

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX111618\
 Data File : VX006057.D
 Acq On : 16 Nov 2018 18:28
 Operator : JC/MD
 Sample : J5968-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2-111418

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\82X110718W.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	119	rVB	247486	354011	9.08%	0.955%
2	2.001	134	139	150	rVB	57682	84822	2.17%	0.229%
3	2.611	229	239	250	rBV	36142	68498	1.76%	0.185%
4	2.843	270	277	281	rBV	66671	148457	3.81%	0.400%
5	2.891	281	285	289	rVB	72572	127185	3.26%	0.343%
6	3.172	322	331	350	rVB	136588	311576	7.99%	0.840%
7	3.812	428	436	446	rVB2	18423	42987	1.10%	0.116%
8	3.922	446	454	461	rBV2	17502	45240	1.16%	0.122%
9	4.397	523	532	545	rVB	146925	419385	10.75%	1.131%
10	5.299	674	680	695	rVB	19538	62304	1.60%	0.168%
11	5.501	699	713	720	rBV2	78080	229094	5.87%	0.618%
12	5.580	720	726	732	rBV3	19413	41359	1.06%	0.112%
13	5.665	732	740	747	rBV	125256	332689	8.53%	0.897%
14	5.927	772	783	796	rBV5	22264	73586	1.89%	0.198%
15	6.068	796	806	819	rVV	83416	219846	5.64%	0.593%
16	6.263	822	838	847	rVV5	38704	149550	3.83%	0.403%
17	6.354	847	853	862	rVV3	42716	121449	3.11%	0.328%
18	6.452	862	869	881	rVB2	26492	73661	1.89%	0.199%
19	6.860	927	936	945	rBV2	189660	488193	12.52%	1.317%
20	6.939	947	949	962	rVB4	28253	57041	1.46%	0.154%
21	7.256	992	1001	1010	rBV	26813	57500	1.47%	0.155%
22	7.458	1020	1034	1046	rBV3	71599	200099	5.13%	0.540%
23	7.957	1105	1116	1123	rBV5	13990	44871	1.15%	0.121%
24	8.177	1144	1152	1154	rBV	21457	39157	1.00%	0.106%
25	8.214	1154	1158	1162	rVB	37479	61429	1.57%	0.166%
26	8.573	1209	1217	1225	rVB2	51815	98205	2.52%	0.265%
27	8.713	1235	1240	1247	rVV	352089	572884	14.69%	1.545%
28	8.787	1247	1252	1257	rVB	183118	279919	7.18%	0.755%
29	9.165	1306	1314	1319	rBV2	27945	52675	1.35%	0.142%
30	9.219	1319	1323	1335	rVB	58551	97735	2.51%	0.264%
31	9.561	1375	1379	1385	rVB3	20341	40630	1.04%	0.110%
32	10.110	1464	1469	1475	rBV	382077	514198	13.18%	1.387%
33	10.256	1487	1493	1503	rVV	2222699	2908538	74.57%	7.846%
34	10.359	1503	1510	1522	rVB	2538220	3417833	87.62%	9.219%

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35	10.701	1560	1566	1581	rVB	991129	1302202	33.38%	3.513%
36	11.018	1612	1618	1626	rBV	403679	496079	12.72%	1.338%
37	11.140	1633	1638	1643	rBV	297064	377112	9.67%	1.017%
38	11.359	1669	1674	1681	rVB	979656	1156210	29.64%	3.119%
39	11.432	1681	1686	1688	rBV	690237	916361	23.49%	2.472%
40	11.506	1694	1698	1705	rVB	755282	890568	22.83%	2.402%
41	11.670	1719	1725	1735	rBV	630632	812714	20.84%	2.192%
42	11.810	1743	1748	1753	rVB	3247820	3900588	100.00%	10.521%
43	11.920	1761	1766	1768	rBV	71960	92086	2.36%	0.248%
44	11.945	1768	1770	1778	rVB	90403	102244	2.62%	0.276%
45	12.024	1778	1783	1787	rBV	49825	62104	1.59%	0.168%
46	12.079	1787	1792	1797	rVB	393055	512821	13.15%	1.383%
47	12.140	1797	1802	1807	rBV	751011	933868	23.94%	2.519%
48	12.298	1821	1828	1836	rBV	1580726	2075397	53.21%	5.598%
49	12.371	1836	1840	1850	rVB3	381968	601717	15.43%	1.623%
50	12.463	1850	1855	1860	rBV2	108980	152297	3.90%	0.411%
51	12.518	1860	1864	1870	rVB	105408	126544	3.24%	0.341%
52	12.585	1870	1875	1878	rBV	240229	339701	8.71%	0.916%
53	12.609	1878	1879	1884	rVB	175901	148173	3.80%	0.400%
54	12.670	1884	1889	1892	rBV	1005983	1170764	30.02%	3.158%
55	12.701	1892	1894	1898	rVV	227585	241997	6.20%	0.653%
56	12.755	1898	1903	1911	rVV	662121	882769	22.63%	2.381%
57	12.902	1922	1927	1932	rVB2	86011	114112	2.93%	0.308%
58	12.975	1932	1939	1943	rBV	512193	653069	16.74%	1.762%
59	13.018	1943	1946	1952	rVB	320707	375707	9.63%	1.013%
60	13.079	1952	1956	1962	rBV3	55845	98449	2.52%	0.266%
61	13.225	1976	1980	1990	rVB	619613	739927	18.97%	1.996%
62	13.316	1990	1995	1996	rVV2	43747	59738	1.53%	0.161%
63	13.347	1996	2000	2006	rVV2	1071141	1526834	39.14%	4.118%
64	13.402	2006	2009	2011	rVV2	58398	78208	2.01%	0.211%
65	13.426	2011	2013	2018	rVV	57413	62530	1.60%	0.169%
66	13.481	2018	2022	2029	rVB	161772	199530	5.12%	0.538%
67	13.560	2029	2035	2038	rBV2	52604	74244	1.90%	0.200%
68	13.603	2038	2042	2045	rVV	117800	172131	4.41%	0.464%
69	13.639	2045	2048	2053	rVV3	132573	218058	5.59%	0.588%
70	13.700	2053	2058	2059	rVV2	57807	85414	2.19%	0.230%
71	13.719	2059	2061	2067	rVB2	114799	174433	4.47%	0.471%

