

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112719\
 Data File : VX013663.D
 Acq On : 27 Nov 2019 19:44
 Operator : JC/SP
 Sample : K6027-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-3

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\82X112619W.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	1.070	26	30	41	rBV	2616396	3071303	9.86%	1.126%
2	2.838	310	320	323	rBV2	163993	371536	1.19%	0.136%
3	2.881	323	327	333	rVB	368766	698128	2.24%	0.256%
4	3.167	364	374	385	rBV	527821	1213965	3.90%	0.445%
5	4.295	550	559	567	rBV2	108213	317912	1.02%	0.117%
6	4.392	567	575	587	rVB	297804	845768	2.71%	0.310%
7	5.295	717	723	734	rVB4	106804	330942	1.06%	0.121%
8	5.490	744	755	763	rBV	390838	1148715	3.69%	0.421%
9	5.654	763	782	788	rVV2	652954	2890711	9.28%	1.060%
10	5.715	788	792	810	rVB	414838	1266973	4.07%	0.465%
11	5.923	815	826	838	rVV2	358790	1093771	3.51%	0.401%
12	6.057	838	848	853	rVV	409728	1036154	3.33%	0.380%
13	6.136	853	861	871	rVV	1814186	4777367	15.33%	1.752%
14	6.252	871	880	888	rVV6	345284	1308582	4.20%	0.480%
15	6.349	889	896	903	rVV3	564190	1742733	5.59%	0.639%
16	6.447	903	912	923	rVB3	268791	981169	3.15%	0.360%
17	6.849	969	978	985	rBV	1006429	2535022	8.14%	0.930%
18	6.929	986	991	1004	rVB3	448828	1314300	4.22%	0.482%
19	7.142	1019	1026	1036	rBV2	136989	334704	1.07%	0.123%
20	7.252	1036	1044	1051	rVB	199088	421608	1.35%	0.155%
21	7.447	1066	1076	1088	rBV3	507520	1536900	4.93%	0.564%
22	7.727	1117	1122	1134	rVB	190012	388816	1.25%	0.143%
23	7.947	1149	1158	1166	rBV5	125456	459191	1.47%	0.168%
24	8.050	1166	1175	1187	rVB2	316046	879228	2.82%	0.322%
25	8.172	1187	1195	1199	rBV	490842	1045977	3.36%	0.384%
26	8.373	1222	1228	1235	rVB4	192170	373213	1.20%	0.137%
27	8.569	1255	1260	1269	rVB5	392872	878529	2.82%	0.322%
28	8.666	1269	1276	1278	rBV2	338258	585968	1.88%	0.215%
29	8.709	1278	1283	1289	rVB	2227295	3492935	11.21%	1.281%
30	8.782	1289	1295	1300	rBV	1902936	2955377	9.49%	1.084%
31	8.983	1323	1328	1332	rBV2	176340	327478	1.05%	0.120%
32	9.032	1332	1336	1341	rVB4	199492	354162	1.14%	0.130%
33	9.160	1350	1357	1362	rVV	308186	532256	1.71%	0.195%
34	9.215	1362	1366	1377	rVB2	765880	1309447	4.20%	0.480%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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 Stop Thrs : 0

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 Title : SW846 8260

35	9.556	1418	1422	1428	rVB3	257932	522067	1.68%	0.191%
36	9.812	1458	1464	1471	rBV3	209847	400496	1.29%	0.147%
37	10.111	1508	1513	1517	rBV	2027358	2693182	8.64%	0.988%
38	10.251	1530	1536	1542	rVB	19110499	25708205	82.51%	9.427%
39	10.355	1545	1553	1559	rVB	22292504	30162054	96.81%	11.060%
40	10.696	1604	1609	1615	rVB	3845231	4882327	15.67%	1.790%
41	11.013	1656	1661	1666	rBV	1486035	1847694	5.93%	0.678%
42	11.135	1676	1681	1686	rBV	1684627	2146345	6.89%	0.787%
43	11.355	1712	1717	1723	rVB	3430043	4072228	13.07%	1.493%
44	11.452	1723	1733	1737	rVV	2823481	5095575	16.35%	1.869%
45	11.501	1737	1741	1753	rVB	1938289	2402949	7.71%	0.881%
46	11.666	1762	1768	1776	rBV	3139185	3872360	12.43%	1.420%
47	11.806	1786	1791	1796	rVB	26129757	31156702	100.00%	11.425%
48	11.916	1804	1809	1811	rBV	362839	484044	1.55%	0.177%
49	11.940	1811	1813	1818	rVB	431039	497670	1.60%	0.182%
50	12.074	1829	1835	1839	rBV2	2218108	2982410	9.57%	1.094%
51	12.141	1839	1846	1850	rVB	6680751	8381906	26.90%	3.074%
52	12.294	1866	1871	1878	rBV	13558101	16821321	53.99%	6.168%
53	12.361	1878	1882	1890	rVB3	2799356	4389046	14.09%	1.609%
54	12.464	1893	1899	1903	rBV2	2121632	2825135	9.07%	1.036%
55	12.513	1903	1907	1913	rVB2	1216338	1777627	5.71%	0.652%
56	12.580	1913	1918	1921	rBV	3160409	4357839	13.99%	1.598%
57	12.604	1921	1922	1927	rVB	1813581	1615395	5.18%	0.592%
58	12.665	1927	1932	1935	rBV	5118399	5946008	19.08%	2.180%
59	12.696	1935	1937	1941	rVB	993655	1081746	3.47%	0.397%
60	12.751	1941	1946	1954	rBV	3874953	5214764	16.74%	1.912%
61	12.897	1965	1970	1975	rVB2	729868	949646	3.05%	0.348%
62	12.976	1975	1983	1986	rBV	4640884	5923230	19.01%	2.172%
63	13.013	1986	1989	1995	rVB	6143616	7090716	22.76%	2.600%
64	13.080	1995	2000	2006	rBV3	557693	1091979	3.50%	0.400%
65	13.190	2015	2018	2020	rBV	361342	393970	1.26%	0.144%
66	13.220	2020	2023	2027	rVV	2992785	3560295	11.43%	1.306%
67	13.263	2027	2030	2033	rVB3	265587	347000	1.11%	0.127%
68	13.306	2033	2037	2039	rBV2	572878	788774	2.53%	0.289%
69	13.342	2039	2043	2048	rVV2	5287168	7495309	24.06%	2.749%
70	13.397	2048	2052	2054	rVV3	494219	683426	2.19%	0.251%
71	13.421	2054	2056	2061	rVB	371260	409391	1.31%	0.150%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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72	13.476	2061	2065	2072	rVB	1140102	1455731	4.67%	0.534%
73	13.555	2072	2078	2081	rBV2	716753	990601	3.18%	0.363%
74	13.598	2081	2085	2088	rVV	1763102	2420614	7.77%	0.888%
75	13.635	2088	2091	2097	rVV4	1561699	2955716	9.49%	1.084%
76	13.696	2097	2101	2102	rVV	824973	1101137	3.53%	0.404%
77	13.720	2102	2105	2110	rVB2	1587621	2347655	7.53%	0.861%
78	13.830	2116	2123	2132	rVB	6823157	8428922	27.05%	3.091%
79	13.921	2132	2138	2142	rVB2	479133	669802	2.15%	0.246%
80	13.976	2142	2147	2152	rBV2	652356	922316	2.96%	0.338%
81	14.025	2152	2155	2159	rVB	551740	615984	1.98%	0.226%
82	14.129	2167	2172	2178	rBV	1294298	1521038	4.88%	0.558%
83	14.269	2190	2195	2201	rBV	1105250	1716080	5.51%	0.629%
84	14.372	2208	2212	2217	rBV2	530235	712780	2.29%	0.261%
85	14.427	2217	2221	2225	rVB	779807	940067	3.02%	0.345%
86	14.689	2261	2264	2269	rVB	835581	1004053	3.22%	0.368%
87	14.836	2283	2288	2297	rVB	1558444	2008130	6.45%	0.736%

