LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002537.D
Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

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\82X060418W.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	1.745	186	190	203	rBV	2320551	2550999	4.62%	0.310%
2	1.934	216	221	236	rVB	4201511	5787029	10.49%	0.704%
3	2.806	357	364	366	rBV	1528696	3036133	5.50%	0.369%
4	2.849	366	371	383	rVV	8389706	18604146	33.71%	2.262%
5	3.135	408	418	441	rVB	5472203	12404513	22.48%	1.508%
6	3.489	465	476	492	rBV	9074489	20424699	37.01%	2.484%
7	3.788	519	525	534	rVB	717048	1727301	3.13%	0.210%
8	4.281	595	606	612	rVV	896108	2753535	4.99%	0.335%
9	4.373	612	621	634	rVV	4281643	12431235	22.53%	1.512%
10	5.598	804	822	831	rVV4	6728222	27910189	50.58%	3.394%
11	5.714	831	841	861	rVB3	2478010	9557941	17.32%	1.162%
12	5.921	861	875	891	rBV	7347259	21815996	39.53%	2.653%
13	6.250	914	929	938	rBV2	2042964	7467374	13.53%	0.908%
14	6.348	938	945	953	rVV2	2216304	6253126	11.33%	0.760%
15	6.446	953	961	972	rVB	2358707	6601149	11.96%	0.803%
16	6.708	989	1004	1011	rBV	12284043	29710270	53.84%	3.613%
17	6.933	1022	1041	1048	rBV5	1722779	6921731	12.54%	0.842%
18	7.256	1085	1094	1100	rBV2	858628	2068596	3.75%	0.252%
19	7.464	1115	1128	1138	rBV	13334408	30743266	55.71%	3.738%
20	7.573	1138	1146	1151	rBV	1738823	3559472	6.45%	0.433%
21	7.634	1151	1156	1166	rVB	2305763	4534123	8.22%	0.551%
22	7.732	1166	1172	1182	rVB	1669348	3229325	5.85%	0.393%
23	7.842	1182	1190	1198	rVB2	1784501	3857075	6.99%	0.469%
24	8.031	1215	1221	1237	rVB4	1092718	4061043	7.36%	0.494%
25	8.171	1237	1244	1247	rBV3	1156746	2341661	4.24%	0.285%
26	8.214	1247	1251	1254	rVV	1441801	2069602	3.75%	0.252%
27	8.256	1254	1258	1262	rVB2	1611319	2476963	4.49%	0.301%
28	8.348	1267	1273	1276	rBV	7940455	13553123	24.56%	1.648%
29	8.384	1276	1279	1285	rVB	4094368	6113384	11.08%	0.743%
30	8.506	1294	1299	1308	rBV	8052665	14914144	27.03%	1.814%
31	8.573	1308	1310	1320	rVB5	1251024	2421993	4.39%	0.295%
32	8.671	1320	1326	1330	rBV3	3532809	7164838	12.98%	0.871%
33	8.841	1343	1354	1358	rBV	1637142	3729454	6.76%	0.454%
34	8.884	1358	1361	1363	rVV	1197408	1767261	3.20%	0.215%

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Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

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

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Title : SW846 8260

35	8.976	1370	1376	1383	rBV	11556516	18659387	33.81%	2.269%
36	9.043	1383	1387	1397	rVB4	1700285	3634886	6.59%	0.442%
37	9.165	1404	1407	1413	rVB2	1111689	1742497	3.16%	0.212%
38	9.226	1413	1417	1421	rBV2	1661991	2445266	4.43%	0.297%
39	9.445	1448	1453	1458	rBV	1693747	2450432	4.44%	0.298%
40	9.561	1467	1472	1479	rBV3	2745173	5337197	9.67%	0.649%
41	9.622	1479	1482	1486	rVB2	1635106	2224674	4.03%	0.271%
42	9.677	1486	1491	1495	rBV4	899061	1958014	3.55%	0.238%
43	9.823	1508	1515	1520	rBV3	1221364	2478920	4.49%	0.301%
44	9.878	1520	1524	1529	rVV2	1248550	2628240	4.76%	0.320%
45	9.927	1529	1532	1534	rVV2	1324362	1932789	3.50%	0.235%
46	9.963	1534	1538	1543	rVB	7217617	10221404	18.52%	1.243%
47	10.067	1548	1555	1560	rVB	6070245	8539817	15.48%	1.038%
48	10.262	1581	1587	1592	rVB	18952394	25777339	46.71%	3.135%
49	10.372	1599	1605	1609	rVB	30619928	42842842	77.64%	5.210%
50	10.420	1609	1613	1624	rVB	8164293	11643360	21.10%	1.416%
51	10.646	1645	1650	1654	rBV	1962119	2644761	4.79%	0.322%
52	10.731	1658	1664	1669	rBV	1172200	2107963	3.82%	0.256%
53	10.841	1674	1682	1685	rBV2	1849359	3338738	6.05%	0.406%
54	10.914	1685	1694	1698	rBV2	1421122	2646128	4.80%	0.322%
55	11.024	1706	1712	1719	rBV	2090920	3428321	6.21%	0.417%
56	11.085	1719	1722	1726	rBV2	1213172	1678327	3.04%	0.204%
57	11.146	1726	1732	1734	rBV3	2513878	4280654	7.76%	0.521%
58	11.170	1734	1736	1742	rVB2	2759861	3879549	7.03%	0.472%
59	11.250	1742	1749	1753	rBV	2606217	3456270	6.26%	0.420%
60	11.365	1764	1768	1775	rVB	7575912	9268404	16.80%	1.127%
61	11.445	1775	1781	1788	rBV	19833717	39892305	72.29%	4.851%
62	11.512	1788	1792	1803	rVB2	14537152	24775275	44.90%	3.013%
63	11.676	1814	1819	1825	rBV	8391827	11232480	20.36%	1.366%
64	11.823	1837	1843	1850	rVB	39948596	55181540	100.00%	6.710%
65	12.085	1881	1886	1891	rBV3	1553219	2696480	4.89%	0.328%
66	12.152	1891	1897	1901	rBV	6740040	9494497	17.21%	1.155%
67	12.304	1917	1922	1925	rBV	9299968	14328662	25.97%	1.742%
68	12.329	1925	1926	1930	rVV	8022992	7180935	13.01%	0.873%
69	12.377	1930	1934	1941	rVB3	13346406	20538026	37.22%	2.497%
70	12.481	1946	1951	1955	rVV3	2851452	4082525	7.40%	0.496%
71	12.524	1955	1958	1965	rVV	2982365	3691673	6.69%	0.449%