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Integration Parameters: RTEINT.P

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72	13.834	2073	2080	2089	rBV	1221894	1473853	37.79%	3.976%
73	13.926	2089	2095	2099	rVB2	41427	59744	1.53%	0.161%
74	13.981	2099	2104	2108	rBV2	49885	71558	1.83%	0.193%
75	14.030	2108	2112	2116	rVB	40663	48000	1.23%	0.129%
76	14.133	2124	2129	2135	rBV	81130	97925	2.51%	0.264%
77	14.273	2147	2152	2162	rBV	64766	115979	2.97%	0.313%
78	14.377	2165	2169	2174	rVV	38802	55179	1.41%	0.149%
79	14.432	2174	2178	2182	rVB	46915	56614	1.45%	0.153%
80	14.694	2217	2221	2232	rVB	693457	833323	21.36%	2.248%
81	14.840	2239	2245	2254	rBV	357220	452412	11.60%	1.220%
82	15.688	2378	2384	2387	rBV	30187	44700	1.15%	0.121%

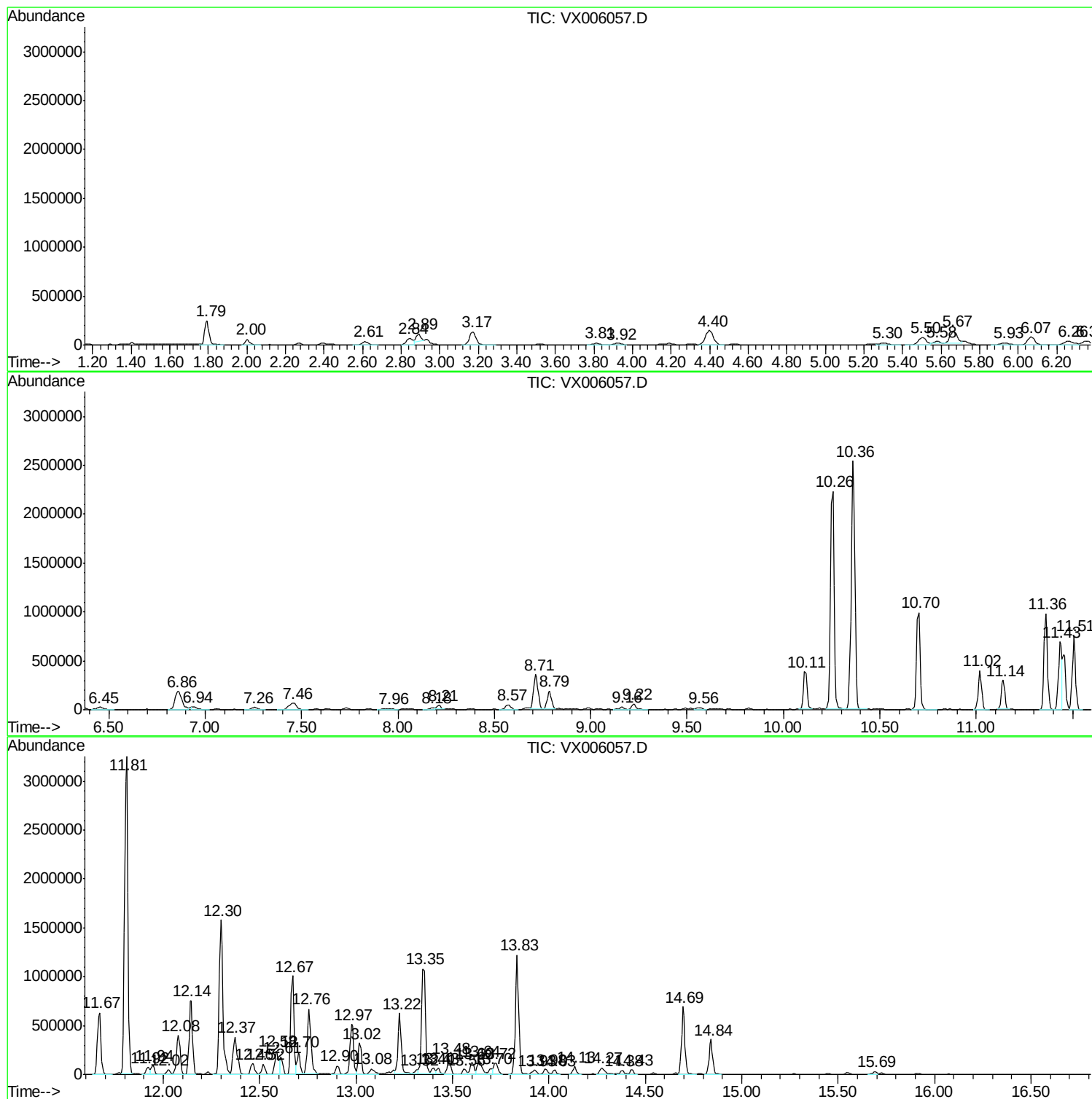
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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



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ALS Vial : 22 Sample Multiplier: 1

