

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112118\
 Data File : VX006130.D
 Acq On : 21 Nov 2018 18:19
 Operator : JC/MD
 Sample : J6016-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

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\82X112018W.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	100	105	112	rVB	137218	196609	3.33%	0.311%
2	2.934	288	292	304	rVB	35281	78754	1.33%	0.125%
3	3.172	324	331	340	rVB	52624	122476	2.07%	0.194%
4	4.397	522	532	545	rVB	74690	214747	3.63%	0.340%
5	5.501	703	713	721	rBV2	205742	601187	10.18%	0.952%
6	5.580	721	726	731	rVV3	46046	130540	2.21%	0.207%
7	5.665	731	740	764	rVB	370057	1118051	18.92%	1.770%
8	5.927	774	783	793	rBV6	17897	60839	1.03%	0.096%
9	6.068	796	806	814	rVV	233411	597470	10.11%	0.946%
10	6.141	814	818	827	rVV	33224	82567	1.40%	0.131%
11	6.257	829	837	843	rVV4	19701	65122	1.10%	0.103%
12	6.354	847	853	861	rVV5	23653	68887	1.17%	0.109%
13	6.860	925	936	959	rBV	551876	1317901	22.31%	2.087%
14	7.458	1024	1034	1045	rBV2	88823	217847	3.69%	0.345%
15	8.665	1225	1232	1235	rBV2	35621	66028	1.12%	0.105%
16	8.713	1235	1240	1250	rVB	1126552	1746189	29.56%	2.765%
17	10.116	1464	1470	1477	rBV	1079368	1497882	25.35%	2.372%
18	10.835	1579	1588	1592	rBV2	38118	62953	1.07%	0.100%
19	11.018	1612	1618	1623	rBV	298953	373836	6.33%	0.592%
20	11.140	1632	1638	1643	rBV	931560	1218804	20.63%	1.930%
21	11.359	1669	1674	1685	rBV	598085	720879	12.20%	1.141%
22	11.670	1720	1725	1732	rVB4	34612	59452	1.01%	0.094%
23	11.920	1759	1766	1767	rBV	68409	84107	1.42%	0.133%
24	11.945	1767	1770	1775	rVB	243176	299516	5.07%	0.474%
25	12.079	1787	1792	1805	rBV	1286841	1550536	26.25%	2.455%
26	12.231	1813	1817	1822	rVB	86178	110410	1.87%	0.175%
27	12.298	1823	1828	1834	rVV	2282414	2790465	47.23%	4.418%
28	12.371	1835	1840	1848	rVB2	257113	487524	8.25%	0.772%
29	12.463	1849	1855	1860	rBV	259895	336372	5.69%	0.533%
30	12.670	1880	1889	1892	rBV	993339	1265664	21.42%	2.004%
31	12.701	1892	1894	1898	rVV	499323	531118	8.99%	0.841%
32	12.755	1898	1903	1910	rVB2	1248187	1779050	30.11%	2.817%
33	12.896	1920	1926	1931	rVB2	170573	262534	4.44%	0.416%
34	12.975	1931	1939	1944	rBV	1205523	1578768	26.72%	2.500%

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 ALS Vial : 20 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Integration Parameters: RTEINT.P

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

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

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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
 Title : SW846 8260

35	13.030	1944	1948	1952	rVB3	41287	65212	1.10%	0.103%
36	13.078	1952	1956	1963	rVB3	204313	296453	5.02%	0.469%
37	13.152	1963	1968	1972	rBV2	123316	207881	3.52%	0.329%
38	13.194	1972	1975	1977	rVV2	317335	383077	6.48%	0.607%
39	13.225	1977	1980	1990	rVB	795589	1050825	17.79%	1.664%
40	13.316	1990	1995	1996	rBV2	143004	199659	3.38%	0.316%
41	13.347	1996	2000	2006	rVV2	3256867	4432959	75.03%	7.019%
42	13.402	2006	2009	2011	rVV2	124924	177493	3.00%	0.281%
43	13.426	2011	2013	2019	rVV2	180792	231072	3.91%	0.366%
44	13.481	2019	2022	2029	rVB2	135967	178506	3.02%	0.283%
45	13.560	2030	2035	2038	rBV2	273608	382136	6.47%	0.605%
46	13.603	2038	2042	2045	rVV	607592	908053	15.37%	1.438%
47	13.639	2045	2048	2052	rVV2	770829	1130382	19.13%	1.790%
48	13.700	2052	2058	2059	rVV2	429625	715822	12.12%	1.133%
49	13.719	2059	2061	2070	rVB2	812610	1177930	19.94%	1.865%
50	13.810	2070	2076	2078	rBV	113877	152294	2.58%	0.241%
51	13.853	2078	2083	2088	rVB3	156983	235680	3.99%	0.373%
52	13.914	2088	2093	2097	rVB	703891	888987	15.05%	1.408%
53	13.987	2097	2105	2109	rBV2	910869	1290275	21.84%	2.043%
54	14.030	2109	2112	2116	rVV	369471	444883	7.53%	0.704%
55	14.060	2116	2117	2124	rVB3	79147	94873	1.61%	0.150%
56	14.133	2124	2129	2132	rBV	718857	868485	14.70%	1.375%
57	14.164	2132	2134	2137	rVB	65766	62773	1.06%	0.099%
58	14.273	2147	2152	2162	rBV3	905958	1852477	31.36%	2.933%
59	14.377	2166	2169	2174	rVV	404047	537864	9.10%	0.852%
60	14.432	2174	2178	2182	rVB	712051	880637	14.91%	1.394%
61	14.475	2182	2185	2191	rVB2	122941	162311	2.75%	0.257%
62	14.542	2191	2196	2200	rBV2	1244037	1665958	28.20%	2.638%
63	14.584	2200	2203	2209	rVB2	65866	97340	1.65%	0.154%
64	14.694	2209	2221	2225	rBV2	1372197	2460833	41.65%	3.896%
65	14.731	2225	2227	2232	rVB	345280	380580	6.44%	0.603%
66	14.786	2232	2236	2240	rBV2	149273	187821	3.18%	0.297%
67	14.840	2240	2245	2249	rVV	4736809	5907889	100.00%	9.354%
68	14.877	2249	2251	2254	rVV2	217832	290374	4.92%	0.460%
69	14.907	2254	2256	2260	rVB2	161842	174690	2.96%	0.277%
70	14.962	2260	2265	2272	rBV3	146215	298752	5.06%	0.473%
71	15.115	2286	2290	2294	rVB2	77372	104960	1.78%	0.166%

