LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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	1.183	24	29	35	rVB	14073	12368	0.25%	0.041%
2	1.399	88	96	108	rBV	299142	306876	6.29%	1.025%
3	1.669	173	180	196	rVV	208671	266860	5.47%	0.892%
4	1.746	197	204	223	rVB	108701	153231	3.14%	0.512%
5	1.939	257	264	279	rVB	73030	101229	2.07%	0.338%
6	2.045	291	297	310	rVB	21738	29545	0.61%	0.099%
7	2.135	320	325	331	rVB3	8758	10073	0.21%	0.034%
8	2.183	334	340	349	rVB8	7170	10483	0.21%	0.035%
9	2.267	356	366	380	rBV	419802	638197	13.08%	2.132%
10	2.633	472	480	485	rBV3	2980	4164	0.09%	0.014%
11	2.701	487	501	504	rBV5	10712	23459	0.48%	0.078%
12	2.990	578	591	612	rVB4	20972	50787	1.04%	0.170%
13	3.080	613	619	622	rBV	1469	1676	0.03%	0.006%
14	3.100	622	625	630	rVB	2217	1764	0.04%	0.006%
15	3.299	673	687	703	rBV4	15942	35611	0.73%	0.119%
16	3.421	719	725	734	rBV3	4398	6776	0.14%	0.023%
17	3.501	738	750	760	rVB5	7193	15168	0.31%	0.051%
18	3.640	788	793	798	rBV2	1685	1932	0.04%	0.006%
19	3.720	809	818	819	rBV	5350	5736	0.12%	0.019%
20	3.887	862	870	872	rBV	2524	2760	0.06%	0.009%
21	4.087	915	932	948	rBV2	64500	174186	3.57%	0.582%
22	4.173	948	959	985	rBV	209924	538157	11.03%	1.798%
23	4.546	1070	1075	1081	rVB2	1405	1706	0.03%	0.006%
24	4.643	1086	1105	1126	rBV	304534	726658	14.89%	2.428%
25	4.990	1198	1213	1233	rBV3	53068	134630	2.76%	0.450%
26	5.183	1265	1273	1275	rBV2	1895	2035	0.04%	0.007%
27	5.218	1275	1284	1293	rBV6	5495	12794	0.26%	0.043%
28	5.334	1298	1320	1333	rBV2	634786	1879905	38.53%	6.281%
29	5.620	1401	1409	1418	rVB5	6865	13032	0.27%	0.044%
30	5.771	1451	1456	1458	rBV2	3591	3270	0.07%	0.011%
31	5.881	1477	1490	1506	rBV	543916	1082607	22.19%	3.617%
32	6.328	1613	1629	1645	rBV	430259	868691	17.81%	2.902%
33	6.640	1720	1726	1736	rVB5	5802	11141	0.23%	0.037%
34	6.813	1776	1780	1783	rBV2	3143	3075	0.06%	0.010%

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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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	6.990	1828	1835	1839	rBV	1671	2159	0.04%	0.007%
36	7.045	1848	1852	1855	rBV	3016	2955	0.06%	0.010%
37	7.132	1872	1879	1884	rBV3	1933	2372	0.05%	0.008%
38	7.225	1890	1908	1924	rBV	230715	427288	8.76%	1.428%
39	7.373	1948	1954	1964	rVB4	10658	17082	0.35%	0.057%
40	7.562	1993	2013	2024	rBV	808315	1423210	29.17%	4.755%
41	7.633	2024	2035	2082	rVB	2759522	4878798	100.00%	16.300%
42	7.845	2092	2101	2120	rVB	159367	268011	5.49%	0.895%
43	8.032	2154	2159	2162	rBV2	2509	2006	0.04%	0.007%
44	8.164	2195	2200	2207	rVB4	2187	2415	0.05%	0.008%
45	8.241	2216	2224	2233	rBV4	6937	13921	0.29%	0.047%
46	8.308	2234	2245	2278	rBV	668757	1221620	25.04%	4.081%
47	8.646	2343	2350	2354	rBV4	2810	4004	0.08%	0.013%
48	8.746	2375	2381	2384	rBV	2653	2517	0.05%	0.008%
49	8.903	2424	2430	2431	rBV2	1533	1814	0.04%	0.006%
50	9.029	2459	2469	2475	rBV5	9321	18052	0.37%	0.060%
51	9.086	2476	2487	2507	rBV	806652	1356703	27.81%	4.533%
52	9.247	2525	2537	2559	rBV	698795	1135333	23.27%	3.793%
53	9.369	2562	2575	2610	rVB	2208000	3708872	76.02%	12.391%
54	9.594	2641	2645	2647	rBV2	2166	2042	0.04%	0.007%
55	9.614	2648	2651	2655	rVB2	2530	1822	0.04%	0.006%
56	9.723	2680	2685	2688	rBV3	1880	2106	0.04%	0.007%
57	9.774	2688	2701	2729	rVV	1125427	1875388	38.44%	6.266%
58	9.942	2747	2753	2756	rBV2	4397	4735	0.10%	0.016%
59	10.164	2814	2822	2836	rBV	23719	39461	0.81%	0.132%
60	10.218	2836	2839	2844	rVB2	2299	1688	0.03%	0.006%
61	10.350	2876	2880	2884	rBV	1855	1718	0.04%	0.006%
62	10.389	2888	2892	2893	rVV	2886	2170	0.04%	0.007%
63	10.427	2893	2904	2925	rVB	597947	954267	19.56%	3.188%
64	10.585	2945	2953	2967	rVB	57620	92295	1.89%	0.308%
65	10.678	2971	2982	3001	rBV	268540	574580	11.78%	1.920%
66	10.768	3001	3010	3022	rVB2	115538	178545	3.66%	0.597%
67	10.967	3060	3072	3092	rBV	120150	199608	4.09%	0.667%
68	11.144	3115	3127	3143	rVV	464254	733163	15.03%	2.449%
69	11.221	3144	3151	3160	rVB7	8183	14510	0.30%	0.048%
70	11.263	3161	3164	3167	rBV2	2220	1870	0.04%	0.006%
71	11.427	3205	3215	3220	rBV4	6304	9605	0.20%	0.032%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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	11.482	3220	3232	3250	rVV	751209	1239228	25.40%	4.140%
73	11.569	3251	3259	3271	rVB	140856	220573	4.52%	0.737%
74	11.762	3305	3319	3330	rBV	153673	267611	5.49%	0.894%
75	11.855	3338	3348	3370	rVB	752525	1266872	25.97%	4.233%
76	11.980	3378	3387	3401	rVB	38166	64102	1.31%	0.214%
77	12.054	3401	3410	3418	rBV4	7354	13077	0.27%	0.044%
78	12.093	3419	3422	3426	rVB	2582	2156	0.04%	0.007%
79	12.147	3430	3439	3444	rBV4	15743	27215	0.56%	0.091%
80	12.250	3459	3471	3479	rBV3	21046	37381	0.77%	0.125%
81	12.286	3479	3482	3488	rVV3	8975	8993	0.18%	0.030%
82	12.353	3495	3503	3518	rVB2	24325	41844	0.86%	0.140%
83	12.549	3557	3564	3576	rVB3	6275	11040	0.23%	0.037%
84	12.652	3587	3596	3602	rBV3	11249	16873	0.35%	0.056%
85	12.700	3604	3611	3626	rVB2	21037	36483	0.75%	0.122%
86	12.762	3626	3630	3632	rBV	2389	1919	0.04%	0.006%
87	12.781	3632	3636	3644	rVB2	2221	2435	0.05%	0.008%
88	12.823	3644	3649	3651	rBV	2097	1835	0.04%	0.006%
89	12.964	3682	3693	3707	rBV3	21010	34897	0.72%	0.117%
90	13.122	3730	3742	3757	rBV4	41095	69256	1.42%	0.231%
91	13.289	3786	3794	3805	rBV6	8624	17235	0.35%	0.058%
92	13.405	3824	3830	3837	rVB3	2189	2857	0.06%	0.010%
93	13.456	3839	3846	3853	rBV2	3046	5292	0.11%	0.018%
94	13.514	3859	3864	3871	rVB4	4676	5981	0.12%	0.020%
95	13.581	3881	3885	3888	rVB	2359	1714	0.04%	0.006%
96	13.742	3924	3935	3961	rBV	99636	216191	4.43%	0.722%
97	14.006	4014	4017	4023	rBV	3195	2890	0.06%	0.010%
98	14.253	4087	4094	4096	rBV	1547	1904	0.04%	0.006%
99	15.244	4399	4402	4406	rBV	2341	2059	0.04%	0.007%
100	15.610	4512	4516	4520	rBV2	2105	2154	0.04%	0.007%

