LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

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\82X061919W.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	1.276	16	20	24	rBV	24176	31559	2.63%	0.228%
2	1.324	25	28	34	rVV	5150	13032	1.09%	0.094%
3	1.788	100	104	111	rVB	23254	33540	2.79%	0.242%
4	2.605	230	238	249	rBV	8529	19528	1.63%	0.141%
5	2.842	270	277	288	rBV	51924	116257	9.69%	0.839%
6	3.172	322	331	339	rVB5	7422	19331	1.61%	0.139%
7	4.312	508	518	528	rBV3	11695	37091	3.09%	0.268%
8	5.501	702	713	724	rBV	125872	360781	30.06%	2.603%
9	5.665	729	740	749	rBV	213257	615901	51.32%	4.443%
10	5.952	775	787	796	rBV5	13440	43091	3.59%	0.311%
11	6.061	796	805	826	rVB	123884	324208	27.02%	2.339%
12	6.348	842	852	865	rVB	50859	148164	12.35%	1.069%
13	6.860	925	936	948	rBV	321982	761908	63.49%	5.496%
14	7.451	1024	1033	1044	rVB5	15347	41604	3.47%	0.300%
15	7.677	1064	1070	1076	rVB2	10452	22495	1.87%	0.162%
16	7.848	1093	1098	1106	rVB3	10783	22699	1.89%	0.164%
17	8.055	1121	1132	1140	rBV	61371	127636	10.64%	0.921%
18	8.177	1144	1152	1160	rBV	142623	275572	22.96%	1.988%
19	8.390	1178	1187	1194	rBV3	10826	22040	1.84%	0.159%
20	8.713	1226	1240	1248	rBV	704476	1118671	93.22%	8.070%
21	8.841	1254	1261	1267	rBV	27698	48072	4.01%	0.347%
22	9.036	1288	1293	1300	rVB	19243	34012	2.83%	0.245%
23	9.165	1306	1314	1319	rVB2	14125	23762	1.98%	0.171%
24	9.567	1375	1380	1385	rVB3	10367	16832	1.40%	0.121%
25	9.677	1393	1398	1402	rBV	15260	23183	1.93%	0.167%
26	9.890	1428	1433	1440	rBV2	8516	14186	1.18%	0.102%
27	10.109	1463	1469	1478	rBV2	685818	1200089	100.00%	8.658%
28	10.183	1478	1481	1486	rVV	25938	37519	3.13%	0.271%
29	10.335	1502	1506	1515	rVB4	28436	55091	4.59%	0.397%
30	10.512	1529	1535	1543	rVB	22559	34642	2.89%	0.250%
31	10.658	1543	1559	1564	rBV3	45584	109448	9.12%	0.790%
32	10.713	1564	1568	1575	rVB3	12404	19695	1.64%	0.142%
33	10.811	1575	1584	1585	rBV3	18288	37012	3.08%	0.267%
34	10.835	1585	1588	1596	rVB	50746	65962	5.50%	0.476%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

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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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Title : SW846 8260

35	10.914	1596	1601	1607	rBV4	24546	47894	3.99%	0.346%
36	11.018	1614	1618	1620	rBV	13178	16743	1.40%	0.121%
37	11.067	1620	1626	1631	rVV5	24606	46816	3.90%	0.338%
38	11.134	1632	1637	1642	rVV	580298	743260	61.93%	5.362%
39	11.182	1642	1645	1648	rVV	51605	67018	5.58%	0.483%
40	11.243	1652	1655	1657	rVV2	15174	19835	1.65%	0.143%
41	11.280	1657	1661	1665	rVB	56185	72800	6.07%	0.525%
42	11.384	1675	1678	1683	rBV4	8742	13223	1.10%	0.095%
43	11.445	1683	1688	1693	rVB2	23668	36801	3.07%	0.265%
44	11.573	1706	1709	1715	rVB5	14909	24199	2.02%	0.175%
45	11.670	1717	1725	1731	rBV2	66807	103908	8.66%	0.750%
46	11.768	1731	1741	1745	rBV	93601	133501	11.12%	0.963%
47	11.804	1745	1747	1751	rVB4	8586	12227	1.02%	0.088%
48	11.920	1758	1766	1768	rBV	204782	303287	25.27%	2.188%
49	11.944	1768	1770	1776	rVB	151154	162851	13.57%	1.175%
50	12.012	1776	1781	1786	rVB9	8532	20772	1.73%	0.150%
51	12.079	1786	1792	1797	rBV	717194	938237	78.18%	6.769%
52	12.170	1803	1807	1812	rBV2	15790	26261	2.19%	0.189%
53	12.231	1812	1817	1823	rVB	316065	387960	32.33%	2.799%
54	12.298	1823	1828	1834	rBV	177445	219642	18.30%	1.585%
55	12.365	1834	1839	1848	rVB2	91298	146060	12.17%	1.054%
56	12.469	1849	1856	1860	rBV4	12798	25044	2.09%	0.181%
57	12.664	1879	1888	1892	rVV	438310	591651	49.30%	4.268%
58	12.700	1892	1894	1899	rVV2	168668	270524	22.54%	1.952%
59	12.755	1899	1903	1909	rVV3	428705	725826	60.48%	5.236%
60	12.816	1909	1913	1917	rVB	71545	95066	7.92%	0.686%
61	12.896	1922	1926	1931	rVV	304244	389081	32.42%	2.807%
62	12.975	1931	1939	1944	rVV2	302725	472209	39.35%	3.407%
63	13.024	1944	1947	1952	rVV	26666	39471	3.29%	0.285%
64	13.085	1952	1957	1963	rVB2	83947	126936	10.58%	0.916%
65	13.145	1963	1967	1971	rBV	55676	79211	6.60%	0.571%
66	13.188	1971	1974	1978	rVV	85726	109441	9.12%	0.790%
67	13.225	1978	1980	1981	rVV	20437	20578	1.71%	0.148%
68	13.249	1981	1984	1988	rVB	71509	84622	7.05%	0.610%
69	13.304	1988	1993	1996	rBV2	31361	56230	4.69%	0.406%
70	13.347	1996	2000	2005	rVV2	263680	372797	31.06%	2.689%
71	13.402	2005	2009	2011	rVV3	40937	60315	5.03%	0.435%

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

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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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Title : SW846 8260