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Integration Parameters: RTEINT.P

Integrator: RTE
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Stop Thrs : 0
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72	15.249	2306	2312	2320	rBV	465853	781831	13.23%	1.238%
73	15.322	2320	2324	2331	rVB4	106908	179111	3.03%	0.284%
74	15.444	2338	2344	2347	rBV	338628	502820	8.51%	0.796%
75	15.480	2347	2350	2354	rVV	339520	451536	7.64%	0.715%
76	15.548	2355	2361	2366	rVV	977629	1577818	26.71%	2.498%
77	15.590	2366	2368	2374	rVV5	136998	230856	3.91%	0.366%
78	15.688	2375	2384	2388	rVV	1538742	2552845	43.21%	4.042%
79	15.724	2388	2390	2398	rVB	902563	1220059	20.65%	1.932%
80	15.810	2399	2404	2411	rBV2	73539	149205	2.53%	0.236%
81	15.889	2411	2417	2418	rVV	290739	384551	6.51%	0.609%
82	15.919	2419	2422	2437	rVB	492258	914537	15.48%	1.448%
83	16.072	2441	2447	2459	rVB	434344	742202	12.56%	1.175%
84	16.279	2478	2481	2487	rVB2	71146	118850	2.01%	0.188%
85	16.352	2488	2493	2495	rBV	37739	62861	1.06%	0.100%
86	16.438	2495	2507	2516	rVV4	96666	351013	5.94%	0.556%
87	16.511	2516	2519	2523	rVB3	37251	60584	1.03%	0.096%
88	16.566	2523	2528	2534	rBV	46645	78064	1.32%	0.124%
89	16.657	2534	2543	2544	rBV5	45471	110282	1.87%	0.175%
90	16.694	2544	2549	2554	rVB	99854	177151	3.00%	0.280%
91	16.755	2554	2559	2564	rBV2	114341	240987	4.08%	0.382%

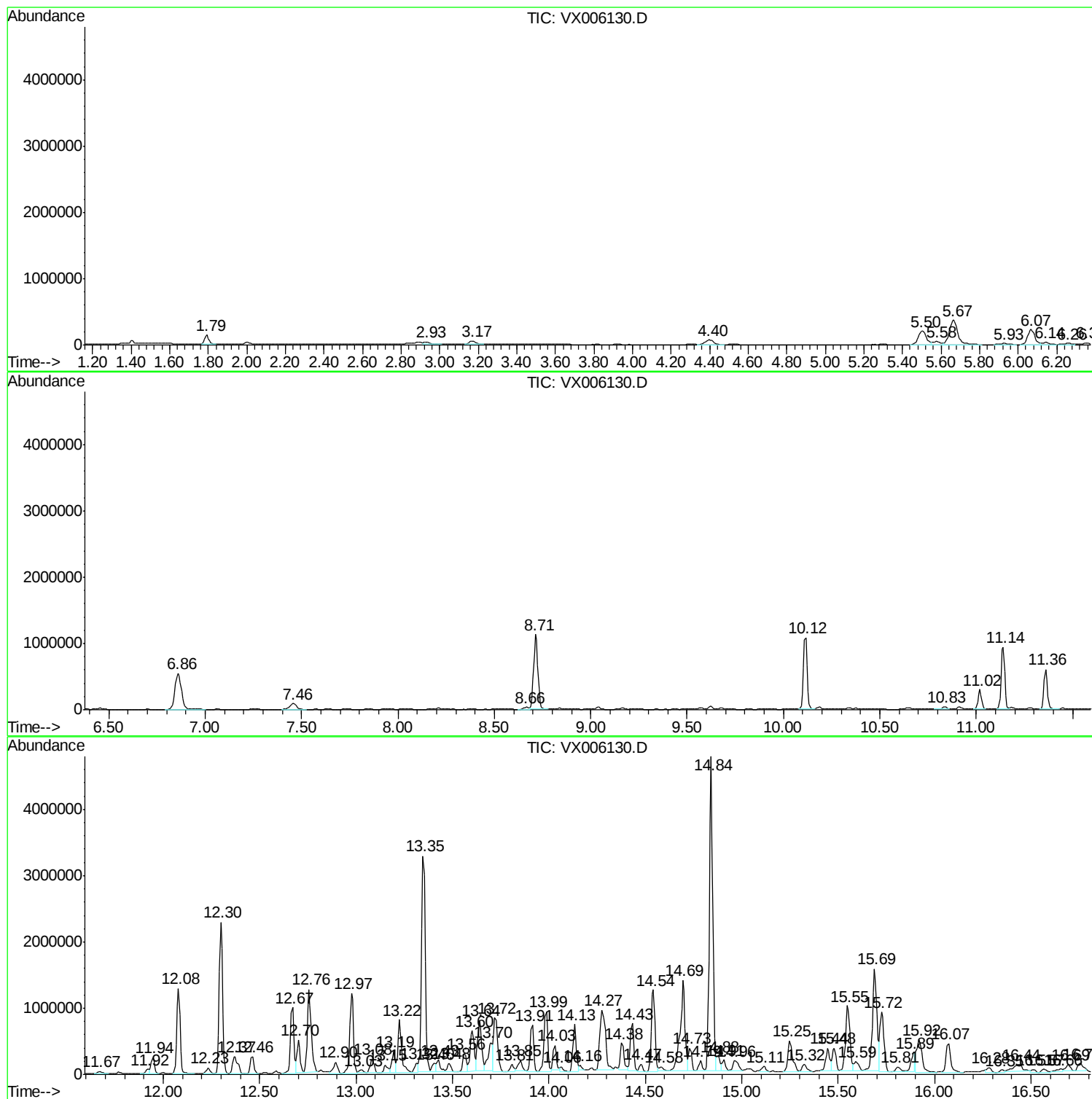
Sum of corrected areas: 63160443

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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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



