

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031519\
 Data File : VX008128.D
 Acq On : 16 Mar 2019 01:17
 Operator : JC/SP
 Sample : K1894-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RFK-RMAB-MW10-GW

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.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	2.843	269	277	285	rBV	26324	58147	1.49%	0.132%
2	5.501	702	713	726	rBV2	142136	409298	10.51%	0.932%
3	5.665	729	740	758	rVB	229606	666233	17.11%	1.517%
4	6.068	796	806	817	rBV	136405	359201	9.22%	0.818%
5	6.860	926	936	946	rBV	353756	827691	21.26%	1.885%
6	8.714	1233	1240	1249	rBV	760853	1154807	29.66%	2.630%
7	9.518	1363	1372	1377	rBV2	38970	62455	1.60%	0.142%
8	10.006	1446	1452	1459	rBV	46034	65688	1.69%	0.150%
9	10.110	1459	1469	1476	rBV	740509	1025933	26.35%	2.336%
10	10.250	1487	1492	1498	rBV	151650	195387	5.02%	0.445%
11	10.360	1503	1510	1516	rVB	45057	66688	1.71%	0.152%
12	10.695	1561	1565	1571	rVB	125089	163951	4.21%	0.373%
13	11.018	1612	1618	1624	rBV	81628	108139	2.78%	0.246%
14	11.134	1632	1637	1643	rBV	677882	853347	21.91%	1.943%
15	11.359	1666	1674	1678	rBV	68745	89786	2.31%	0.204%
16	11.451	1684	1689	1696	rBV	53756	71069	1.83%	0.162%
17	11.664	1719	1724	1732	rBV	779418	966248	24.81%	2.200%
18	11.804	1743	1747	1753	rVB	509942	603437	15.50%	1.374%
19	11.920	1761	1766	1768	rBV	26005	40643	1.04%	0.093%
20	12.024	1779	1783	1787	rBV	111448	127153	3.27%	0.290%
21	12.079	1787	1792	1796	rVV	827730	1046156	26.87%	2.382%
22	12.140	1796	1802	1807	rVB	286830	347457	8.92%	0.791%
23	12.231	1813	1817	1823	rVB	68150	83326	2.14%	0.190%
24	12.298	1823	1828	1835	rBV	680837	857002	22.01%	1.952%
25	12.365	1835	1839	1846	rVB	165055	233808	6.00%	0.532%
26	12.457	1850	1854	1859	rVV2	269704	354942	9.12%	0.808%
27	12.518	1859	1864	1870	rVB	430056	561918	14.43%	1.280%
28	12.585	1870	1875	1877	rBV	925425	1134675	29.14%	2.584%
29	12.609	1877	1879	1884	rVV	784080	833303	21.40%	1.898%
30	12.670	1884	1889	1893	rVV	2473320	3037445	78.00%	6.917%
31	12.701	1893	1894	1898	rVV	105573	90307	2.32%	0.206%
32	12.755	1898	1903	1911	rVV2	1223472	1921983	49.36%	4.377%
33	12.853	1915	1919	1922	rBV	69749	82368	2.12%	0.188%
34	12.902	1922	1927	1932	rVB2	511058	654690	16.81%	1.491%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
 Title : SW846 8260

35	12.975	1932	1939	1943	rBV	2357930	2929900	75.24%	6.672%
36	13.018	1943	1946	1951	rVB	3193026	3646837	93.65%	8.305%
37	13.103	1952	1960	1965	rBV2	399502	719119	18.47%	1.638%
38	13.152	1965	1968	1970	rVV	50790	63183	1.62%	0.144%
39	13.194	1971	1975	1977	rVV2	235897	315080	8.09%	0.718%
40	13.225	1977	1980	1983	rVV	566023	697593	17.91%	1.589%
41	13.261	1983	1986	1990	rVV3	173078	267962	6.88%	0.610%
42	13.310	1990	1994	1996	rVV2	397698	561304	14.41%	1.278%
43	13.347	1996	2000	2006	rVV2	2709523	3894037	100.00%	8.868%
44	13.426	2010	2013	2018	rVV	252603	339900	8.73%	0.774%
45	13.481	2018	2022	2028	rVB	331849	436987	11.22%	0.995%
46	13.554	2029	2034	2038	rBV2	189299	304828	7.83%	0.694%
47	13.597	2038	2041	2044	rVV	528062	731445	18.78%	1.666%
48	13.639	2044	2048	2054	rVV3	935620	1931732	49.61%	4.399%
49	13.700	2054	2058	2059	rVV	492979	654958	16.82%	1.492%
50	13.725	2059	2062	2067	rVV2	760192	1187149	30.49%	2.703%
51	13.774	2067	2070	2073	rVB	48028	51798	1.33%	0.118%
52	13.810	2073	2076	2078	rBV	44986	54588	1.40%	0.124%
53	13.853	2078	2083	2088	rVB3	130846	234619	6.03%	0.534%
54	13.914	2088	2093	2098	rVB4	115065	188112	4.83%	0.428%
55	13.987	2098	2105	2108	rBV2	313021	464516	11.93%	1.058%
56	14.030	2108	2112	2116	rVV	279462	333927	8.58%	0.760%
57	14.060	2116	2117	2124	rVB3	50083	63362	1.63%	0.144%
58	14.133	2124	2129	2134	rBV	597551	736150	18.90%	1.676%
59	14.188	2134	2138	2141	rBV2	27205	47674	1.22%	0.109%
60	14.274	2146	2152	2158	rBV	565033	888080	22.81%	2.022%
61	14.377	2165	2169	2174	rVB	389822	432172	11.10%	0.984%
62	14.432	2174	2178	2182	rVB	429636	539470	13.85%	1.229%
63	14.475	2182	2185	2191	rVB	43511	62527	1.61%	0.142%
64	14.542	2191	2196	2200	rBV2	156774	232615	5.97%	0.530%
65	14.578	2200	2202	2208	rVB2	28694	39165	1.01%	0.089%
66	14.664	2208	2216	2222	rBV3	116680	266213	6.84%	0.606%
67	14.731	2225	2227	2231	rVB	60996	70301	1.81%	0.160%
68	14.786	2231	2236	2239	rBV	44406	59649	1.53%	0.136%
69	14.840	2240	2245	2254	rBV	587143	814726	20.92%	1.855%
70	15.249	2306	2312	2320	rBV	75463	139370	3.58%	0.317%
71	15.481	2346	2350	2356	rVB3	26305	39089	1.00%	0.089%

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ClientSampleId :
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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

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72	15.688	2379	2384	2387	rVV	118738	180055	4.62%	0.410%
73	15.724	2387	2390	2397	rVB	52335	82488	2.12%	0.188%
74	15.889	2412	2417	2420	rBV	30949	54640	1.40%	0.124%

