

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
 Data File : VX011254.D
 Acq On : 31 Jul 2019 19:37
 Operator : JC/SP
 Sample : K4048-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OWL-C1-GW

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\82X071619W.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	2.440	206	211	222	rBV	19433	36686	1.52%	0.137%
2	5.500	702	713	724	rBV	117789	330264	13.71%	1.238%
3	5.665	730	740	755	rVB	225427	631424	26.20%	2.366%
4	6.061	793	805	820	rBV	122504	321391	13.34%	1.204%
5	6.860	925	936	950	rBV	312319	735285	30.51%	2.756%
6	7.463	1025	1035	1044	rBV2	22336	56912	2.36%	0.213%
7	8.713	1234	1240	1250	rVB	657514	994135	41.25%	3.726%
8	9.036	1286	1293	1300	rVB3	17950	34542	1.43%	0.129%
9	9.621	1385	1389	1394	rVB	18051	25344	1.05%	0.095%
10	10.109	1463	1469	1476	rBV	635036	846334	35.12%	3.172%
11	10.640	1547	1556	1563	rBV4	14515	40709	1.69%	0.153%
12	10.835	1575	1588	1594	rBV3	24807	49195	2.04%	0.184%
13	10.908	1594	1600	1612	rBV5	17897	37077	1.54%	0.139%
14	11.018	1612	1618	1622	rBV	191949	244587	10.15%	0.917%
15	11.133	1632	1637	1642	rBV	537742	674248	27.98%	2.527%
16	11.274	1656	1660	1665	rVB4	20070	28974	1.20%	0.109%
17	11.359	1668	1674	1683	rBV	237290	307322	12.75%	1.152%
18	11.670	1719	1725	1733	rVB4	25368	44641	1.85%	0.167%
19	11.914	1758	1765	1767	rBV2	25487	40124	1.67%	0.150%
20	11.944	1767	1770	1776	rVB	92716	114023	4.73%	0.427%
21	12.072	1786	1791	1797	rBV	658449	845380	35.08%	3.168%
22	12.231	1813	1817	1822	rVB	51267	68512	2.84%	0.257%
23	12.298	1822	1828	1833	rBV	679308	827682	34.35%	3.102%
24	12.365	1835	1839	1840	rVV	50204	58789	2.44%	0.220%
25	12.383	1840	1842	1847	rVB	61609	72428	3.01%	0.271%
26	12.456	1849	1854	1861	rBV	119186	156805	6.51%	0.588%
27	12.664	1880	1888	1891	rBV	641708	757845	31.45%	2.840%
28	12.700	1891	1894	1898	rVV	251936	308098	12.79%	1.155%
29	12.755	1898	1903	1910	rVV2	795897	1183528	49.11%	4.435%
30	12.810	1910	1912	1918	rVB2	26700	38206	1.59%	0.143%
31	12.895	1918	1926	1930	rVB	80404	130584	5.42%	0.489%
32	12.975	1930	1939	1944	rBV	566758	708048	29.38%	2.653%
33	13.029	1944	1948	1952	rVV2	20358	32121	1.33%	0.120%
34	13.078	1952	1956	1963	rVB3	104756	158539	6.58%	0.594%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
 Data File : VX011254.D
 Acq On : 31 Jul 2019 19:37
 Operator : JC/SP
 Sample : K4048-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OWL-C1-GW

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\82X071619W.M
 Title : SW846 8260

35	13.145	1963	1967	1971	rBV	66631	90175	3.74%	0.338%
36	13.194	1971	1975	1977	rVV2	137844	198645	8.24%	0.744%
37	13.225	1977	1980	1989	rVV	508413	650356	26.99%	2.437%
38	13.310	1989	1994	1995	rVV2	70244	80969	3.36%	0.303%
39	13.346	1995	2000	2006	rVV2	1713879	2409783	100.00%	9.031%
40	13.401	2006	2009	2010	rVV	62658	72609	3.01%	0.272%
41	13.420	2010	2012	2016	rVV	87102	121107	5.03%	0.454%
42	13.474	2018	2021	2029	rVV	62990	97407	4.04%	0.365%
43	13.560	2030	2035	2038	rVV	106216	147243	6.11%	0.552%
44	13.596	2038	2041	2044	rVV	282453	371744	15.43%	1.393%
45	13.639	2044	2048	2051	rVV2	339239	489394	20.31%	1.834%
46	13.694	2051	2057	2058	rVV2	213723	328377	13.63%	1.231%
47	13.718	2058	2061	2070	rVV2	396196	617109	25.61%	2.313%
48	13.804	2070	2075	2079	rVV	64677	91608	3.80%	0.343%
49	13.852	2079	2083	2088	rVV2	100930	147627	6.13%	0.553%
50	13.907	2088	2092	2097	rVV	371142	454612	18.87%	1.704%
51	13.980	2097	2104	2108	rVV2	468496	635112	26.36%	2.380%
52	14.023	2108	2111	2115	rVV2	164006	224923	9.33%	0.843%
53	14.060	2115	2117	2121	rVV	40889	45742	1.90%	0.171%
54	14.127	2123	2128	2132	rVV	256569	339438	14.09%	1.272%
55	14.157	2132	2133	2140	rVV3	40710	69195	2.87%	0.259%
56	14.218	2140	2143	2146	rVV3	28556	42757	1.77%	0.160%
57	14.285	2146	2154	2161	rVV3	551487	1116392	46.33%	4.184%
58	14.377	2165	2169	2173	rVV	191812	252778	10.49%	0.947%
59	14.432	2173	2178	2182	rVV2	326751	497515	20.65%	1.864%
60	14.474	2182	2185	2190	rVB	59524	81317	3.37%	0.305%
61	14.535	2190	2195	2200	rBV2	606037	798787	33.15%	2.994%
62	14.578	2200	2202	2208	rVV3	34997	58753	2.44%	0.220%
63	14.669	2208	2217	2223	rVV3	224615	457760	19.00%	1.715%
64	14.730	2223	2227	2231	rVV	152465	210245	8.72%	0.788%
65	14.779	2231	2235	2240	rVV3	61381	95647	3.97%	0.358%
66	14.846	2241	2246	2248	rVV3	76262	128908	5.35%	0.483%
67	14.871	2248	2250	2253	rVV	100643	145541	6.04%	0.545%
68	14.901	2253	2255	2259	rVV2	84291	104412	4.33%	0.391%
69	14.962	2259	2265	2272	rVV4	76707	171252	7.11%	0.642%
70	15.035	2272	2277	2283	rVB6	19486	45409	1.88%	0.170%
71	15.108	2283	2289	2294	rBV	38399	62547	2.60%	0.234%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
 Data File : VX011254.D
 Acq On : 31 Jul 2019 19:37
 Operator : JC/SP
 Sample : K4048-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OWL-C1-GW

