LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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_X\METHOD\82X042220W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.064	24	29	56	rBV2	18434046	56644019	100.00%	15.058%
2	2.826	305	318	321	rBV2	383615	863492	1.52%	0.230%
3	2.875	321	326	331	rVB	669904	1240525	2.19%	0.330%
4	3.155	360	372	385	rBV3	1402571	3597533	6.35%	0.956%
5	3.503	421	429	440	rBV	556448	1251849	2.21%	0.333%
6	3.795	468	477	486	rBV	621875	1496715	2.64%	0.398%
7	3.905	486	495	502	rBV3	354220	935294	1.65%	0.249%
8	4.173	525	539	547	rVB3	380033	1129266	1.99%	0.300%
9	4.289	548	558	564	rVV	585333	1747409	3.08%	0.465%
10	4.380	564	573	585	rVV	3365997	9703833	17.13%	2.580%
11	5.277	694	720	736	rVB	1111731	4166109	7.35%	1.108%
12	5.472	739	752	757	rBV2	603235	1915444	3.38%	0.509%
13	5.557	757	766	776	rBV3	2465703	9469631	16.72%	2.517%
14	5.703	783	790	810	rVB	2316882	7990176	14.11%	2.124%
15	5.905	812	823	837	rVV2	1669487	4991643	8.81%	1.327%
16	6.039	837	845	851	rVV	552715	1395357	2.46%	0.371%
17	6.118	851	858	865	rVV	455977	1211433	2.14%	0.322%
18	6.240	866	878	881	rVV	1160930	3454900	6.10%	0.918%
19	6.325	884	892	902	rVV2	4118559	13711961	24.21%	3.645%
20	6.435	902	910	922	rVB	1072938	3106374	5.48%	0.826%
21	6.679	937	950	960	rBV2	779691	1952858	3.45%	0.519%
22	6.837	964	976	980	rBV	1499108	3805708	6.72%	1.012%
23	7.045	1001	1010	1017	rBV3	263135	619214	1.09%	0.165%
24	7.130	1017	1024	1033	rVB	360581	776041	1.37%	0.206%
25	7.234	1033	1041	1049	rBV	1143606	2571567	4.54%	0.684%
26	7.441	1061	1075	1087	rBV5	4121515	10841072	19.14%	2.882%
27	7.557	1087	1094	1099	rBV2	305168	586744	1.04%	0.156%
28	7.618	1099	1104	1109	rBV	415987	777855	1.37%	0.207%
29	7.715	1114	1120	1132	rVB	865741	1723957	3.04%	0.458%
30	7.910	1146	1152	1153	rBV	573788	867579	1.53%	0.231%
31	7.935	1154	1156	1163	rVV	696006	1215766	2.15%	0.323%
32	8.038	1163	1173	1184	rVB	1278045	3026113	5.34%	0.804%
33	8.160	1184	1193	1195	rBV	1305617	2454540	4.33%	0.653%
34	8.197	1195	1199	1203	rVV	2196283	3909215	6.90%	1.039%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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_X\METHOD\82X042220W.M
Title : SW846 8260

35	8.246	1203	1207	1215	rVV3	727939	1469364	2.59%	0.391%
36	8.361	1222	1226	1233	rVB4	340189	643363	1.14%	0.171%
37	8.483	1241	1246	1250	rBV2	359073	569696	1.01%	0.151%
38	8.557	1250	1258	1268	rVB	2046567	3737380	6.60%	0.994%
39	8.654	1268	1274	1276	rBV3	477893	845442	1.49%	0.225%
40	8.697	1276	1281	1288	rVB	3035514	4894161	8.64%	1.301%
41	8.886	1305	1312	1319	rVB3	313218	789685	1.39%	0.210%
42	8.971	1320	1326	1331	rVV2	416791	692476	1.22%	0.184%
43	9.209	1360	1365	1375	rVB2	950027	1665844	2.94%	0.443%
44	10.099	1507	1511	1515	rVB	3029842	3954249	6.98%	1.051%
45	10.239	1529	1534	1538	rBV	2050112	2703133	4.77%	0.719%
46	10.343	1545	1551	1557	rVB	1492424	2082713	3.68%	0.554%
47	11.001	1652	1659	1670	rBV	12262308	16052830	28.34%	4.268%
48	11.123	1674	1679	1684	rBV	2836765	3475015	6.13%	0.924%
49	11.343	1710	1715	1722	rVB	14878755	18597398	32.83%	4.944%
50	11.440	1726	1731	1735	rBV	1166081	1407490	2.48%	0.374%
51	11.489	1735	1739	1745	rVB	848526	1204939	2.13%	0.320%
52	11.654	1760	1766	1774	rBV	2672136	3295835	5.82%	0.876%
53	11.794	1784	1789	1793	rVB	1619153	1881726	3.32%	0.500%
54	11.910	1802	1808	1809	rBV2	626097	825415	1.46%	0.219%
55	11.928	1809	1811	1818	rVB2	727830	896584	1.58%	0.238%
56	12.062	1828	1833	1839	rBV	3768132	4630226	8.17%	1.231%
57	12.129	1839	1844	1848	rBV	1385567	1662128	2.93%	0.442%
58	12.288	1864	1870	1876	rBV	32260883	43437169	76.68%	11.547%
59	12.355	1876	1881	1889	rVB2	1432168	2668790	4.71%	0.709%
60	12.446	1891	1896	1902	rBV2	569208	681578	1.20%	0.181%
61	12.653	1925	1930	1933	rBV	7303233	8564775	15.12%	2.277%
62	12.684	1933	1935	1940	rVV	2289123	2596454	4.58%	0.690%
63	12.739	1940	1944	1951	rVB	6962046	9090553	16.05%	2.417%
64	12.964	1973	1981	1985	rBV	4062739	4939654	8.72%	1.313%
65	13.001	1985	1987	1992	rVV	510868	586985	1.04%	0.156%
66	13.208	2017	2021	2031	rVB	5223836	6419646	11.33%	1.707%
67	13.336	2036	2042	2047	rVB2	9397097	12832108	22.65%	3.411%
68	13.464	2058	2063	2071	rVB	1604928	1981073	3.50%	0.527%
69	13.586	2079	2083	2086	rVV	693756	955326	1.69%	0.254%
70	13.623	2086	2089	2094	rVV2	680312	1058853	1.87%	0.281%
71	13.684	2094	2099	2100	rVV2	432395	585190	1.03%	0.156%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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_X\METHOD\82X042220W.M
Title : SW846 8260