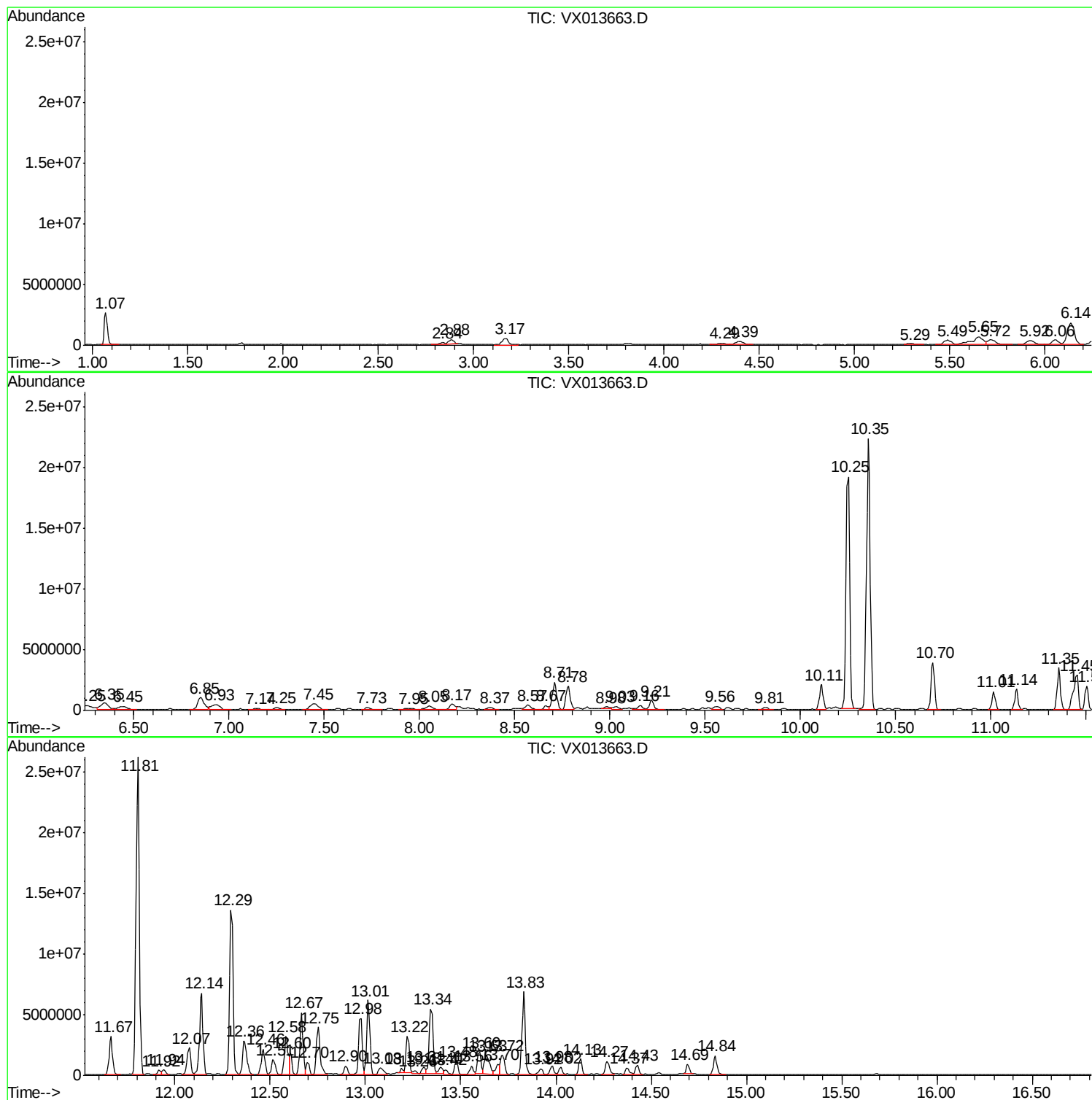
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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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



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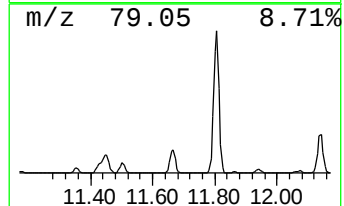
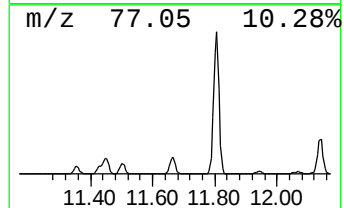
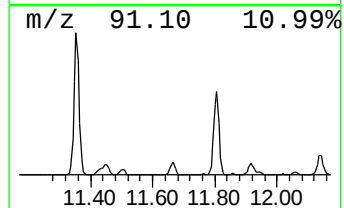
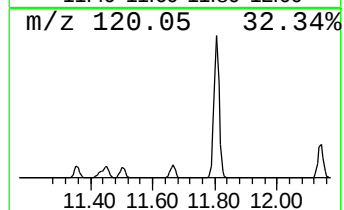
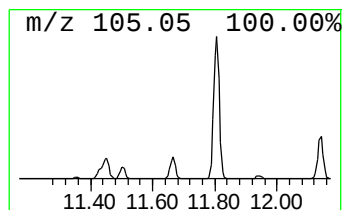
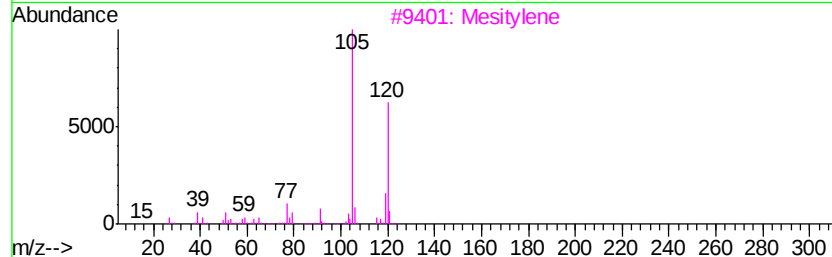
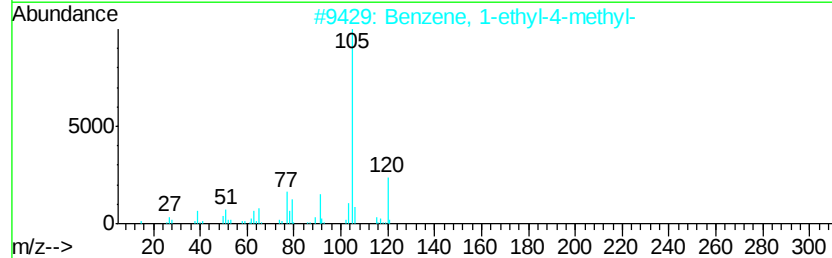
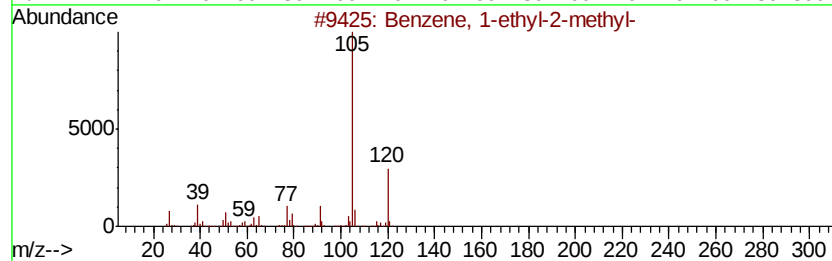
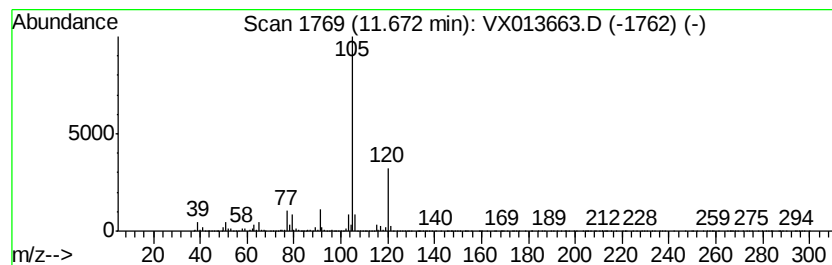
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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.
11.67	64.92 ug/l	3872360	1,4-Dichlorobenzene-d4	12.07