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
 Title : SW846 8260

72	15.242	2306	2311	2319	rVB2	186934	333274	13.83%	1.249%
73	15.322	2319	2324	2331	rVB4	34511	67404	2.80%	0.253%
74	15.438	2338	2343	2346	rVV	144792	219070	9.09%	0.821%
75	15.474	2346	2349	2354	rVV	101947	146547	6.08%	0.549%
76	15.547	2354	2361	2365	rVV	457296	749016	31.08%	2.807%
77	15.590	2365	2368	2374	rVV5	66013	129559	5.38%	0.486%
78	15.681	2374	2383	2388	rVV	523957	853116	35.40%	3.197%
79	15.718	2388	2389	2397	rVB3	47360	62651	2.60%	0.235%
80	15.803	2398	2403	2410	rBV4	26506	64691	2.68%	0.242%
81	15.913	2411	2421	2436	rVV	185255	522035	21.66%	1.956%
82	16.065	2440	2446	2457	rVB	111231	203991	8.47%	0.764%
83	16.273	2477	2480	2486	rVB3	17890	30819	1.28%	0.115%
84	16.425	2494	2505	2515	rVV3	40504	163447	6.78%	0.613%
85	16.504	2515	2518	2523	rVB2	14711	24985	1.04%	0.094%
86	16.559	2523	2527	2532	rBV2	16429	28149	1.17%	0.105%
87	16.651	2532	2542	2543	rBV7	16438	41313	1.71%	0.155%
88	16.681	2543	2547	2553	rVV	44590	85045	3.53%	0.319%
89	16.742	2553	2557	2564	rVB2	35821	65840	2.73%	0.247%

