LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\METHOD\SOMULM032319WMA.M
Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.181	25	28	38	rVB	16437	16591	0.19%	0.044%
2	1.396	88	95	105	rBV	286457	294763	3.42%	0.785%
3	1.663	170	178	189	rBV	202754	254666	2.95%	0.678%
4	2.129	317	323	331	rVV3	9446	11800	0.14%	0.031%
5	2.180	334	339	347	rVB3	12647	15993	0.19%	0.043%
6	2.271	355	367	380	rBV3	718170	1313487	15.22%	3.499%
7	2.621	473	476	479	rBV	2365	1796	0.02%	0.005%
8	2.676	483	493	502	rBV3	20765	41091	0.48%	0.109%
9	2.785	519	527	542	rVB5	17232	36075	0.42%	0.096%
10	2.991	578	591	604	rBV5	11439	28834	0.33%	0.077%
11	3.184	647	651	652	rBV	2526	1648	0.02%	0.004%
12	3.274	675	679	680	rBV	2764	1918	0.02%	0.005%
13	3.299	681	687	699	rVB4	6175	12311	0.14%	0.033%
14	3.489	741	746	757	rVV4	5099	8343	0.10%	0.022%
15	3.643	790	794	802	rVB3	2292	2988	0.03%	0.008%
16	3.724	814	819	823	rBV2	5581	6877	0.08%	0.018%
17	3.920	875	880	884	rVB	1791	1596	0.02%	0.004%
18	4.087	917	932	947	rBV5	29740	80600	0.93%	0.215%
19	4.174	947	959	994	rBV	212458	556798	6.45%	1.483%
20	4.425	1033	1037	1040	rBV2	1768	1718	0.02%	0.005%
21	4.640	1089	1104	1124	rBV	304676	720664	8.35%	1.920%
22	4.846	1163	1168	1171	rBV	1417	1576	0.02%	0.004%
23	4.872	1172	1176	1180	rVV2	2305	2256	0.03%	0.006%
24	4.897	1181	1184	1190	rVB	3343	3564	0.04%	0.009%
25	4.991	1200	1213	1230	rVV4	31960	75359	0.87%	0.201%
26	5.219	1274	1284	1295	rBV6	4301	10630	0.12%	0.028%
27	5.335	1298	1320	1326	rBV2	658715	1807753	20.95%	4.816%
28	5.383	1328	1335	1354	rVB	1593392	3222433	37.35%	8.584%
29	5.701	1432	1434	1439	rVB	2433	1650	0.02%	0.004%
30	5.881	1477	1490	1517	rBV	549204	1117674	12.95%	2.977%
31	6.116	1560	1563	1569	rVB	3609	2964	0.03%	0.008%
32	6.180	1569	1583	1612	rBV	424740	852713	9.88%	2.271%
33	6.325	1615	1628	1645	rBV	433359	876505	10.16%	2.335%
34	6.817	1774	1781	1782	rBV3	4384	4562	0.05%	0.012%

LSC Area Percent Report

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Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\METHOD\SOMULM032319WMA.M
Title : VOC Analysis

35	6.891	1800	1804	1808	rBV	2311	1860	0.02%	0.005%
36	6.936	1814	1818	1825	rVB2	1628	2069	0.02%	0.006%
37	7.061	1847	1857	1868	rVV5	10898	21851	0.25%	0.058%
38	7.113	1868	1873	1879	rVV4	2015	2738	0.03%	0.007%
39	7.174	1886	1892	1894	rBV2	1989	2262	0.03%	0.006%
40	7.225	1894	1908	1924	rBV	236831	425246	4.93%	1.133%
41	7.370	1944	1953	1964	rBV3	20892	36710	0.43%	0.098%
42	7.560	1992	2012	2023	rBV	822688	1444702	16.74%	3.848%
43	7.634	2023	2035	2057	rVB	5007933	8627679	100.00%	22.982%
44	7.846	2091	2101	2116	rVB	165736	279268	3.24%	0.744%
45	8.097	2171	2179	2187	rVB5	4190	6632	0.08%	0.018%
46	8.193	2204	2209	2215	rVB2	2190	2636	0.03%	0.007%
47	8.254	2222	2228	2234	rVB4	2299	3340	0.04%	0.009%
48	8.309	2234	2245	2280	rBV	706181	1278504	14.82%	3.406%
49	8.502	2300	2305	2312	rVB4	7785	9937	0.12%	0.026%
50	8.585	2322	2331	2341	rVB3	16117	27518	0.32%	0.073%
51	8.637	2344	2347	2355	rVB5	3318	3715	0.04%	0.010%
52	8.830	2403	2407	2413	rVB	1556	1645	0.02%	0.004%
53	8.997	2454	2459	2460	rBV	2508	1962	0.02%	0.005%
54	9.029	2462	2469	2475	rBV2	8981	12467	0.14%	0.033%
55	9.087	2475	2487	2491	rBV	848931	1373041	15.91%	3.658%
56	9.245	2526	2536	2561	rVB	571255	926227	10.74%	2.467%
57	9.370	2562	2575	2611	rVB	1903199	3131979	36.30%	8.343%
58	9.775	2689	2701	2732	rVB	974610	1596519	18.50%	4.253%
59	9.923	2743	2747	2749	rBV	2440	1598	0.02%	0.004%
60	9.939	2749	2752	2755	rBV	2597	2187	0.03%	0.006%
61	10.007	2770	2773	2781	rVB4	2228	2430	0.03%	0.006%
62	10.058	2784	2789	2792	rVV2	2039	2203	0.03%	0.006%
63	10.080	2792	2796	2802	rVB4	3752	4303	0.05%	0.011%
64	10.132	2808	2812	2814	rVV	2898	2444	0.03%	0.007%
65	10.164	2814	2822	2834	rVB3	23990	40400	0.47%	0.108%
66	10.283	2856	2859	2864	rBV	2061	2194	0.03%	0.006%
67	10.354	2876	2881	2883	rBV	2256	1857	0.02%	0.005%
68	10.428	2889	2904	2925	rBV	613546	983379	11.40%	2.620%
69	10.585	2945	2953	2966	rVB	64825	102306	1.19%	0.273%
70	10.679	2970	2982	3000	rBV	311772	660594	7.66%	1.760%
71	10.769	3001	3010	3025	rVB	138651	220967	2.56%	0.589%

LSC Area Percent Report

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Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\METHOD\SOMULM032319WMA.M
Title : VOC Analysis

