LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

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\82X092618W.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	2.879	278	283	295	rVB	1085838	2333281	3.48%	0.357%
2	3.160	320	329	349	rVB	780139	1816109	2.71%	0.278%
3	3.513	378	387	401	rBV	760383	1758584	2.63%	0.269%
4	4.300	506	516	523	rVV	691359	2075414	3.10%	0.317%
5	4.391	523	531	543	rVB3	779171	2210979	3.30%	0.338%
6	5.293	652	679	696	rVB4	289193	1260842	1.88%	0.193%
7	5.610	719	731	738	rVV2	1602671	5186041	7.74%	0.793%
8	5.720	739	749	770	rVB	2258739	8064862	12.04%	1.233%
9	5.921	770	782	797	rBV4	1781606	5251355	7.84%	0.803%
10	6.257	824	837	841	rBV2	536411	1580412	2.36%	0.242%
11	6.348	842	852	863	rVV2	6059257	18781432	28.04%	2.871%
12	6.458	863	870	879	rVB4	497812	1395199	2.08%	0.213%
13	6.702	898	910	918	rBV	1519018	3683097	5.50%	0.563%
14	6.933	940	948	956	rVB3	773576	2527180	3.77%	0.386%
15	7.256	991	1001	1008	rBV4	520750	1246031	1.86%	0.190%
16	7.458	1022	1034	1045	rBV5	1644012	4455434	6.65%	0.681%
17	7.573	1045	1053	1058	rBV	1349834	2735719	4.08%	0.418%
18	7.634	1058	1063	1069	rBV	1505103	2872997	4.29%	0.439%
19	7.732	1075	1079	1090	rVB	567592	1079248	1.61%	0.165%
20	7.848	1090	1098	1105	rVB2	519251	1093803	1.63%	0.167%
21	8.055	1122	1132	1144	rVB	4847839	10131245	15.13%	1.549%
22	8.177	1144	1152	1161	rBV2	4068486	9755995	14.57%	1.491%
23	8.256	1161	1165	1174	rVB2	2160846	3841963	5.74%	0.587%
24	8.342	1174	1179	1183	rBV	2615423	4346794	6.49%	0.664%
25	8.378	1183	1185	1191	rVB2	1358680	1873050	2.80%	0.286%
26	8.500	1201	1205	1214	rBV	2765158	4999075	7.46%	0.764%
27	8.573	1214	1217	1227	rVB4	660991	1378290	2.06%	0.211%
28	8.677	1227	1234	1238	rBV3	2328441	4872944	7.28%	0.745%
29	8.787	1246	1252	1258	rBV	12153217	18936289	28.27%	2.894%
30	8.835	1258	1260	1264	rVV3	648211	1045315	1.56%	0.160%
31	8.884	1264	1268	1276	rVV2	756596	1797695	2.68%	0.275%
32	8.976	1276	1283	1290	rVV2	2824947	5208598	7.78%	0.796%
33	9.049	1290	1295	1304	rVV4	638602	1668630	2.49%	0.255%
34	9.165	1311	1314	1320	rVV2	655191	1257546	1.88%	0.192%

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ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Title : SW846 8260

35	9.220	1320	1323	1327	rVV2	721753	1226906	1.83%	0.188%
36	9.262	1327	1330	1340	rVB5	538699	1146461	1.71%	0.175%
37	9.445	1355	1360	1365	rBV2	752551	1188165	1.77%	0.182%
38	9.555	1374	1378	1385	rBV3	1368147	2717341	4.06%	0.415%
39	9.817	1415	1421	1427	rBV4	499797	981475	1.47%	0.150%
40	9.878	1427	1431	1435	rBV2	572682	1097620	1.64%	0.168%
41	9.963	1441	1445	1450	rVB	3414277	4587743	6.85%	0.701%
42	10.067	1457	1462	1466	rVB	2728478	3634660	5.43%	0.556%
43	10.116	1466	1470	1473	rBV	1020124	1271666	1.90%	0.194%
44	10.158	1473	1477	1480	rBV	767494	984246	1.47%	0.150%
45	10.256	1488	1493	1498	rVB	13178145	17345617	25.90%	2.651%
46	10.372	1505	1512	1516	rVV2	43281957	66977948	100.00%	10.237%
47	10.414	1516	1519	1531	rVB2	2948228	5630119	8.41%	0.861%
48	10.646	1552	1557	1560	rBV	1215874	1600871	2.39%	0.245%
49	10.707	1560	1567	1575	rVB	21226816	29148289	43.52%	4.455%
50	10.835	1580	1588	1592	rBV	962554	1746363	2.61%	0.267%
51	10.914	1592	1601	1605	rBV3	781900	1460932	2.18%	0.223%
52	11.018	1612	1618	1623	rBV	2153342	3010792	4.50%	0.460%
53	11.085	1625	1629	1632	rVV	707369	961377	1.44%	0.147%
54	11.140	1632	1638	1640	rVV2	2016821	3075476	4.59%	0.470%
55	11.170	1640	1643	1649	rVB2	2202559	3601160	5.38%	0.550%
56	11.250	1649	1656	1659	rBV	1376275	1909961	2.85%	0.292%
57	11.365	1664	1675	1682	rVV	7071724	10266039	15.33%	1.569%
58	11.445	1682	1688	1695	rVV	29167635	56138620	83.82%	8.581%
59	11.512	1695	1699	1710	rVB	15624881	21995909	32.84%	3.362%
60	11.670	1720	1725	1732	rBV	12646171	15713176	23.46%	2.402%
61	11.817	1743	1749	1755	rVB	45270357	59439303	88.74%	9.085%
62	11.951	1768	1771	1774	rVB	968955	1162509	1.74%	0.178%
63	12.030	1780	1784	1787	rVB	1802911	2059129	3.07%	0.315%
64	12.079	1787	1792	1797	rVB4	1312631	2291810	3.42%	0.350%
65	12.146	1797	1803	1808	rBV	12190273	15550823	23.22%	2.377%
66	12.304	1823	1829	1831	rBV	11550832	15620384	23.32%	2.388%
67	12.329	1831	1833	1836	rVV	8089399	8261758	12.34%	1.263%
68	12.378	1836	1841	1847	rVB3	11891725	18526895	27.66%	2.832%
69	12.475	1847	1857	1861	rBV3	1784852	3286266	4.91%	0.502%
70	12.524	1861	1865	1871	rVB2	3720948	5415853	8.09%	0.828%
71	12.591	1871	1876	1878	rBV	5972036	8073552	12.05%	1.234%

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ClientSampleId : B-7(8-8.5)

