

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062419\
 Data File : VX010408.D
 Acq On : 24 Jun 2019 13:17
 Operator : JC/SP
 Sample : K3500-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 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	1.794	99	105	117	rVB	559161	804261	7.17%	0.616%
2	2.001	133	139	148	rBB	245002	337313	3.01%	0.258%
3	2.849	270	278	280	rBV	56924	114308	1.02%	0.088%
4	2.891	280	285	296	rVB	125071	271296	2.42%	0.208%
5	3.172	323	331	343	rVB	87600	205124	1.83%	0.157%
6	4.403	522	533	545	rBV2	52802	160042	1.43%	0.123%
7	5.500	699	713	722	rBV	170889	496127	4.42%	0.380%
8	5.659	731	739	757	rVB	288466	823484	7.34%	0.631%
9	6.061	794	805	819	rBV	155549	408701	3.64%	0.313%
10	6.860	925	936	947	rBV	424540	1012969	9.03%	0.776%
11	7.464	1024	1035	1046	rVB3	63041	158232	1.41%	0.121%
12	8.713	1234	1240	1249	rVB	899557	1389006	12.38%	1.064%
13	10.109	1464	1469	1475	rBV	879799	1167520	10.41%	0.894%
14	11.018	1612	1618	1626	rBV	397787	506182	4.51%	0.388%
15	11.134	1632	1637	1642	rBV	742517	950858	8.47%	0.728%
16	11.359	1665	1674	1681	rBV	2252914	2719211	24.24%	2.082%
17	11.432	1681	1686	1688	rBV	4478098	5716947	50.95%	4.378%
18	11.506	1694	1698	1707	rVB	3471858	4155050	37.03%	3.182%
19	11.664	1719	1724	1734	rVB	1569777	1970145	17.56%	1.509%
20	11.804	1742	1747	1753	rVB	8881167	11219931	100.00%	8.593%
21	11.920	1760	1766	1768	rBV	389915	531890	4.74%	0.407%
22	11.944	1768	1770	1775	rVB	488739	519492	4.63%	0.398%
23	12.024	1775	1783	1786	rBV	1018991	1201813	10.71%	0.920%
24	12.073	1786	1791	1797	rVB2	1006012	1576330	14.05%	1.207%
25	12.140	1797	1802	1808	rBV	1071164	1309139	11.67%	1.003%
26	12.231	1813	1817	1821	rVB	108330	130814	1.17%	0.100%
27	12.304	1821	1829	1830	rBV	2621888	3400636	30.31%	2.604%
28	12.322	1830	1832	1836	rVV	3603879	3780946	33.70%	2.896%
29	12.371	1836	1840	1847	rVB3	6743499	10004122	89.16%	7.661%
30	12.457	1850	1854	1859	rVB	469919	596530	5.32%	0.457%
31	12.518	1859	1864	1870	rVB	1974088	2392281	21.32%	1.832%
32	12.585	1870	1875	1877	rBV	3926807	4920861	43.86%	3.769%
33	12.609	1877	1879	1884	rVB	3825726	4096150	36.51%	3.137%
34	12.670	1884	1889	1892	rBV	6691687	7985070	71.17%	6.115%

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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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 Title : SW846 8260

35	12.700	1892	1894	1898	rVB	714943	745930	6.65%	0.571%
36	12.755	1898	1903	1911	rBV3	1989636	3349830	29.86%	2.565%
37	12.853	1914	1919	1922	rVB2	184776	255870	2.28%	0.196%
38	12.902	1922	1927	1932	rVB	1383928	1712207	15.26%	1.311%
39	12.975	1932	1939	1943	rBV	3753194	4842911	43.16%	3.709%
40	13.017	1943	1946	1951	rVB	5429798	6233995	55.56%	4.774%
41	13.078	1951	1956	1957	rBV	634997	714740	6.37%	0.547%
42	13.097	1957	1959	1962	rVV2	631743	838396	7.47%	0.642%
43	13.121	1962	1963	1966	rVB	423424	361603	3.22%	0.277%
44	13.225	1972	1980	1984	rBV2	2509625	3589359	31.99%	2.749%
45	13.267	1984	1987	1990	rVB	437419	512445	4.57%	0.392%
46	13.310	1990	1994	1996	rBV2	780009	1077228	9.60%	0.825%
47	13.347	1996	2000	2009	rVV2	4640779	6743019	60.10%	5.164%
48	13.426	2009	2013	2018	rVB	623094	794873	7.08%	0.609%
49	13.481	2018	2022	2029	rVB	374166	495014	4.41%	0.379%
50	13.560	2029	2035	2038	rBV2	672300	952827	8.49%	0.730%
51	13.597	2038	2041	2044	rVV	1253287	1704403	15.19%	1.305%
52	13.639	2044	2048	2054	rVV3	1674708	3387526	30.19%	2.594%
53	13.719	2058	2061	2067	rVB2	1369894	2191499	19.53%	1.678%
54	13.847	2077	2082	2089	rVB3	154709	276578	2.47%	0.212%
55	13.914	2089	2093	2097	rVB2	159988	230506	2.05%	0.177%
56	13.981	2097	2104	2108	rBV2	694808	1046923	9.33%	0.802%
57	14.029	2108	2112	2115	rVV	542049	673184	6.00%	0.516%
58	14.060	2115	2117	2124	rVB2	75121	115167	1.03%	0.088%
59	14.133	2124	2129	2135	rBV	1194166	1474114	13.14%	1.129%
60	14.273	2147	2152	2158	rBV	947388	1599963	14.26%	1.225%
61	14.377	2165	2169	2174	rBV	506037	659082	5.87%	0.505%
62	14.426	2174	2177	2182	rVB	679135	850578	7.58%	0.651%
63	14.542	2190	2196	2200	rBV2	170677	276384	2.46%	0.212%
64	14.694	2208	2221	2226	rBV	1241742	1667883	14.87%	1.277%
65	14.834	2239	2244	2254	rVB	1869577	2450895	21.84%	1.877%
66	15.267	2304	2315	2320	rBV2	122179	226261	2.02%	0.173%
67	15.444	2333	2344	2347	rBV	134346	223272	1.99%	0.171%
68	15.547	2354	2361	2374	rVB	252684	443275	3.95%	0.339%
69	15.682	2374	2383	2387	rVV	278250	467216	4.16%	0.358%
70	15.724	2387	2390	2398	rVB	140102	200282	1.79%	0.153%
71	15.913	2411	2421	2433	rVB2	61118	159094	1.42%	0.122%