Instrument :
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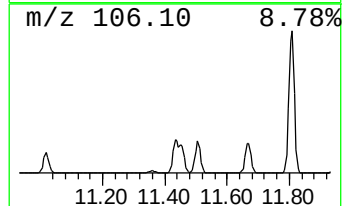
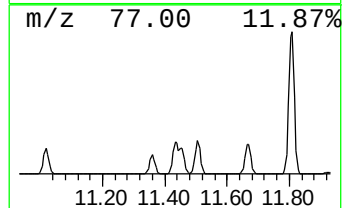
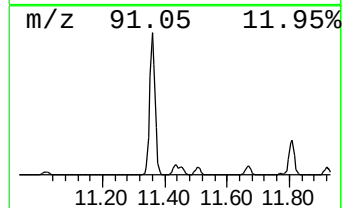
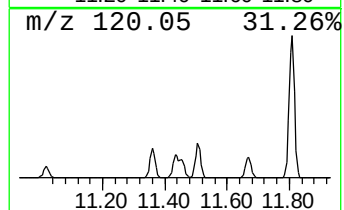
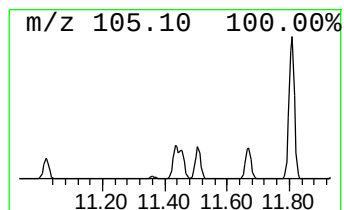
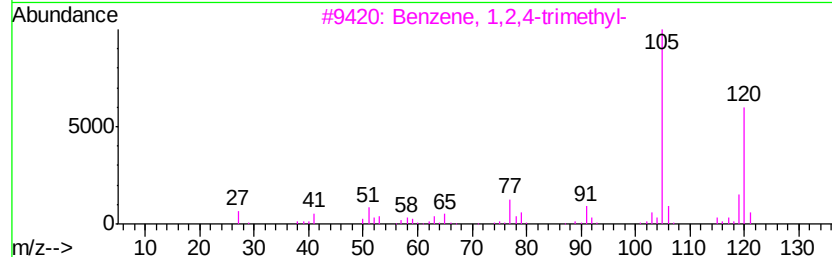
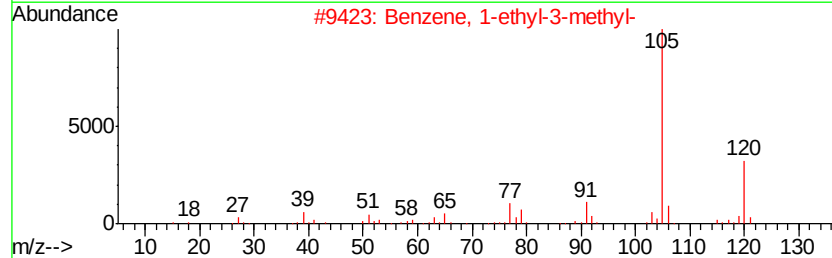
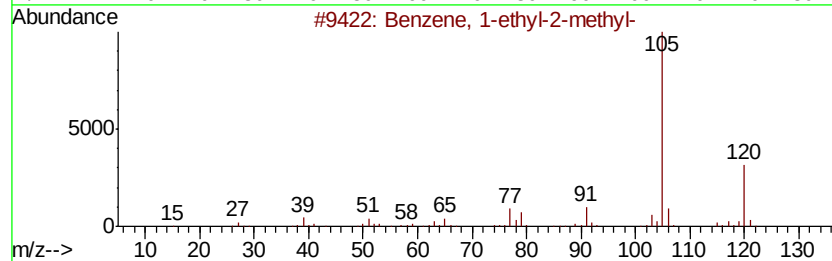
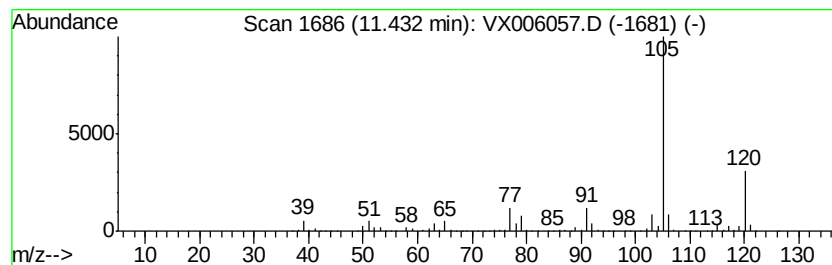
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	89.35 ug/l	916361	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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



