LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

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_V\METHOD\SOMVTR111519WMA.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	1.315	62	70	80	rBV	259980	266083	12.93%	1.303%
2	1.582	145	153	164	rVB	198882	222366	10.80%	1.089%
3	1.650	164	174	184	rBV	649026	797745	38.75%	3.907%
4	1.823	220	228	238	rVB	174998	228239	11.09%	1.118%
5	1.920	251	258	270	rVB	41401	55003	2.67%	0.269%
6	2.045	290	297	306	rVB	45155	57177	2.78%	0.280%
7	2.129	312	323	344	rVB	373281	609442	29.60%	2.985%
8	2.531	435	448	450	rBV2	58647	96499	4.69%	0.473%
9	2.605	465	471	489	rVB2	56988	109627	5.33%	0.537%
10	2.791	518	529	542	rBV2	66845	133438	6.48%	0.653%
11	2.978	577	587	598	rVB5	14423	31270	1.52%	0.153%
12	3.074	607	617	627	rBV5	14779	29298	1.42%	0.143%
13	3.257	662	674	683	rBV5	15393	32067	1.56%	0.157%
14	3.312	683	691	706	rVB2	22979	47132	2.29%	0.231%
15	3.409	709	721	728	rBV4	15019	30707	1.49%	0.150%
16	3.624	776	788	800	rBV3	21647	45469	2.21%	0.223%
17	3.807	825	845	862	rBV2	97836	253522	12.31%	1.242%
18	3.936	873	885	924	rBV	157274	479637	23.30%	2.349%
19	4.396	1009	1028	1050	rBV2	262002	695363	33.78%	3.405%
20	4.505	1053	1062	1078	rVB3	15988	38469	1.87%	0.188%
21	4.724	1114	1130	1148	rBV4	51372	130684	6.35%	0.640%
22	5.093	1224	1245	1276	rBV2	594672	1594000	77.43%	7.806%
23	5.270	1292	1300	1316	rVB2	11089	31388	1.52%	0.154%
24	5.659	1408	1421	1457	rBV	381521	814118	39.55%	3.987%
25	6.113	1547	1562	1578	rBV	337804	703140	34.16%	3.444%
26	6.849	1780	1791	1798	rVB3	12303	24367	1.18%	0.119%
27	7.026	1834	1846	1862	rBV	157235	290595	14.12%	1.423%
28	7.357	1937	1949	1961	rBV	727508	1242487	60.35%	6.085%
29	7.428	1961	1971	1992	rVB	363467	634891	30.84%	3.109%
30	7.662	2029	2044	2057	rBV2	95991	166529	8.09%	0.816%
31	8.132	2177	2190	2222	rBV2	544540	1235725	60.03%	6.052%
32	8.891	2412	2426	2444	rBV	558490	917136	44.55%	4.492%
33	9.051	2464	2476	2490	rBV	392858	619246	30.08%	3.033%
34	9.177	2502	2515	2540	rBV	1302274	2058649	100.00%	10.082%

LSC Area Percent Report

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Operator : SY/MD
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ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

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_V\METHOD\SOMVTR111519WMA.M
Title : TRACE VOA SOM01.0

35	9.582	2626	2641	2661	rBV	302135	483939	23.51%	2.370%
36	9.971	2751	2762	2772	rBV	24918	39855	1.94%	0.195%
37	10.257	2836	2851	2863	rBV2	239456	381533	18.53%	1.869%
38	10.392	2880	2893	2909	rBV	82022	139715	6.79%	0.684%
39	10.489	2911	2923	2941	rVV	264254	580615	28.20%	2.843%
40	10.579	2941	2951	2967	rVB	132236	207780	10.09%	1.018%
41	10.778	3001	3013	3025	rBV	113393	186695	9.07%	0.914%
42	10.955	3057	3068	3086	rBV	532033	807270	39.21%	3.954%
43	11.289	3160	3172	3189	rBV	618600	957634	46.52%	4.690%
44	11.379	3189	3200	3211	rVB	111431	173633	8.43%	0.850%
45	11.572	3247	3260	3270	rBV2	84988	158884	7.72%	0.778%
46	11.666	3278	3289	3310	rVB	741587	1192768	57.94%	5.841%
47	11.955	3369	3379	3383	rBV2	20448	30609	1.49%	0.150%
48	11.984	3383	3388	3398	rVB2	17219	26358	1.28%	0.129%
49	12.058	3398	3411	3418	rBV	29554	49025	2.38%	0.240%
50	12.157	3433	3442	3451	rBV2	26098	43356	2.11%	0.212%
51	12.457	3526	3535	3542	rBV2	18311	26889	1.31%	0.132%
52	12.508	3542	3551	3562	rVB2	28359	48422	2.35%	0.237%
53	12.768	3622	3632	3650	rVB4	23404	40773	1.98%	0.200%
54	12.926	3672	3681	3693	rVB3	42560	68450	3.32%	0.335%
55	13.550	3859	3875	3890	rBV2	27958	53377	2.59%	0.261%

Sum of corrected areas: 20419088

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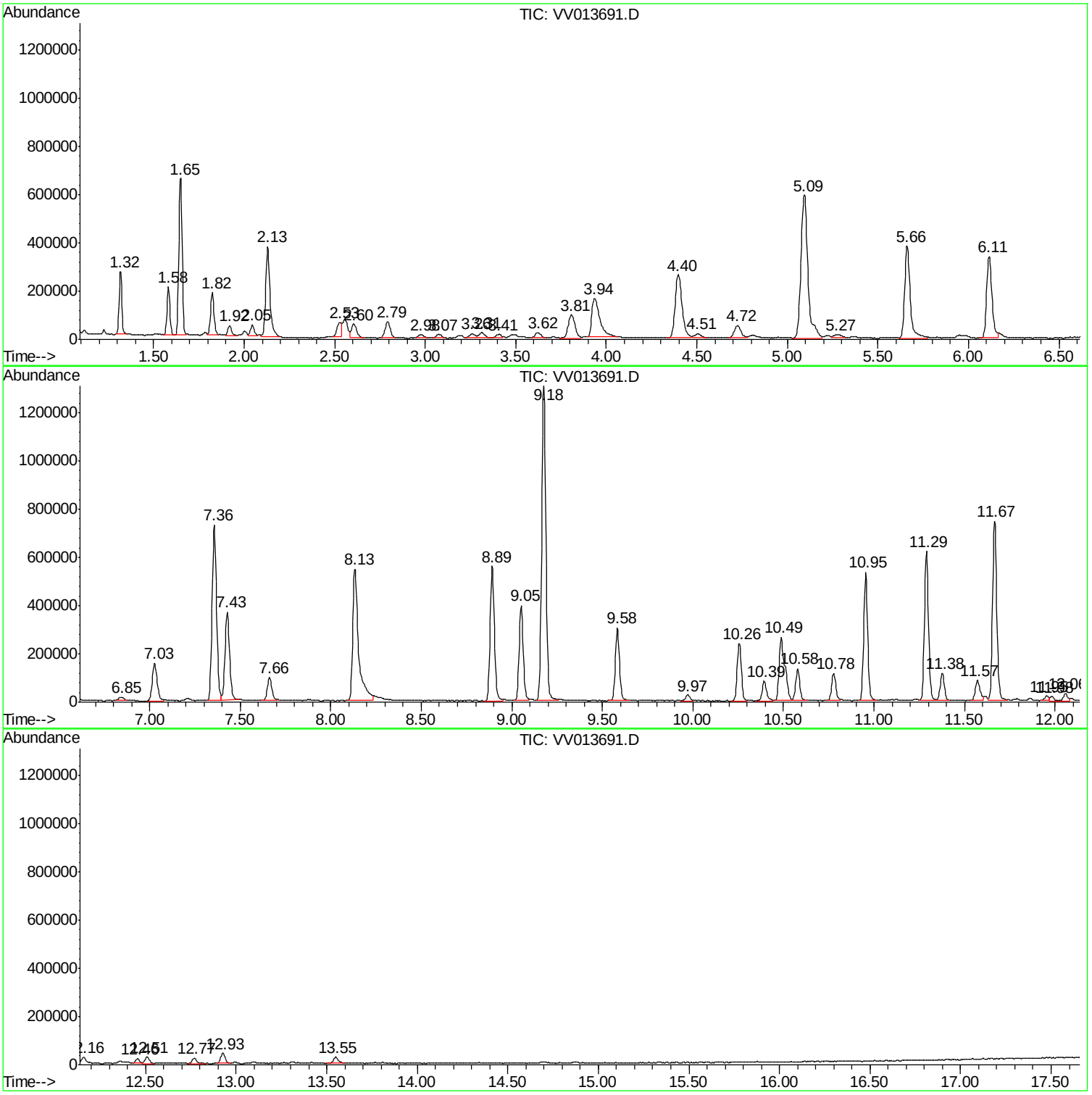
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.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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Quant Title : TRACE VOA SOM01.0