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ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

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

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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Title : SW846 8260

72	12.591	1965	1969	1971	rVV	5559809	6853772	12.42%	0.833%
73	12.615	1971	1973	1979	rVV	5702119	6985128	12.66%	0.849%
74	12.676	1979	1983	1986	rVV	10148299	11916479	21.60%	1.449%
75	12.707	1986	1988	1992	rVV	1680790	1996936	3.62%	0.243%
76	12.762	1992	1997	2005	rVB3	4696200	7715503	13.98%	0.938%
77	12.908	2018	2021	2025	rVB	2041375	2397939	4.35%	0.292%
78	12.987	2025	2034	2037	rBV	5262376	7790208	14.12%	0.947%
79	13.024	2037	2040	2046	rVB	7848590	9385009	17.01%	1.141%
80	13.085	2046	2050	2051	rBV	1774908	1877793	3.40%	0.228%
81	13.109	2051	2054	2056	rVV3	2140513	3082893	5.59%	0.375%
82	13.133	2056	2058	2061	rVV	1920751	1891506	3.43%	0.230%
83	13.231	2067	2074	2078	rVV3	5652502	8881534	16.10%	1.080%
84	13.274	2078	2081	2084	rVV2	1847088	2116533	3.84%	0.257%
85	13.322	2084	2089	2091	rVV3	2409686	3720215	6.74%	0.452%
86	13.359	2091	2095	2100	rVV2	9043690	12986584	23.53%	1.579%
87	13.432	2104	2107	2112	rVB	2023691	2398654	4.35%	0.292%
88	13.487	2112	2116	2124	rVB3	1735917	2826822	5.12%	0.344%
89	13.566	2124	2129	2132	rBV2	1767242	2413708	4.37%	0.294%
90	13.609	2132	2136	2139	rVV	2772443	3586097	6.50%	0.436%
91	13.646	2139	2142	2147	rVV3	3185422	5541333	10.04%	0.674%
92	13.731	2153	2156	2161	rVV2	2673211	3788400	6.87%	0.461%
93	13.841	2170	2174	2179	rVB	6171297	7095529	12.86%	0.863%
94	13.926	2184	2188	2192	rVB3	1171058	1844568	3.34%	0.224%
95	13.987	2192	2198	2203	rBV4	1761627	3494596	6.33%	0.425%
96	14.139	2219	2223	2227	rBV2	2607148	3106657	5.63%	0.378%
97	14.280	2241	2246	2255	rBV	2373605	4148686	7.52%	0.504%
98	14.438	2268	2272	2276	rVB	1546556	1882583	3.41%	0.229%
99	14.700	2311	2315	2326	rVB	6255884	8571167	15.53%	1.042%
100	14.847	2335	2339	2348	rVB	2351048	3157999	5.72%	0.384%

Sum of corrected areas: 822358097

LSC Report - Integrated Chromatogram

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Acq On : 14 Jun 2018 04:54
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Sample : J3383-01
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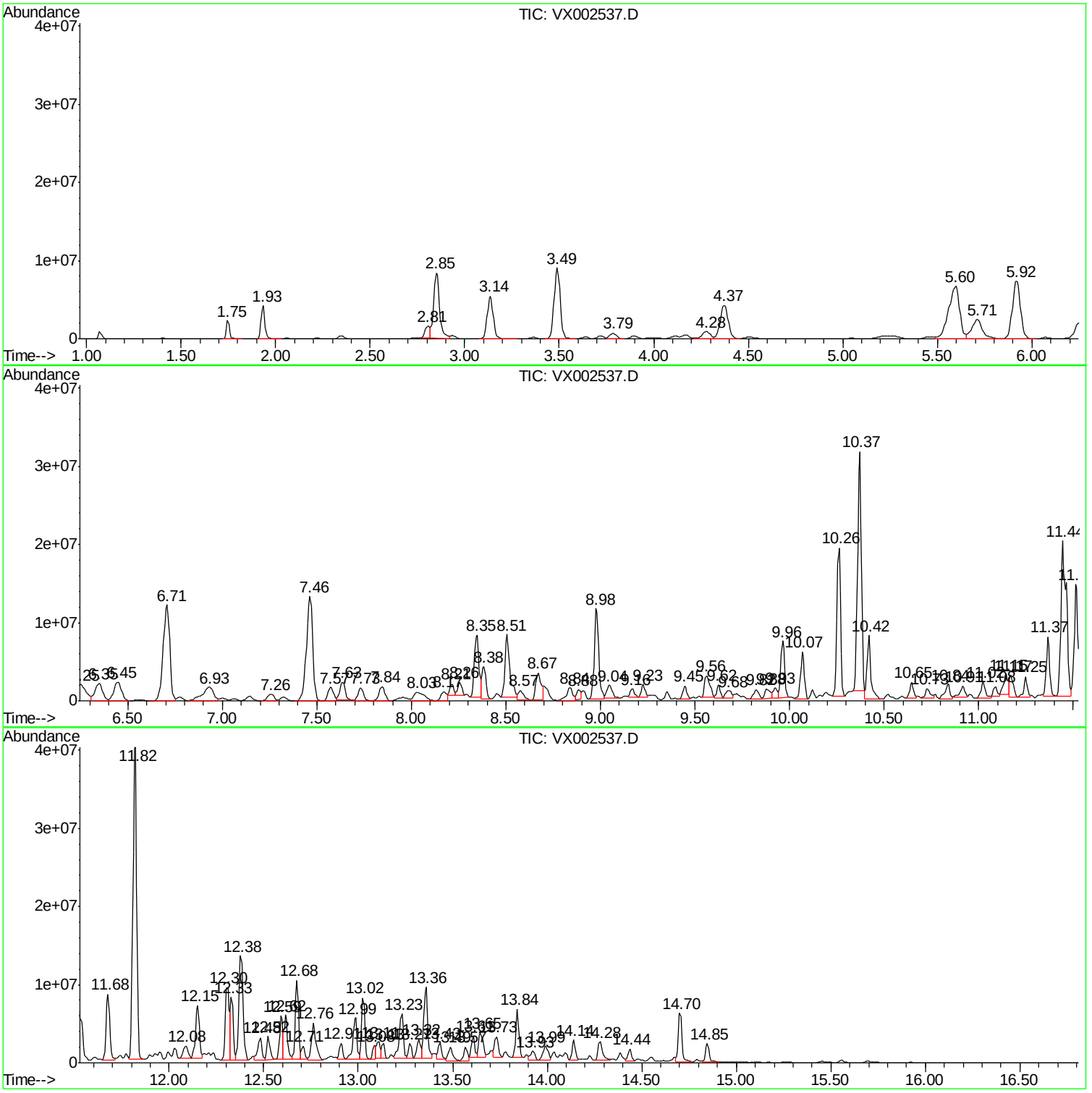
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061318\
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Instrument : MSVOA_X
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

