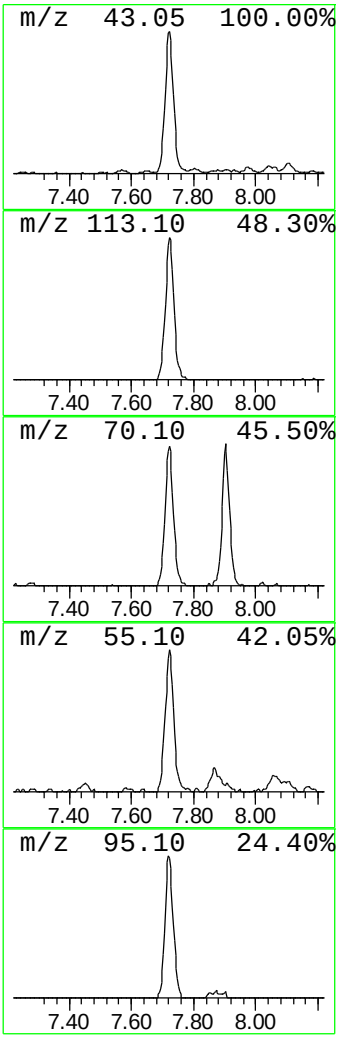
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Furan, tetrahydro-2,2,5,5-t... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	2.00 ug/L	135824	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, tetrahydro-2,2,5,5-tetram...	128	C8H16O	015045-43-9	64
2	Furan, tetrahydro-2,2,5,5-tetram...	128	C8H16O	015045-43-9	64
3	Furan, tetrahydro-2,2,4,4-tetram...	128	C8H16O	003358-28-9	53
4	4-Piperidinone, 1-methyl-	113	C6H11NO	001445-73-4	50
5	Pyrrolidine, 1-acetyl-	113	C6H11NO	004030-18-6	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

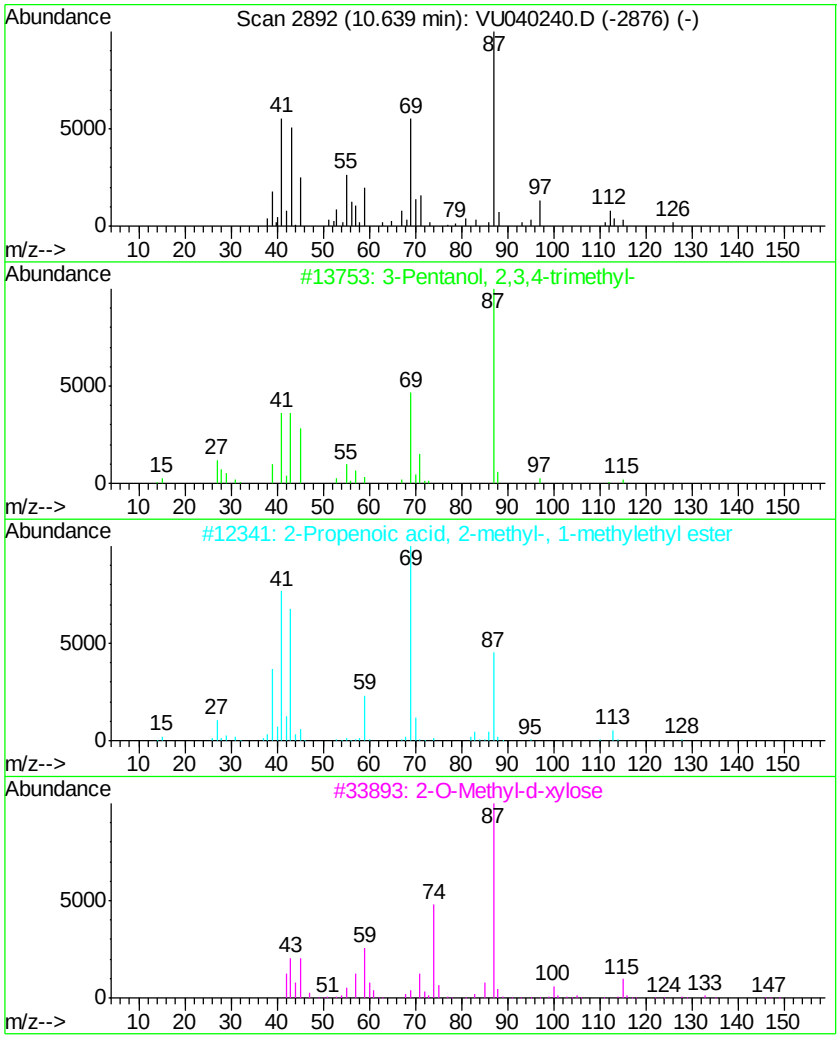
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

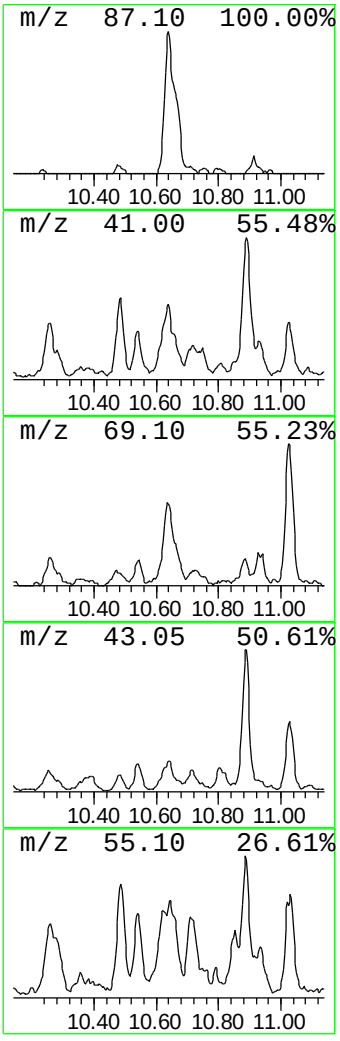
****

Peak Number 9 3-Pentanol, 2,3,4-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.64	1.22 ug/L	108132	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 2,3,4-trimethyl-	130	C8H18O	003054-92-0	59
2	2-Propenoic acid, 2-methyl-, 1-m...	128	C7H12O2	004655-34-9	47
3	2-O-Methyl-d-xylose	164	C6H12O5	1000130-01-3	45
4	3-Pentanol, 2,3,4-trimethyl-	130	C8H18O	003054-92-0	45
5	2,3,4-Trimethyl-5-hexen-3-ol	142	C9H18O	028638-29-1	40





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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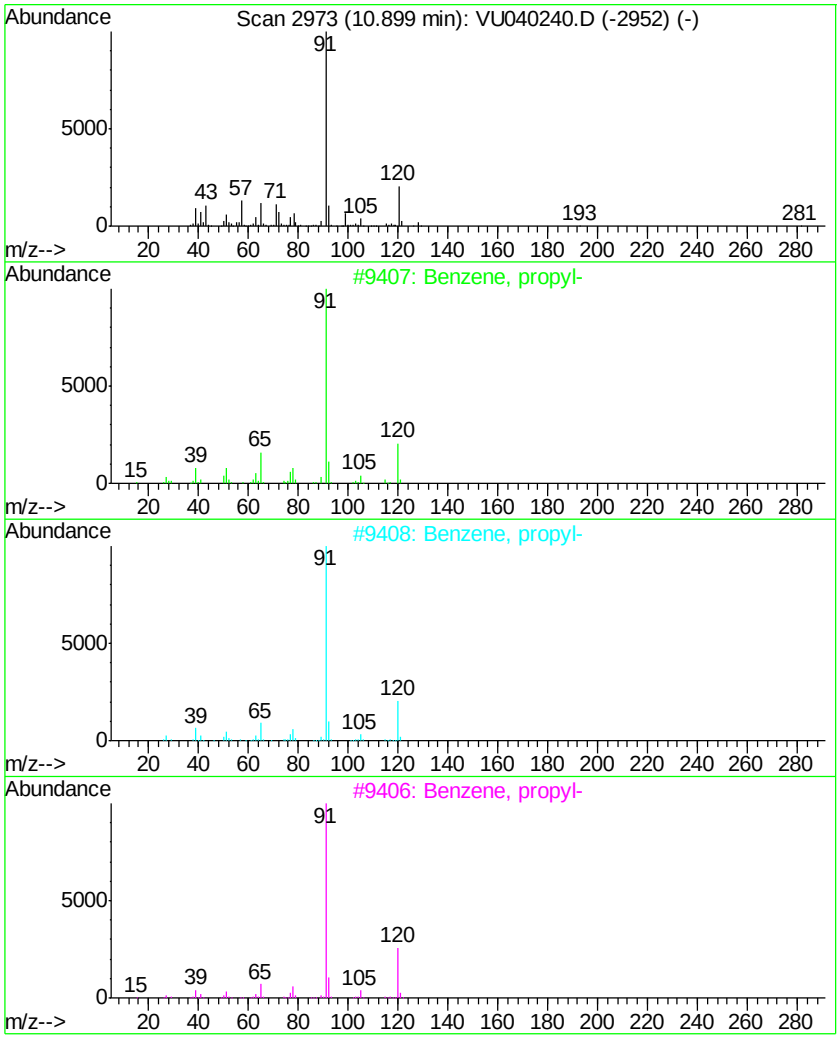
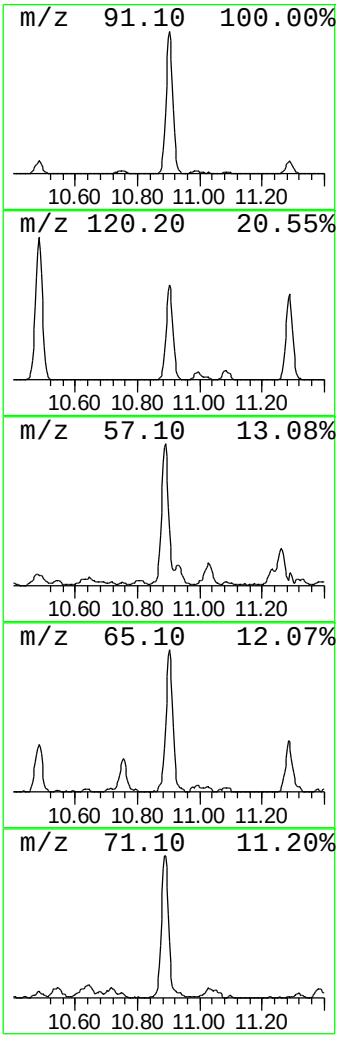
Peak Number 10 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	5.29 ug/L	467390	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	60
2	Benzene, propyl-	120	C9H12	000103-65-1	60
3	Benzene, propyl-	120	C9H12	000103-65-1	53
4	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	47
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	47