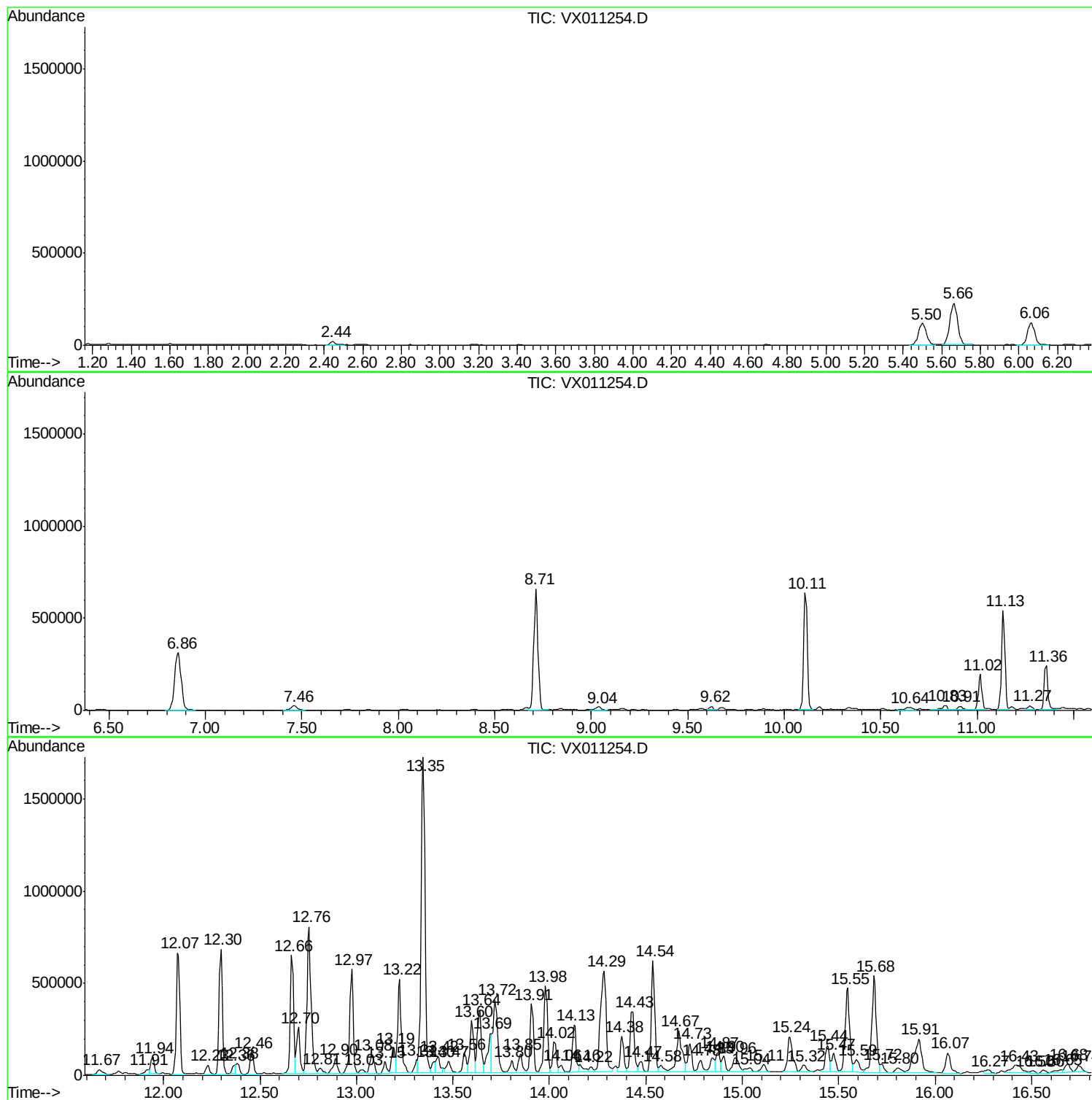
Sum of corrected areas: 26683929

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

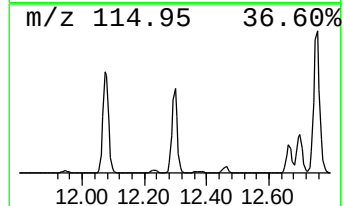
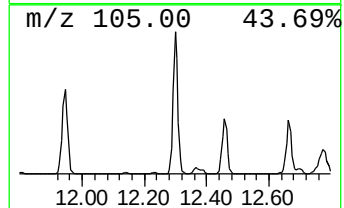
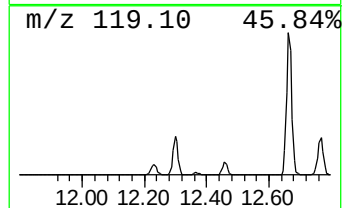
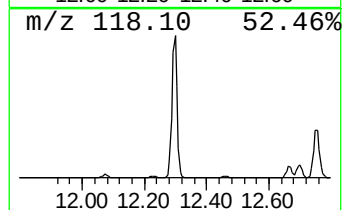
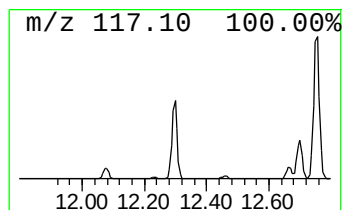
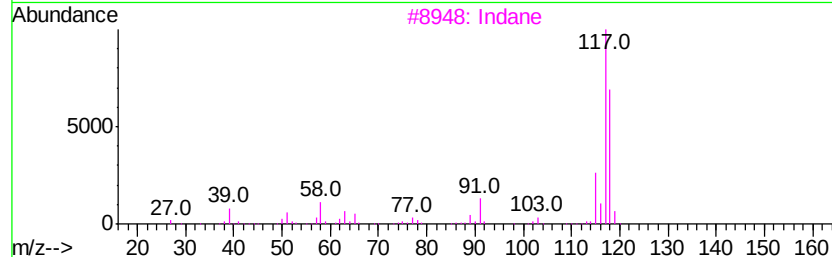
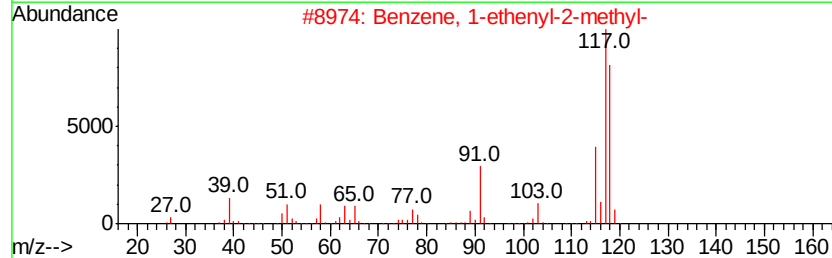
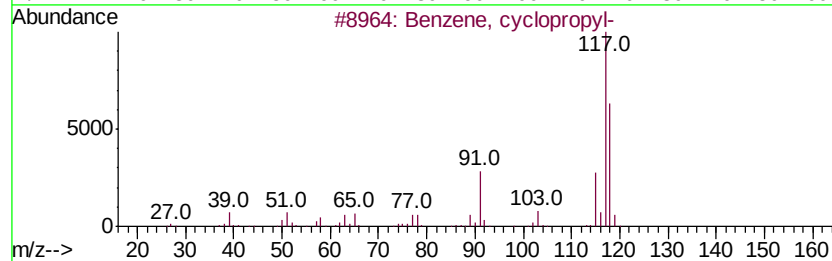
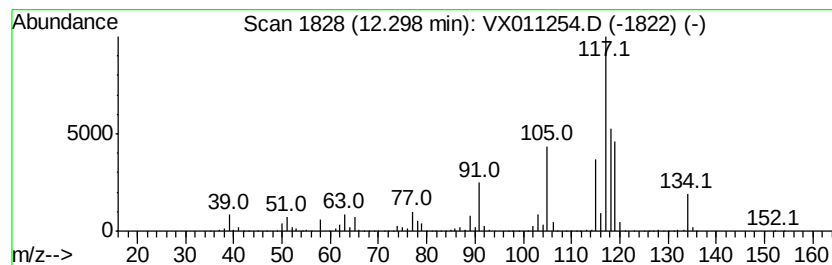
Instrument :
MSVOA_X
ClientSampled :
OWL-C1-GW