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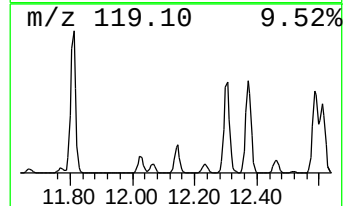
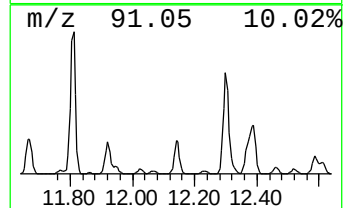
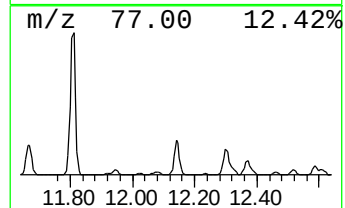
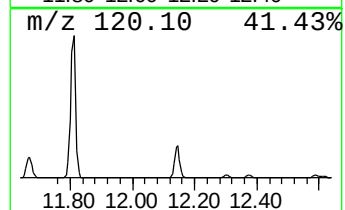
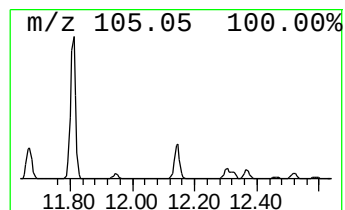
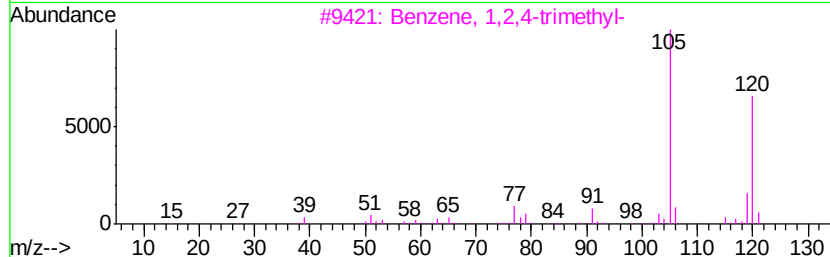
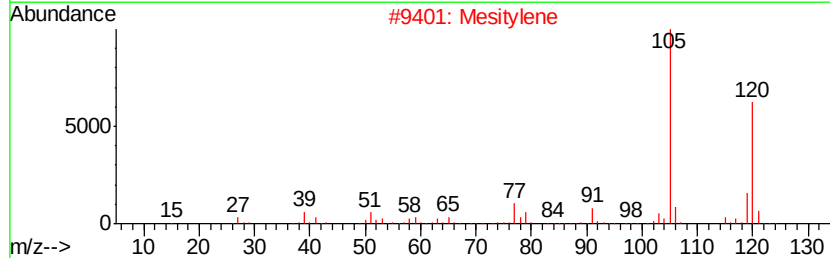
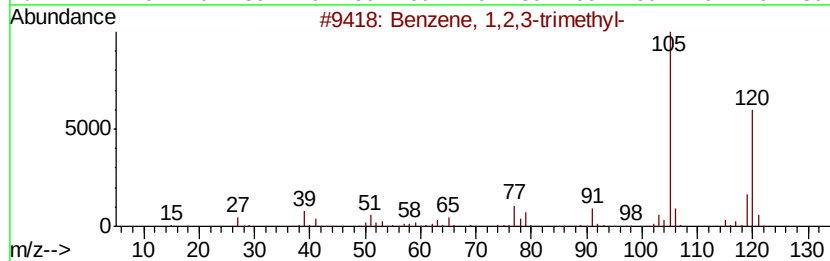
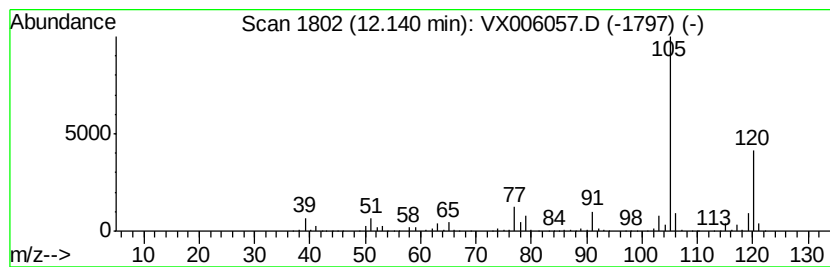
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.14	91.05 ug/l	933868	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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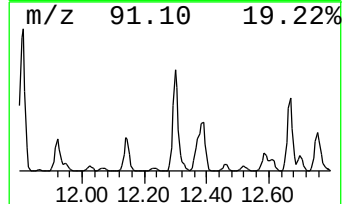
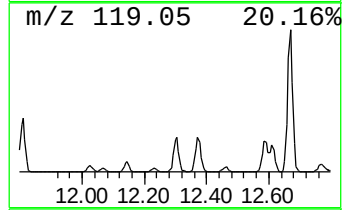
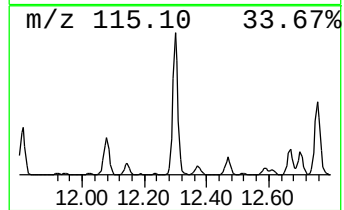
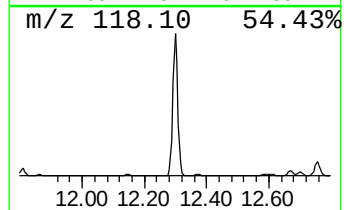
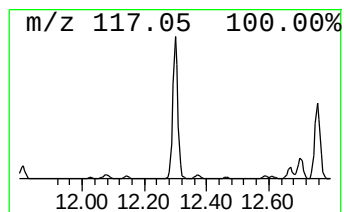
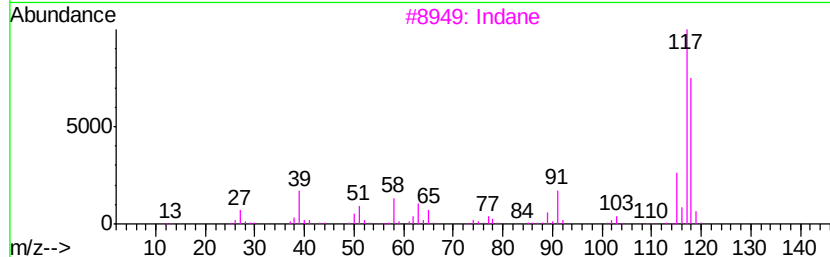
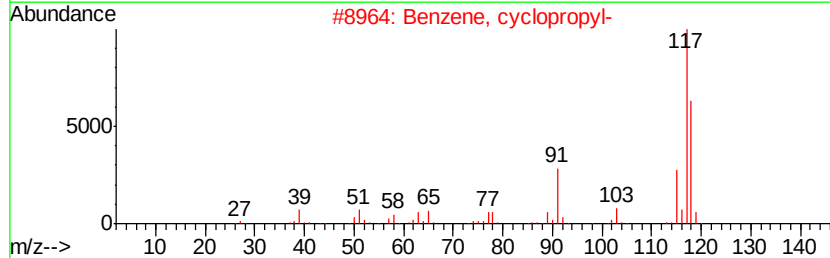
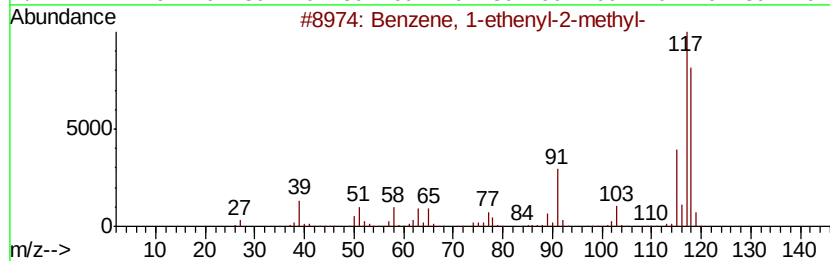
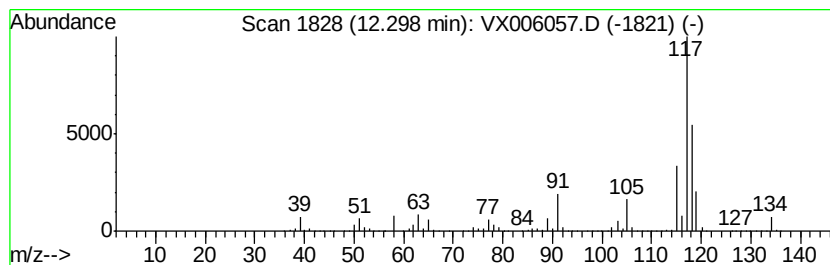
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	202.35 ug/l	2075400	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 86
2		Benzene, cyclopropyl-	118 C9H10	000873-49-4 70
3		Indane	118 C9H10	000496-11-7 70
4		Benzene, 2-propenyl-	118 C9H10	000300-57-2 70
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 58