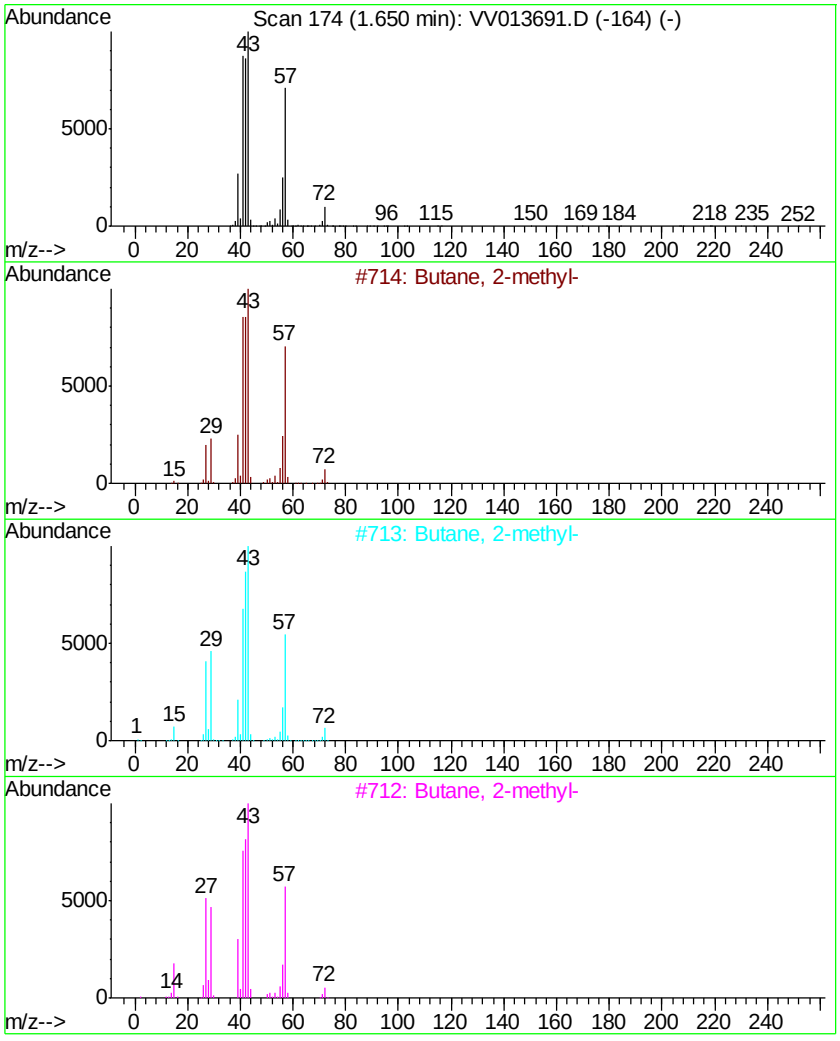
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

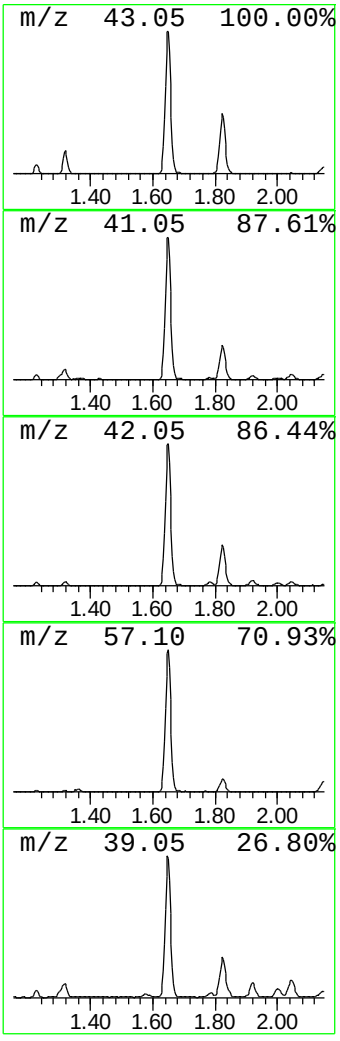
****

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.65	4.90 ug/L	797745	1,4-Difluorobenzene	5.66

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	90
4	3-Buten-1-ol	72	C4H8O	000627-27-0	38
5	Pentane	72	C5H12	000109-66-0	37