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-4-methyl-	120	C9H12	000622-96-8	93
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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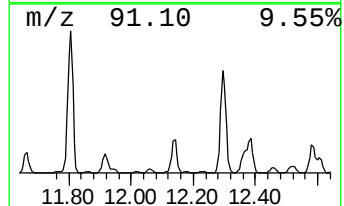
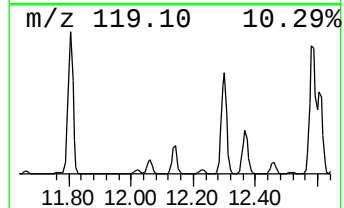
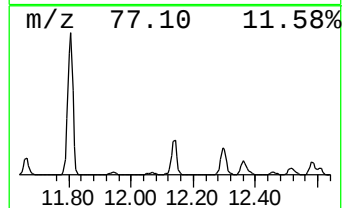
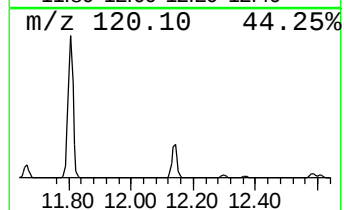
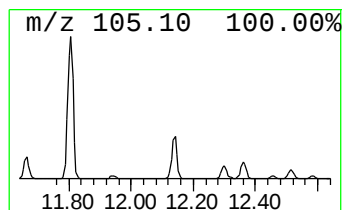
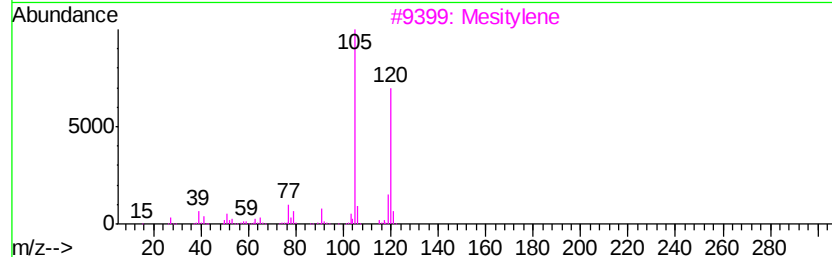
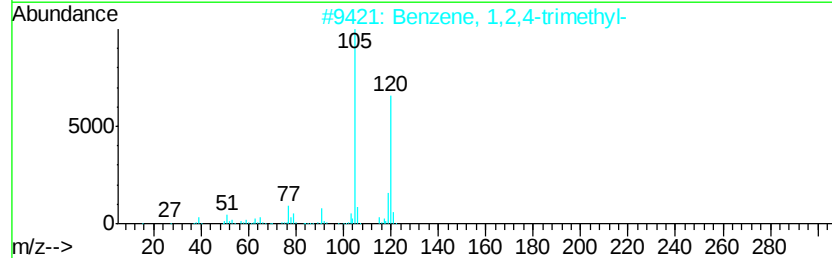
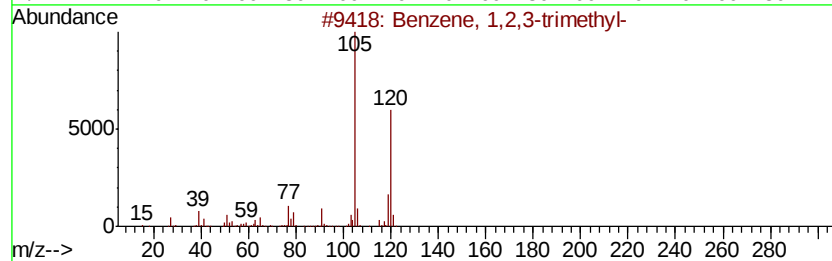
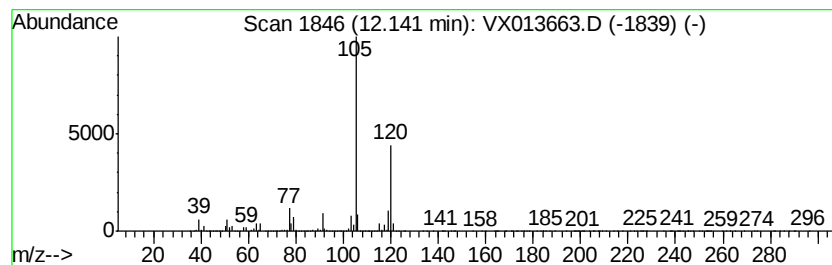
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	140.52 ug/l	8381910	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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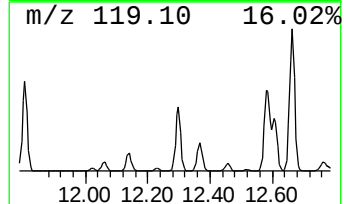
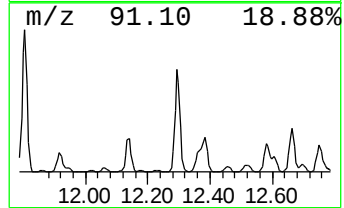
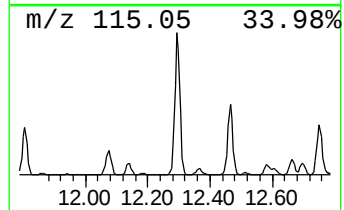
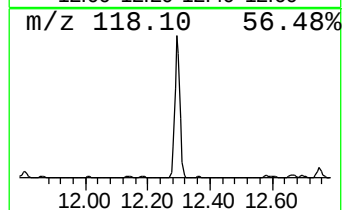
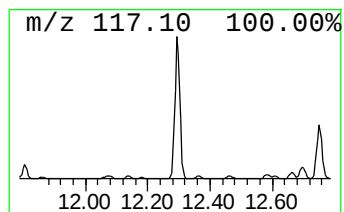
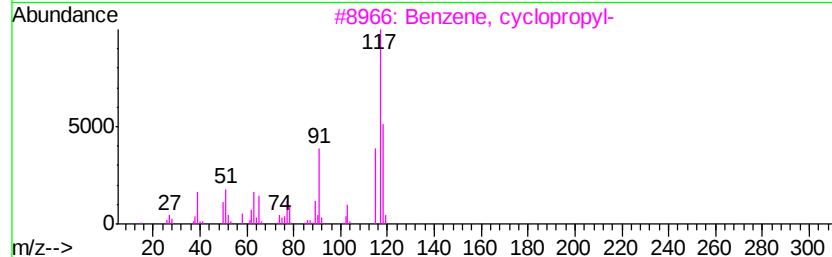
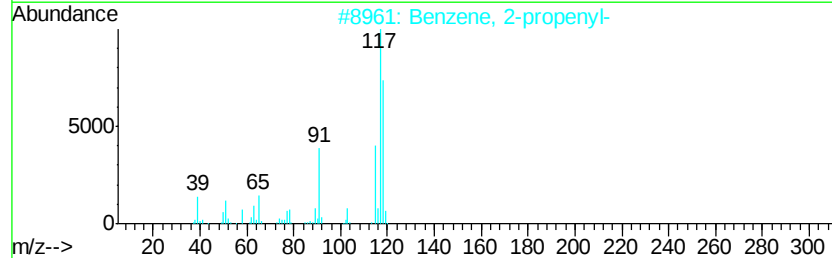
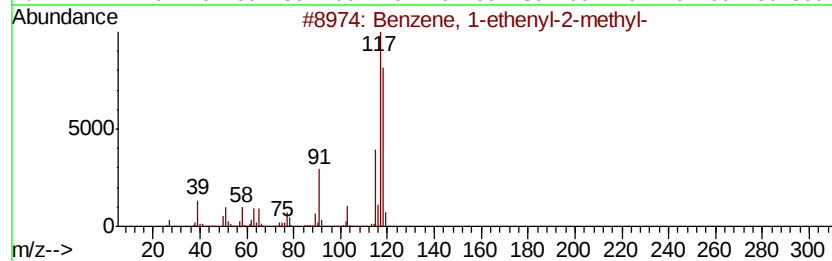
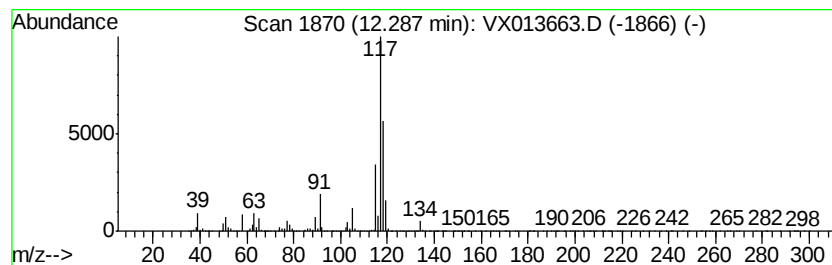
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	282.01 ug/l	16821300	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4		Indane	118	C9H10	000496-11-7	70
5		(Z)-1-Phenylpropene	118	C9H10	000766-90-5	62