Sum of corrected areas: 29931354

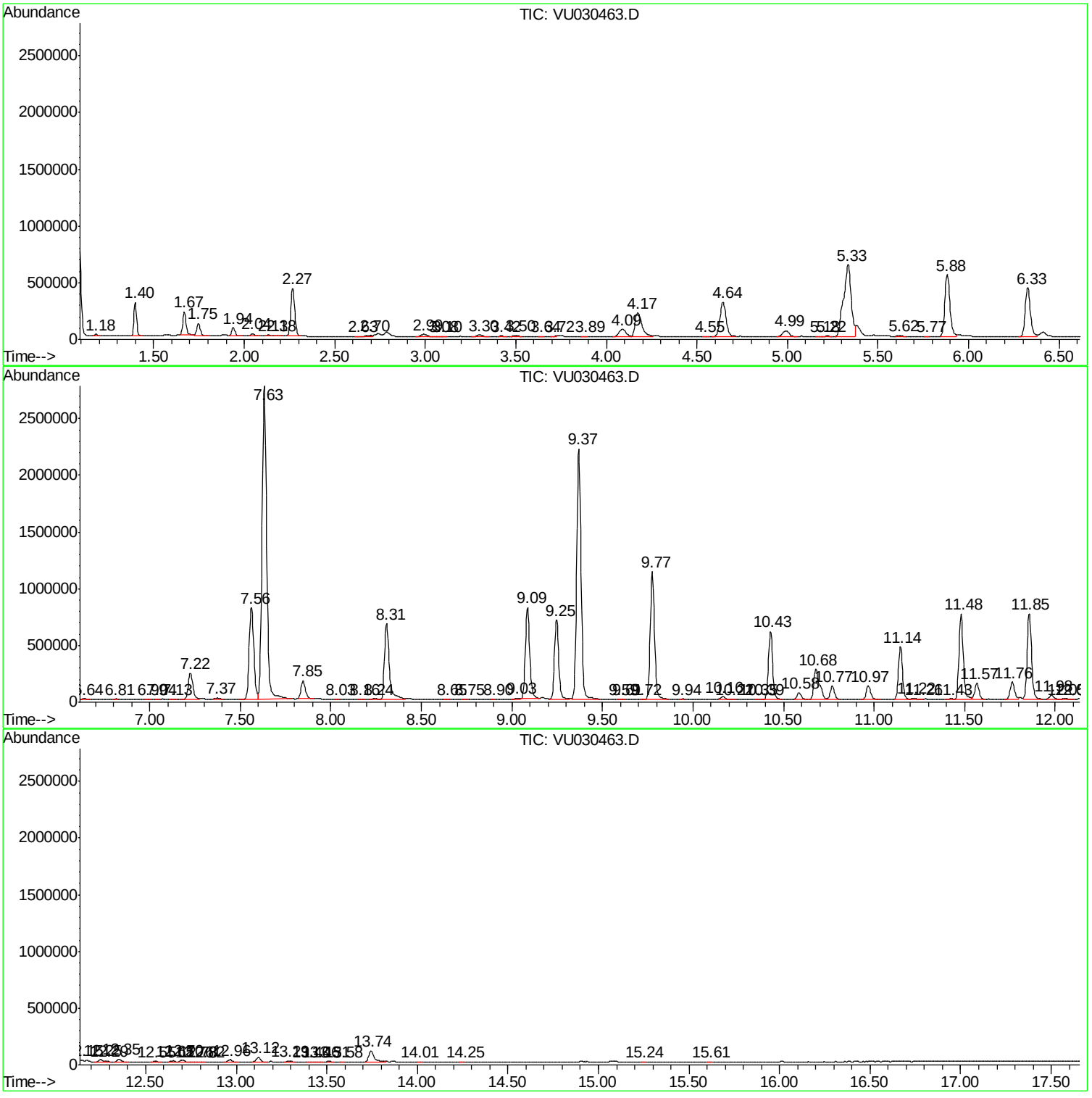
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

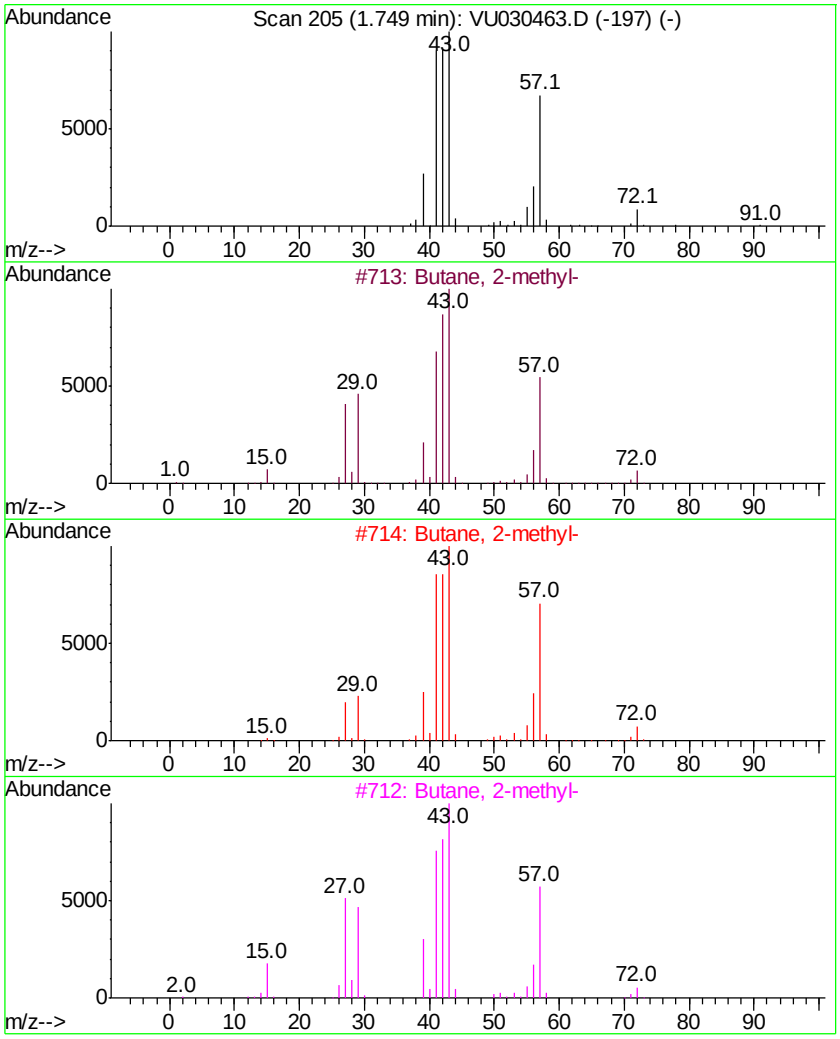
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