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

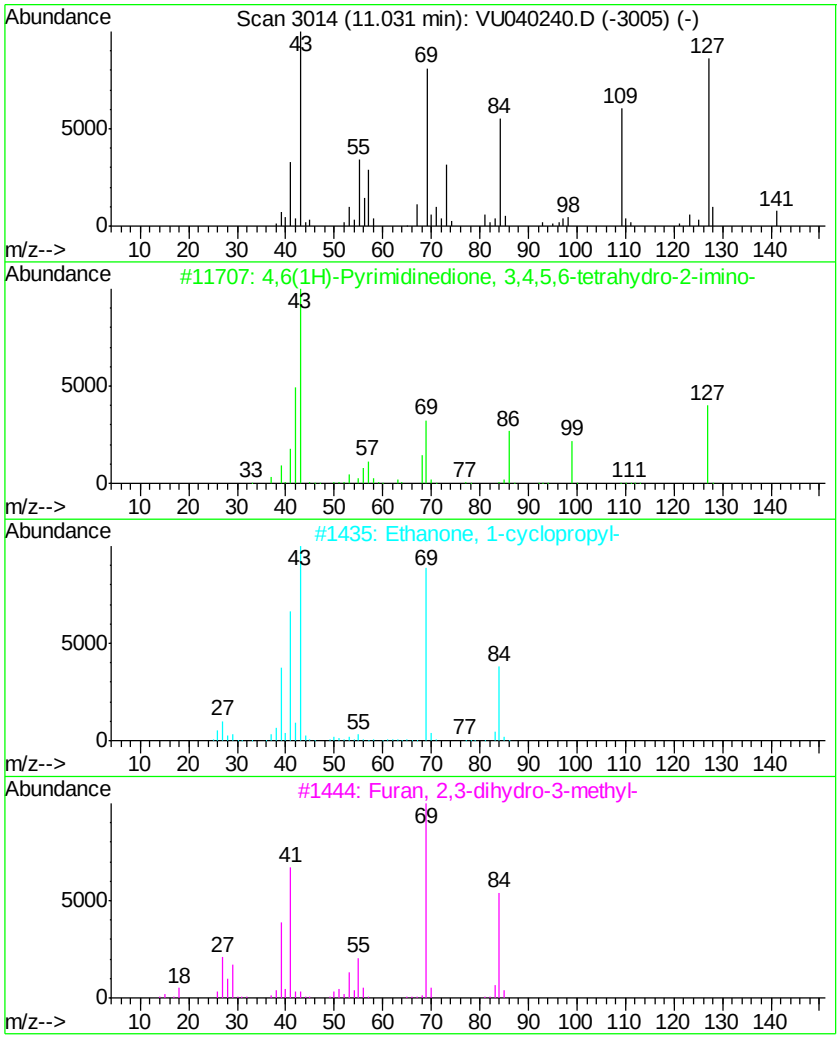
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

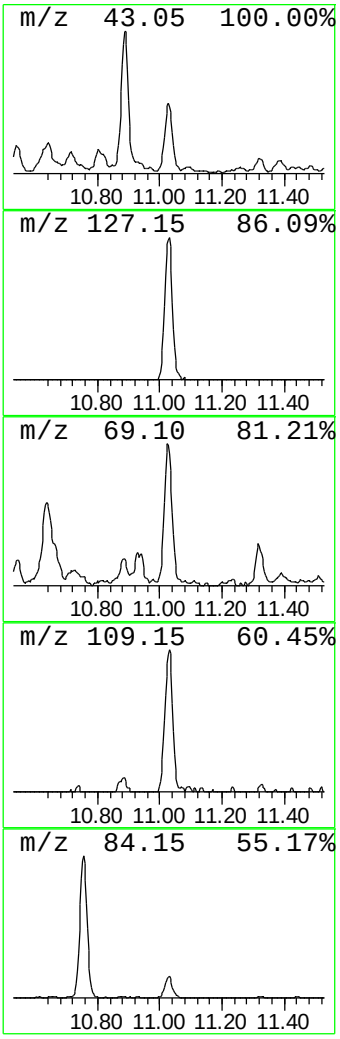
****

Peak Number 11 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.03	1.07 ug/L	94486	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,6(1H)-Pyrimidinedione, 3,4,5,6...	127	C4H5N3O2	004425-67-6	43
2	Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	27
3	Furan, 2,3-dihydro-3-methyl-	84	C5H8O	001708-27-6	27
4	2,5-Pyrrolidinedione, 1-ethyl-	127	C6H9NO2	002314-78-5	25
5	2-Thiophenecarboxaldehyde, oxime	127	C5H5NOS	029683-84-9	25





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

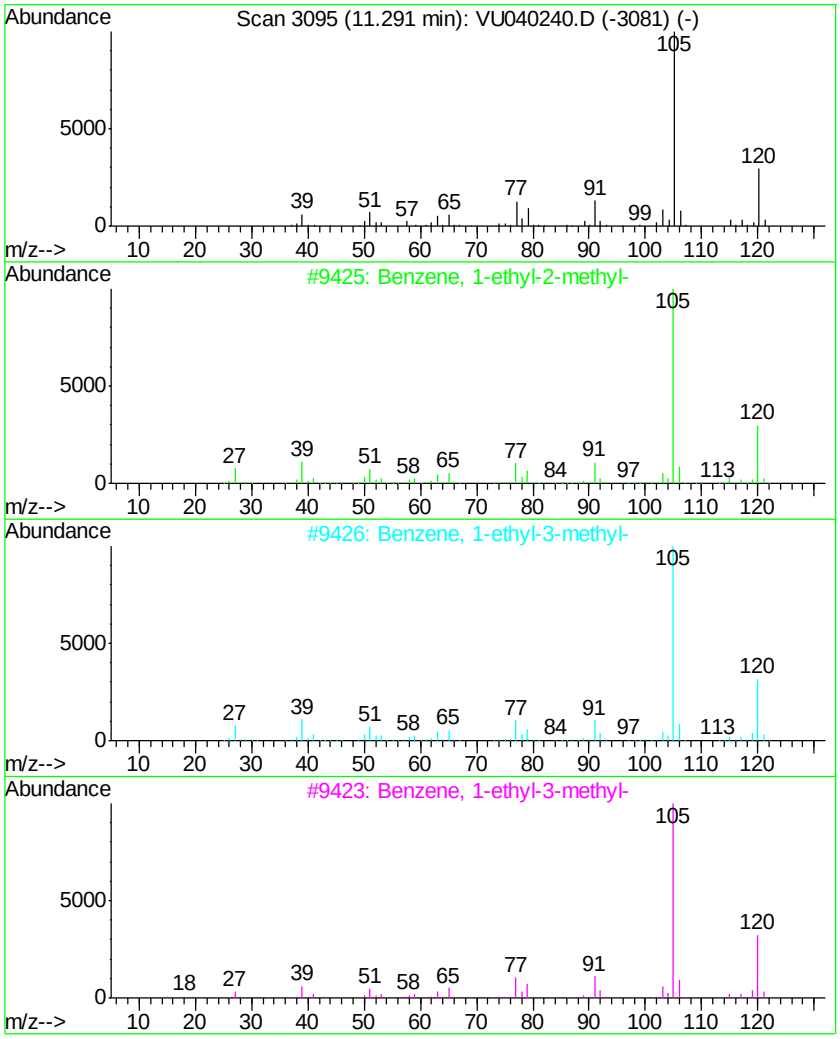
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

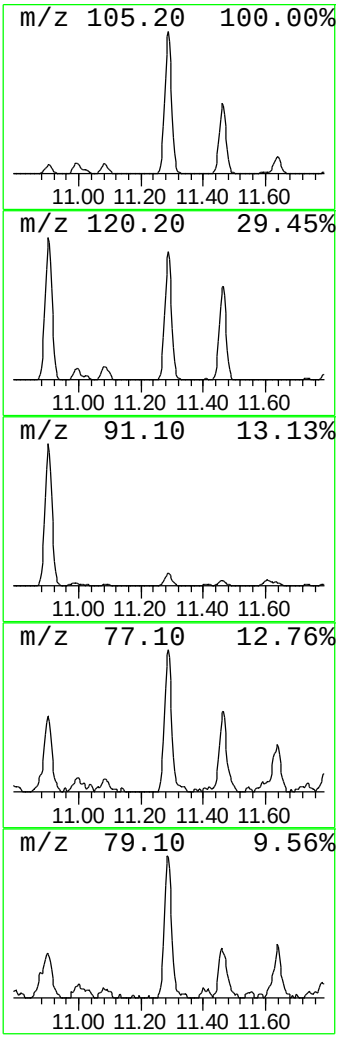
****

Peak Number 12 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	2.44 ug/L	215363	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

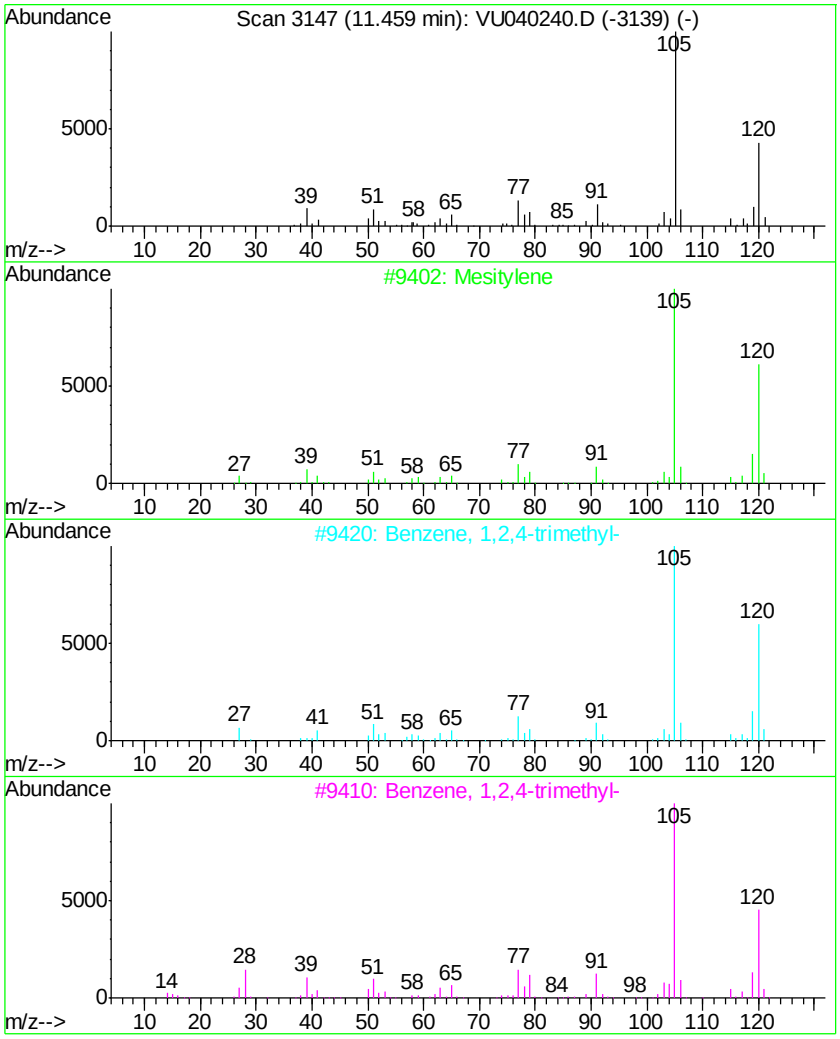
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