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MW-3

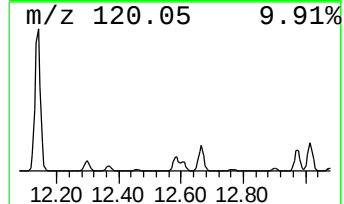
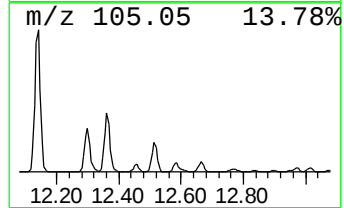
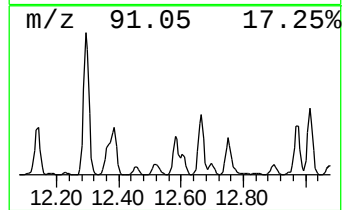
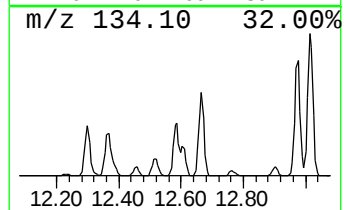
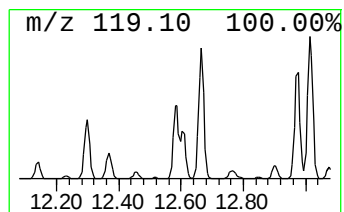
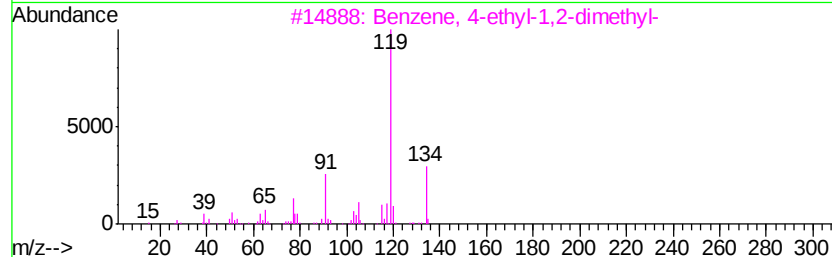
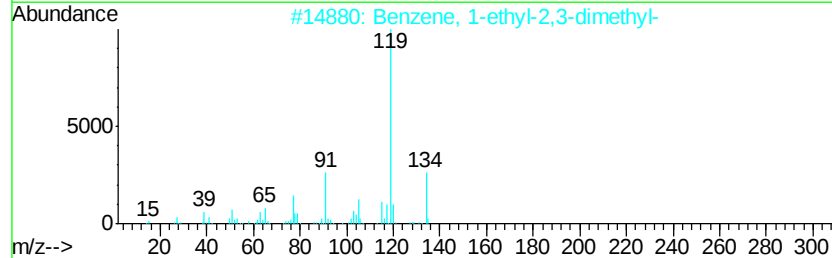
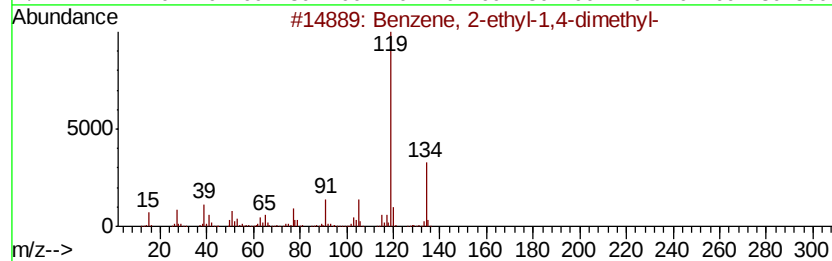
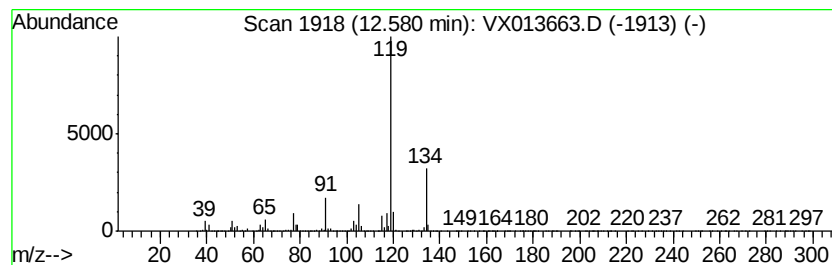
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	73.06 ug/l	4357840	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-3