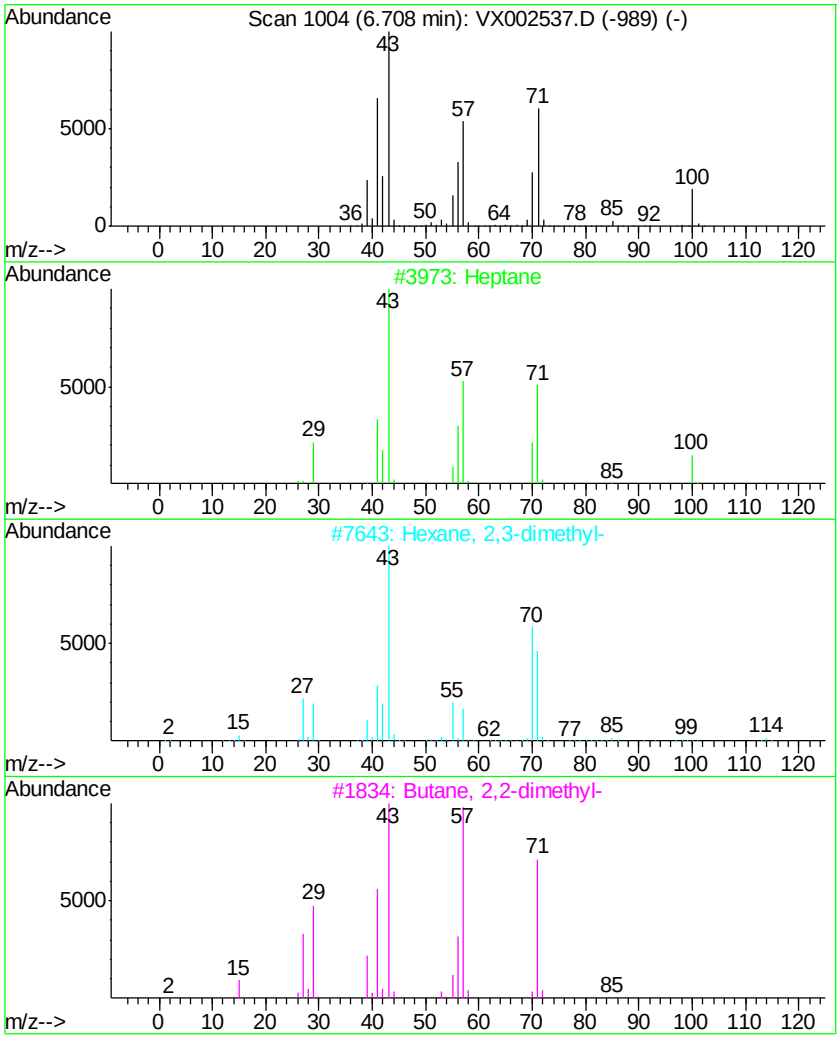
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

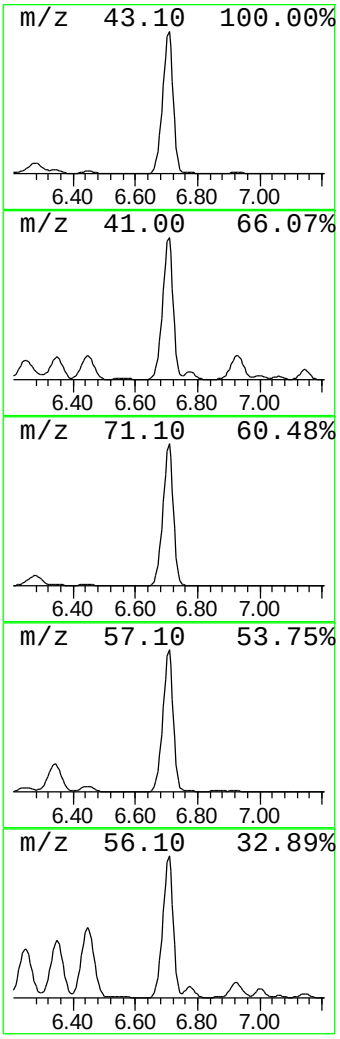
****

Peak Number 1 Heptane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	214.62 ug/l	29710300	1,4-Difluorobenzene	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	91
2	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40
3	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	38
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	38
5	Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	38