72	13.456	2015	2018	2027	rVB6	21960	51639	4.30%	0.373%
73	13.560	2032	2035	2037	rVV	16879	19081	1.59%	0.138%
74	13.597	2037	2041	2044	rVV2	40624	62124	5.18%	0.448%
75	13.639	2044	2048	2051	rVV2	97650	132931	11.08%	0.959%
76	13.676	2051	2054	2055	rVV2	41359	50704	4.23%	0.366%
77	13.694	2055	2057	2059	rVV	63285	81741	6.81%	0.590%
78	13.719	2059	2061	2066	rVV2	75204	92007	7.67%	0.664%
79	13.804	2070	2075	2078	rBV	28993	39152	3.26%	0.282%
80	13.853	2080	2083	2088	rVB	18093	24361	2.03%	0.176%
81	13.908	2088	2092	2097	rBV	26226	35345	2.95%	0.255%
82	13.981	2097	2104	2108	rBV2	34578	53854	4.49%	0.389%
83	14.023	2108	2111	2116	rBV3	14803	20337	1.69%	0.147%
84	14.127	2124	2128	2131	rBV3	13063	19171	1.60%	0.138%
85	14.158	2131	2133	2138	rVB2	14658	16146	1.35%	0.116%
86	14.267	2147	2151	2157	rBV2	23908	43889	3.66%	0.317%
87	14.377	2165	2169	2173	rVB2	9863	13062	1.09%	0.094%
88	14.432	2173	2178	2181	rBV4	10160	14238	1.19%	0.103%
89	14.535	2190	2195	2200	rBV3	32606	47810	3.98%	0.345%
90	14.731	2223	2227	2231	rVB5	9920	13352	1.11%	0.096%
91	14.834	2240	2244	2254	rBV2	45472	72133	6.01%	0.520%
92	15.688	2375	2384	2387	rBV3	11230	21716	1.81%	0.157%

Sum of corrected areas: 13861703

LSC Report - Integrated Chromatogram

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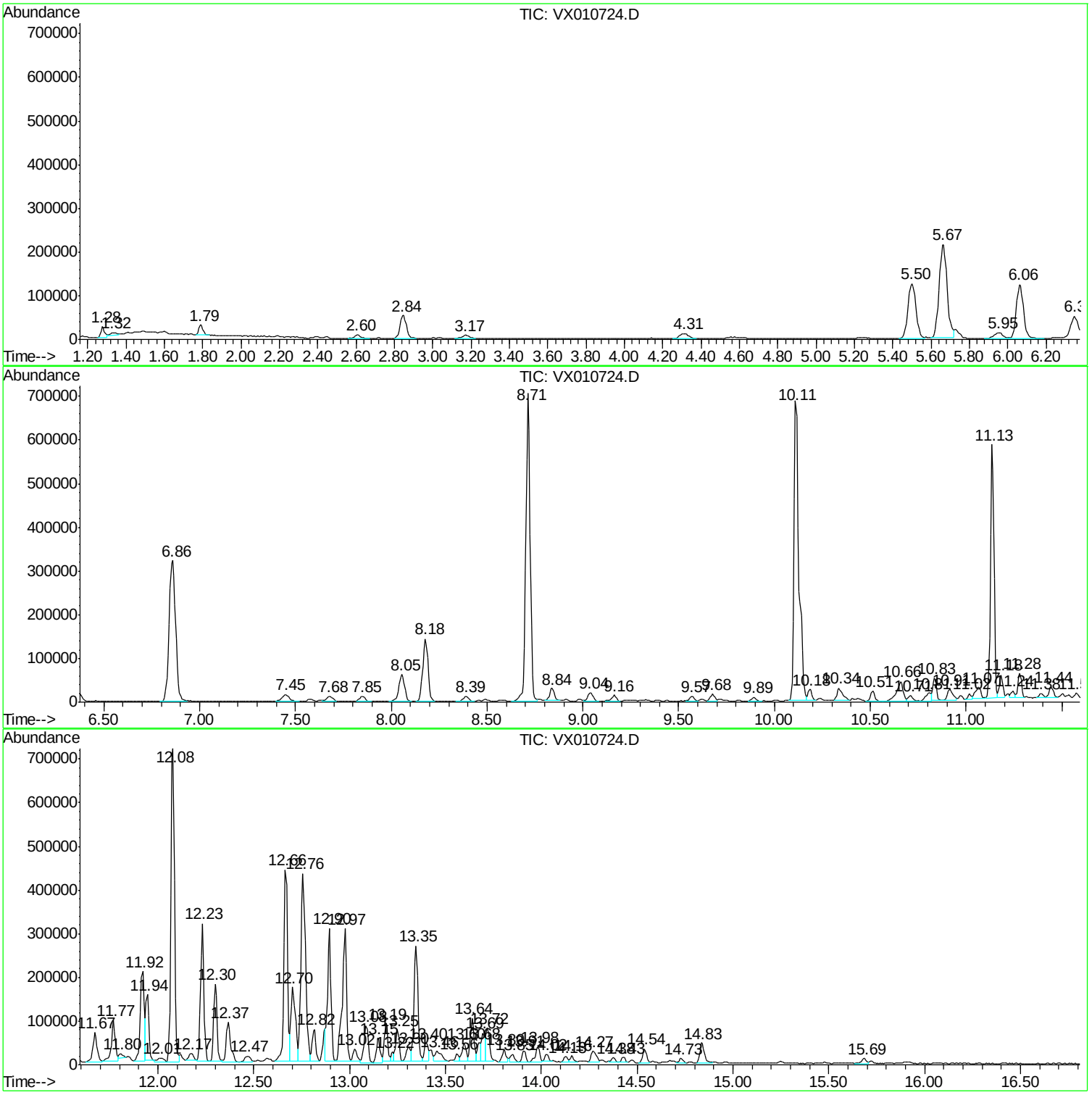
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
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Quant Title : SW846 8260

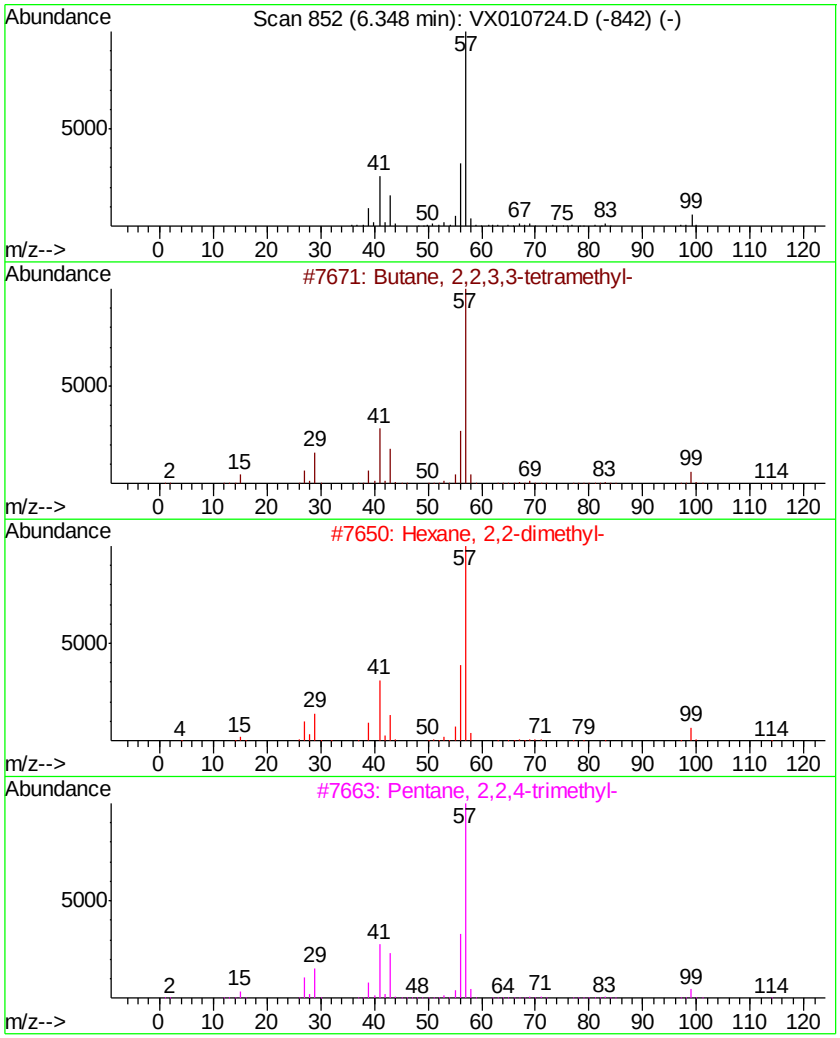
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