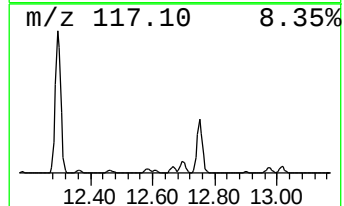
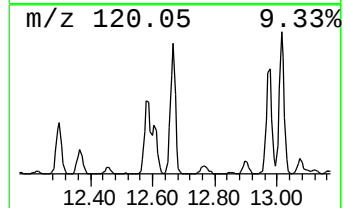
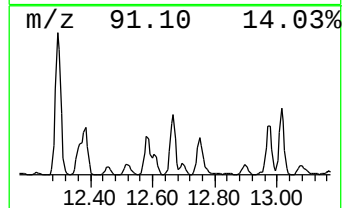
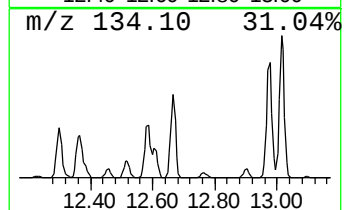
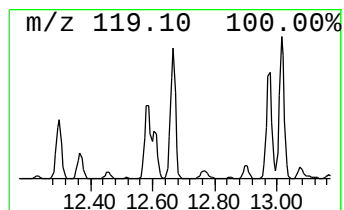
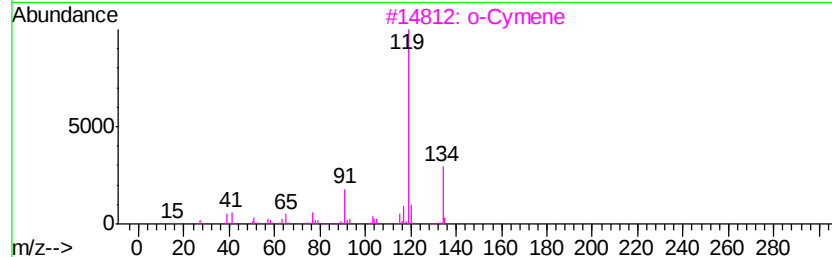
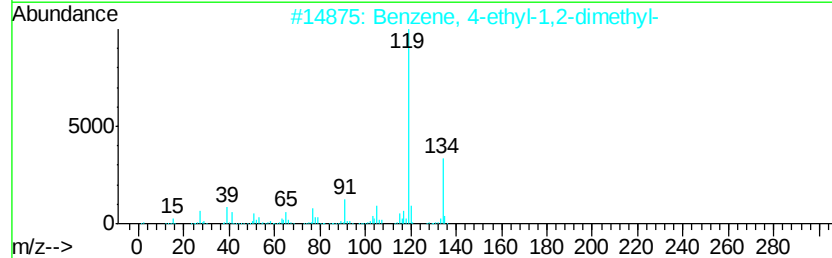
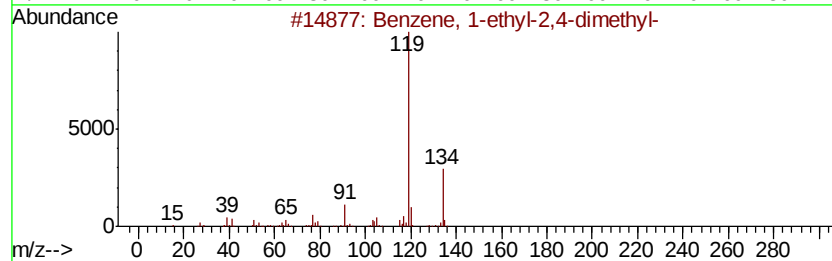
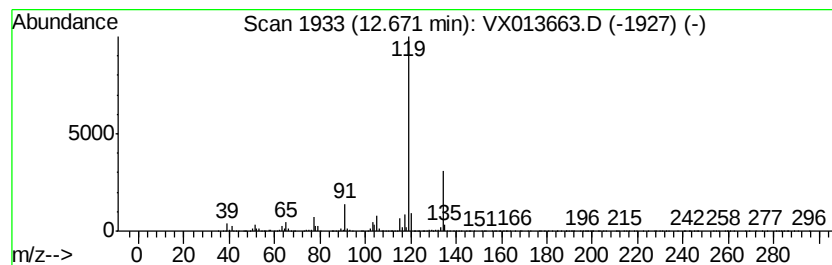
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	99.68 ug/l	5946010	1,4-Dichlorobenzene-d4	12.07

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	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	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

