

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
 Data File : VX002012.D
 Acq On : 26 May 2018 07:41
 Operator : JC/MD
 Sample : J3064-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-18

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.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	1.075	76	80	94	rBV	3153632	3712056	39.53%	4.433%
2	3.081	402	409	418	rVV	128584	278547	2.97%	0.333%
3	3.215	419	431	441	rVB	87183	235851	2.51%	0.282%
4	5.513	795	808	816	rBV	253457	739973	7.88%	0.884%
5	5.672	825	834	853	rVB	283545	796648	8.48%	0.951%
6	6.074	890	900	911	rBV	281627	732025	7.79%	0.874%
7	6.769	1006	1014	1022	rBV2	63861	158730	1.69%	0.190%
8	6.867	1022	1030	1041	rVV	437312	1019327	10.85%	1.217%
9	7.464	1119	1128	1139	rVB	68334	161154	1.72%	0.192%
10	8.659	1318	1324	1328	rBV2	65367	114521	1.22%	0.137%
11	8.720	1328	1334	1341	rVV	1323335	2048219	21.81%	2.446%
12	8.787	1341	1345	1351	rVB	80646	127555	1.36%	0.152%
13	8.885	1356	1361	1374	rVB	148066	248788	2.65%	0.297%
14	9.525	1460	1466	1471	rBV2	77007	122489	1.30%	0.146%
15	9.866	1517	1522	1526	rBV	94536	139742	1.49%	0.167%
16	10.012	1540	1546	1554	rBV	102629	148967	1.59%	0.178%
17	10.116	1554	1563	1568	rBV2	996423	1627430	17.33%	1.943%
18	10.165	1568	1571	1580	rVB4	72588	157336	1.68%	0.188%
19	10.281	1580	1590	1592	rBV2	66301	131933	1.40%	0.158%
20	10.335	1596	1599	1604	rVB	109017	143501	1.53%	0.171%
21	10.610	1640	1644	1648	rBV	87570	132252	1.41%	0.158%
22	10.915	1690	1694	1695	rBV2	110348	138478	1.47%	0.165%
23	10.963	1695	1702	1706	rVB3	347912	761901	8.11%	0.910%
24	11.018	1706	1711	1717	rBV2	562993	916028	9.75%	1.094%
25	11.079	1717	1721	1726	rVB2	88758	149981	1.60%	0.179%
26	11.140	1726	1731	1737	rBV	1189439	1643219	17.50%	1.962%
27	11.232	1742	1746	1752	rBV2	292421	452941	4.82%	0.541%
28	11.287	1752	1755	1760	rVB3	89128	134910	1.44%	0.161%
29	11.366	1764	1768	1772	rBV	623269	734322	7.82%	0.877%
30	11.439	1776	1780	1786	rBV7	68023	123331	1.31%	0.147%
31	11.567	1795	1801	1805	rBV4	462353	767272	8.17%	0.916%
32	11.701	1815	1823	1826	rBV	318339	463233	4.93%	0.553%
33	11.732	1826	1828	1831	rVV2	131631	153558	1.64%	0.183%
34	11.774	1831	1835	1837	rVV	229258	327579	3.49%	0.391%

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
 Data File : VX002012.D
 Acq On : 26 May 2018 07:41
 Operator : JC/MD
 Sample : J3064-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-18

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
 Title : SW846 8260

35	11.799	1837	1839	1846	rVB2	190472	260457	2.77%	0.311%
36	11.951	1861	1864	1866	rVV	229675	277496	2.95%	0.331%
37	11.982	1866	1869	1872	rVB3	119472	183640	1.96%	0.219%
38	12.085	1881	1886	1891	rBV	1194534	1523844	16.23%	1.820%
39	12.140	1891	1895	1898	rVB3	154982	198759	2.12%	0.237%
40	12.213	1903	1907	1910	rVB2	196124	271833	2.89%	0.325%
41	12.305	1918	1922	1926	rBV	2031778	2627083	27.97%	3.137%
42	12.347	1926	1929	1932	rVB2	491254	602184	6.41%	0.719%
43	12.469	1942	1949	1953	rBV3	451474	723514	7.70%	0.864%
44	12.536	1956	1960	1964	rVB2	163057	261000	2.78%	0.312%
45	12.603	1964	1971	1974	rBV5	225233	576821	6.14%	0.689%
46	12.640	1974	1977	1979	rVV3	214125	339247	3.61%	0.405%
47	12.677	1979	1983	1986	rVV	1169879	1540721	16.41%	1.840%
48	12.707	1986	1988	1991	rVV	295536	382856	4.08%	0.457%
49	12.762	1993	1997	2001	rVV2	1090142	1900560	20.24%	2.270%
50	12.805	2001	2004	2010	rVB2	1072092	1691686	18.01%	2.020%
51	12.884	2010	2017	2018	rBV5	193751	309989	3.30%	0.370%
52	12.902	2018	2020	2024	rVB3	197313	231788	2.47%	0.277%
53	12.981	2025	2033	2037	rBV	1201633	1693413	18.03%	2.022%
54	13.042	2040	2043	2046	rVB2	206476	256006	2.73%	0.306%
55	13.079	2046	2049	2056	rVB6	338891	523522	5.57%	0.625%
56	13.152	2056	2061	2065	rBV3	350559	552152	5.88%	0.659%
57	13.207	2066	2070	2073	rBV2	644689	917680	9.77%	1.096%
58	13.256	2076	2078	2083	rVV3	155200	206134	2.19%	0.246%
59	13.323	2085	2089	2090	rVV2	429921	540626	5.76%	0.646%
60	13.353	2090	2094	2098	rVV2	3535022	5157676	54.92%	6.159%
61	13.402	2098	2102	2106	rVV	7868106	9391475	100.00%	11.215%
62	13.481	2110	2115	2118	rVV2	719003	1244060	13.25%	1.486%
63	13.536	2122	2124	2132	rVB3	570075	1086512	11.57%	1.297%
64	13.609	2132	2136	2139	rBV	1335255	1577289	16.79%	1.883%
65	13.646	2139	2142	2146	rVV3	500883	698997	7.44%	0.835%
66	13.707	2147	2152	2153	rVV3	463133	685298	7.30%	0.818%
67	13.725	2153	2155	2161	rVB5	735394	1242130	13.23%	1.483%
68	13.817	2167	2170	2172	rBV	192431	257532	2.74%	0.308%
69	13.914	2182	2186	2189	rBV2	535386	823992	8.77%	0.984%
70	13.945	2189	2191	2193	rVV2	395001	506127	5.39%	0.604%
71	13.987	2195	2198	2202	rVV2	708952	1128745	12.02%	1.348%