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

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

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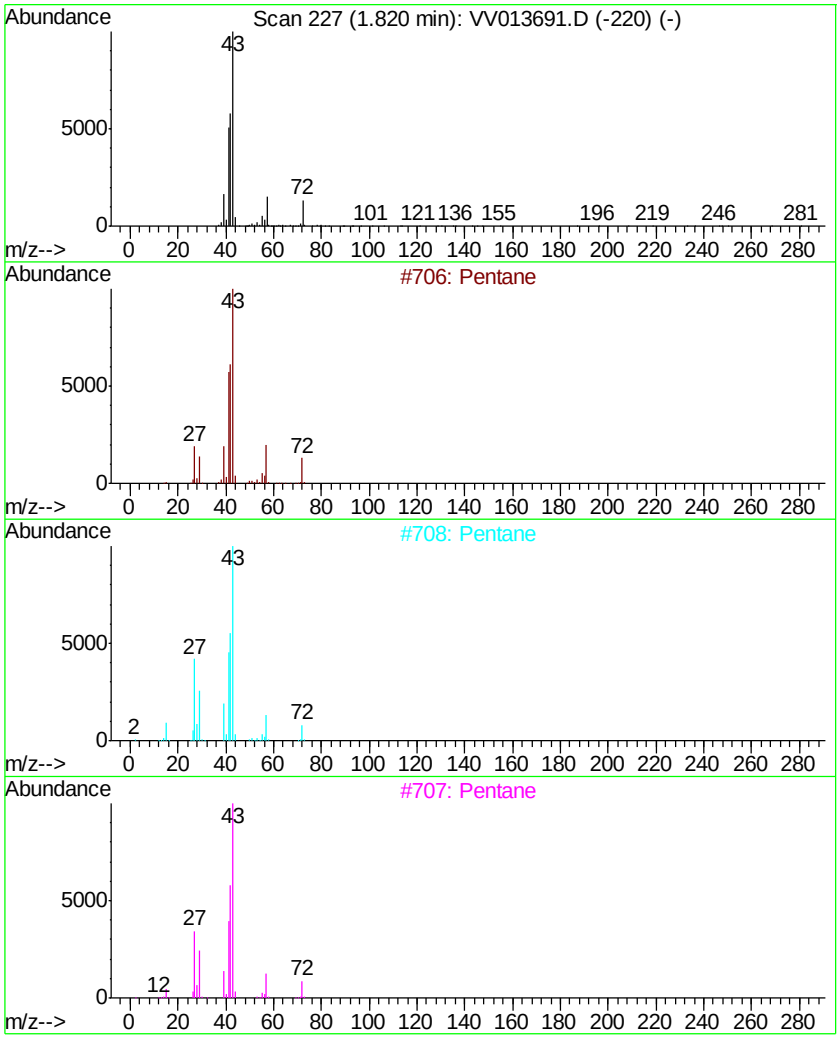
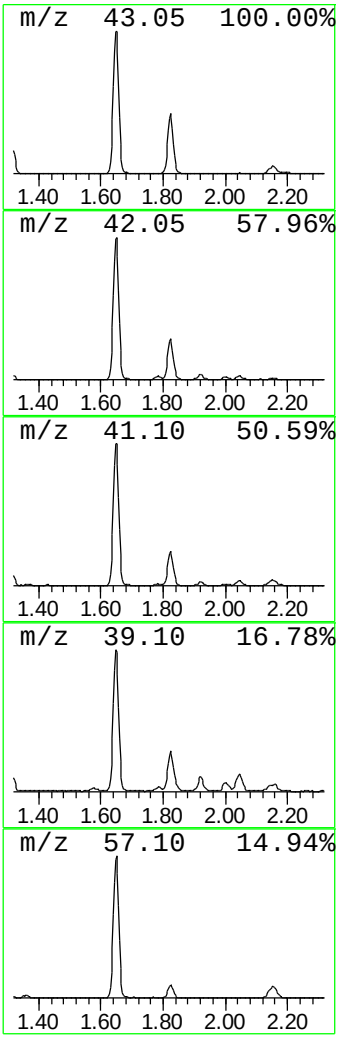
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.82	1.40 ug/L	228239	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Pentane	72	C5H12	000109-66-0	86
3	Pentane	72	C5H12	000109-66-0	86
4	Pentane	72	C5H12	000109-66-0	72
5	Oxirane, ethyl-	72	C4H8O	000106-88-7	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

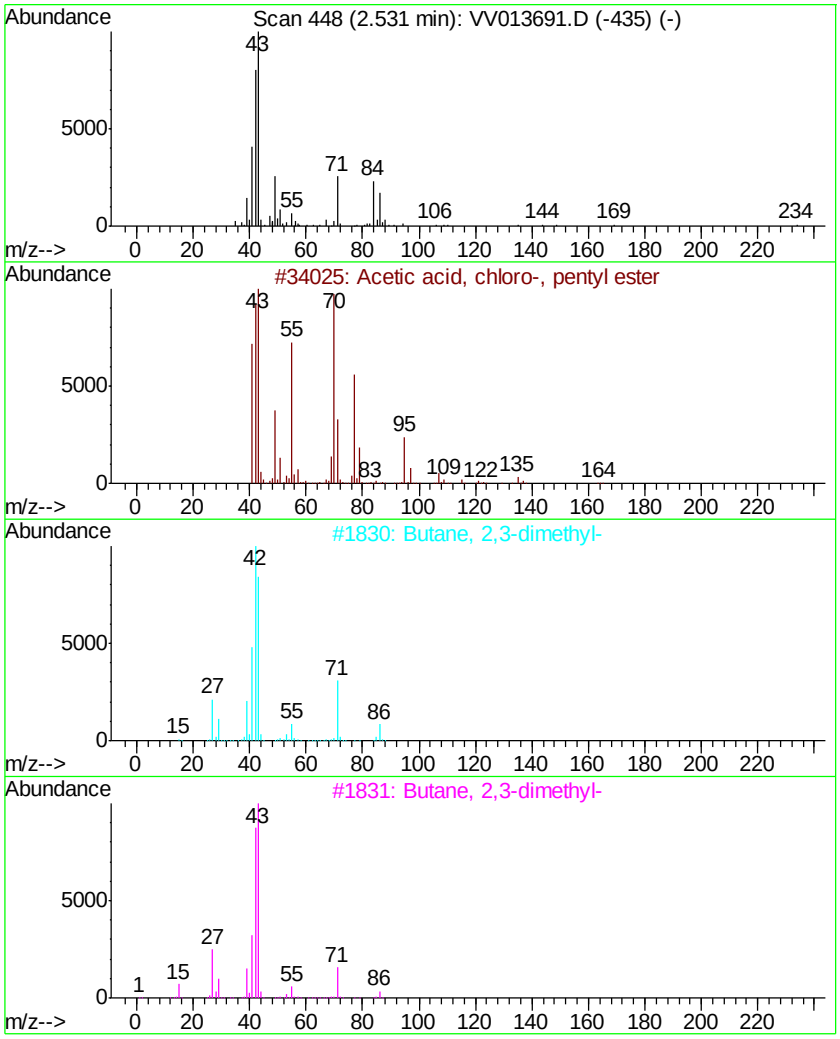
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

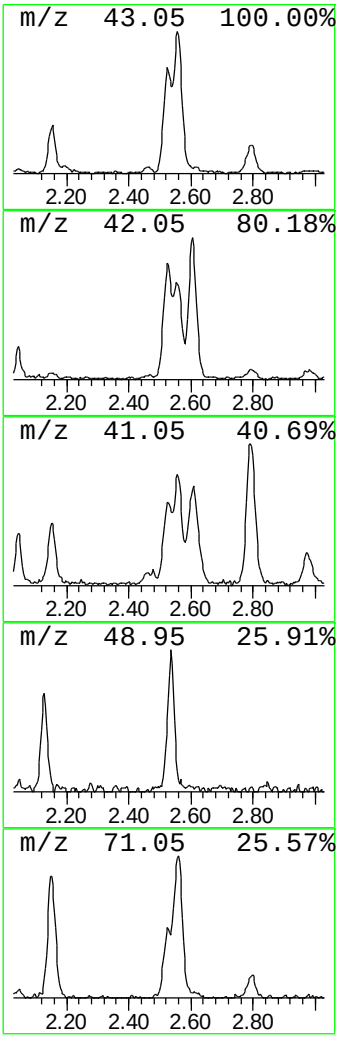
****

Peak Number 3 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.53	0.59 ug/L	96499	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, chloro-, pentyl ester	164	C7H13ClO2	005411-55-2	39
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	35
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	35
4	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	32
5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	32