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Title : SW846 8260

72	12.615	1878	1880	1884	rVV	5537075	5588461	8.34%	0.854%
73	12.670	1884	1889	1893	rVV	9945392	11971429	17.87%	1.830%
74	12.707	1893	1895	1899	rVV	1093474	1159472	1.73%	0.177%
75	12.762	1899	1904	1914	rVB2	3337103	6471085	9.66%	0.989%
76	12.908	1923	1928	1932	rVB	2171498	2711648	4.05%	0.414%
77	12.981	1932	1940	1943	rBV	5072508	6721424	10.04%	1.027%
78	13.024	1943	1947	1952	rVB	7037473	8548847	12.76%	1.307%
79	13.085	1952	1957	1958	rBV	1245730	1640096	2.45%	0.251%
80	13.103	1958	1960	1962	rVV	1399669	1619185	2.42%	0.247%
81	13.127	1962	1964	1967	rVB	1190651	1167470	1.74%	0.178%
82	13.225	1973	1980	1984	rVB3	4570560	6699947	10.00%	1.024%
83	13.268	1984	1987	1991	rVB	1051248	1187036	1.77%	0.181%
84	13.316	1991	1995	1997	rBV3	1398777	2072619	3.09%	0.317%
85	13.353	1997	2001	2006	rVV2	7898402	10923586	16.31%	1.670%
86	13.402	2006	2009	2011	rVV3	769385	1034224	1.54%	0.158%
87	13.426	2011	2013	2018	rVB	1312240	1457567	2.18%	0.223%
88	13.481	2018	2022	2031	rVB	1536822	2186187	3.26%	0.334%
89	13.560	2031	2035	2038	rBV2	824760	1153373	1.72%	0.176%
90	13.603	2038	2042	2045	rBV	1299968	1648953	2.46%	0.252%
91	13.639	2045	2048	2054	rVB3	1920394	3257610	4.86%	0.498%
92	13.725	2059	2062	2067	rVB2	1426996	2250593	3.36%	0.344%
93	13.835	2075	2080	2086	rVB	6860208	8827619	13.18%	1.349%
94	13.987	2098	2105	2109	rBV4	992621	1857418	2.77%	0.284%
95	14.133	2125	2129	2133	rBV	1529718	1857803	2.77%	0.284%
96	14.274	2148	2152	2157	rBV	1415822	2151397	3.21%	0.329%
97	14.377	2165	2169	2175	rBV	798831	1165467	1.74%	0.178%
98	14.432	2175	2178	2182	rVB	1027174	1186003	1.77%	0.181%
99	14.700	2217	2222	2232	rVB	7445684	9641226	14.39%	1.474%
100	14.840	2241	2245	2254	rVB	3452453	4459416	6.66%	0.682%

Sum of corrected areas: 654245823

LSC Report - Integrated Chromatogram

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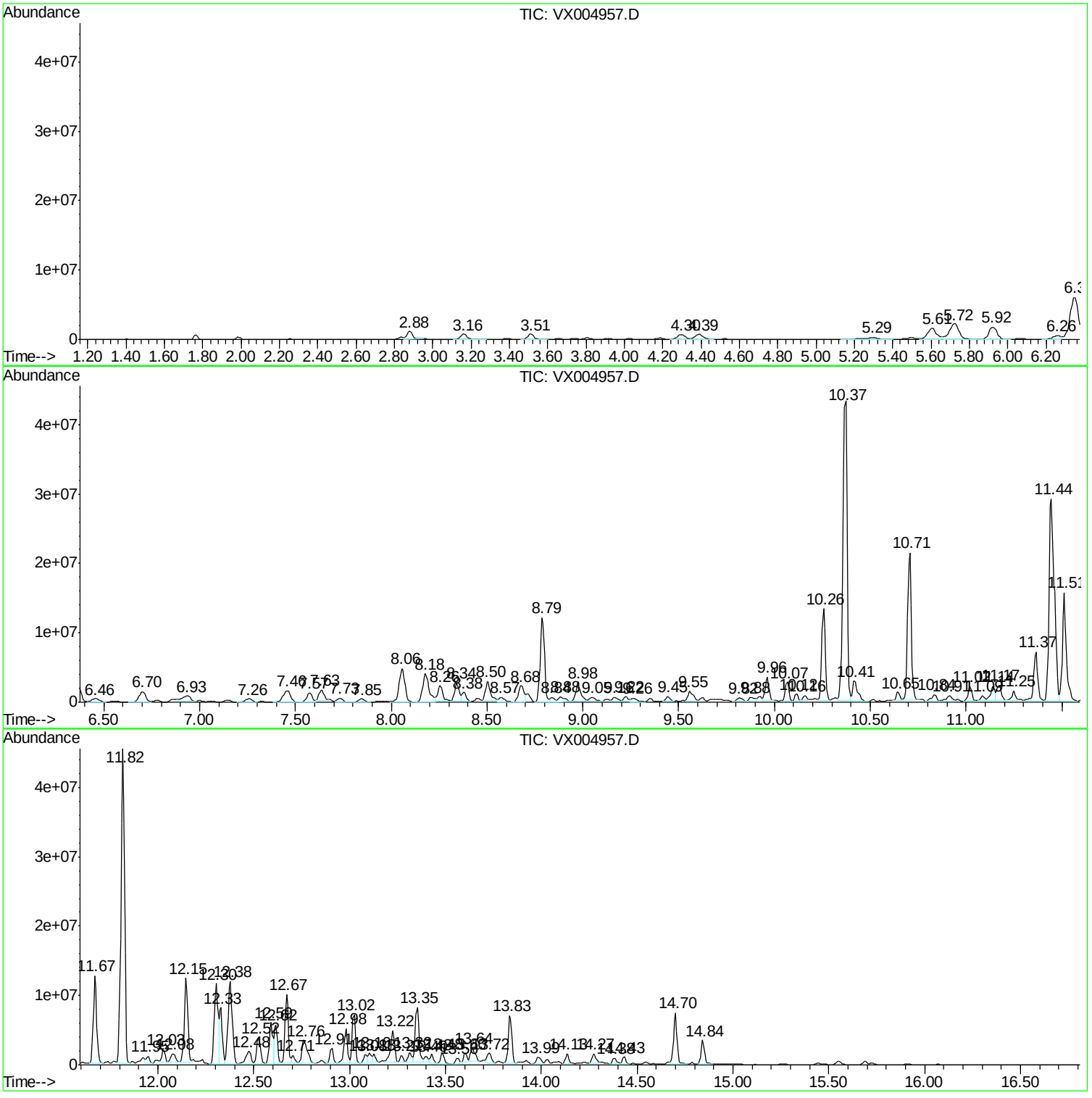
Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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\VX100318\
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ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