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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
Title : SW846 8260

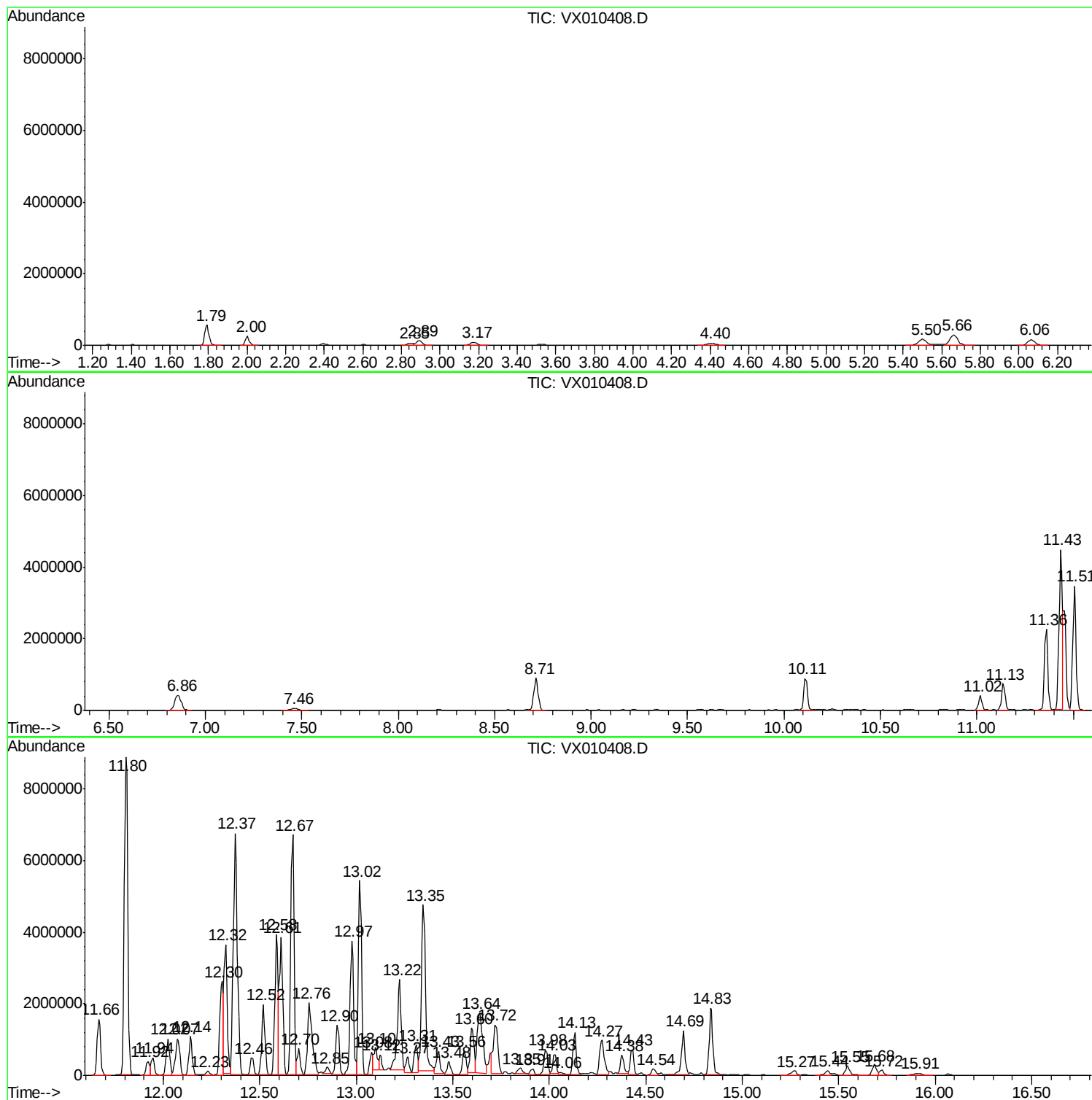
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

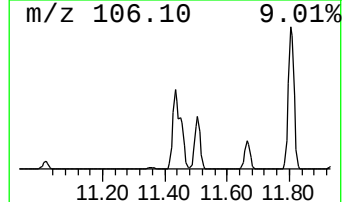
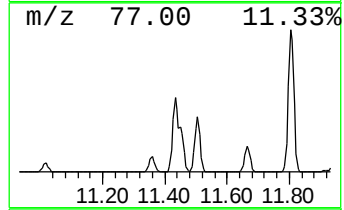
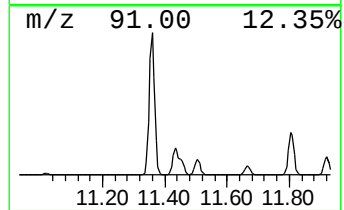
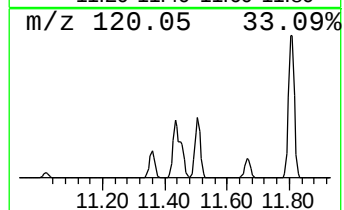
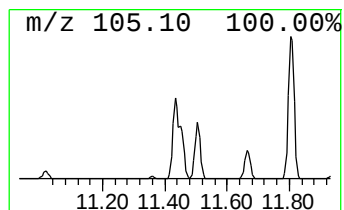
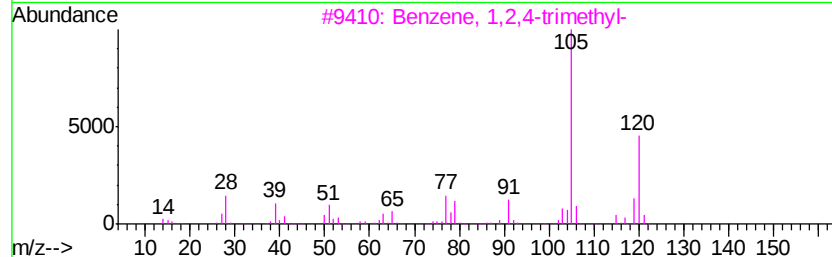
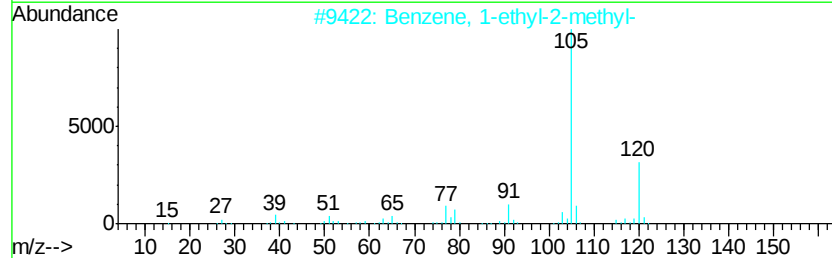
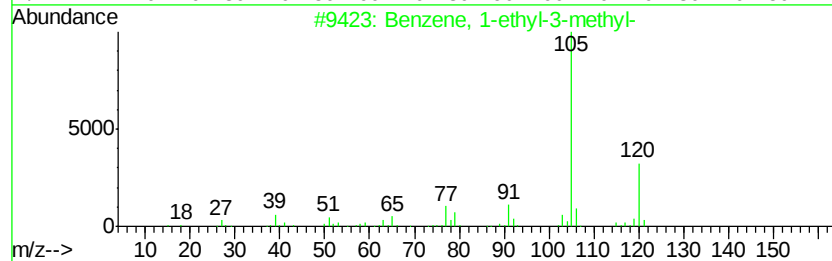
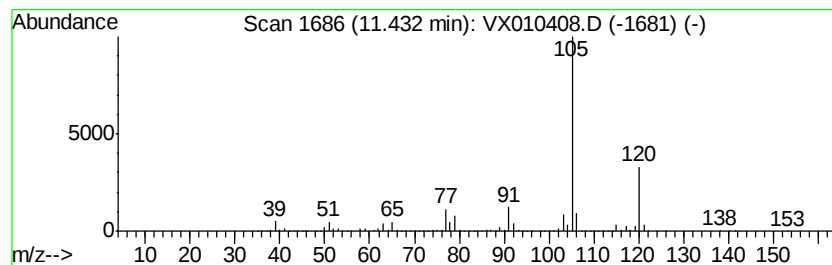
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	181.34 ug/l	5716950	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampled :
GW-B-061919