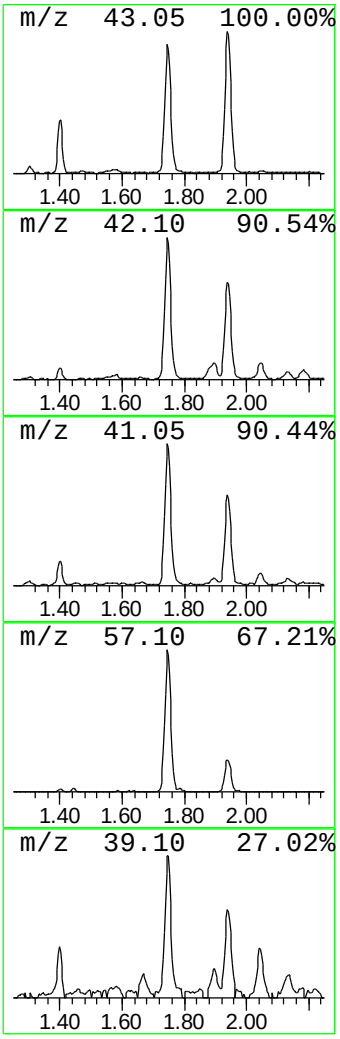
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.75	7.08 ug/L	153231	1,4-Difluorobenzene	5.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	91
3	Butane, 2-methyl-	72	C5H12	000078-78-4	91
4	Pentane	72	C5H12	000109-66-0	45
5	3-Buten-1-ol	72	C4H8O	000627-27-0	25





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

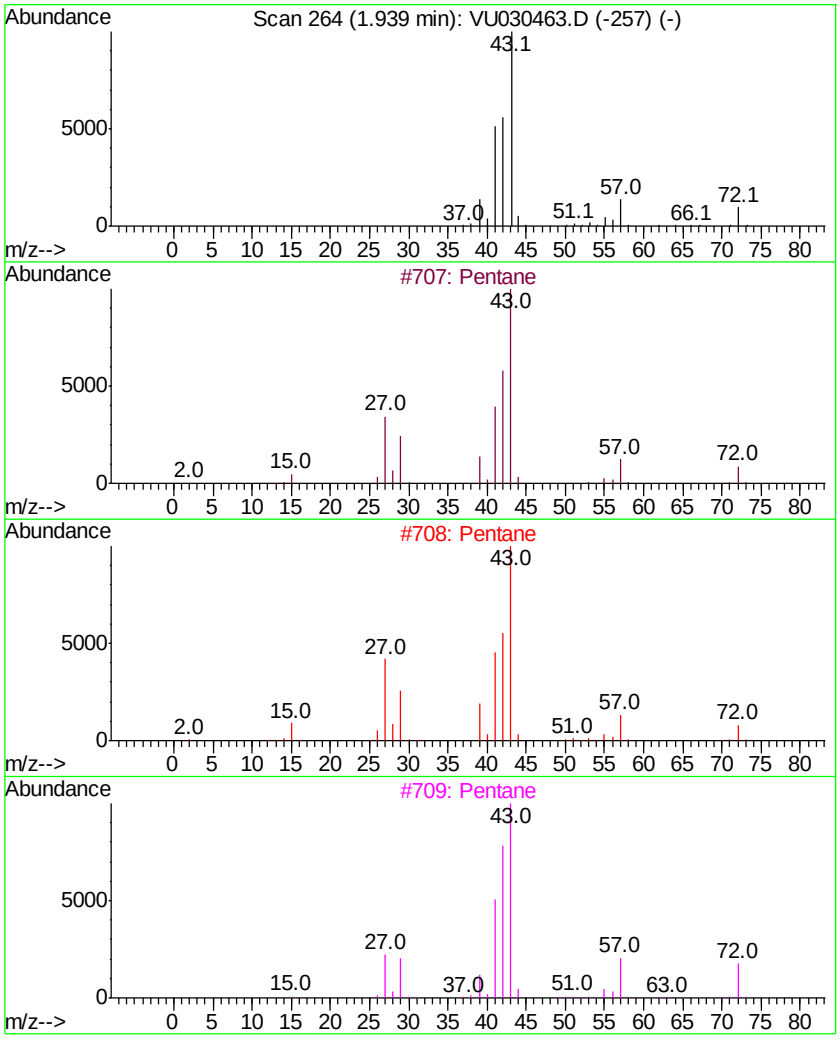
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

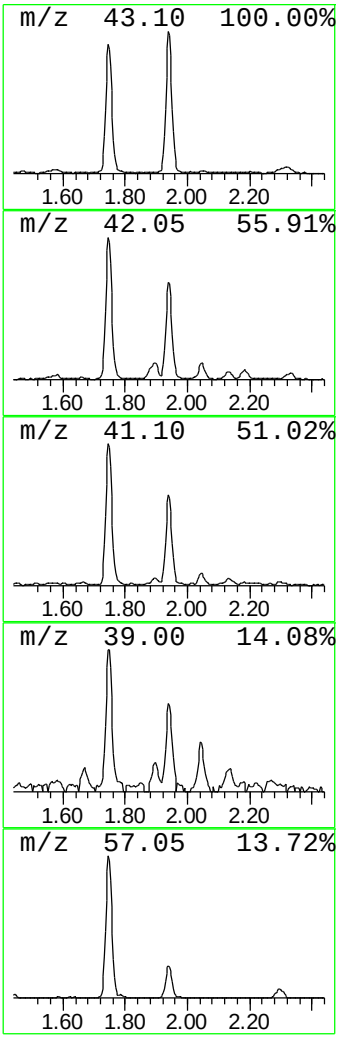
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.94	4.68 ug/L	101229	1,4-Difluorobenzene	5.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	80
2	Pentane	72	C5H12	000109-66-0	80
3	Pentane	72	C5H12	000109-66-0	72
4	Pentane	72	C5H12	000109-66-0	43
5	Hydrazine, 2-propenyl-	72	C3H8N2	007422-78-8	9





