

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062519\
 Data File : VX010428.D
 Acq On : 25 Jun 2019 10:58
 Operator : JC/SP
 Sample : K3500-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GW-B-061919

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\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	1.788	99	104	117	rVB	489686	715656	6.97%	0.579%
2	2.001	133	139	147	rVB	208492	296316	2.89%	0.240%
3	2.843	270	277	280	rBV	53796	105407	1.03%	0.085%
4	2.891	280	285	295	rVB	111568	237984	2.32%	0.193%
5	3.172	323	331	345	rVB	80815	186510	1.82%	0.151%
6	4.403	521	533	546	rVB	45572	133603	1.30%	0.108%
7	5.501	700	713	721	rBV	153729	444216	4.33%	0.360%
8	5.665	731	740	759	rVB	257557	751014	7.32%	0.608%
9	6.062	796	805	819	rBV	143770	371159	3.62%	0.301%
10	6.860	924	936	946	rBV	380081	905636	8.83%	0.733%
11	7.464	1022	1035	1043	rBV3	59469	149674	1.46%	0.121%
12	8.714	1234	1240	1249	rVB	814031	1260476	12.28%	1.021%
13	10.110	1464	1469	1475	rBV	789838	1054861	10.28%	0.854%
14	11.018	1612	1618	1627	rBV	348148	458083	4.46%	0.371%
15	11.134	1632	1637	1642	rBV	685815	872865	8.51%	0.707%
16	11.359	1668	1674	1681	rBV	1997196	2480494	24.18%	2.009%
17	11.433	1681	1686	1693	rVV	4075282	7617878	74.24%	6.168%
18	11.506	1693	1698	1707	rVB	3150245	3850550	37.53%	3.118%
19	11.664	1719	1724	1732	rVB	1463093	1798845	17.53%	1.457%
20	11.804	1742	1747	1753	rVB	8369828	10260498	100.00%	8.308%
21	11.920	1760	1766	1768	rBV	344375	508173	4.95%	0.411%
22	11.945	1768	1770	1775	rVB	441999	473649	4.62%	0.384%
23	12.024	1775	1783	1786	rBV	926607	1116351	10.88%	0.904%
24	12.073	1786	1791	1797	rVB2	914132	1456379	14.19%	1.179%
25	12.140	1797	1802	1808	rBV	978227	1180048	11.50%	0.956%
26	12.231	1812	1817	1822	rVB	104828	123983	1.21%	0.100%
27	12.304	1822	1829	1830	rBV	2444212	3214109	31.33%	2.603%
28	12.323	1830	1832	1835	rVV	3267315	3351264	32.66%	2.714%
29	12.371	1835	1840	1847	rVB3	6543239	9367300	91.29%	7.585%
30	12.457	1850	1854	1859	rVB	459203	554316	5.40%	0.449%
31	12.518	1859	1864	1870	rBV	1845098	2238534	21.82%	1.813%
32	12.585	1870	1875	1877	rBV	3717027	4643462	45.26%	3.760%
33	12.609	1877	1879	1884	rVB	3569658	3713029	36.19%	3.007%
34	12.670	1884	1889	1892	rBV	5984864	7407824	72.20%	5.998%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
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35	12.701	1892	1894	1898	rVB	636503	668605	6.52%	0.541%
36	12.756	1898	1903	1911	rBV2	1801452	3099740	30.21%	2.510%
37	12.847	1914	1918	1922	rVB2	178306	241449	2.35%	0.196%
38	12.902	1922	1927	1932	rVB	1281366	1575653	15.36%	1.276%
39	12.975	1932	1939	1943	rBV	3586455	4496716	43.83%	3.641%
40	13.018	1943	1946	1951	rVB	5115050	5794101	56.47%	4.692%
41	13.079	1951	1956	1957	rBV	592655	671734	6.55%	0.544%
42	13.097	1957	1959	1962	rVV2	578675	775856	7.56%	0.628%
43	13.121	1962	1963	1966	rVB	418892	341649	3.33%	0.277%
44	13.225	1972	1980	1984	rBV2	2301218	3315115	32.31%	2.684%
45	13.268	1984	1987	1990	rVB2	386319	456790	4.45%	0.370%
46	13.310	1990	1994	1996	rBV2	715888	1007136	9.82%	0.815%
47	13.347	1996	2000	2009	rVV2	4269076	6262477	61.03%	5.071%
48	13.426	2009	2013	2018	rVB	561555	720935	7.03%	0.584%
49	13.481	2018	2022	2029	rVB	334839	457843	4.46%	0.371%
50	13.560	2029	2035	2038	rBV2	618590	883587	8.61%	0.715%
51	13.597	2038	2041	2044	rVV	1195124	1601274	15.61%	1.297%
52	13.633	2044	2047	2054	rVV3	1557245	3144959	30.65%	2.547%
53	13.694	2054	2057	2058	rVV2	546384	620496	6.05%	0.502%
54	13.719	2058	2061	2067	rVB2	1257679	2025506	19.74%	1.640%
55	13.853	2077	2083	2089	rVB3	144930	220769	2.15%	0.179%
56	13.908	2089	2092	2097	rVB2	144994	215513	2.10%	0.175%
57	13.981	2097	2104	2108	rBV2	677036	989571	9.64%	0.801%
58	14.024	2108	2111	2115	rVV	514982	633085	6.17%	0.513%
59	14.133	2124	2129	2134	rBV	1055543	1358834	13.24%	1.100%
60	14.267	2147	2151	2161	rBV	899389	1541202	15.02%	1.248%
61	14.377	2165	2169	2174	rVV	491043	667830	6.51%	0.541%
62	14.426	2174	2177	2182	rVB	635853	790165	7.70%	0.640%
63	14.536	2190	2195	2199	rBV2	154910	244622	2.38%	0.198%
64	14.694	2209	2221	2226	rBV	1093390	1517877	14.79%	1.229%
65	14.834	2239	2244	2254	rVB	1714416	2260257	22.03%	1.830%
66	15.267	2307	2315	2320	rBV2	114075	207375	2.02%	0.168%
67	15.444	2332	2344	2347	rBV	118391	206959	2.02%	0.168%
68	15.548	2354	2361	2373	rVB	238426	412289	4.02%	0.334%
69	15.682	2373	2383	2387	rVV	258879	437402	4.26%	0.354%
70	15.725	2387	2390	2398	rVB	127346	187694	1.83%	0.152%
71	15.907	2411	2420	2432	rVB2	55679	146493	1.43%	0.119%

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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