72	13.702	2100	2102	2109	rVB2	760288	1027716	1.81%	0.273%
73	13.818	2114	2121	2129	rBV	21825316	27671508	48.85%	7.356%
74	14.257	2187	2193	2203	rBV	350491	613939	1.08%	0.163%
75	14.677	2258	2262	2272	rVB	3785892	4684315	8.27%	1.245%
76	14.824	2281	2286	2294	rBV	2159816	2638165	4.66%	0.701%

Sum of corrected areas: 376162081

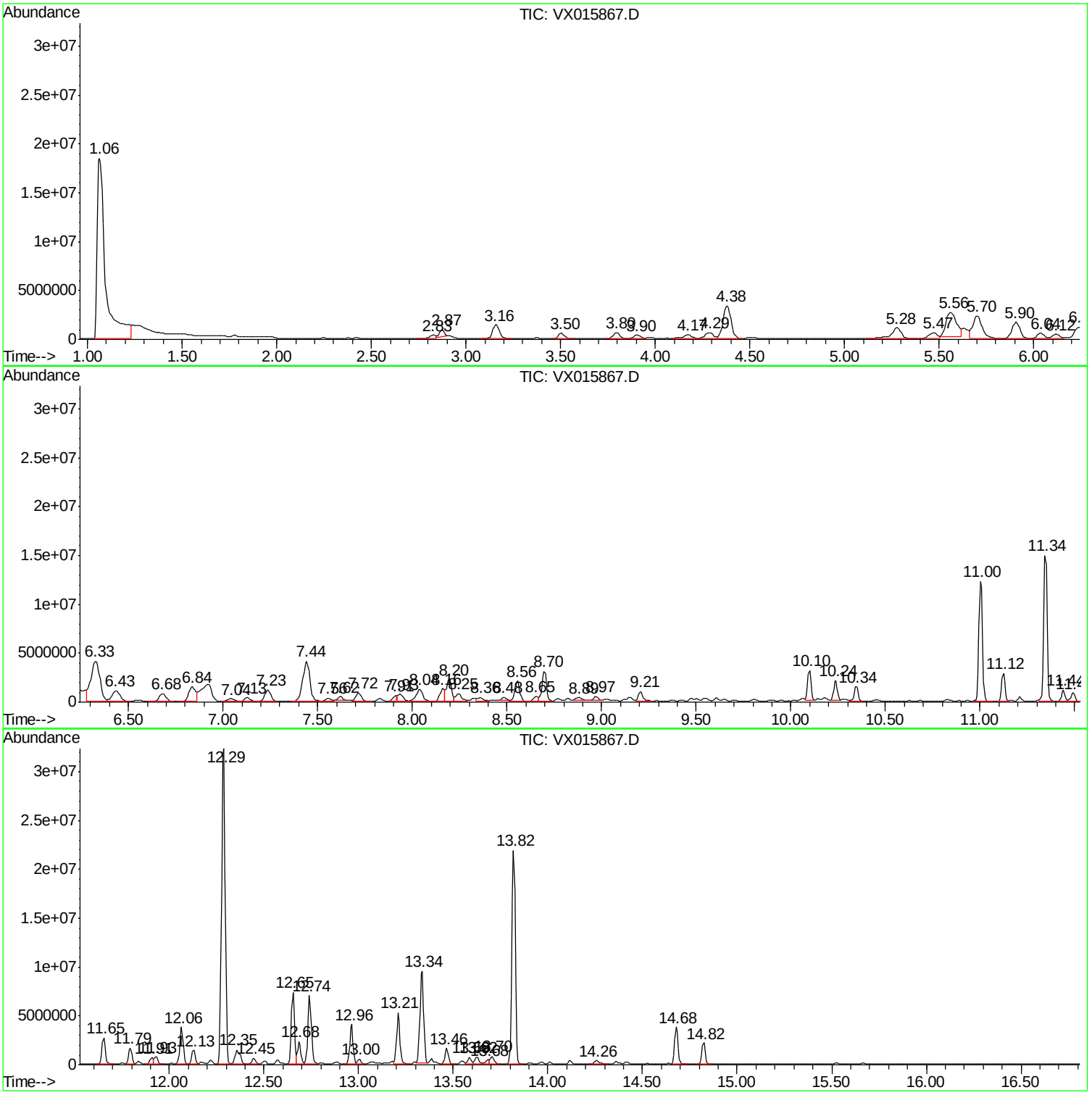
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

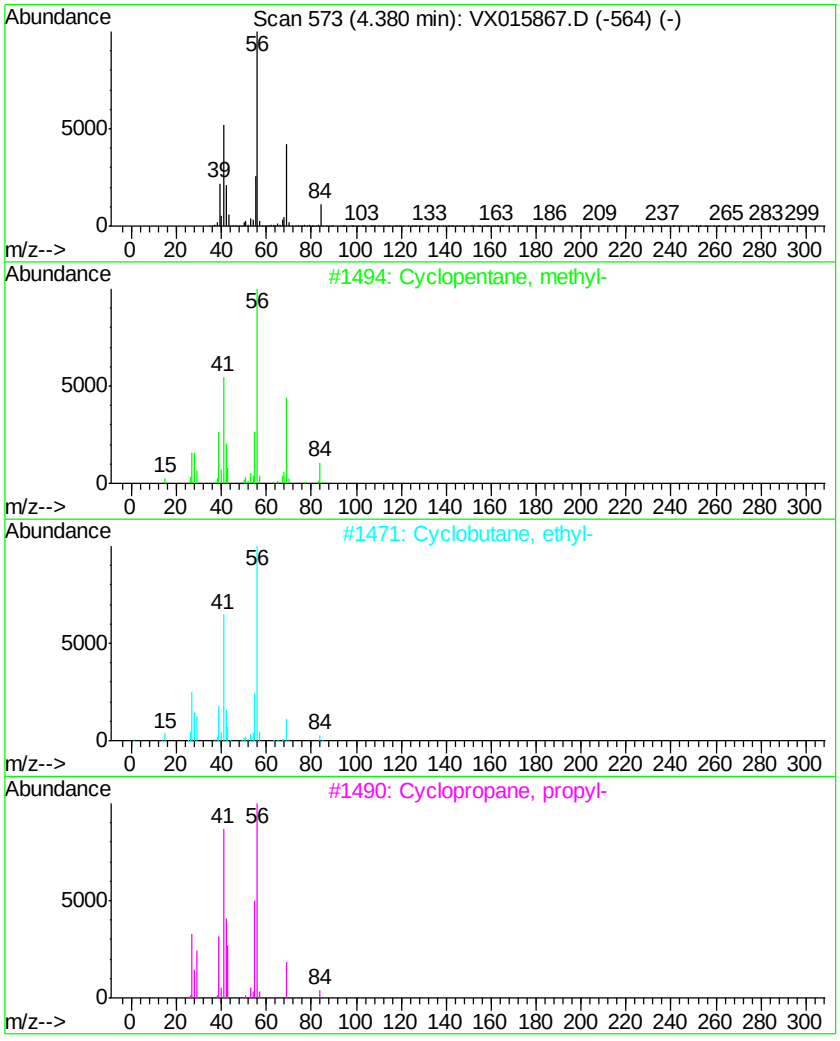
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