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
 Data File : VX002012.D
 Acq On : 26 May 2018 07:41
 Operator : JC/MD
 Sample : J3064-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-18

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
 Title : SW846 8260

72	14.036	2202	2206	2214	rVV6	708521	1506430	16.04%	1.799%
73	14.103	2214	2217	2219	rVV	208663	235257	2.51%	0.281%
74	14.134	2220	2222	2224	rVV2	207794	272195	2.90%	0.325%
75	14.164	2224	2227	2231	rVV2	342964	538032	5.73%	0.642%
76	14.225	2231	2237	2242	rVV6	350597	893576	9.51%	1.067%
77	14.274	2242	2245	2251	rVV3	718387	1859635	19.80%	2.221%
78	14.323	2251	2253	2258	rVV3	689204	983599	10.47%	1.175%
79	14.384	2258	2263	2267	rVV4	523838	939816	10.01%	1.122%
80	14.438	2267	2272	2276	rVB3	625055	855913	9.11%	1.022%
81	14.548	2286	2290	2293	rBV2	766799	1051142	11.19%	1.255%
82	14.579	2293	2295	2299	rVB3	409088	438250	4.67%	0.523%
83	14.640	2302	2305	2308	rVV4	175890	262369	2.79%	0.313%
84	14.707	2308	2316	2324	rVB2	1655969	3395540	36.16%	4.055%
85	14.847	2334	2339	2343	rBV	1828894	2473335	26.34%	2.953%
86	14.963	2353	2358	2366	rBV7	323635	783698	8.34%	0.936%
87	15.036	2366	2370	2377	rVB6	297046	637371	6.79%	0.761%
88	15.127	2381	2385	2389	rBV	361913	510859	5.44%	0.610%
89	15.255	2400	2406	2410	rBV5	226912	476271	5.07%	0.569%
90	15.402	2425	2430	2434	rBV	298281	417088	4.44%	0.498%
91	15.450	2434	2438	2441	rVV2	153618	226515	2.41%	0.270%
92	15.487	2441	2444	2451	rVB4	143596	264101	2.81%	0.315%
93	15.554	2451	2455	2460	rBV	337594	508666	5.42%	0.607%
94	15.694	2472	2478	2482	rBV	419057	733377	7.81%	0.876%
95	15.731	2482	2484	2490	rVB	299757	408573	4.35%	0.488%
96	15.816	2492	2498	2504	rBV4	108180	220790	2.35%	0.264%
97	15.932	2513	2517	2523	rVB3	118792	236550	2.52%	0.282%
98	16.084	2537	2542	2553	rVB2	173831	373286	3.97%	0.446%
99	16.182	2554	2558	2568	rVB9	45727	104066	1.11%	0.124%