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, cyclopropyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.30	48.95 ug/l	827682	1,4-Dichlorobenzene-d4		12.08	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	cvclopropyl-	118	C9H10	000873-49-4	87
2	Benzene,	1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
3	Indane		118	C9H10	000496-11-7	55
4	Benzene,	2-propenyl-	118	C9H10	000300-57-2	55
5	Benzene,	1-propenyl-	118	C9H10	000637-50-3	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

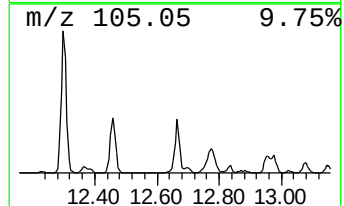
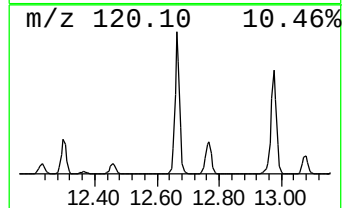
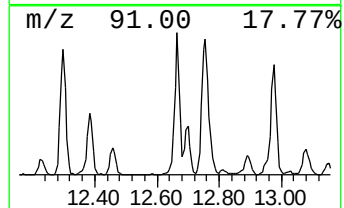
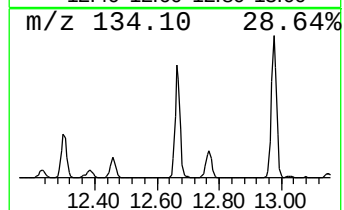
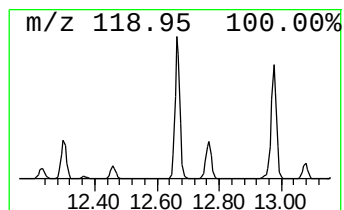
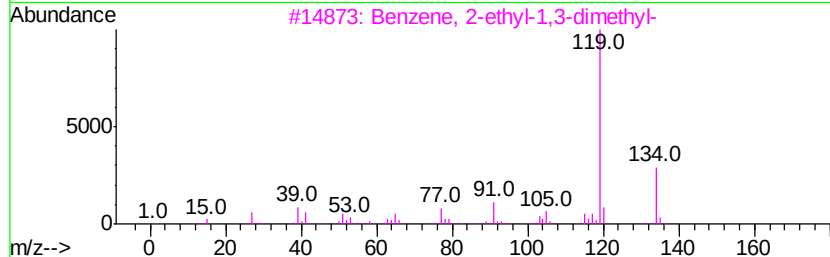
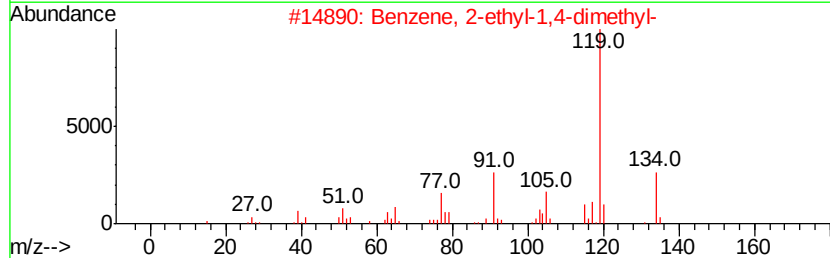
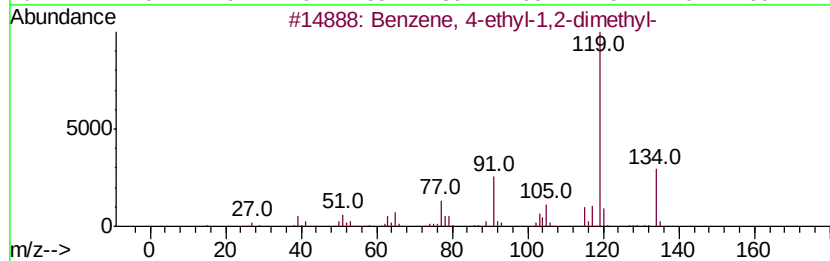
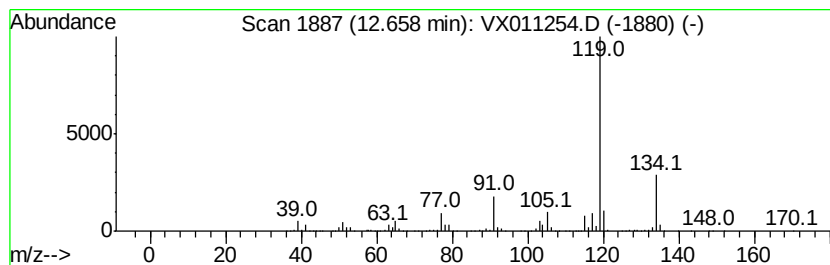
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	44.82 ug/l	757845	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