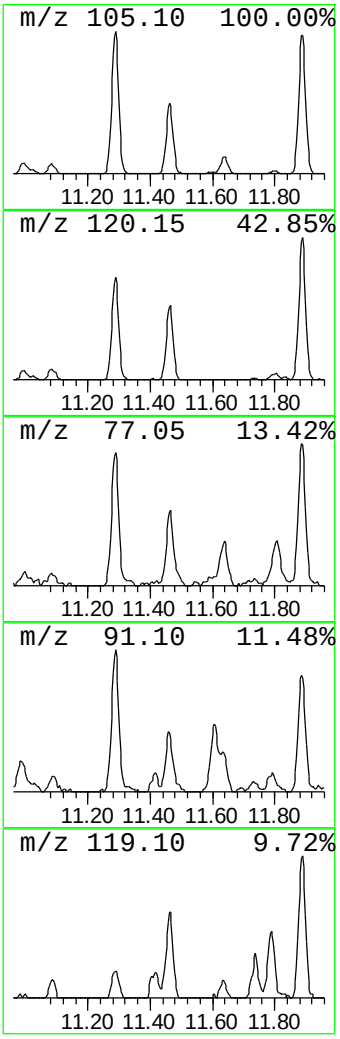
****

Peak Number 13 Mesitylene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	1.41 ug/L	124613	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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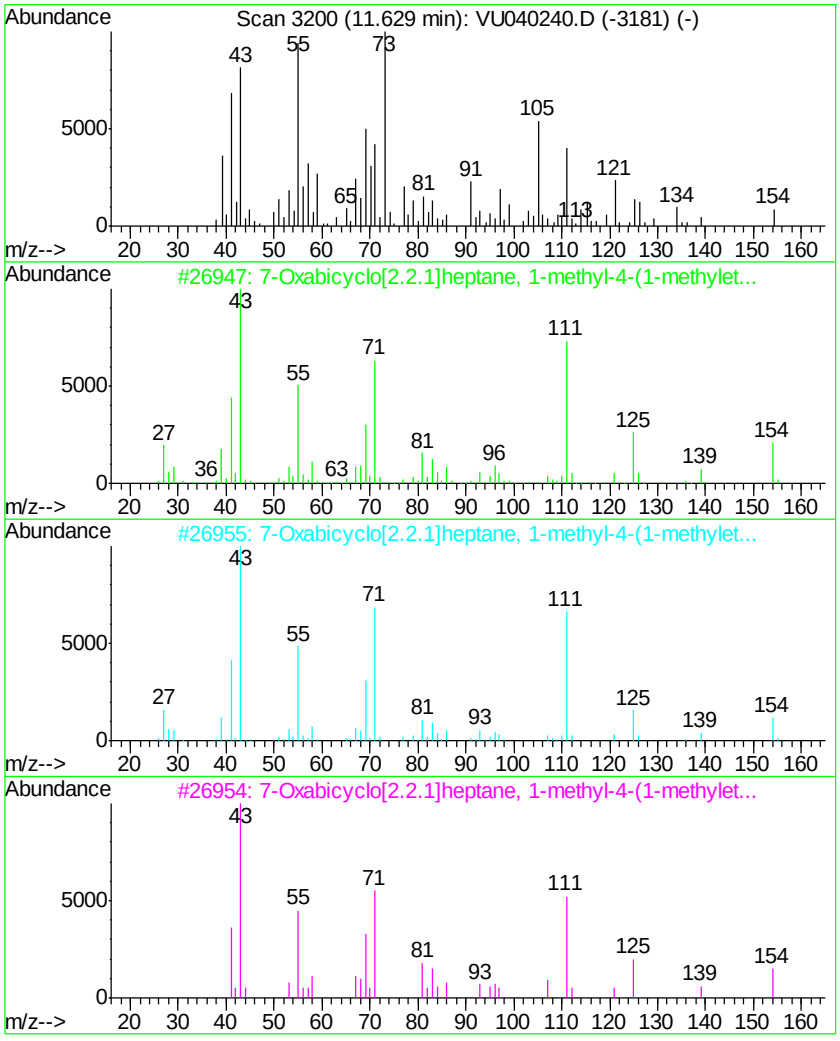
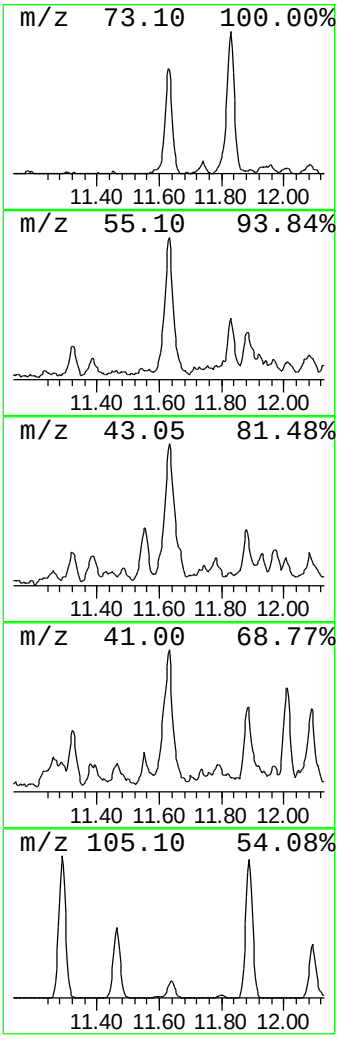
Peak Number 14 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.63	2.97 ug/L	262907	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	46
2	7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	43
3	7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	42
4	Ethanone, 1-(3-ethylcyclobutyl)-	126	C8H14O	056335-71-8	42
5	3-Octen-2-one, (E)-	126	C8H14O	018402-82-9	38





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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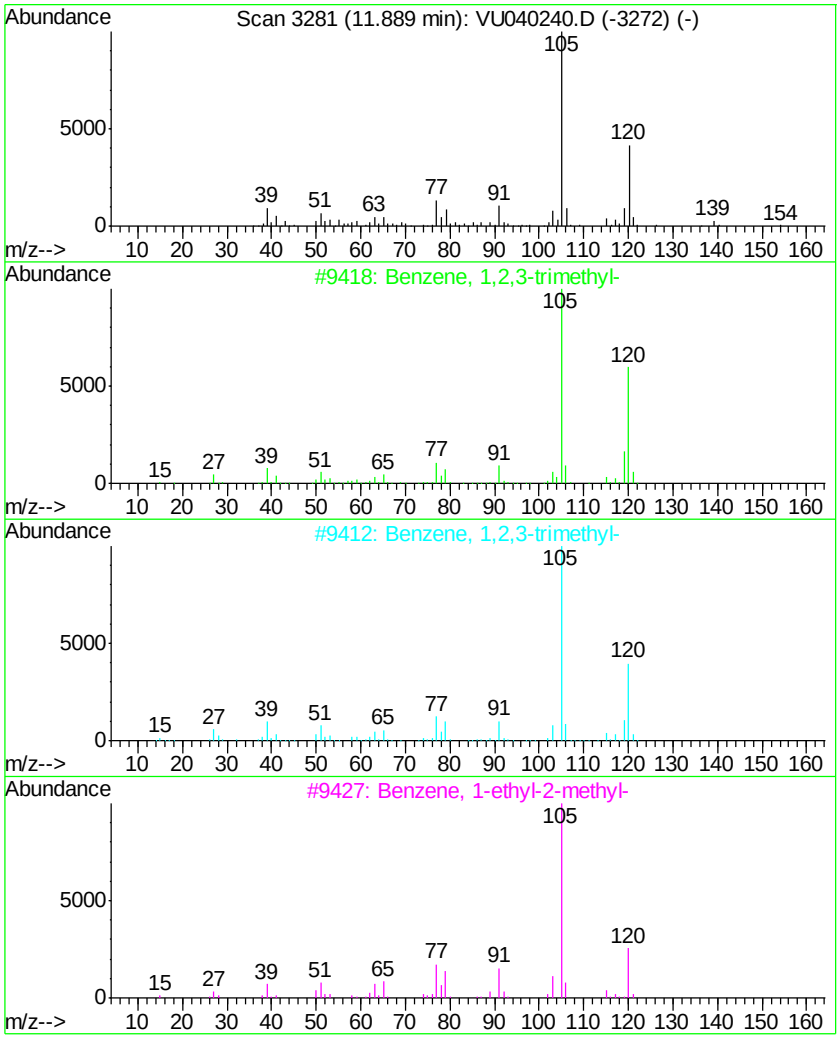
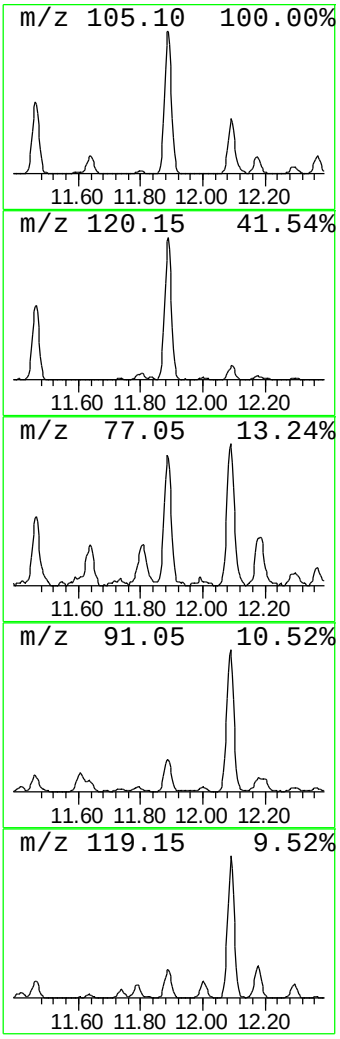
Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	3.06 ug/L	270564	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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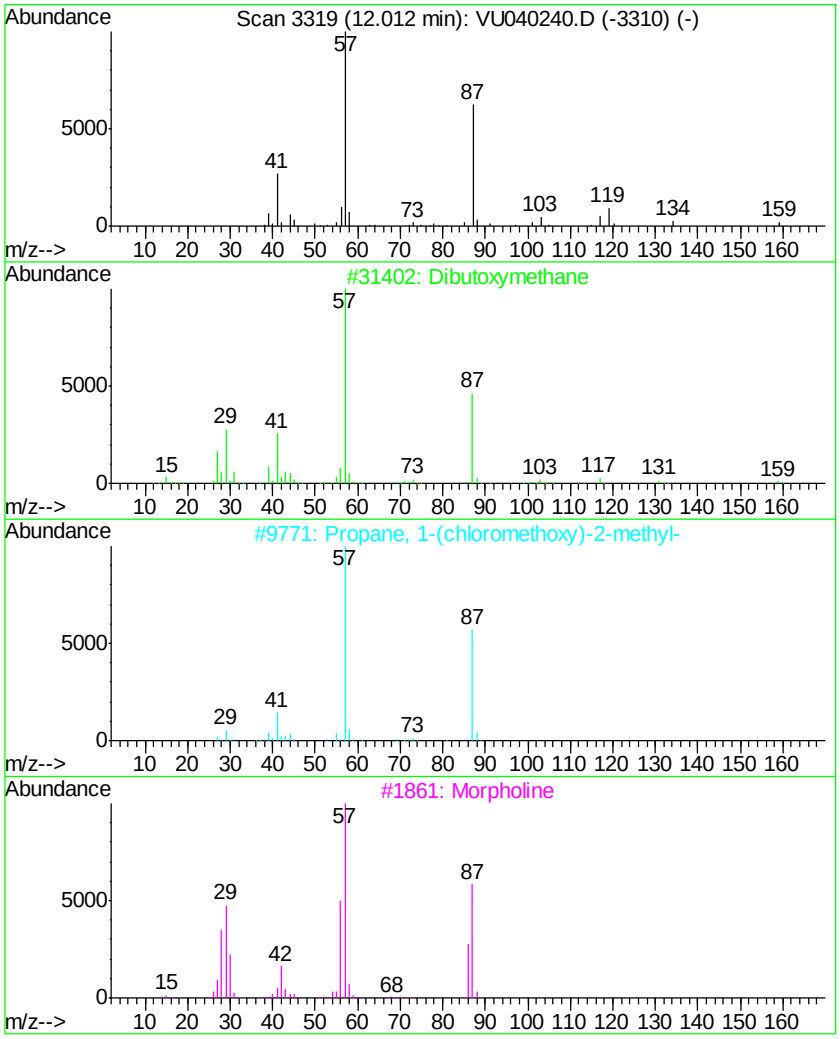
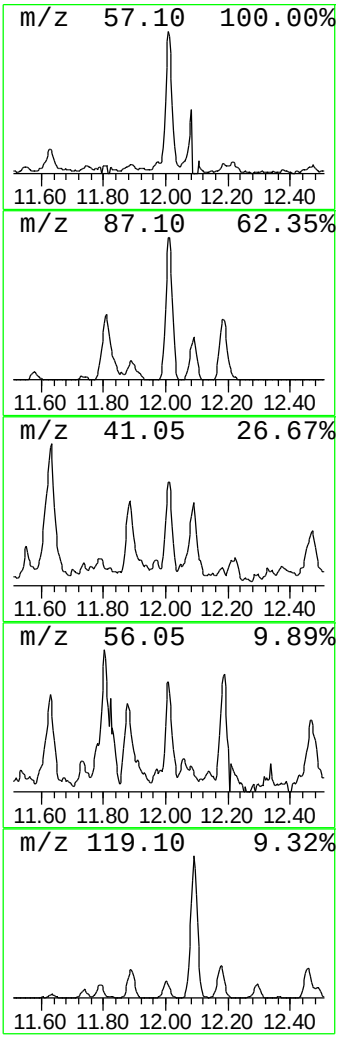
Peak Number 16 unknown-03 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	0.84 ug/L	74112	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibutoxymethane	160	C9H20O2	002568-90-3	45
2	Propane, 1-(chloromethoxy)-2-met...	122	C5H11ClO	034180-11-5	38
3	Morpholine	87	C4H9NO	000110-91-8	38
4	N-Thio-valero-morpholine	187	C9H17NOS	002832-99-7	28
5	4-Heptanol, 2,4-dimethyl-	144	C9H20O	019549-77-0	23