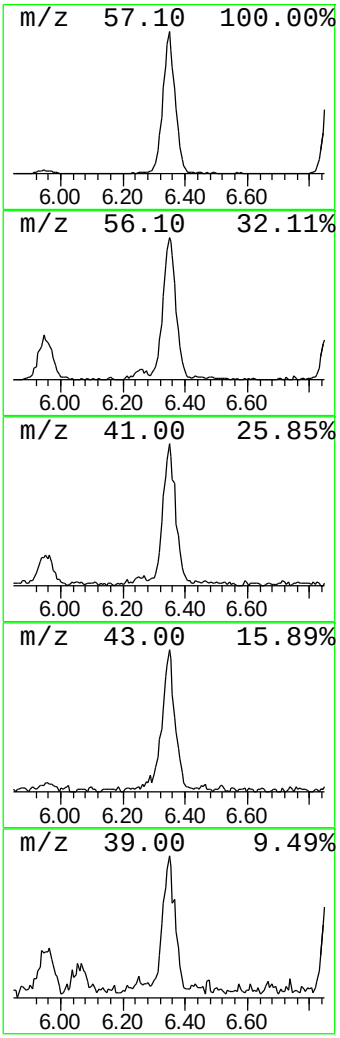
****

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	9.72 ug/l	148164	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	72
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

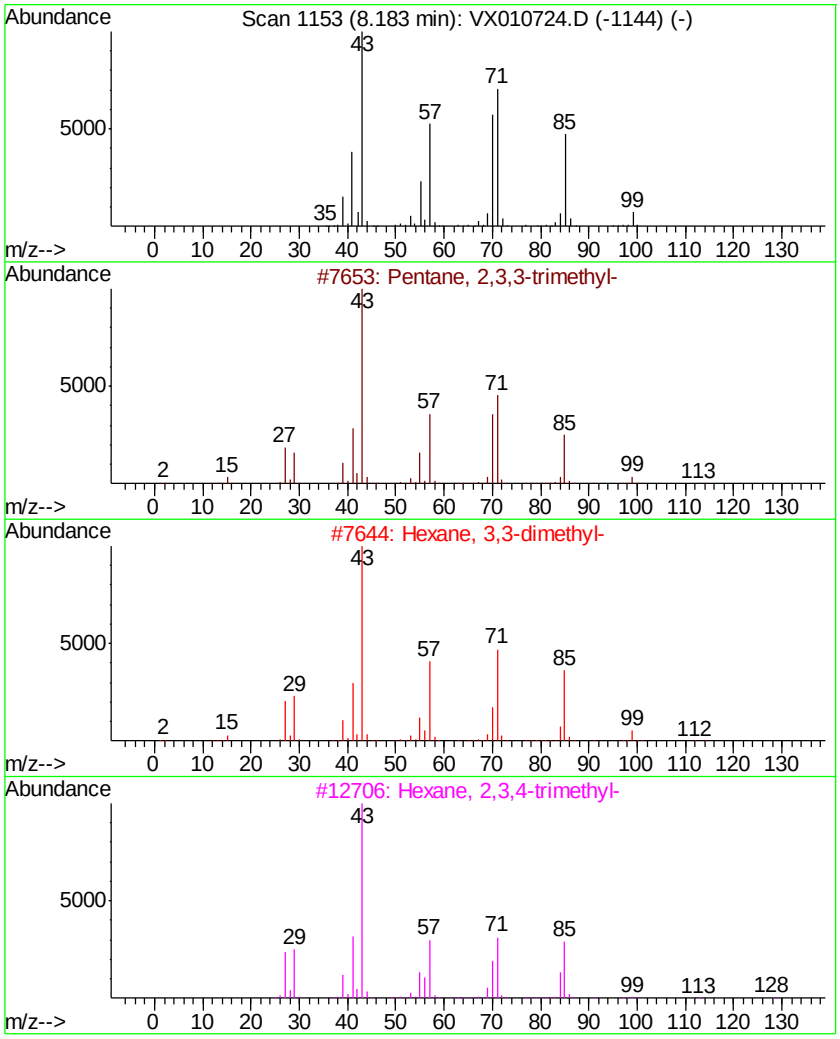
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

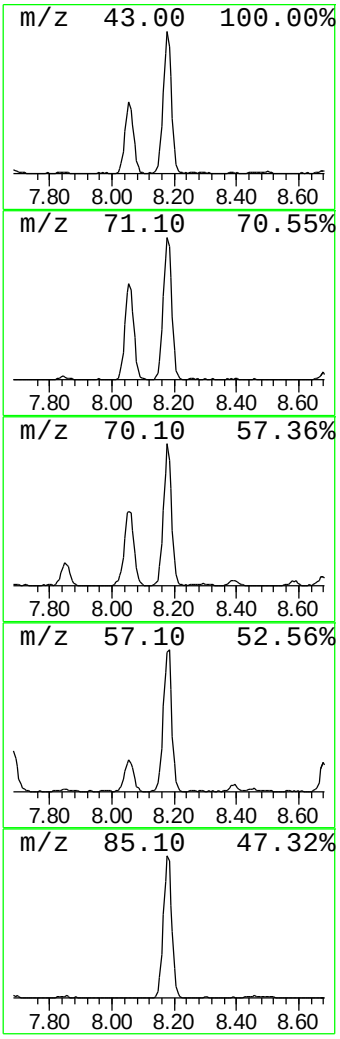
****

Peak Number 2 Pentane, 2,3,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	18.08 ug/l	275572	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5	Propanoic acid, 2-methyl-, 2-eth...	200	C12H24O2	035061-61-1	50