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
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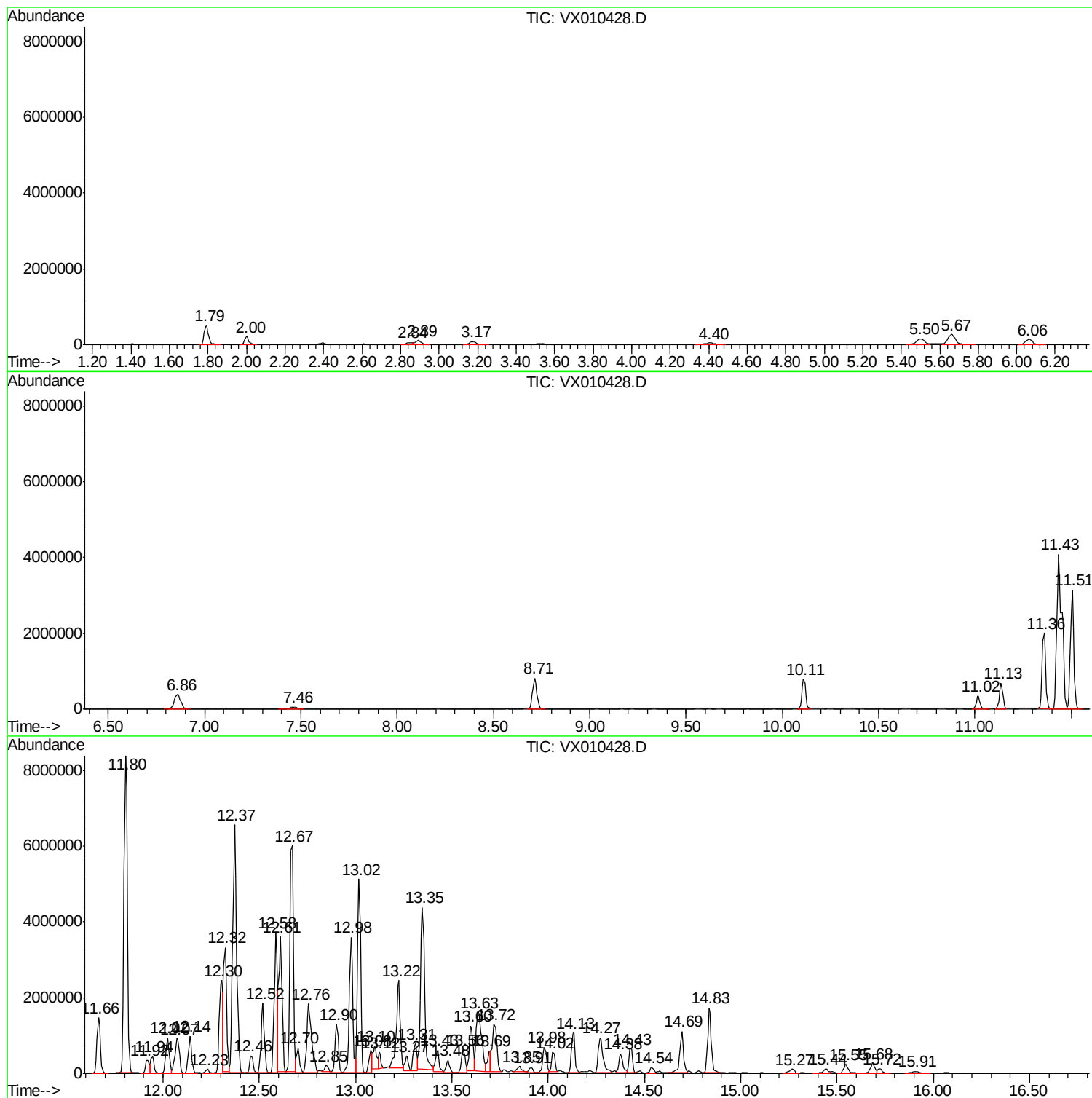
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Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

Quant Method : Z:\V0ASRV\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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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
GW-B-061919

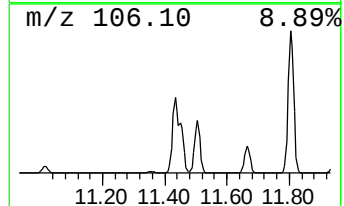
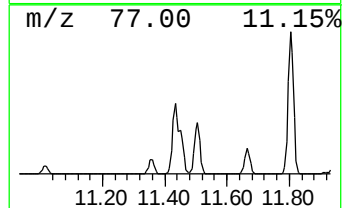
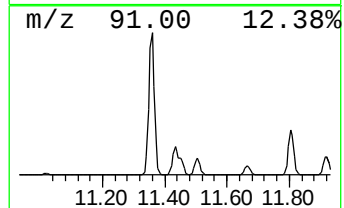
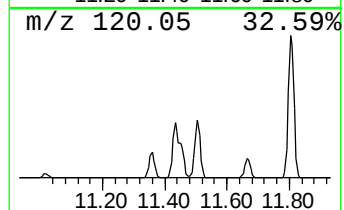
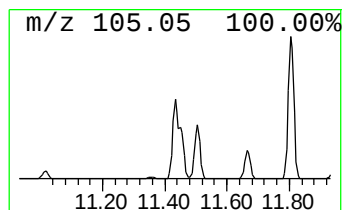
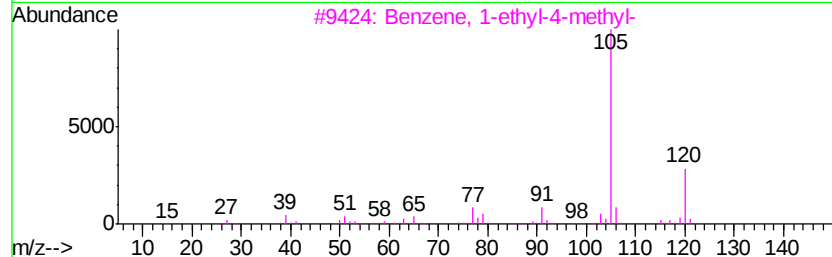
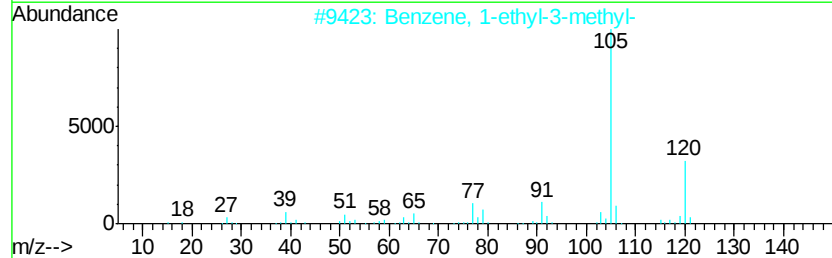
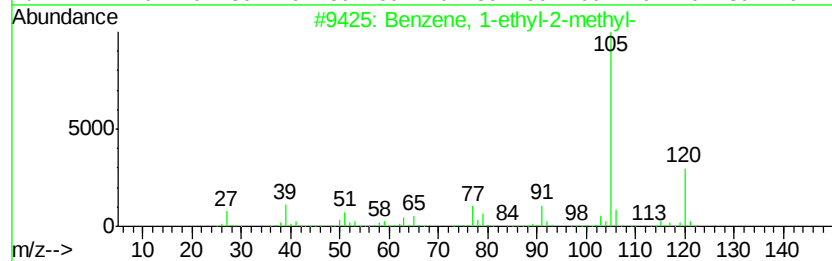
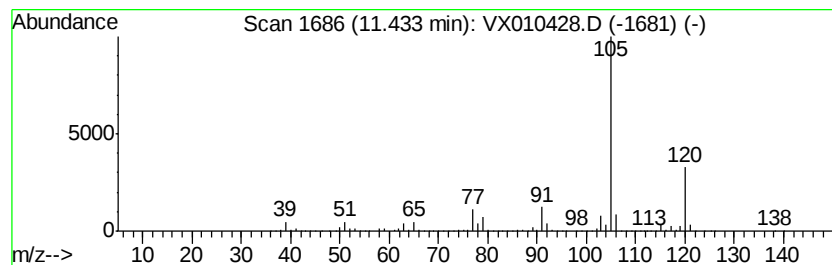
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	261.54 ug/l	7617880	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	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Mesitylene	120	C9H12	000108-67-8	91