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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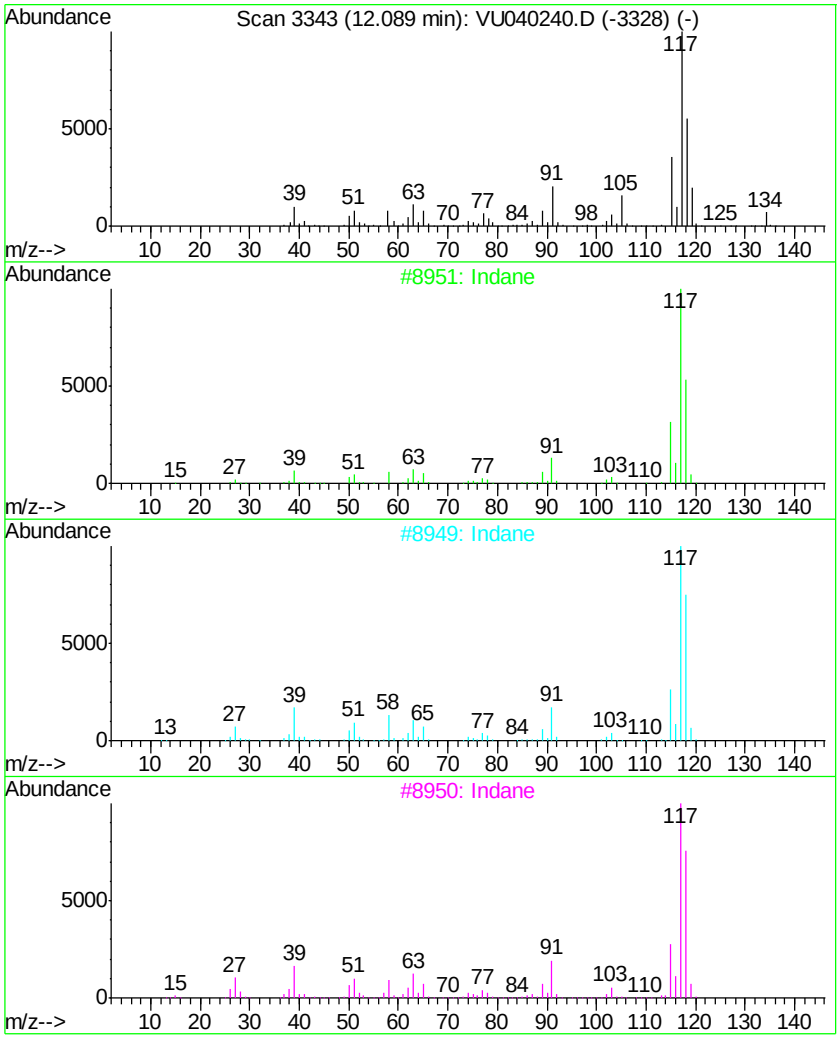
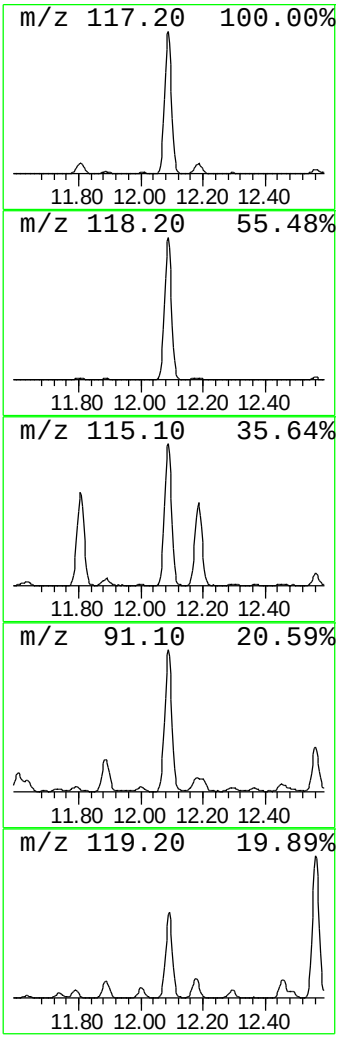
Peak Number 17 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	8.58 ug/L	758289	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	89
2	Indane	118	C9H10	000496-11-7	76
3	Indane	118	C9H10	000496-11-7	76
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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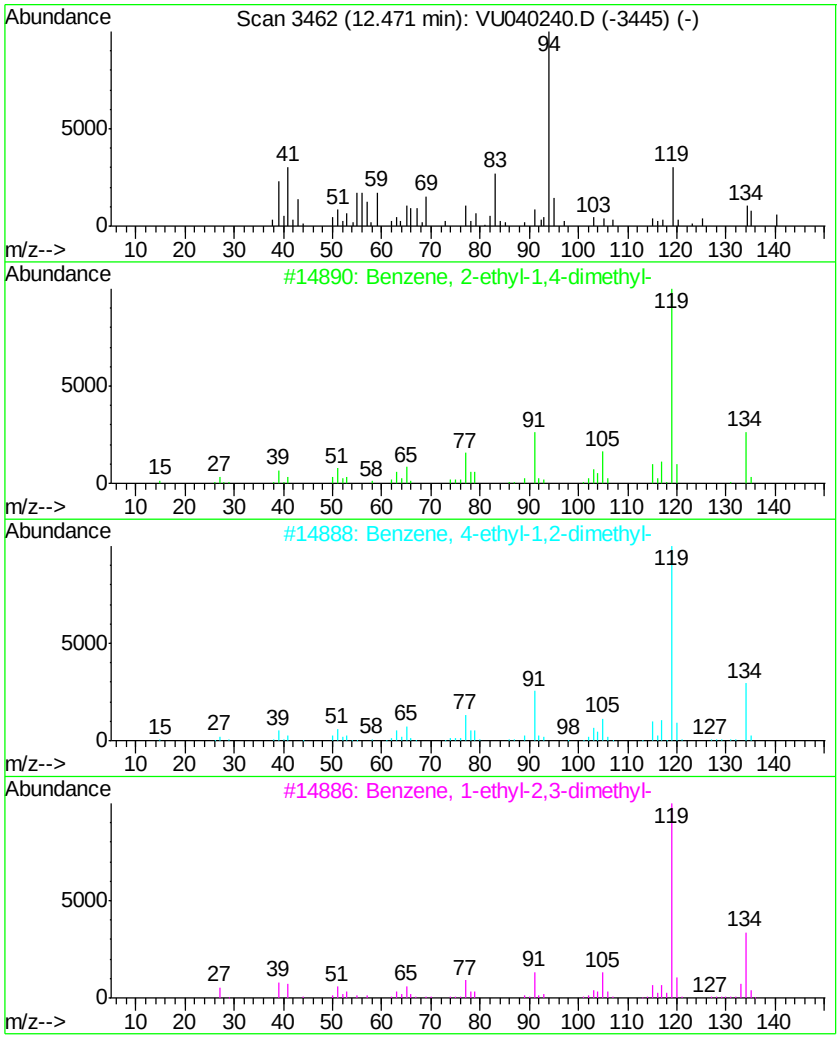
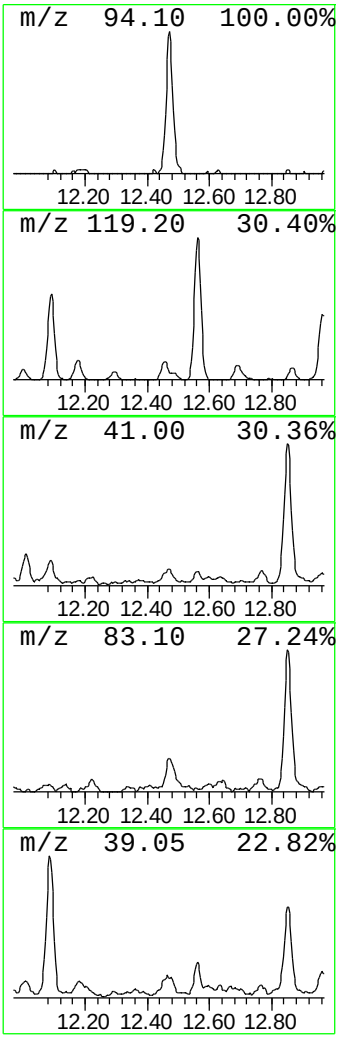
Peak Number 18 unknown-04 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	1.06 ug/L	94000	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	25
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	25
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	25
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	25
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	25