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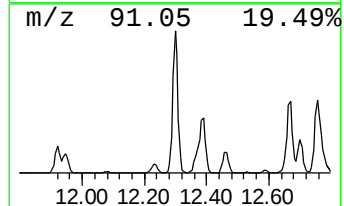
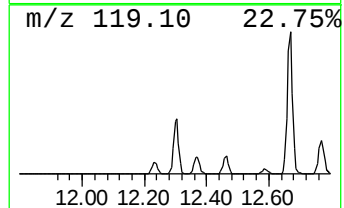
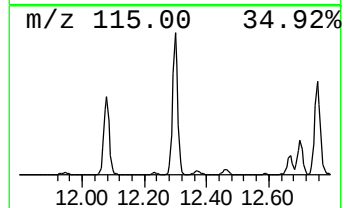
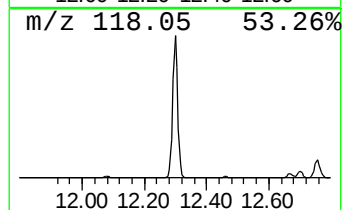
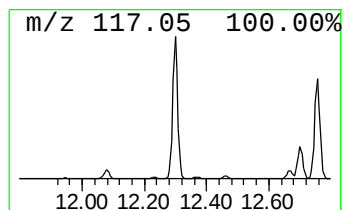
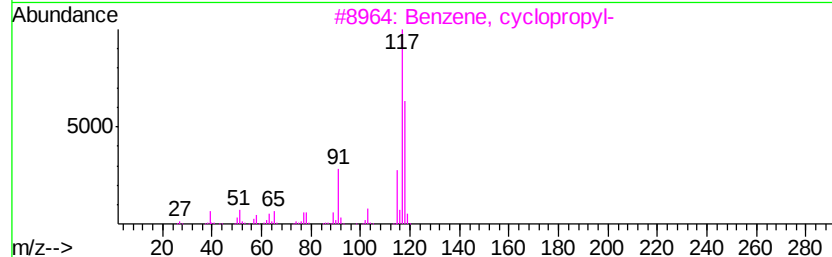
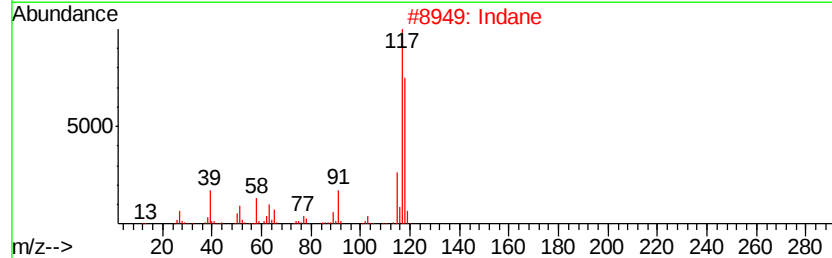
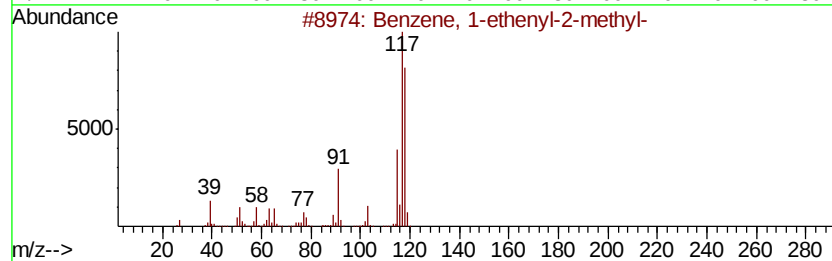
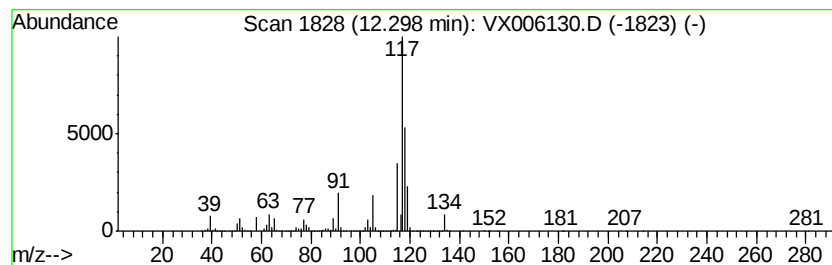
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	89.98 ug/l	2790470	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
2		Indane	118	C9H10	000496-11-7	76
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	62