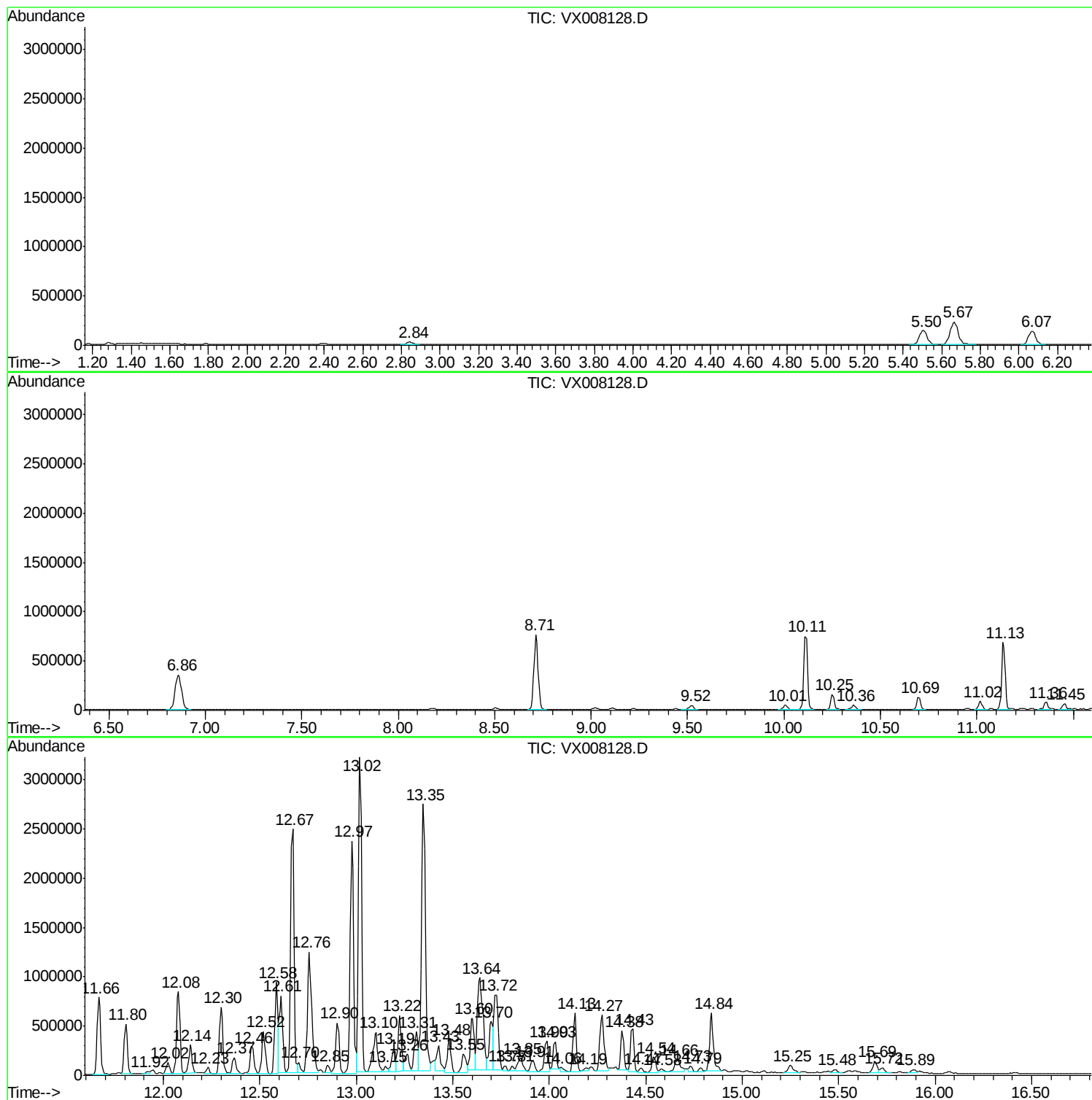
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Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
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RFK-RMAB-MW10-GW

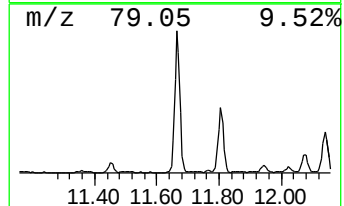
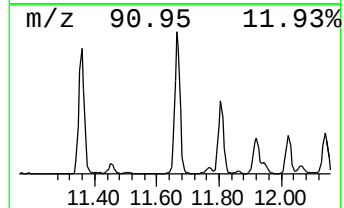
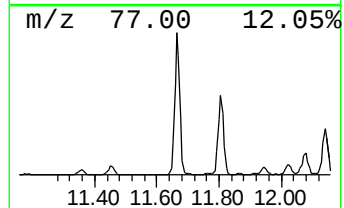
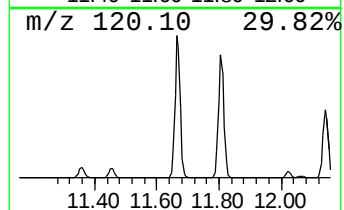
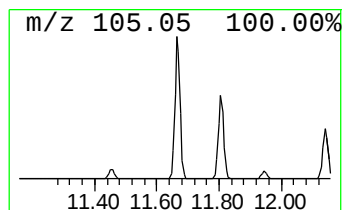
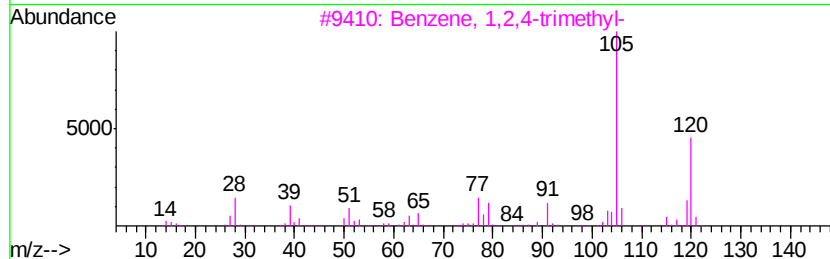
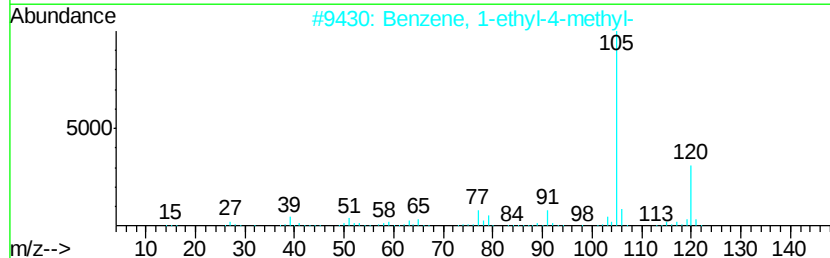
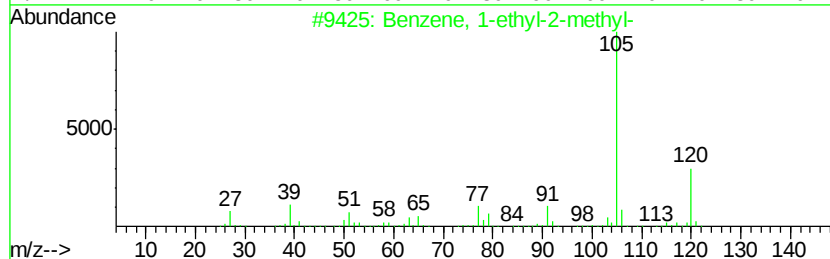
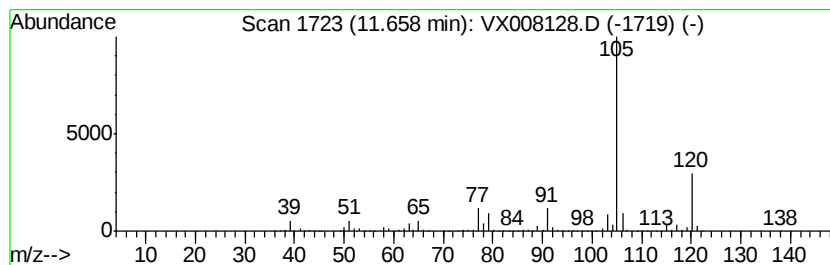
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	46.18 ug/l	966248	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5		Mesitylene	120	C9H12	000108-67-8	90