Library Search Compound Report

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ClientSampleId : DB6R8

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

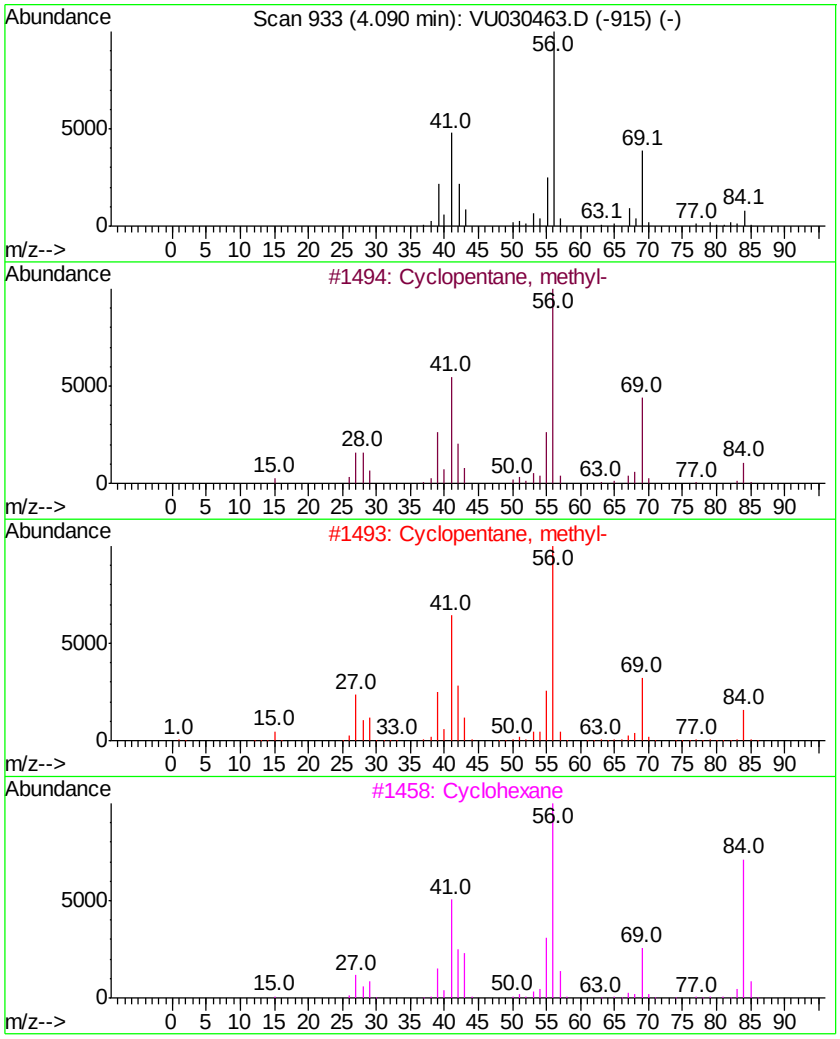
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

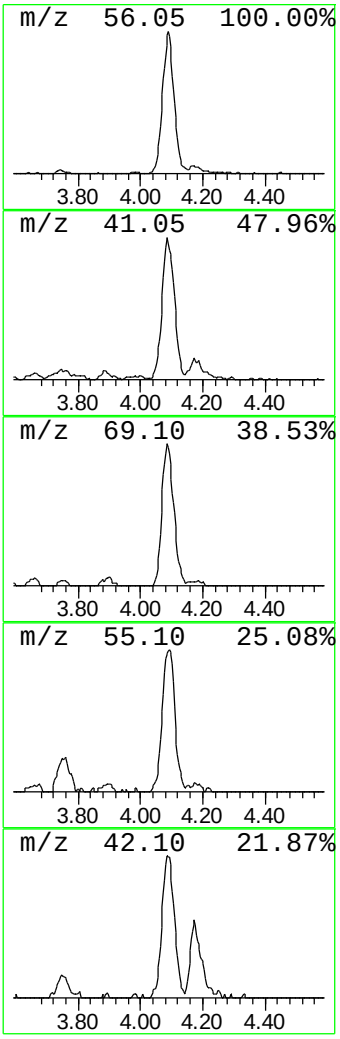
****

Peak Number 3 (DEL) Alkane: Cyclic4.09 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	8.04 ug/L	174186	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	86
3	Cyclohexane	84	C6H12	000110-82-7	72
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	59





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Instrument : MSVOA_U
ClientSampleId : DB6R8

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

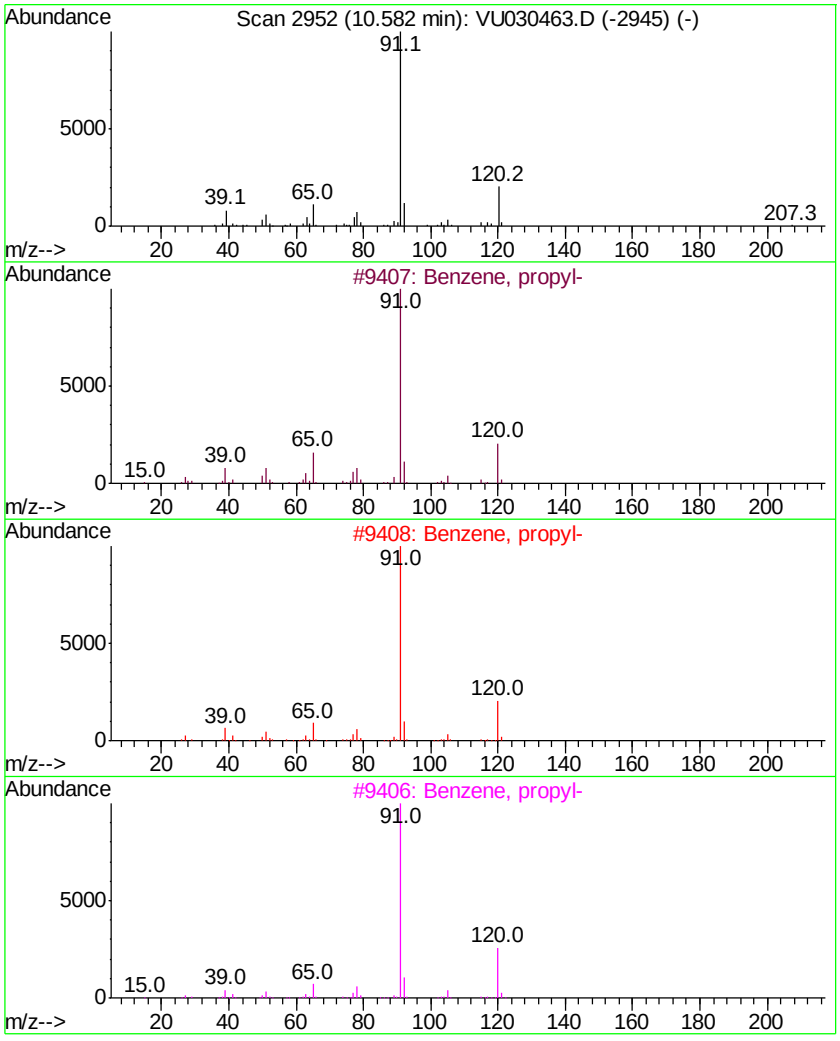
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