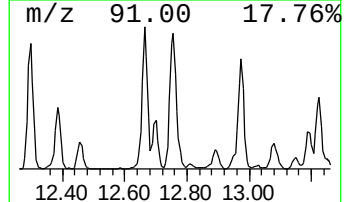
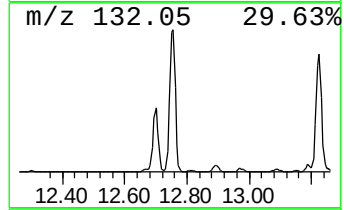
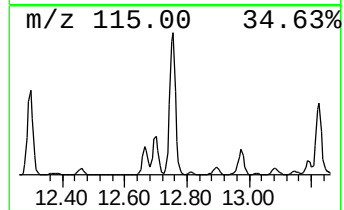
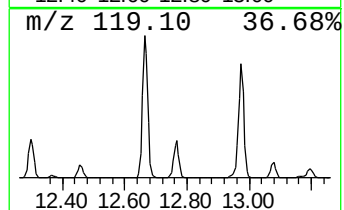
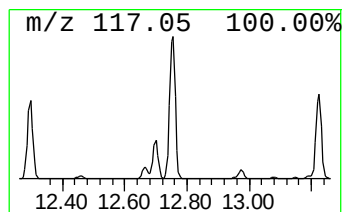
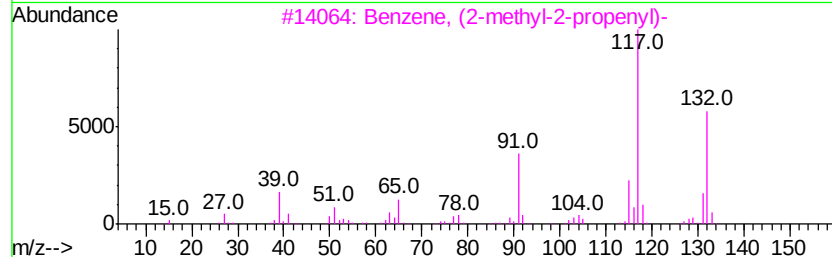
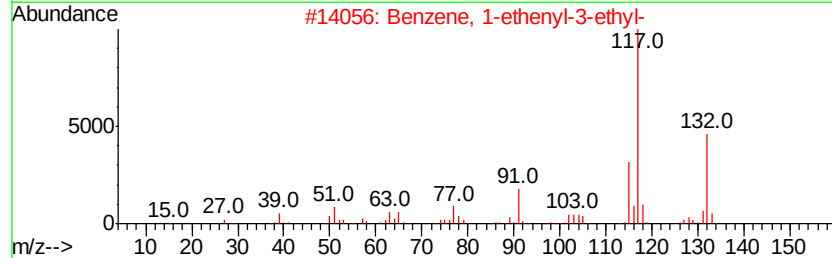
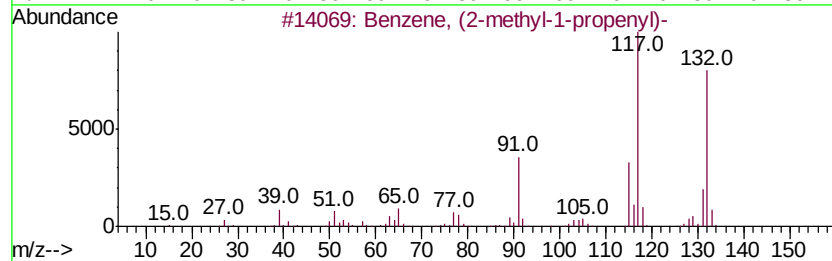
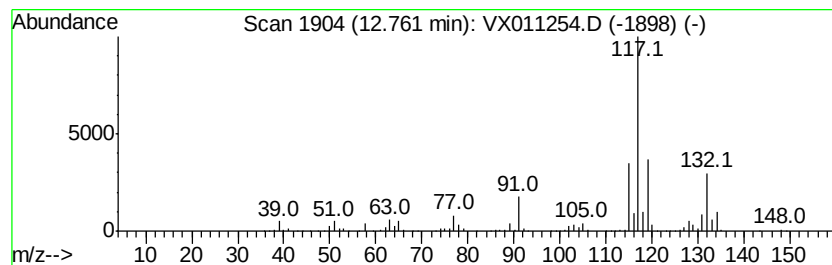
Instrument :
MSVOA_X
ClientSampled :
OWL-C1-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, (2-methyl-1-propen... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	70.00 ug/l	1183530	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 91
2		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 90
3		Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7 87
4		Benzene, 2-butenyl-	132 C10H12	001560-06-1 87
5		Indan, 1-methyl-	132 C10H12	000767-58-8 87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