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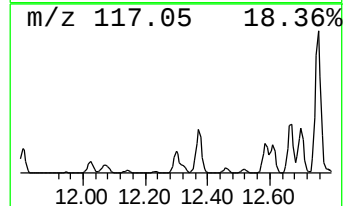
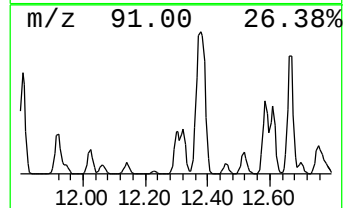
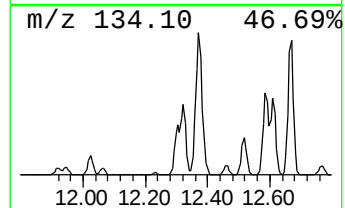
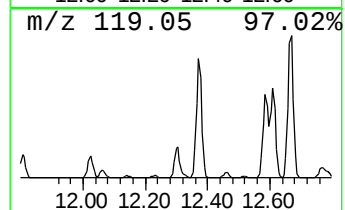
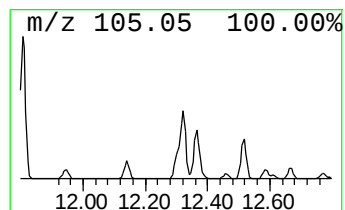
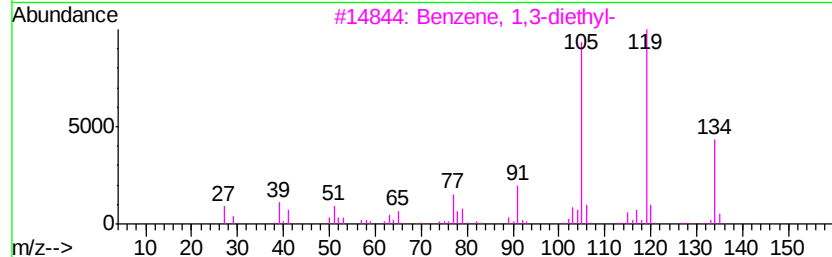
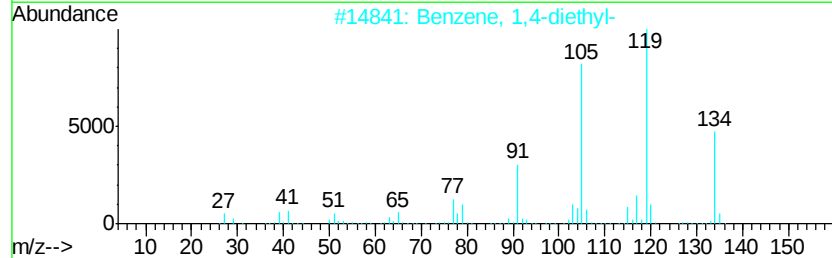
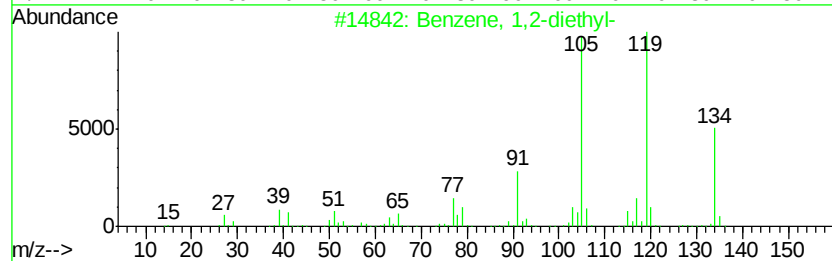
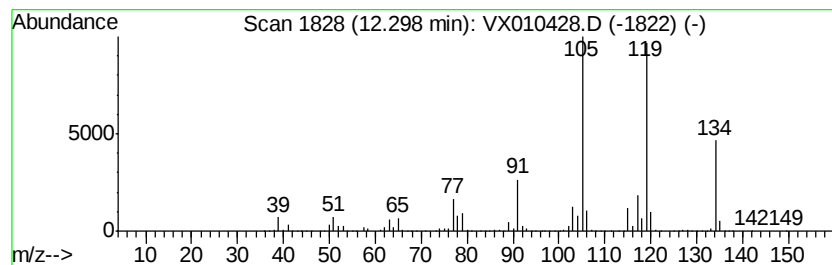
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	110.35 ug/l	3214110	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	83