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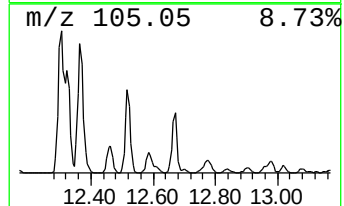
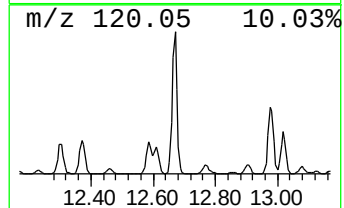
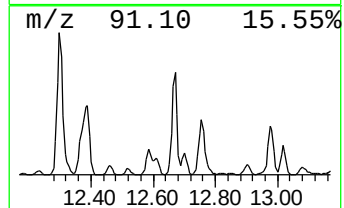
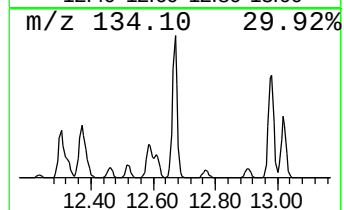
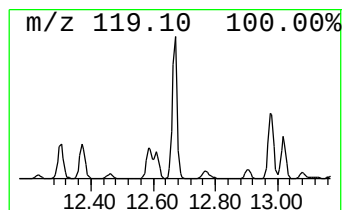
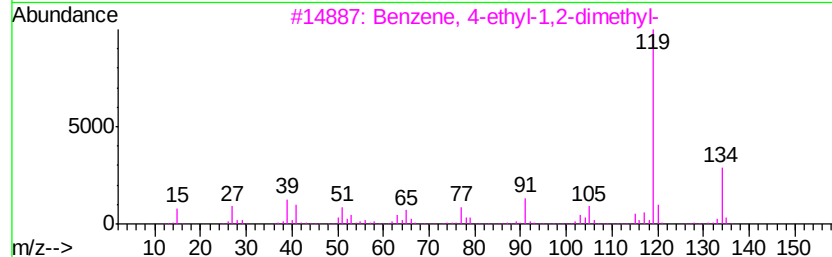
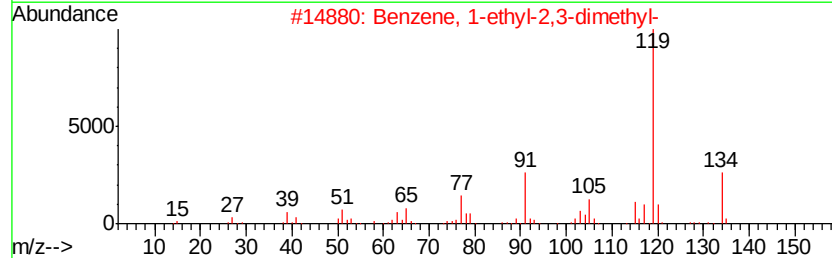
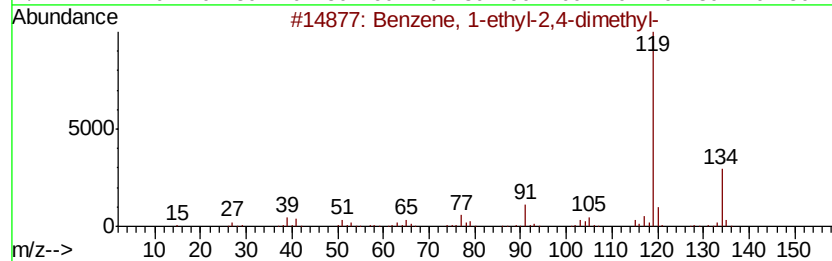
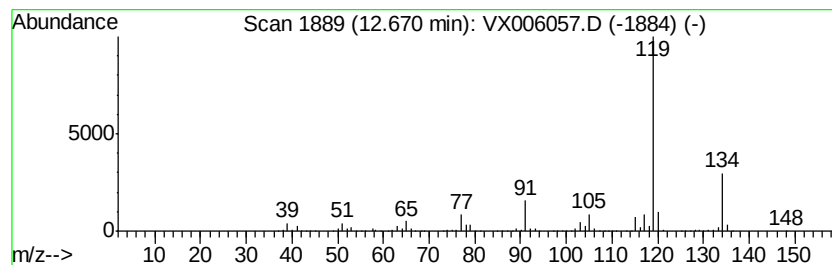
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	114.15 ug/l	1170760	1,4-Dichlorobenzene-d4	12.08

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	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\VX111618\
Data File : VX006057.D
Acq On : 16 Nov 2018 18:28
Operator : JC/MD
Sample : J5968-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