72	10.968	3061	3072	3093	rVB	132545	215336	2.50%	0.574%
73	11.145	3115	3127	3143	rVV	525178	828182	9.60%	2.206%
74	11.219	3143	3150	3160	rVB5	9532	15317	0.18%	0.041%
75	11.424	3204	3214	3220	rBV2	8137	12604	0.15%	0.034%
76	11.479	3220	3231	3244	rBV	779788	1253421	14.53%	3.339%
77	11.569	3251	3259	3273	rVB	153476	244061	2.83%	0.650%
78	11.762	3308	3319	3329	rBV	162334	273990	3.18%	0.730%
79	11.855	3338	3348	3373	rVB	798939	1350225	15.65%	3.597%
80	11.977	3378	3386	3400	rVV2	42065	72203	0.84%	0.192%
81	12.055	3404	3410	3420	rVB2	10274	14655	0.17%	0.039%
82	12.145	3430	3438	3443	rBV3	21265	33524	0.39%	0.089%
83	12.251	3459	3471	3477	rBV2	31413	51580	0.60%	0.137%
84	12.350	3493	3502	3520	rVB2	29953	54090	0.63%	0.144%
85	12.550	3555	3564	3573	rVV4	7971	12912	0.15%	0.034%
86	12.650	3585	3595	3602	rBV2	17007	26422	0.31%	0.070%
87	12.701	3603	3611	3622	rVB2	26256	40086	0.46%	0.107%
88	12.961	3685	3692	3704	rVB3	26582	40538	0.47%	0.108%
89	13.122	3726	3742	3753	rBV3	53523	91752	1.06%	0.244%
90	13.186	3758	3762	3768	rVB5	4401	4618	0.05%	0.012%
91	13.286	3785	3793	3810	rVB5	10027	20585	0.24%	0.055%
92	13.450	3840	3844	3847	rBV3	2731	2702	0.03%	0.007%
93	13.511	3855	3863	3868	rBV6	5754	8825	0.10%	0.024%
94	13.743	3925	3935	3957	rBV	107808	219294	2.54%	0.584%
95	13.961	3996	4003	4004	rBV2	2049	2197	0.03%	0.006%
96	14.267	4096	4098	4102	rVV3	1786	1633	0.02%	0.004%
97	14.338	4113	4120	4124	rBV4	2892	3946	0.05%	0.011%
98	14.887	4285	4291	4292	rBV	3853	3259	0.04%	0.009%
99	15.456	4465	4468	4475	rBV3	2024	2352	0.03%	0.006%
100	15.916	4608	4611	4615	rVB	3145	2382	0.03%	0.006%

Sum of corrected areas: 37540256

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
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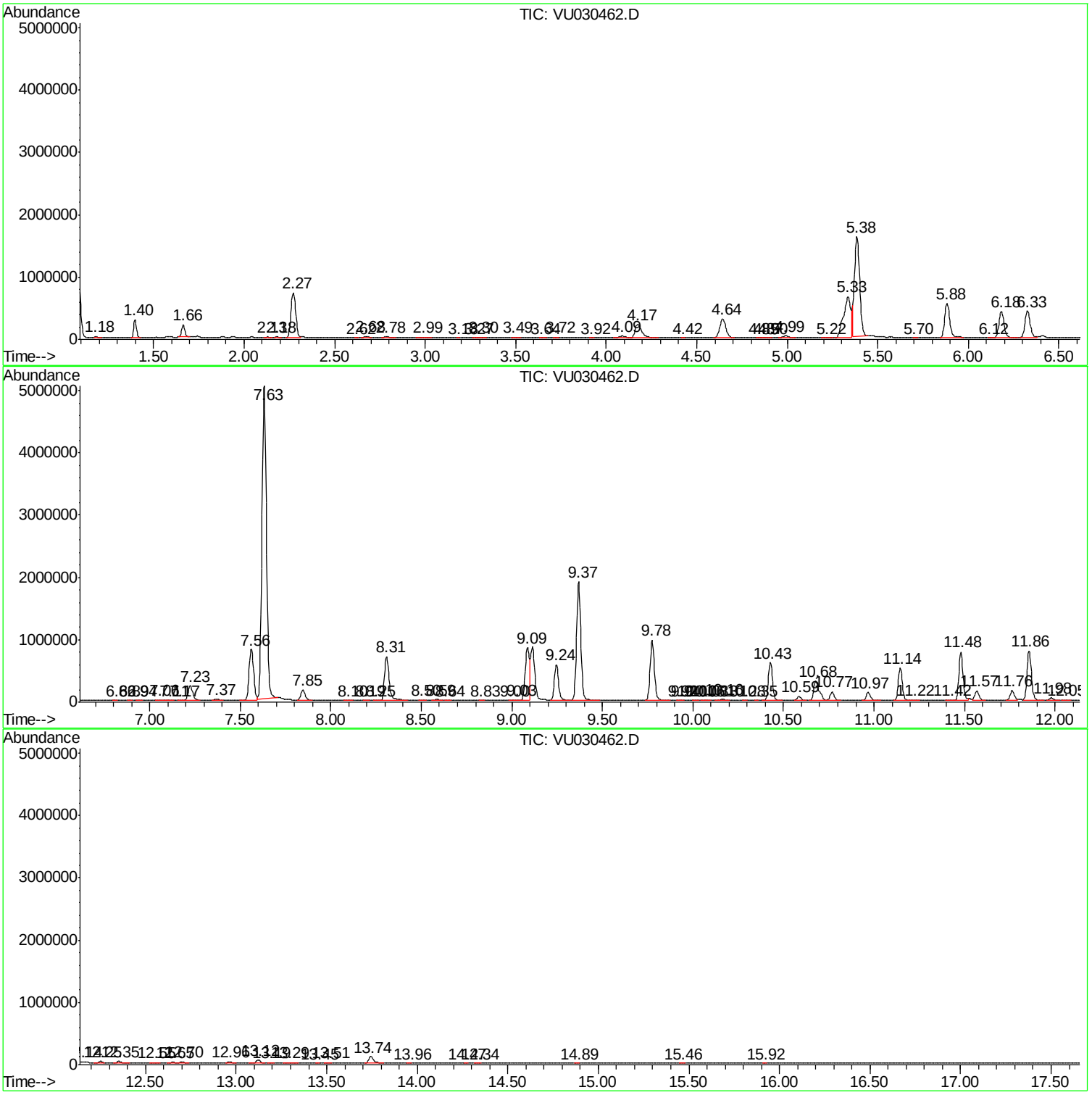
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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