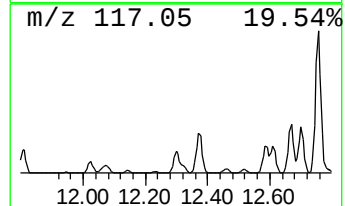
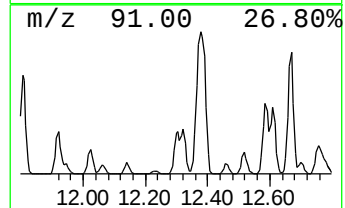
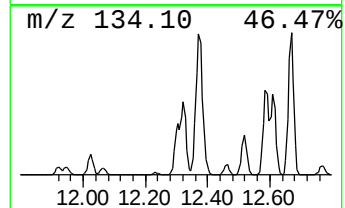
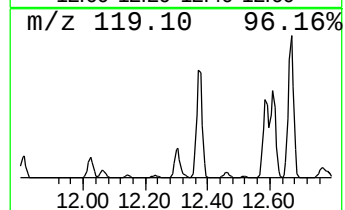
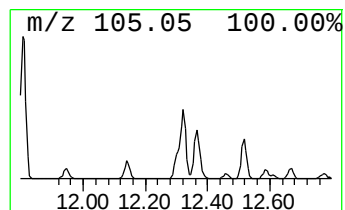
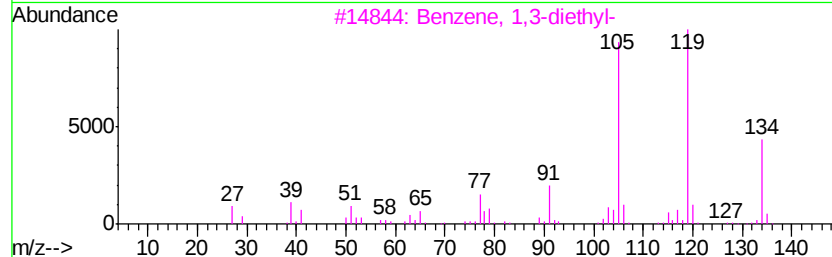
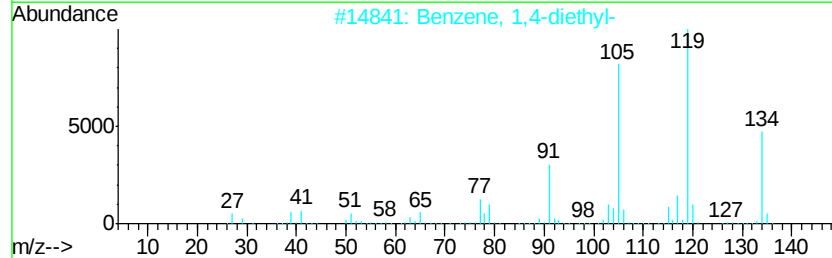
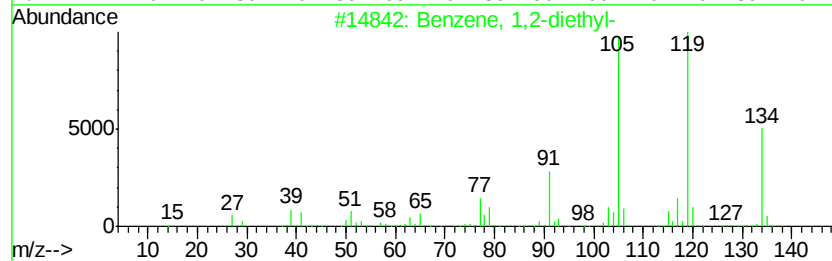
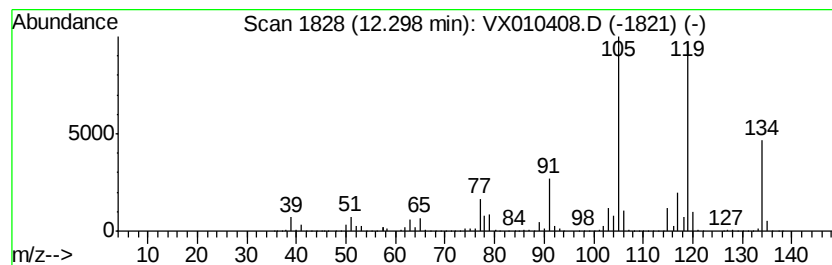
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	107.87 ug/l	3400640	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	91
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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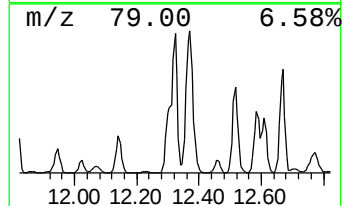
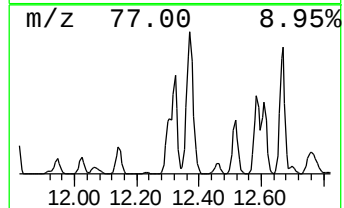
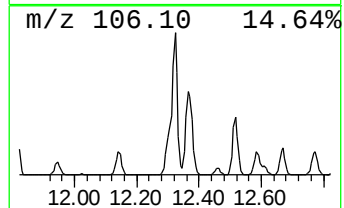
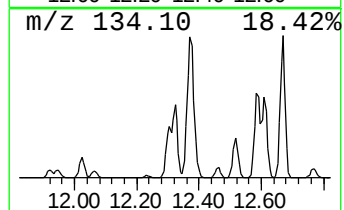
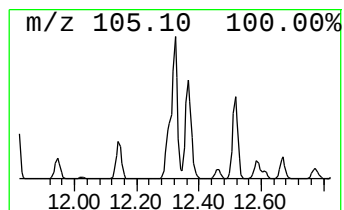
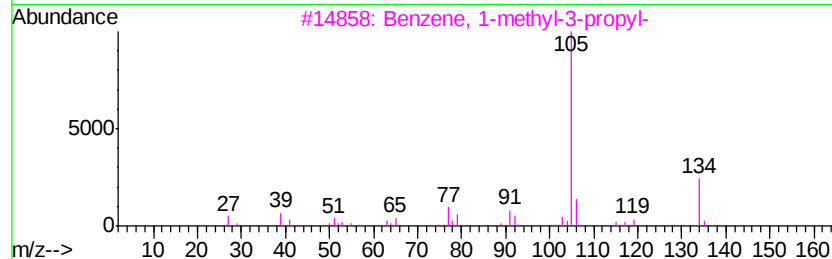
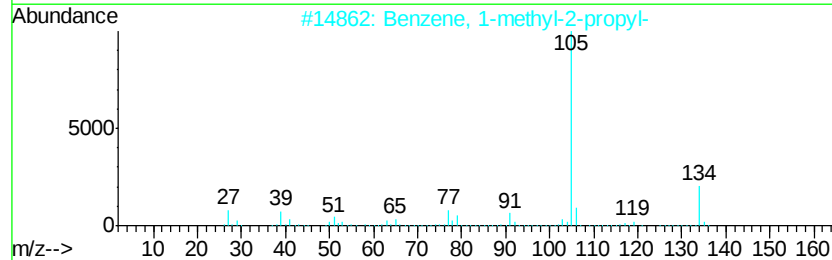
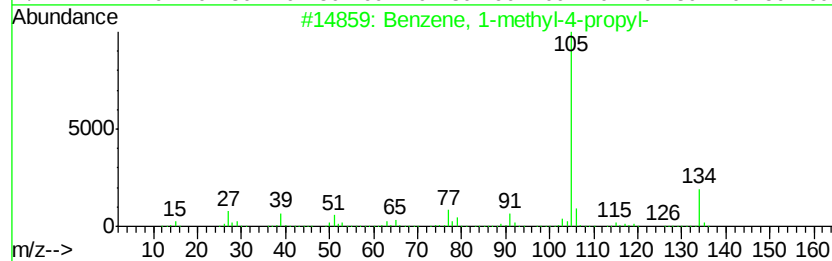
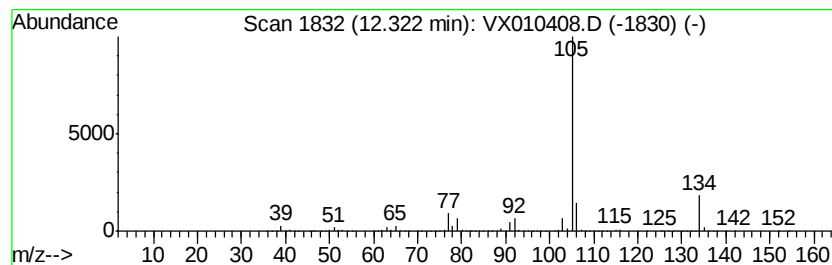
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-methyl-4-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	119.93 ug/l	3780950	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	80
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	80
3		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	74
4		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	50
5		2-Indanol	134	C9H10O	004254-29-9	50