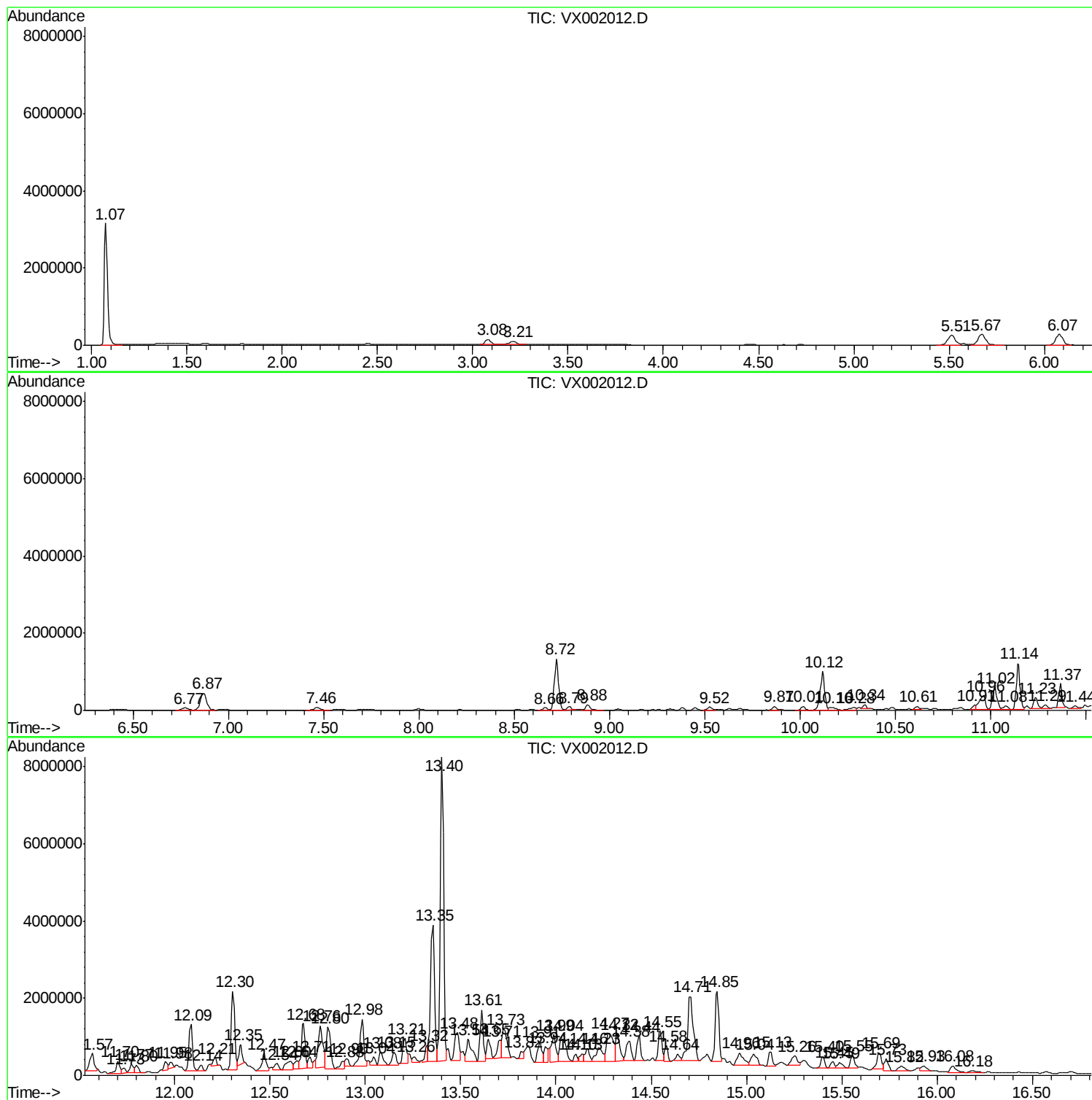
Sum of corrected areas: 83742941

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

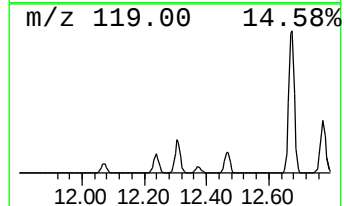
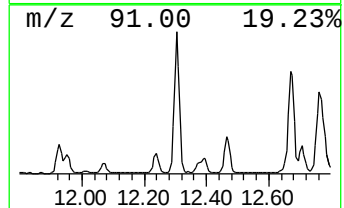
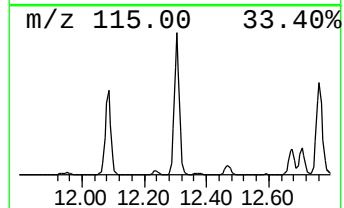
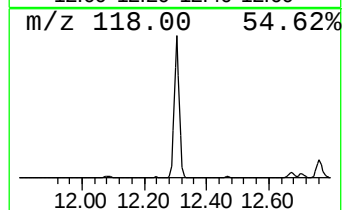
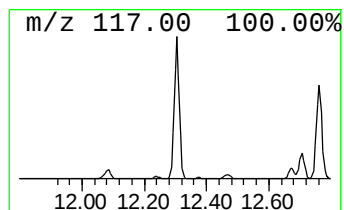
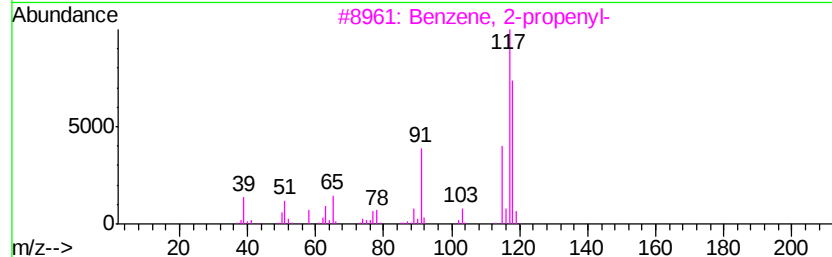
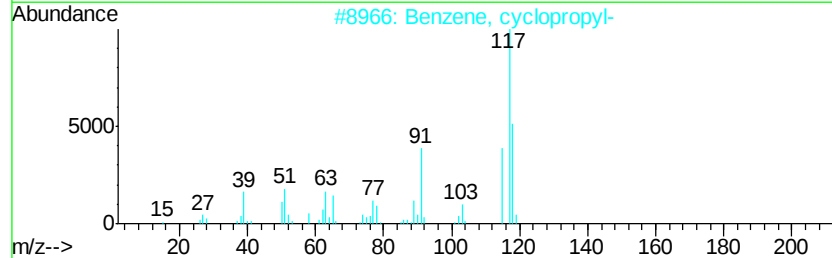
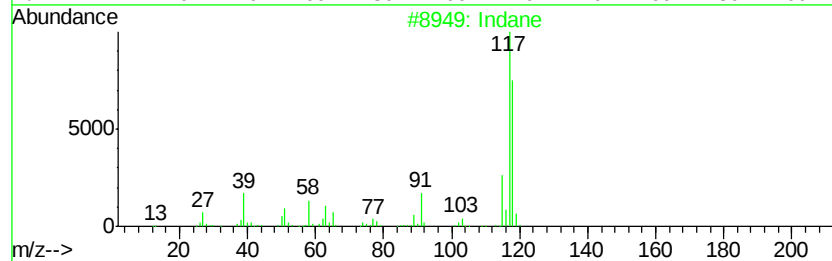
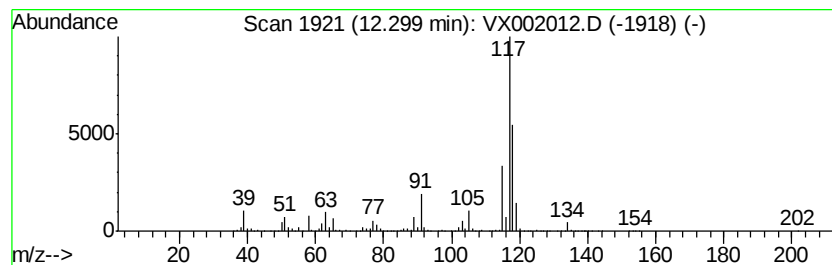
Instrument :
MSVOA_X
ClientSampled :
MW-18

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.30	86.20 ug/l	2627080	1,4-Dichlorobenzene-d4	12.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	50
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	50