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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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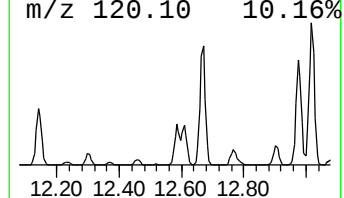
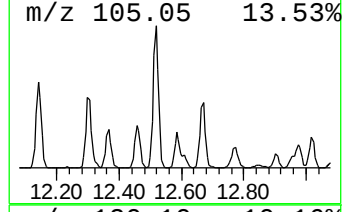
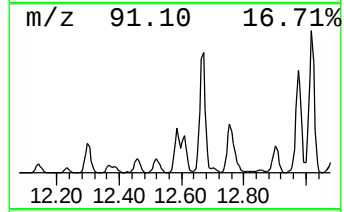
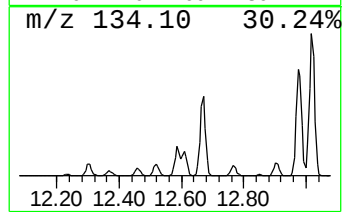
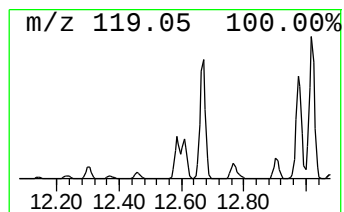
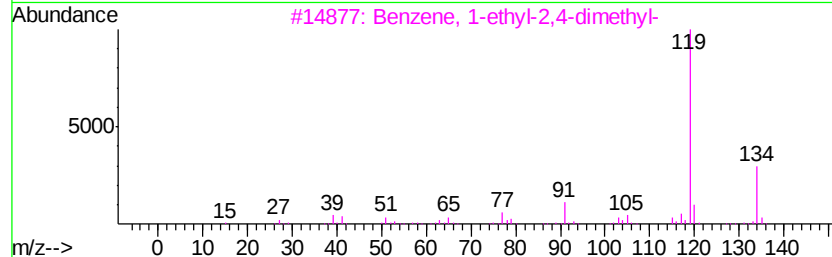
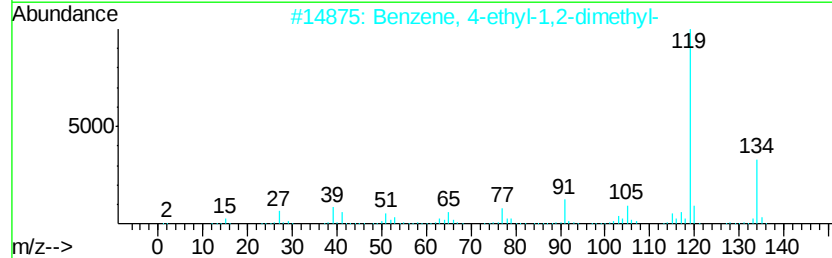
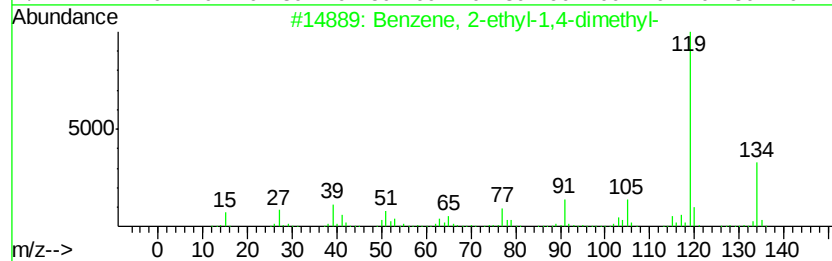
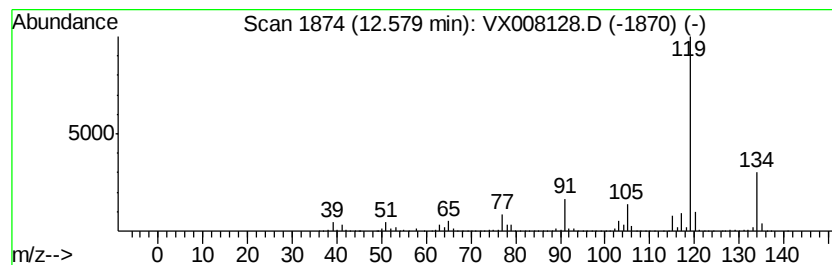
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	54.23 ug/l	1134680	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