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Instrument :
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ClientSampled :
GW-B-061919

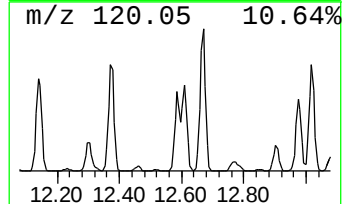
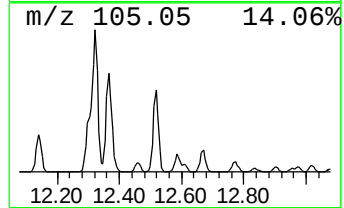
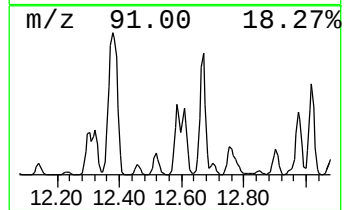
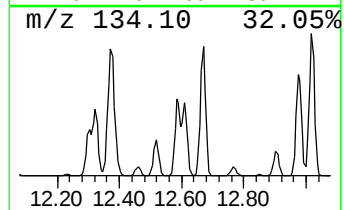
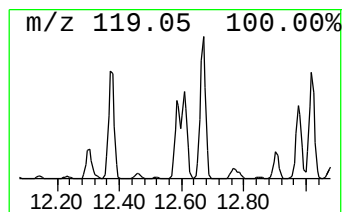
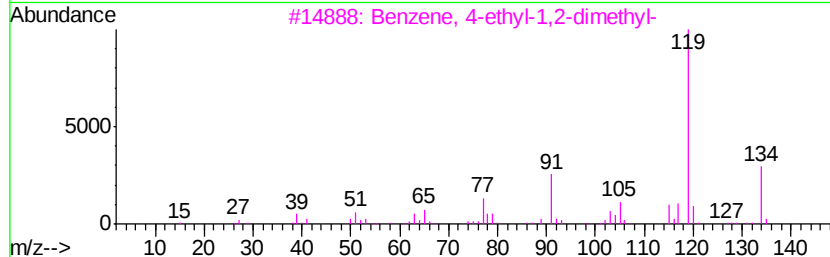
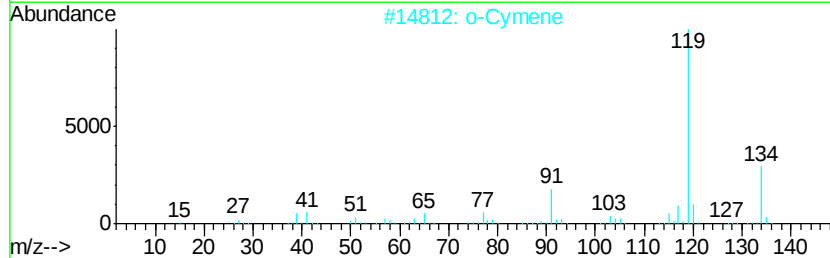
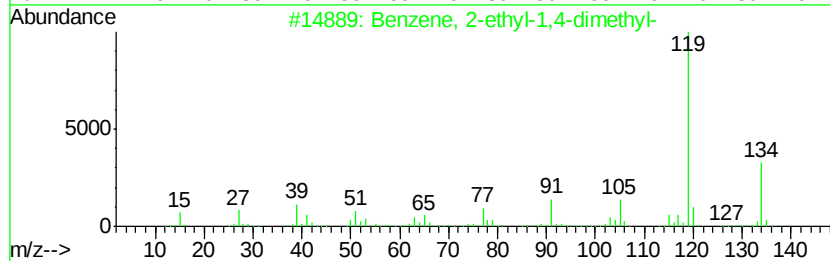
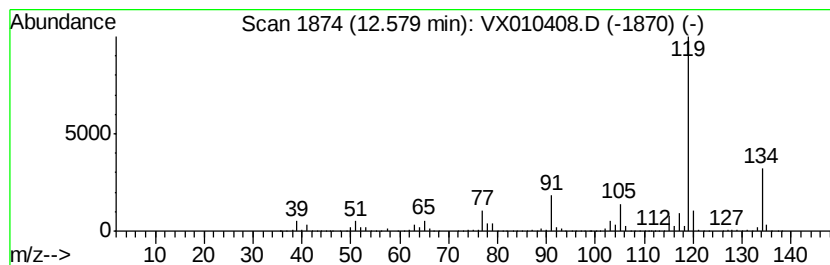
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Quant Title : SW846 8260

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	156.09 ug/l	4920860	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		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	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\VX062419\