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Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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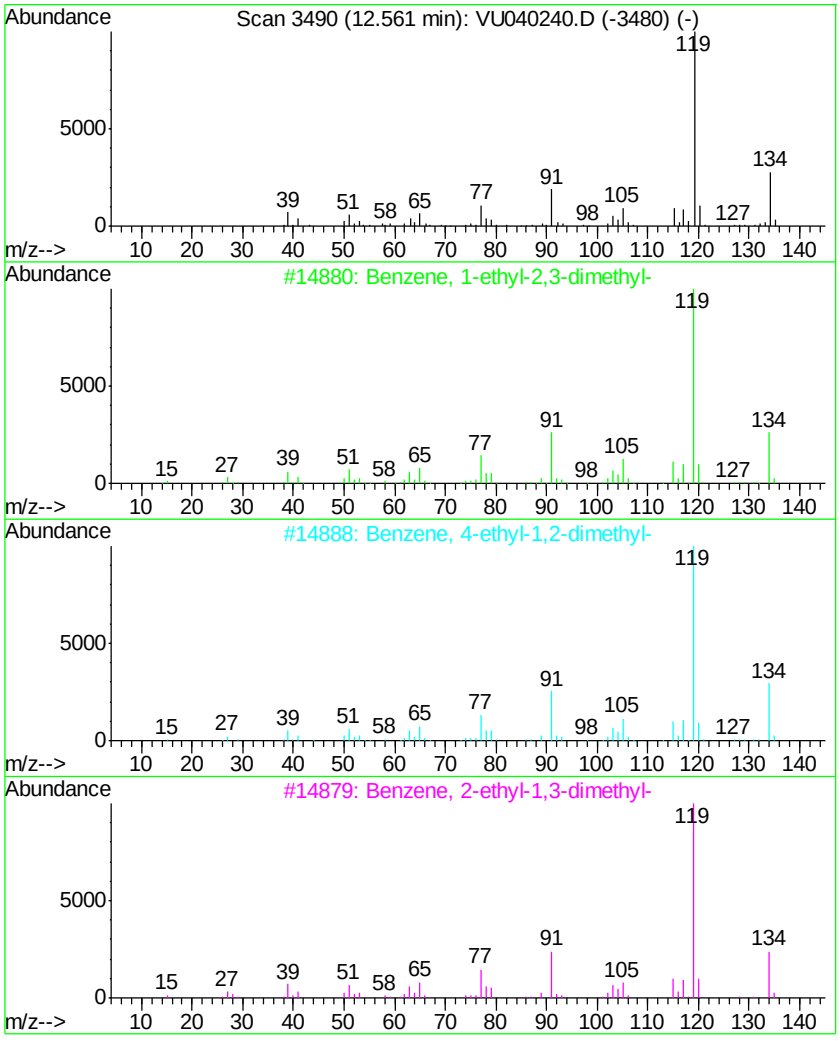
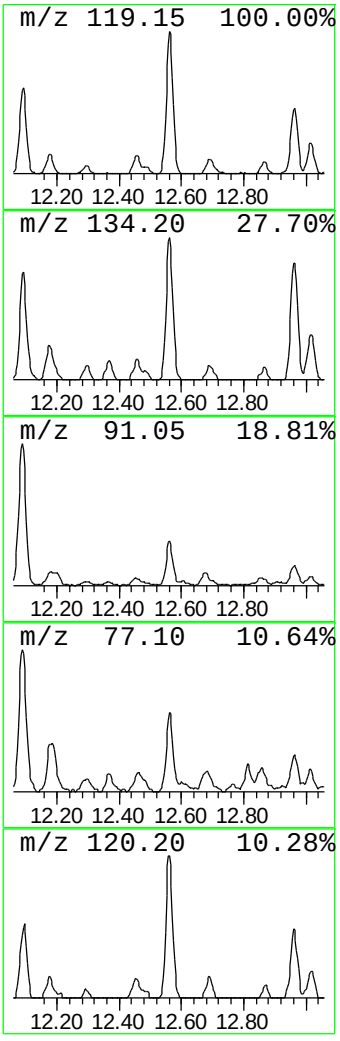
Peak Number 19 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	1.91 ug/L	168613	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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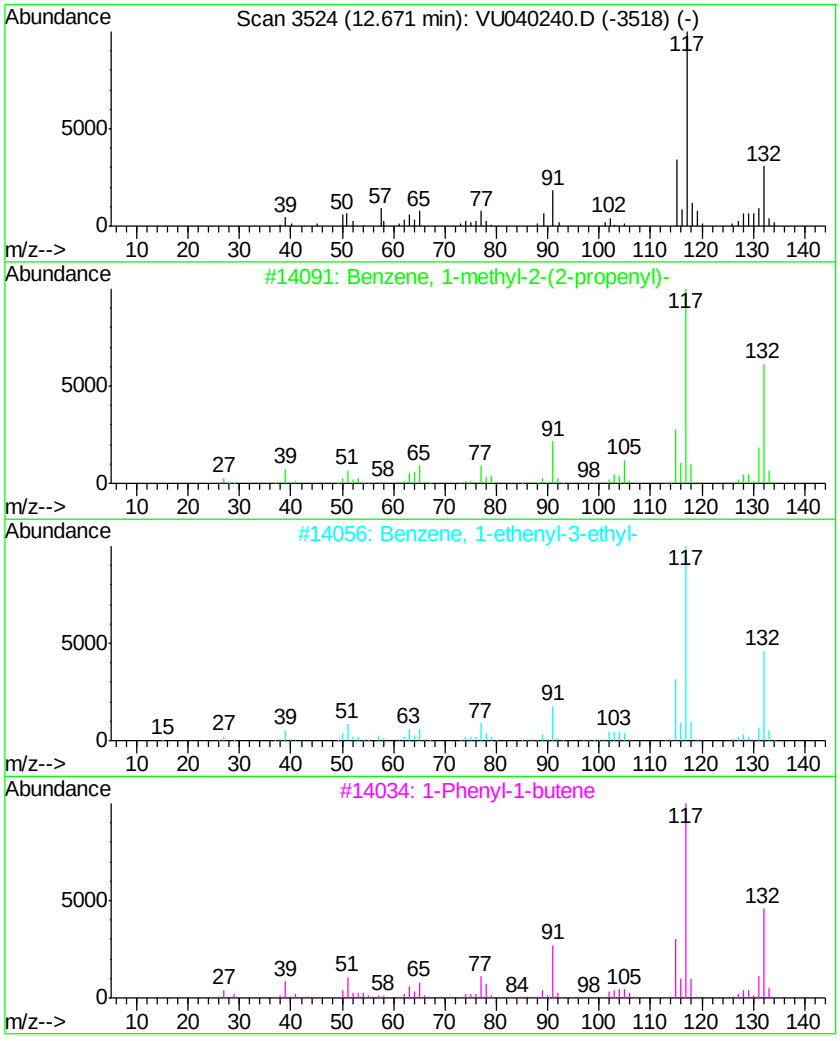
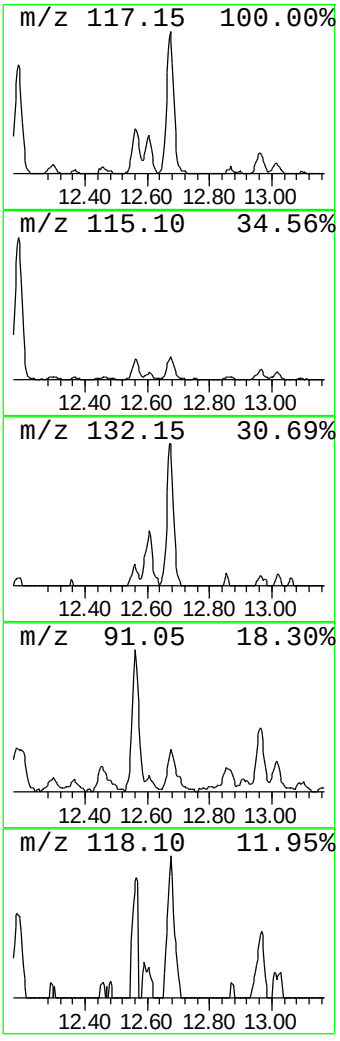
Peak Number 20 Benzene, 1-methyl-2-(2-prop... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	0.85 ug/L	75207	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Indan, 1-methyl-	132	C10H12	000767-58-8	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	86





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

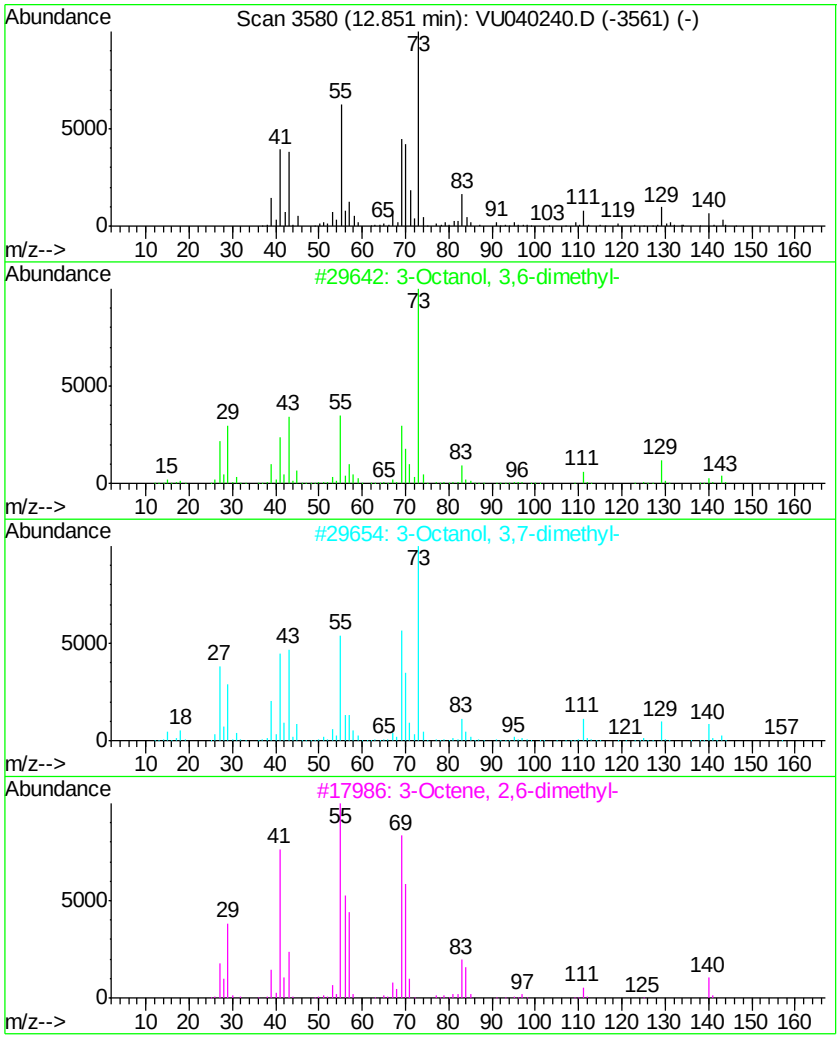
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

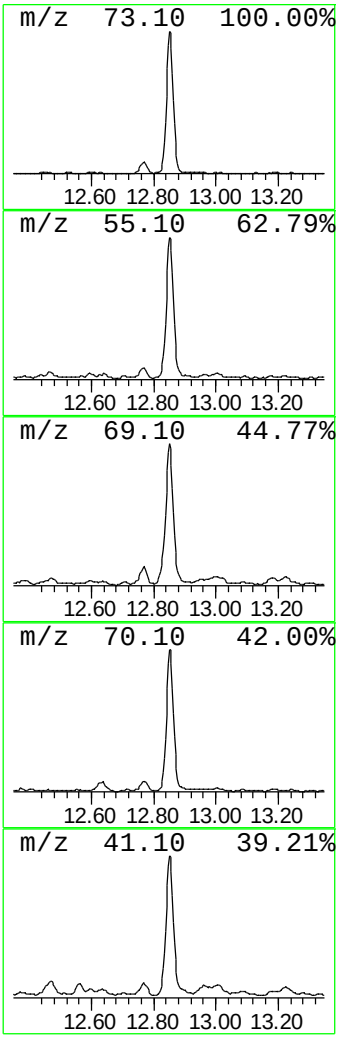
****

Peak Number 21 3-Octanol, 3,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.85	5.53 ug/L	488539	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	53
2	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	47
3	3-Octene, 2,6-dimethyl-	140	C10H20	006874-28-8	46
4	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	43
5	1-Octyn-3-ol	126	C8H14O	000818-72-4	37