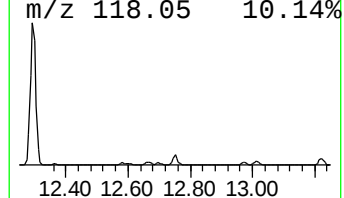
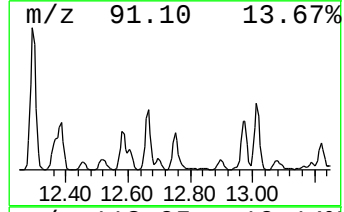
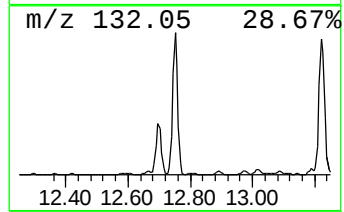
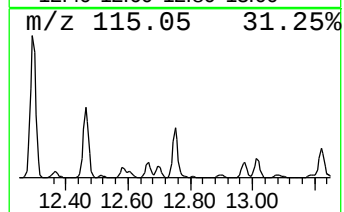
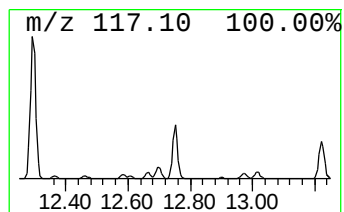
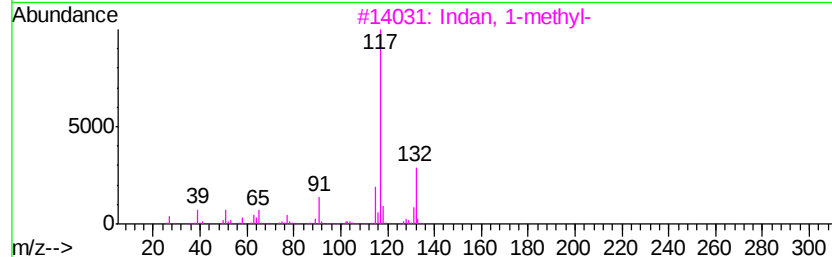
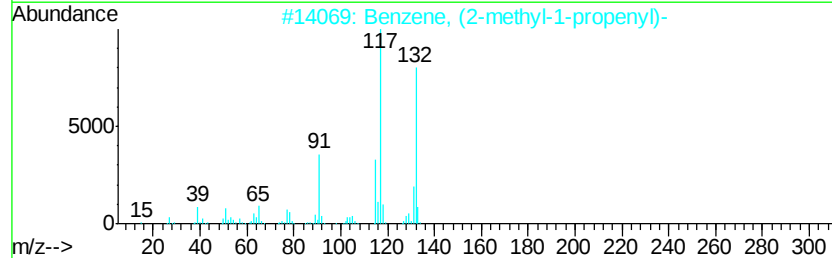
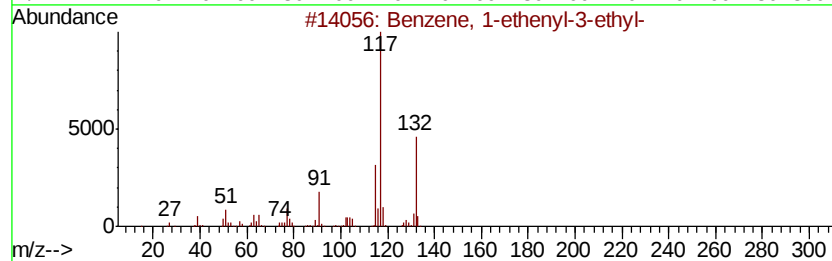
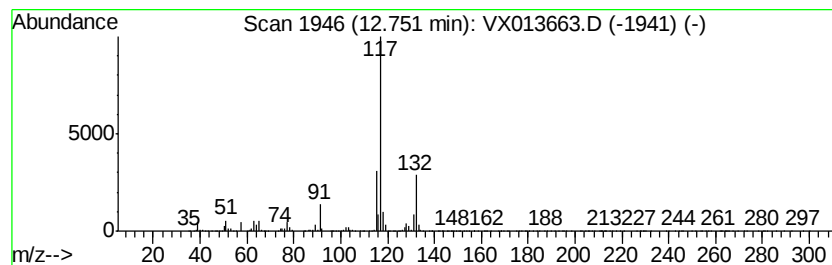
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	87.43 ug/l	5214760	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

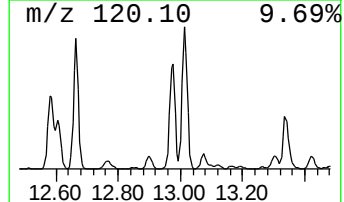
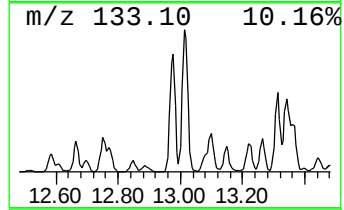
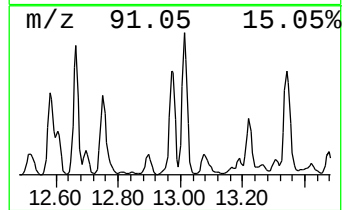
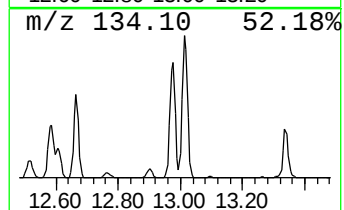
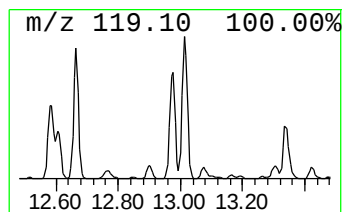
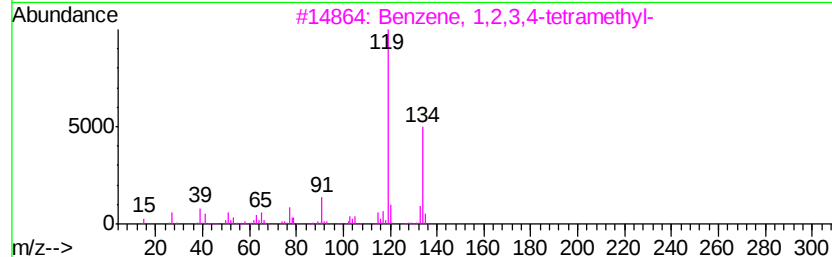
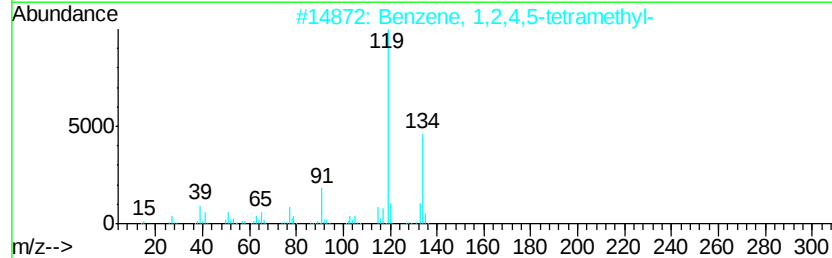
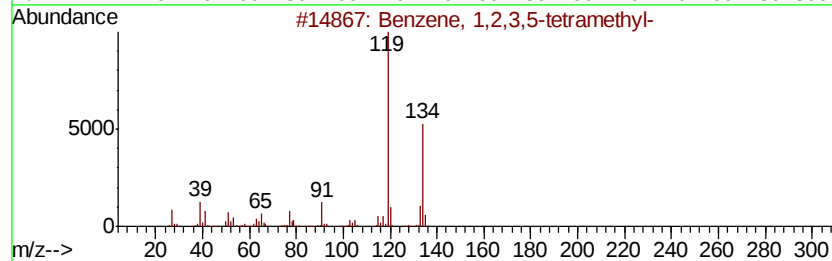
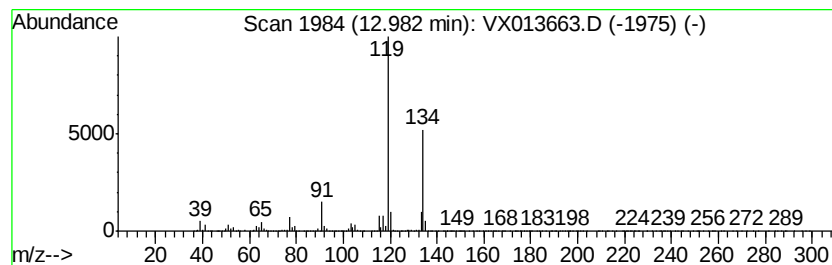
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	99.30 ug/l	5923230	1,4-Dichlorobenzene-d4	12.07