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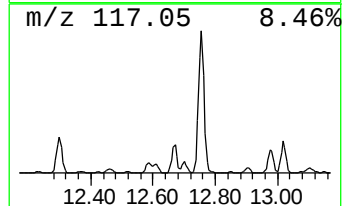
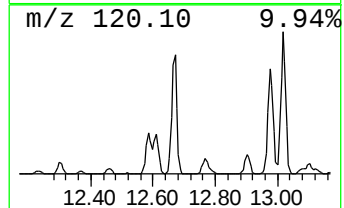
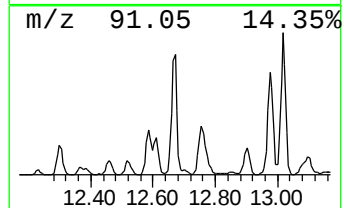
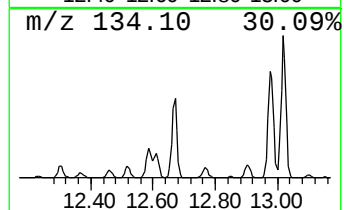
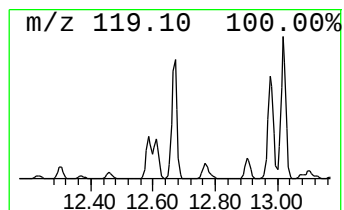
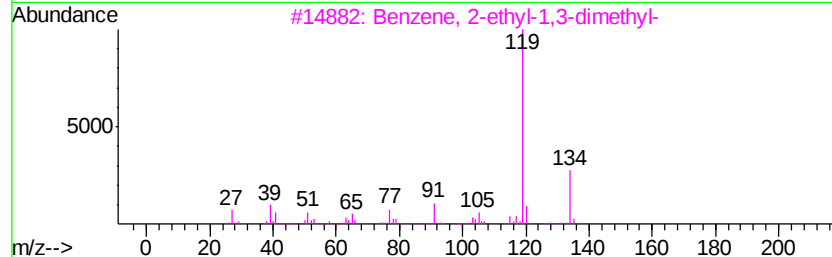
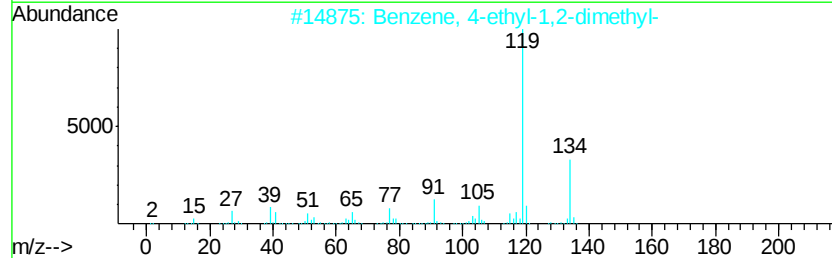
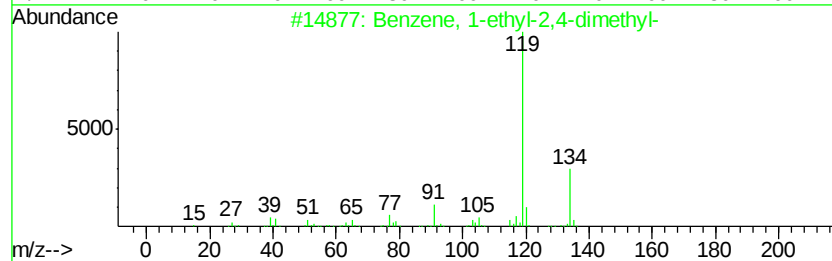
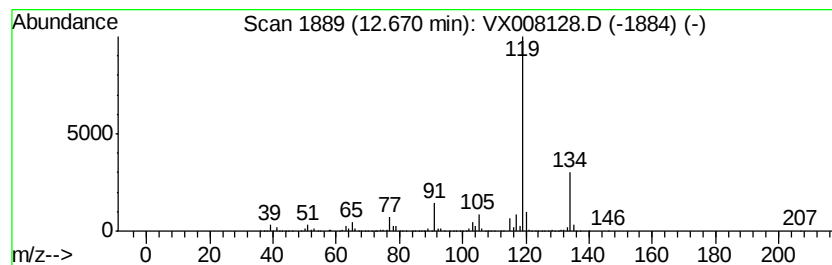
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	145.17 ug/l	3037450	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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ClientSampleId :
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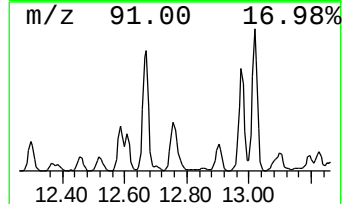
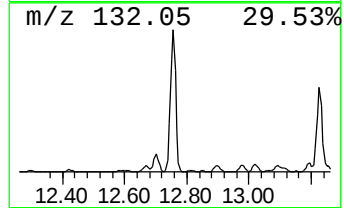
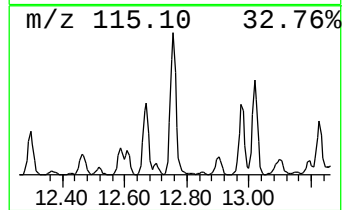
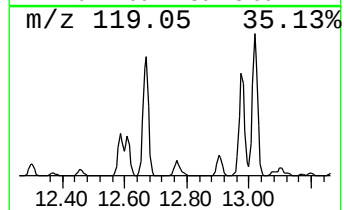
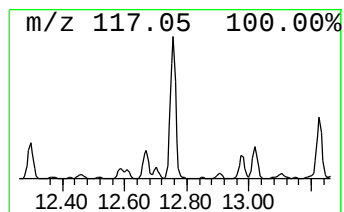
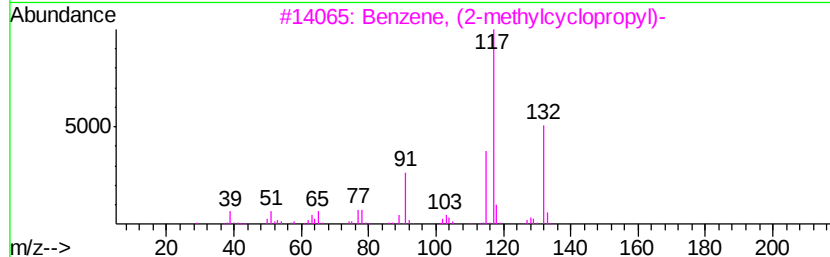
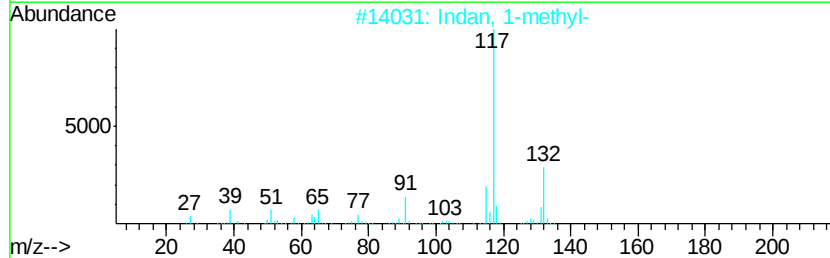
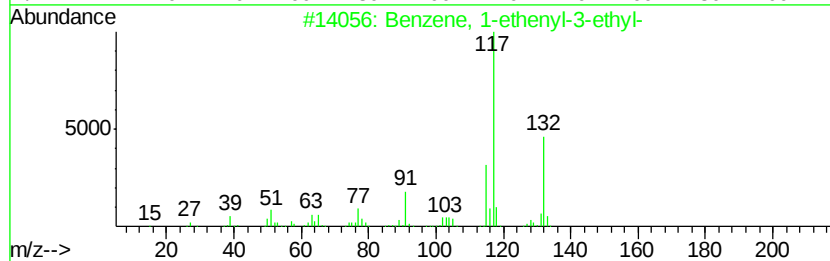
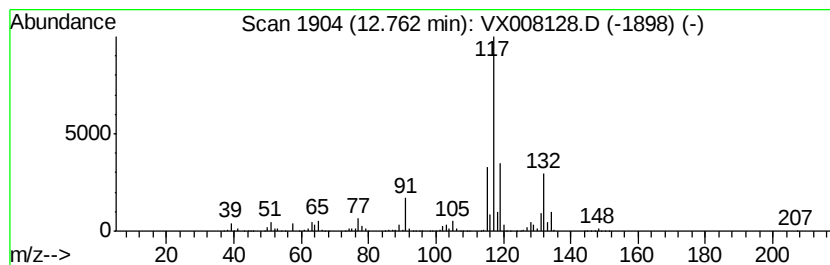
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	91.86 ug/l	1921980	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