Library Search Compound Report

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Quant Title : VOC Analysis

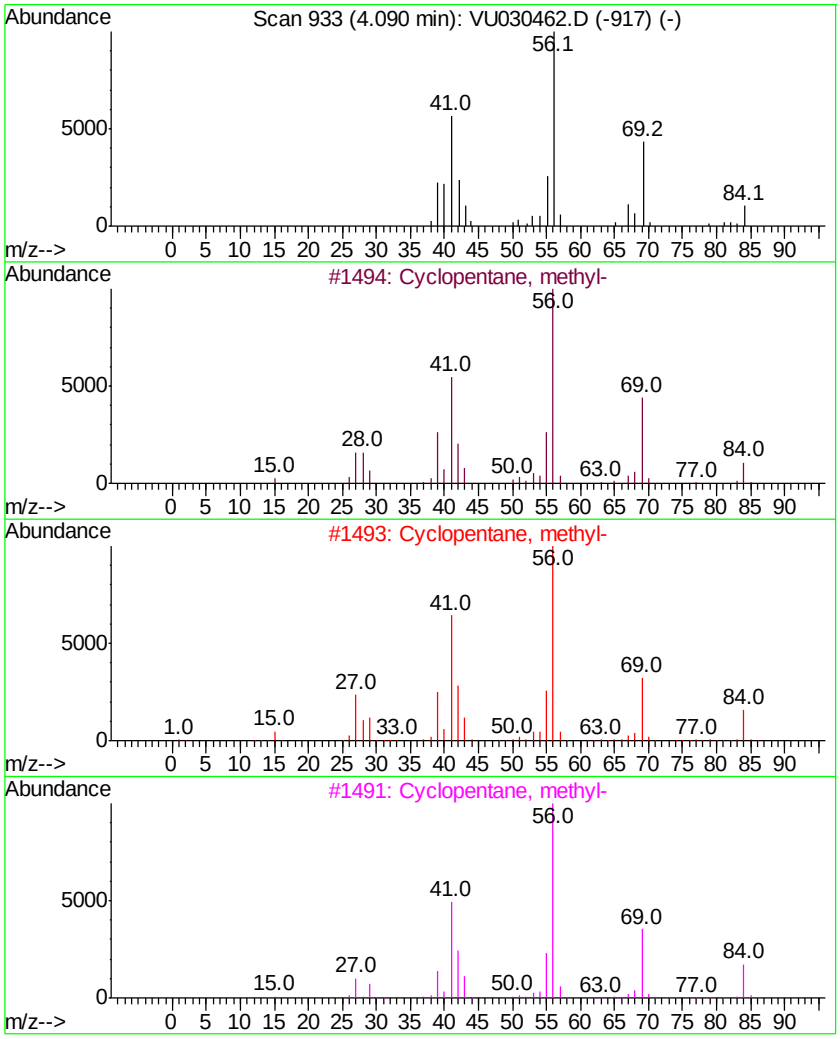
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

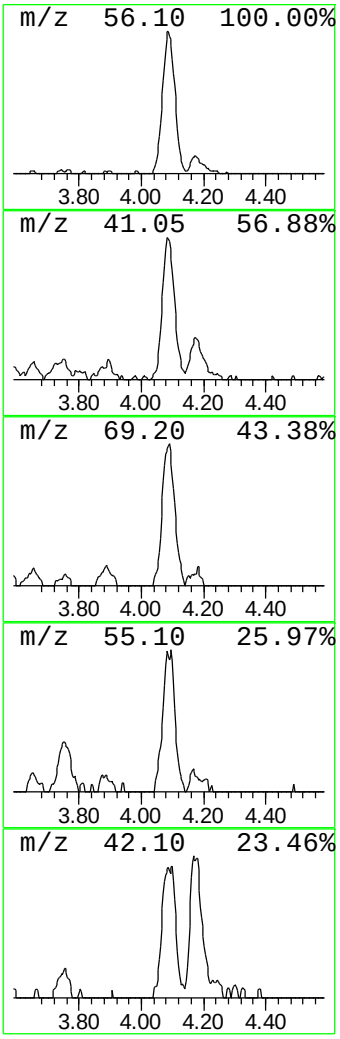
****

Peak Number 1 Cyclopentane, methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	3.61 ug/L	80600	1,4-Difluorobenzene	5.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

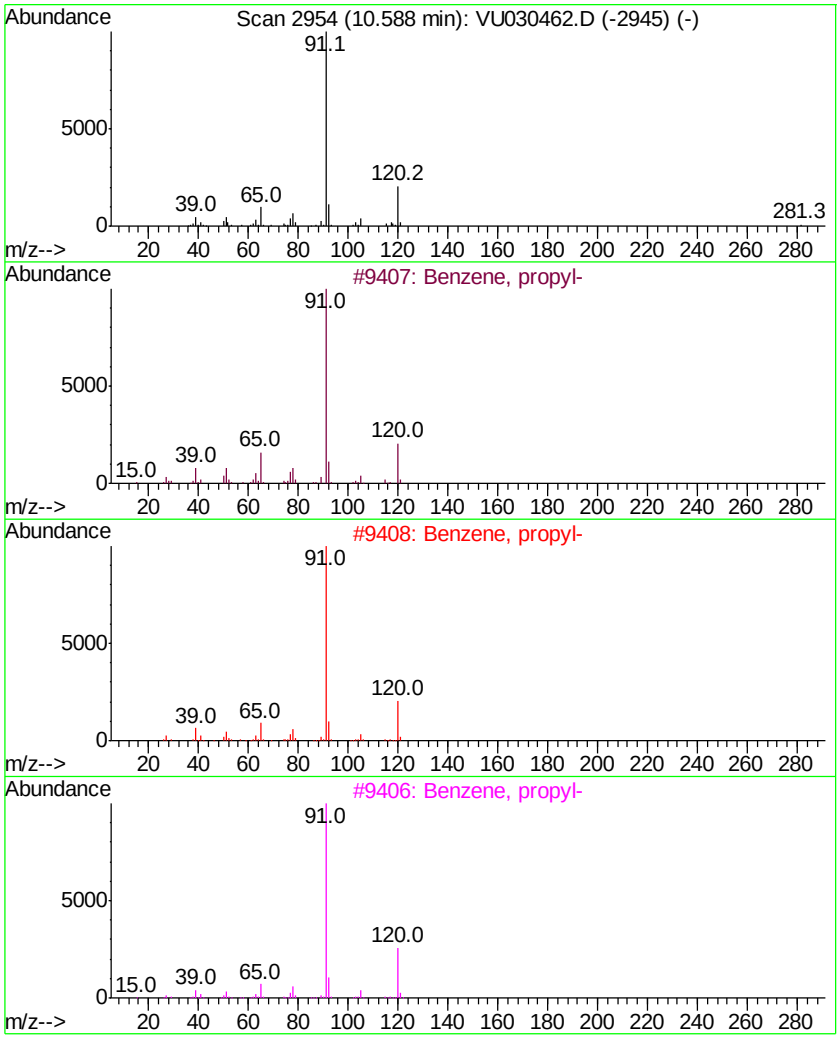
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

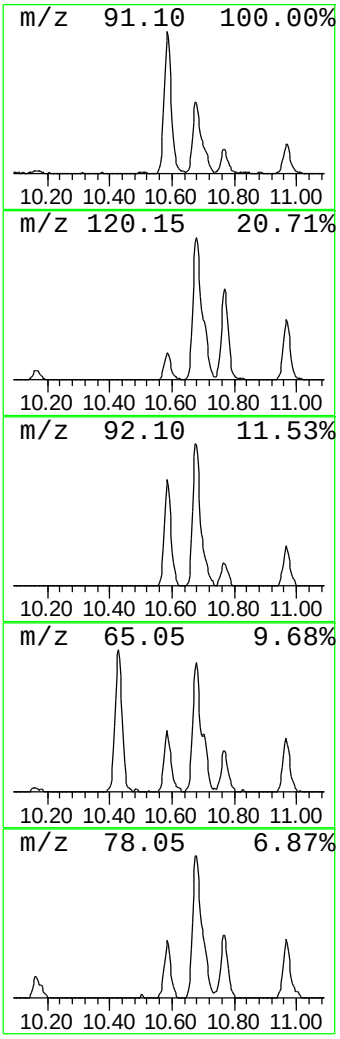
****

Peak Number 3 Benzene, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.59	4.08 ug/L	102306	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	90
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	83
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78