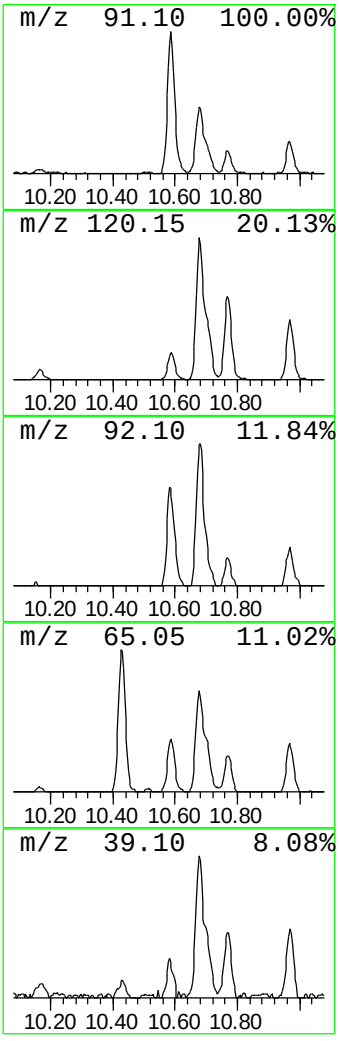
****

Peak Number 5 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	3.72 ug/L	92295	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	87
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	83
4	n-Butylbenzylamine	163	C11H17N	002403-22-7	64
5	2,4,6-Cycloheptatrien-1-one, 4-m...	120	C8H8O	1000327-34-9	50





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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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

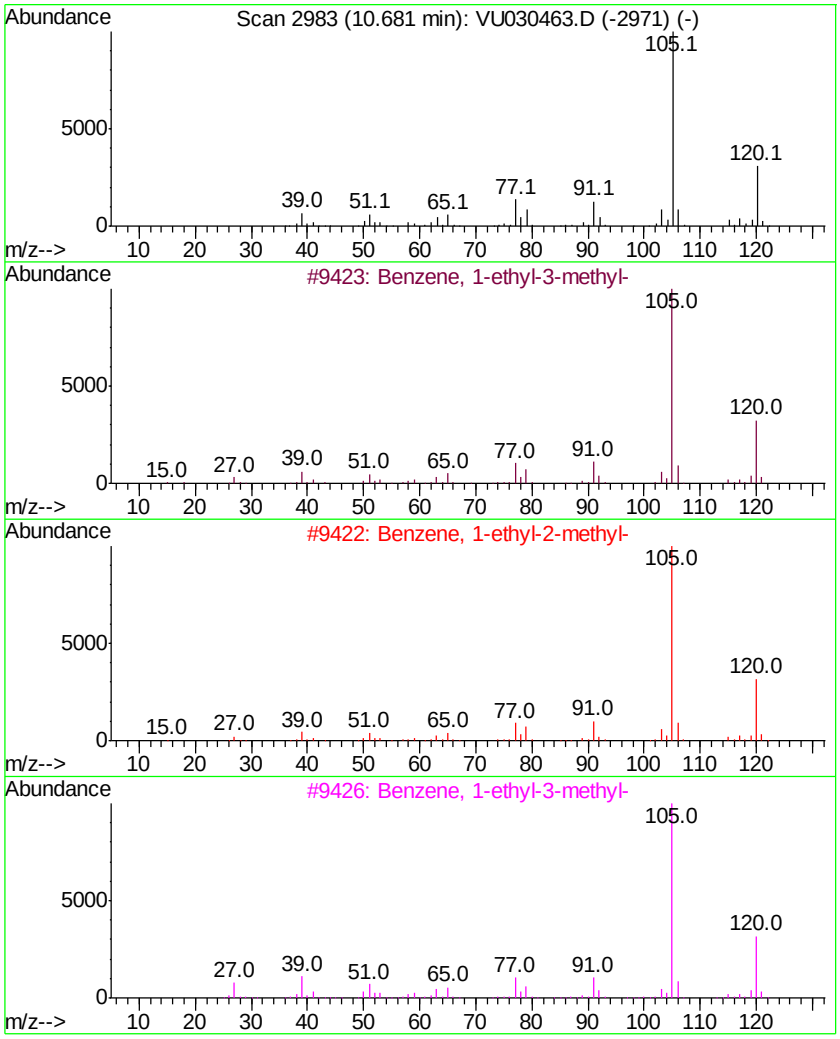
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