Library Search Compound Report

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Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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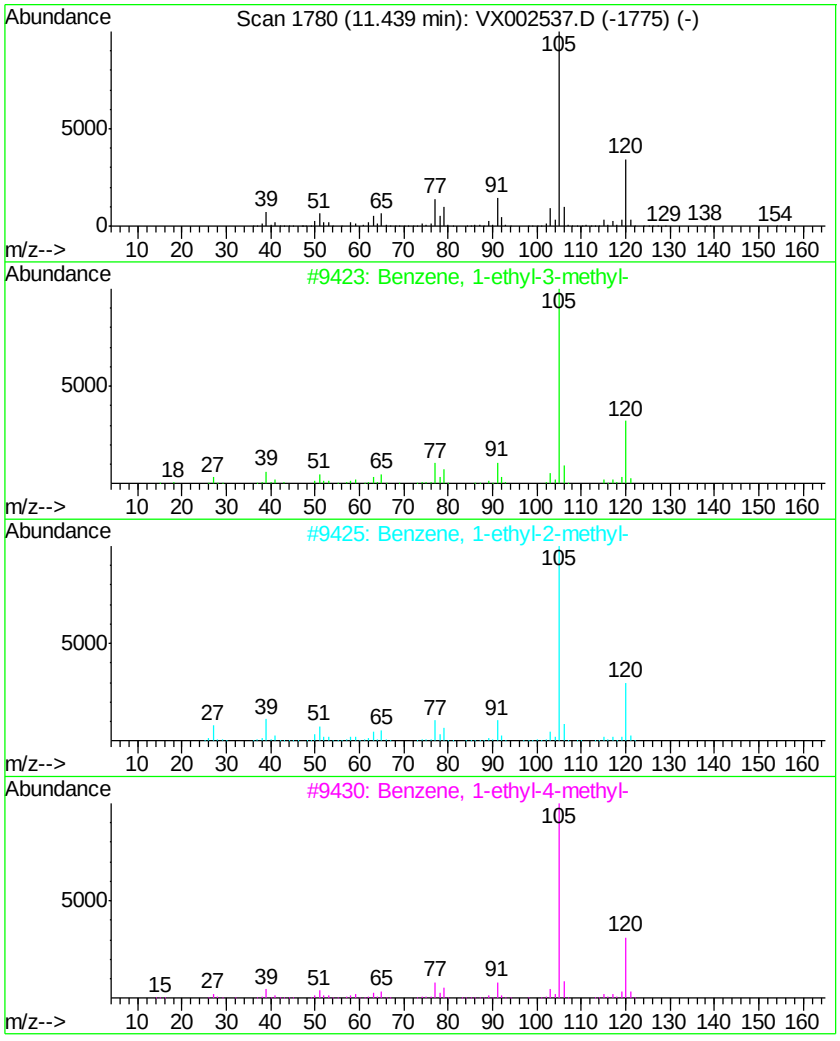
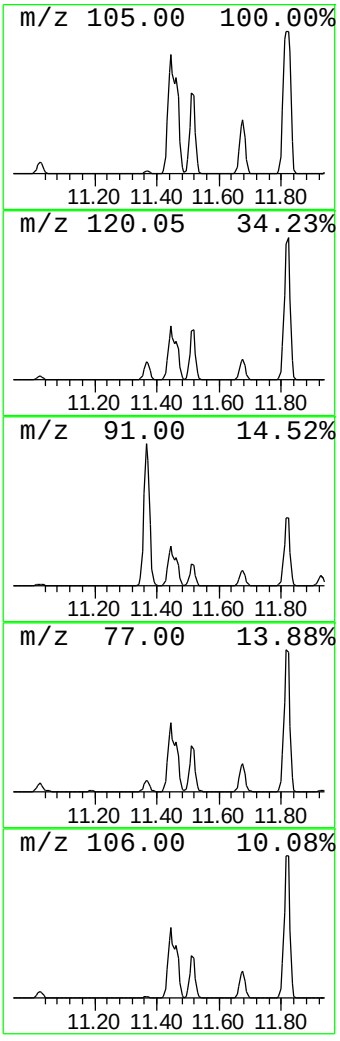
Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	739.71 ug/l	39892300	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90





Library Search Compound Report

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Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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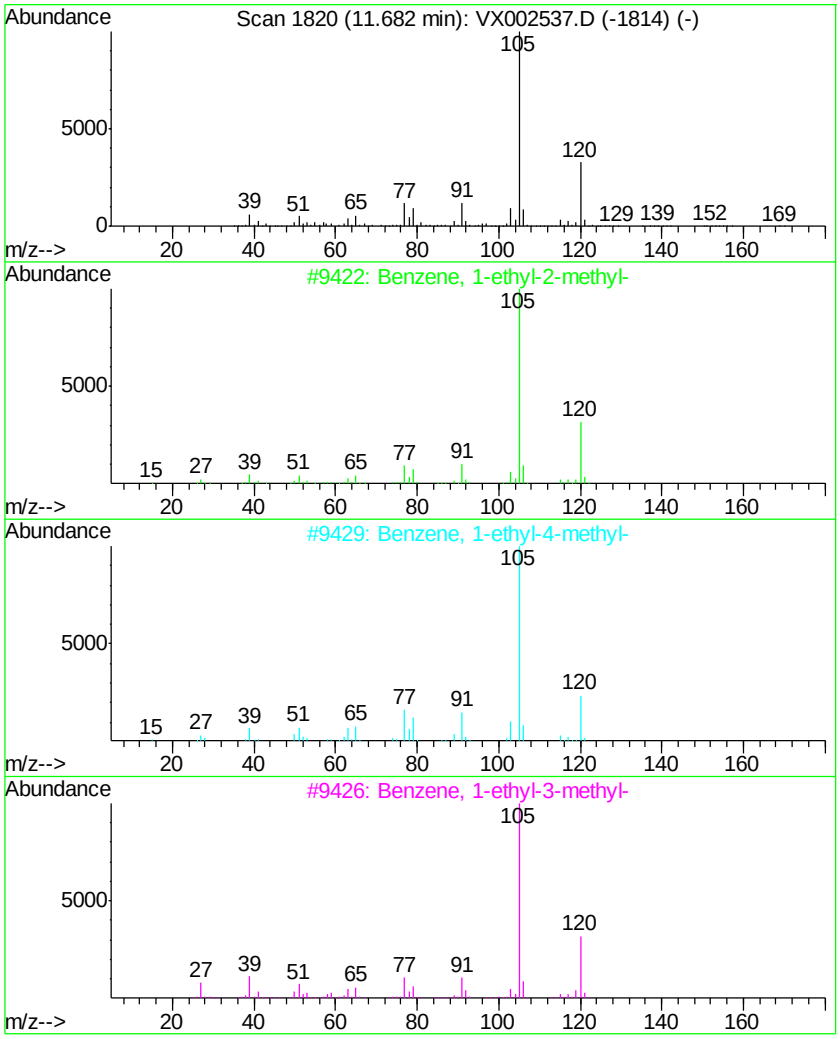
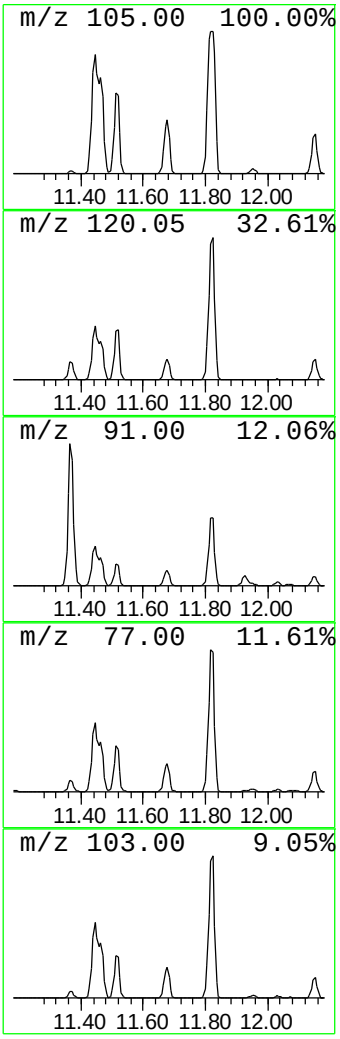
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.68	208.28 ug/l	11232500	1,4-Dichlorobenzene-d4	12.08