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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampled :
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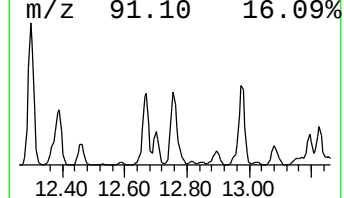
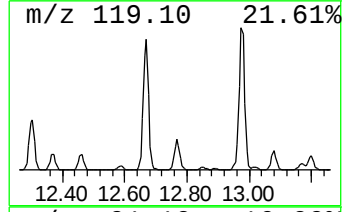
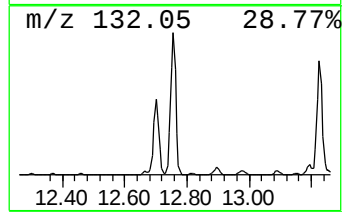
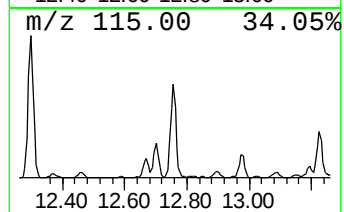
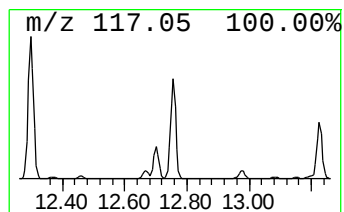
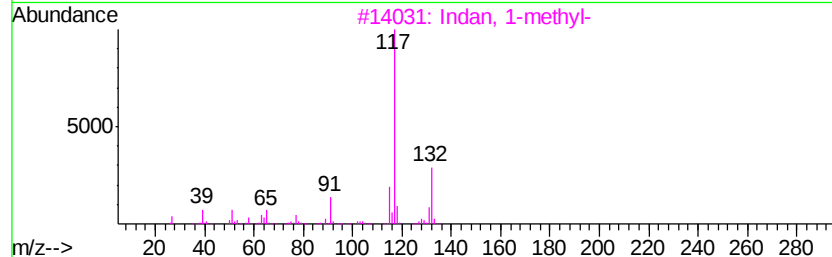
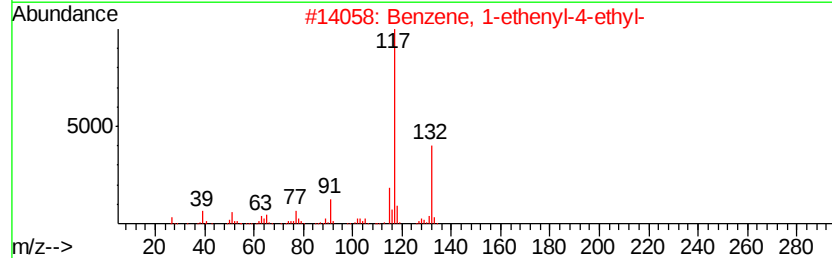
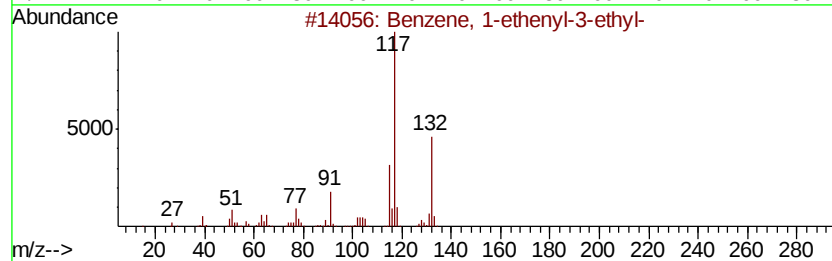
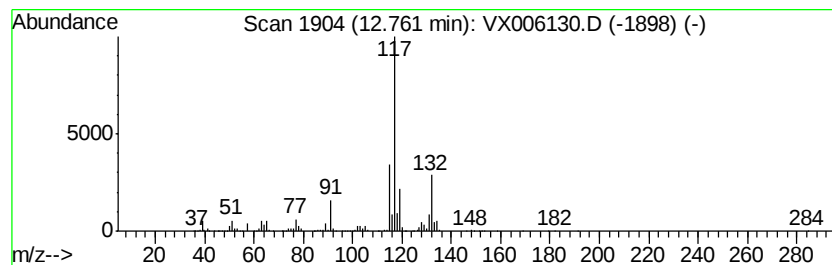
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	57.37 ug/l	1779050	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	83



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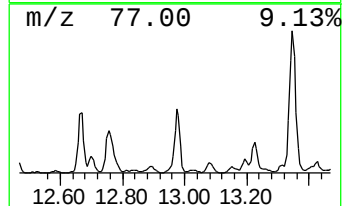
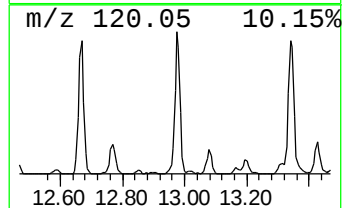
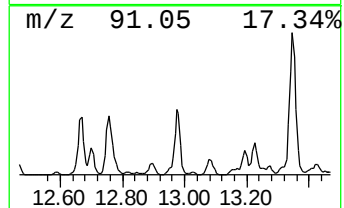
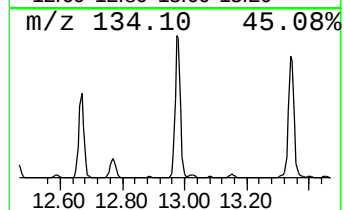
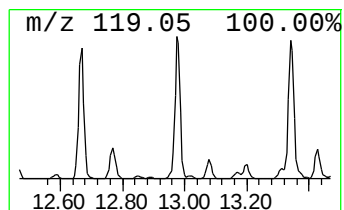
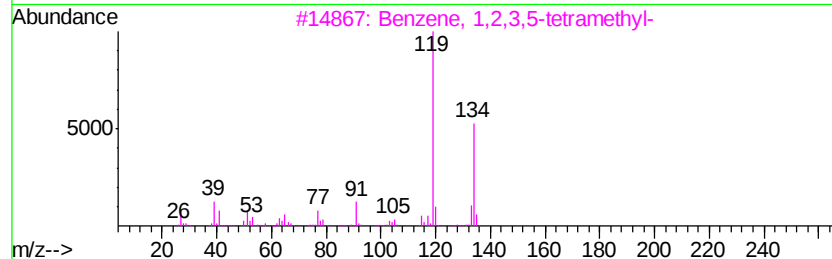
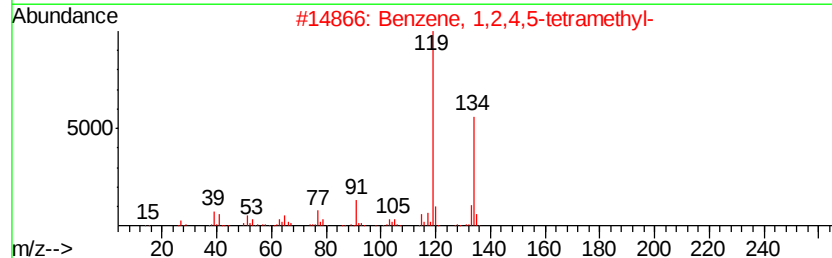
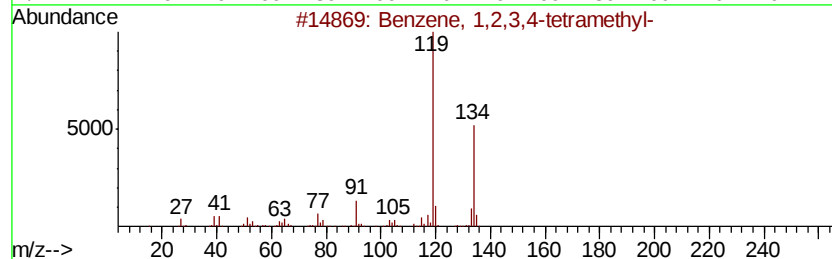
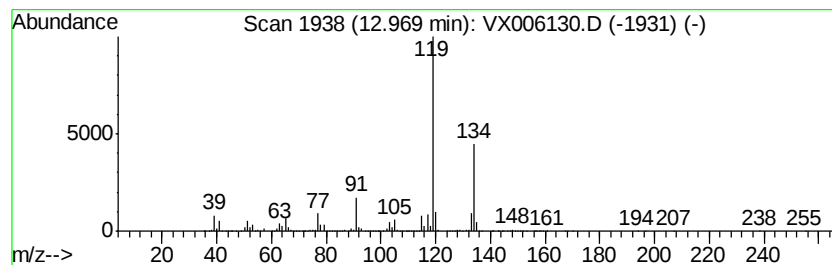
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	50.91 ug/l	1578770	1,4-Dichlorobenzene-d4	12.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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