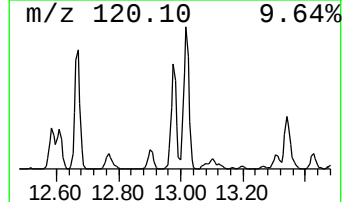
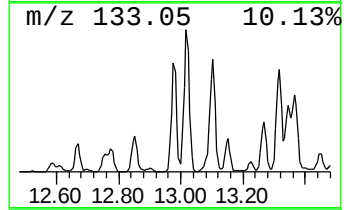
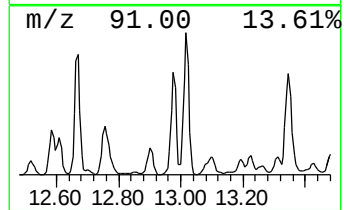
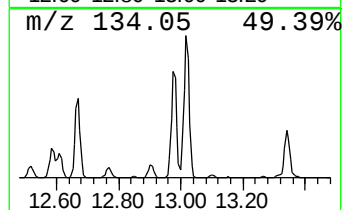
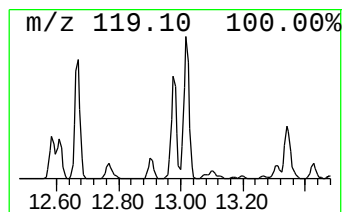
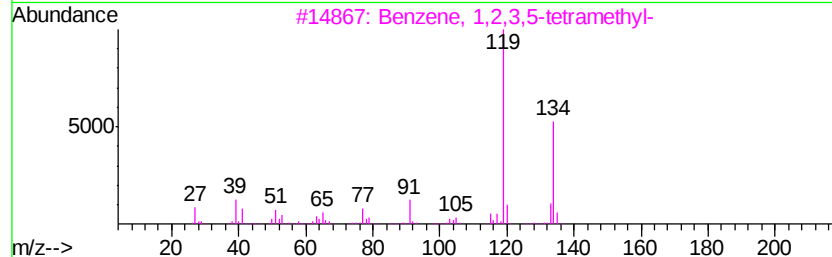
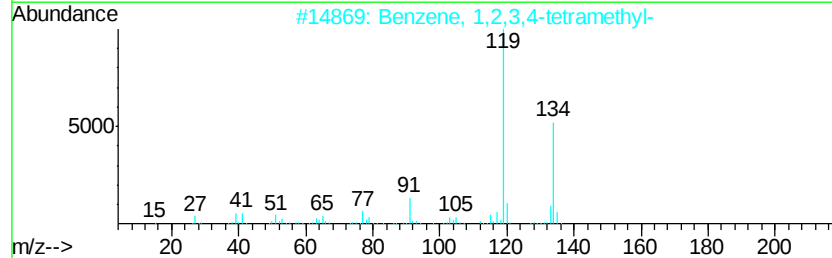
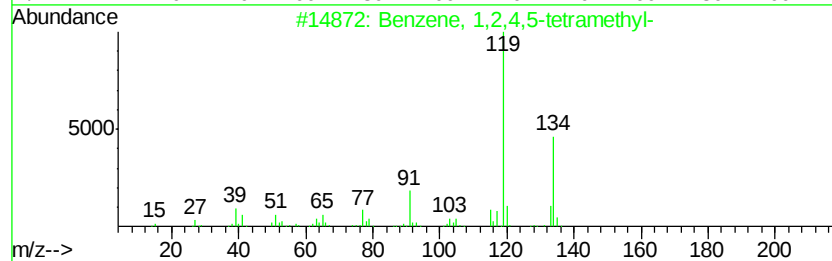
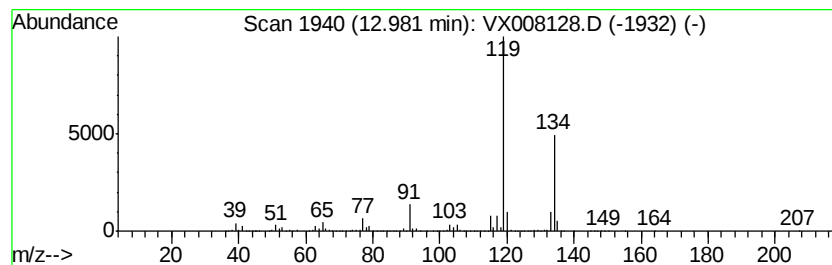
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MSVOA_X
ClientSampleId :
RFK-RMAB-MW10-GW

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.98	140.03 ug/l	2929900	1,4-Dichlorobenzene-d4	12.08	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RFK-RMAB-MW10-GW

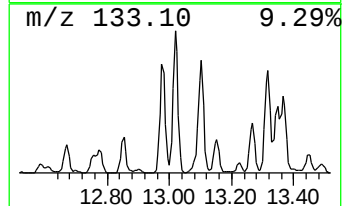
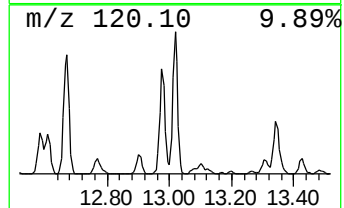
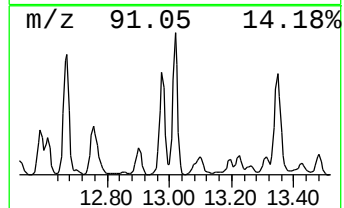
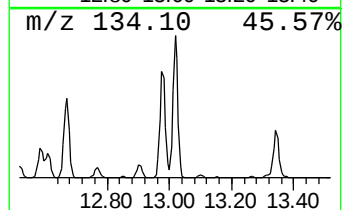
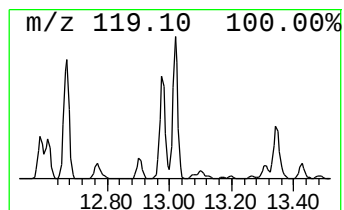
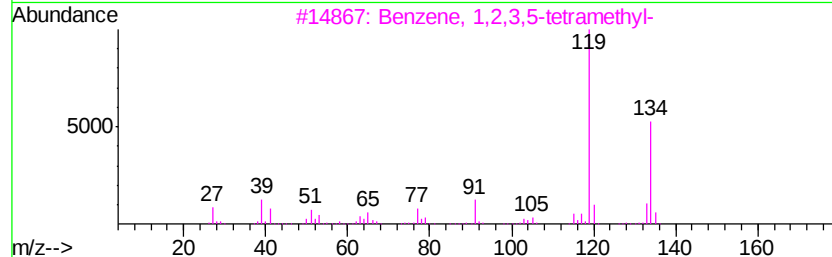
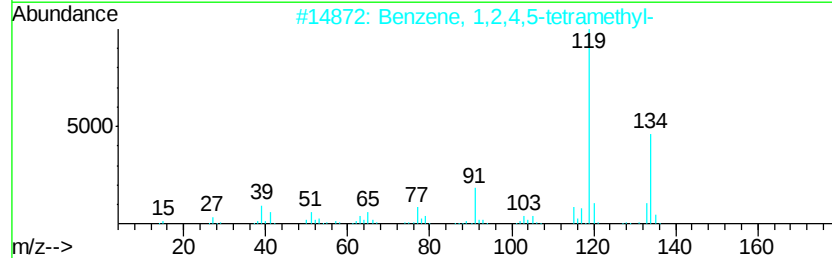
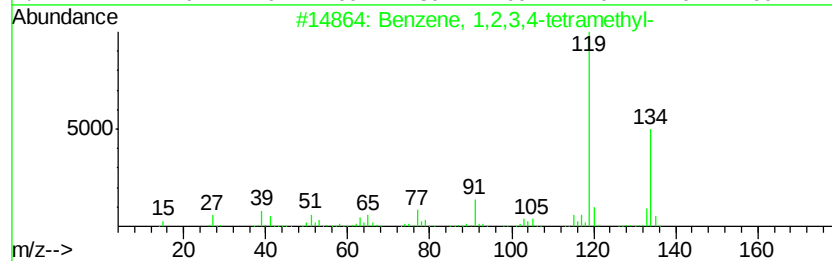
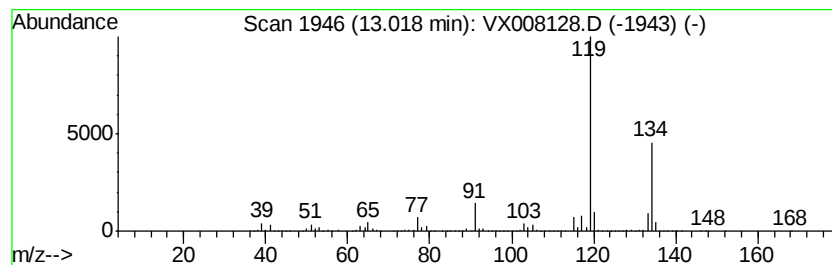
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	174.30 ug/l	3646840	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4		o-Cymene	134	C10H14	000527-84-4	96
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RFK-RMAB-MW10-GW