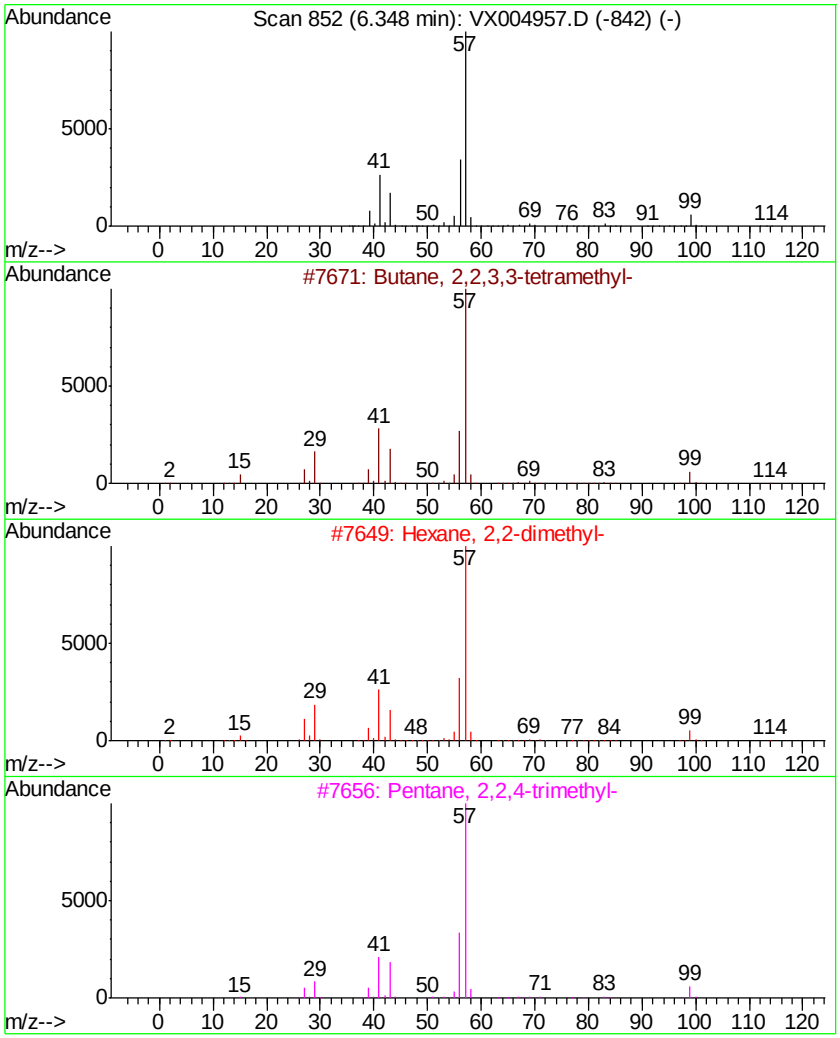
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

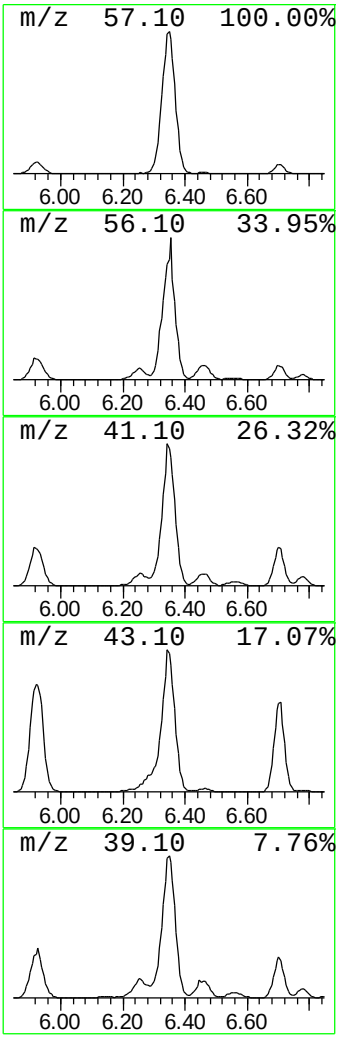
****

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	371.59 ug/l	18781400	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
5	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64





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Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

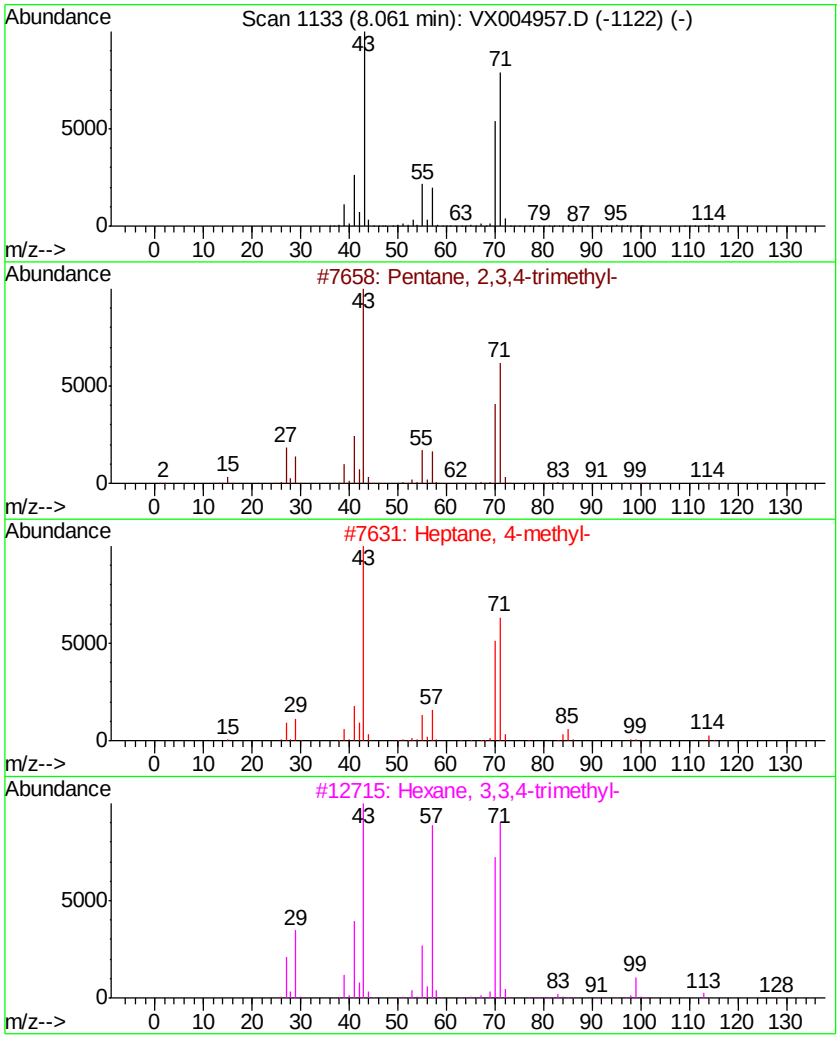
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

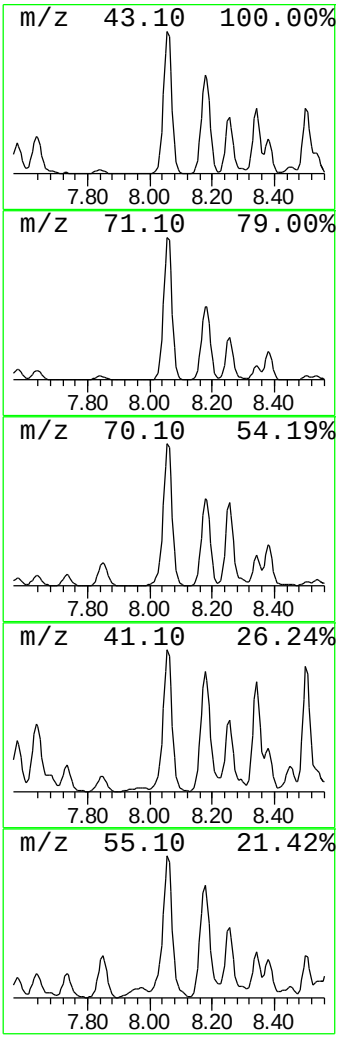
****

Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	200.45 ug/l	10131200	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
4	Pentane, 3-ethyl-	100	C7H16	000617-78-7	74
5	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64