Data File : VX010408.D
Acq On : 24 Jun 2019 13:17
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

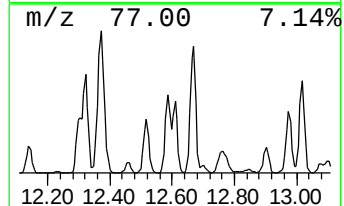
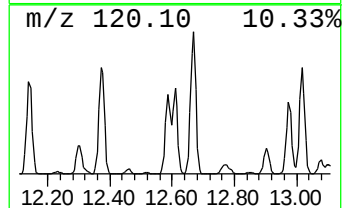
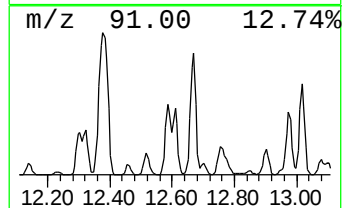
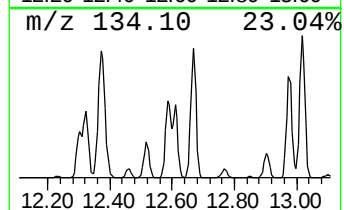
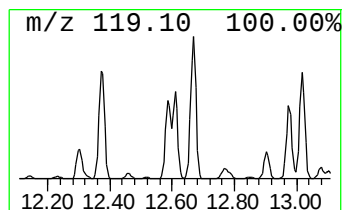
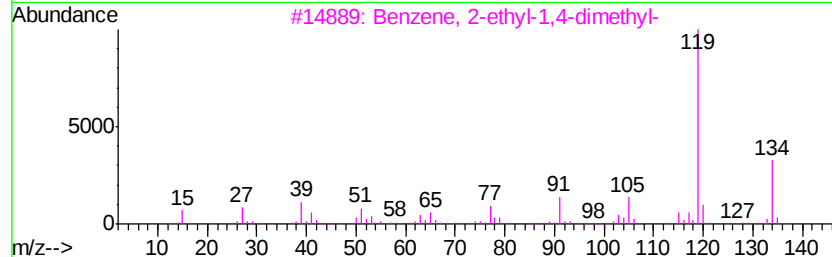
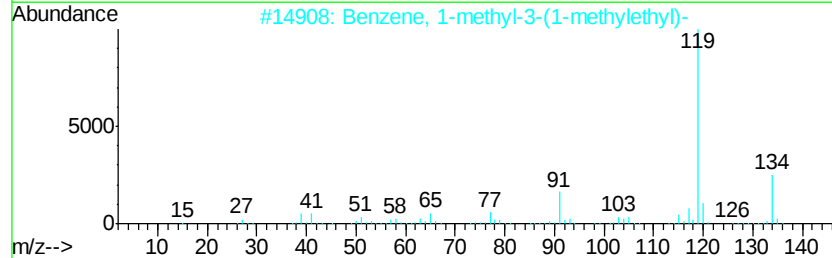
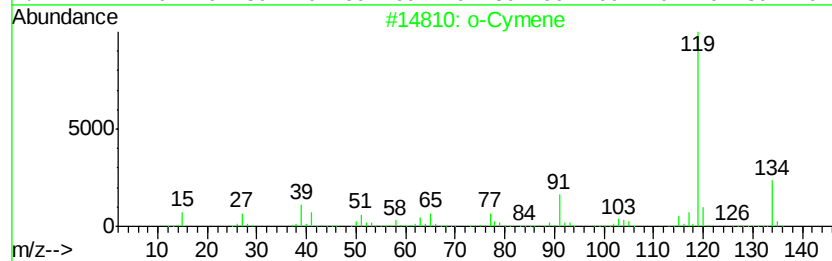
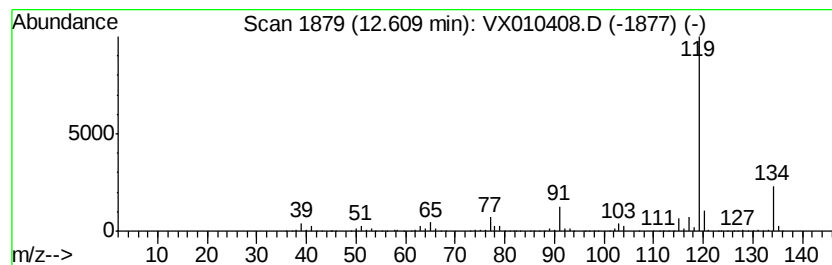
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Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	129.93 ug/l	4096150	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	94
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062419\
Data File : VX010408.D
Acq On : 24 Jun 2019 13:17
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