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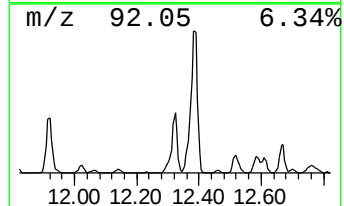
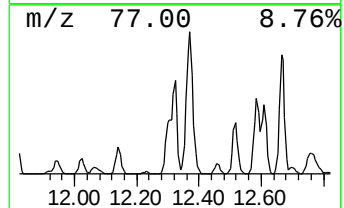
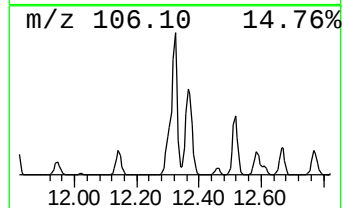
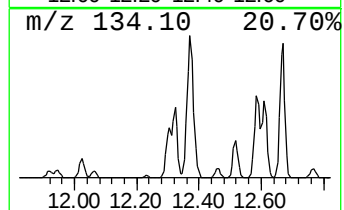
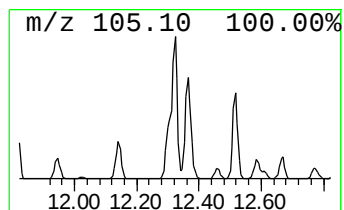
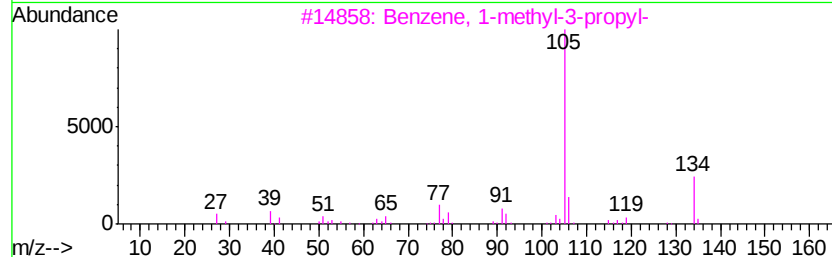
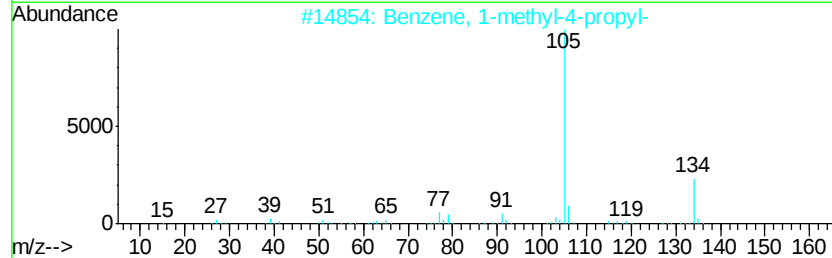
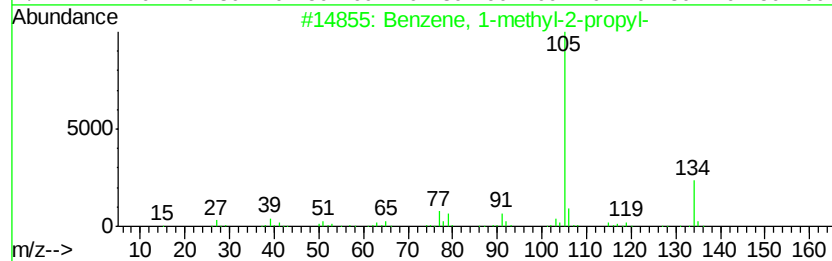
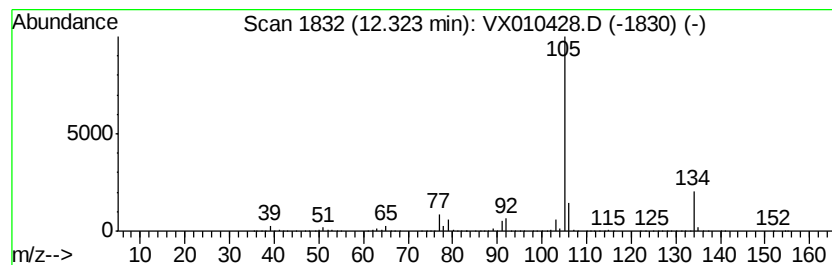
Instrument :
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-methyl-2-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	115.06 ug/l	3351260	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-2-propyl-	134 C10H14	001074-17-5 87
2		Benzene, 1-methyl-4-propyl-	134 C10H14	001074-55-1 86
3		Benzene, 1-methyl-3-propyl-	134 C10H14	001074-43-7 74
4		Benzeneacetaldehyde, .alpha.-met...	134 C9H10O	000093-53-8 64
5		2-Indanol	134 C9H10O	004254-29-9 64



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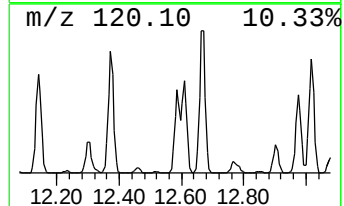
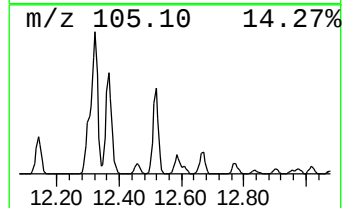
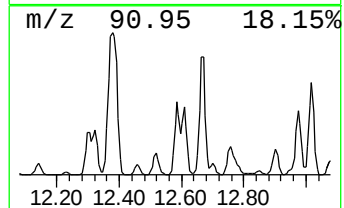
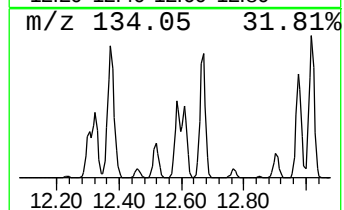
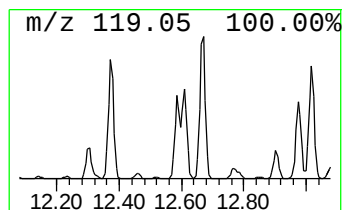
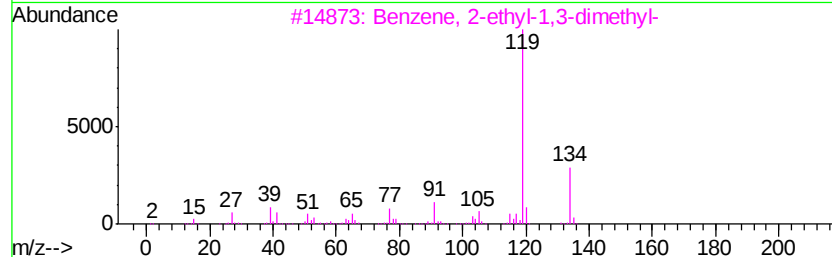
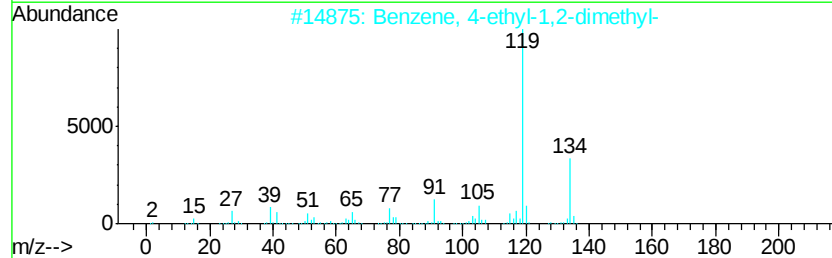
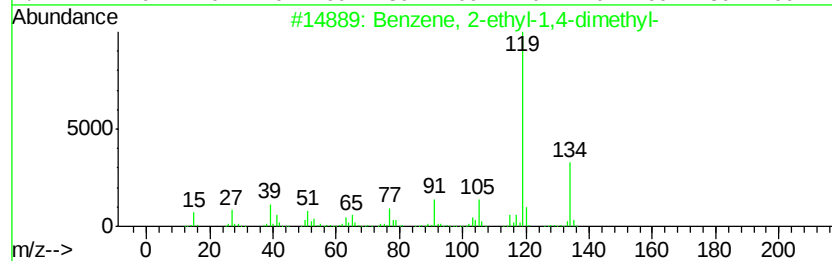
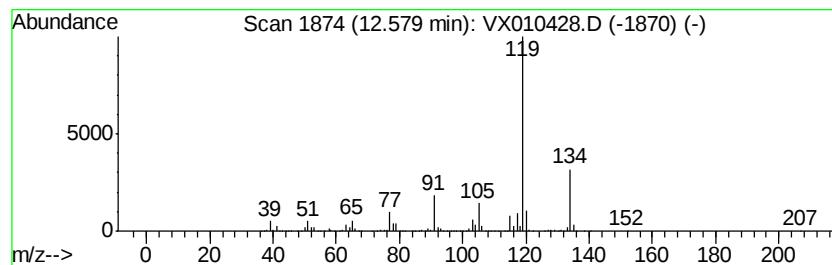
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	159.42 ug/l	4643460	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, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062519\
Data File : VX010428.D
Acq On : 25 Jun 2019 10:58
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