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

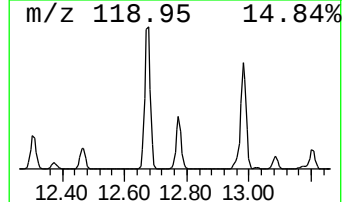
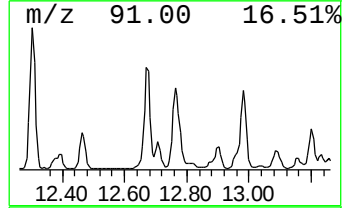
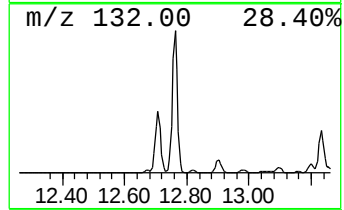
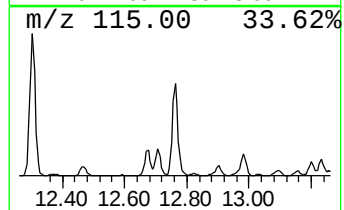
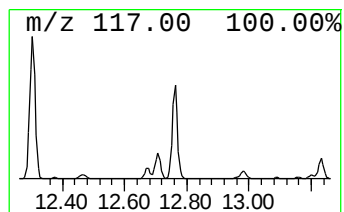
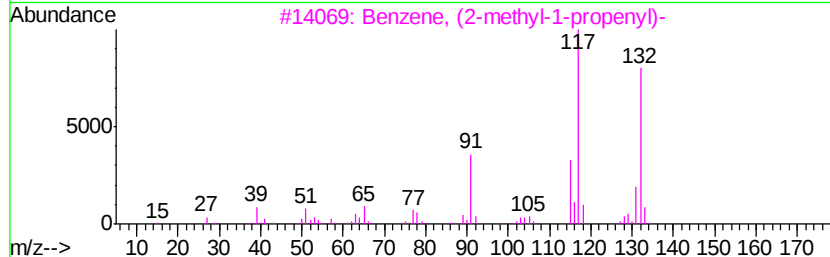
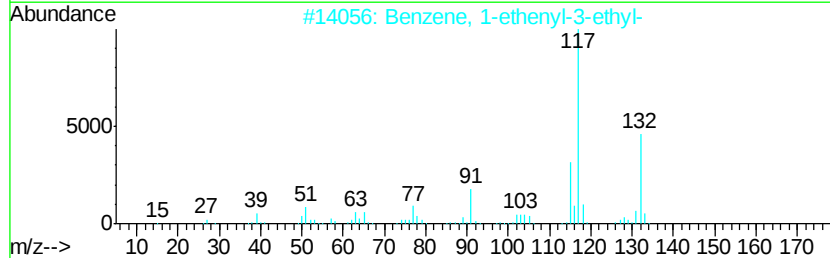
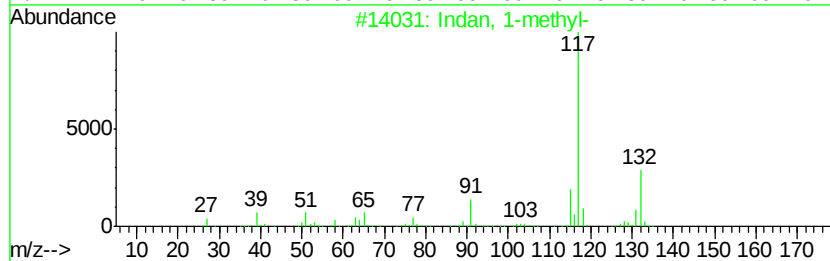
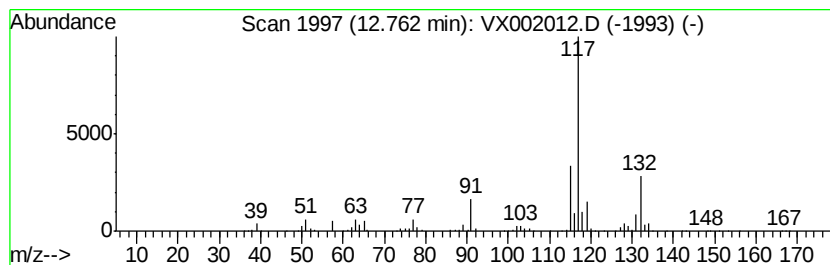
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 3 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	62.36 ug/l	1900560	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

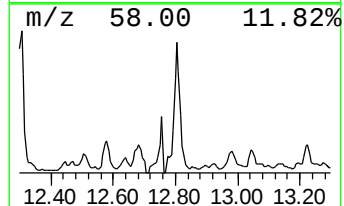
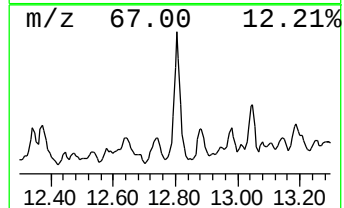
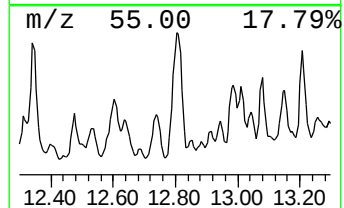
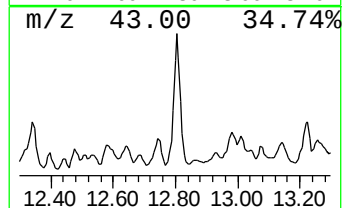
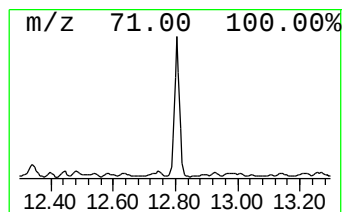
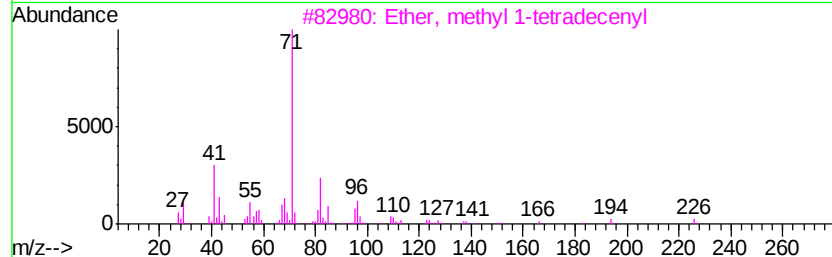
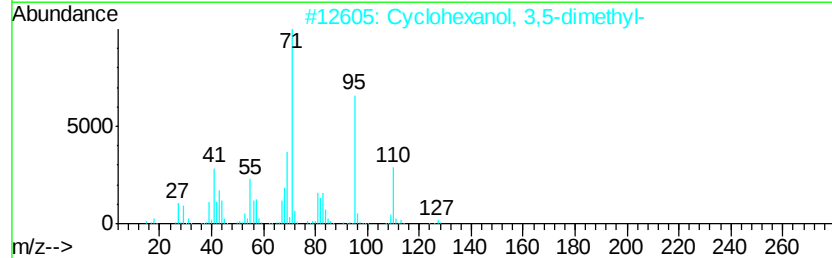
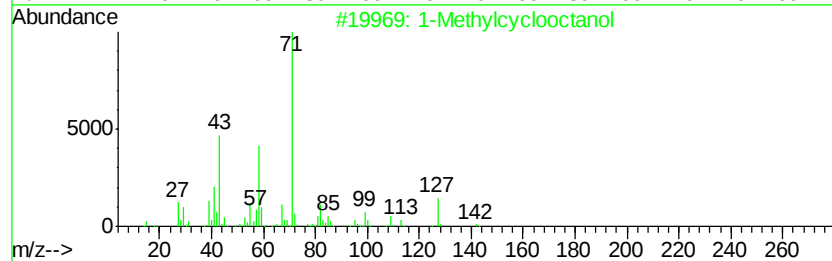
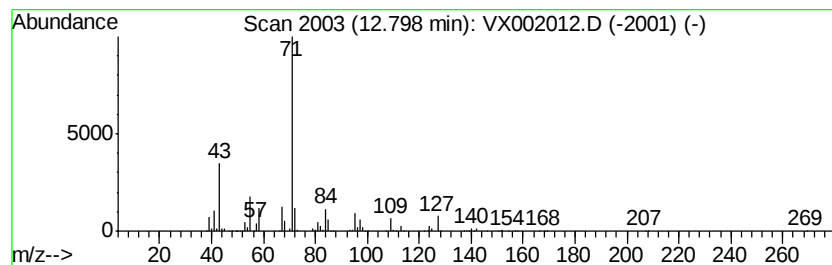
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 4 1-Methylcyclooctanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.80	55.51 ug/l	1691690	1,4-Dichlorobenzene-d4	12.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Methylcyclooctanol	142	C9H18O	059123-41-0	59
2		Cyclohexanol, 3,5-dimethyl-	128	C8H16O	005441-52-1	59
3		Ether, methyl 1-tetradecenyl	226	C15H30O	026537-05-3	50
4		(2E)-Dodec-2-en-1-yl methyl ether	198	C13H26O	1000334-06-3	50
5		2-Furanmethanol, tetrahydro-, ac...	144	C7H12O3	000637-64-9	47