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

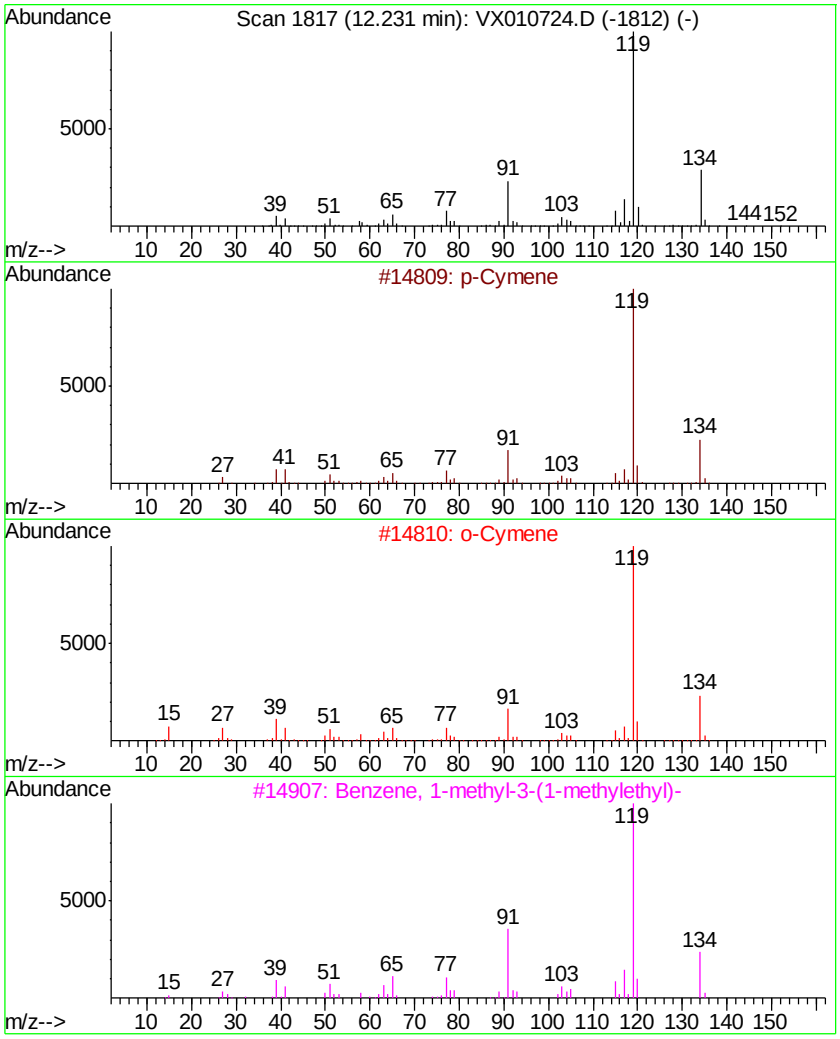
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

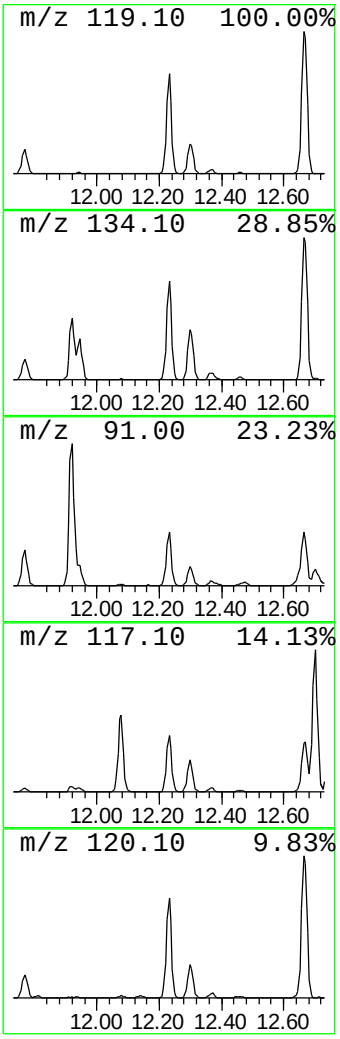
****

Peak Number 3 p-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	20.67 ug/l	387960	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91





Library Search Compound Report

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

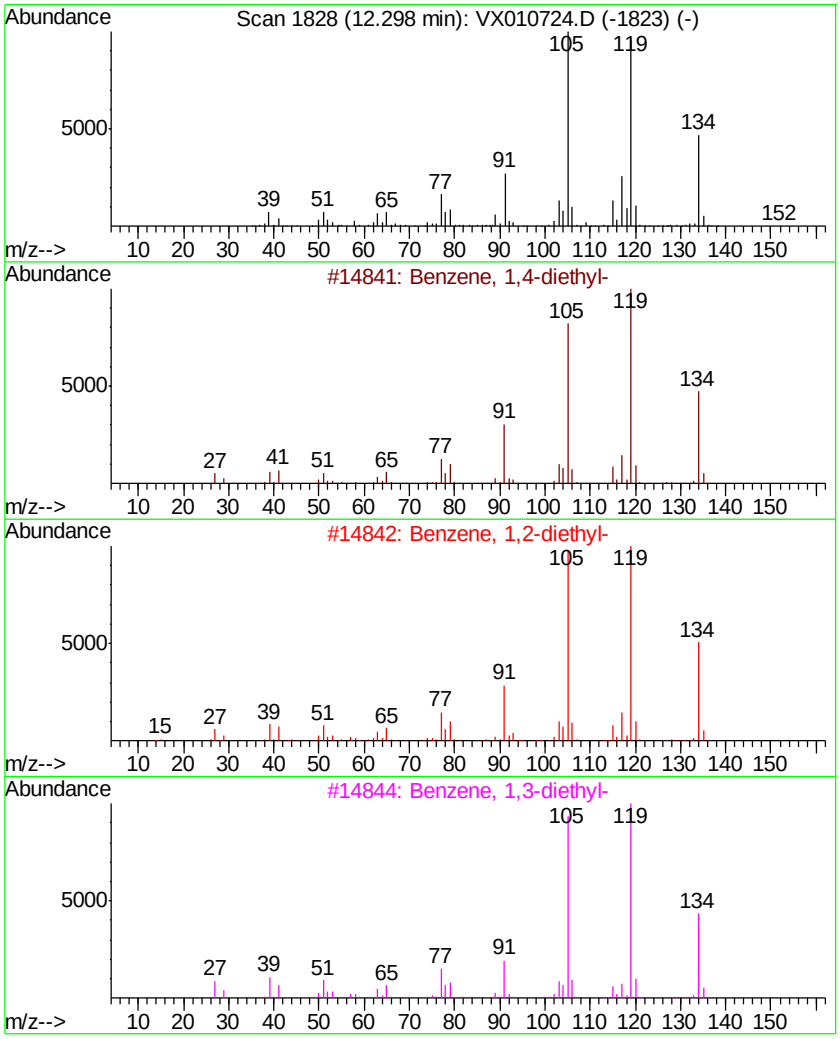
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