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

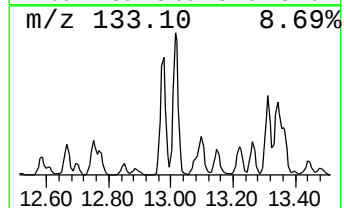
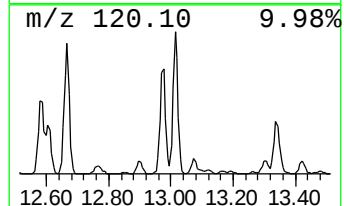
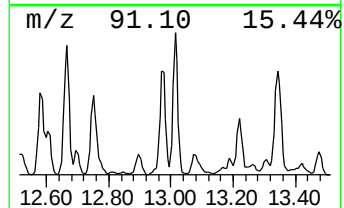
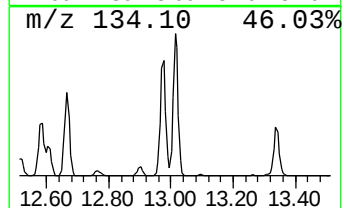
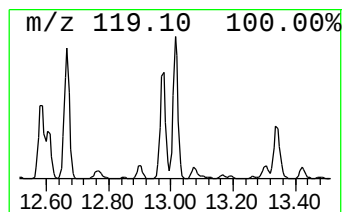
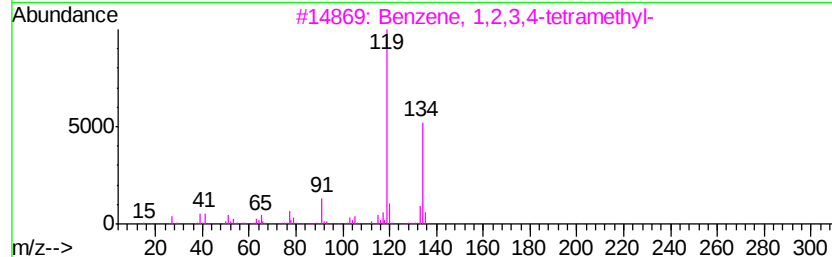
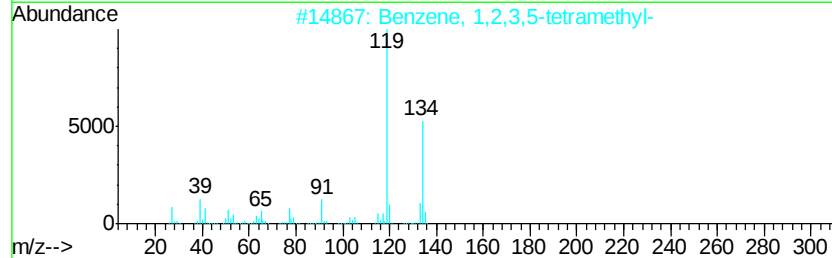
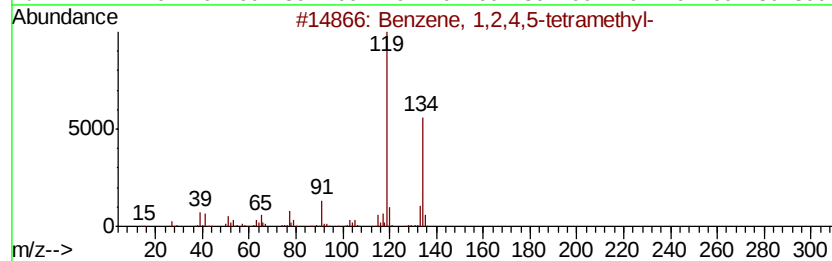
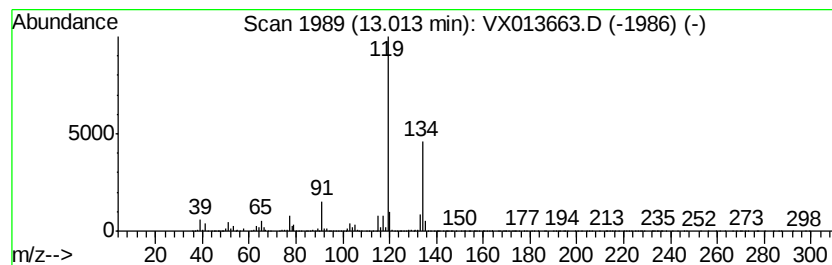
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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.
13.01	118.88 ug/l	7090720	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

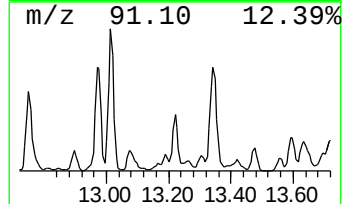
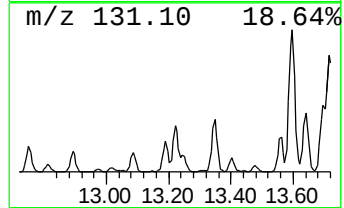
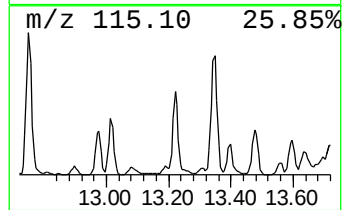
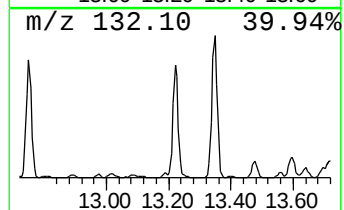
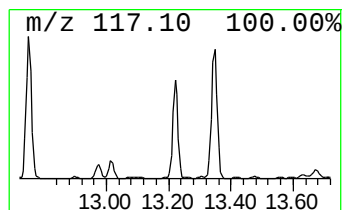
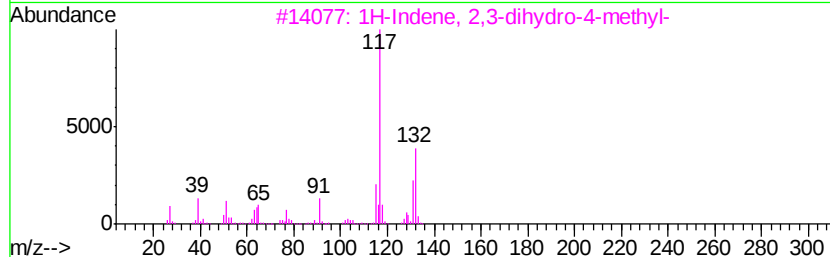
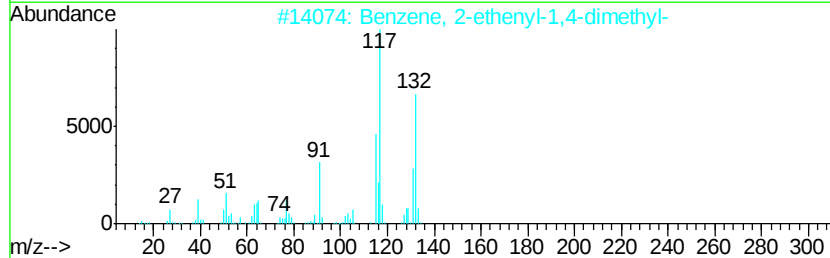
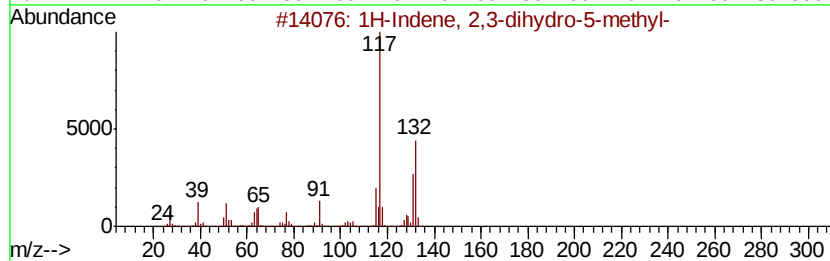
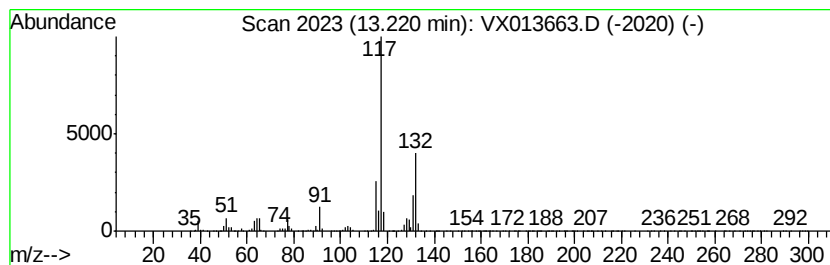
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	59.69 ug/l	3560300	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-3