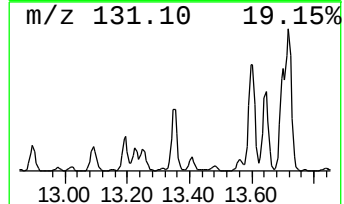
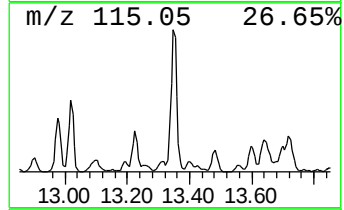
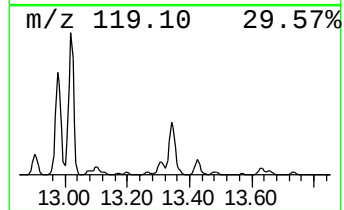
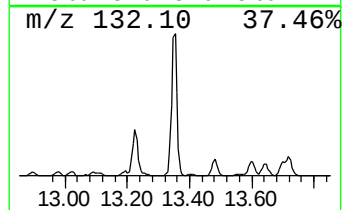
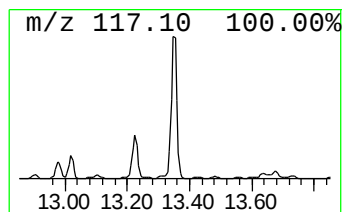
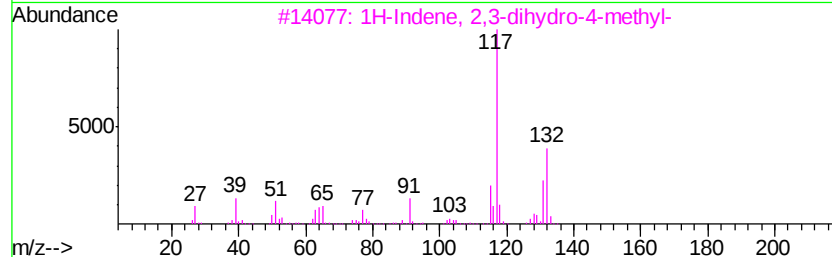
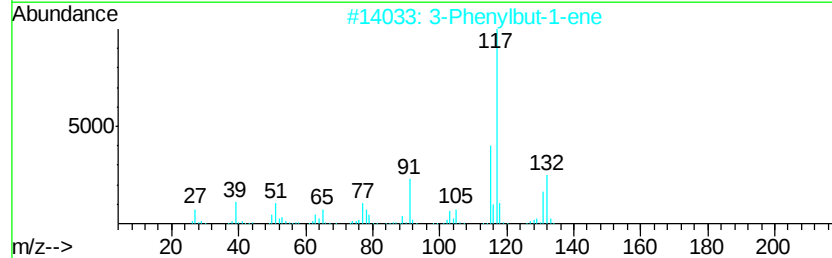
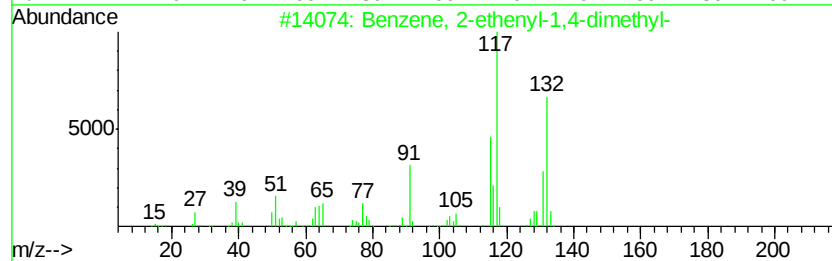
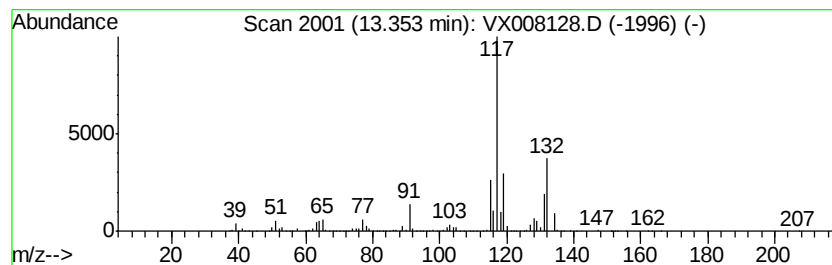
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	186.11 ug/l	3894040	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
4		Indan, 1-methyl-	132	C10H12	000767-58-8	64
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RFK-RMAB-MW10-GW