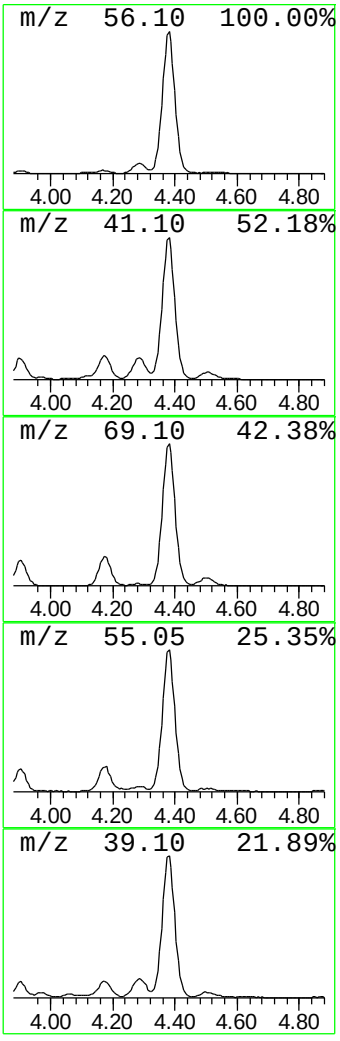
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	51.24 ug/l	9703830	Pentafluorobenzene	5.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	72
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
5	1-Hexene	84	C6H12	000592-41-6	43





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

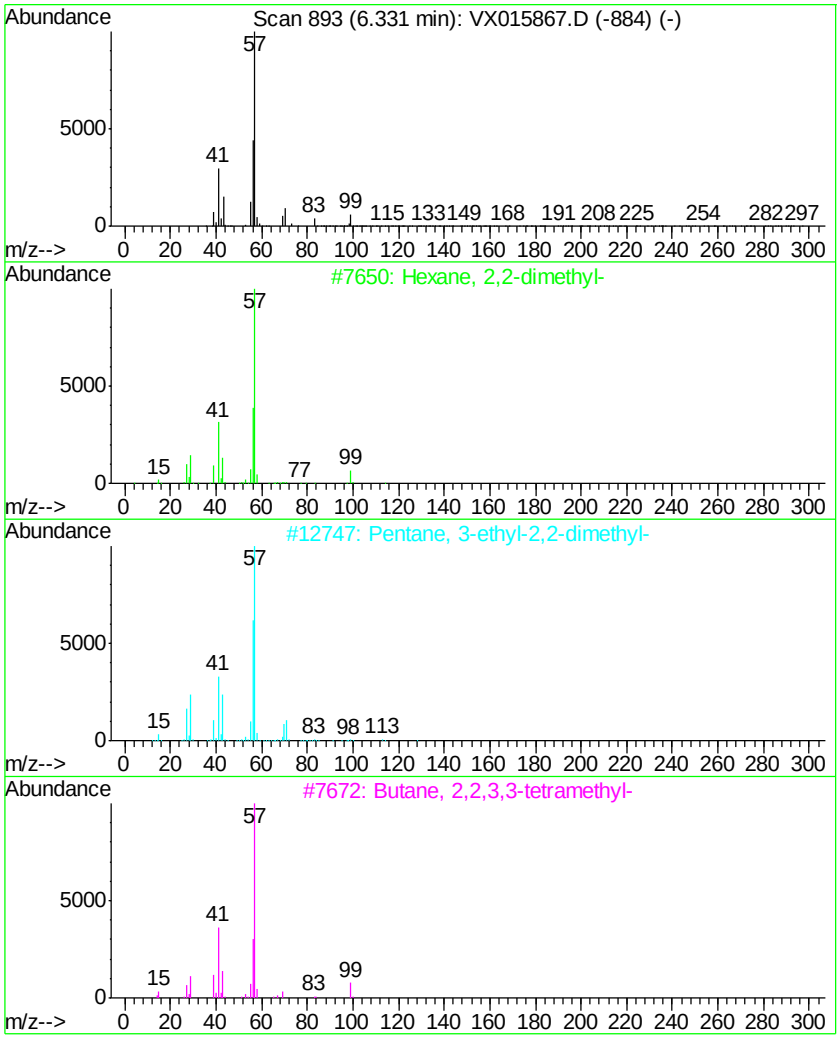
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