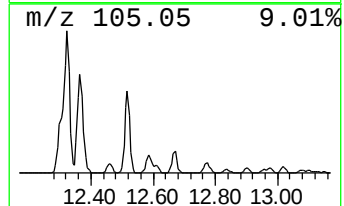
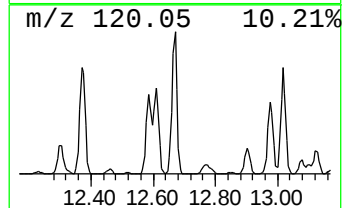
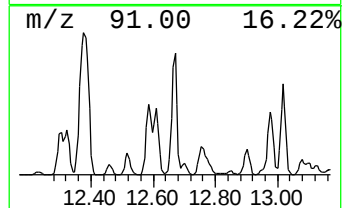
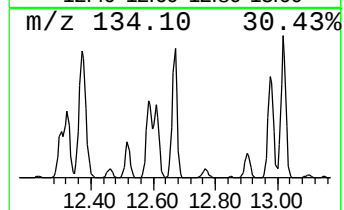
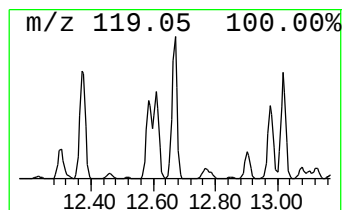
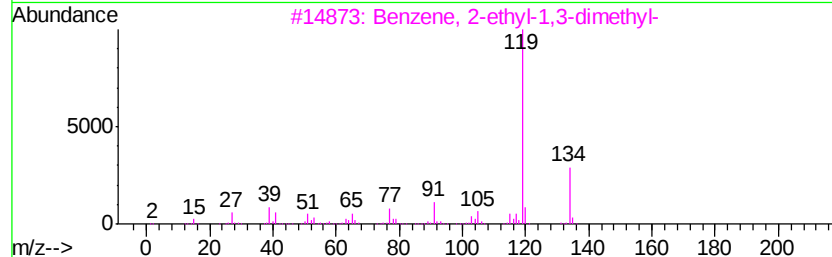
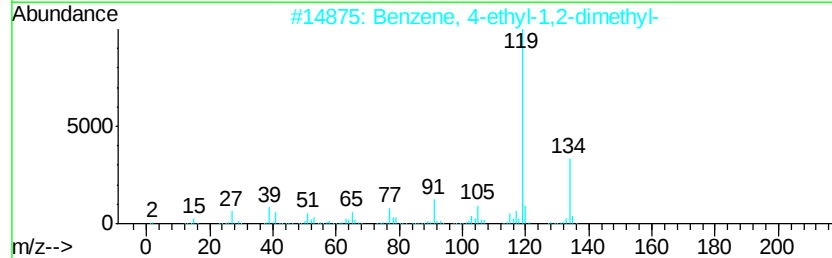
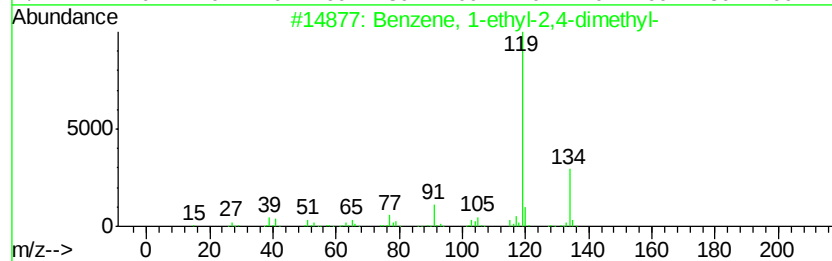
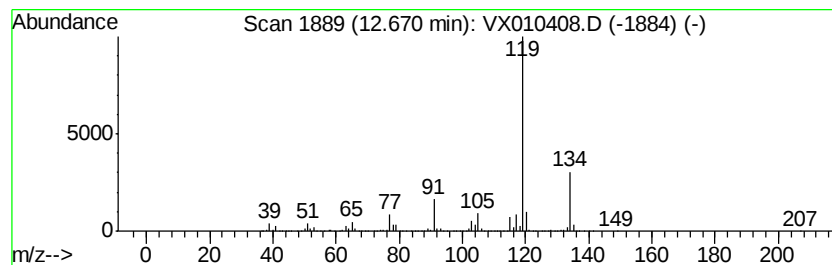
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	253.28 ug/l	7985070	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, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062419\
Data File : VX010408.D
Acq On : 24 Jun 2019 13:17
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