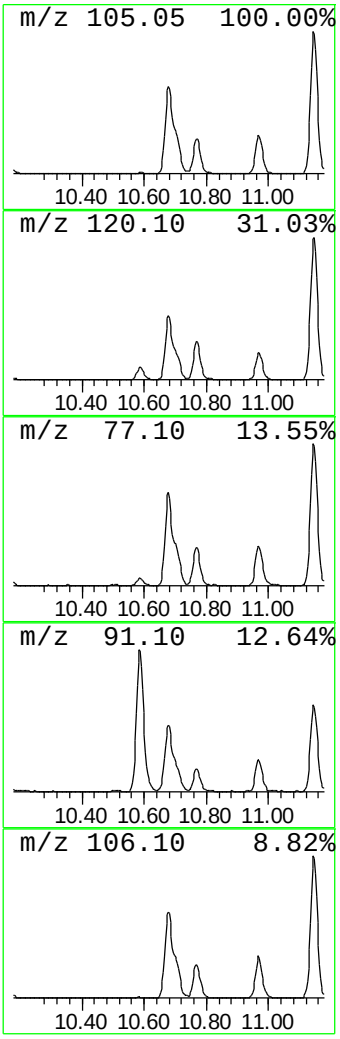
****

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	23.18 ug/L	574580	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-3-methyl-	120	C9H12	000620-14-4	94
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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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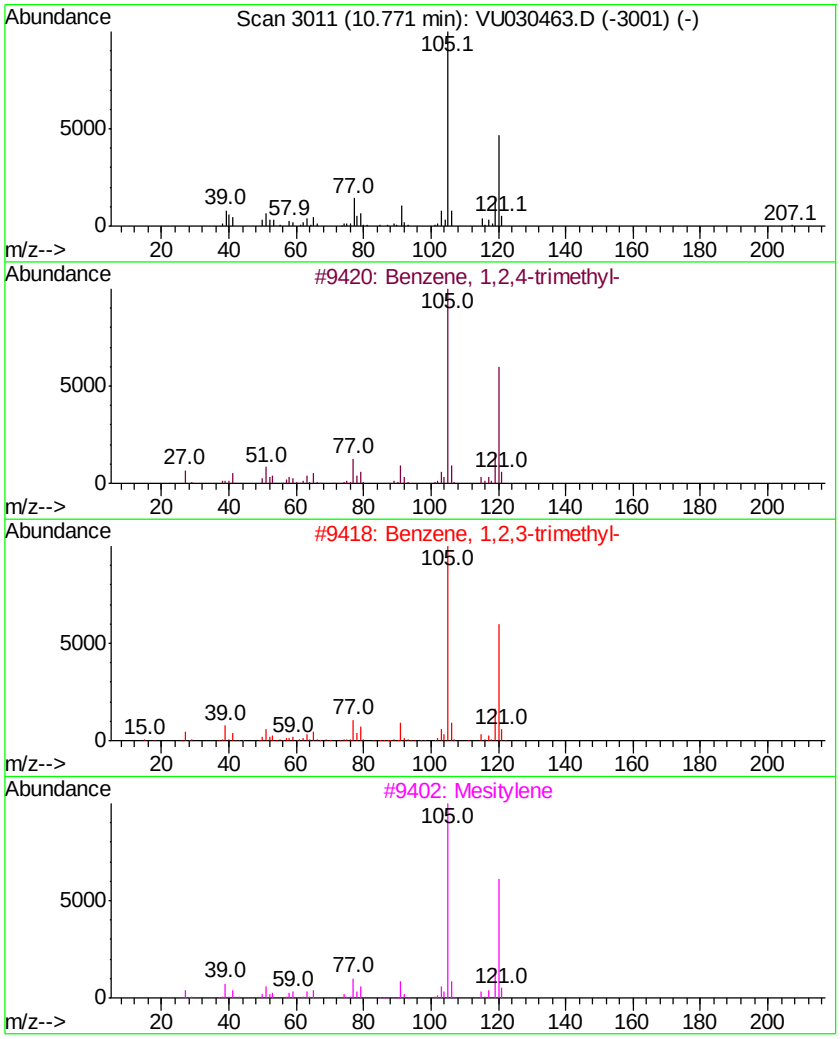
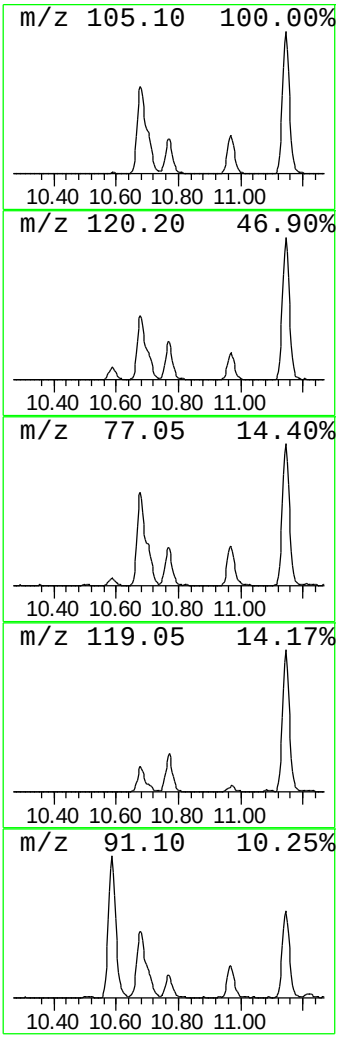
Peak Number 7 Benzene, 1,2,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	7.20 ug/L	178545	1,4-Dichlorobenzene-d4	11.48