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

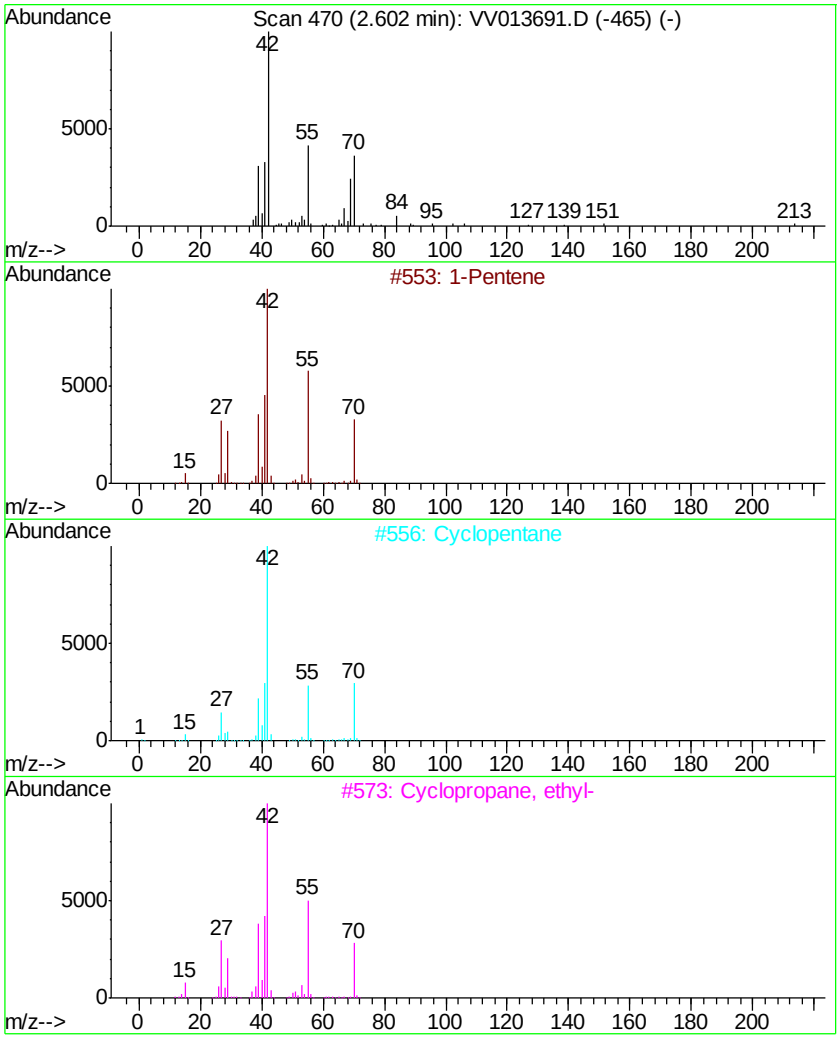
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

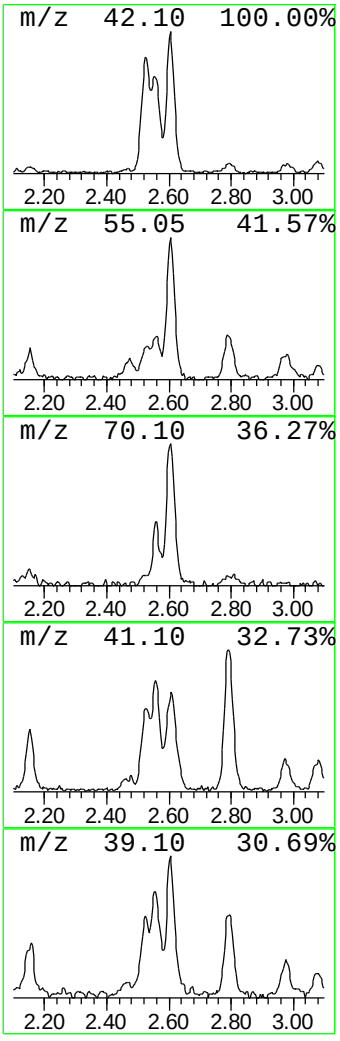
****

Peak Number 4 (DEL) Alkane: Cyclic2.60 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.60	0.67 ug/L	109627	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene	70	C5H10	000109-67-1	64
2	Cyclopentane	70	C5H10	000287-92-3	64
3	Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
4	1-Pentene	70	C5H10	000109-67-1	53
5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	50





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Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

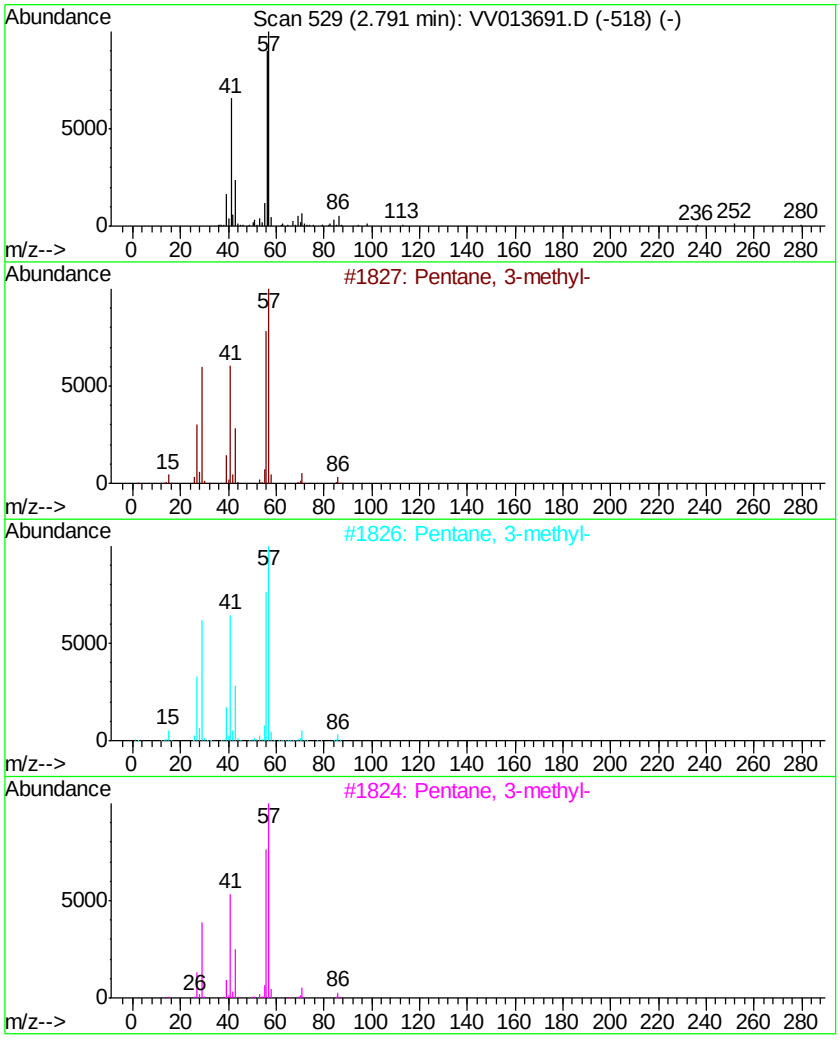
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

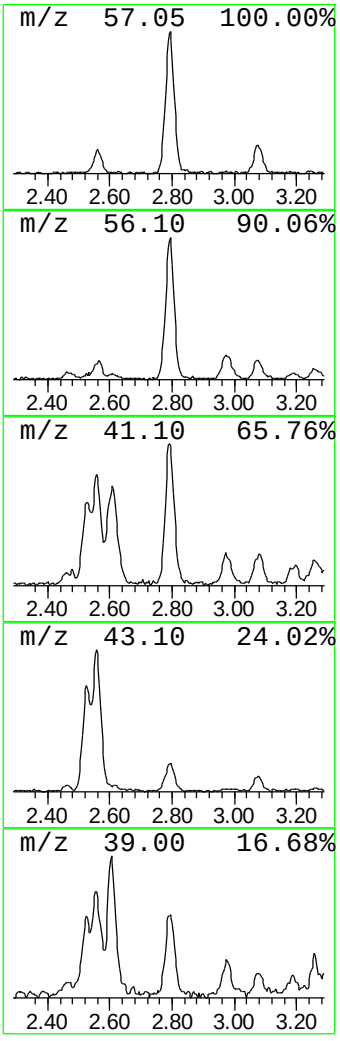
****

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	0.82 ug/L	133438	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
5	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	56





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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