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-18

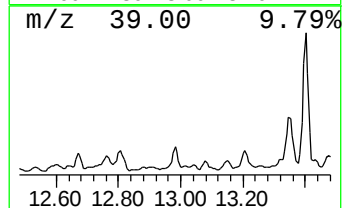
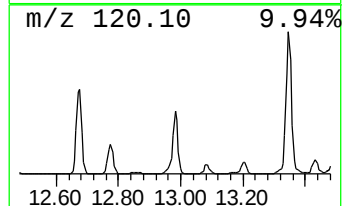
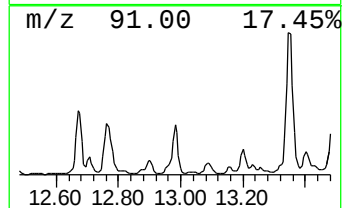
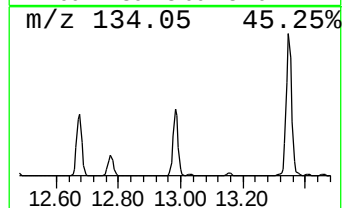
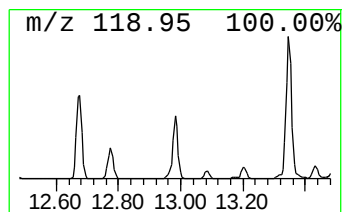
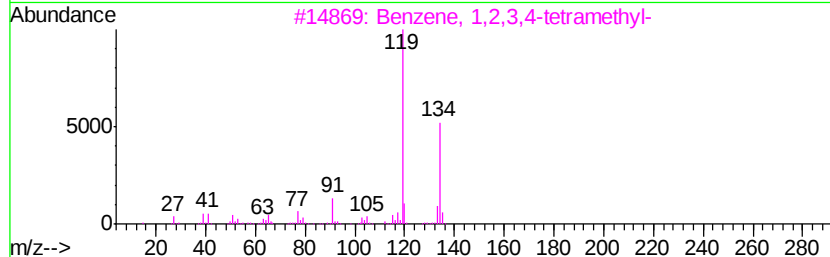
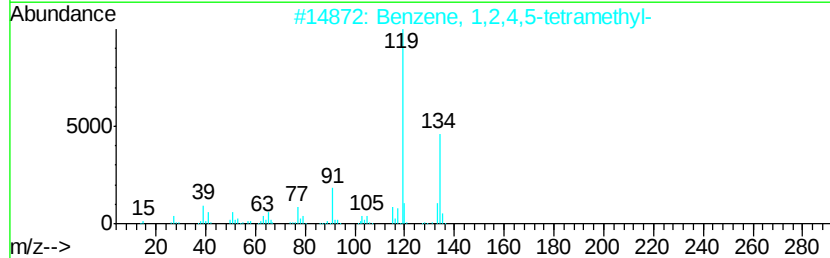
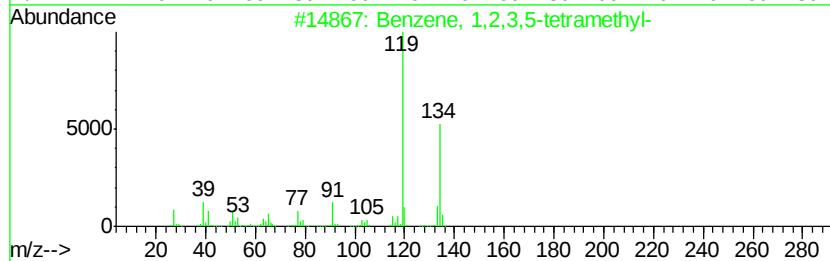
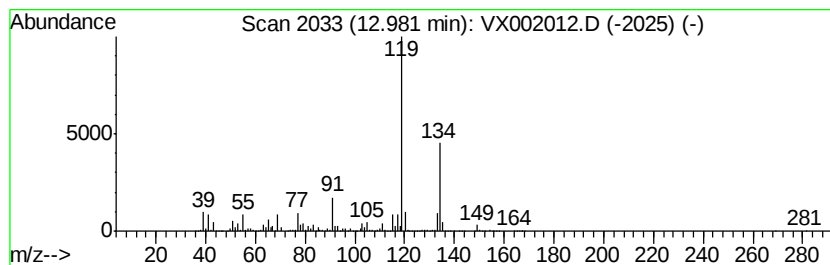
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	55.56 ug/l	1693410	1,4-Dichlorobenzene-d4	12.09