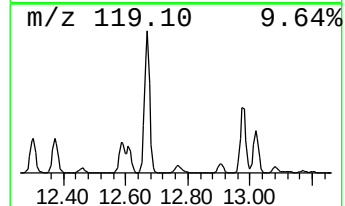
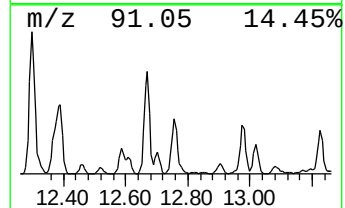
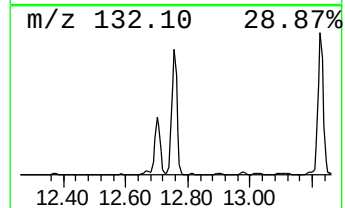
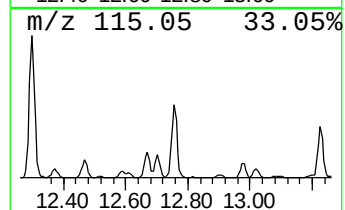
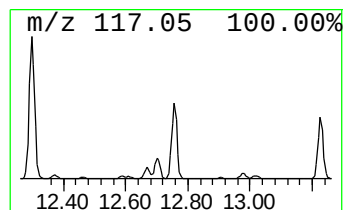
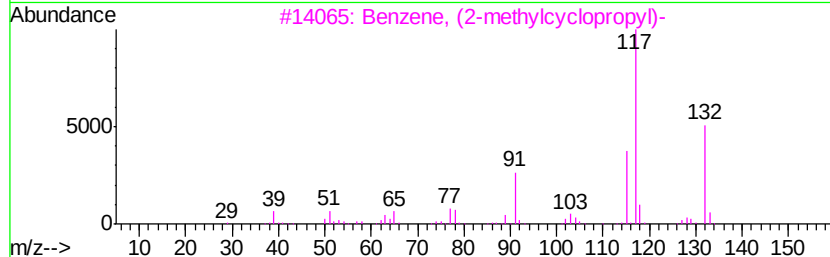
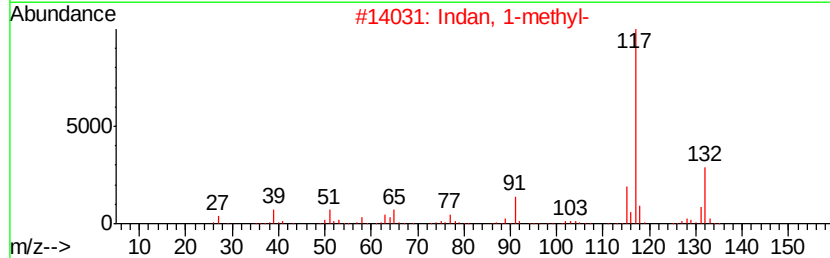
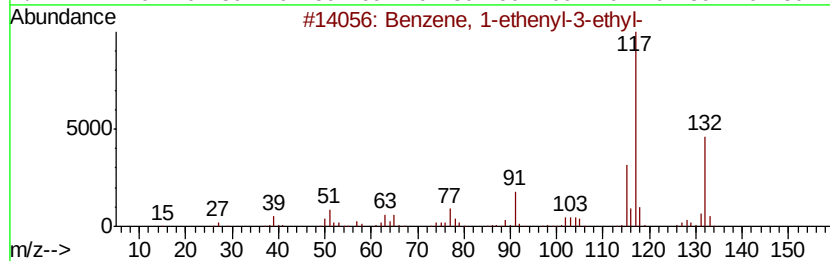
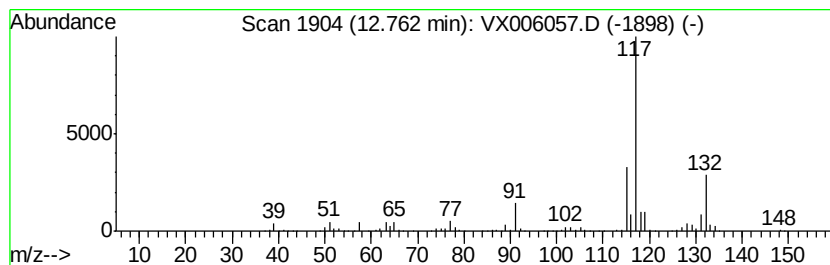
Instrument :
MSVOA_X
ClientSampleId :
MW-2-111418

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
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 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	86.07 ug/l	882769	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 91
2		Indan, 1-methyl-	132 C10H12	000767-58-8 90
3		Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0 90
4		1H-Indene, 2,3-dihydro-2-methyl-	132 C10H12	000824-63-5 87
5		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX111618\
Data File : VX006057.D
Acq On : 16 Nov 2018 18:28
Operator : JC/MD
Sample : J5968-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-2-111418

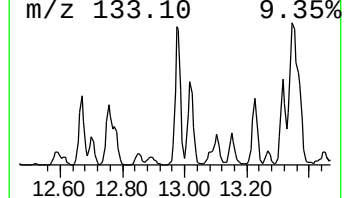
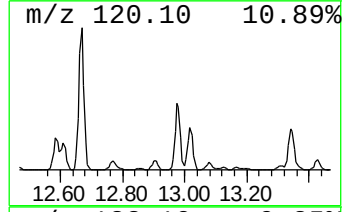
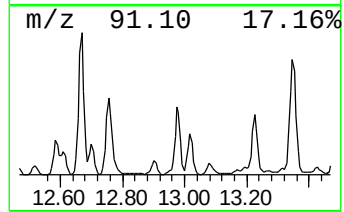
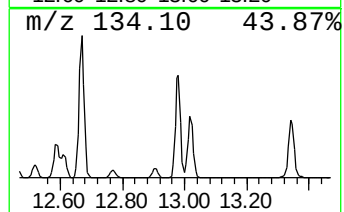
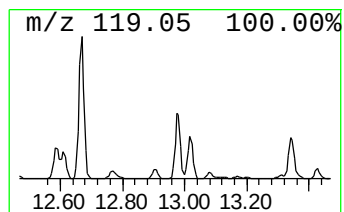
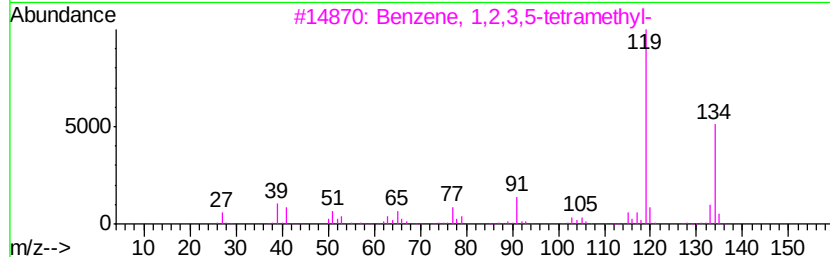
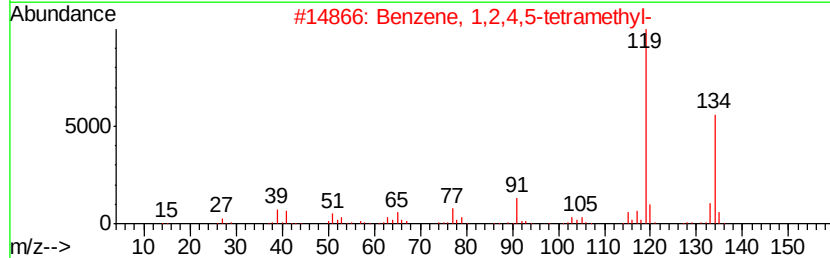
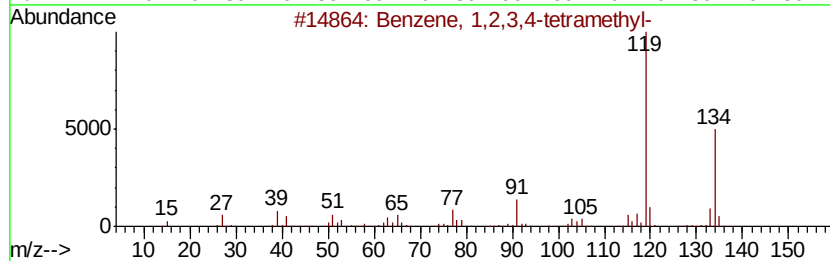
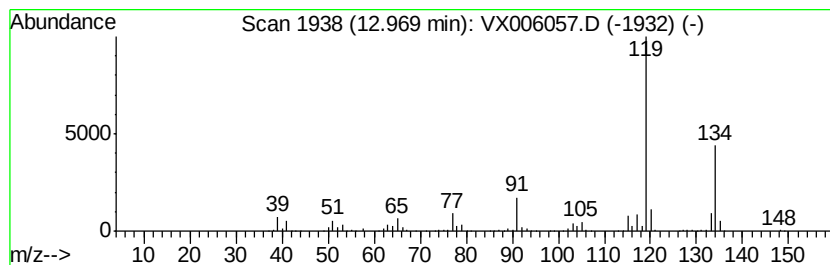
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	63.67 ug/l	653069	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	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX111618\
Data File : VX006057.D
Acq On : 16 Nov 2018 18:28
Operator : JC/MD
Sample : J5968-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