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Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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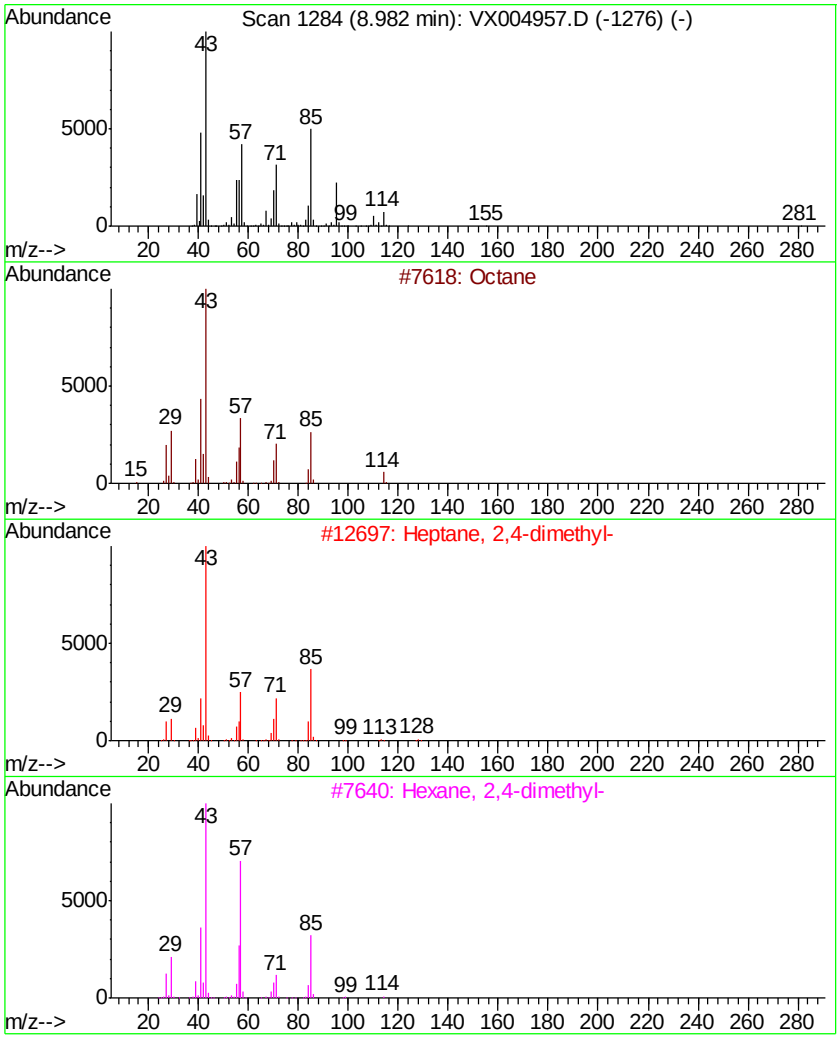
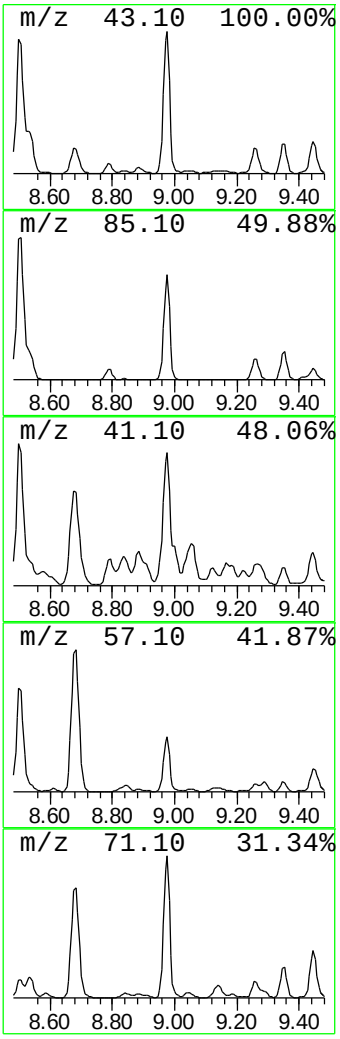
Peak Number 3 Octane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	204.79 ug/l	5208600	Chlorobenzene-d5	10.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	94
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
4	Oxalic acid, diisohexyl ester	258	C14H26O4	1000309-32-9	59
5	Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	53





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

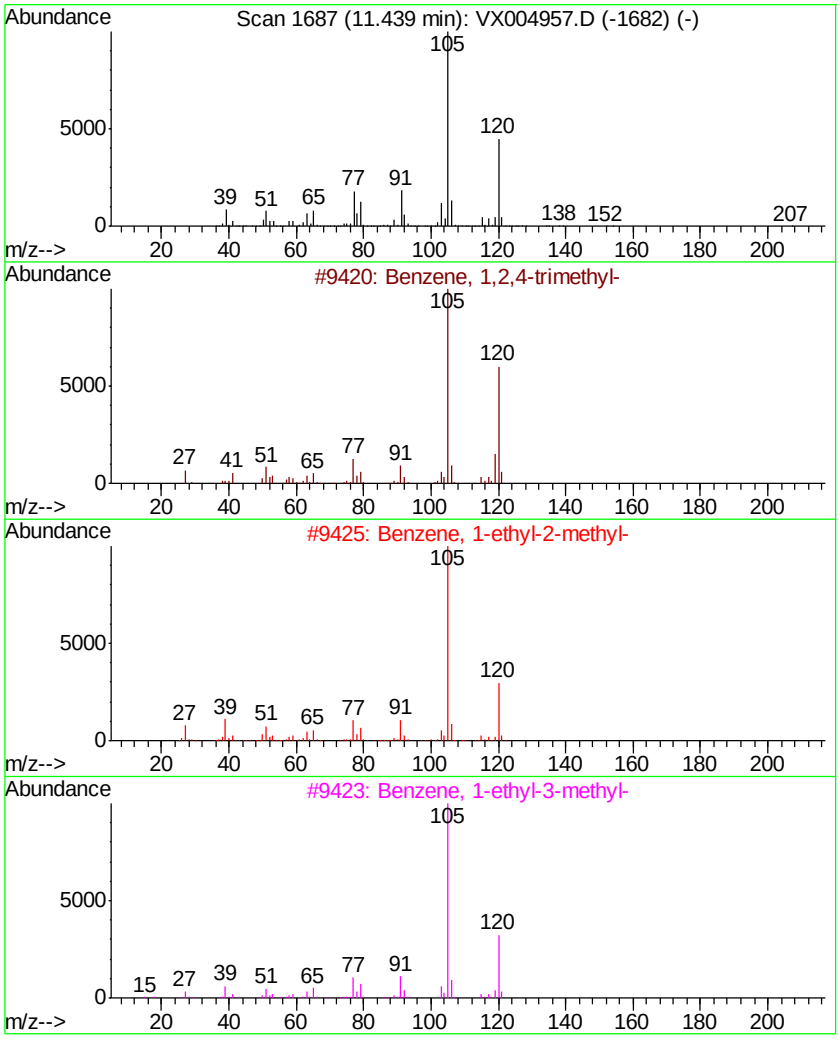
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