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

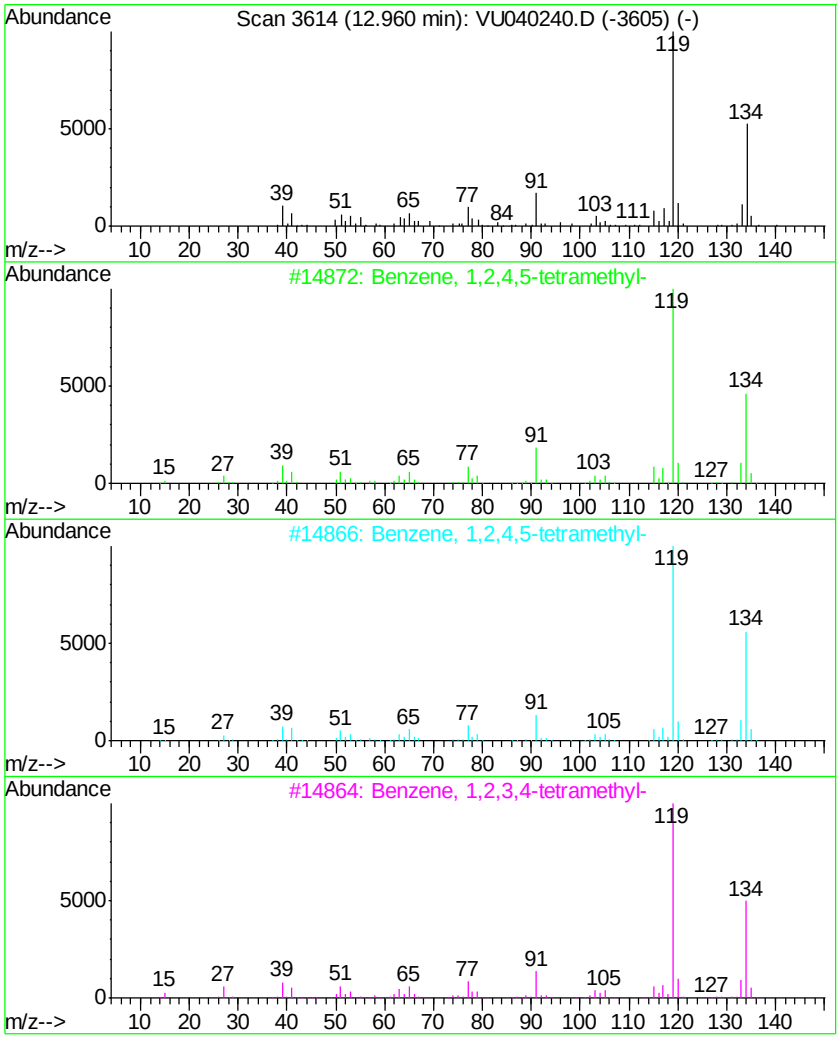
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

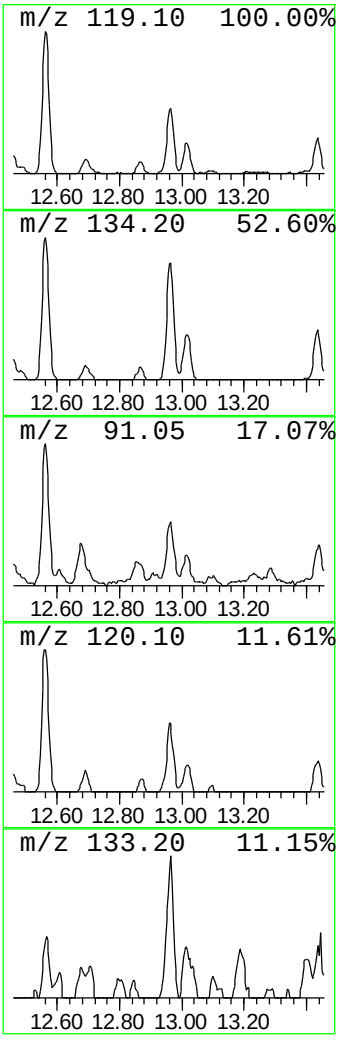
****

Peak Number 22 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	1.18 ug/L	104518	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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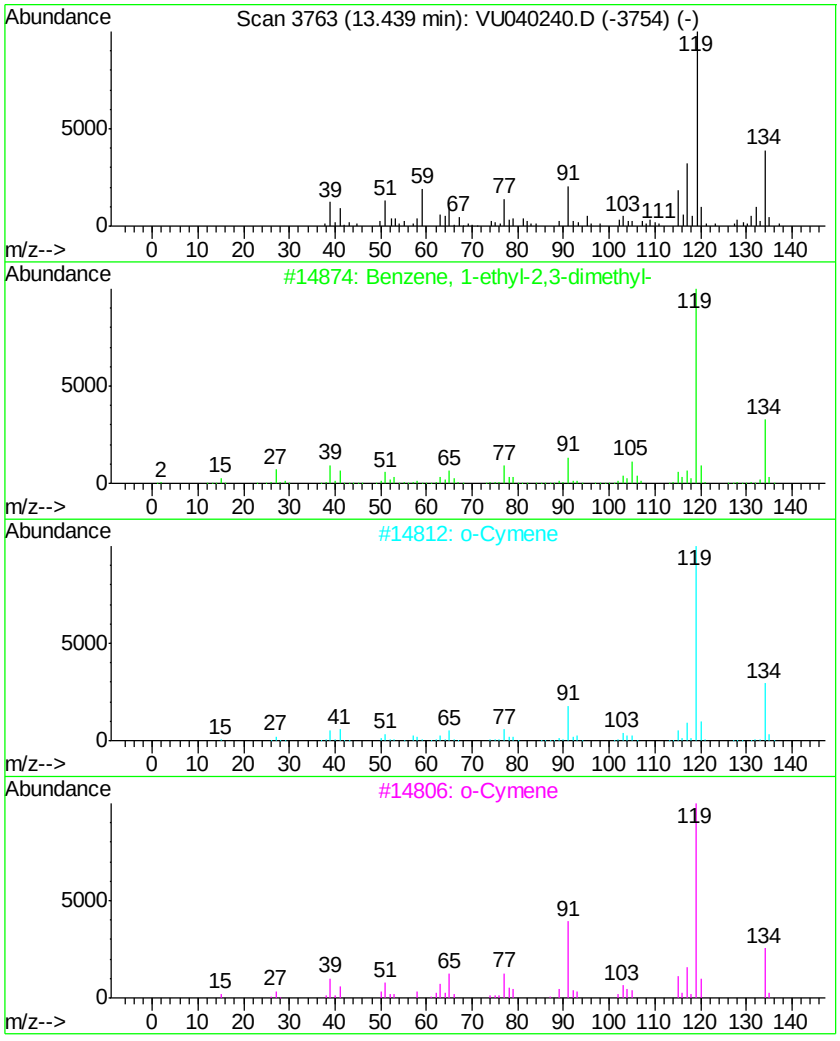
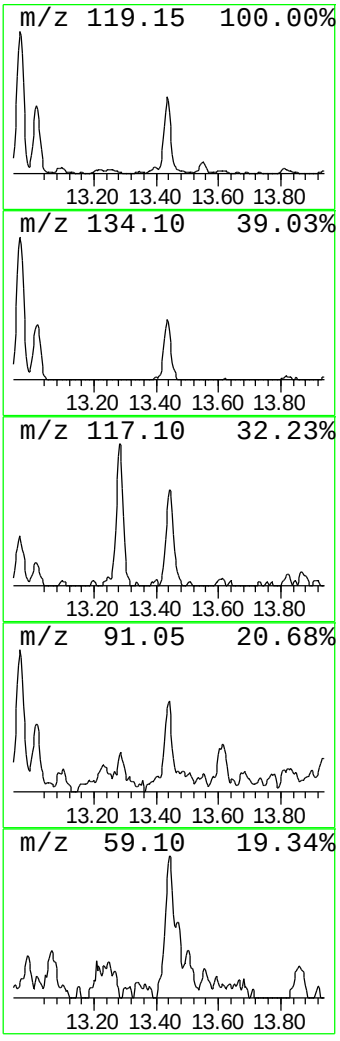
Peak Number 23 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	0.82 ug/L	72760	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76
2	o-Cymene	134	C10H14	000527-84-4	76
3	o-Cymene	134	C10H14	000527-84-4	76
4	o-Cymene	134	C10H14	000527-84-4	70
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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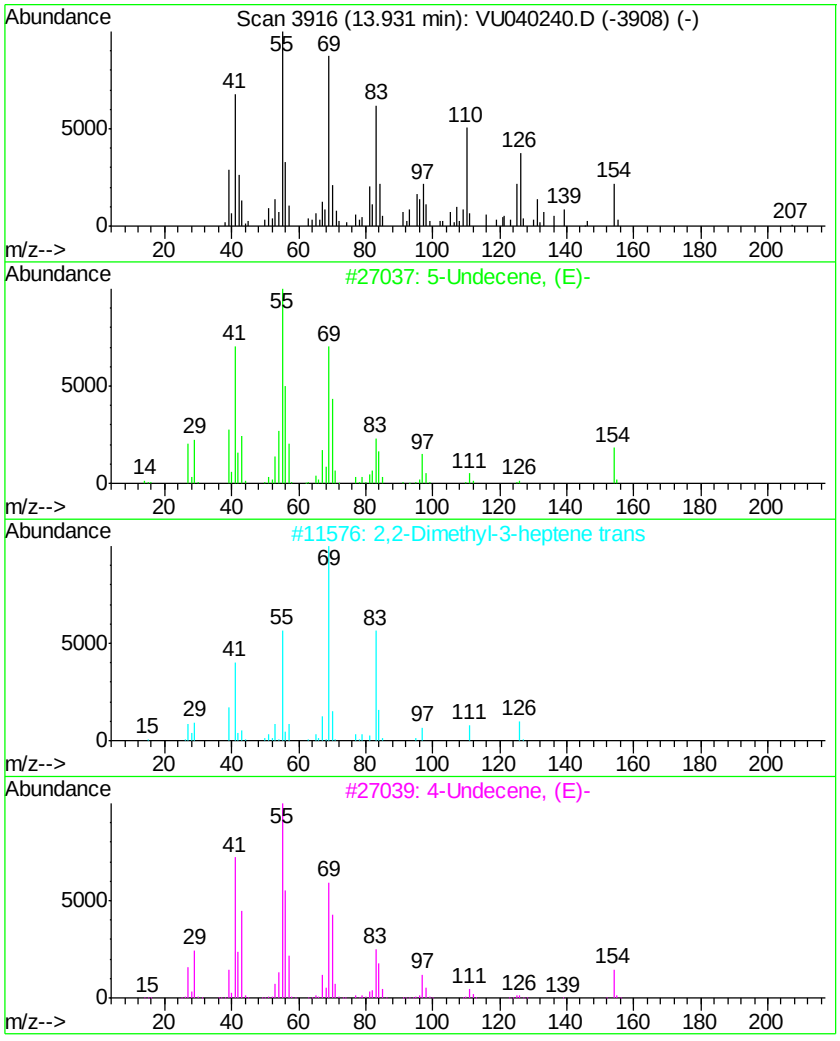
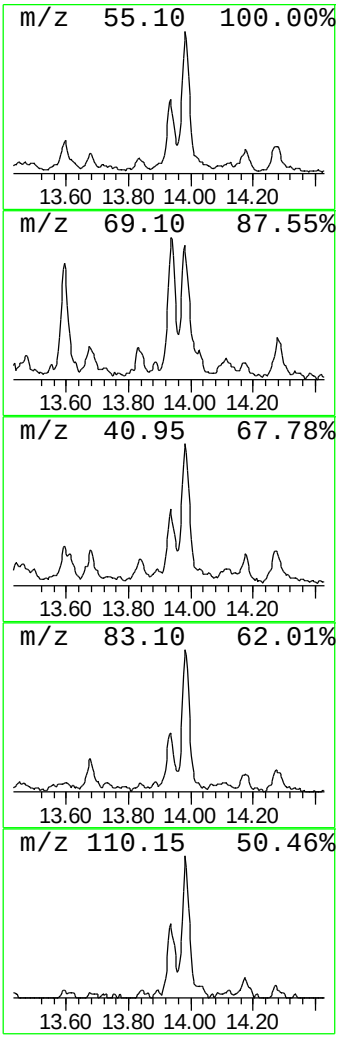
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	0.94 ug/L	83483	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Undecene, (E)-	154	C11H22	000764-97-6	46
2	2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	46
3	4-Undecene, (E)-	154	C11H22	000693-62-9	43
4	1-Decene, 5-methyl-	154	C11H22	054244-79-0	43
5	Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	41