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ClientSampleId :
MW-1

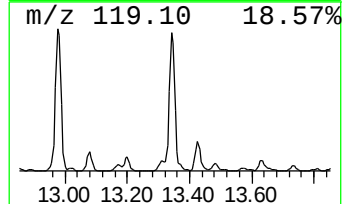
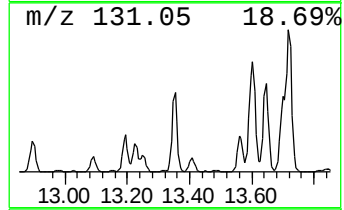
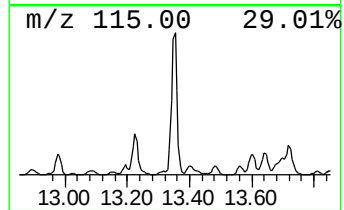
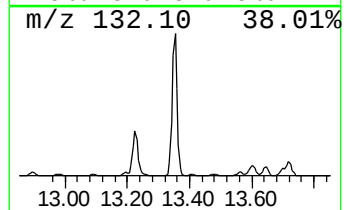
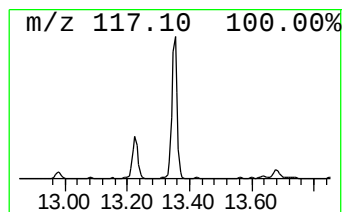
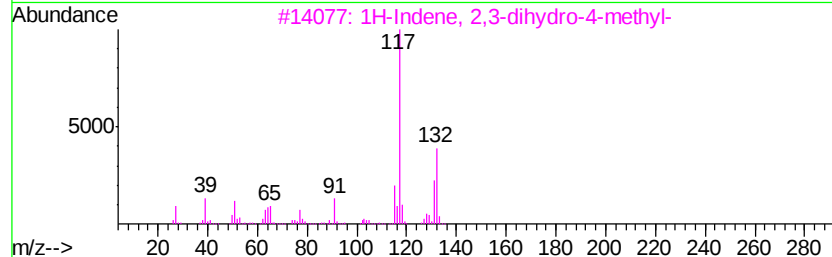
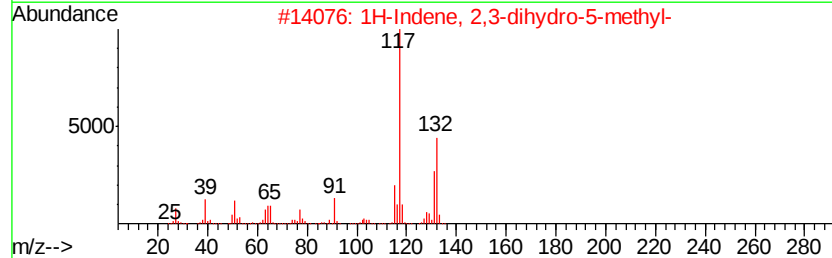
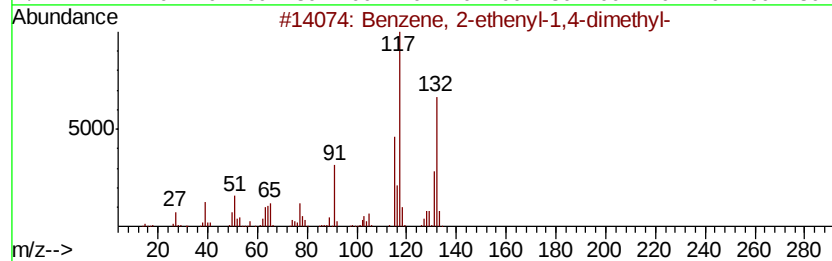
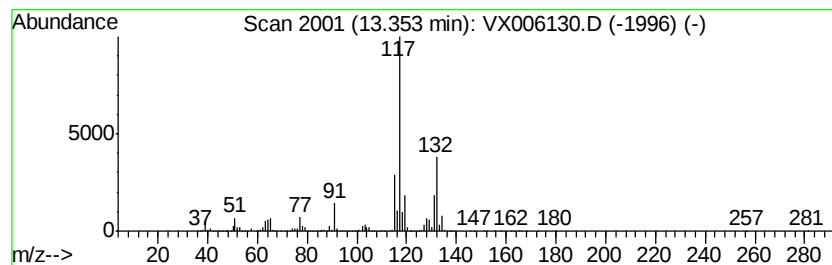
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	142.95 ug/l	4432960	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	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112118\
Data File : VX006130.D
Acq On : 21 Nov 2018 18:19
Operator : JC/MD
Sample : J6016-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