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





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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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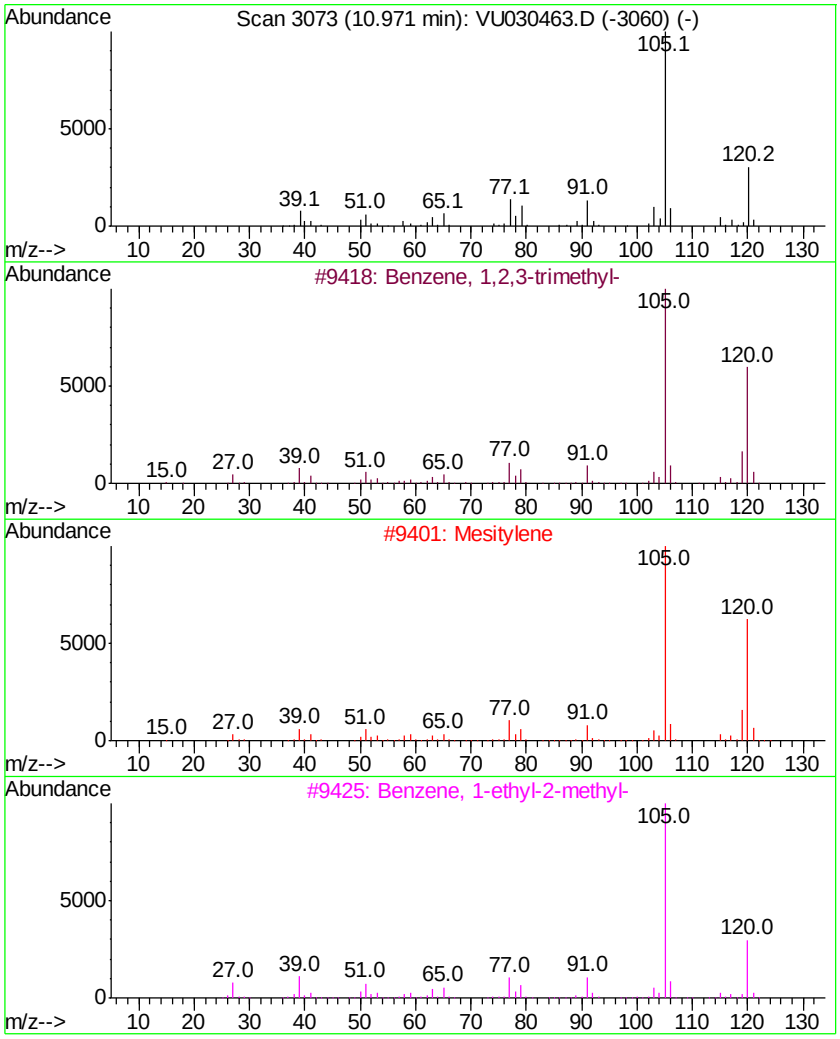
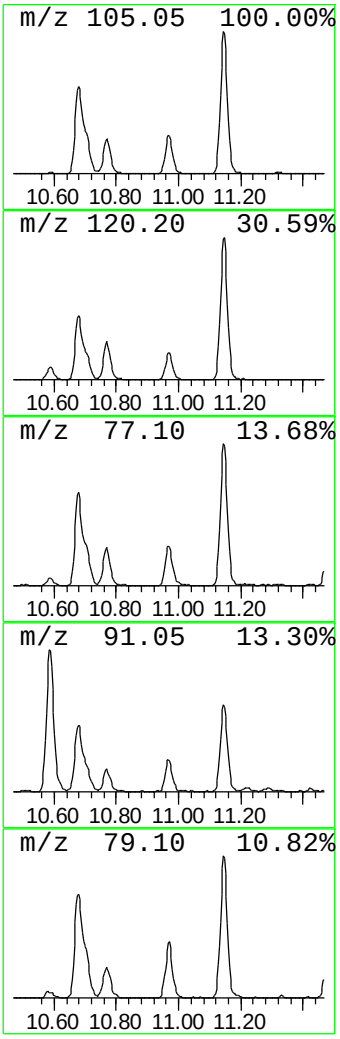
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	8.05 ug/L	199608	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	91
2	Mesitylene	120	C9H12	000108-67-8	91
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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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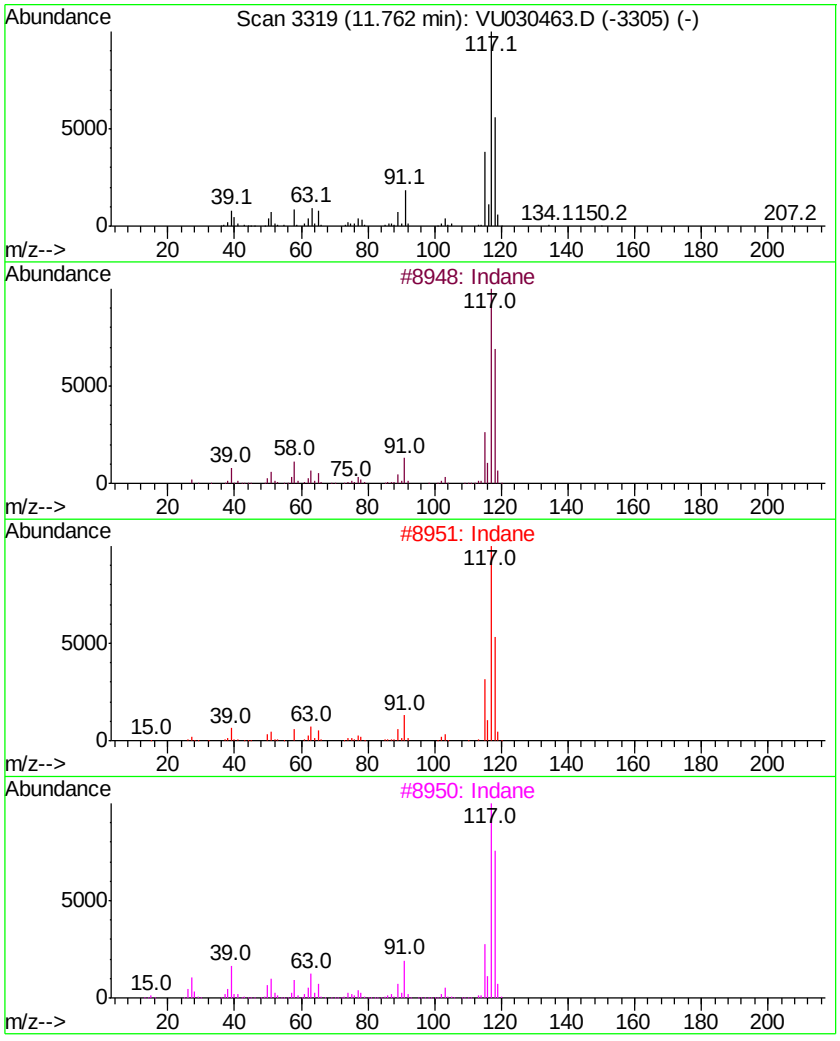
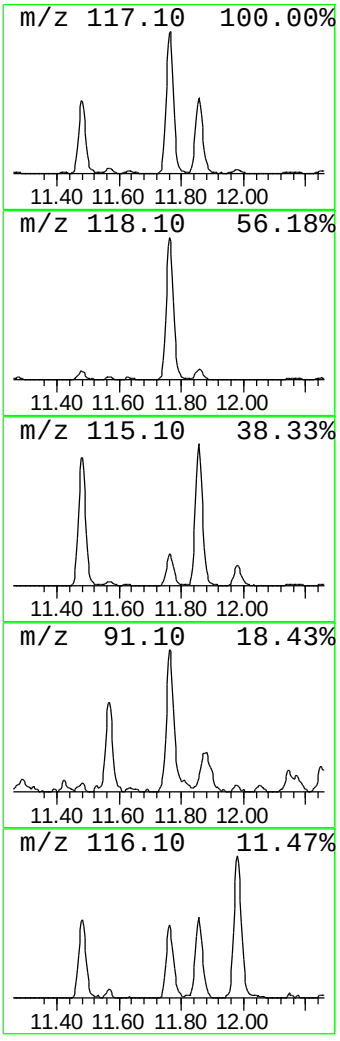
Peak Number 11 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	10.80 ug/L	267611	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	68
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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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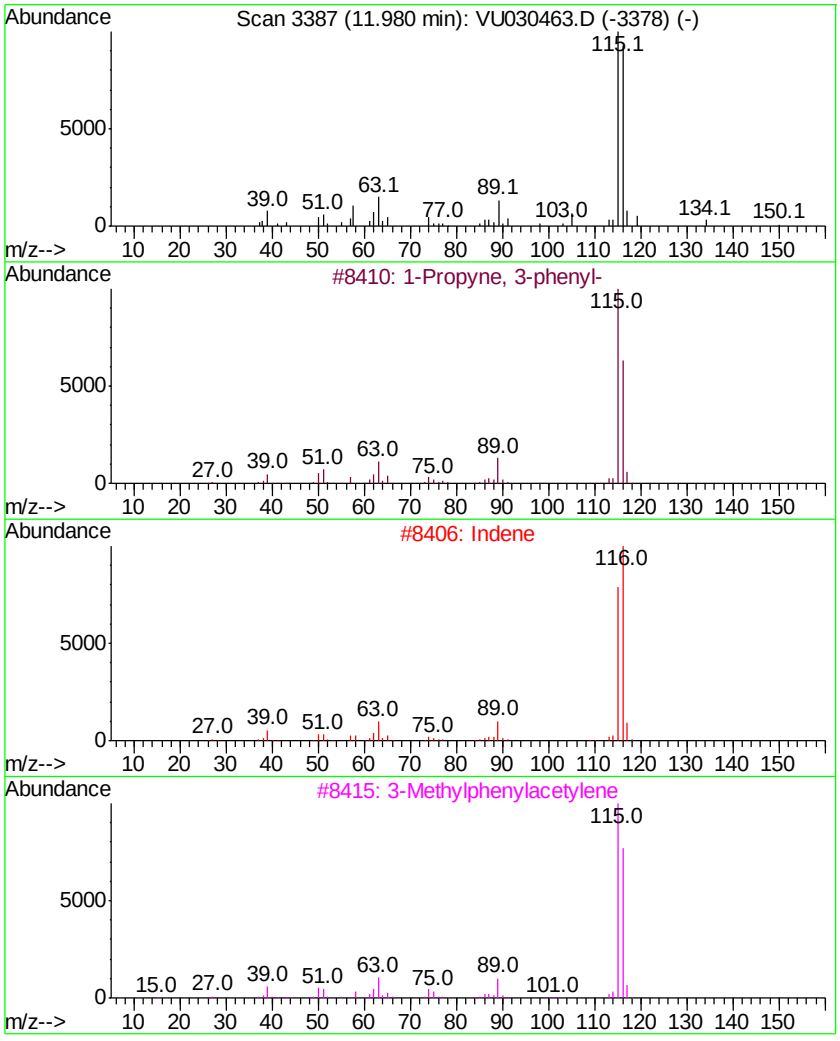
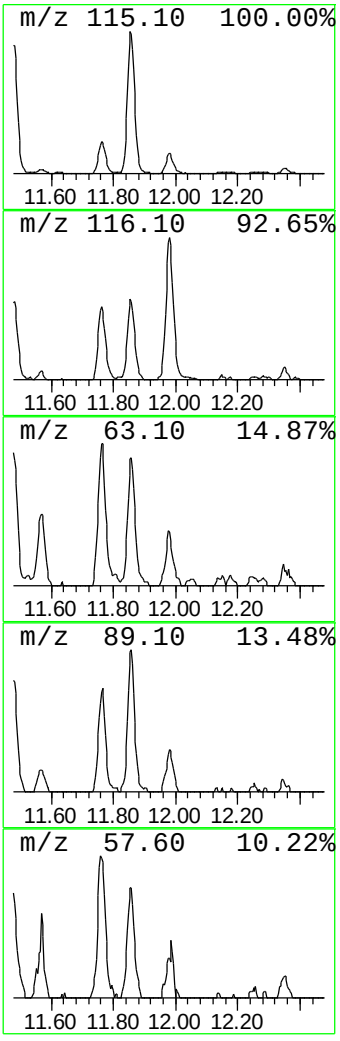
Peak Number 12 1-Propyne, 3-phenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	2.59 ug/L	64102	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
2	Indene	116	C9H8	000095-13-6	93