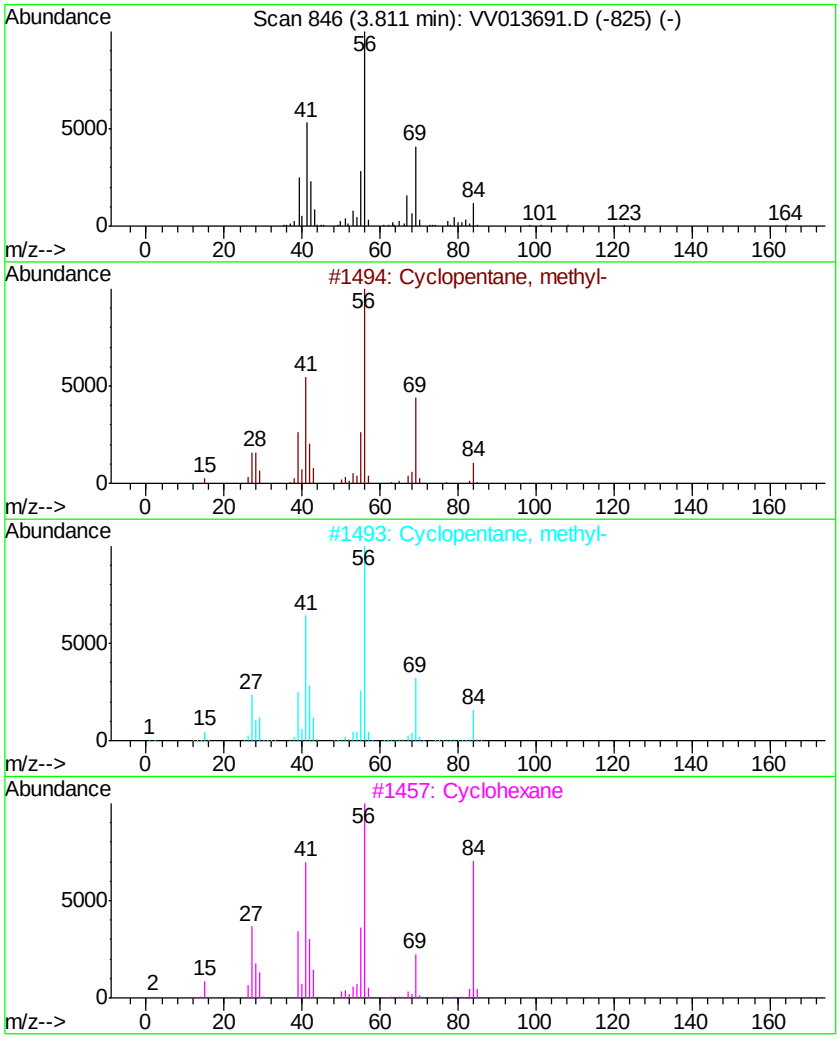
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

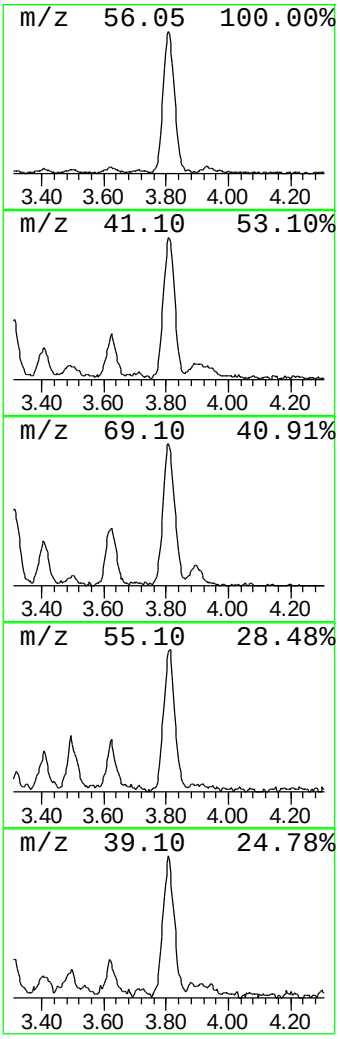
****

Peak Number 6 (DEL) Alkane: Cyclic3.81 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.81	1.56 ug/L	253522	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
3	Cyclohexane	84	C6H12	000110-82-7	72
4	Cyclohexane	84	C6H12	000110-82-7	64
5	Cyclohexane	84	C6H12	000110-82-7	64





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ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

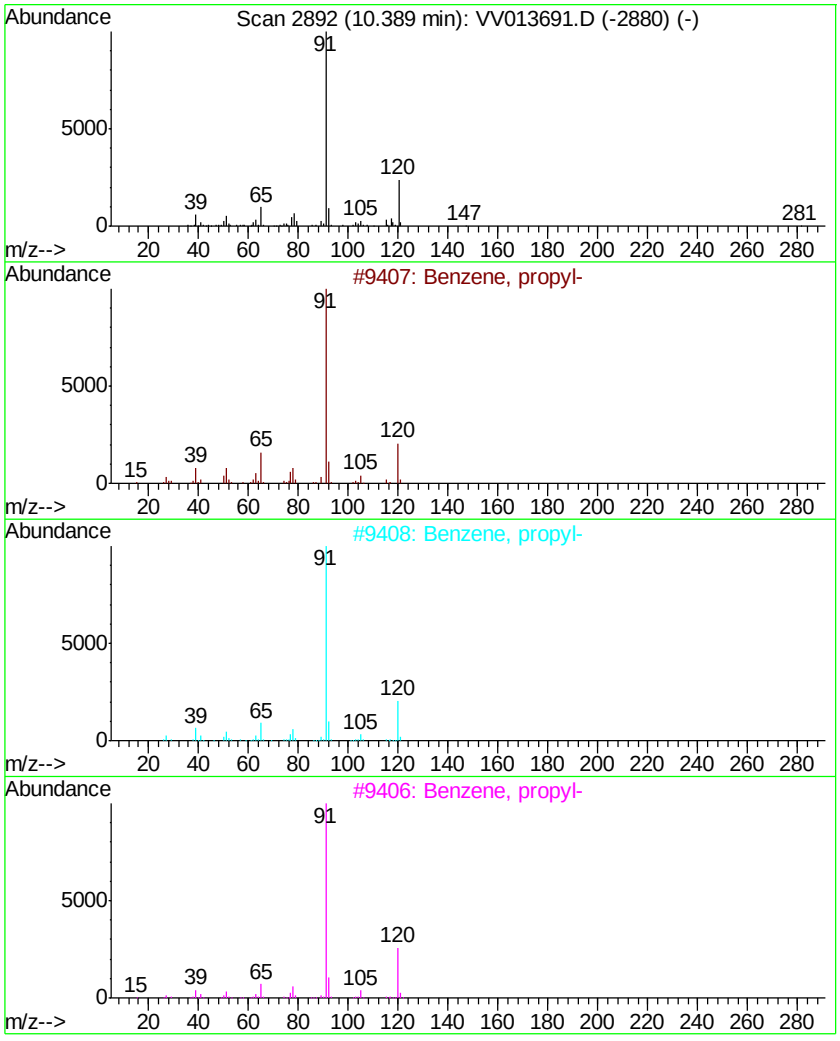
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

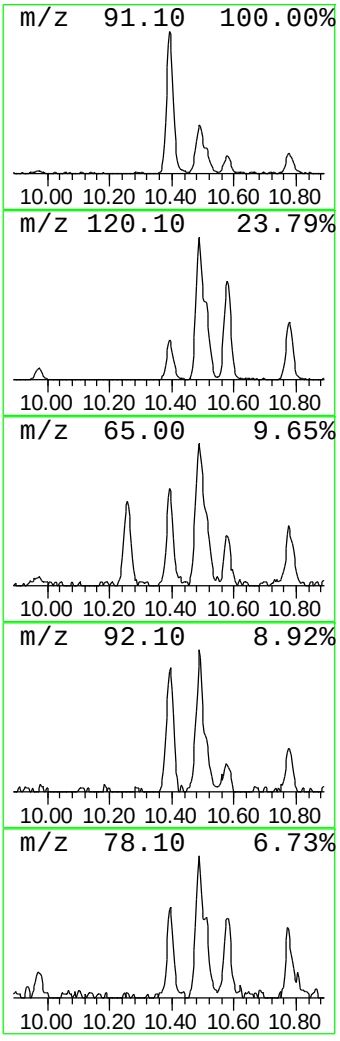
****

Peak Number 8 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	0.73 ug/L	139715	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	81
2	Benzene, propyl-	120	C9H12	000103-65-1	81
3	Benzene, propyl-	120	C9H12	000103-65-1	74
4	N,N-Dibenzylethylenediamine diac...	324	C20H24N2O2	000122-75-8	72
5	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
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Instrument : MSVOA_V
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