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Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

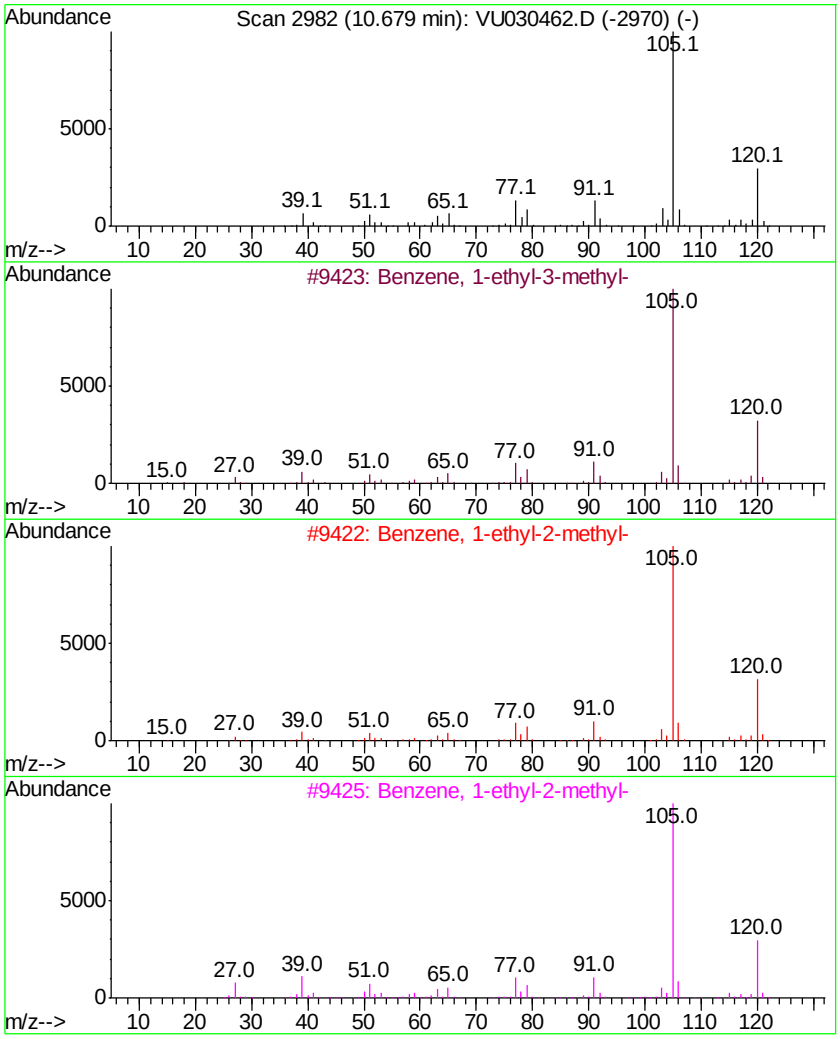
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

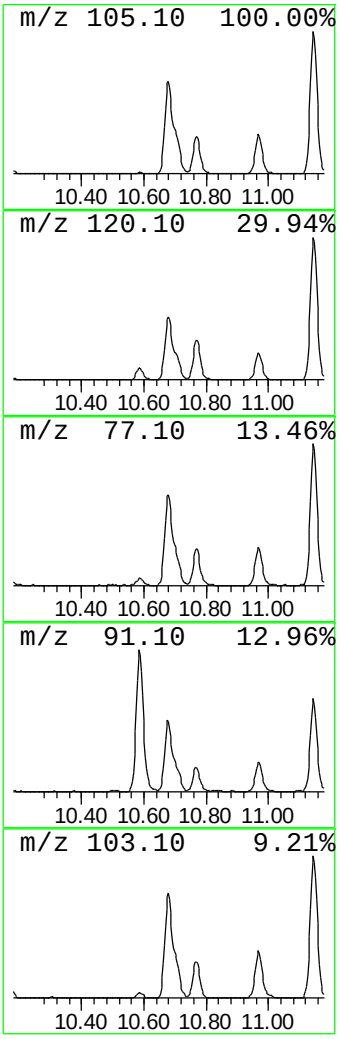
****

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	26.35 ug/L	660594	1,4-Dichlorobenzene-d4	11.48

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-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030462.D
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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

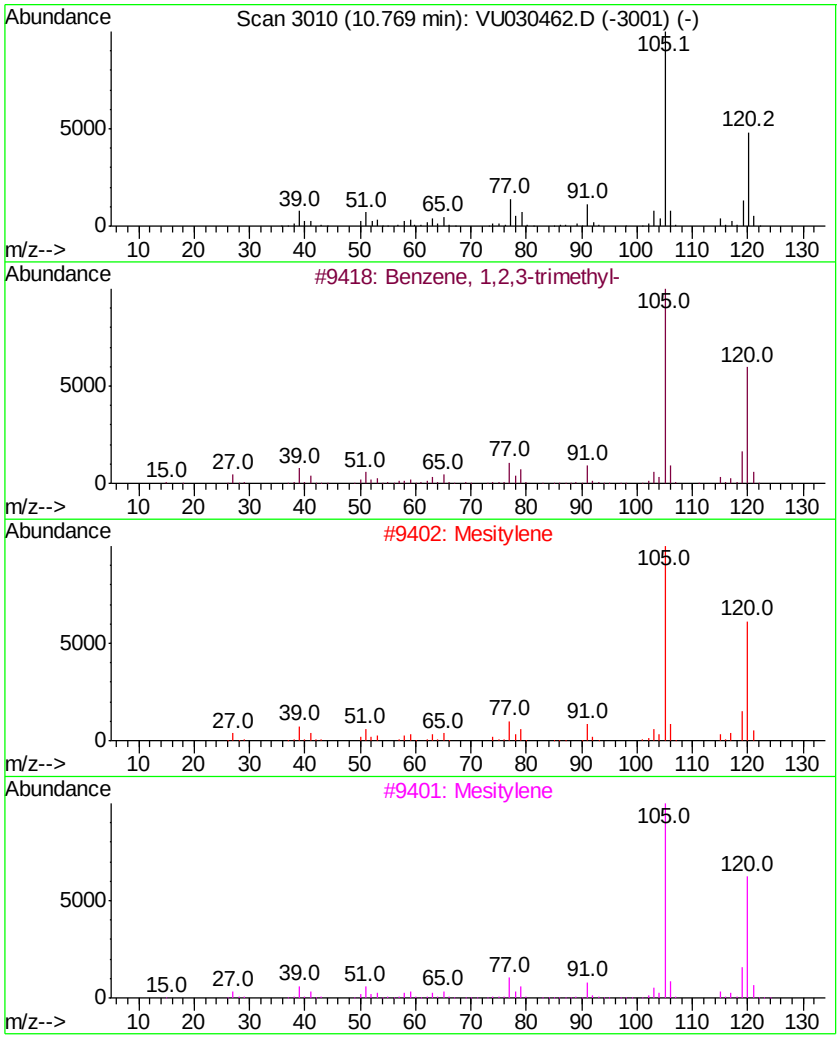
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