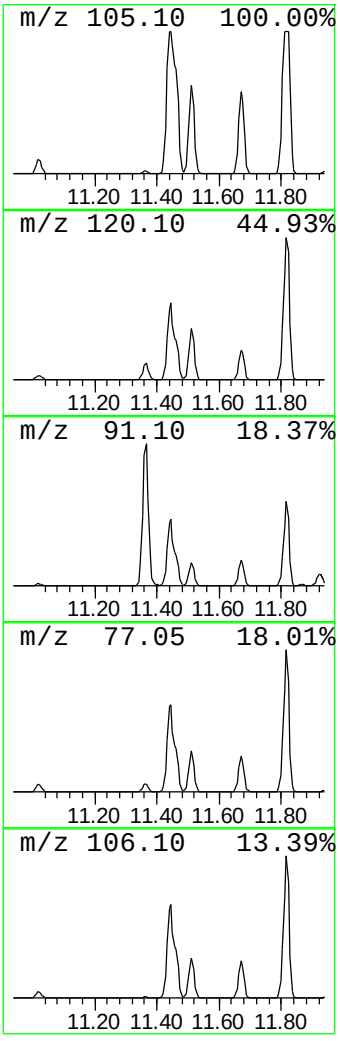
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	1224.77 ug/l	56138600	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87





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Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

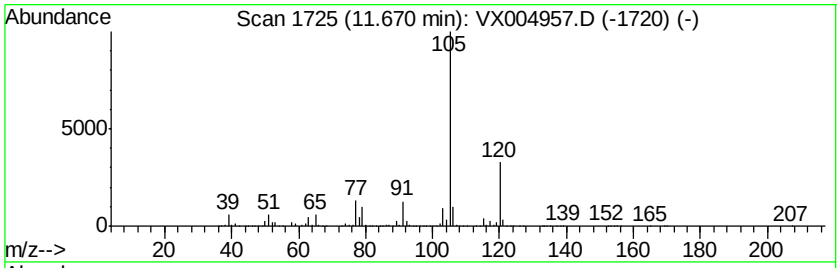
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

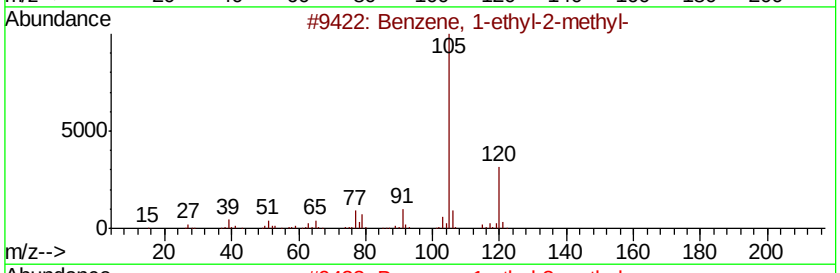
****

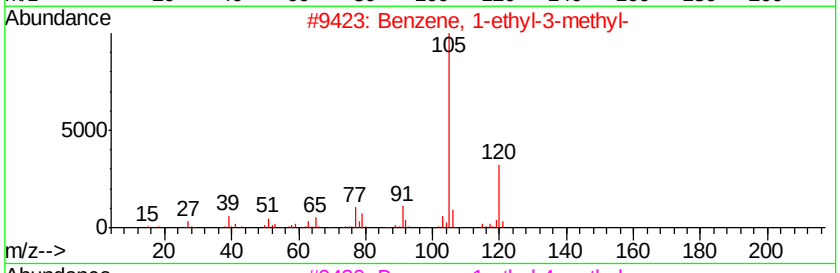
Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

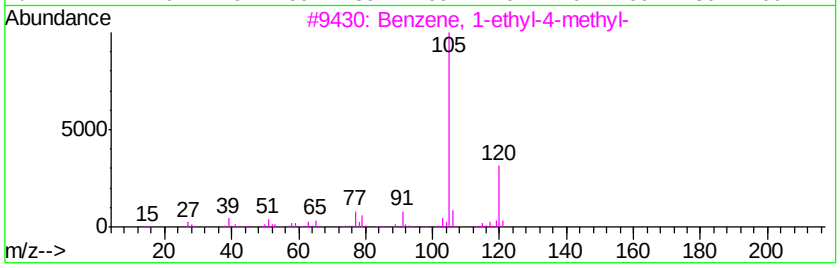
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	342.81 ug/l	15713200	1,4-Dichlorobenzene-d4	12.08

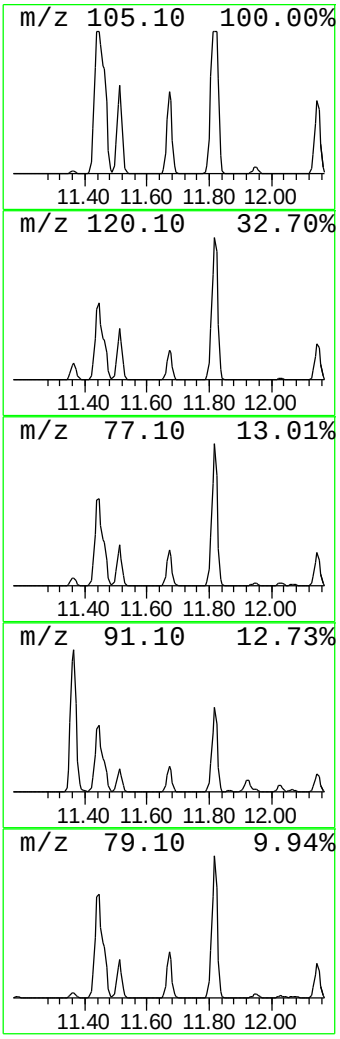
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91











Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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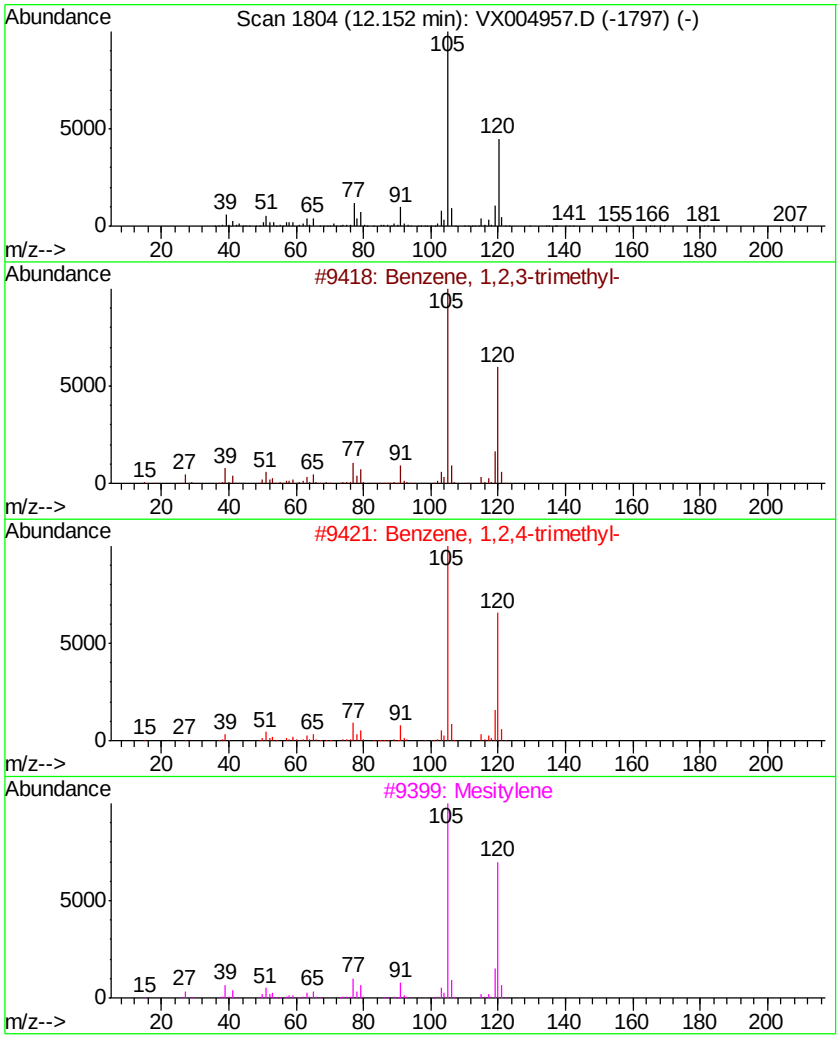
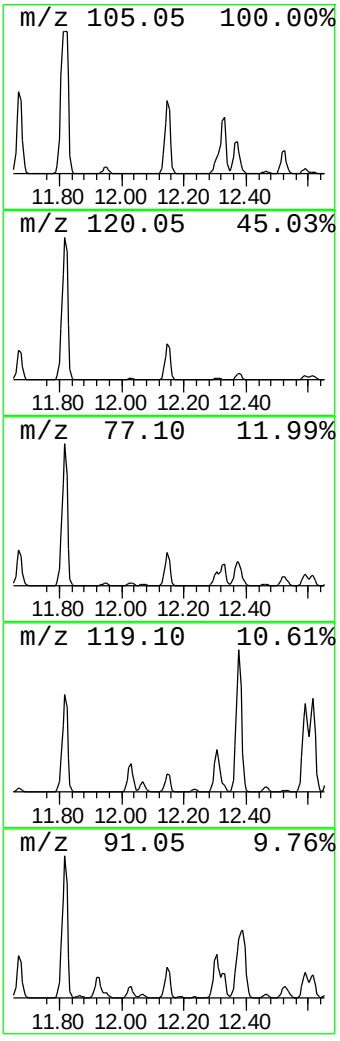
Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	339.27 ug/l	15550800	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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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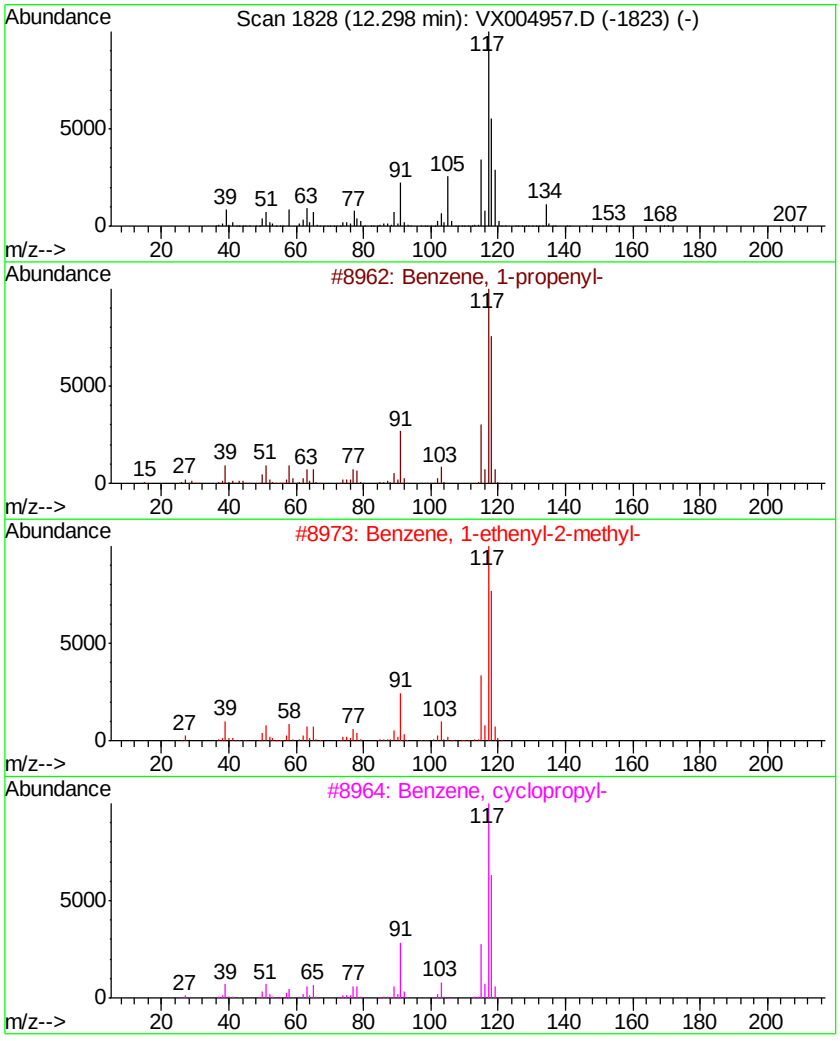
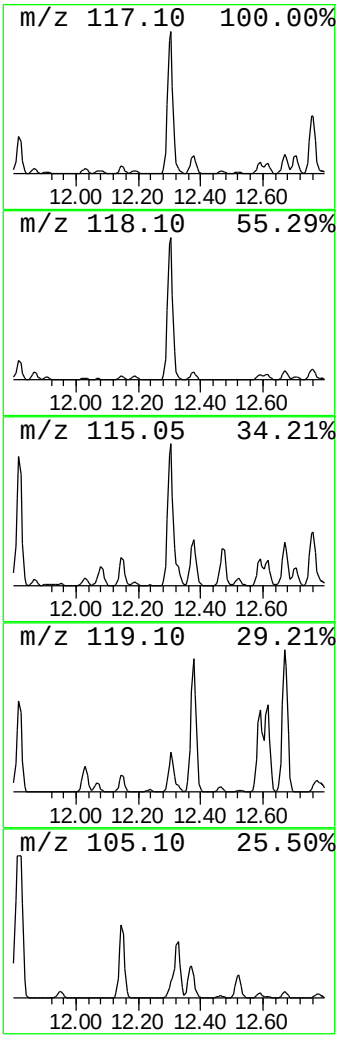
Peak Number 7 Benzene, 1-propenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	340.79 ug/l	15620400	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, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	90
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
4	Indane	118	C9H10	000496-11-7	60
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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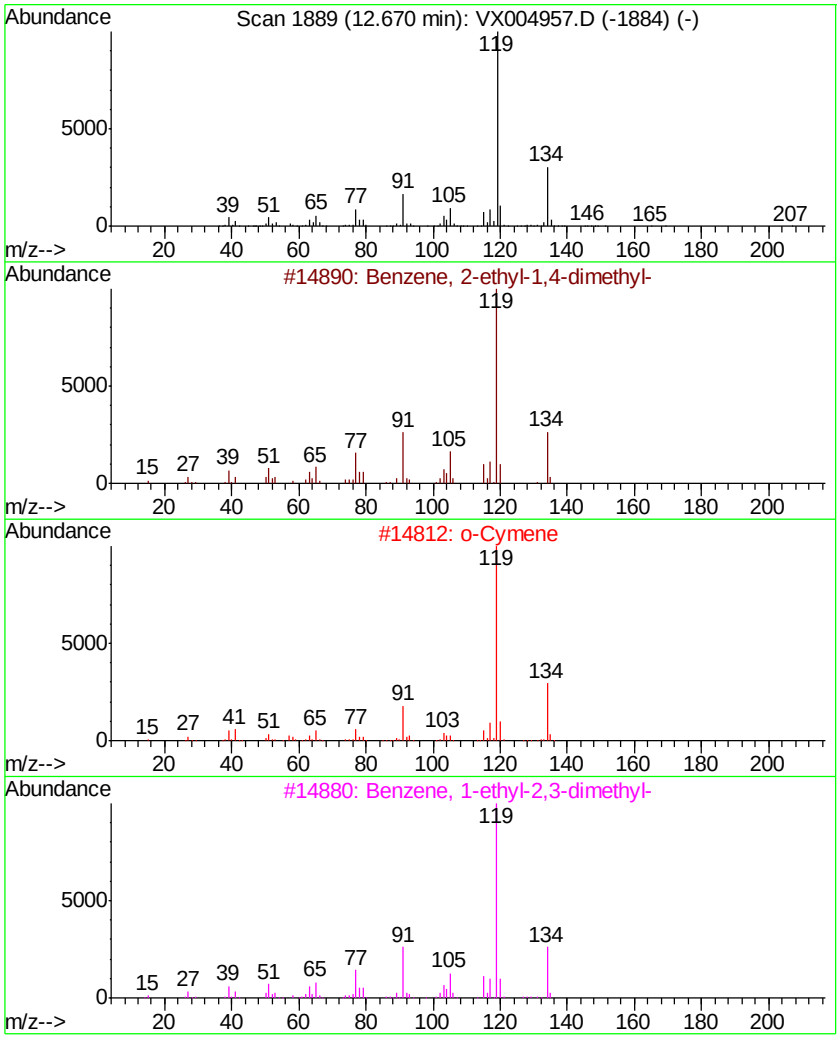