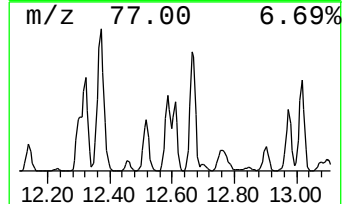
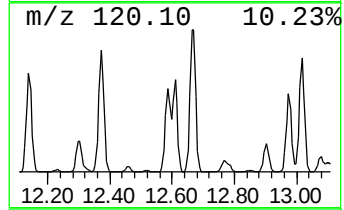
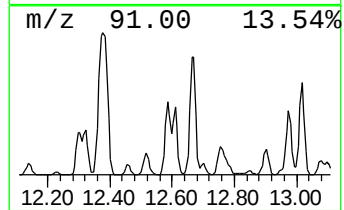
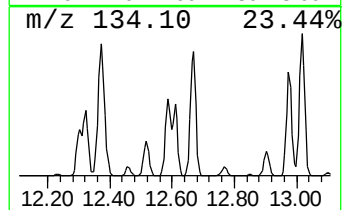
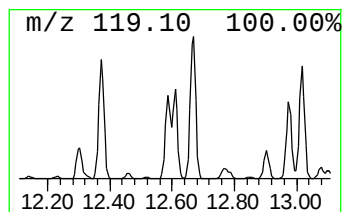
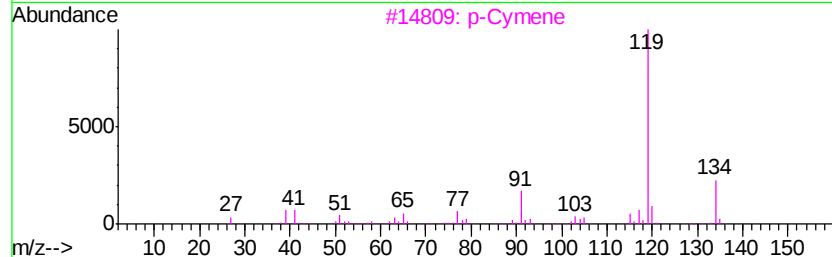
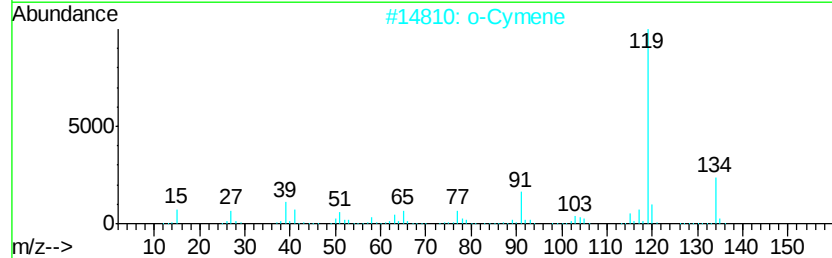
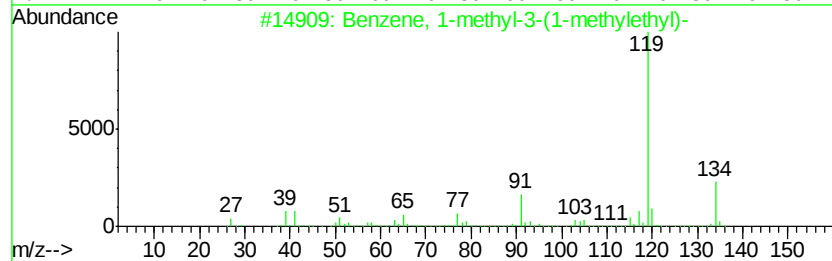
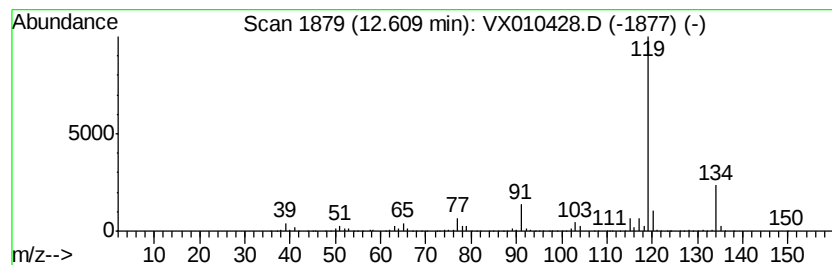
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MSVOA_X
ClientSampleId :
GW-B-061919

Quant Method : Z:\V0ASRV\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, 1-methyl-3-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	127.48 ug/l	3713030	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 94
2		o-Cymene	134 C10H14	000527-84-4 94
3		p-Cymene	134 C10H14	000099-87-6 94
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 91
5		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062519\
Data File : VX010428.D
Acq On : 25 Jun 2019 10:58
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

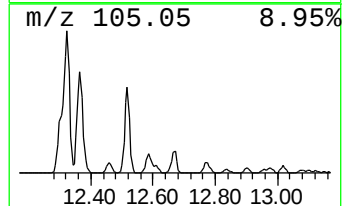
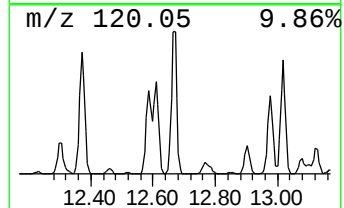
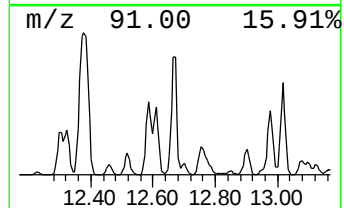
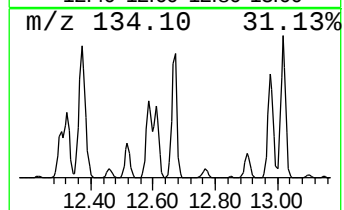
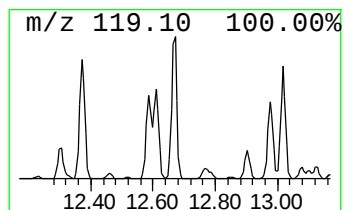
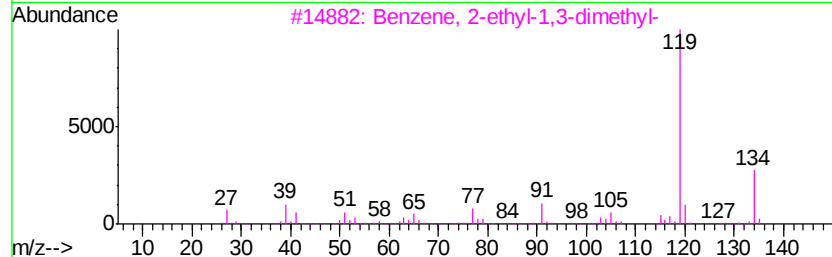
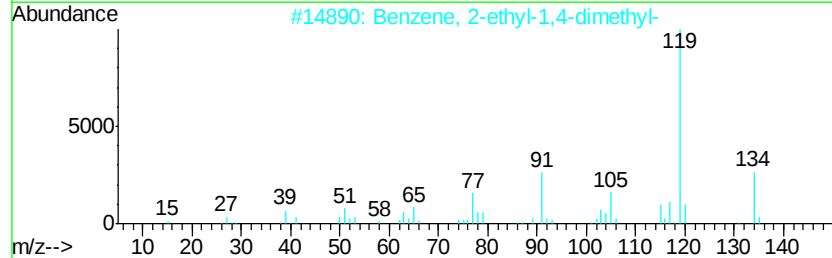
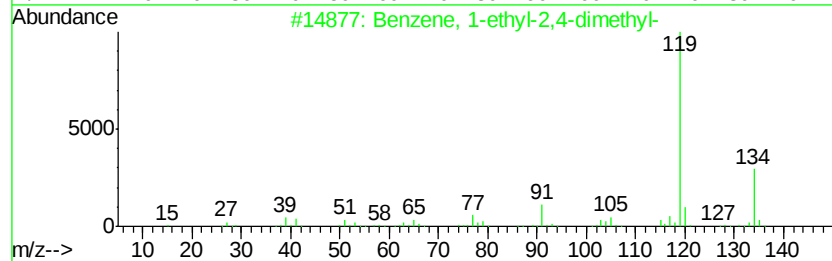
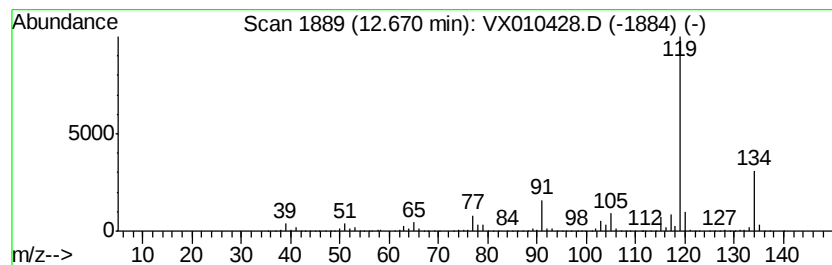
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	254.32 ug/l	7407820	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, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062519\
Data File : VX010428.D
Acq On : 25 Jun 2019 10:58
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