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-4-methyl-	120	C9H12	000622-96-8	93
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	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
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Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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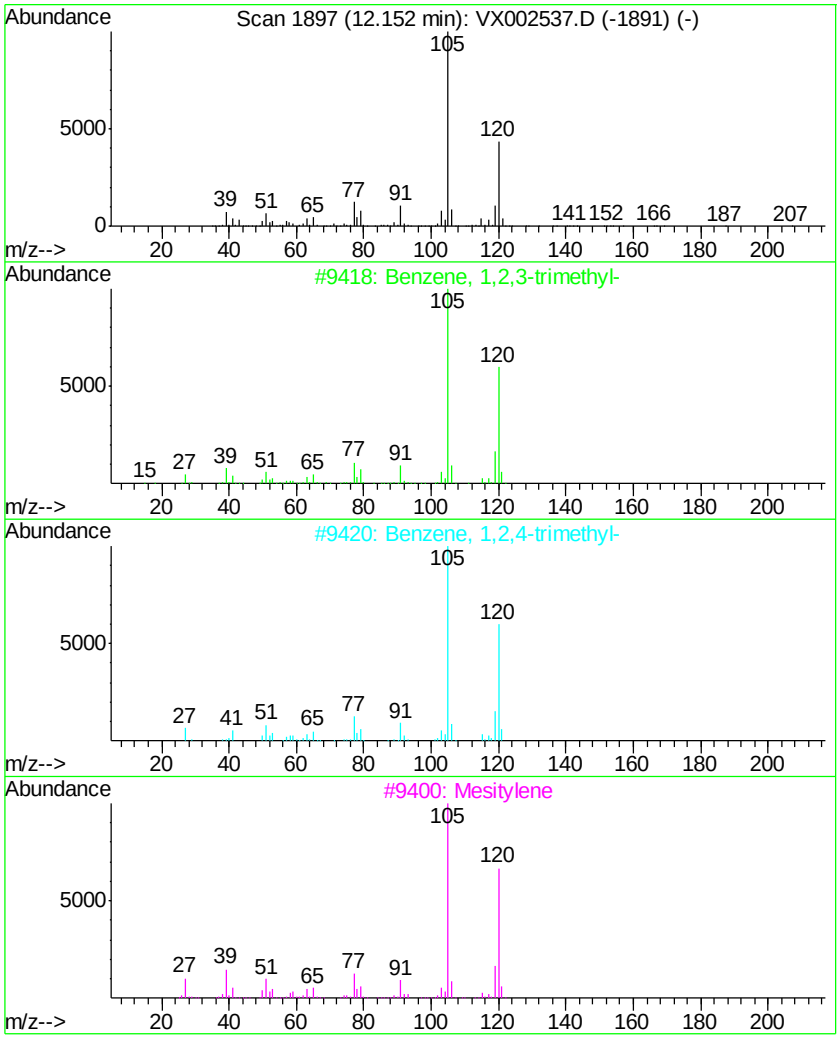
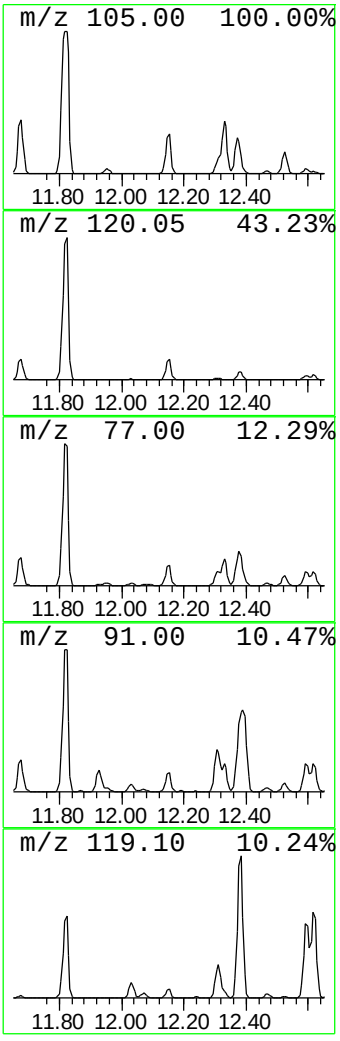
Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	176.05 ug/l	9494500	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Library Search Compound Report

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Data File : VX002537.D
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ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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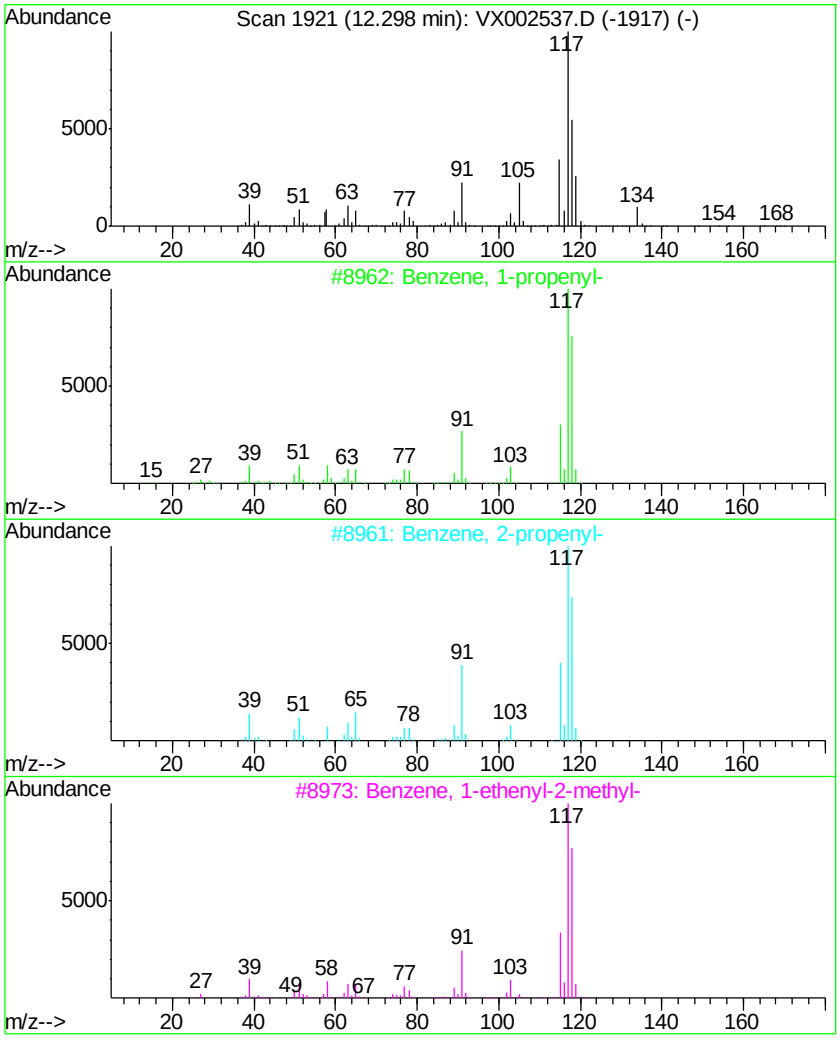
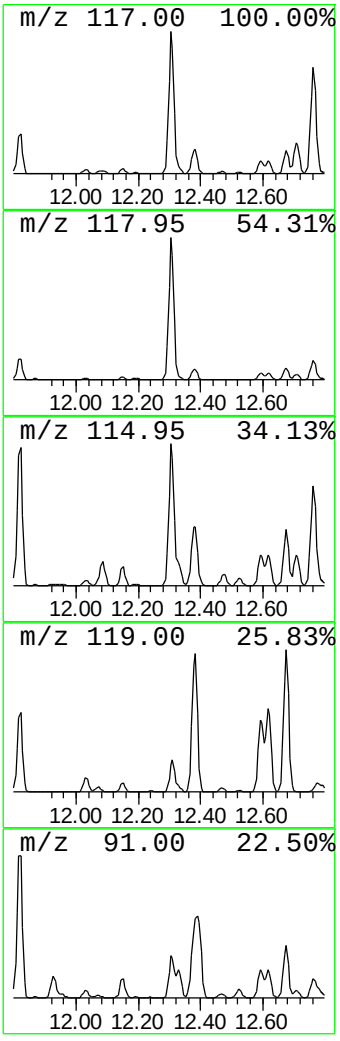
Peak Number 5 Benzene, 1-propenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	265.69 ug/l	14328700	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5	Indane	118	C9H10	000496-11-7	60