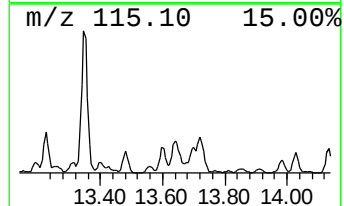
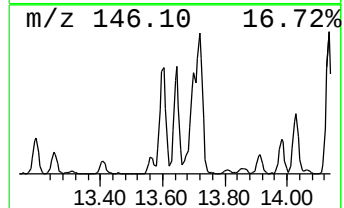
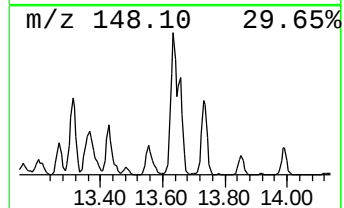
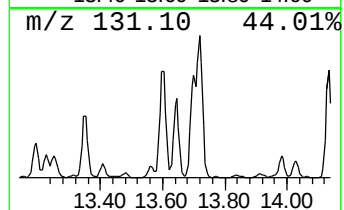
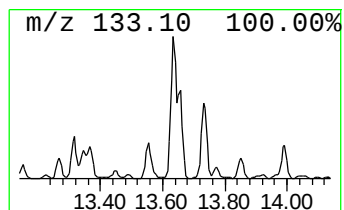
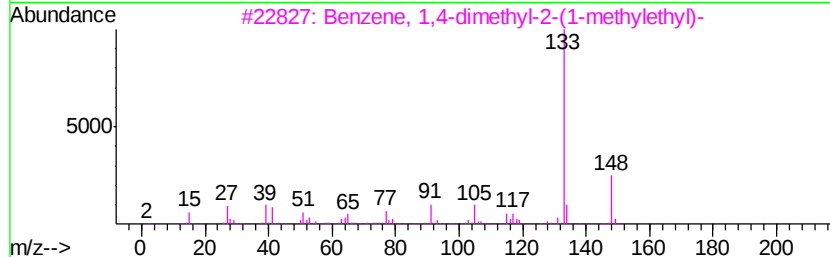
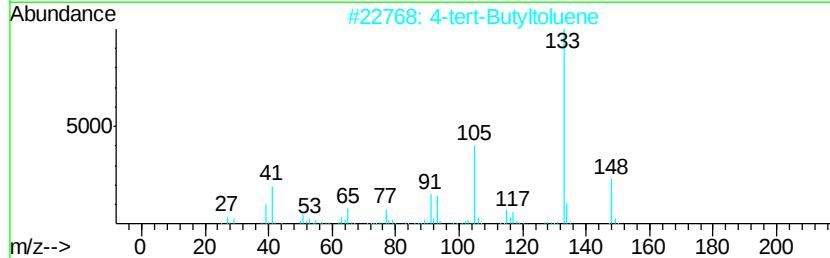
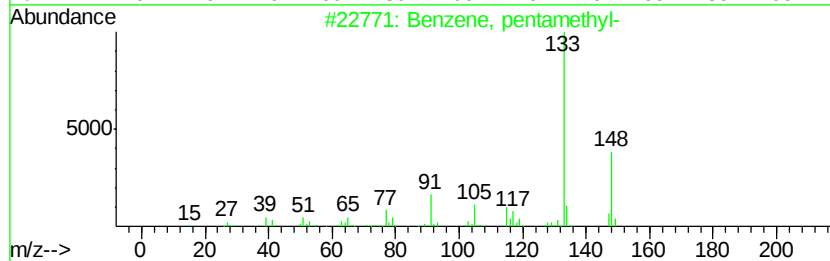
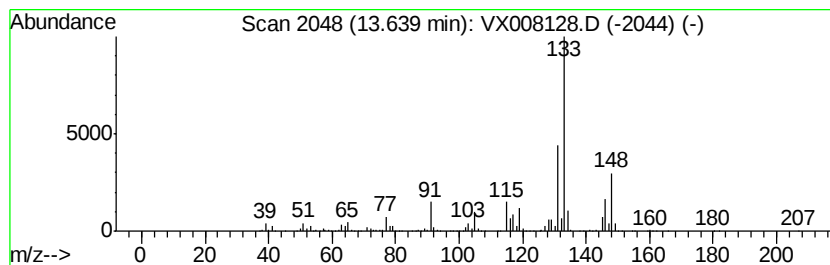
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, pentamethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	92.33 ug/l	1931730	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, pentamethyl-	148	C11H16	000700-12-9	64
2		4-tert-Butyltoluene	148	C11H16	000098-51-1	53
3		Benzene, 1,4-dimethyl-2-(1-methyl-...	148	C11H16	004132-72-3	53
4		1,5,6,7-Tetramethylbicyclo[3.2.0]...	148	C11H16	134329-46-7	52
5		Benzene, 1,3-dimethyl-5-(1-methyl-...	148	C11H16	004706-90-5	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RFK-RMAB-MW10-GW

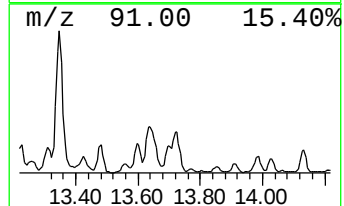
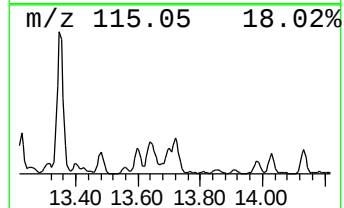
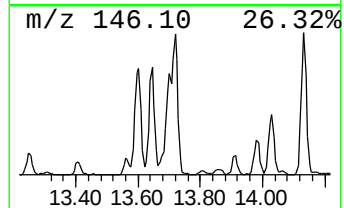
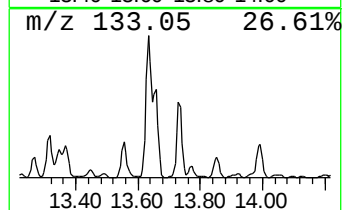
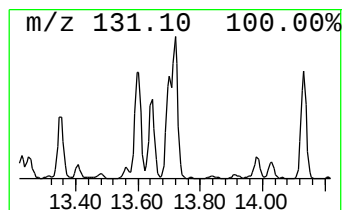
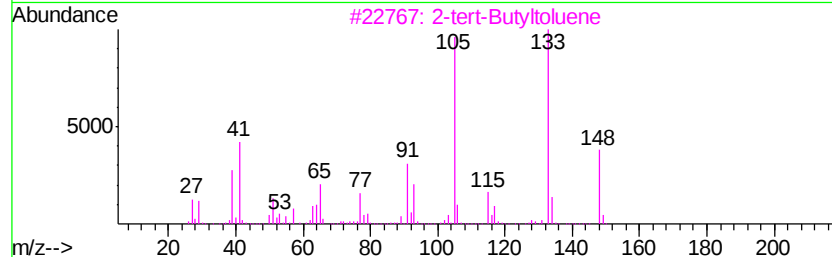
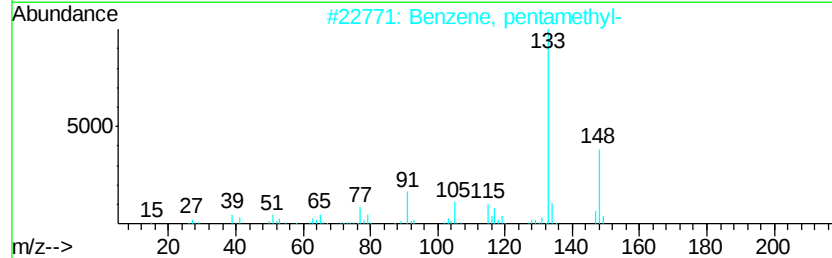
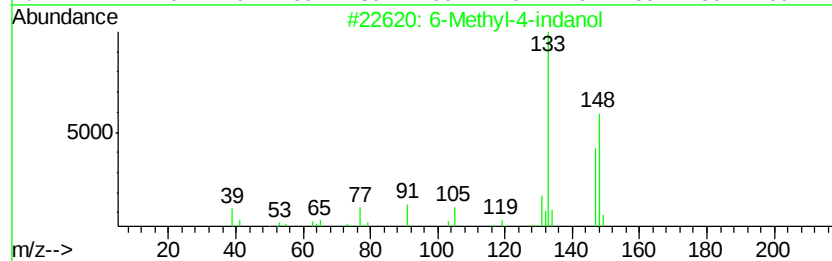
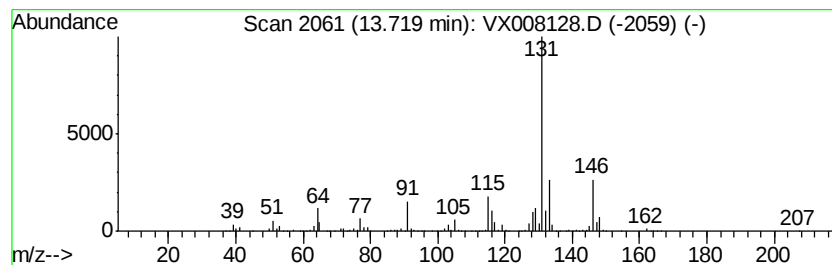
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Quant Title : SW846 8260