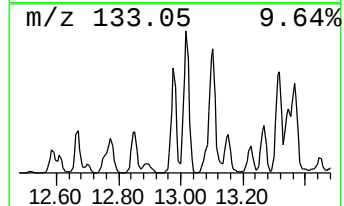
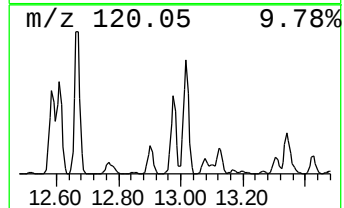
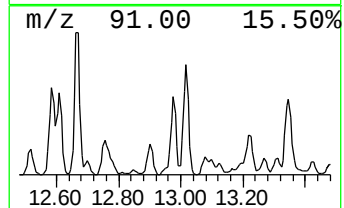
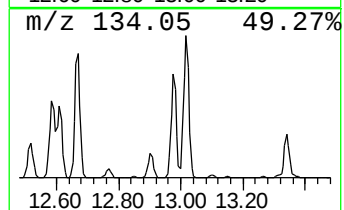
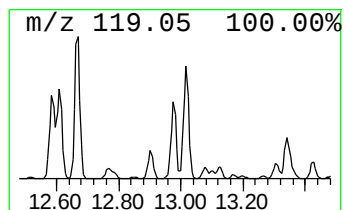
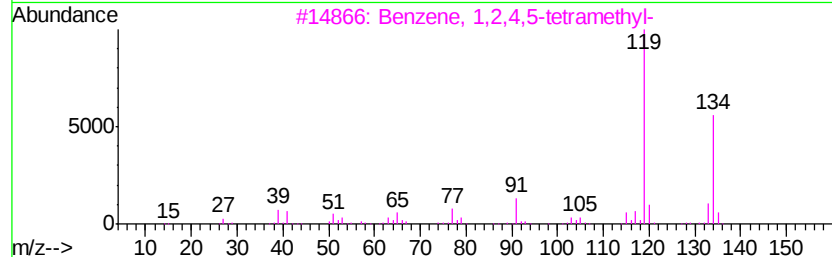
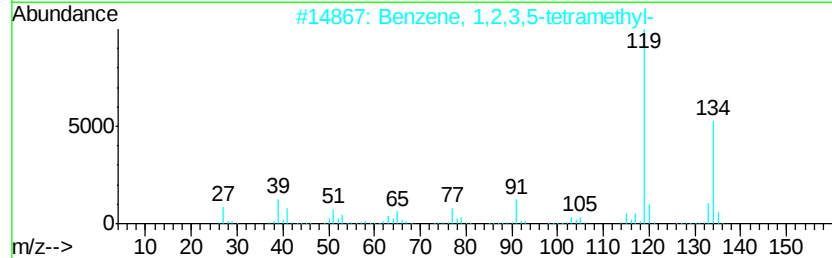
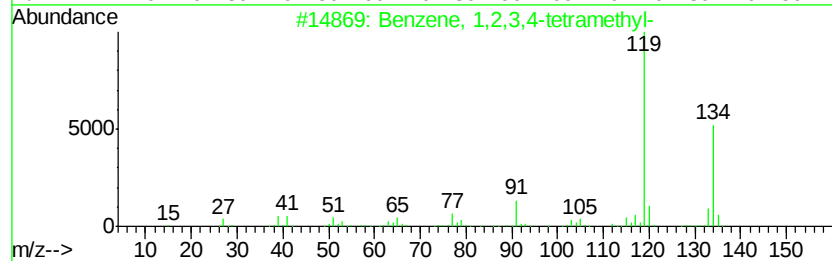
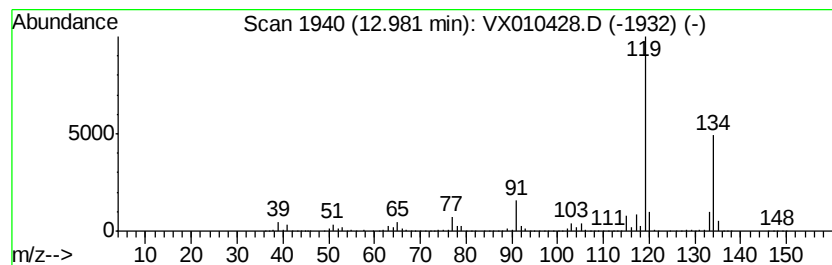
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	154.38 ug/l	4496720	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062519\
Data File : VX010428.D
Acq On : 25 Jun 2019 10:58
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