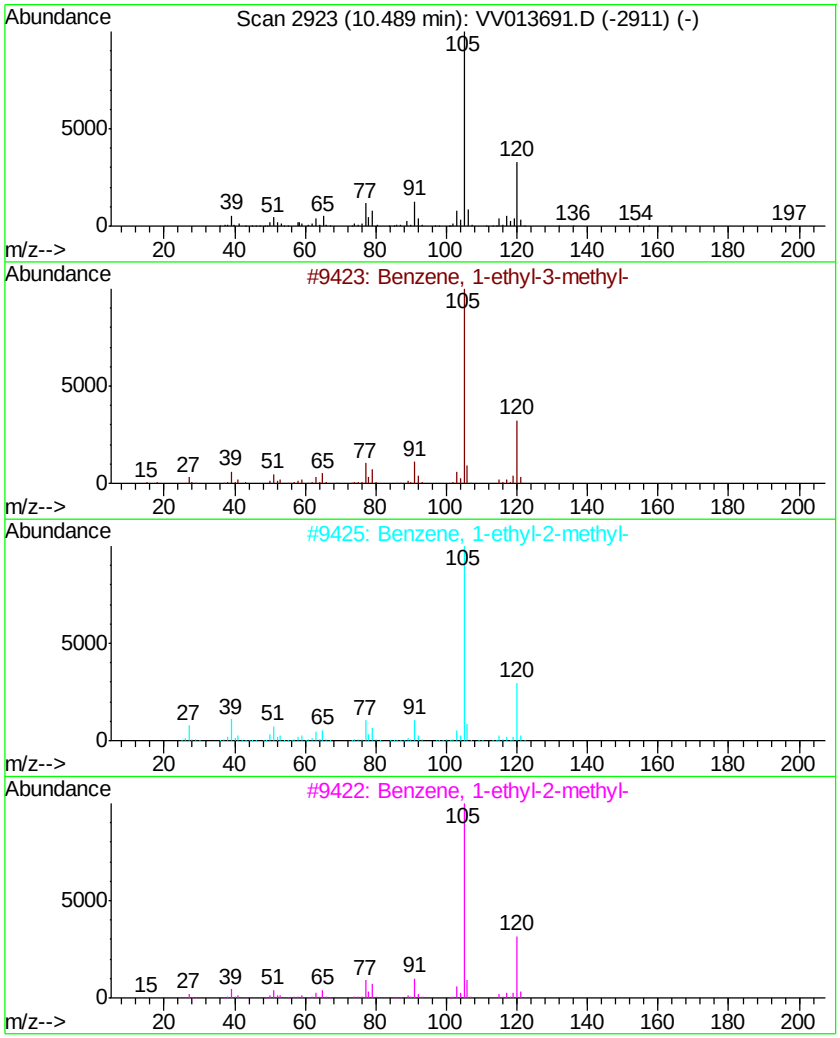
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

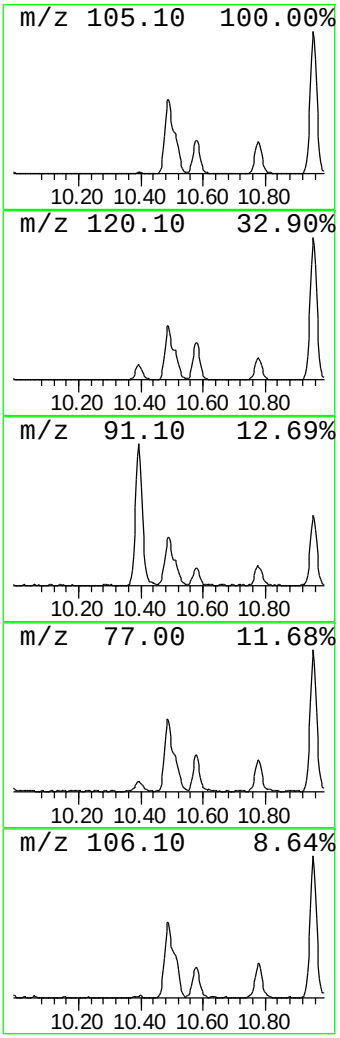
****

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.49	3.03 ug/L	580615	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

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

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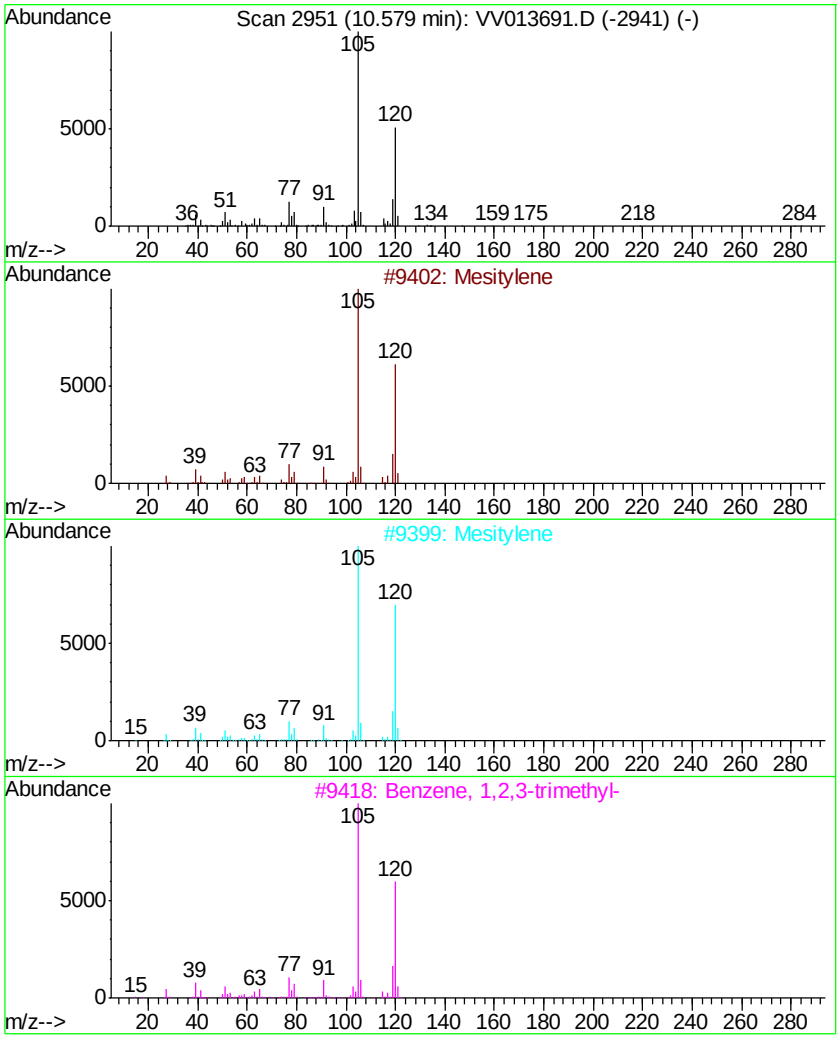
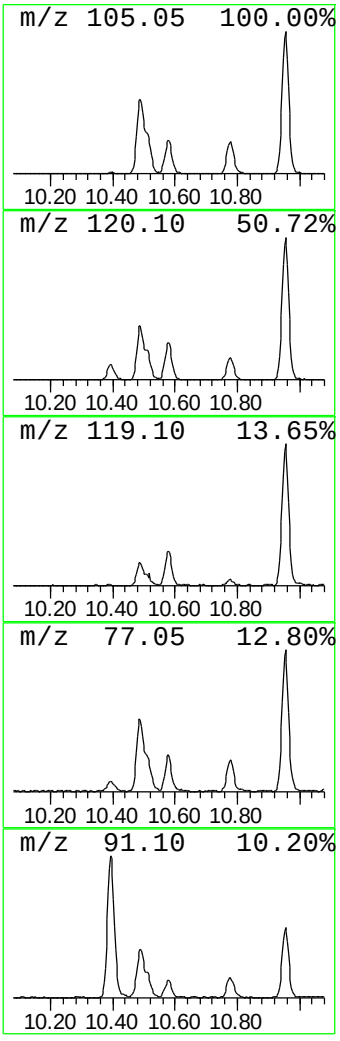
Peak Number 10 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	1.08 ug/L	207780	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Mesitylene	120	C9H12	000108-67-8	97
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
4	Mesitylene	120	C9H12	000108-67-8	95
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