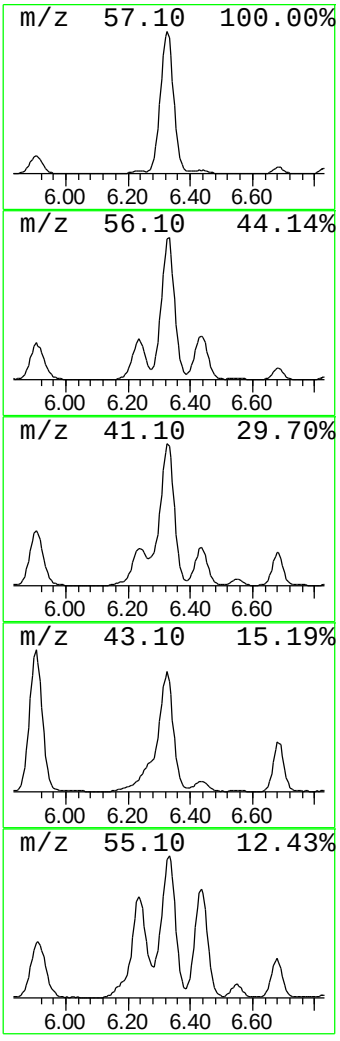
****

Peak Number 3 Hexane, 2,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	180.15 ug/l	13712000	1,4-Difluorobenzene	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
2	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
3	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	56
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	50
5	1-Butanol, 3,3-dimethyl-	102	C6H14O	000624-95-3	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

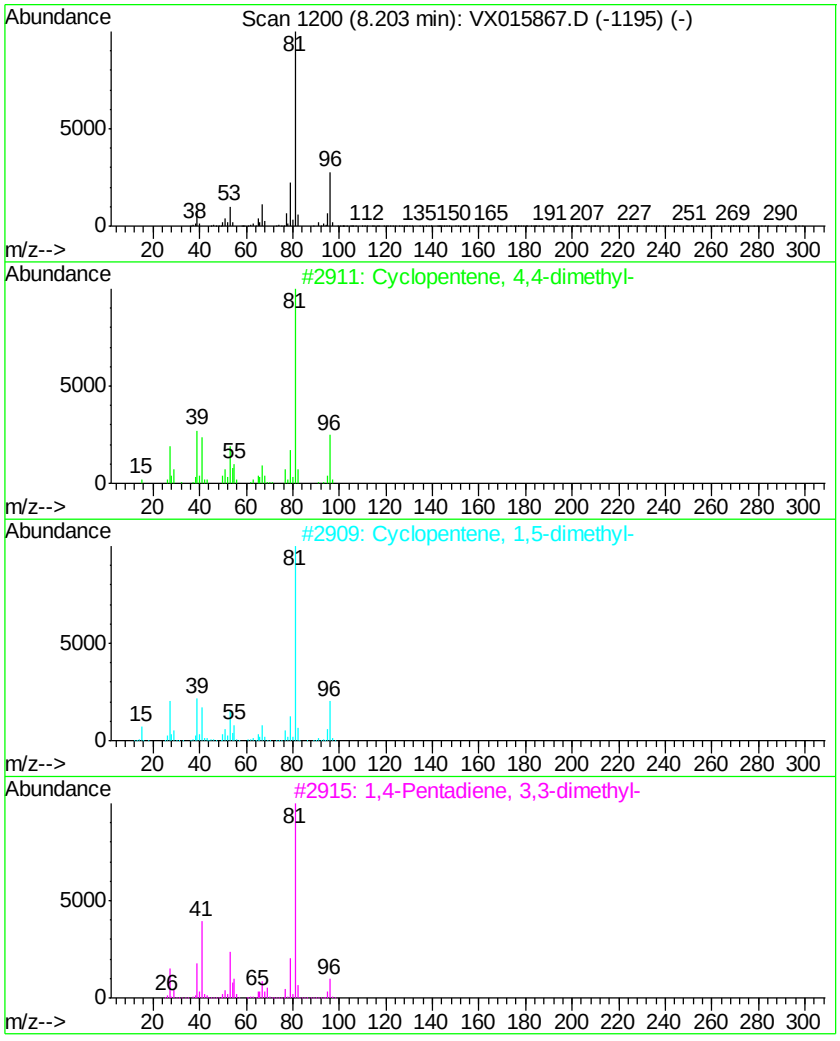
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

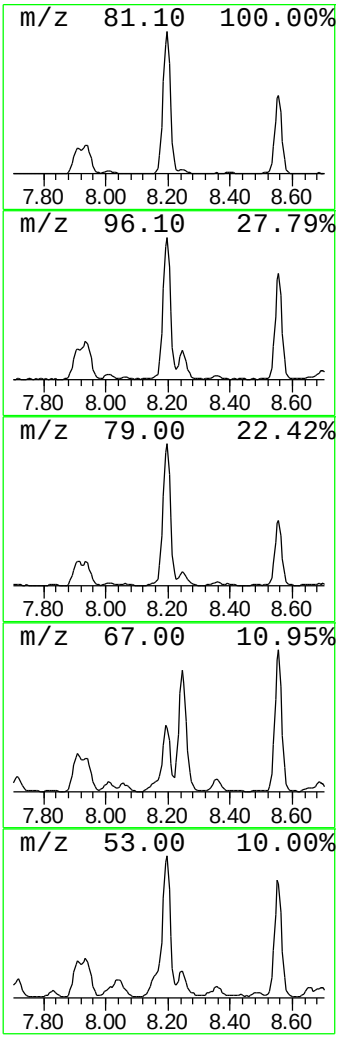
****

Peak Number 4 Cyclopentene, 4,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	51.36 ug/l	3909220	1,4-Difluorobenzene	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	87
2	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	80
3	1,4-Pentadiene, 3,3-dimethyl-	96	C7H12	001112-35-2	78
4	3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	78
5	Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