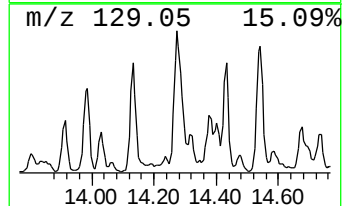
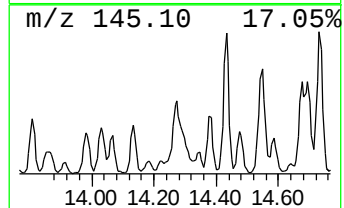
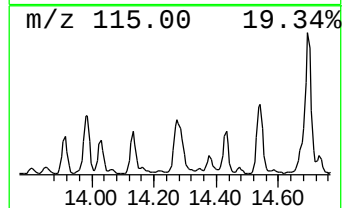
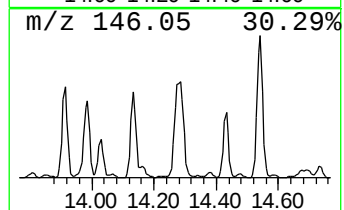
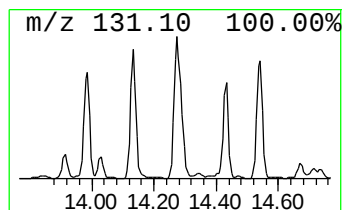
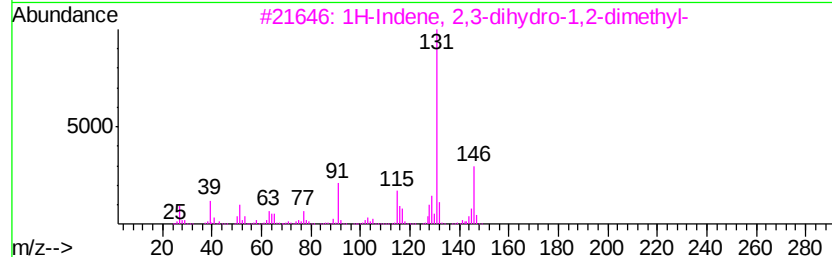
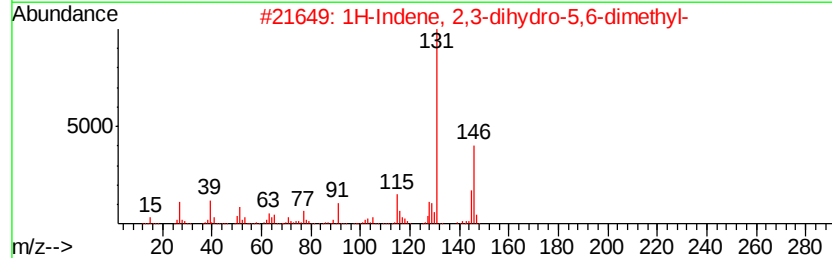
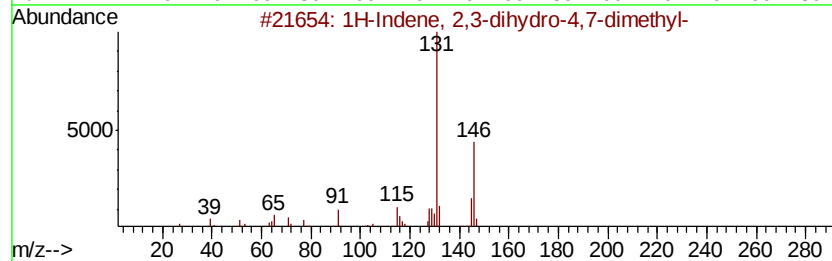
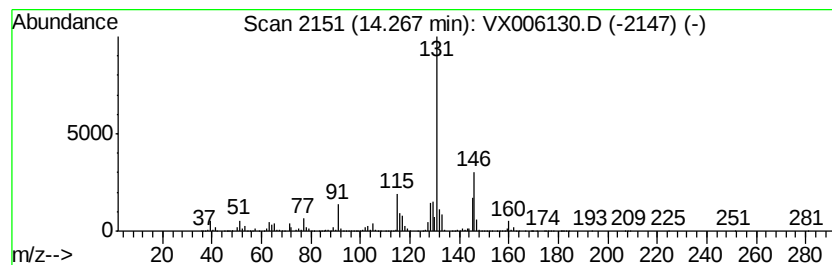
Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	59.74 ug/l	1852480	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	95
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146 C11H14	001075-22-5	94
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	93
4	Benzene, 1-methyl-3-(1-methyl-2-...	146 C11H14	052161-57-6	90
5	Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112118\
Data File : VX006130.D
Acq On : 21 Nov 2018 18:19
Operator : JC/MD
Sample : J6016-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