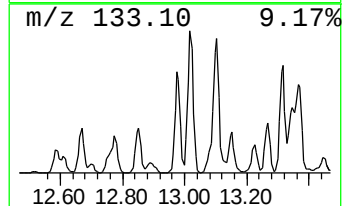
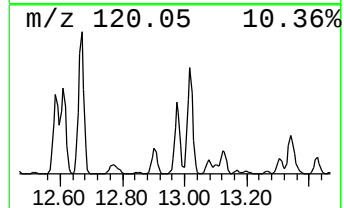
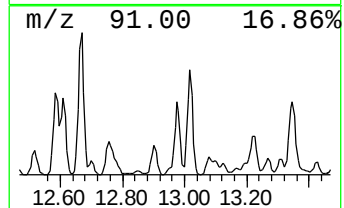
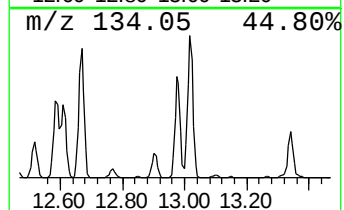
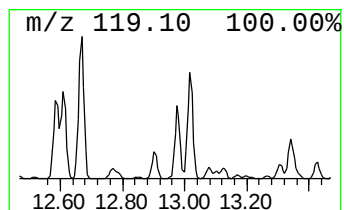
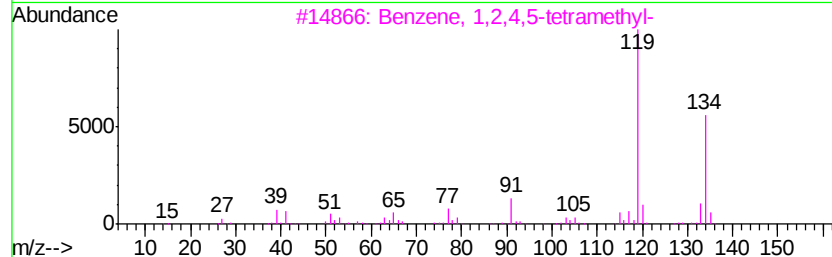
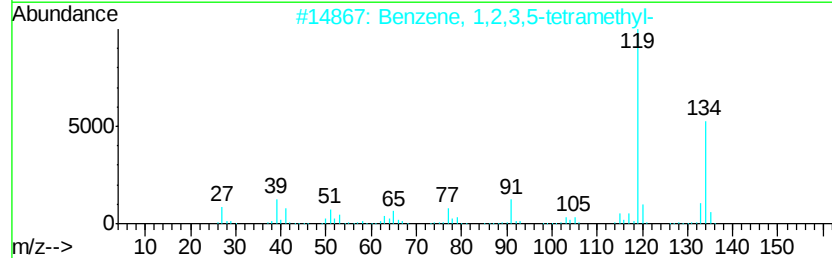
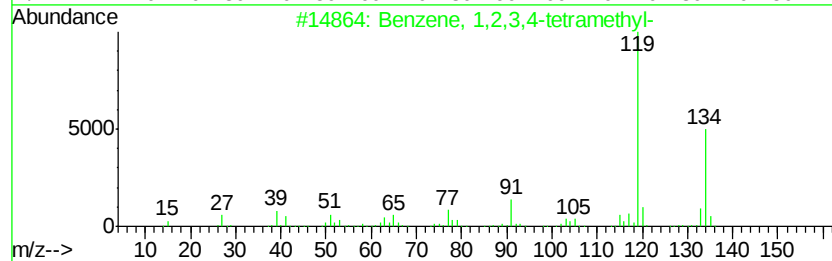
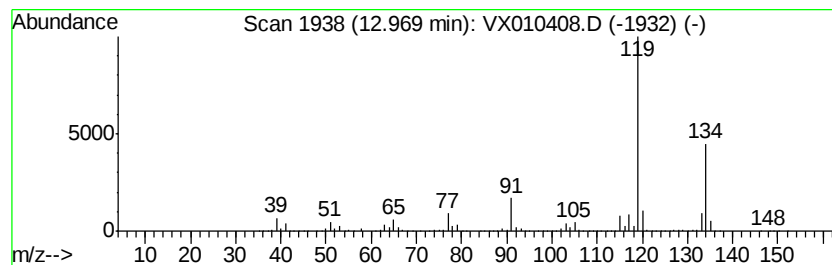
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	153.61 ug/l	4842910	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\VX062419\
Data File : VX010408.D
Acq On : 24 Jun 2019 13:17
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

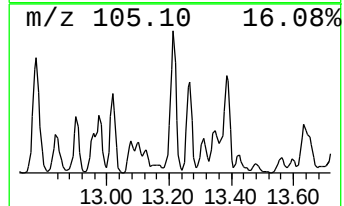
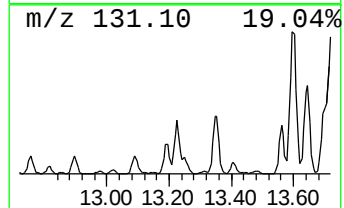
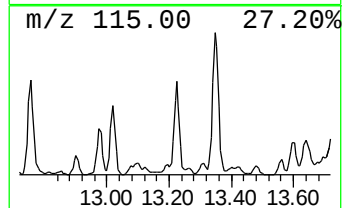
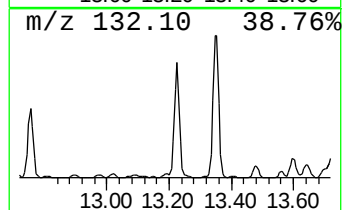
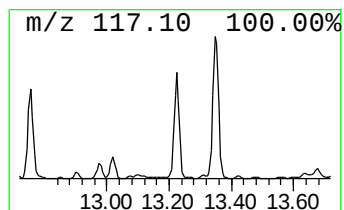
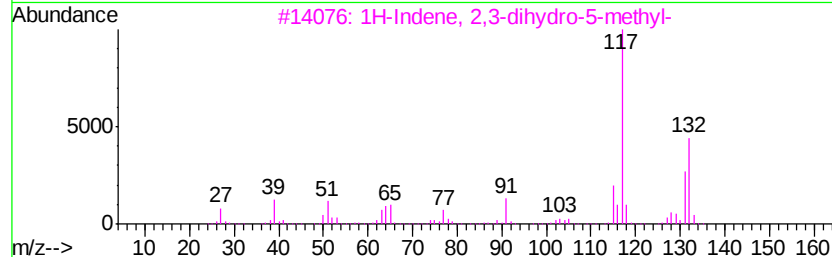
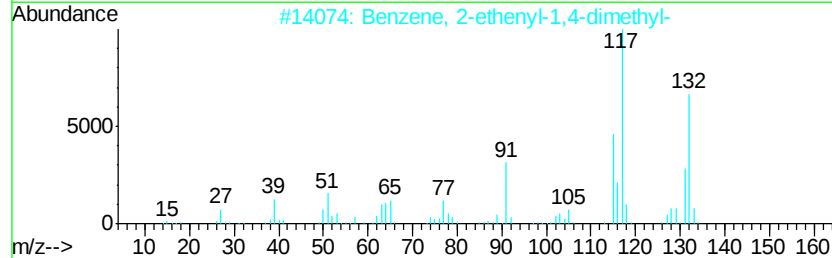
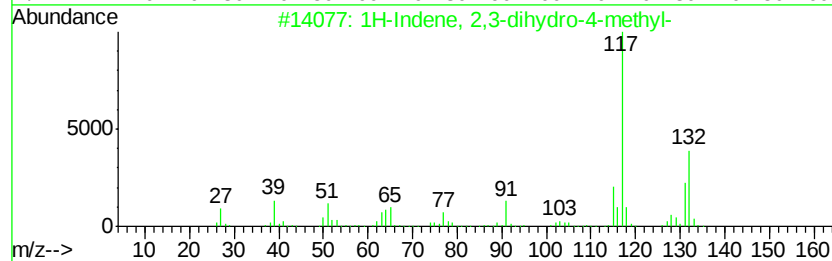
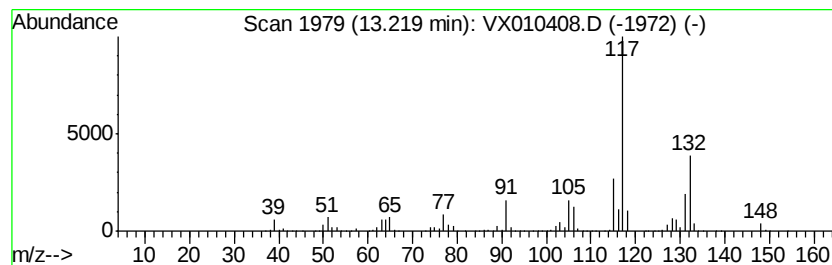
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	113.85 ug/l	3589360	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		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062419\
Data File : VX010408.D
Acq On : 24 Jun 2019 13:17
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-B-061919