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-18

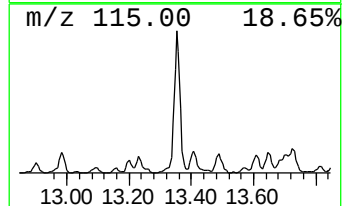
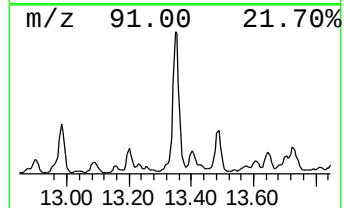
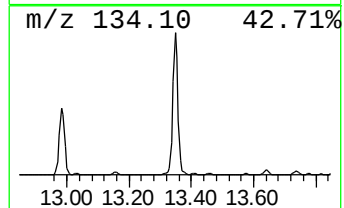
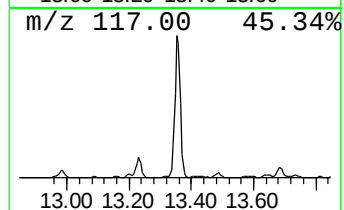
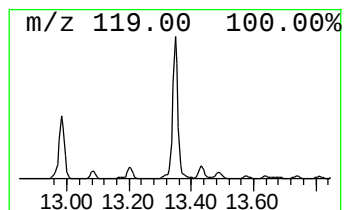
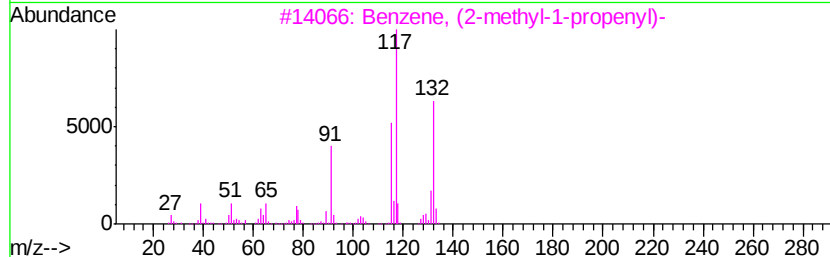
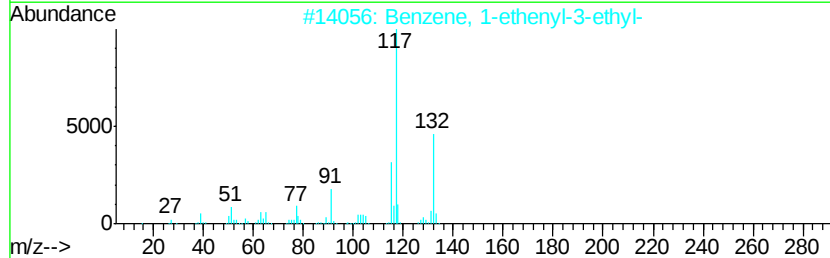
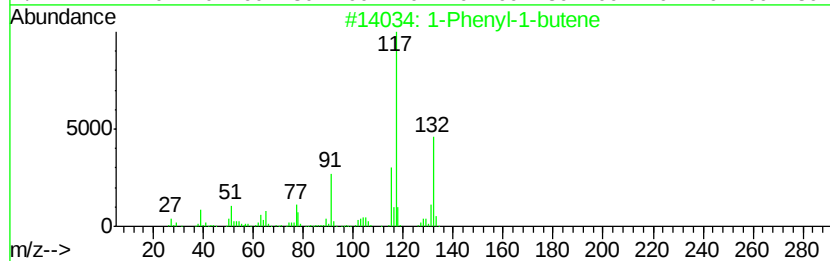
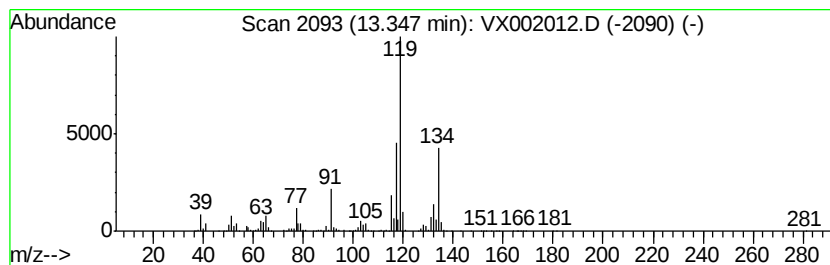
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 6 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	169.23 ug/l	5157680	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	70
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	64