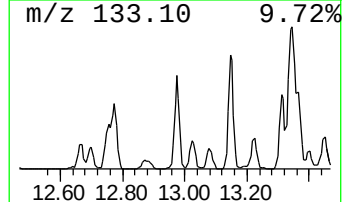
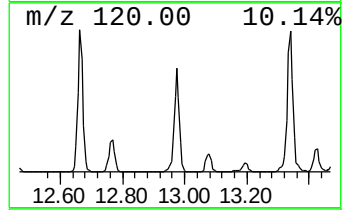
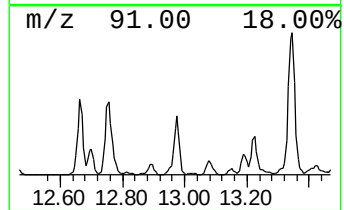
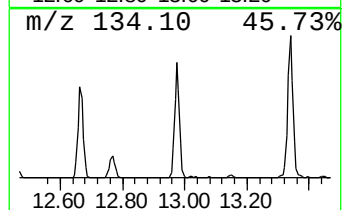
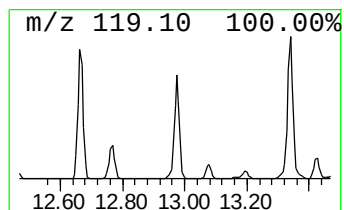
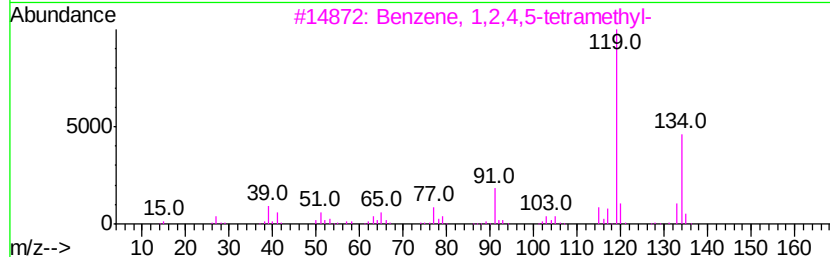
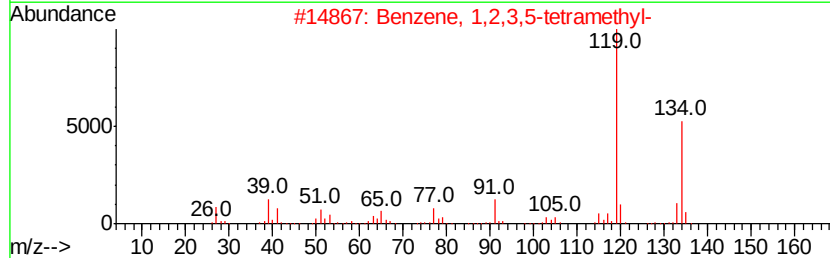
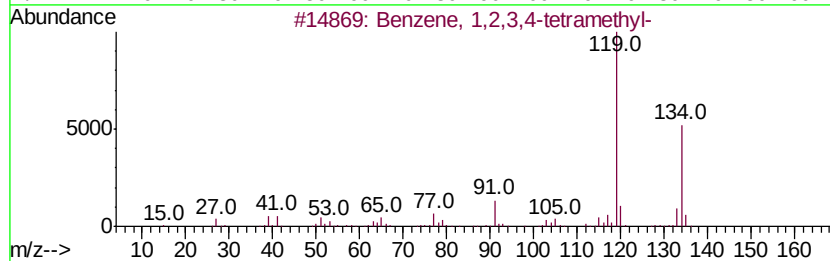
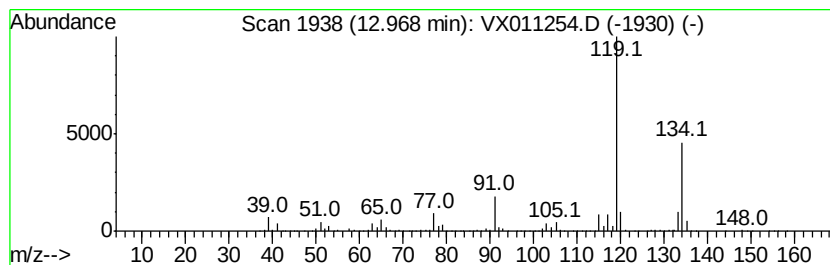
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	41.88 ug/l	708048	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