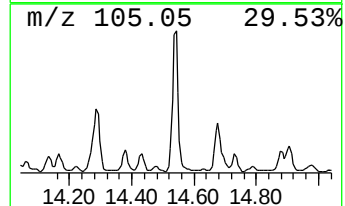
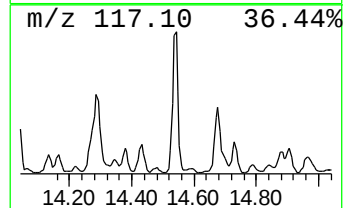
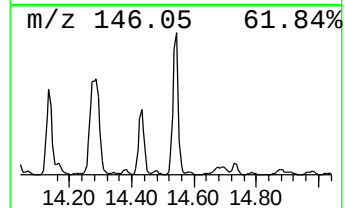
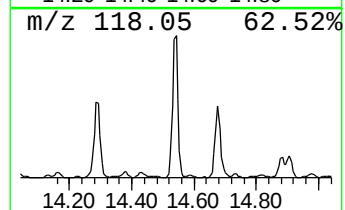
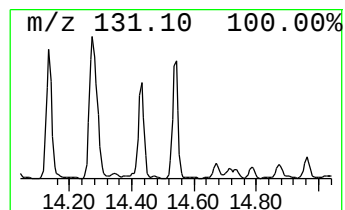
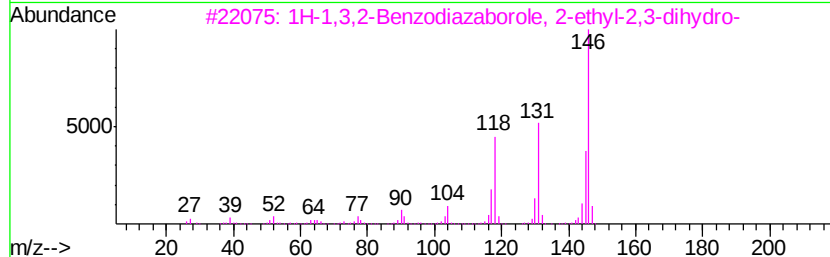
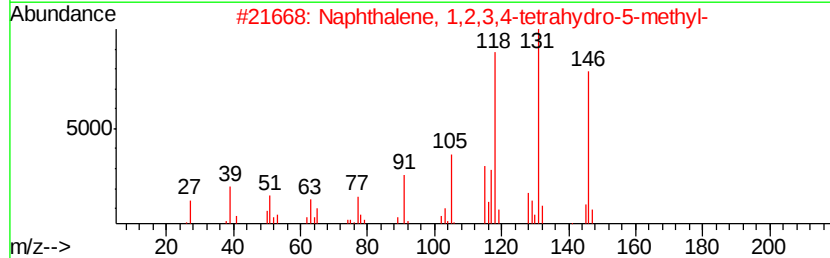
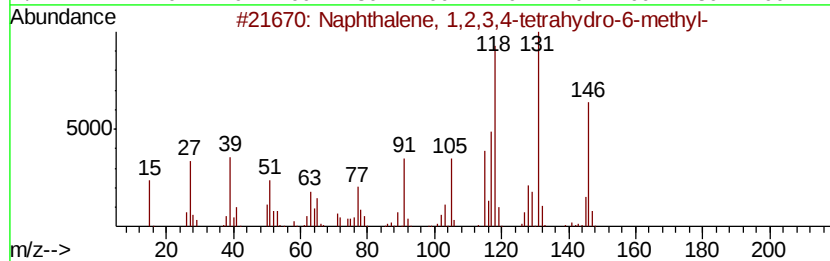
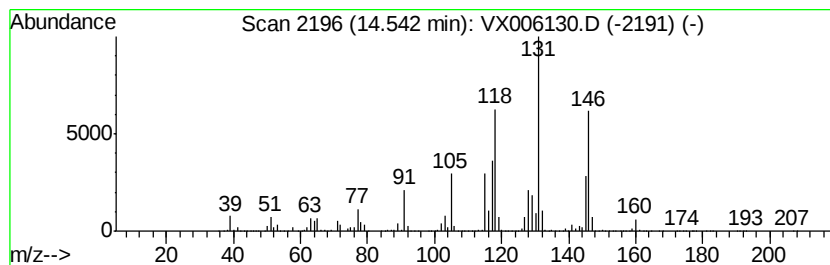
Instrument :
MSVOA_X
ClientSampled :
MW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	53.72 ug/l	1665960	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 95
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 91
4		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 70
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112118\
Data File : VX006130.D
Acq On : 21 Nov 2018 18:19
Operator : JC/MD
Sample : J6016-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

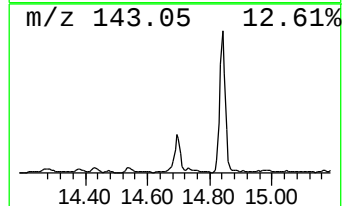
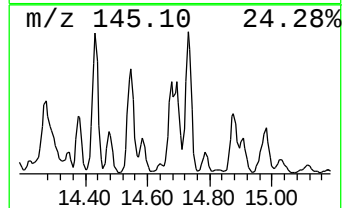
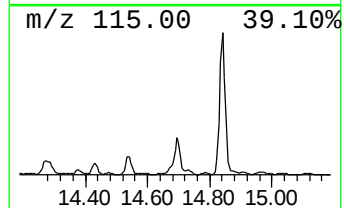
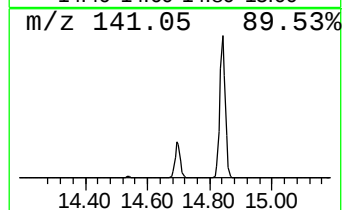
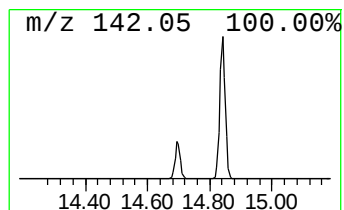
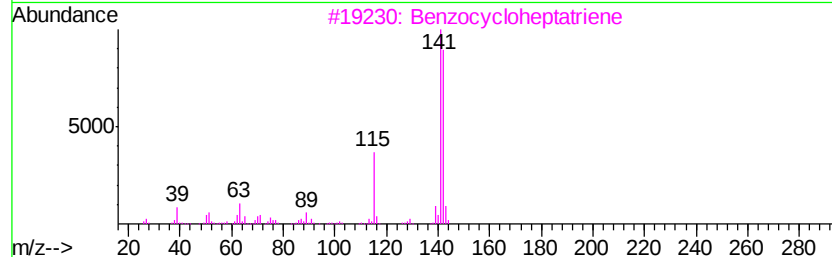
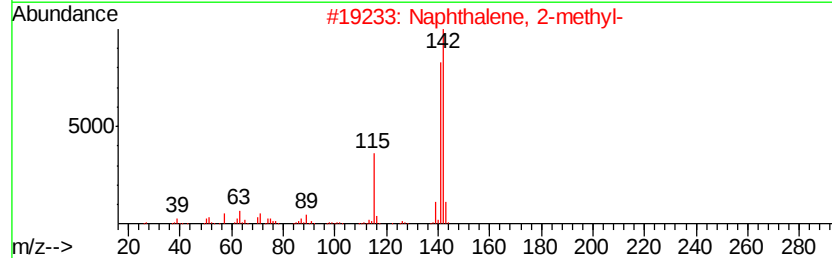
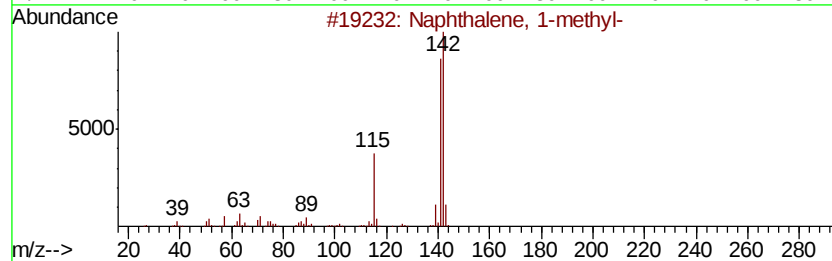
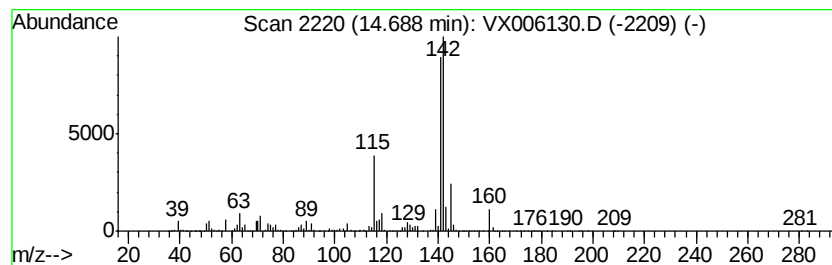
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	79.35 ug/l	2460830	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	90
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112118\
Data File : VX006130.D
Acq On : 21 Nov 2018 18:19
Operator : JC/MD
Sample : J6016-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