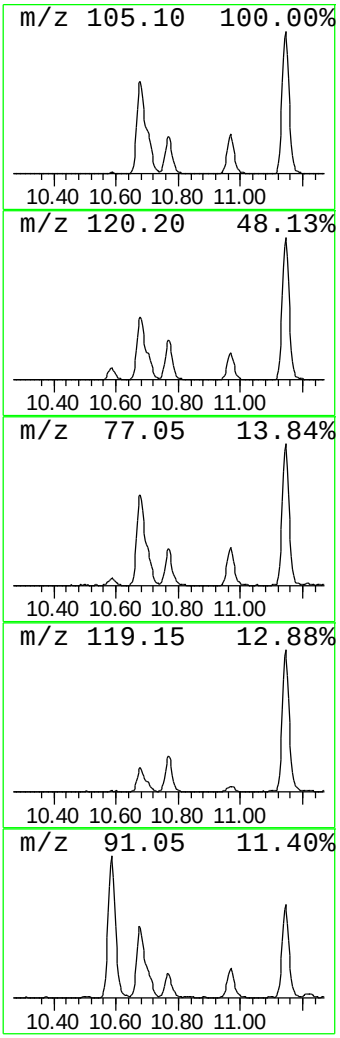
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	8.81 ug/L	220967	1,4-Dichlorobenzene-d4	11.48

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





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

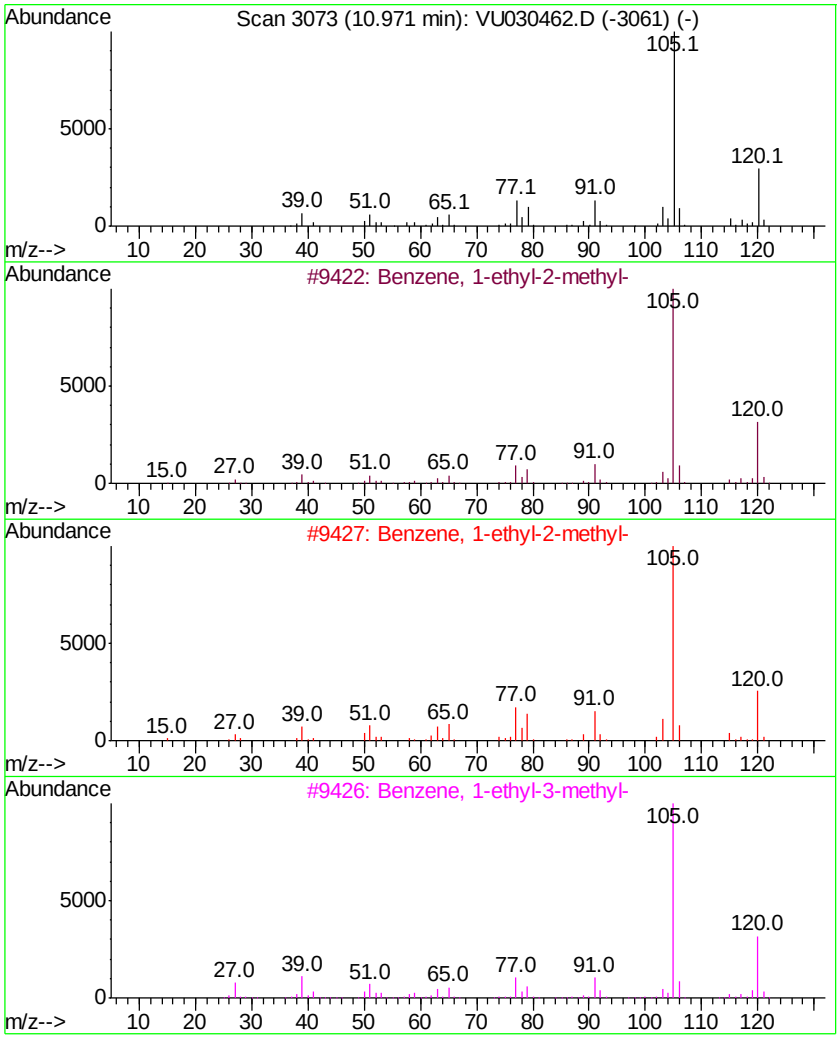
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

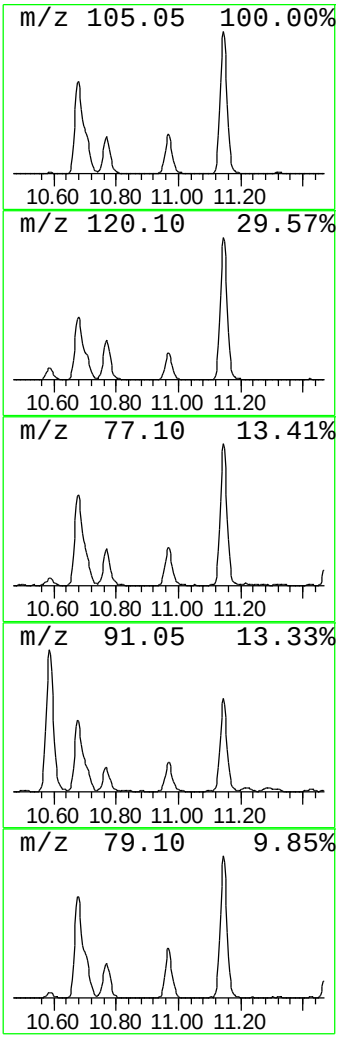
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	8.59 ug/L	215336	1,4-Dichlorobenzene-d4	11.48

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-2-methyl-	120	C9H12	000611-14-3	94
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	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