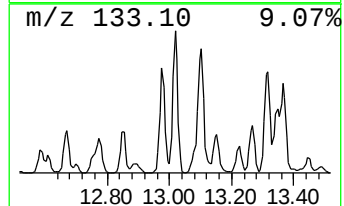
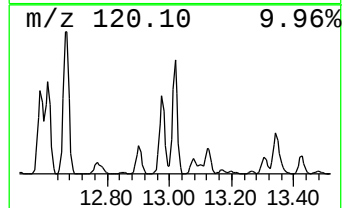
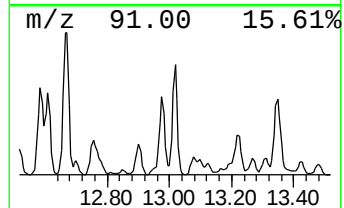
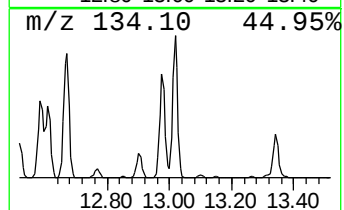
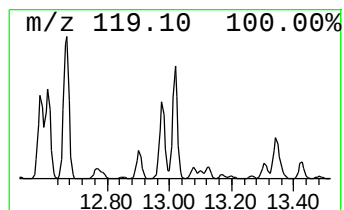
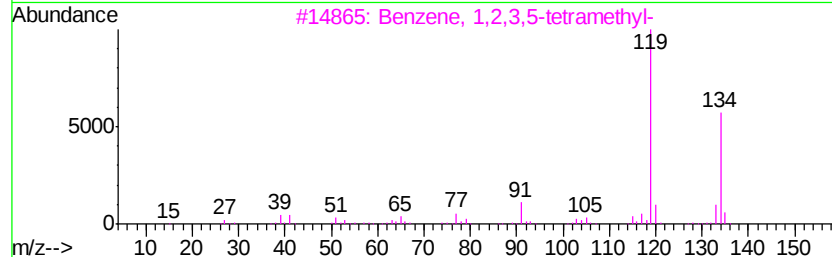
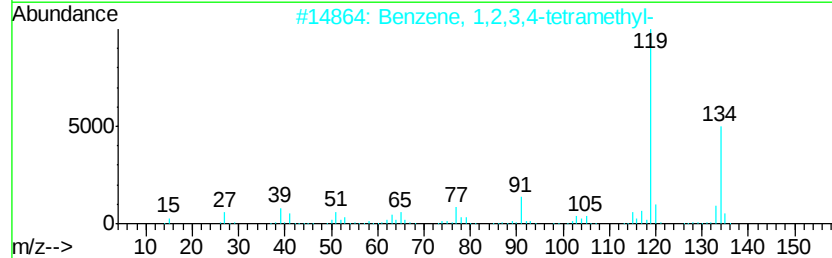
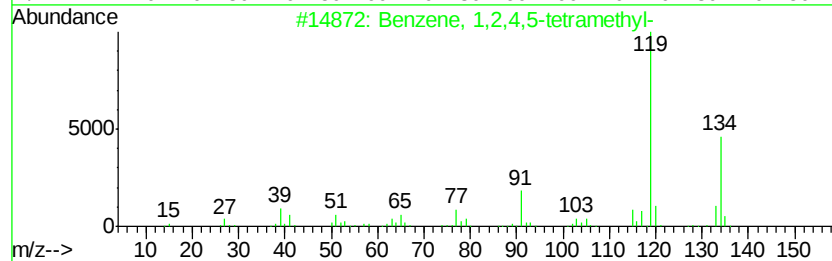
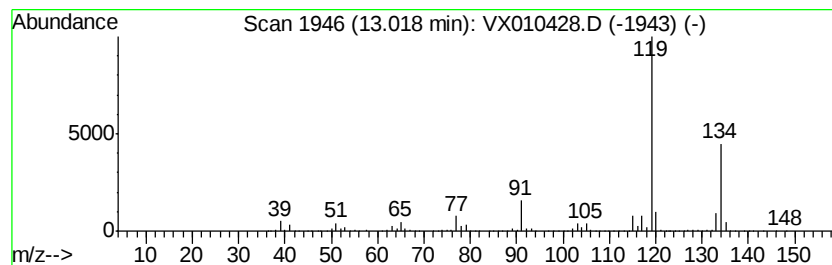
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	198.92 ug/l	5794100	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	95
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\VX062519\
Data File : VX010428.D
Acq On : 25 Jun 2019 10:58
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

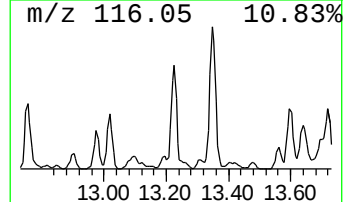
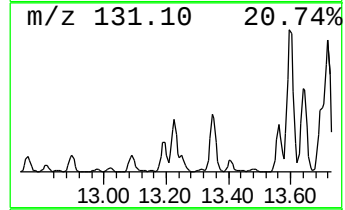
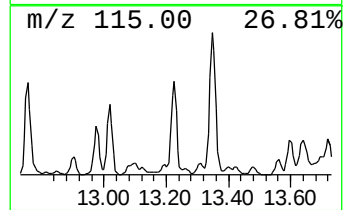
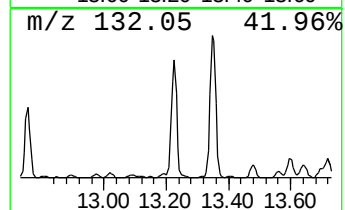
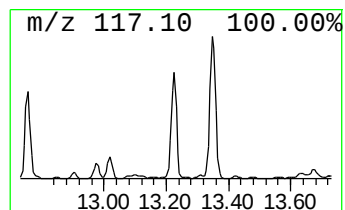
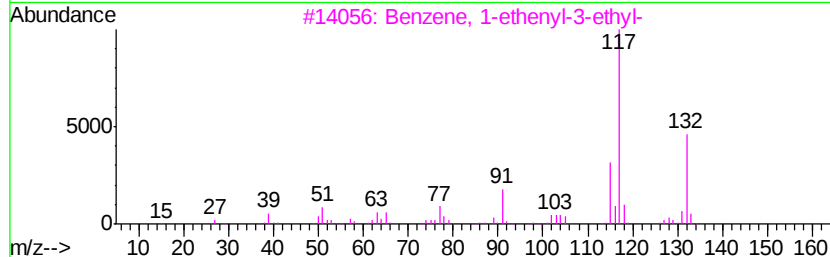
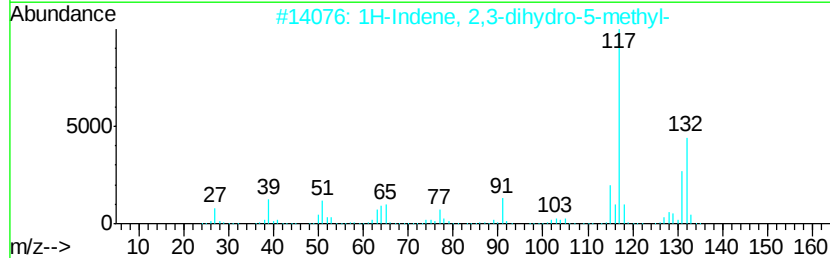
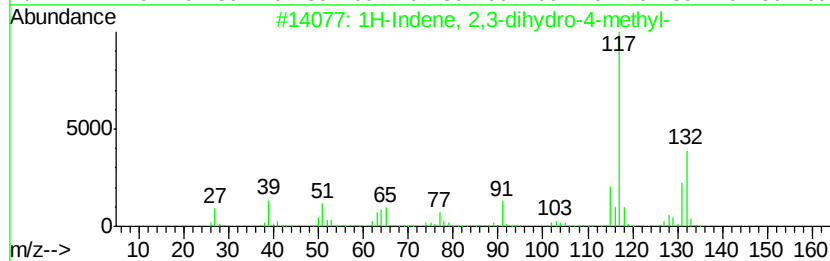
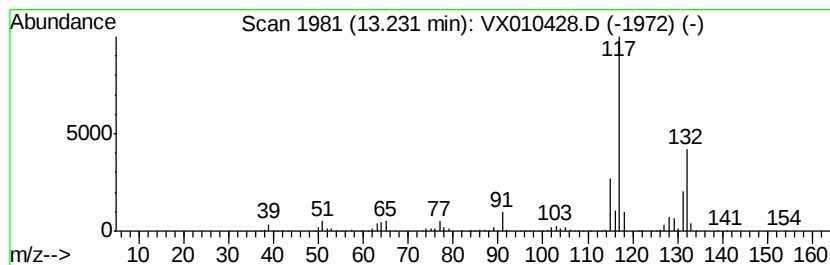
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.23	113.81 ug/l	3315120	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
4		Indan, 1-methyl-	132	C10H12	000767-58-8	90
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062519\
 Data File : VX010428.D
 Acq On : 25 Jun 2019 10:58
 Operator : JC/SP
 Sample : K3500-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 GW-B-061919