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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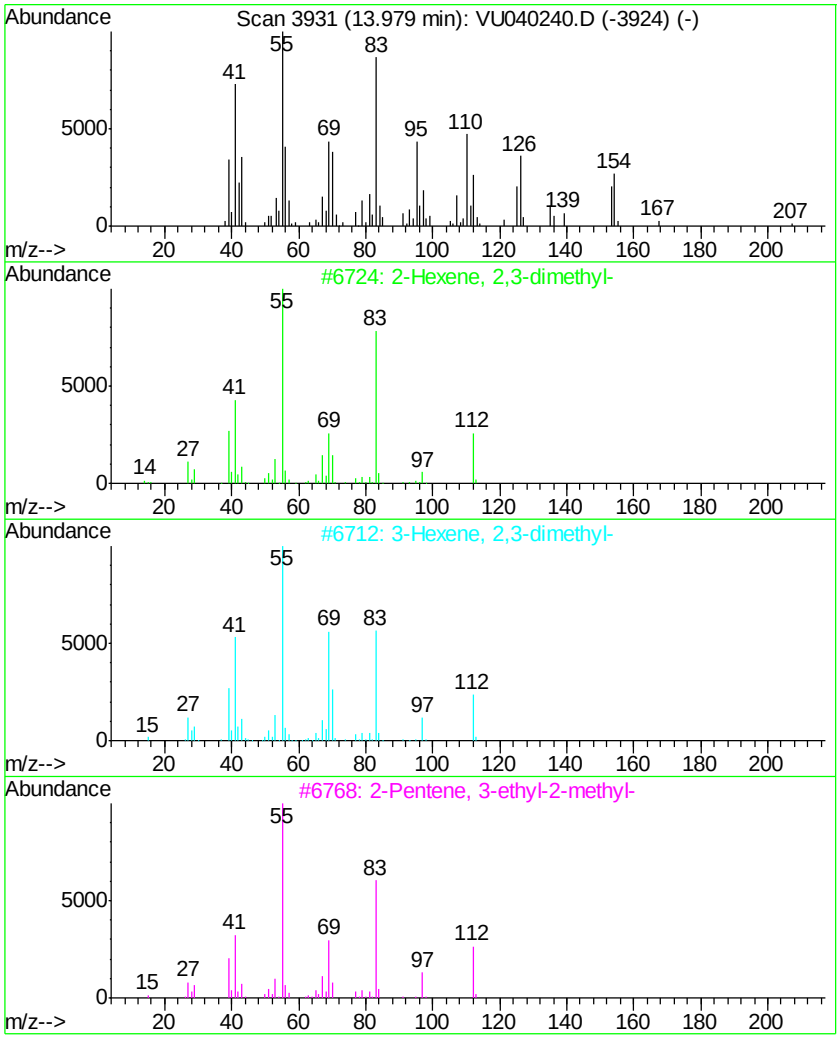
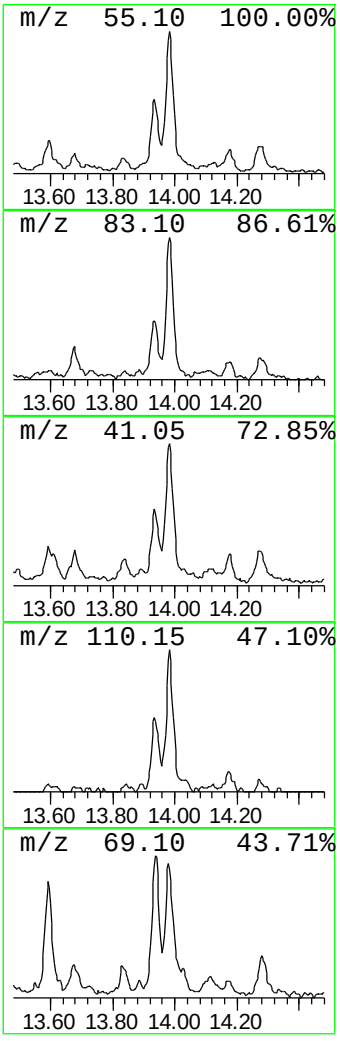
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	1.94 ug/L	171206	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	38
2	3-Hexene, 2,3-dimethyl-	112	C8H16	007145-23-5	38
3	2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	38
4	3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	38
5	2-Octene, 2,3,7-trimethyl-	154	C11H22	033933-75-4	30





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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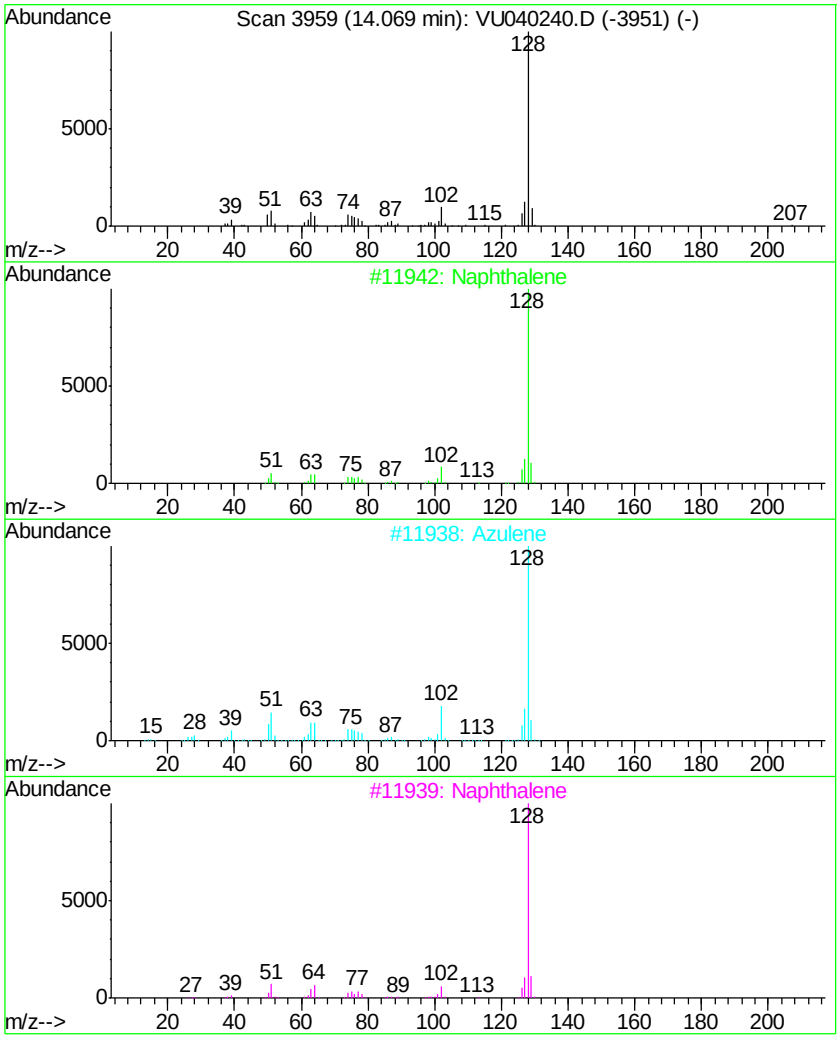
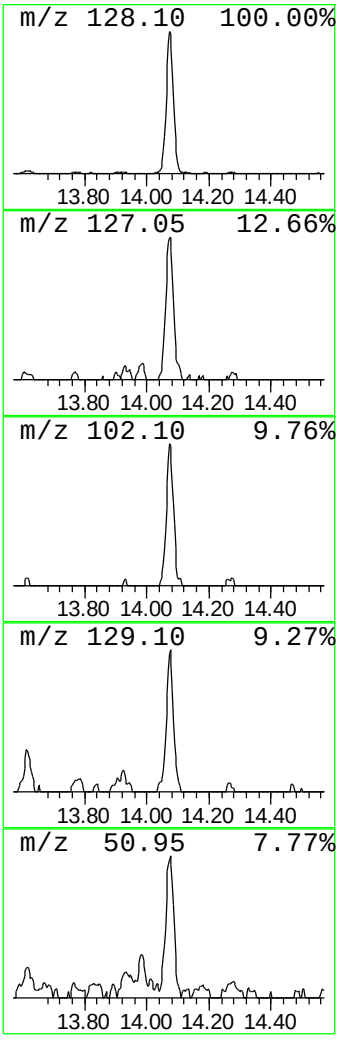
Peak Number 26 Azulene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.07	1.08 ug/L	95872	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Azulene	128	C10H8	000275-51-4	91
3	Naphthalene	128	C10H8	000091-20-3	91
4	Azulene	128	C10H8	000275-51-4	91
5	Azulene	128	C10H8	000275-51-4	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