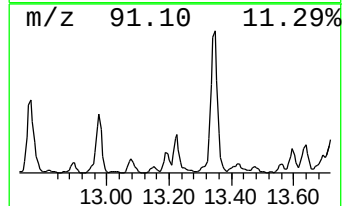
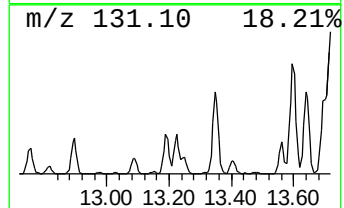
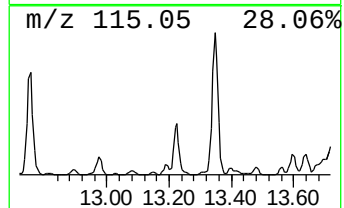
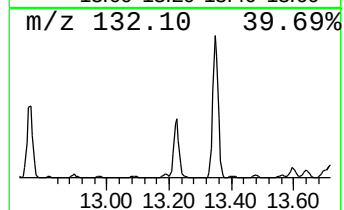
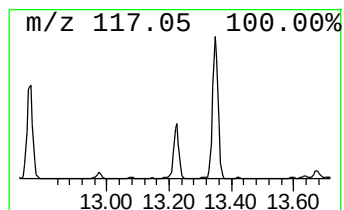
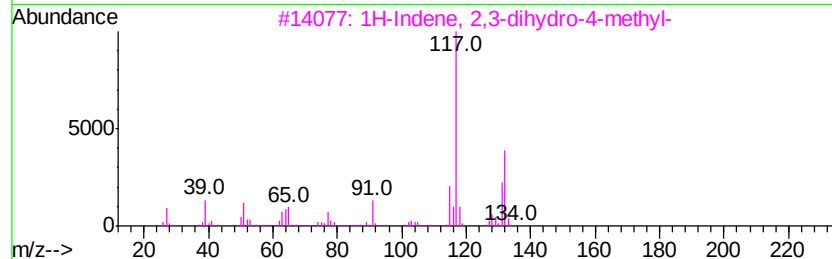
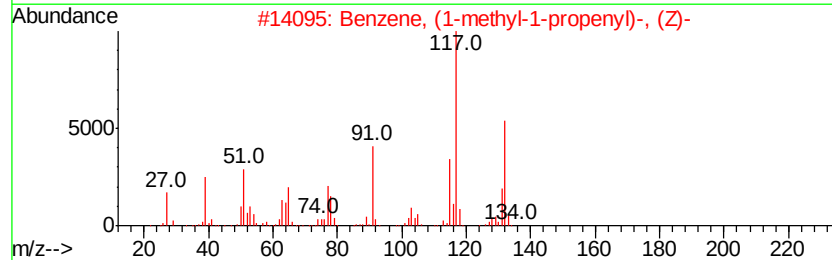
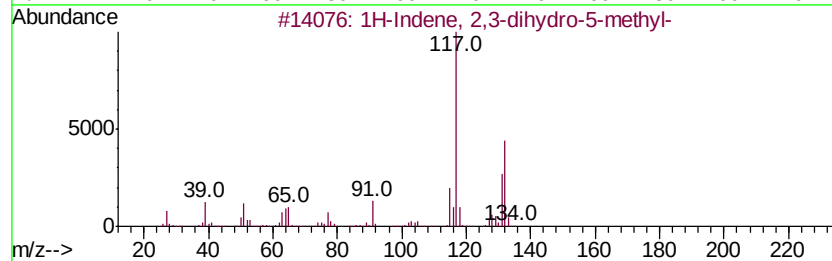
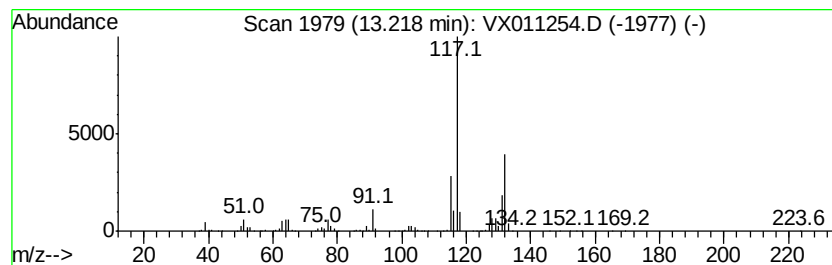
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	38.47 ug/l	650356	1,4-Dichlorobenzene-d4	12.08

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
5	Indan, 1-methyl-	132	C10H12	000767-58-8	83