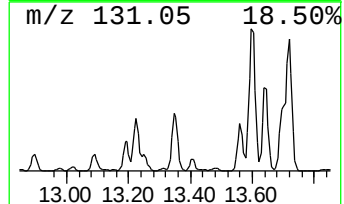
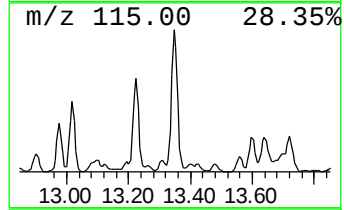
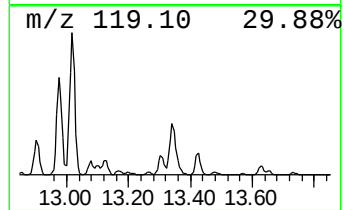
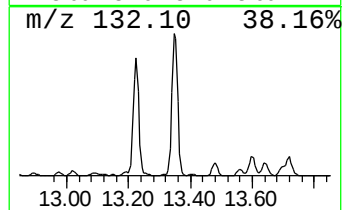
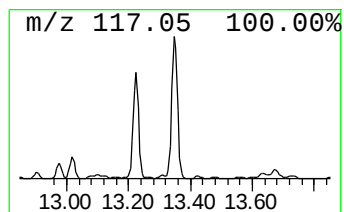
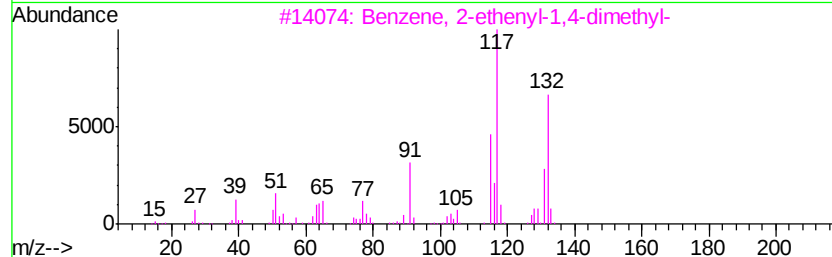
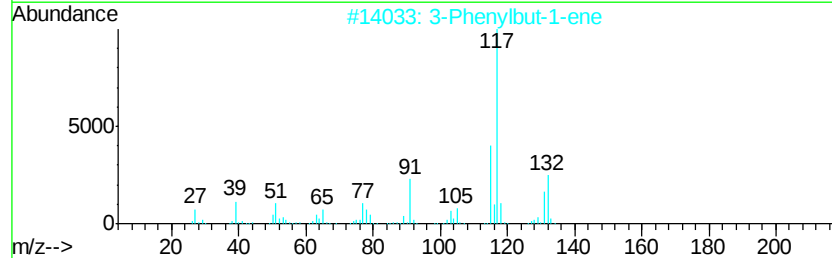
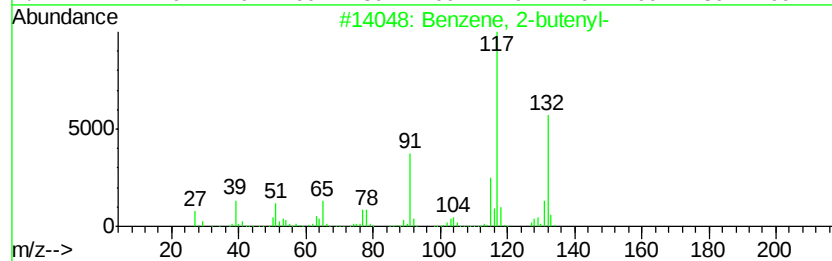
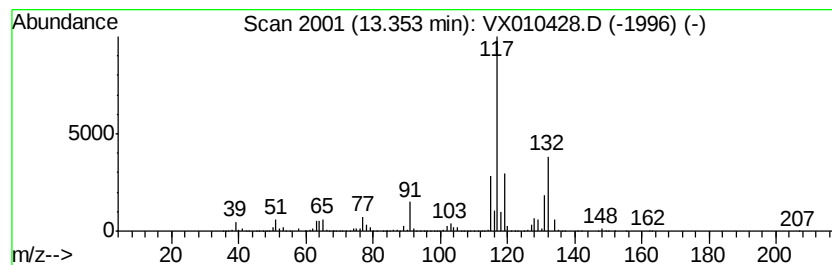
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 Quant Title : SW846 8260

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

 Peak Number 10 Benzene, 2-butenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	215.00 ug/l	6262480	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	70
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062519\
Data File : VX010428.D
Acq On : 25 Jun 2019 10:58
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	11.43	261.5	ug/l	7617880	4	12.08	1456380	50.0
Benzene, 1,2-diet...	12.30	110.3	ug/l	3214110	4	12.08	1456380	50.0
Benzene, 1-methyl...	12.32	115.1	ug/l	3351260	4	12.08	1456380	50.0
Benzene, 2-ethyl-...	12.58	159.4	ug/l	4643460	4	12.08	1456380	50.0
Benzene, 1-methyl...	12.61	127.5	ug/l	3713030	4	12.08	1456380	50.0
Benzene, 1-ethyl-...	12.67	254.3	ug/l	7407820	4	12.08	1456380	50.0
Benzene, 1,2,3,4-...	12.98	154.4	ug/l	4496720	4	12.08	1456380	50.0
Benzene, 1,2,4,5-...	13.02	198.9	ug/l	5794100	4	12.08	1456380	50.0
1H-Indene, 2,3-di...	13.23	113.8	ug/l	3315120	4	12.08	1456380	50.0
Benzene, 2-butenyl-	13.35	215.0	ug/l	6262480	4	12.08	1456380	50.0