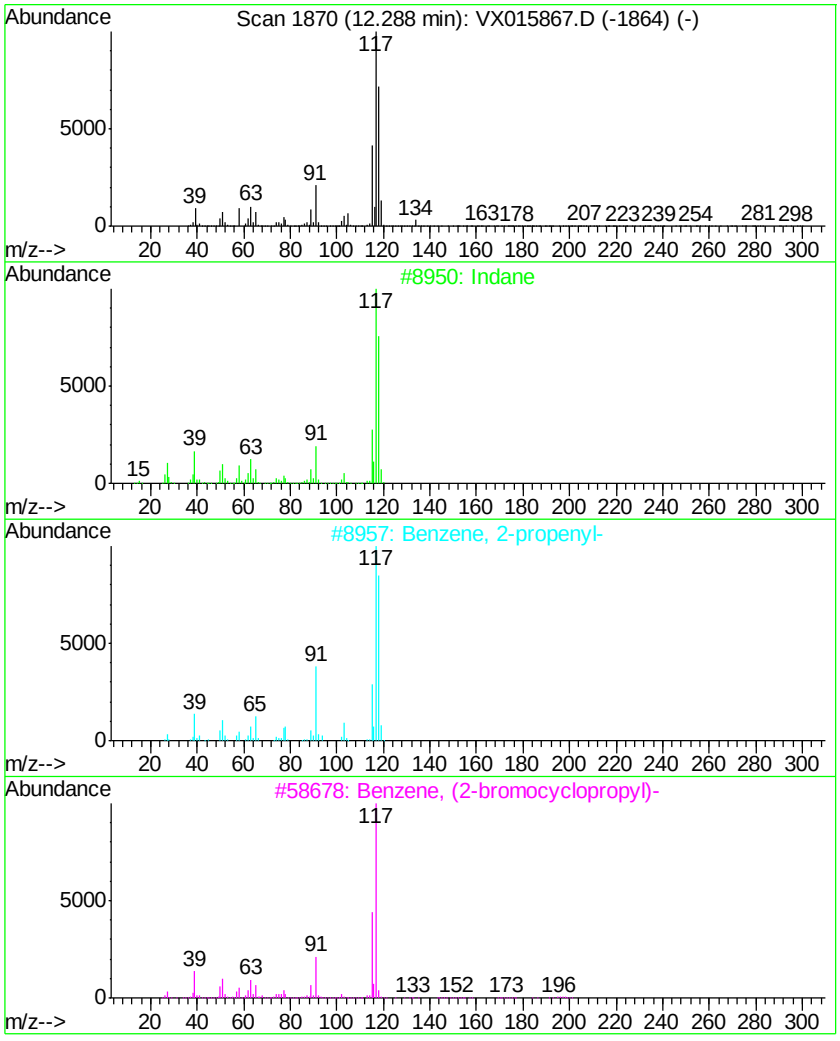
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

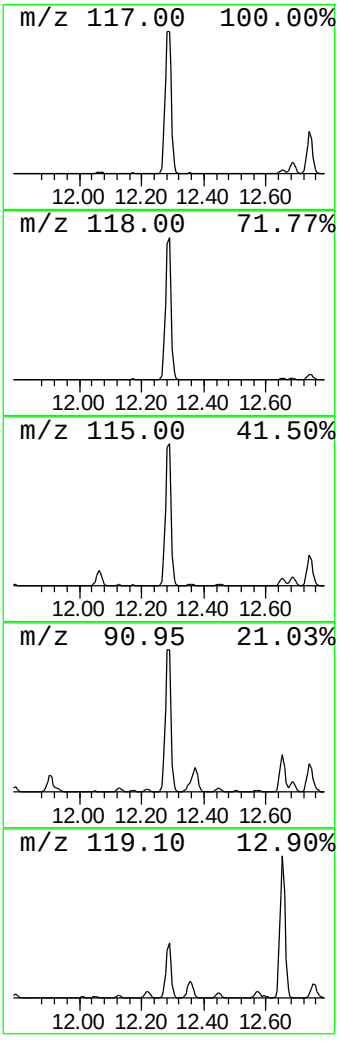
****

Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	469.06 ug/l	43437200	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	64
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	62





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

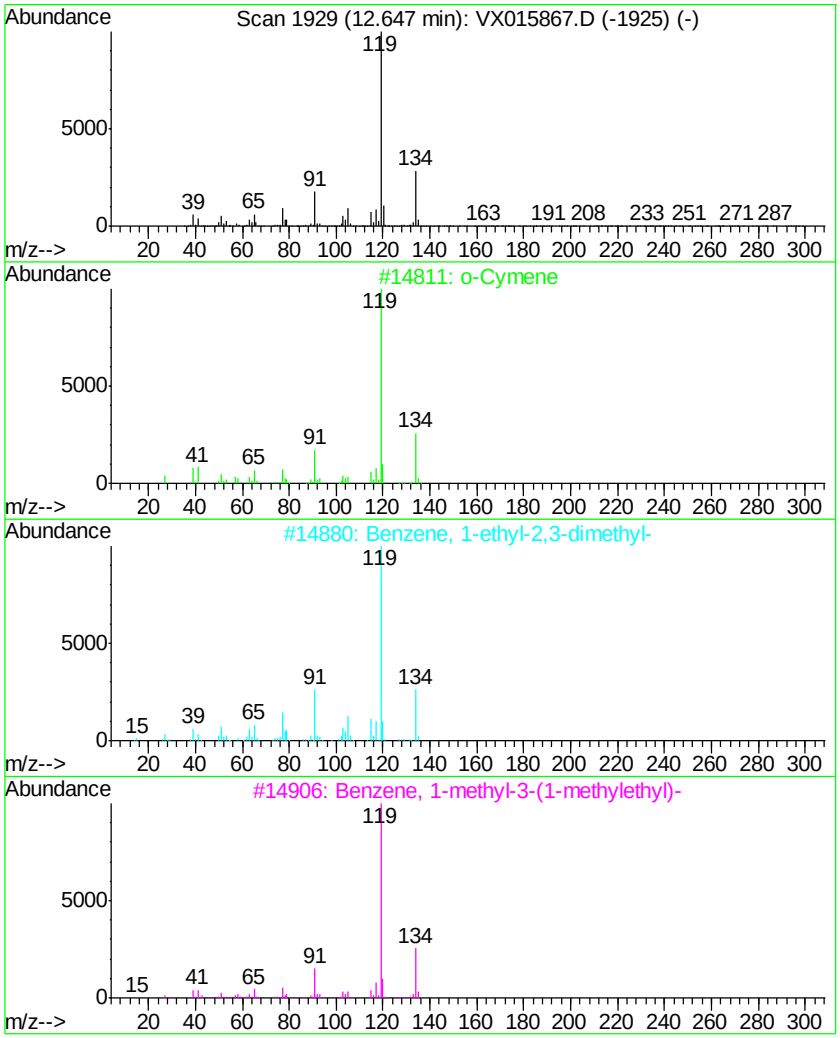
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