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

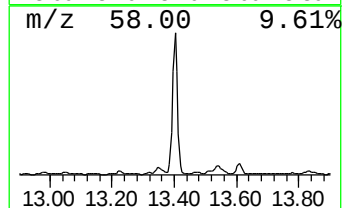
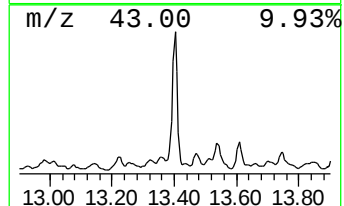
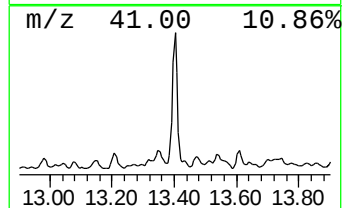
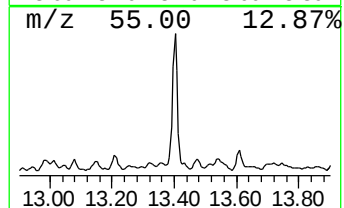
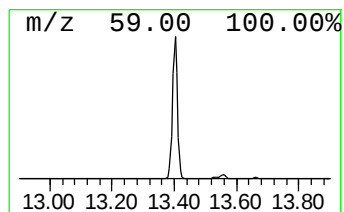
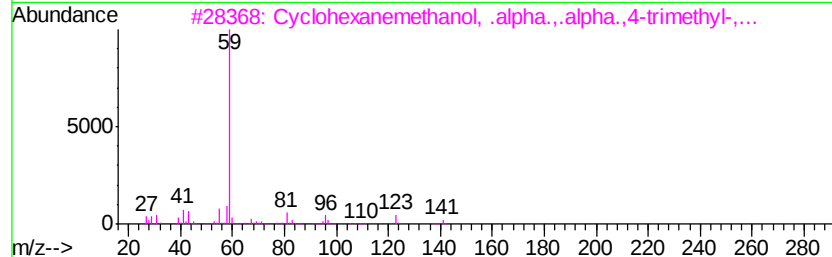
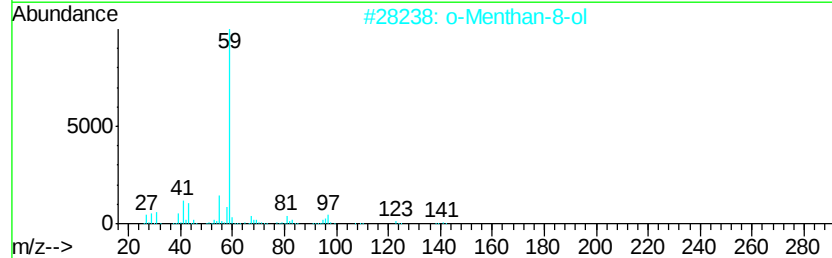
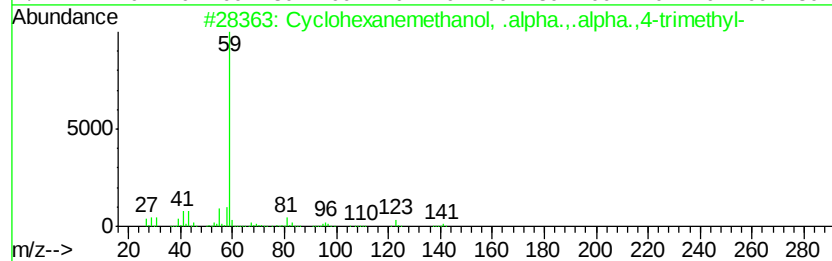
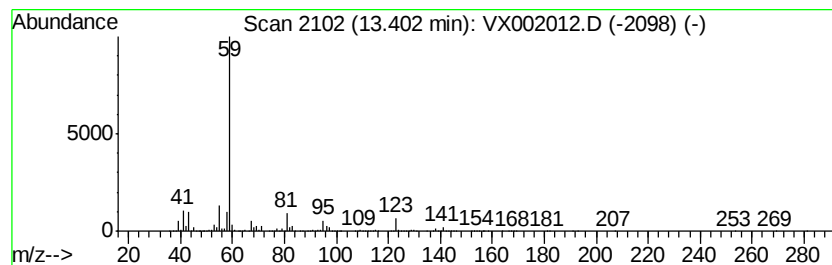
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclohexanemethanol, .alpha... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.40	308.15 ug/l	9391480	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	83
2		o-Menthan-8-ol	156	C10H20O	343855-44-7	78
3		Cyclohexanemethanol, .alpha...al...	156	C10H20O	007322-63-6	78
4		1-Cyclohexylethanol, methyl ether	142	C9H18O	1000365-12-8	72
5		Cyclohexanemethanol, .alpha...al...	156	C10H20O	005114-00-1	64



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

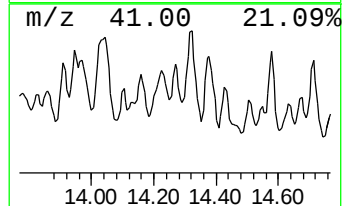
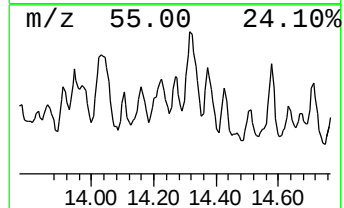
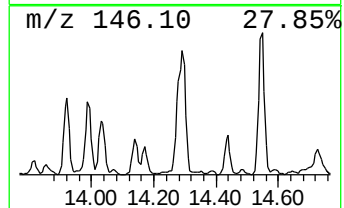
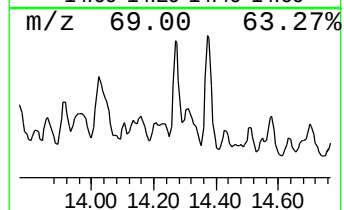
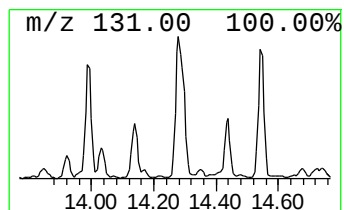
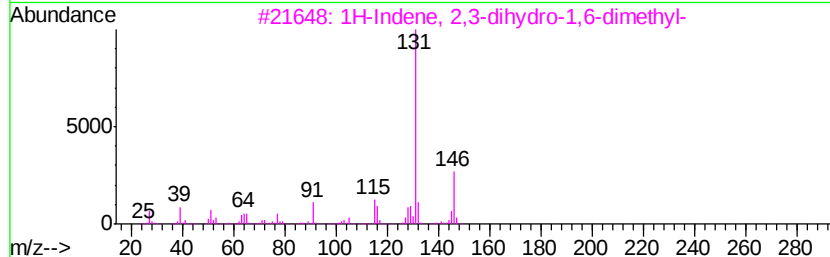
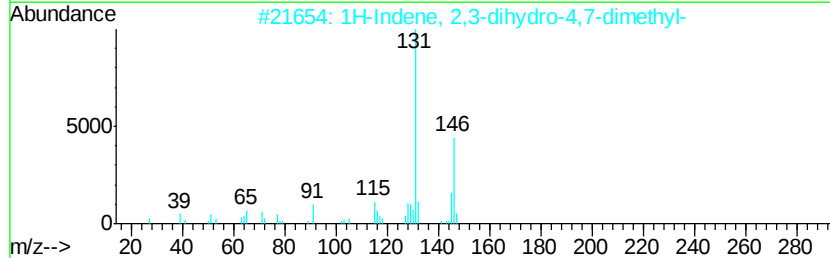
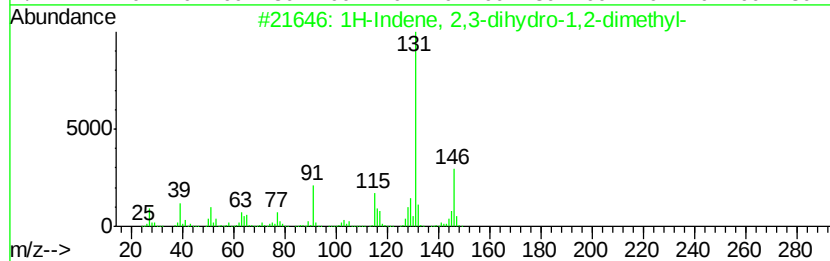
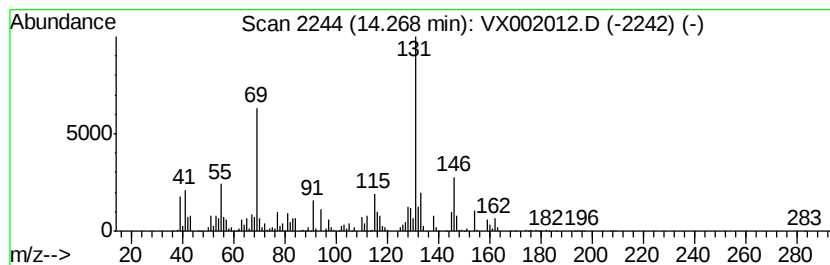
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	61.02 ug/l	1859640	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87
4		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	87
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