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

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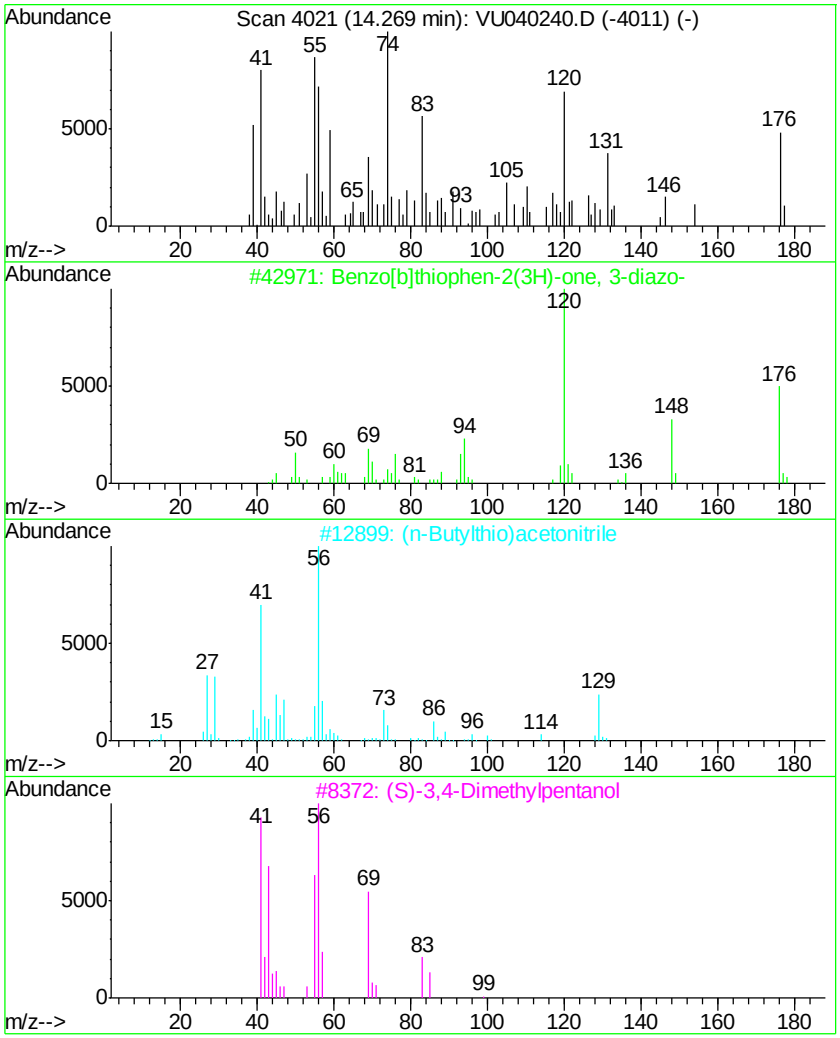
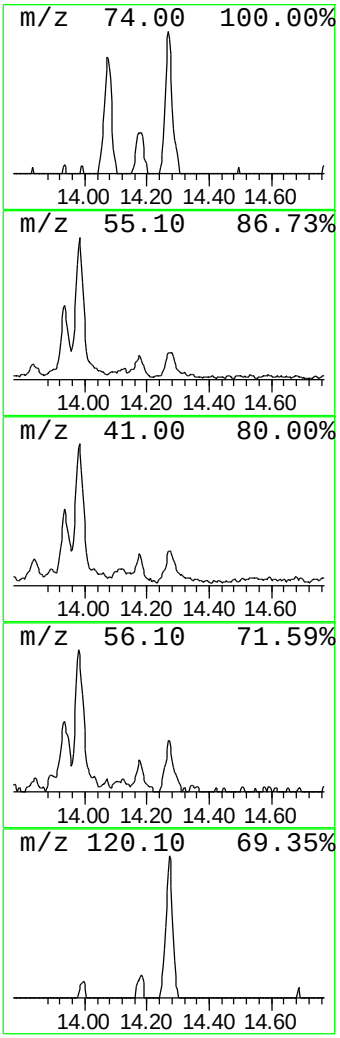
Peak Number 27 unknown-05 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	0.72 ug/L	63604	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophen-2(3H)-one, 3-di...	176	C8H4N2OS	054518-09-1	11
2	(n-Butylthio)acetonitrile	129	C6H11NS	071037-08-6	10
3	(S)-3,4-Dimethylpentanol	116	C7H16O	1000143-83-9	10
4	Thiirane, methyl-	74	C3H6S	001072-43-1	10
5	Formic acid, butyl ester	102	C5H10O2	000592-84-7	10





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.M
Quant Title : TRACE VOA SOM01.0

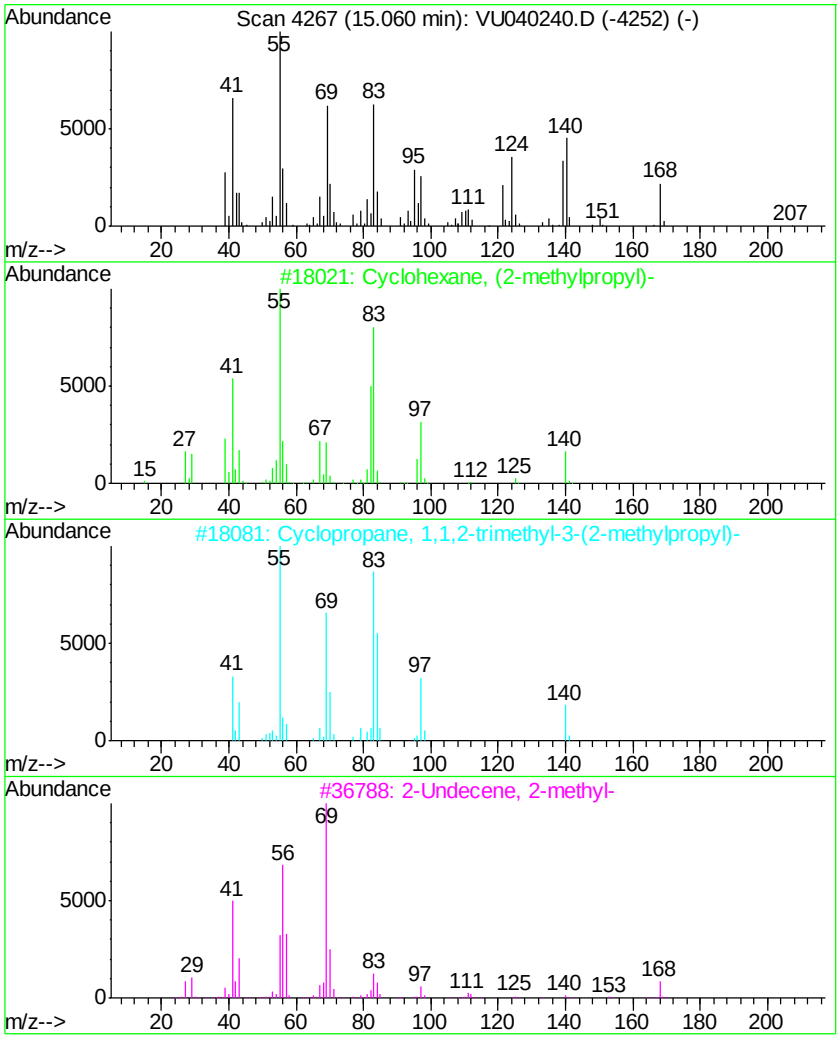
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

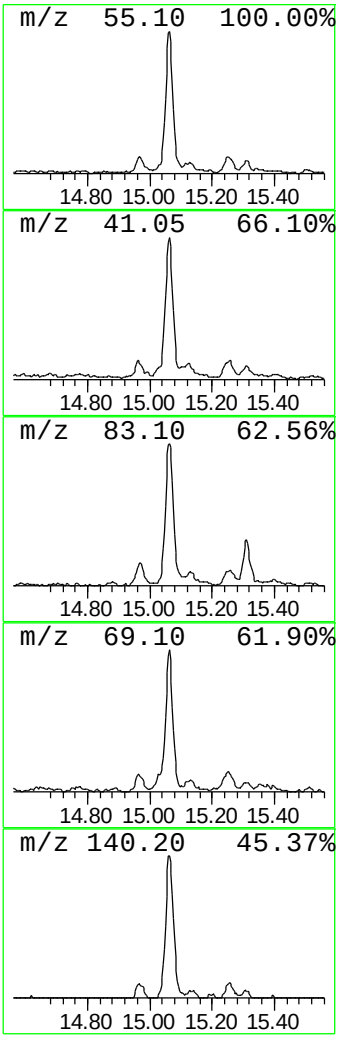
****

Peak Number 28 (DEL) Alkane: Cyclic15.06 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	2.91 ug/L	257113	1,4-Dichlorobenzene-d4	11.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	49
2	Cyclopropane, 1,1,2-trimethyl-3-...	140	C10H20	041977-43-9	43
3	2-Undecene, 2-methyl-	168	C12H24	056888-88-1	42
4	2,3-Dimethyl-2-octene	140	C10H20	019781-18-1	35
5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	30





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU092120\
Data File : VU040240.D
Acq On : 21 Sep 2020 17:45
Operator : SY/MD
Sample : L4046-16
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : BG0Q2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR083120WMA.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...	2.23	1.3	ug/L	85280	1	6.26	338985	5.0
Ethyl ether	2.39	20.2	ug/L	1368860	1	6.26	338985	5.0
(DEL) Alkane: Str...	3.10	1.9	ug/L	127174	1	6.26	338985	5.0
(DEL) Alkane: Str...	3.72	2.8	ug/L	192711	1	6.26	338985	5.0
Diisopropyl ether	4.01	75.8	ug/L	5139160	1	6.26	338985	5.0
(DEL) Alkane: Cyc...	4.55	15.1	ug/L	1025530	1	6.26	338985	5.0
Furan, tetrahydro...	7.72	2.0	ug/L	135824	1	6.26	338985	5.0
3-Pentanol, 2,3,4...	10.64	1.2	ug/L	108132	3	11.81	441885	5.0
Benzene, propyl-	10.90	5.3	ug/L	467390	3	11.81	441885	5.0
unknown-01	11.03	1.1	ug/L	94486	3	11.81	441885	5.0
Benzene, 1-ethyl-...	11.29	2.4	ug/L	215363	3	11.81	441885	5.0
Mesitylene	11.46	1.4	ug/L	124613	3	11.81	441885	5.0
unknown-02	11.63	3.0	ug/L	262907	3	11.81	441885	5.0
Benzene, 1,2,3-tr...	11.89	3.1	ug/L	270564	3	11.81	441885	5.0
unknown-03	12.01	0.8	ug/L	74112	3	11.81	441885	5.0
Indane	12.09	8.6	ug/L	758289	3	11.81	441885	5.0
unknown-04	12.47	1.1	ug/L	94000	3	11.81	441885	5.0
Benzene, 2-ethyl-...	12.56	1.9	ug/L	168613	3	11.81	441885	5.0
Benzene, 1-methyl...	12.67	0.8	ug/L	75207	3	11.81	441885	5.0
3-Octanol, 3,6-di...	12.85	5.5	ug/L	488539	3	11.81	441885	5.0
Benzene, 1,2,4,5-...	12.96	1.2	ug/L	104518	3	11.81	441885	5.0
Benzene, 1-ethyl-...	13.44	0.8	ug/L	72760	3	11.81	441885	5.0
(DEL) Alkane: Str...	13.93	0.9	ug/L	83483	3	11.81	441885	5.0
(DEL) Alkane: Str...	13.98	1.9	ug/L	171206	3	11.81	441885	5.0
Azulene	14.07	1.1	ug/L	95872	3	11.81	441885	5.0
unknown-05	14.27	0.7	ug/L	63604	3	11.81	441885	5.0
(DEL) Alkane: Cyc...	15.06	2.9	ug/L	257113	3	11.81	441885	5.0