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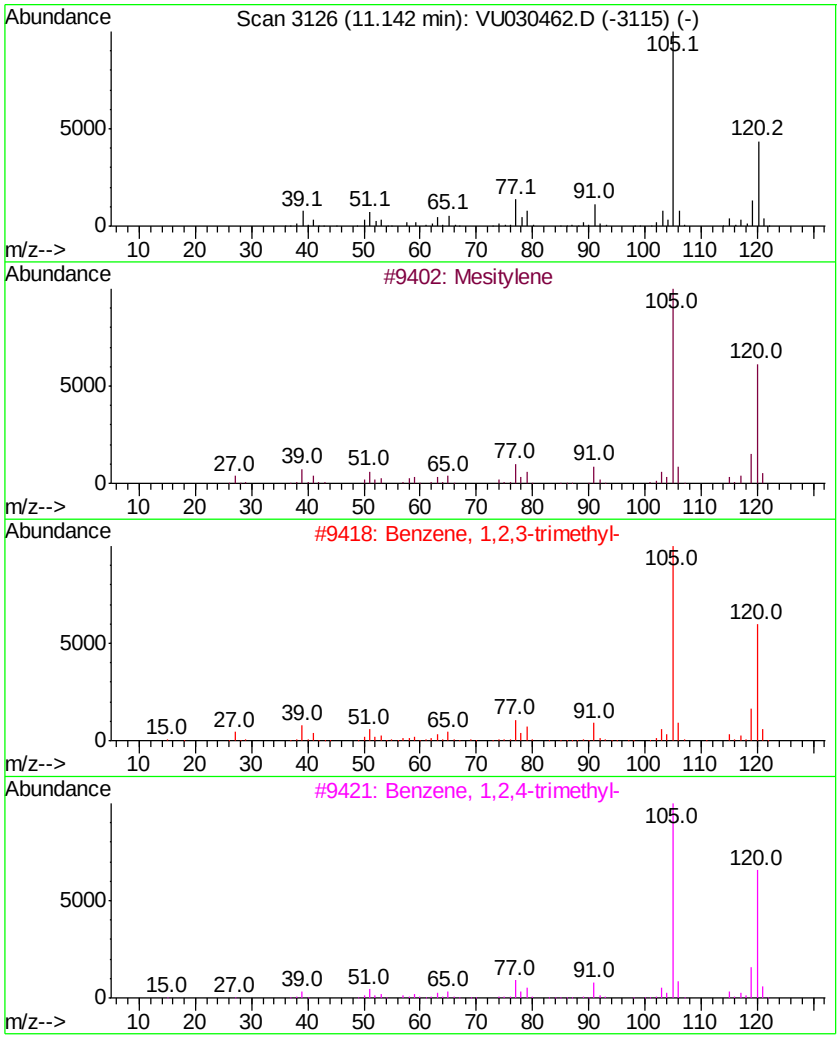
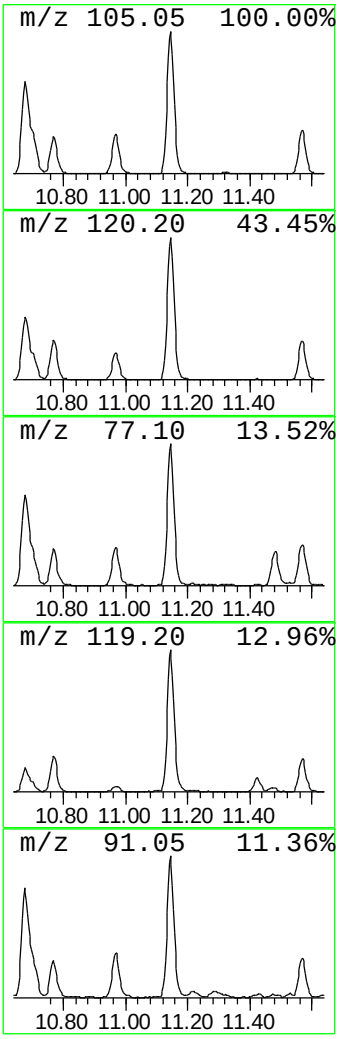
Peak Number 7 Mesitylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.14	33.04 ug/L	828182	1,4-Dichlorobenzene-d4	11.48

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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

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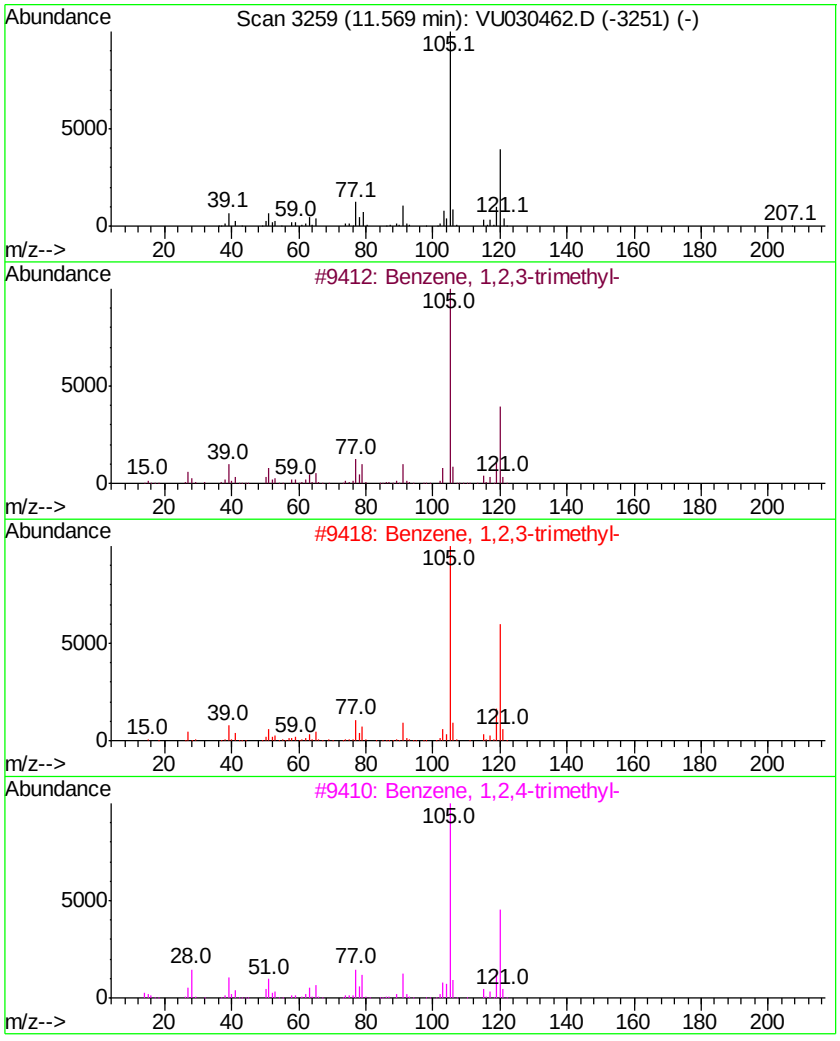
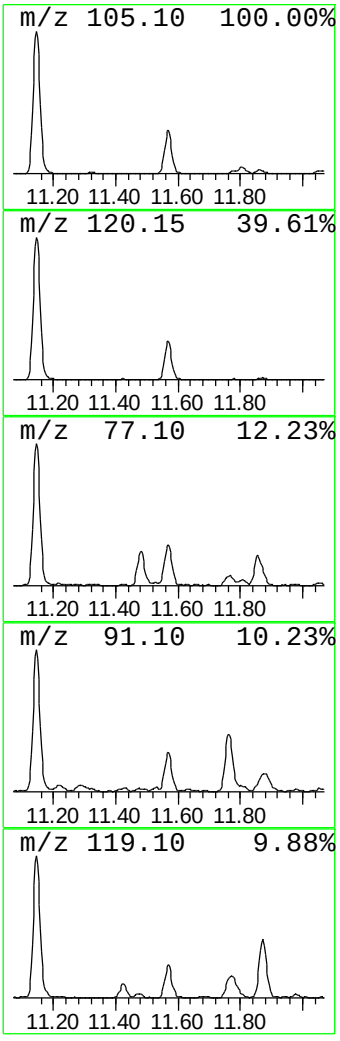
Peak Number 8 Benzene, 1,2,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	9.74 ug/L	244061	1,4-Dichlorobenzene-d4	11.48