3	3-Methylphenylacetylene	116	C9H8	000766-82-5	93
4	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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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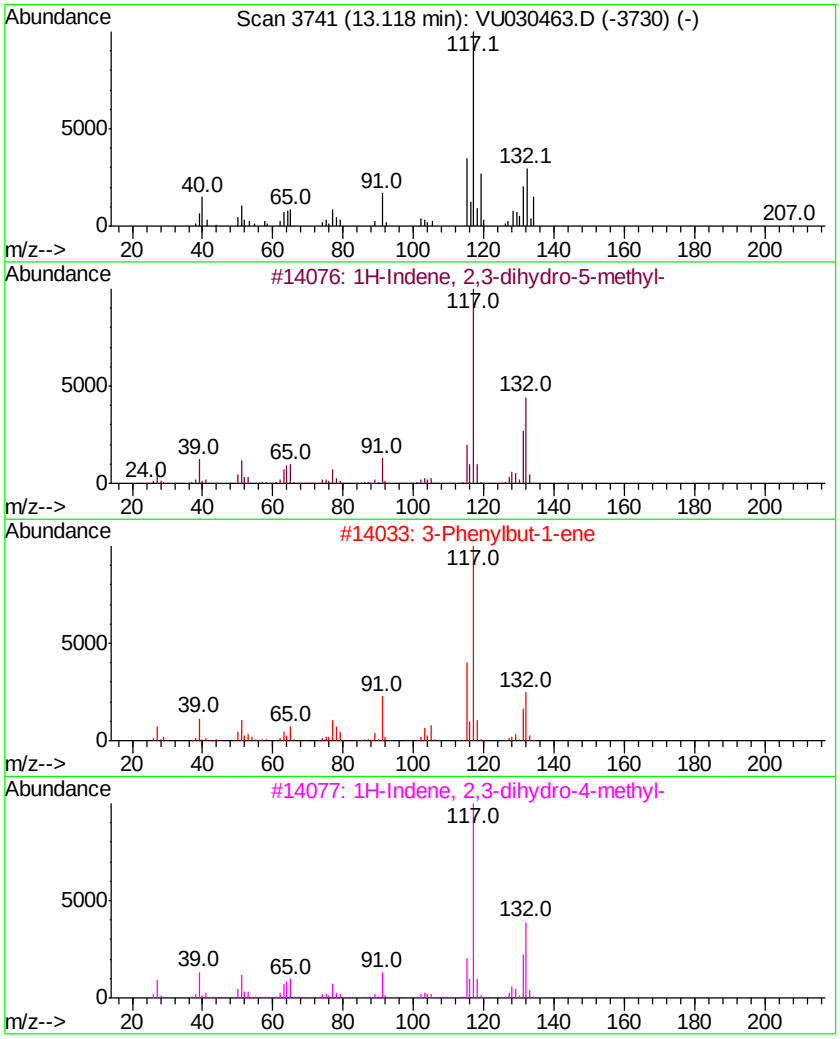
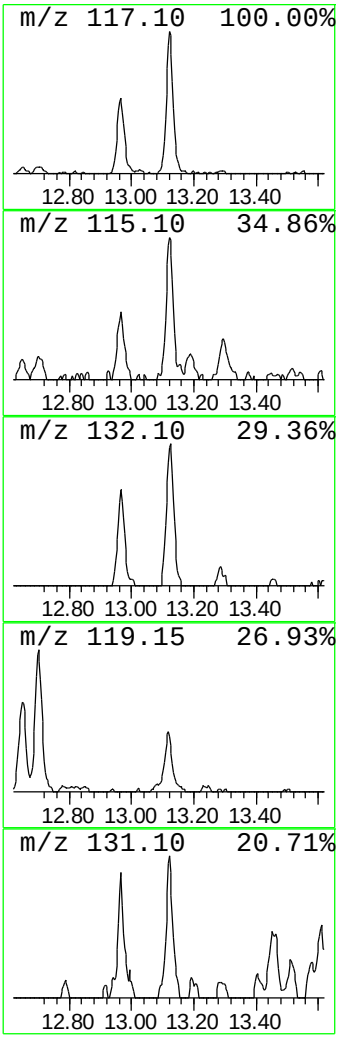
Peak Number 13 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	2.79 ug/L	69256	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
5	Indan, 1-methyl-	132	C10H12	000767-58-8	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030463.D
Acq On : 26 Mar 2019 23:02
Operator : JC/SP
Sample : K2063-08 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DB6R8

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
(DEL) Alkane: Str...	1.75	7.1	ug/L	153231	1	5.88	1082610	50.0
(DEL) Alkane: Str...	1.94	4.7	ug/L	101229	1	5.88	1082610	50.0
(DEL) Alkane: Cyc...	4.09	8.0	ug/L	174186	1	5.88	1082610	50.0
Benzene, propyl-	10.58	3.7	ug/L	92295	3	11.48	1239230	50.0
Benzene, 1-ethyl-...	10.68	23.2	ug/L	574580	3	11.48	1239230	50.0
Benzene, 1,2,4-tr...	10.77	7.2	ug/L	178545	3	11.48	1239230	50.0
Benzene, 1,2,3-tr...	10.97	8.1	ug/L	199608	3	11.48	1239230	50.0
Indane	11.76	10.8	ug/L	267611	3	11.48	1239230	50.0
1-Propyne, 3-phenyl-	11.98	2.6	ug/L	64102	3	11.48	1239230	50.0
1H-Indene, 2,3-di...	13.12	2.8	ug/L	69256	3	11.48	1239230	50.0