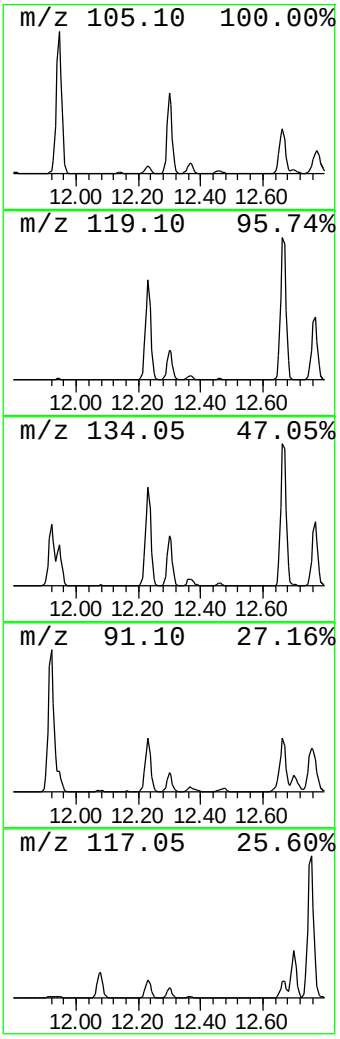
****

Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	11.71 ug/l	219642	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	68





Library Search Compound Report

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Acq On : 09 Jul 2019 03:06
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Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

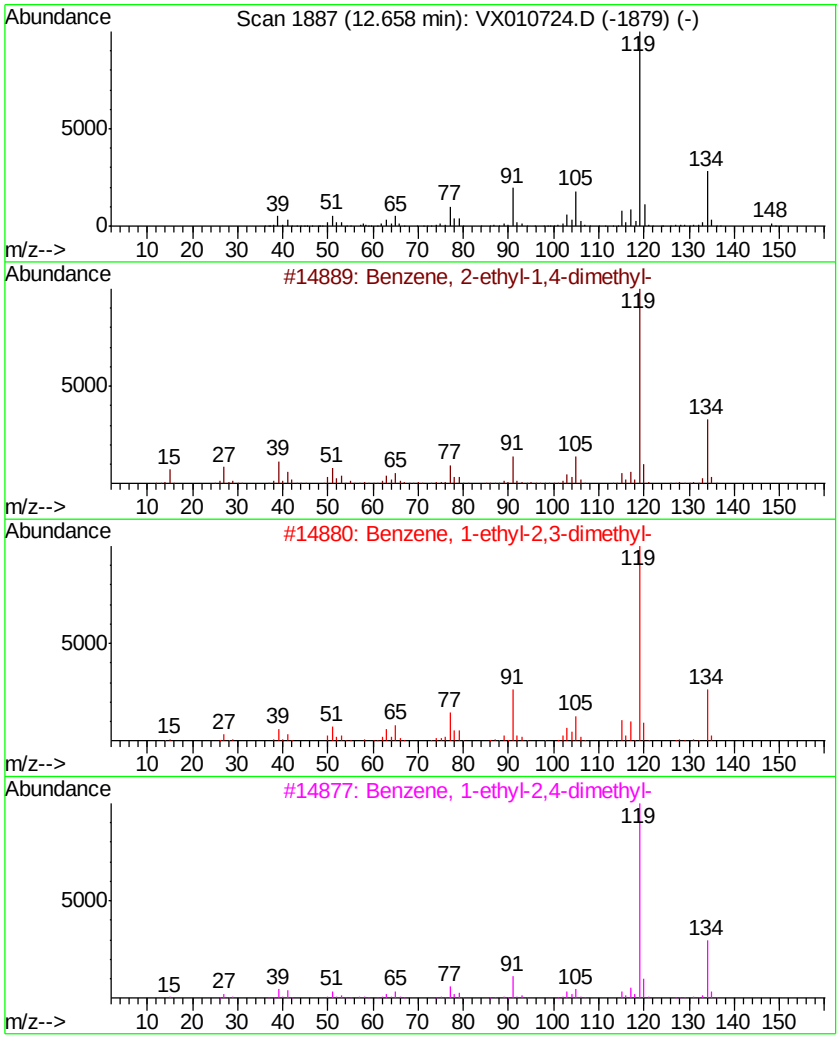
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

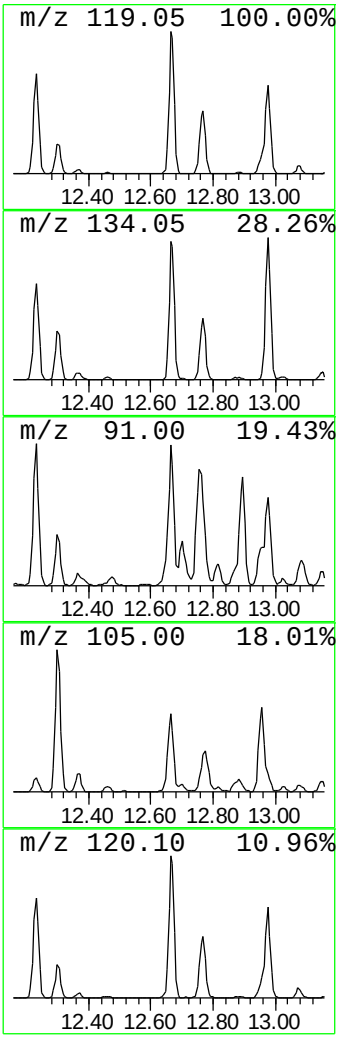
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	31.53 ug/l	591651	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	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





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Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

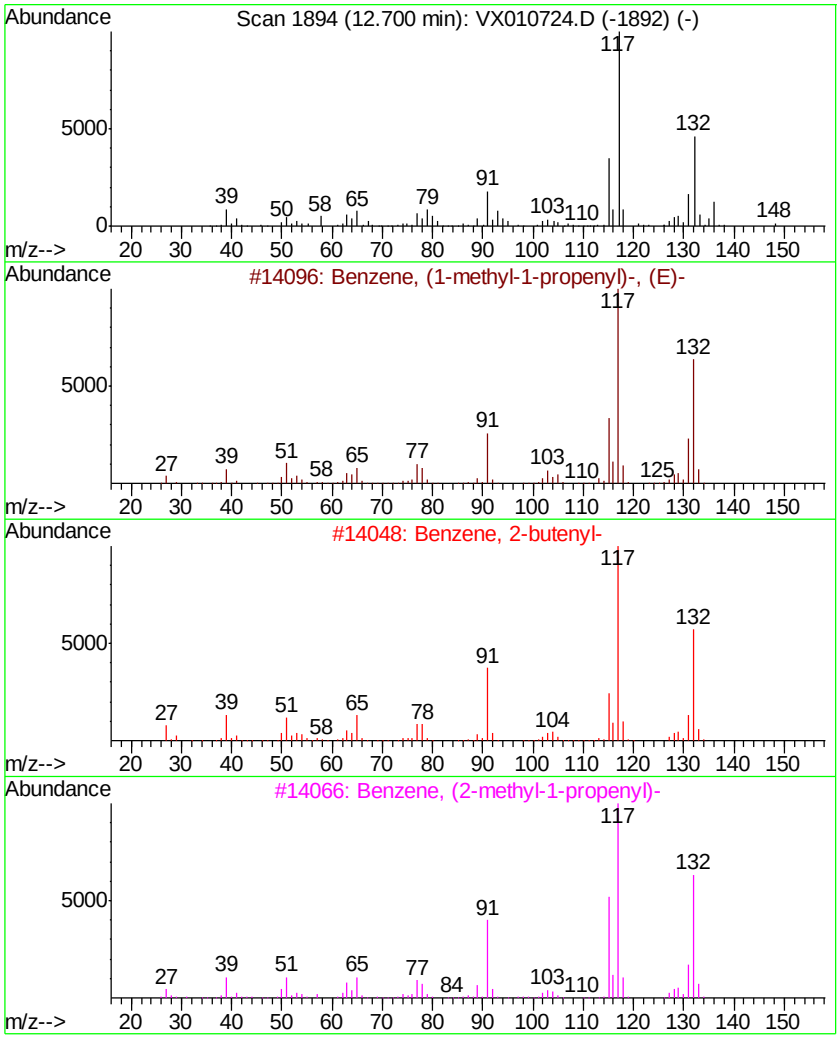
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