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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

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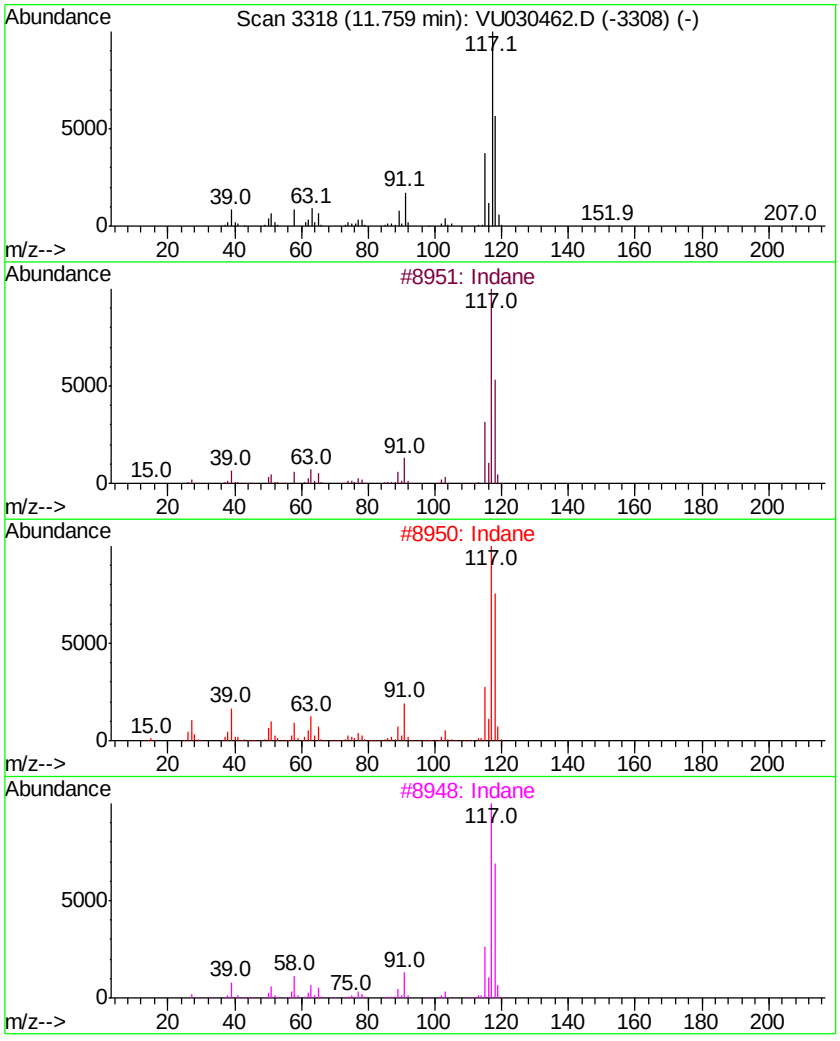
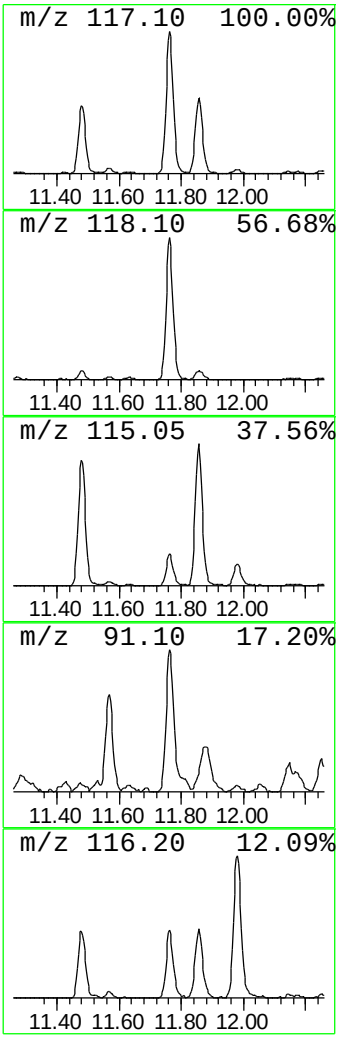
Peak Number 9 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	10.93 ug/L	273990	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	90
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

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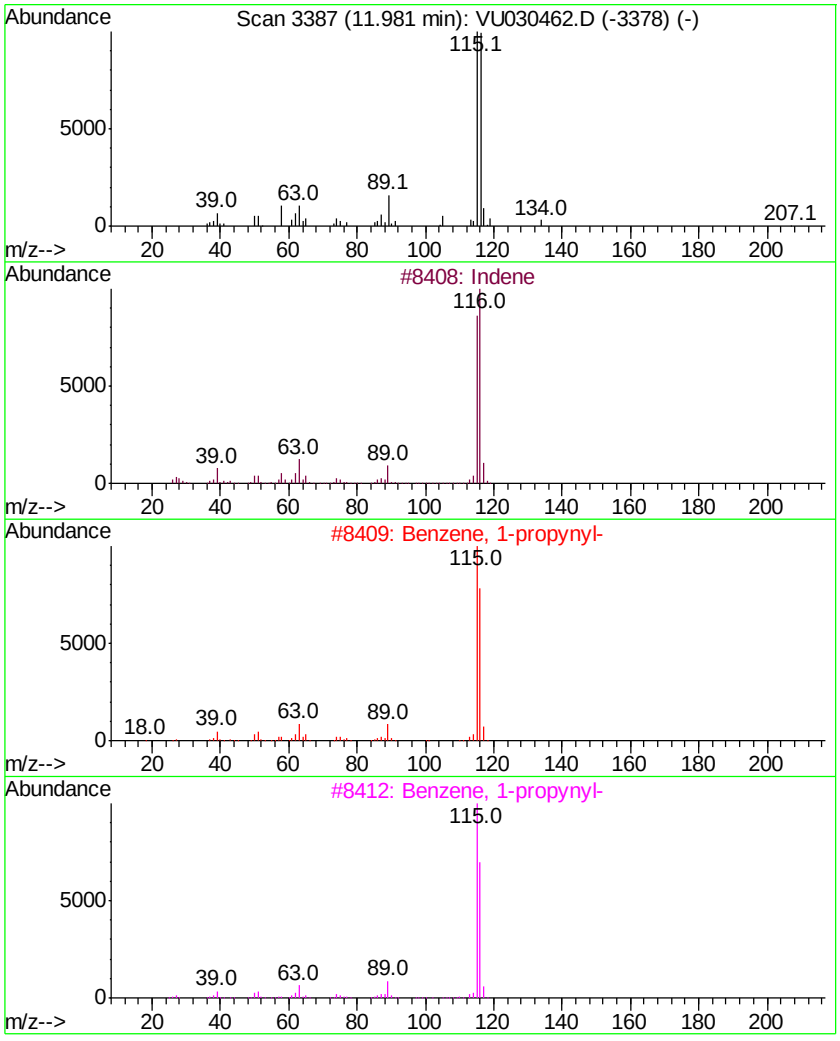
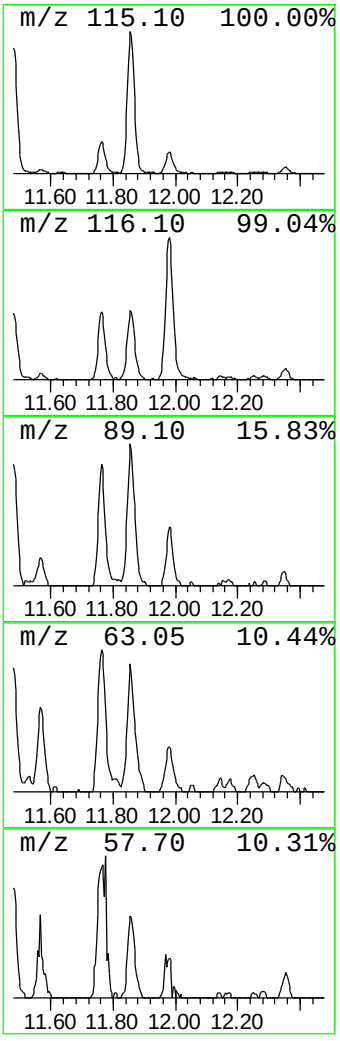
Peak Number 10 Indene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	2.88 ug/L	72203	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	93
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
4	Indene	116	C9H8	000095-13-6	90
5	3-Methylphenylacetylene	116	C9H8	000766-82-5	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