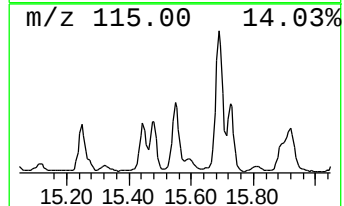
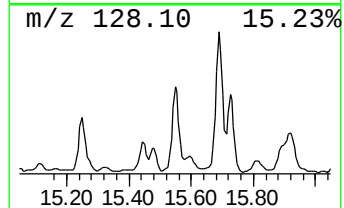
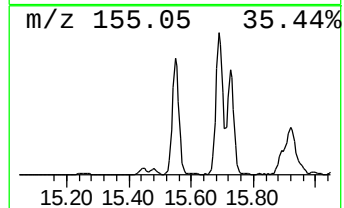
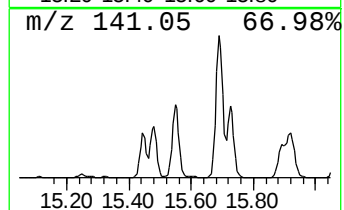
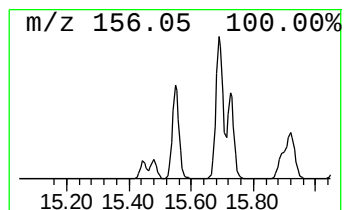
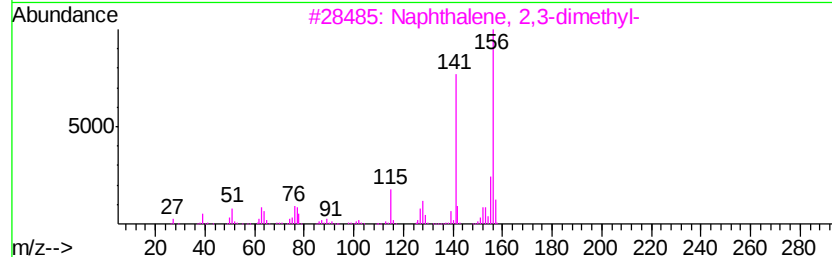
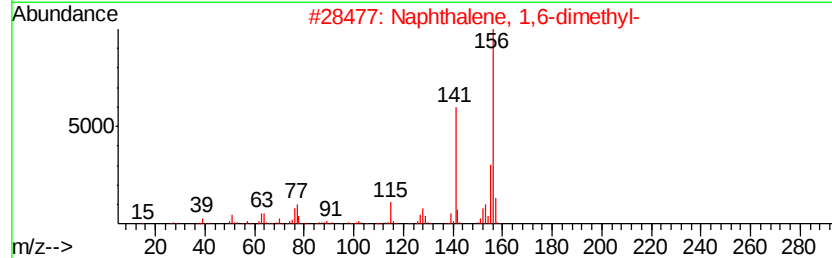
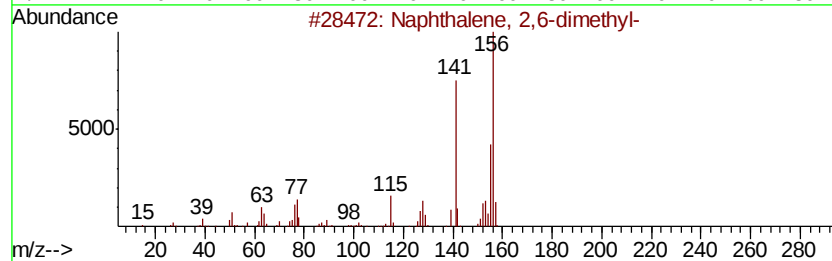
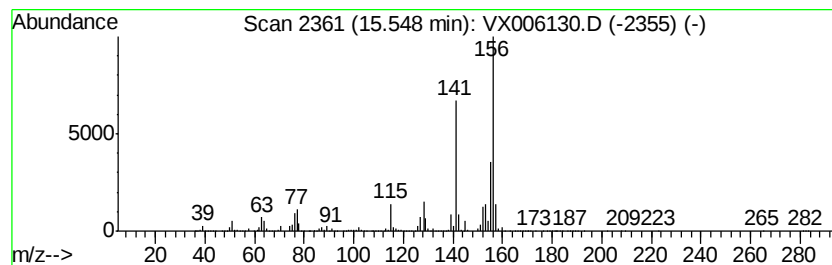
Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	50.88 ug/l	1577820	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 98
2		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
3		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
4		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 96
5		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 96



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112118\
Data File : VX006130.D
Acq On : 21 Nov 2018 18:19
Operator : JC/MD
Sample : J6016-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.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-etheny...	12.30	90.0	ug/l	2790470	4	12.08	1550540	50.0
Benzene, 1-etheny...	12.76	57.4	ug/l	1779050	4	12.08	1550540	50.0
Benzene, 1,2,3,4-...	12.97	50.9	ug/l	1578770	4	12.08	1550540	50.0
Benzene, 2-etheny...	13.35	142.9	ug/l	4432960	4	12.08	1550540	50.0
1H-Indene, 2,3-di...	14.27	59.7	ug/l	1852480	4	12.08	1550540	50.0
Naphthalene, 1,2,...	14.54	53.7	ug/l	1665960	4	12.08	1550540	50.0
Naphthalene, 1-me...	14.69	79.3	ug/l	2460830	4	12.08	1550540	50.0
Naphthalene, 2,6-...	15.55	50.9	ug/l	1577820	4	12.08	1550540	50.0