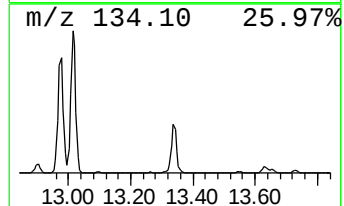
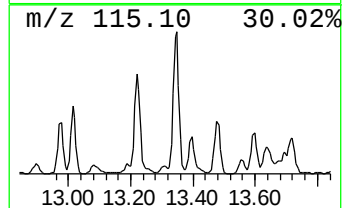
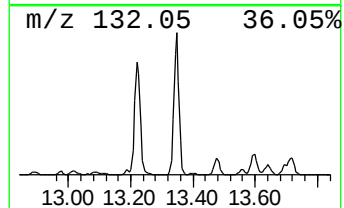
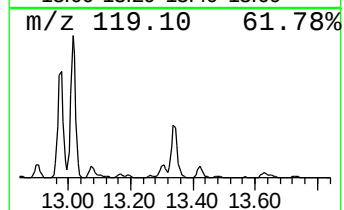
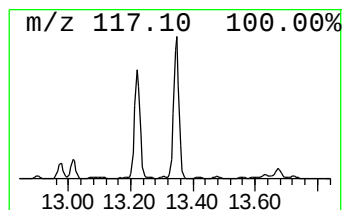
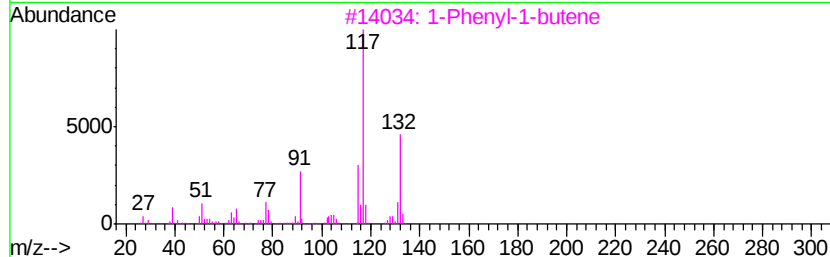
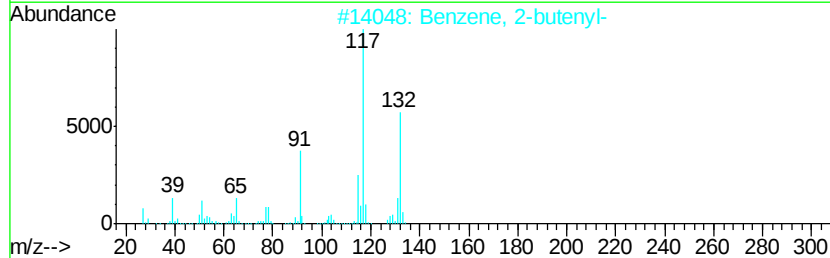
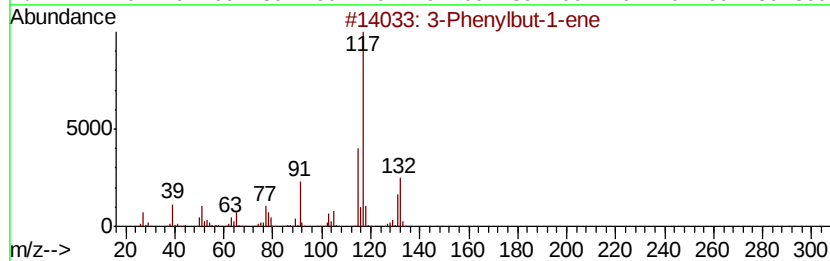
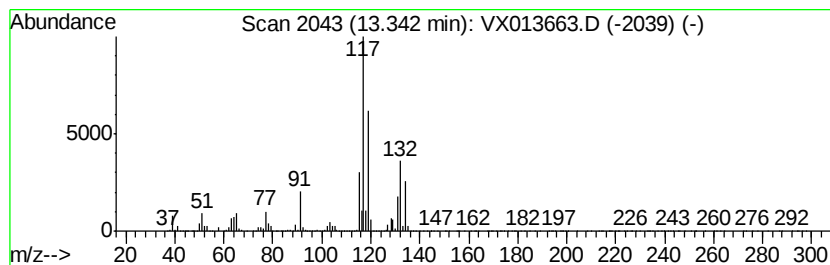
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

Peak Number 10 3-Phenylbut-1-ene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	125.66 ug/l	7495310	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112719\
Data File : VX013663.D
Acq On : 27 Nov 2019 19:44
Operator : JC/SP
Sample : K6027-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.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-...	11.67	64.9	ug/l	3872360	4	12.07	2982410	50.0
Benzene, 1,2,3-tr...	12.14	140.5	ug/l	8381910	4	12.07	2982410	50.0
Benzene, 1-etheny...	12.29	282.0	ug/l	16821300	4	12.07	2982410	50.0
Benzene, 2-ethyl-...	12.58	73.1	ug/l	4357840	4	12.07	2982410	50.0
Benzene, 1-ethyl-...	12.67	99.7	ug/l	5946010	4	12.07	2982410	50.0
Benzene, 1-etheny...	12.75	87.4	ug/l	5214760	4	12.07	2982410	50.0
Benzene, 1,2,3,5-...	12.98	99.3	ug/l	5923230	4	12.07	2982410	50.0
Benzene, 1,2,4,5-...	13.01	118.9	ug/l	7090720	4	12.07	2982410	50.0
1H-Indene, 2,3-di...	13.22	59.7	ug/l	3560300	4	12.07	2982410	50.0
3-Phenylbut-1-ene	13.34	125.7	ug/l	7495310	4	12.07	2982410	50.0