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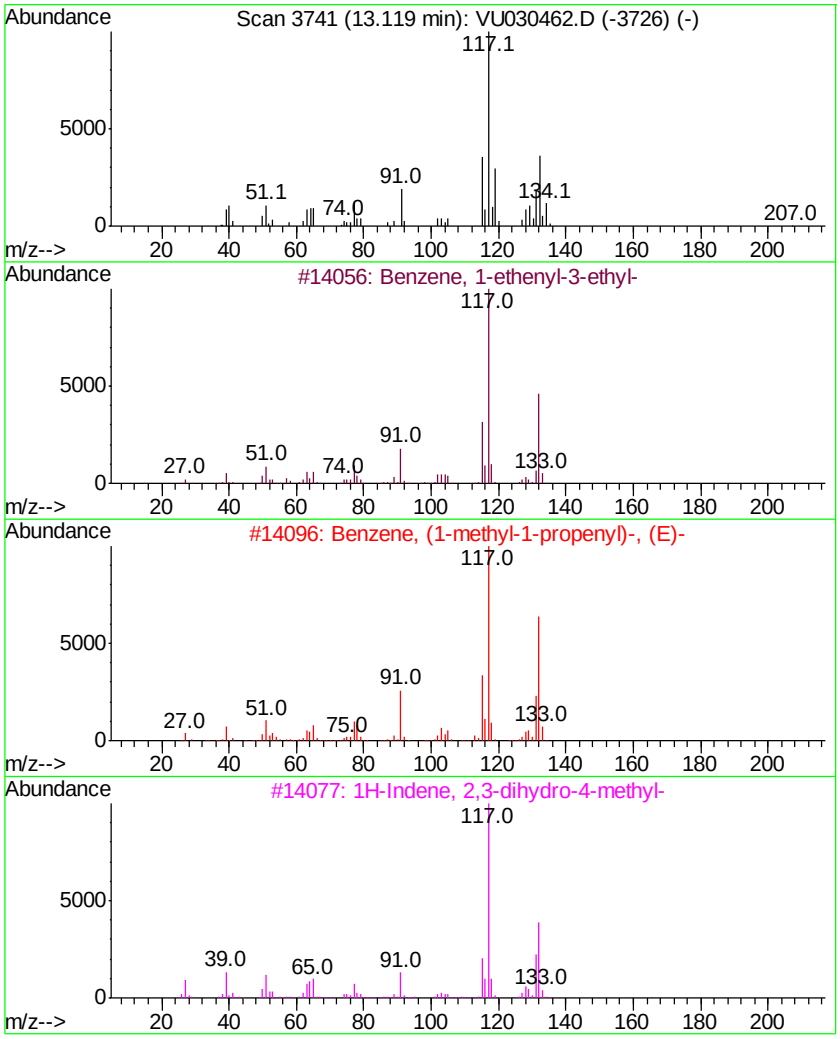
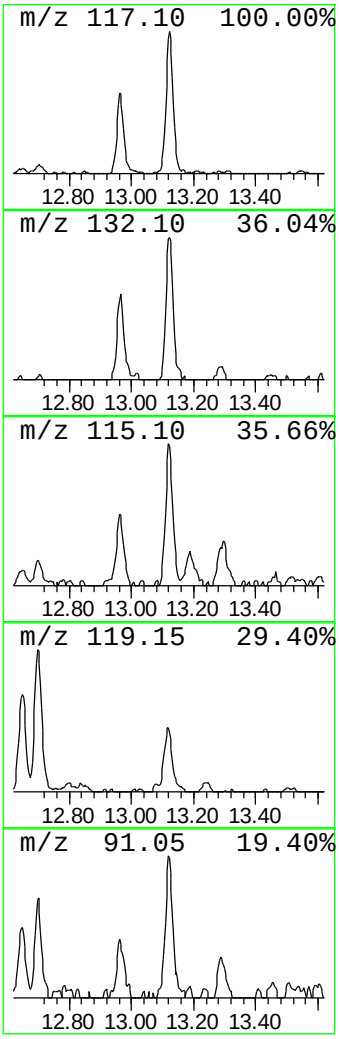
Peak Number 11 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	3.66 ug/L	91752	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	76
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030462.D
Acq On : 26 Mar 2019 22:38
Operator : JC/SP
Sample : K2063-07MSD 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R7MSD

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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...	4.09	3.6	ug/L	80600	1	5.88	1117670	50.0
Benzene, propyl-	10.59	4.1	ug/L	102306	3	11.48	1253420	50.0
Benzene, 1-ethyl-...	10.68	26.4	ug/L	660594	3	11.48	1253420	50.0
Benzene, 1,2,3-tr...	10.77	8.8	ug/L	220967	3	11.48	1253420	50.0
Benzene, 1-ethyl-...	10.97	8.6	ug/L	215336	3	11.48	1253420	50.0
Mesitylene	11.14	33.0	ug/L	828182	3	11.48	1253420	50.0
Benzene, 1,2,4-tr...	11.57	9.7	ug/L	244061	3	11.48	1253420	50.0
Indane	11.76	10.9	ug/L	273990	3	11.48	1253420	50.0
Indene	11.98	2.9	ug/L	72203	3	11.48	1253420	50.0
Benzene, 1-etheny...	13.12	3.7	ug/L	91752	3	11.48	1253420	50.0