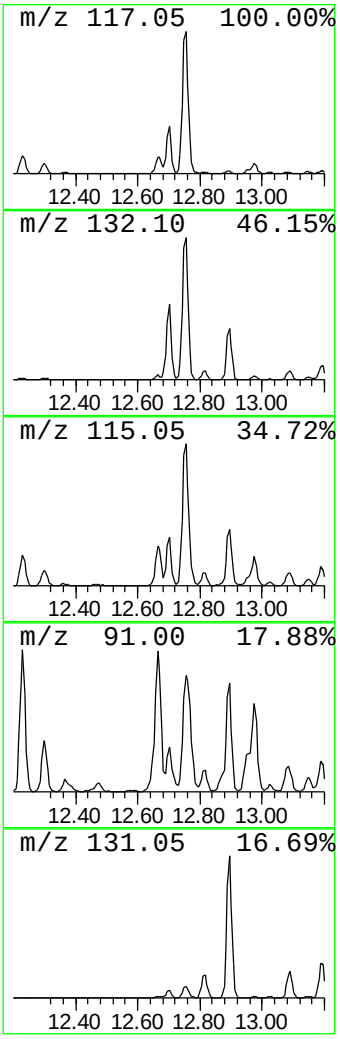
****

Peak Number 6 Benzene, (1-methyl-1-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	14.42 ug/l	270524	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	93
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	89
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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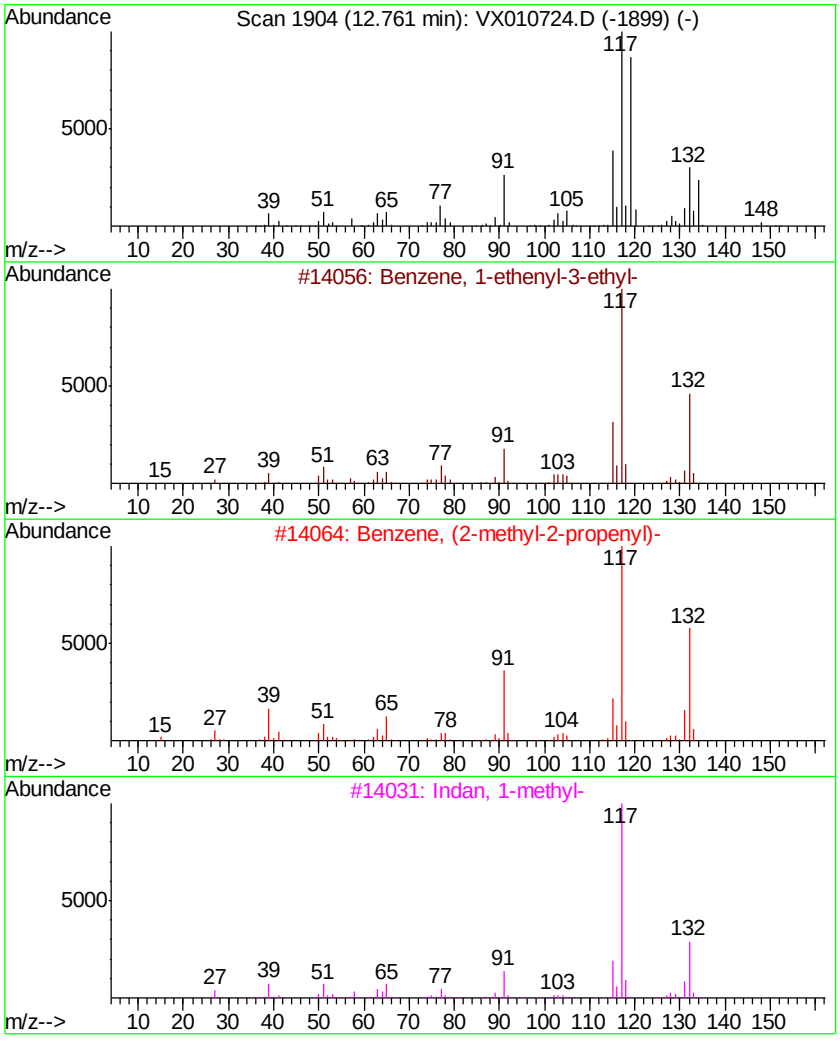
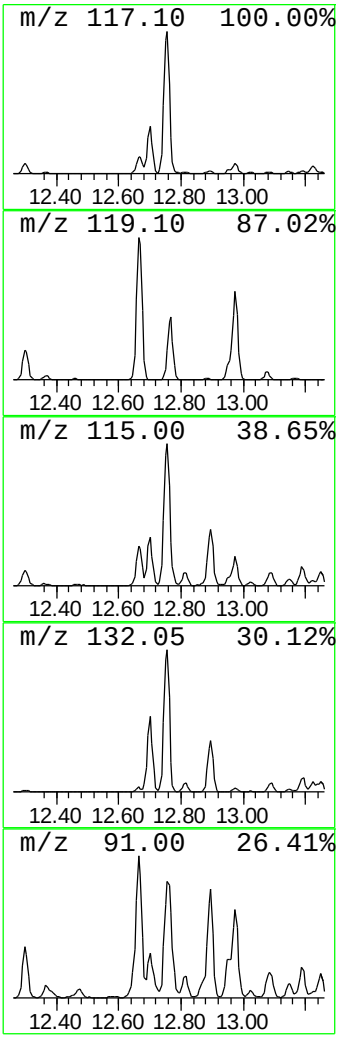
Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	38.68 ug/l	725826	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
2	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
3	Indan, 1-methyl-	132	C10H12	000767-58-8	76
4	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	74