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

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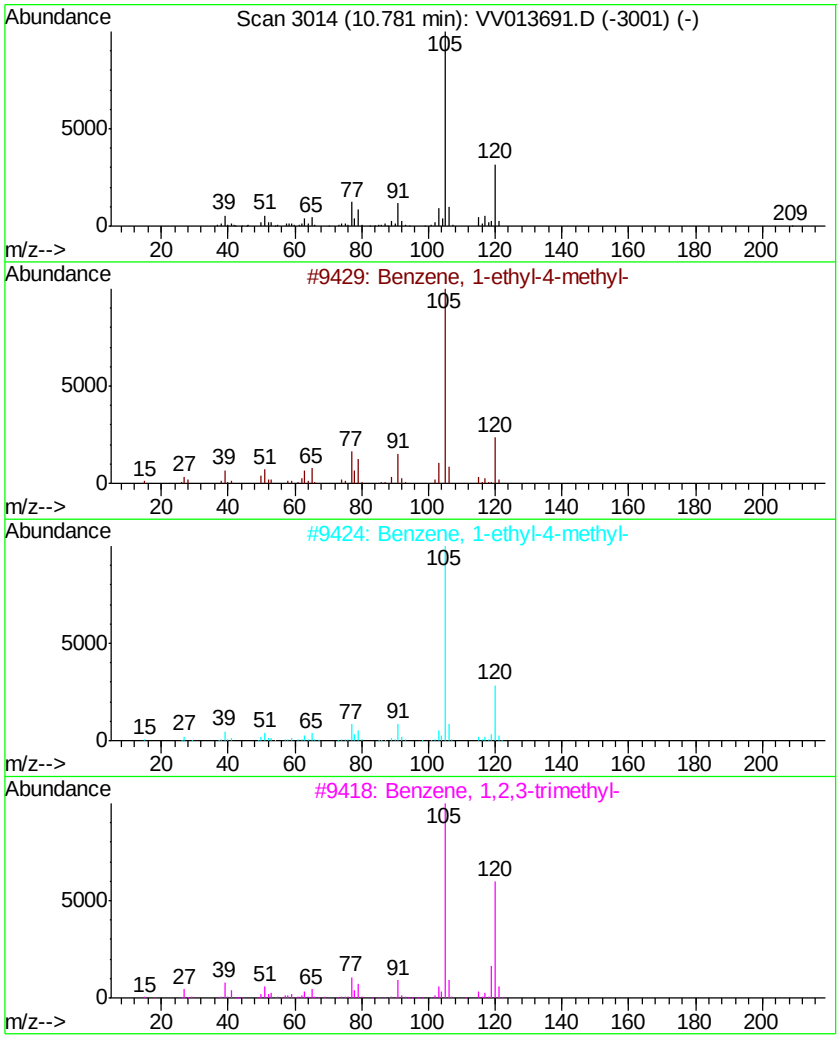
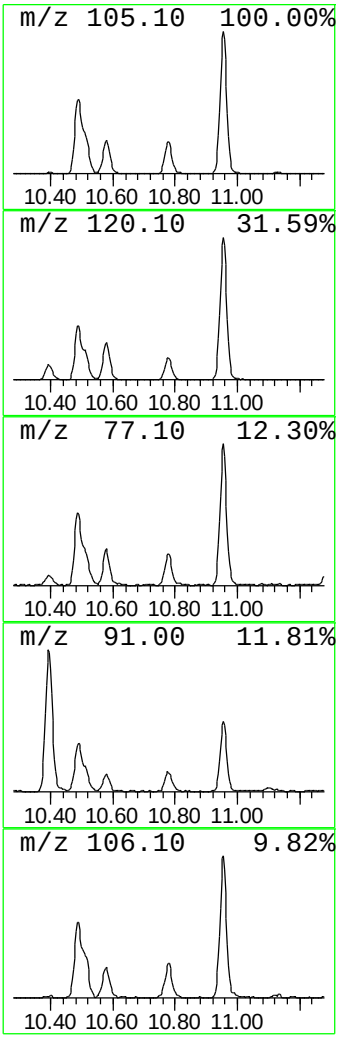
Peak Number 11 Benzene, 1-ethyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	0.97 ug/L	186695	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

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

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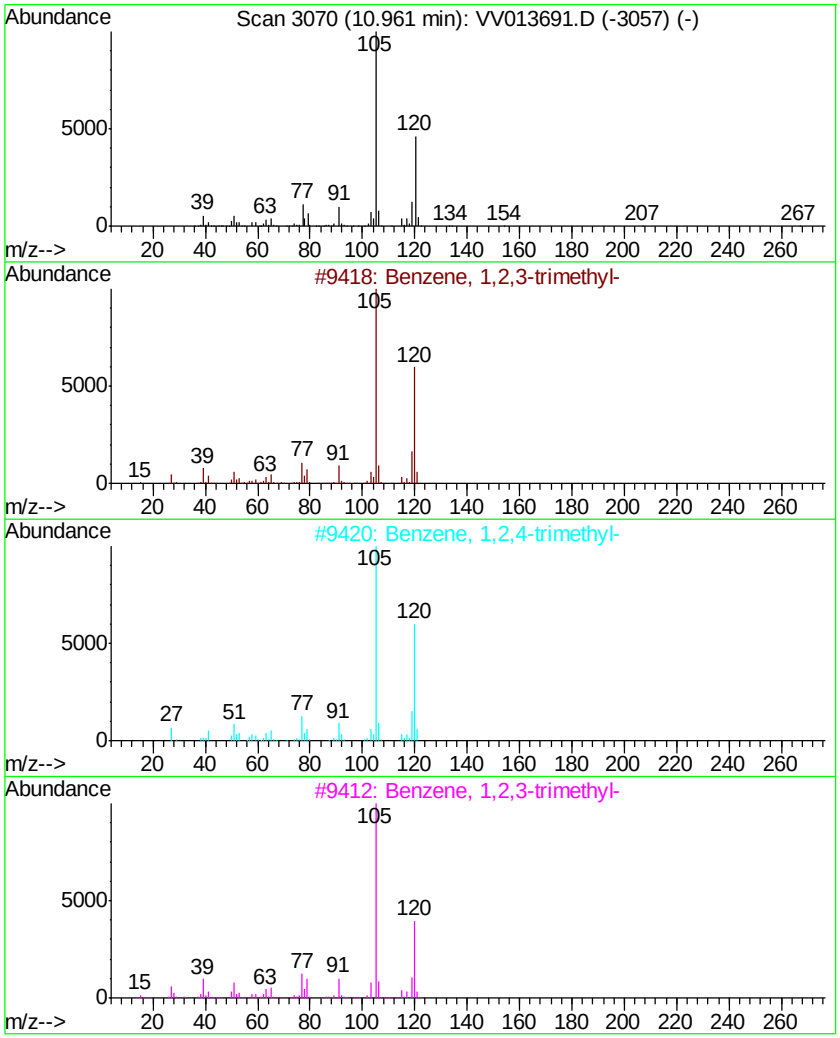
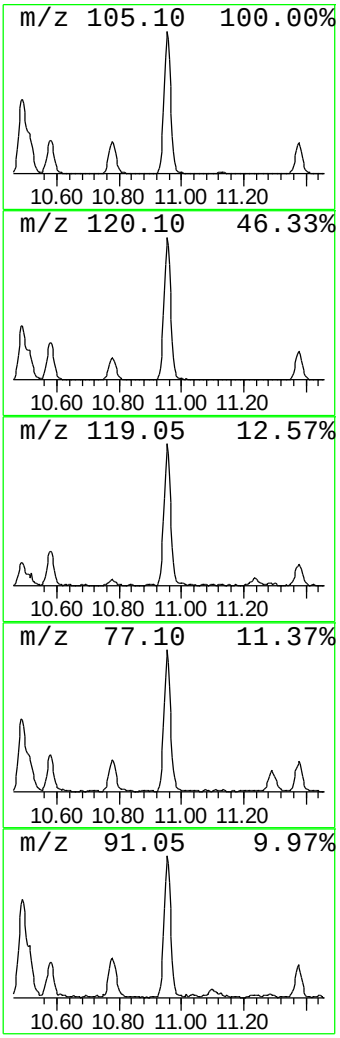
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	4.21 ug/L	807270	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Mesitylene	120	C9H12	000108-67-8	93