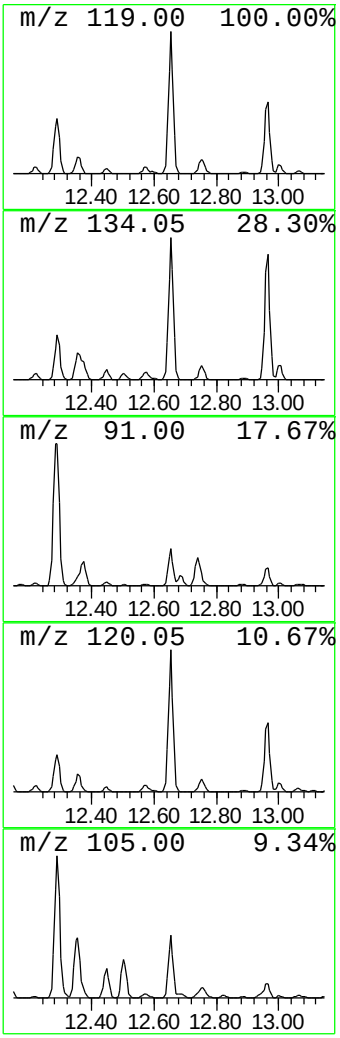
****

Peak Number 6 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	92.49 ug/l	8564780	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

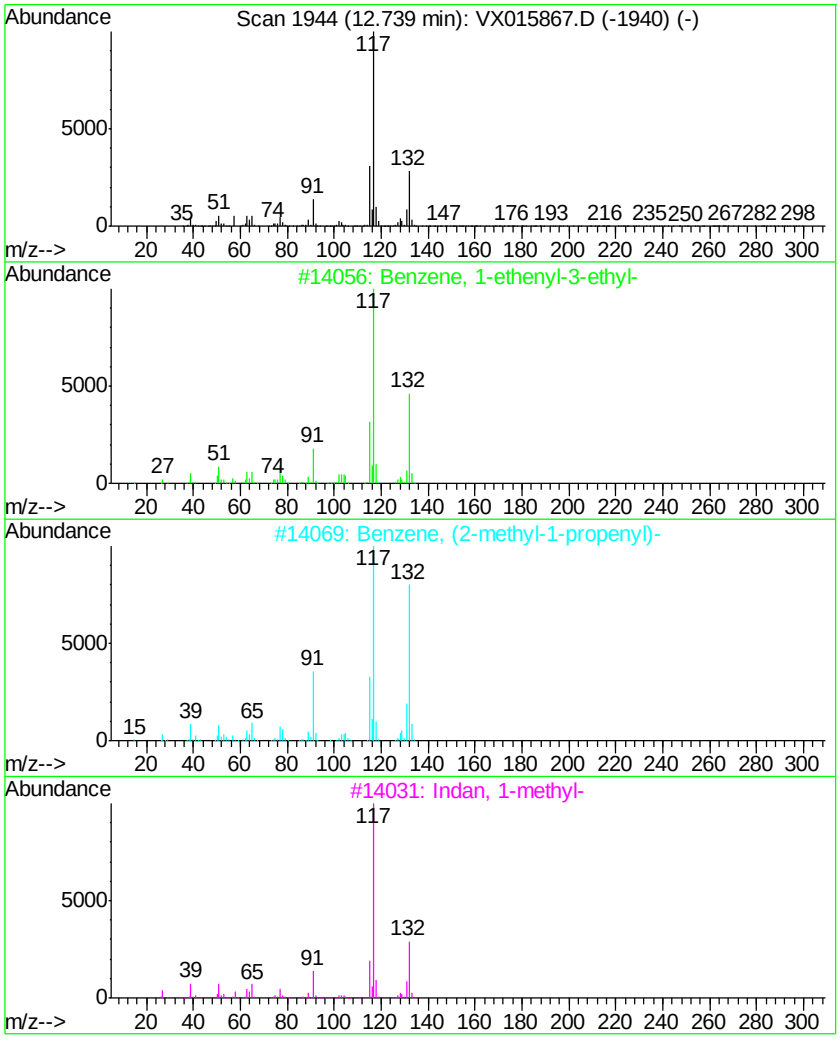
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

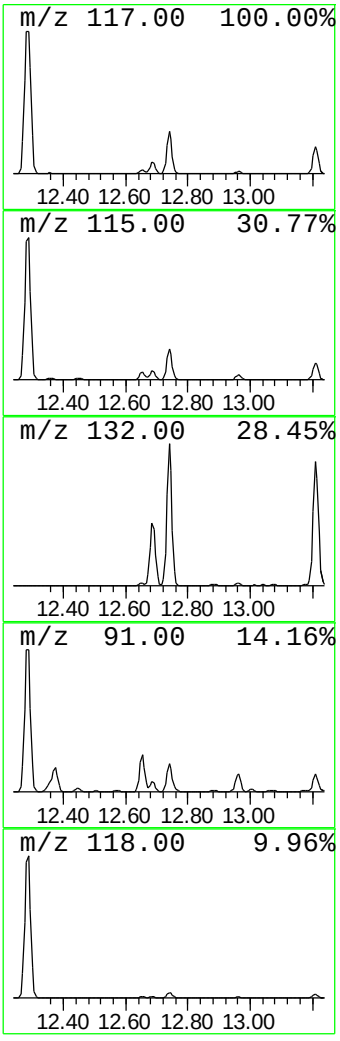
****

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	98.17 ug/l	9090550	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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