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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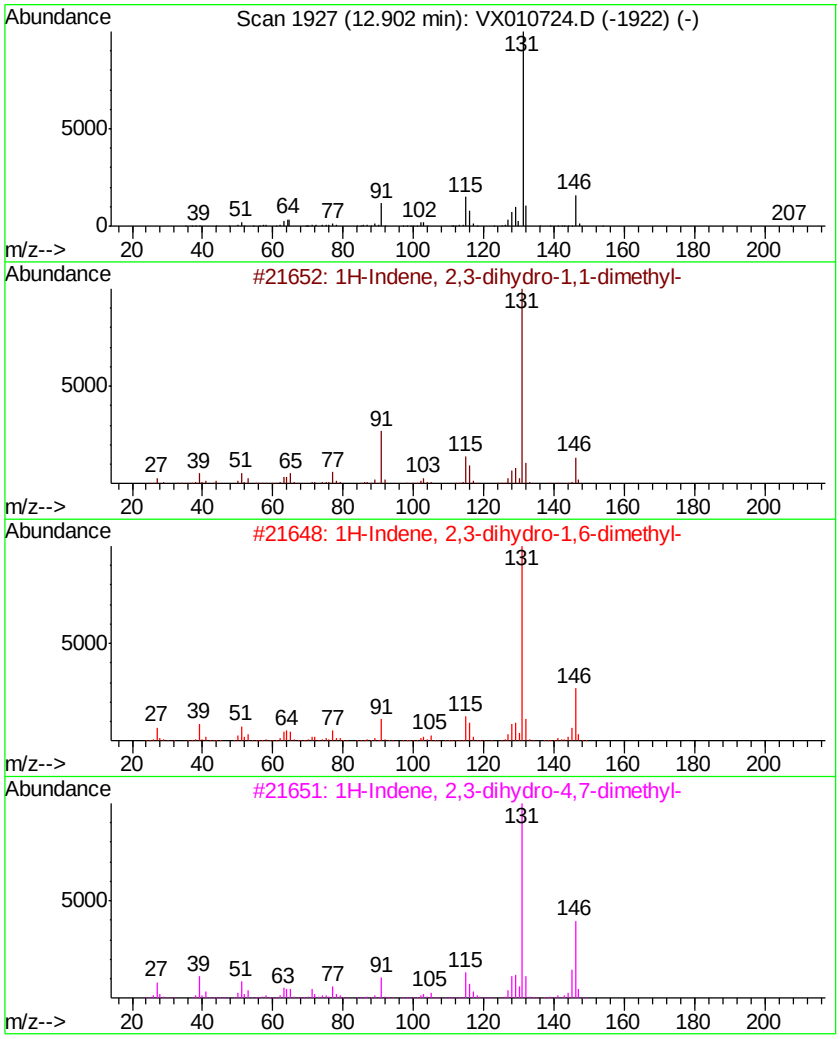
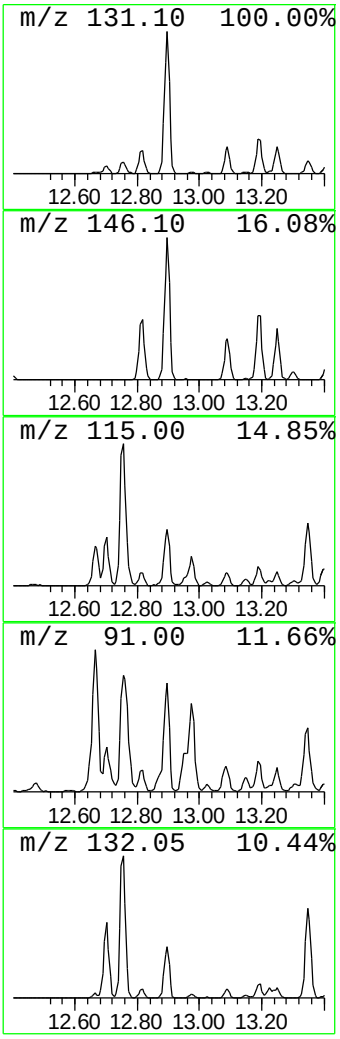
Peak Number 8 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	20.73 ug/l	389081	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	93
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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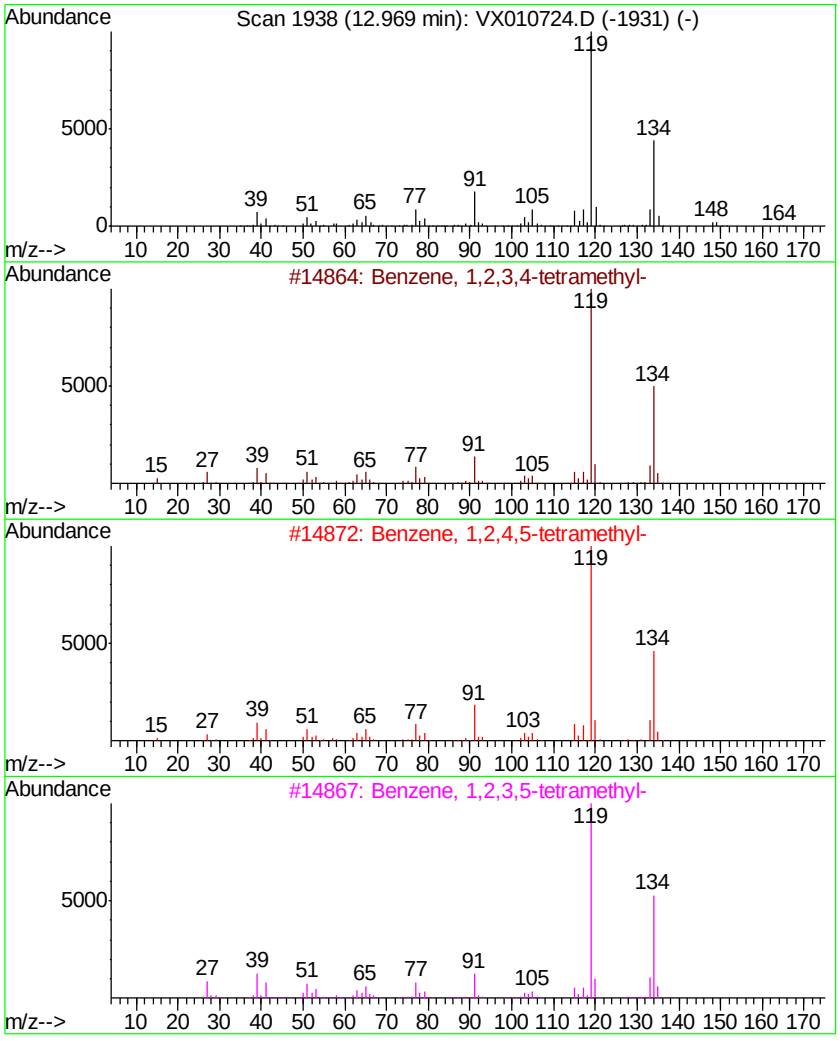
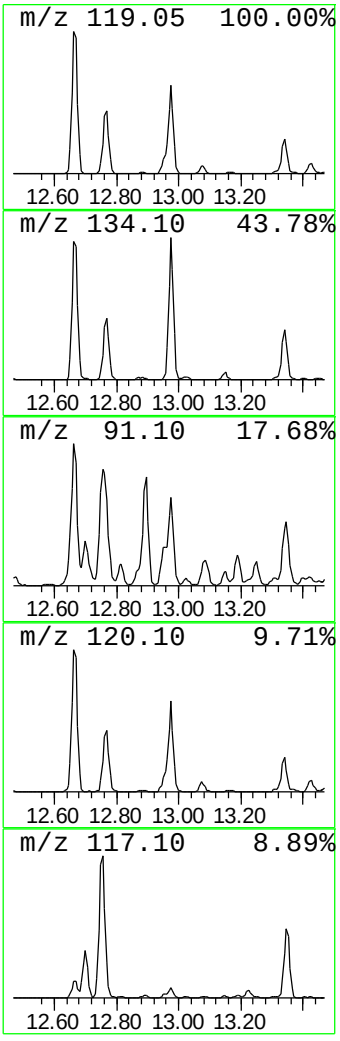
Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	25.16 ug/l	472209	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
5	o-Cymene	134	C10H14	000527-84-4	93