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Data File : VX002537.D
Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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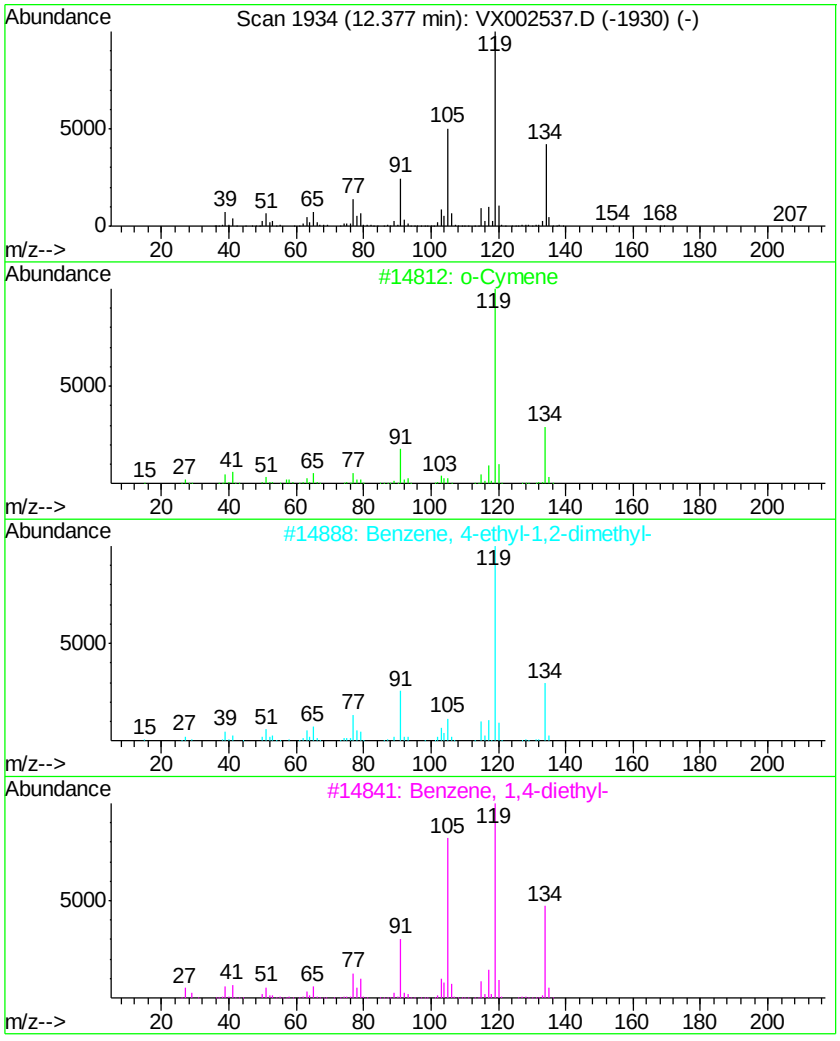
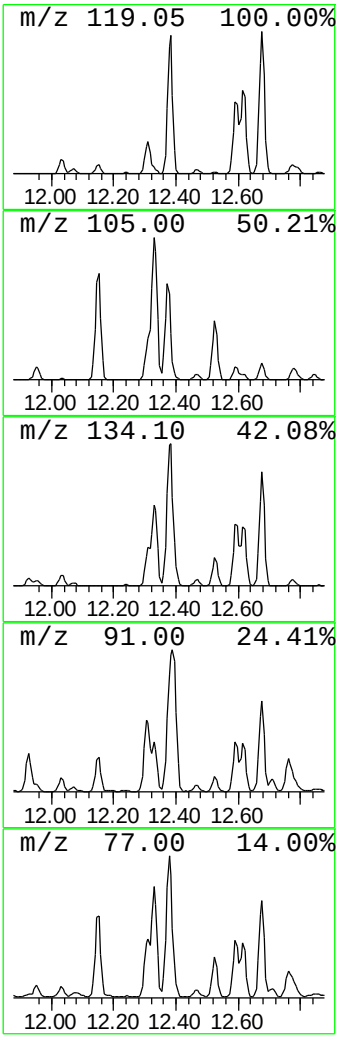
Peak Number 6 o-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	380.83 ug/l	20538000	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	94
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
Data File : VX002537.D
Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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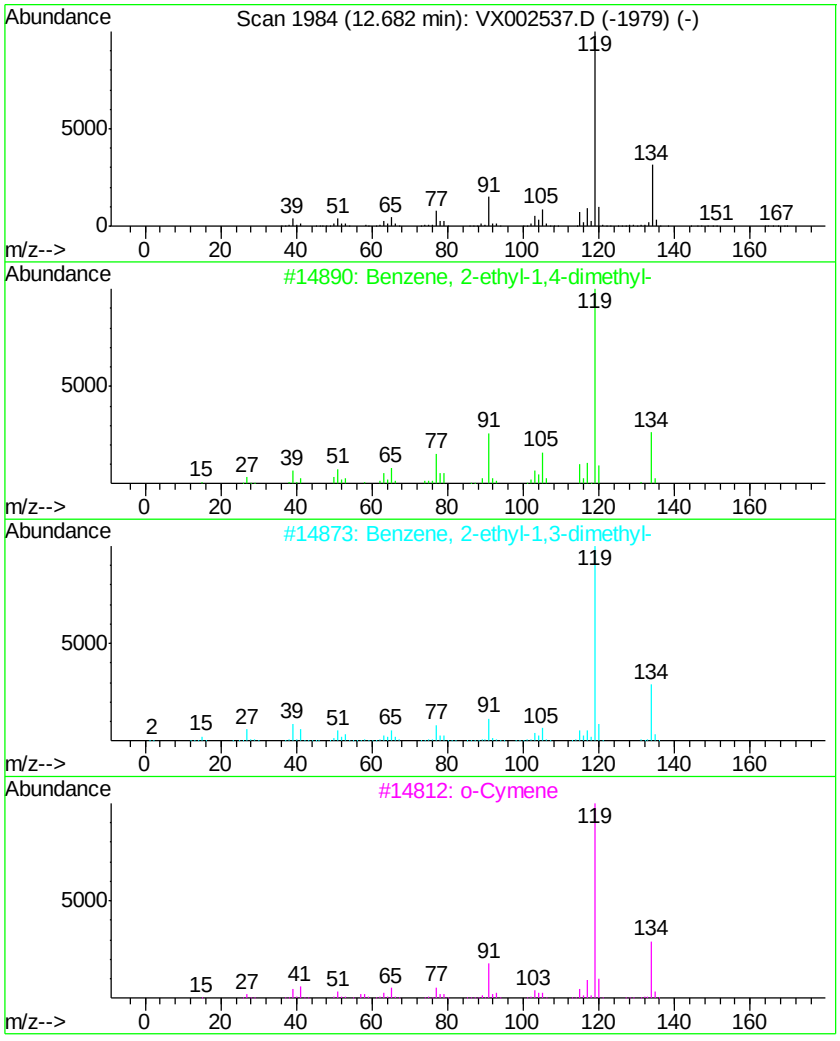
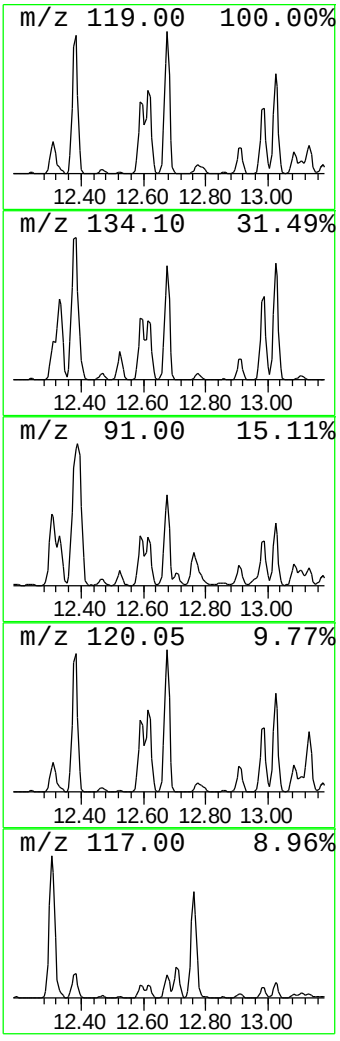
Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	220.96 ug/l	11916500	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
Data File : VX002537.D
Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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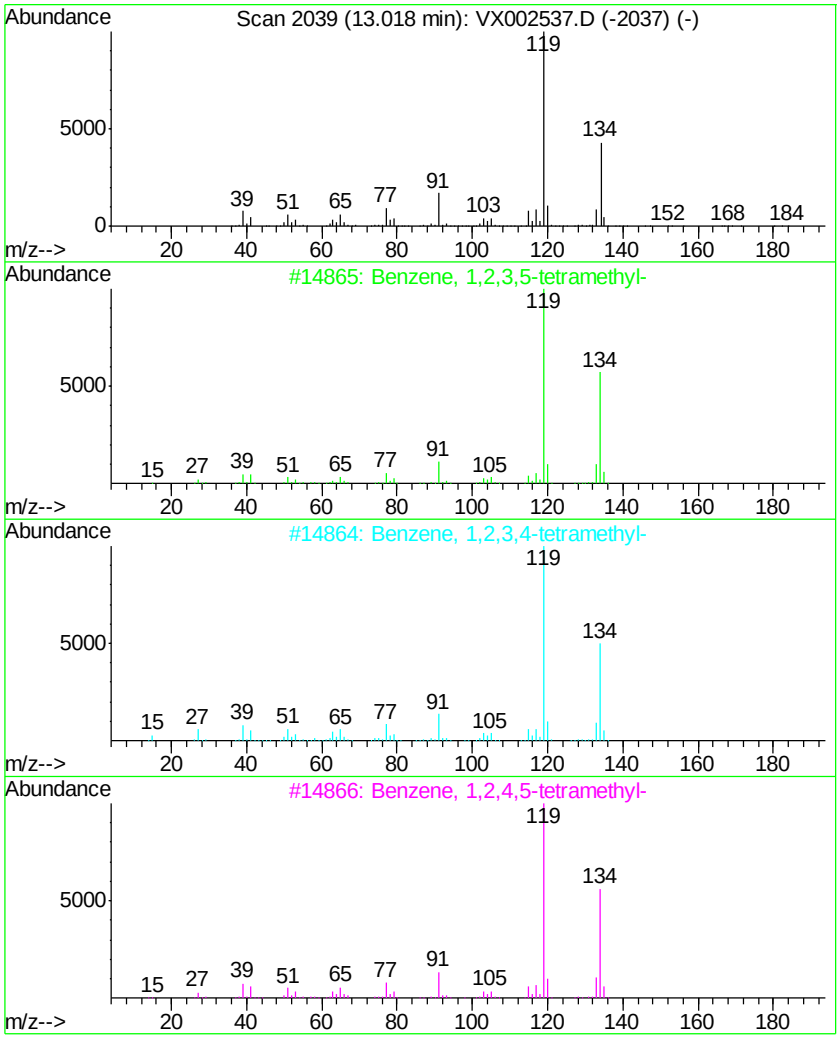
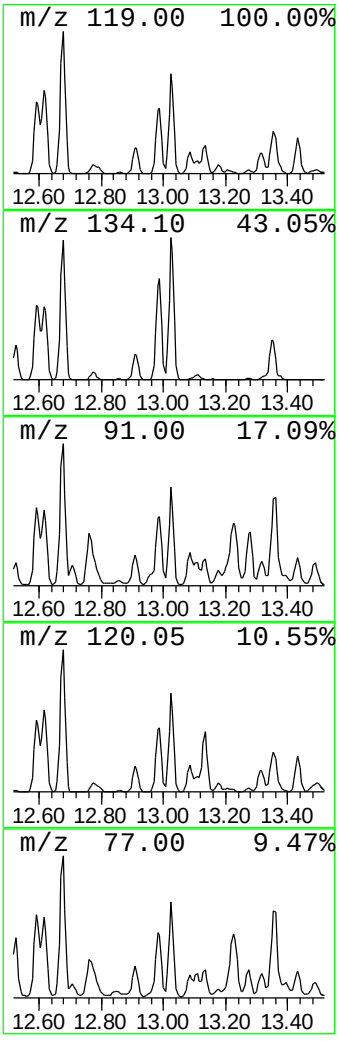
Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	174.02 ug/l	9385010	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002537.D
Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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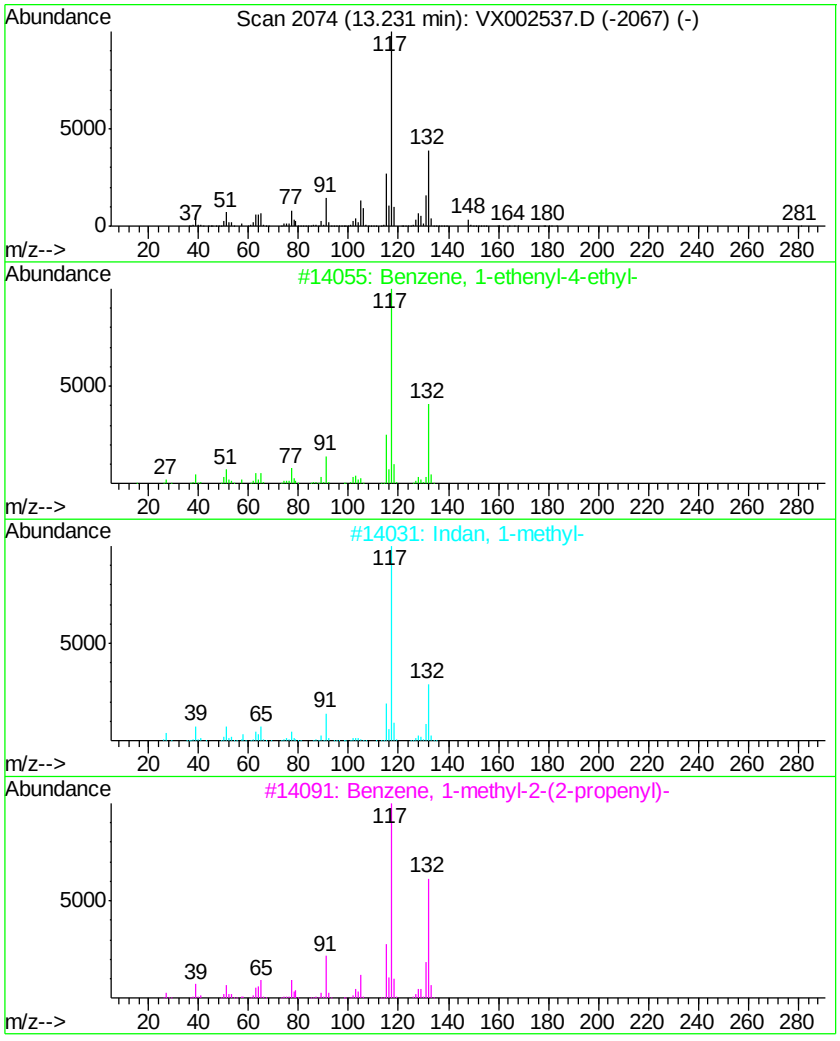
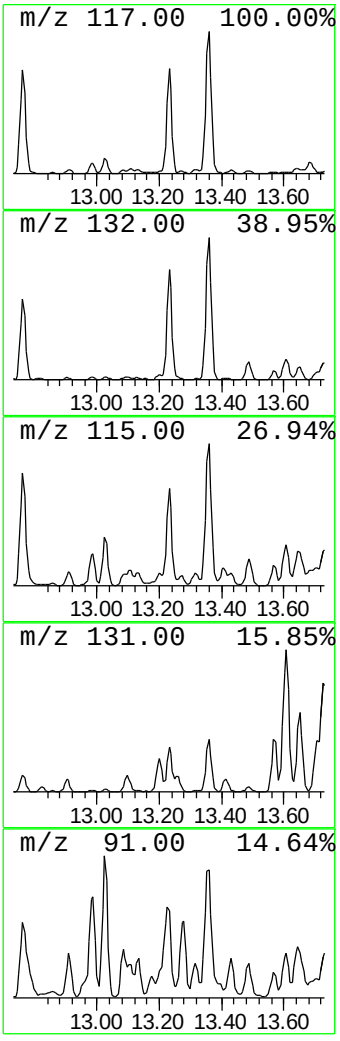
Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.23	164.69 ug/l	8881530	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Indan, 1-methyl-	132	C10H12	000767-58-8	93
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	93
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
Data File : VX002537.D
Acq On : 14 Jun 2018 04:54
Operator : JC/MD
Sample : J3383-01
Misc : 6.86g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-2RR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.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
Heptane	6.71	214.6	ug/l	29710300	2	6.87	6921730	50.0
Benzene, 1-ethyl-...	11.44	739.7	ug/l	39892300	4	12.08	2696480	50.0
Benzene, 1-ethyl-...	11.68	208.3	ug/l	11232500	4	12.08	2696480	50.0
Benzene, 1,2,3-tr...	12.15	176.1	ug/l	9494500	4	12.08	2696480	50.0
Benzene, 1-propenyl-	12.30	265.7	ug/l	14328700	4	12.08	2696480	50.0
o-Cymene	12.38	380.8	ug/l	20538000	4	12.08	2696480	50.0
Benzene, 2-ethyl-...	12.68	221.0	ug/l	11916500	4	12.08	2696480	50.0
Benzene, 1,2,3,5-...	13.02	174.0	ug/l	9385010	4	12.08	2696480	50.0
Benzene, 1-etheny...	13.23	164.7	ug/l	8881530	4	12.08	2696480	50.0