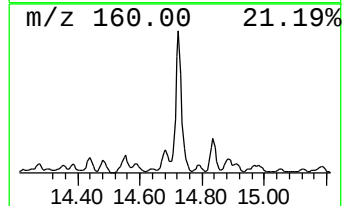
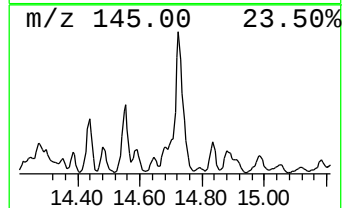
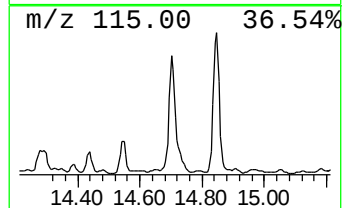
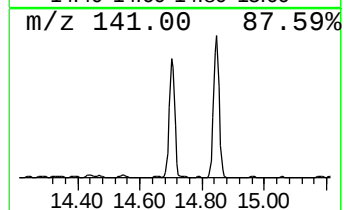
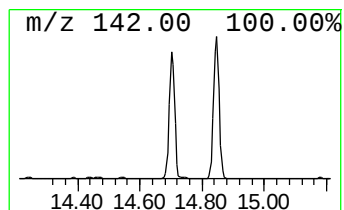
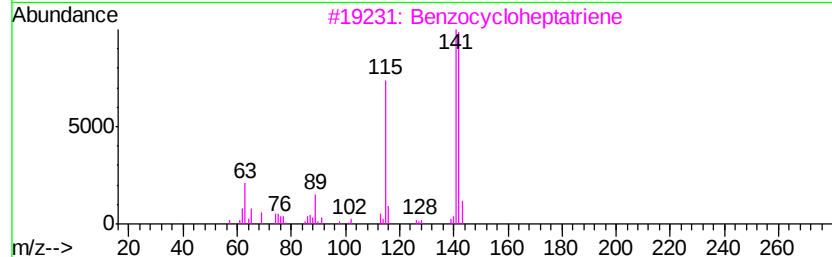
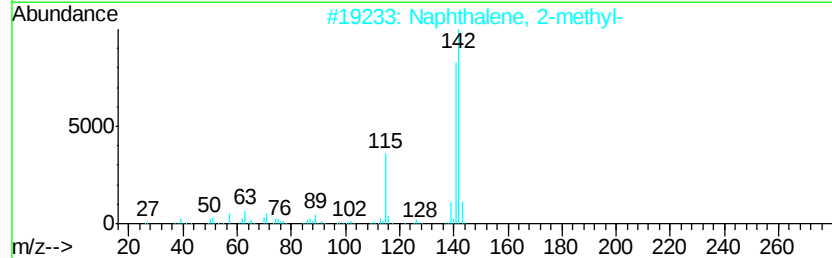
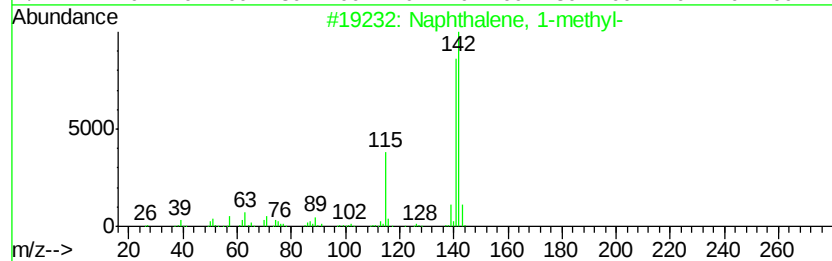
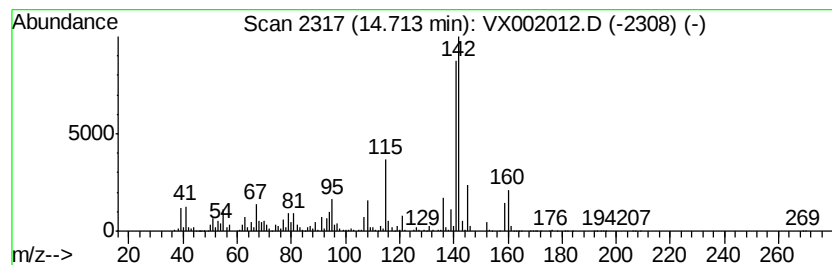
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	111.41 ug/l	3395540	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	53



Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX052518\
Data File : VX002012.D
Acq On : 26 May 2018 07:41
Operator : JC/MD
Sample : J3064-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-18

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.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
Indane	12.30	86.2	ug/l	2627080	4	12.09	1523840	50.0
Indan, 1-methyl-	12.76	62.4	ug/l	1900560	4	12.09	1523840	50.0
1-Methylcyclooctanol	12.80	55.5	ug/l	1691690	4	12.09	1523840	50.0
Benzene, 1,2,3,5-...	12.98	55.6	ug/l	1693410	4	12.09	1523840	50.0
1-Phenyl-1-butene	13.35	169.2	ug/l	5157680	4	12.09	1523840	50.0
Cyclohexanemethan...	13.40	308.1	ug/l	9391480	4	12.09	1523840	50.0
1H-Indene, 2,3-di...	14.27	61.0	ug/l	1859640	4	12.09	1523840	50.0
Naphthalene, 1-me...	14.71	111.4	ug/l	3395540	4	12.09	1523840	50.0