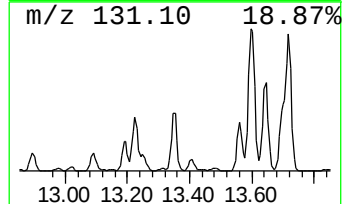
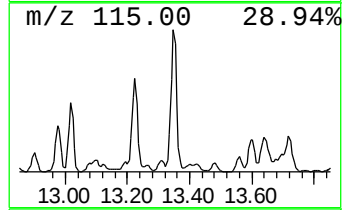
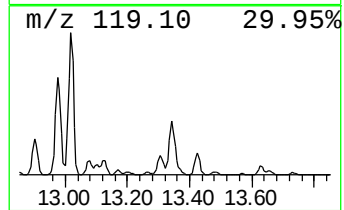
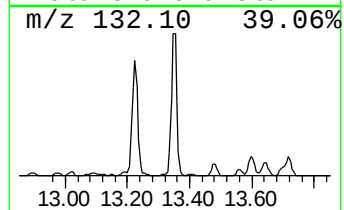
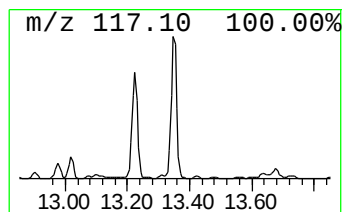
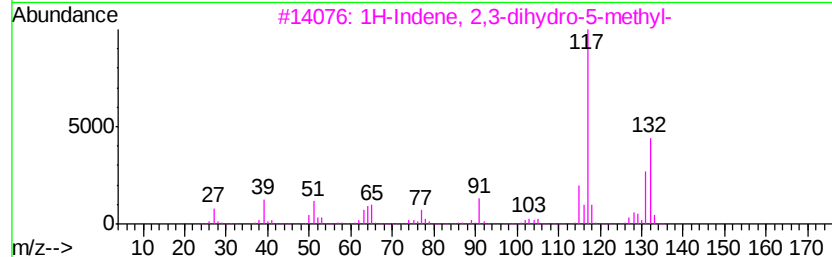
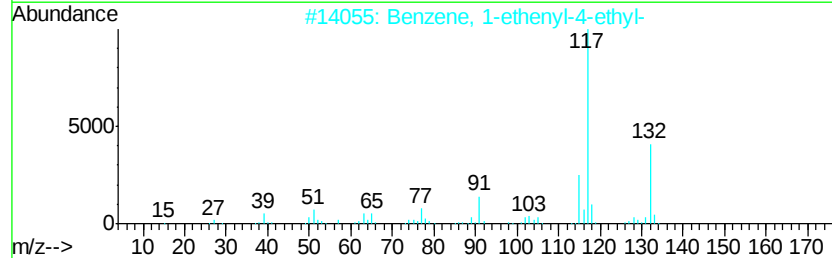
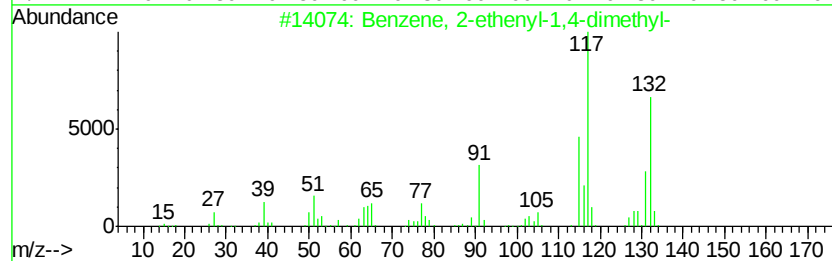
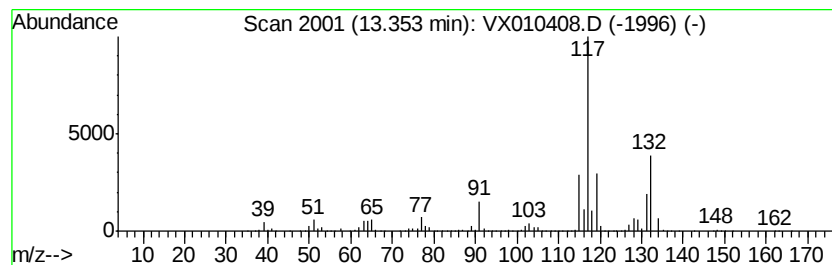
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	213.88 ug/l	6743020	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	91
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062419\
Data File : VX010408.D
Acq On : 24 Jun 2019 13:17
Operator : JC/SP
Sample : K3500-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 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	181.3	ug/l	5716950	4	12.08	1576330	50.0
Benzene, 1,2-diet...	12.30	107.9	ug/l	3400640	4	12.08	1576330	50.0
Benzene, 1-methyl...	12.32	119.9	ug/l	3780950	4	12.08	1576330	50.0
Benzene, 2-ethyl-...	12.58	156.1	ug/l	4920860	4	12.08	1576330	50.0
o-Cymene	12.61	129.9	ug/l	4096150	4	12.08	1576330	50.0
Benzene, 1-ethyl-...	12.67	253.3	ug/l	7985070	4	12.08	1576330	50.0
Benzene, 1,2,3,4-...	12.97	153.6	ug/l	4842910	4	12.08	1576330	50.0
1H-Indene, 2,3-di...	13.22	113.8	ug/l	3589360	4	12.08	1576330	50.0
Benzene, 2-etheny...	13.35	213.9	ug/l	6743020	4	12.08	1576330	50.0