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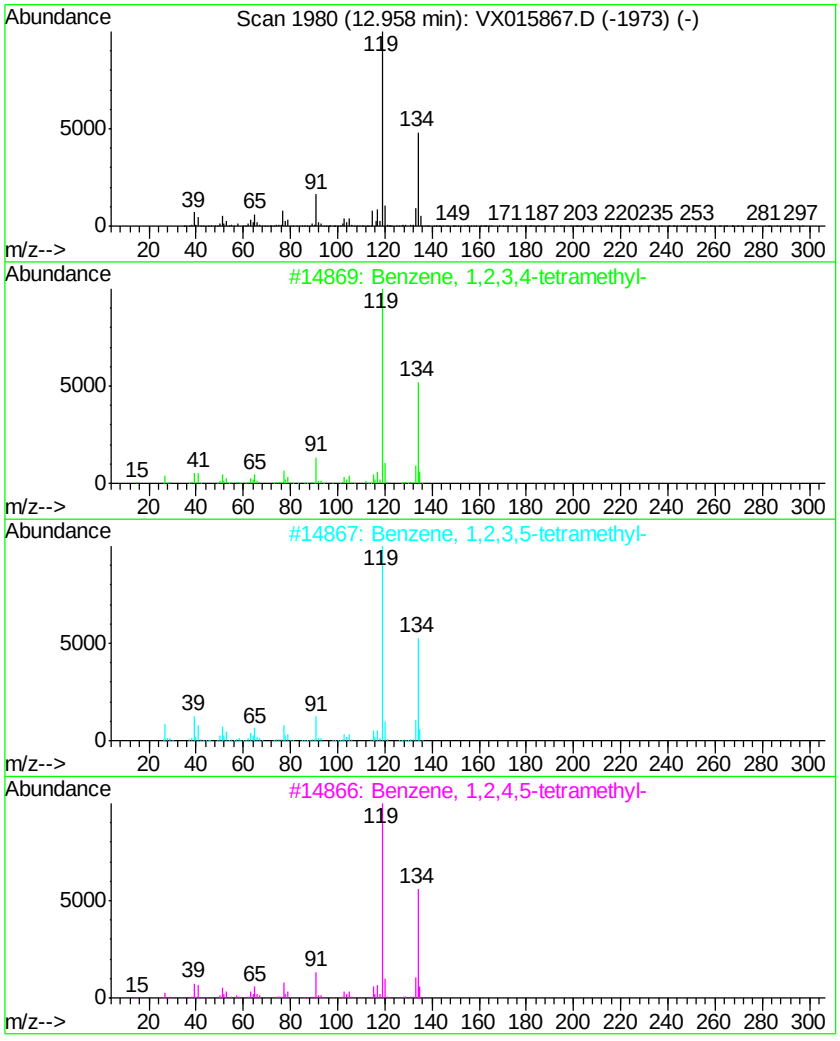
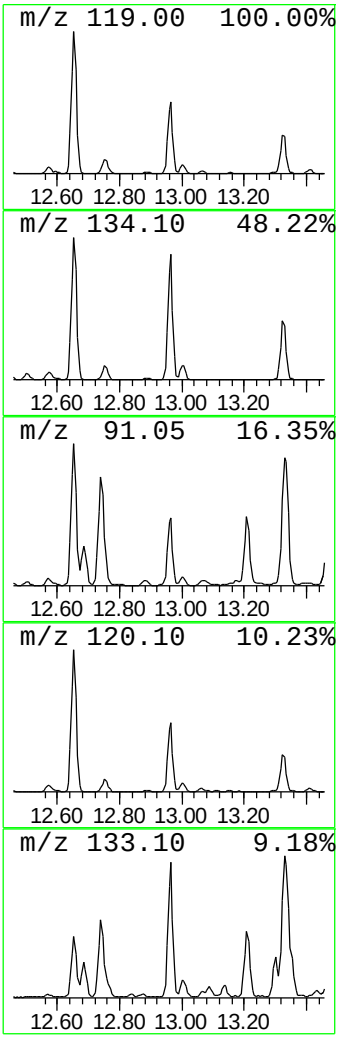
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	53.34 ug/l	4939650	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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

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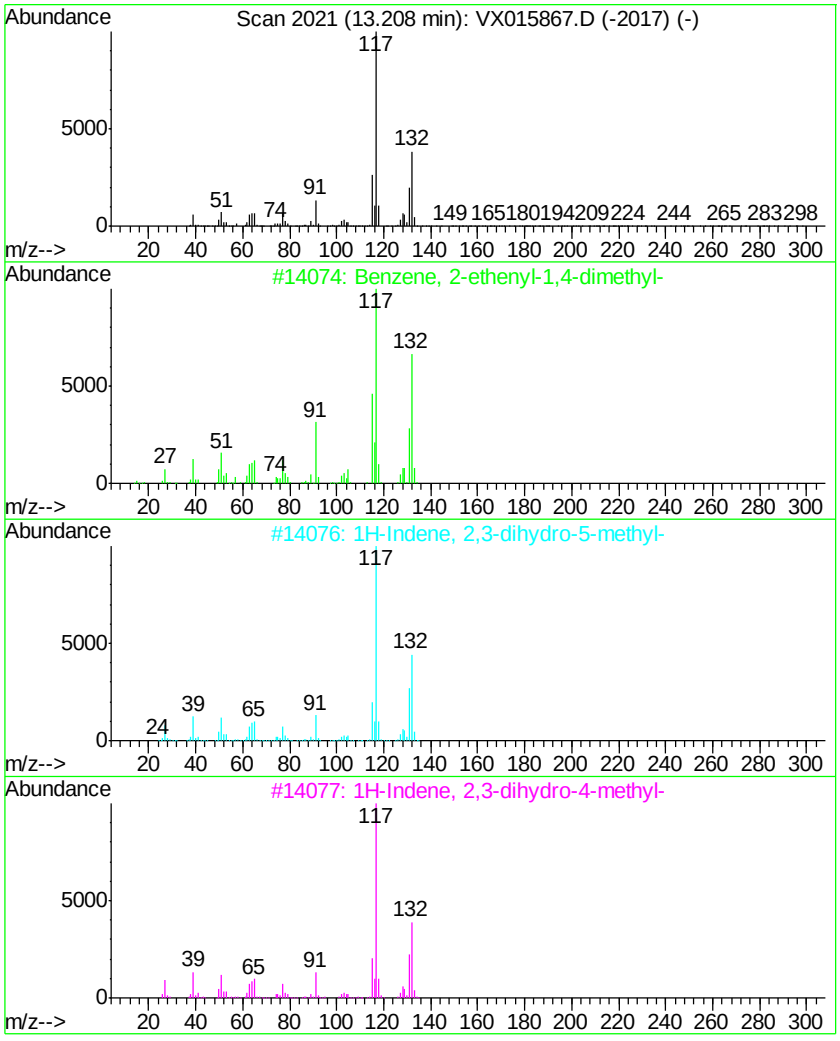
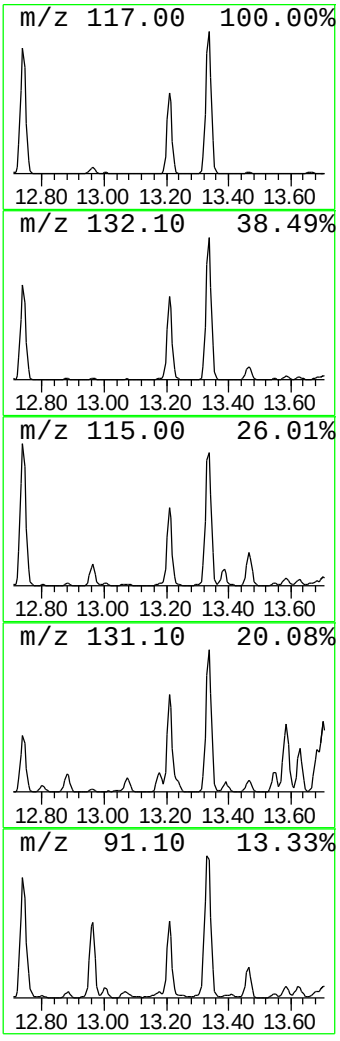
Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	69.32 ug/l	6419650	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
5	Indan, 1-methyl-	132	C10H12	000767-58-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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