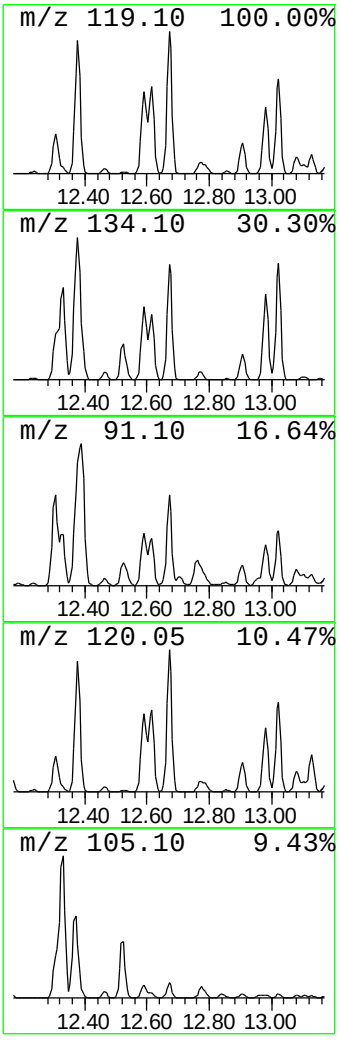
Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	261.18 ug/l	11971400	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	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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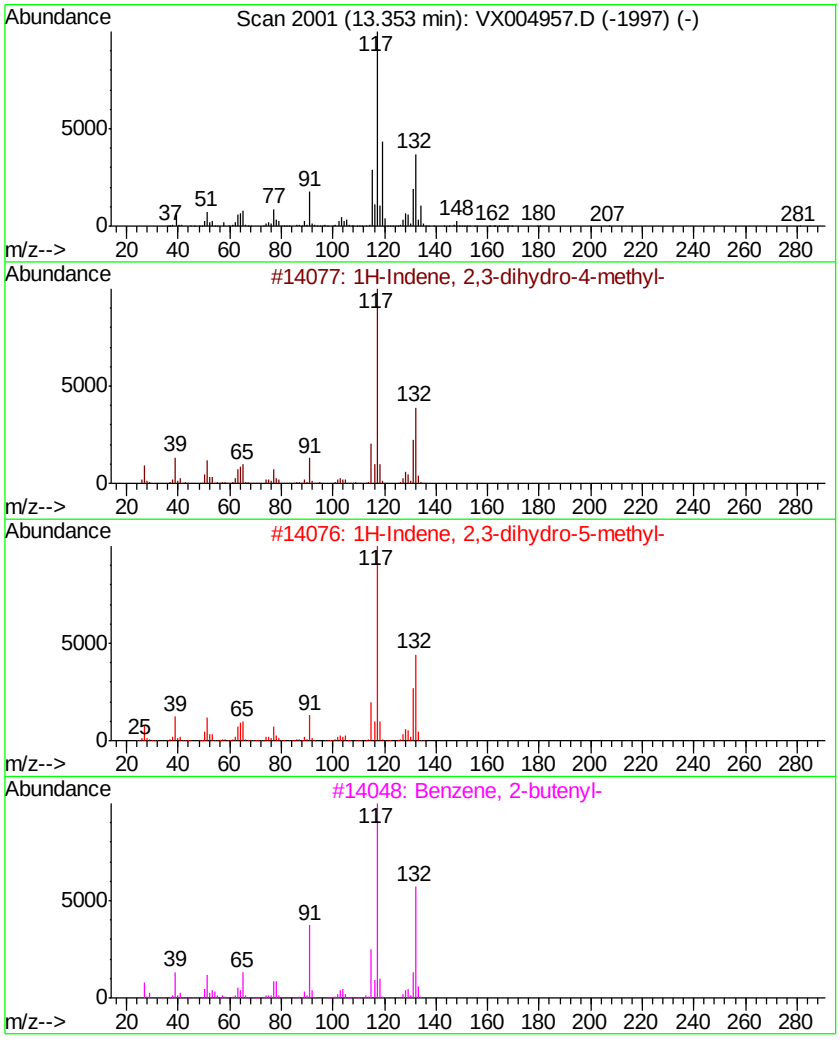
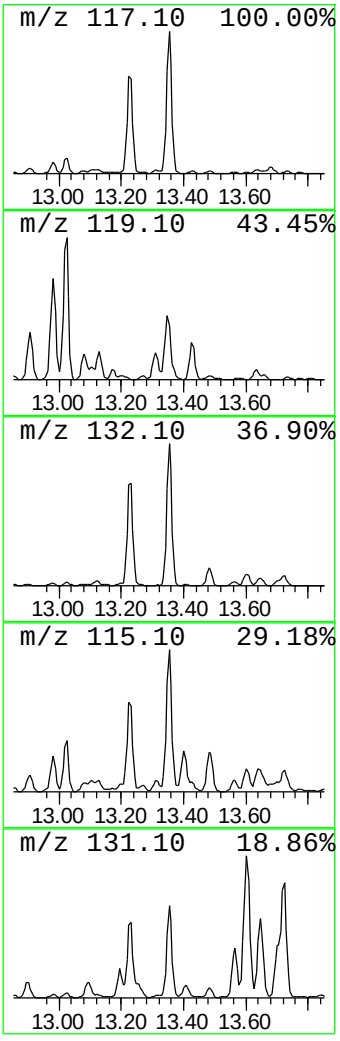
Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	238.32 ug/l	10923600	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	86
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4	Indan, 1-methyl-	132	C10H12	000767-58-8	70
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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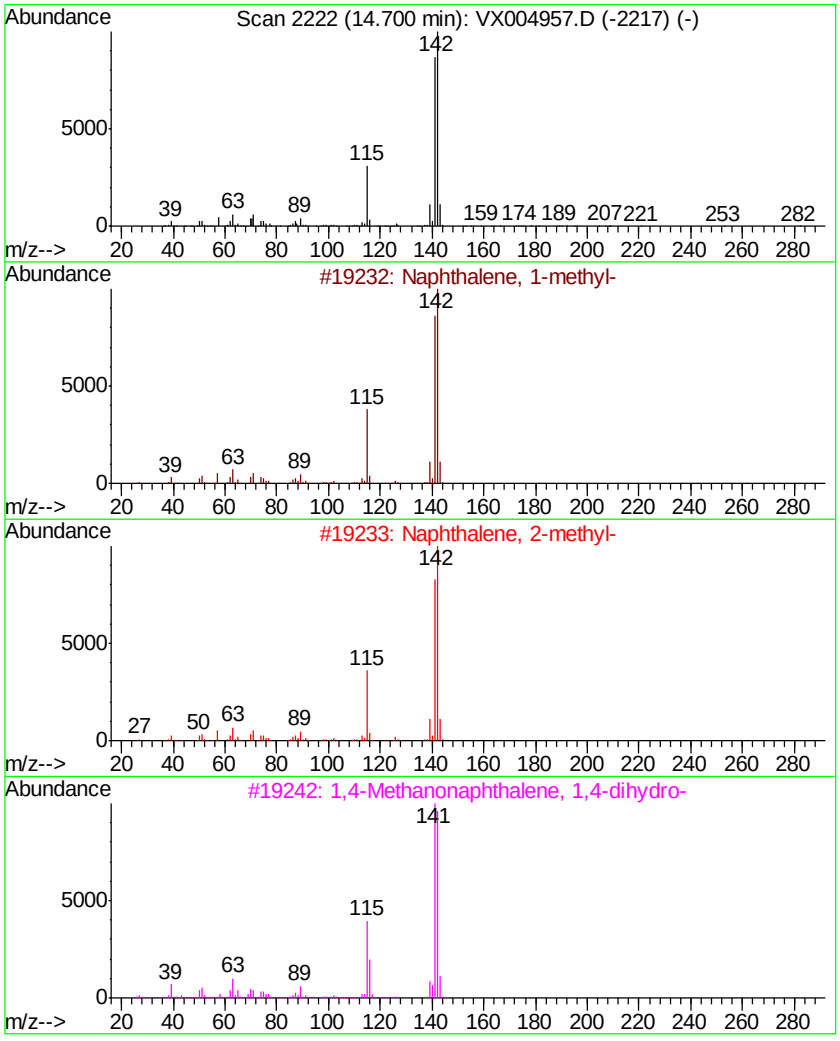
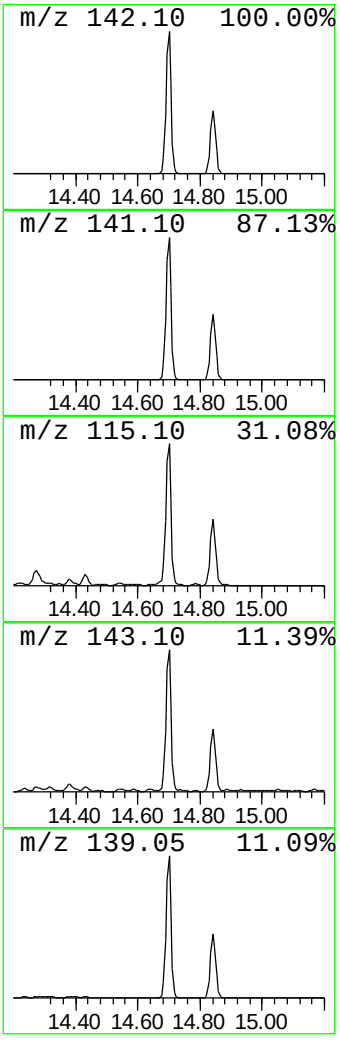
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	210.34 ug/l	9641230	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004957.D
Acq On : 03 Oct 2018 09:46
Operator : JC/MD
Sample : J5084-05 10X
Misc : 5.61g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 39 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : B-7(8-8.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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
Butane, 2,2,3,3-t...	6.35	371.6	ug/l	18781400	2	6.86	2527180	50.0
Pentane, 2,3,4-tr...	8.06	200.4	ug/l	10131200	2	6.86	2527180	50.0
Octane	8.98	204.8	ug/l	5208600	3	10.12	1271670	50.0
Benzene, 1,2,4-tr...	11.44	1224.8	ug/l	56138600	4	12.08	2291810	50.0
Benzene, 1-ethyl-...	11.67	342.8	ug/l	15713200	4	12.08	2291810	50.0
Benzene, 1,2,3-tr...	12.15	339.3	ug/l	15550800	4	12.08	2291810	50.0
Benzene, 1-propenyl-	12.30	340.8	ug/l	15620400	4	12.08	2291810	50.0
Benzene, 2-ethyl-...	12.67	261.2	ug/l	11971400	4	12.08	2291810	50.0
1H-Indene, 2,3-di...	13.35	238.3	ug/l	10923600	4	12.08	2291810	50.0
Naphthalene, 1-me...	14.70	210.3	ug/l	9641230	4	12.08	2291810	50.0