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
Acq On : 09 Jul 2019 03:06
Operator : JC/SP
Sample : K3690-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW016-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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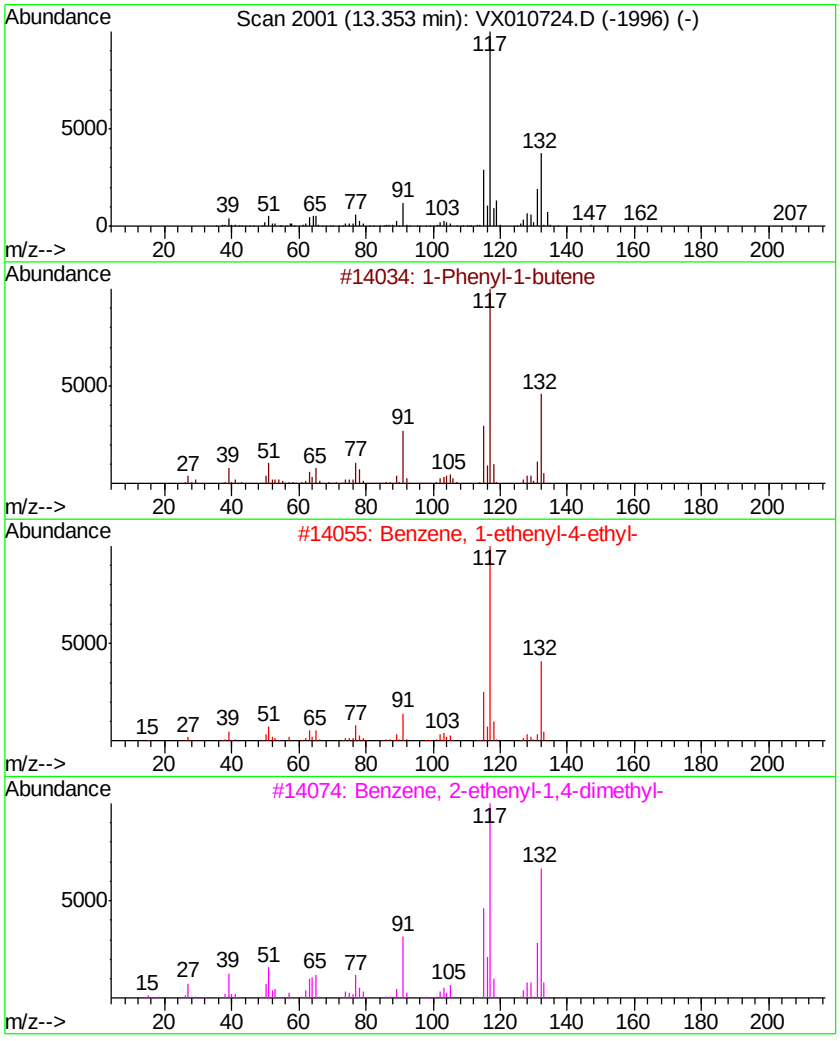
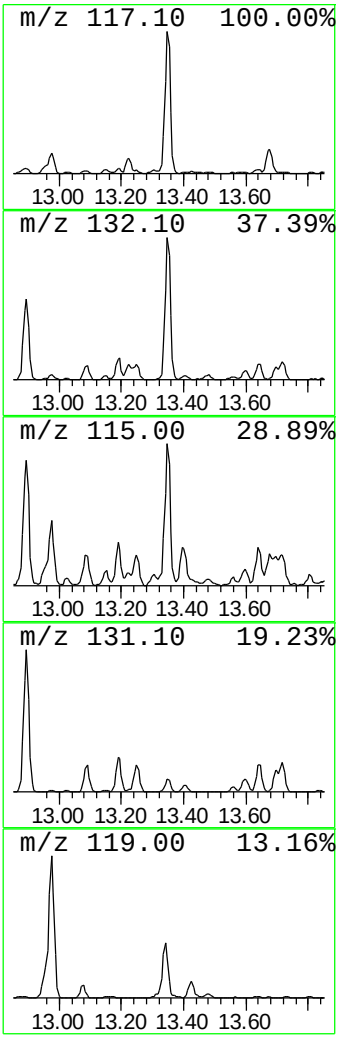
Peak Number 10 1-Phenyl-1-butene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	19.87 ug/l	372797	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	70
4	Indan, 1-methyl-	132	C10H12	000767-58-8	64
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	62





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010724.D
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ALS Vial : 43 Sample Multiplier: 1

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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	9.7	ug/l	148164	2	6.86	761908	50.0
Pentane, 2,3,3-tr...	8.18	18.1	ug/l	275572	2	6.86	761908	50.0
p-Cymene	12.23	20.7	ug/l	387960	4	12.08	938237	50.0
Benzene, 1,4-diet...	12.30	11.7	ug/l	219642	4	12.08	938237	50.0
Benzene, 2-ethyl-...	12.66	31.5	ug/l	591651	4	12.08	938237	50.0
Benzene, (1-methy...	12.70	14.4	ug/l	270524	4	12.08	938237	50.0
Benzene, 1-etheny...	12.76	38.7	ug/l	725826	4	12.08	938237	50.0
1H-Indene, 2,3-di...	12.90	20.7	ug/l	389081	4	12.08	938237	50.0
Benzene, 1,2,3,4-...	12.97	25.2	ug/l	472209	4	12.08	938237	50.0
1-Phenyl-1-butene	13.35	19.9	ug/l	372797	4	12.08	938237	50.0