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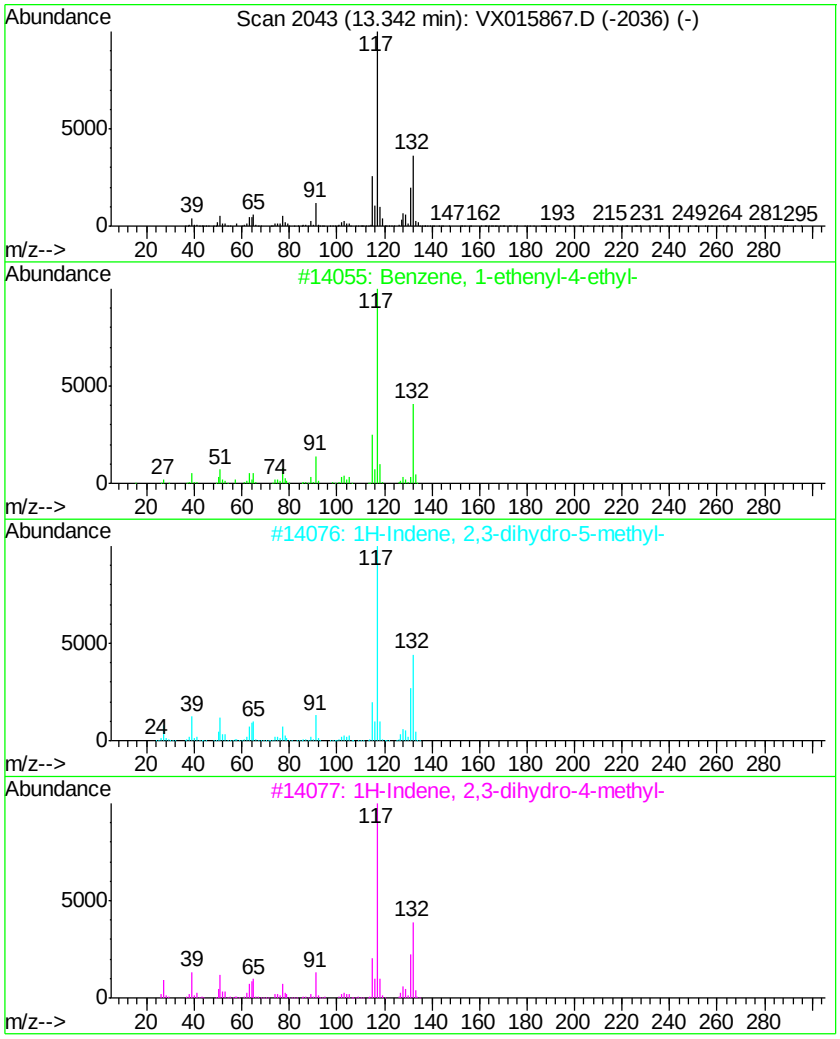
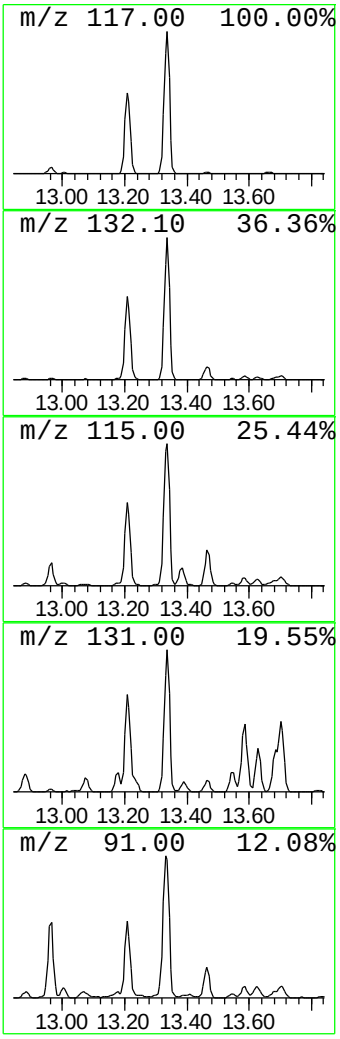
Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	138.57 ug/l	12832100	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015867.D
Acq On : 23 Apr 2020 02:27
Operator : JC/SP
Sample : L2380-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : KY001MW038-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Cyclopentane, met...	4.38	51.2	ug/l	9703830	1	5.63	9469630	50.0
Hexane, 2,2-dimet...	6.33	180.2	ug/l	13712000	2	6.84	3805710	50.0
Cyclopentene, 4,4...	8.20	51.4	ug/l	3909220	2	6.84	3805710	50.0
Indane	12.29	469.1	ug/l	43437200	4	12.06	4630230	50.0
o-Cymene	12.65	92.5	ug/l	8564780	4	12.06	4630230	50.0
Benzene, 1-etheny...	12.74	98.2	ug/l	9090550	4	12.06	4630230	50.0
Benzene, 1,2,3,4-...	12.96	53.3	ug/l	4939650	4	12.06	4630230	50.0
Benzene, 2-etheny...	13.21	69.3	ug/l	6419650	4	12.06	4630230	50.0
Benzene, 1-etheny...	13.34	138.6	ug/l	12832100	4	12.06	4630230	50.0