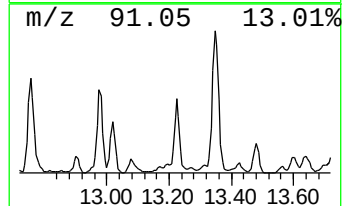
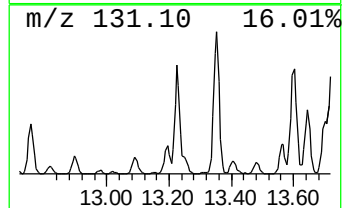
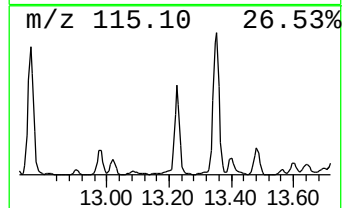
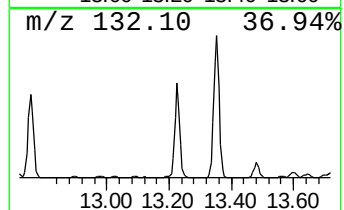
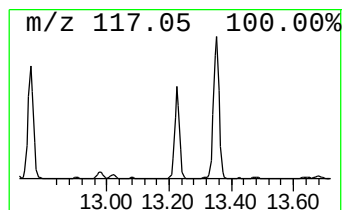
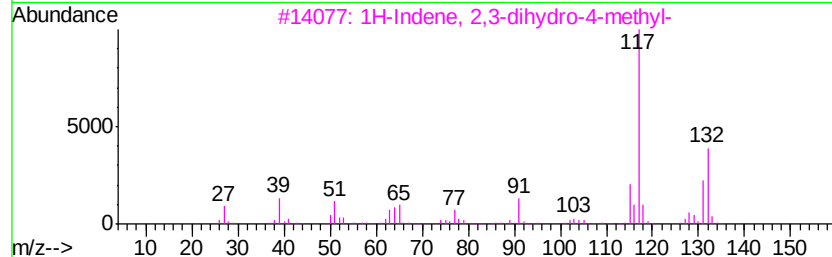
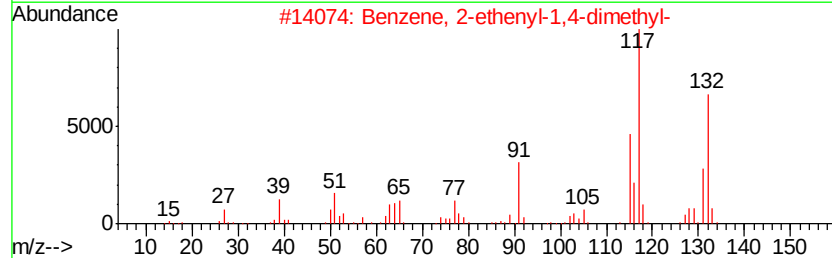
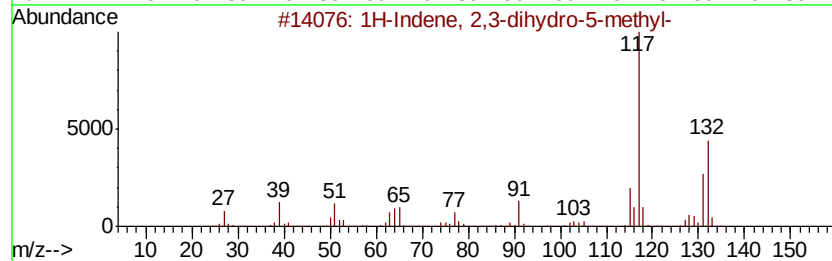
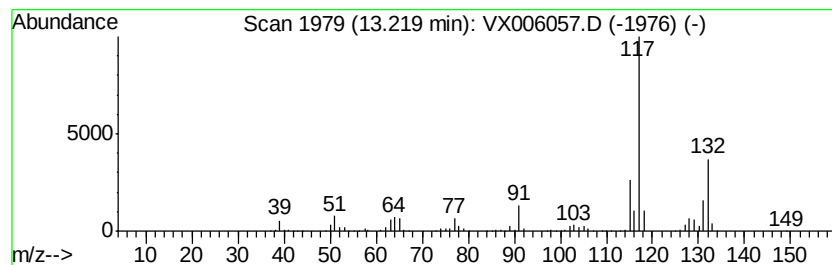
Instrument :
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ClientSampleId :
MW-2-111418