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.M
Quant Title : TRACE VOA SOM01.0

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

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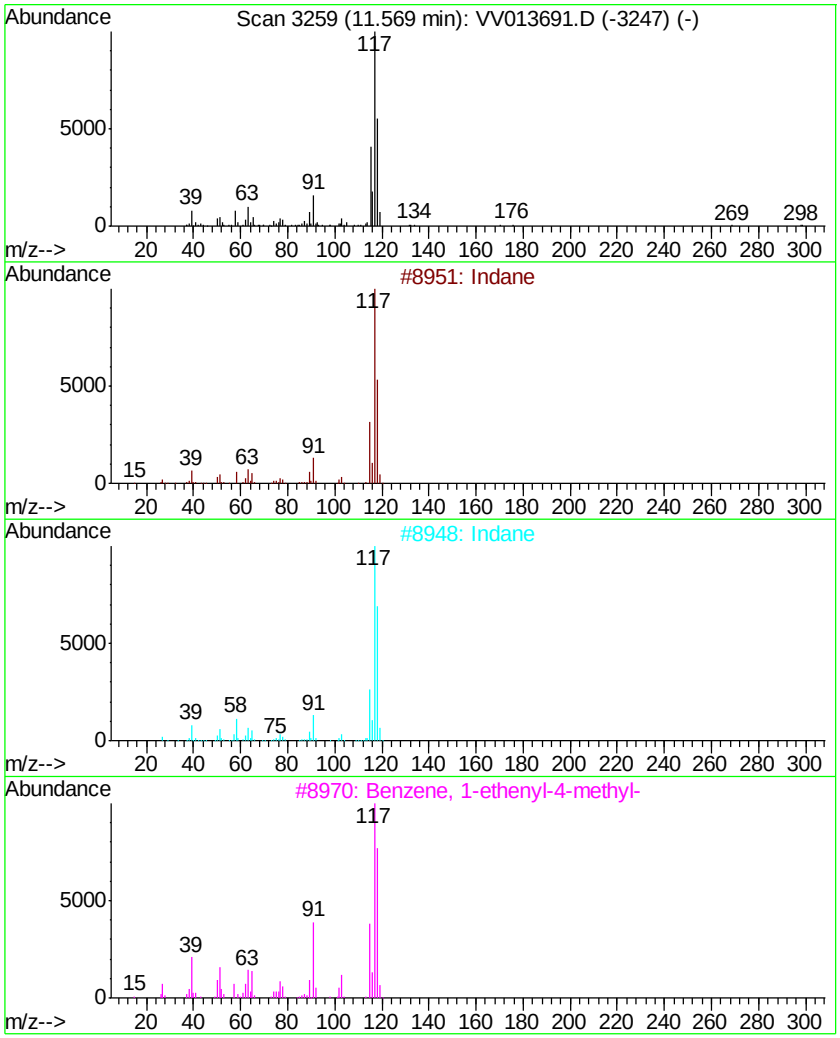
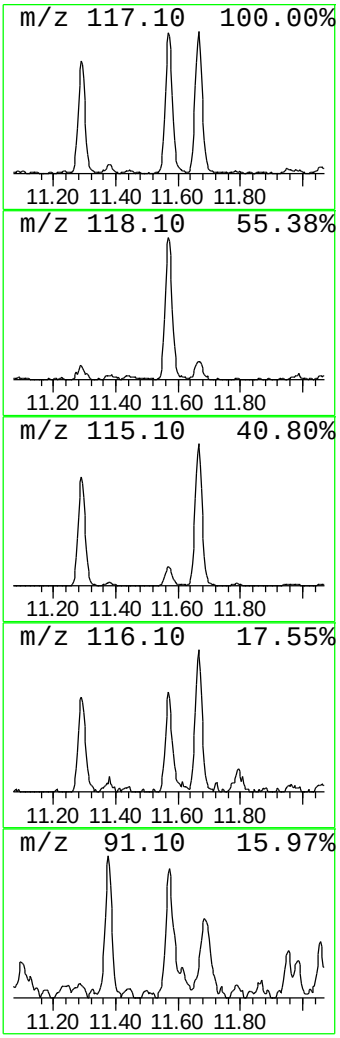
Peak Number 14 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	0.83 ug/L	158884	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Indane	118	C9H10	000496-11-7	76
3	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	76
4	Indane	118	C9H10	000496-11-7	76
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112019\
Data File : VV013691.D
Acq On : 20 Nov 2019 16:21
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR111519WMA.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: Str...	1.65	4.9	ug/L	797745	1	5.66	814118	5.0
(DEL) Alkane: Str...	1.82	1.4	ug/L	228239	1	5.66	814118	5.0
unknown-01	2.53	0.6	ug/L	96499	1	5.66	814118	5.0
(DEL) Alkane: Cyc...	2.60	0.7	ug/L	109627	1	5.66	814118	5.0
(DEL) Alkane: Str...	2.79	0.8	ug/L	133438	1	5.66	814118	5.0
(DEL) Alkane: Cyc...	3.81	1.6	ug/L	253522	1	5.66	814118	5.0
Benzene, propyl-	10.39	0.7	ug/L	139715	3	11.29	957634	5.0
Benzene, 1-ethyl-...	10.49	3.0	ug/L	580615	3	11.29	957634	5.0
Mesitylene	10.58	1.1	ug/L	207780	3	11.29	957634	5.0
Benzene, 1-ethyl-...	10.78	1.0	ug/L	186695	3	11.29	957634	5.0
Benzene, 1,2,3-tr...	10.96	4.2	ug/L	807270	3	11.29	957634	5.0
Indane	11.57	0.8	ug/L	158884	3	11.29	957634	5.0