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

Peak Number 9 6-Methyl-4-indanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	56.74 ug/l	1187150	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Methyl-4-indanol	148	C10H12O	020294-32-0	68
2		Benzene, pentamethyl-	148	C11H16	000700-12-9	64
3		2-tert-Butyltoluene	148	C11H16	001074-92-6	58
4		4-tert-Butyltoluene	148	C11H16	000098-51-1	53
5		Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RFK-RMAB-MW10-GW

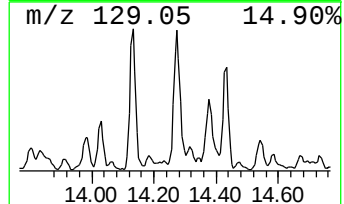
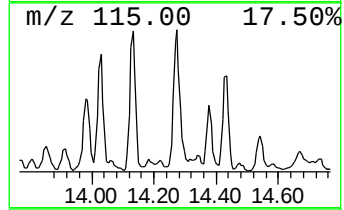
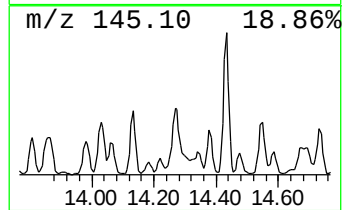
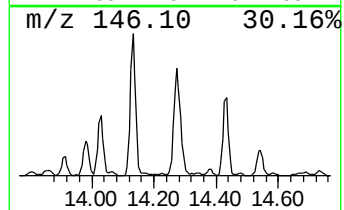
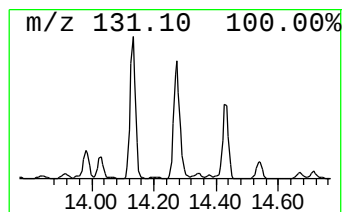
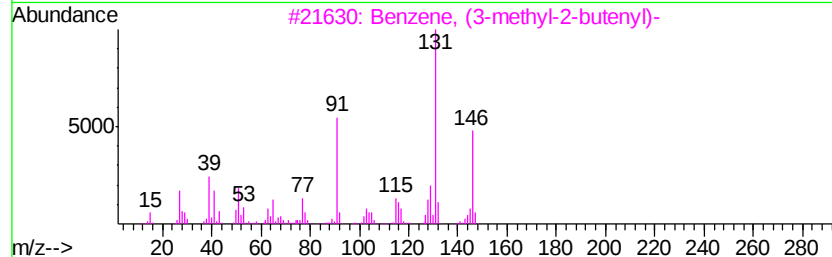
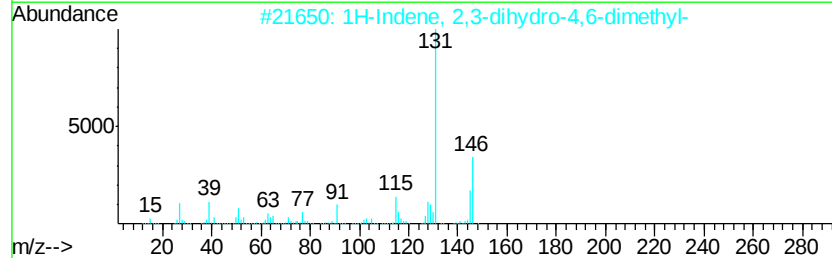
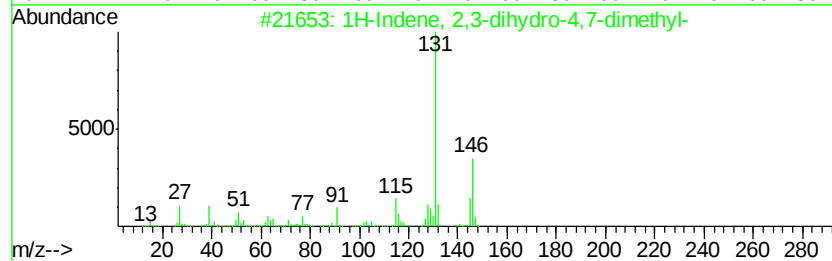
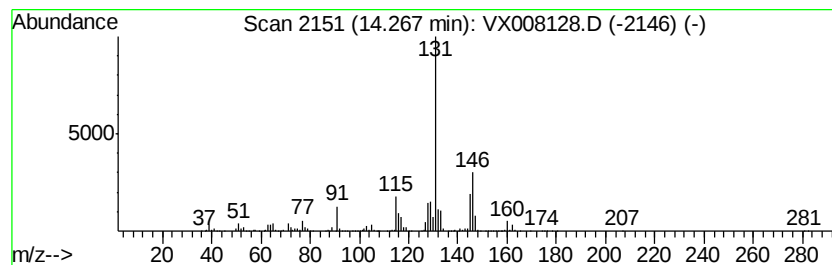
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	42.44 ug/l	888080	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	94
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	93
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
5		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031519\
Data File : VX008128.D
Acq On : 16 Mar 2019 01:17
Operator : JC/SP
Sample : K1894-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RFK-RMAB-MW10-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	11.66	46.2	ug/l	966248	4	12.08	1046160	50.0
Benzene, 2-ethyl-...	12.58	54.2	ug/l	1134680	4	12.08	1046160	50.0
Benzene, 1-ethyl-...	12.67	145.2	ug/l	3037450	4	12.08	1046160	50.0
Benzene, 1-etheny...	12.76	91.9	ug/l	1921980	4	12.08	1046160	50.0
Benzene, 1,2,4,5-...	12.98	140.0	ug/l	2929900	4	12.08	1046160	50.0
Benzene, 1,2,3,4-...	13.02	174.3	ug/l	3646840	4	12.08	1046160	50.0
Benzene, 2-etheny...	13.35	186.1	ug/l	3894040	4	12.08	1046160	50.0
Benzene, pentamet...	13.64	92.3	ug/l	1931730	4	12.08	1046160	50.0
6-Methyl-4-indanol	13.72	56.7	ug/l	1187150	4	12.08	1046160	50.0
1H-Indene, 2,3-di...	14.27	42.4	ug/l	888080	4	12.08	1046160	50.0