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.22	72.14 ug/l	739927	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX111618\
Data File : VX006057.D
Acq On : 16 Nov 2018 18:28
Operator : JC/MD
Sample : J5968-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

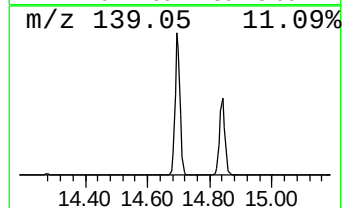
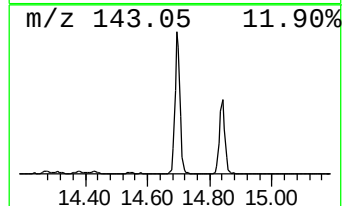
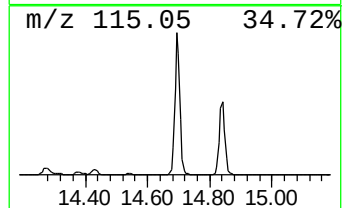
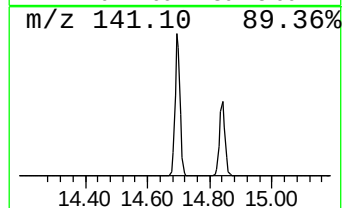
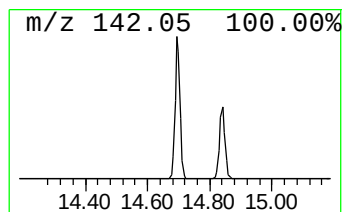
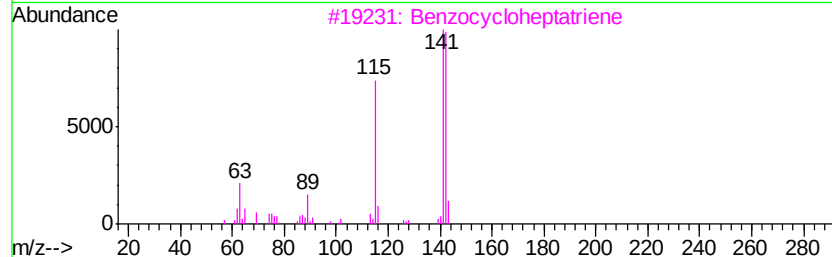
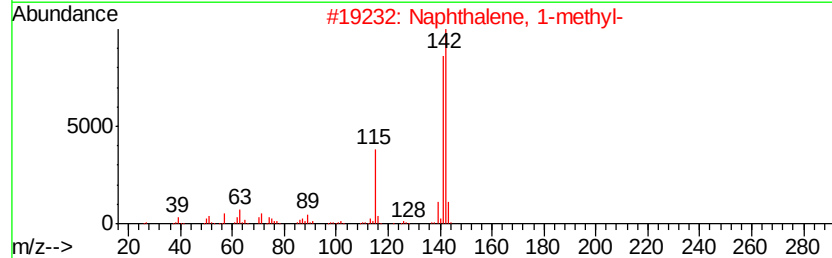
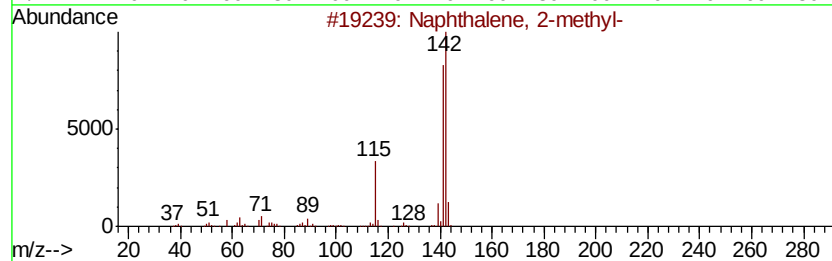
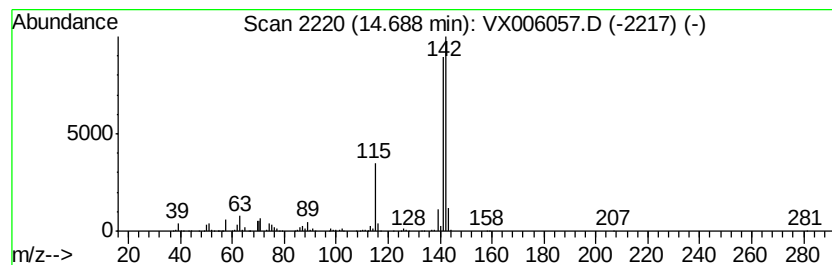
Instrument :
MSVOA_X
ClientSampleId :
MW-2-111418

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	81.25 ug/l	833323	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		Benzocycloheptatriene	142 C11H10	000264-09-5 91
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 47



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111618\
Data File : VX006057.D
Acq On : 16 Nov 2018 18:28
Operator : JC/MD
Sample : J5968-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2-111418

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.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	89.3	ug/l	916361	4	12.08	512821	50.0
Benzene, 1,2,3-tr...	12.14	91.0	ug/l	933868	4	12.08	512821	50.0
Benzene, 1-etheny...	12.30	202.3	ug/l	2075400	4	12.08	512821	50.0
Benzene, 1-ethyl-...	12.67	114.2	ug/l	1170760	4	12.08	512821	50.0
Benzene, 1-etheny...	12.76	86.1	ug/l	882769	4	12.08	512821	50.0
Benzene, 1,2,3,4-...	12.97	63.7	ug/l	653069	4	12.08	512821	50.0
1H-Indene, 2,3-di...	13.22	72.1	ug/l	739927	4	12.08	512821	50.0
Naphthalene, 1-me...	14.69	81.3	ug/l	833323	4	12.08	512821	50.0