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

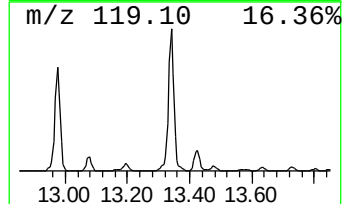
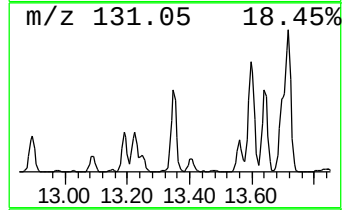
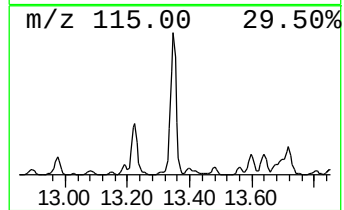
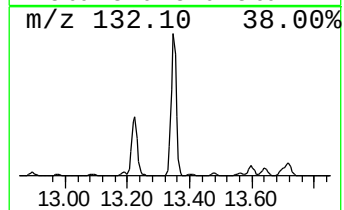
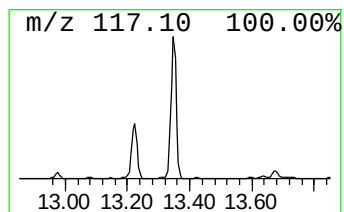
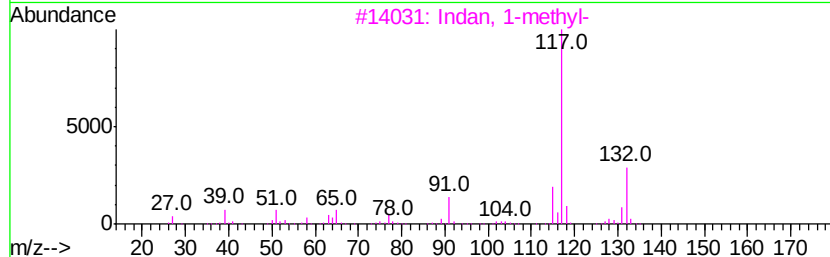
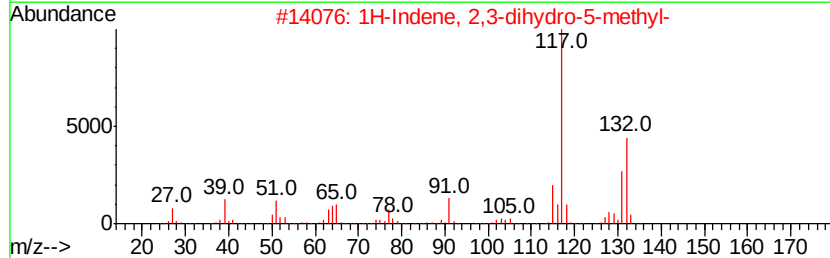
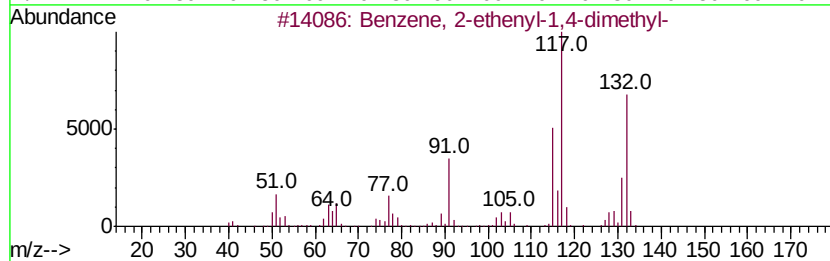
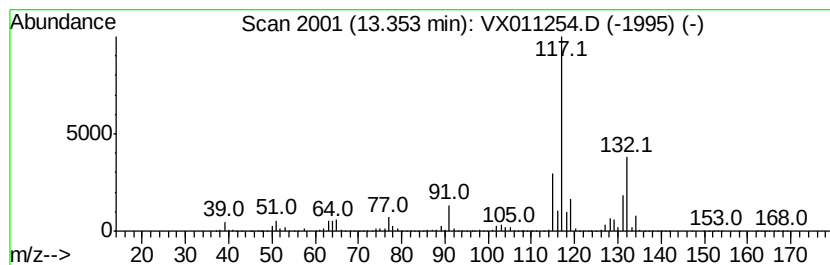
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	142.53 ug/l	2409780	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	87
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	86
3		Indan, 1-methyl-	132	C10H12	000767-58-8	76
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

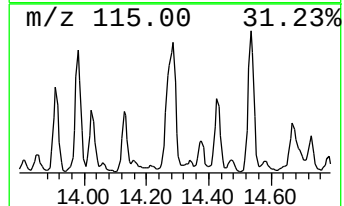
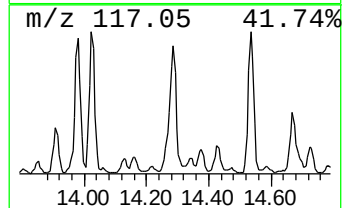
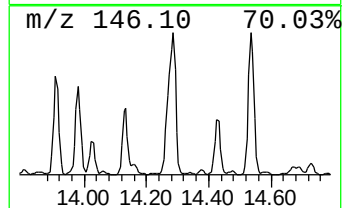
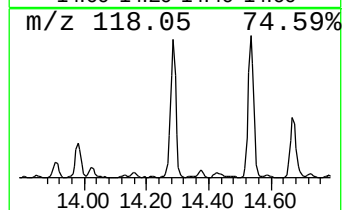
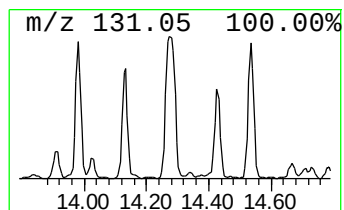
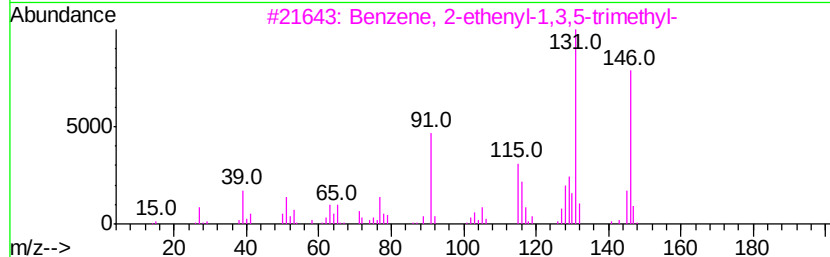
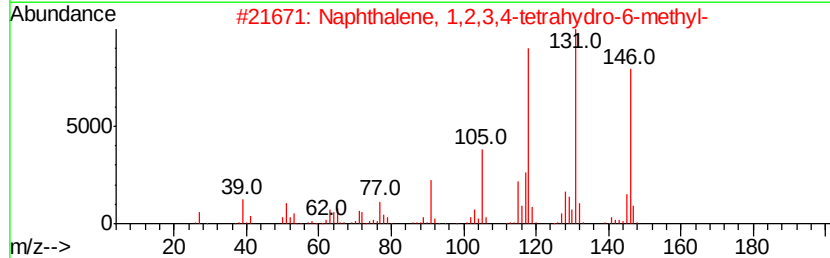
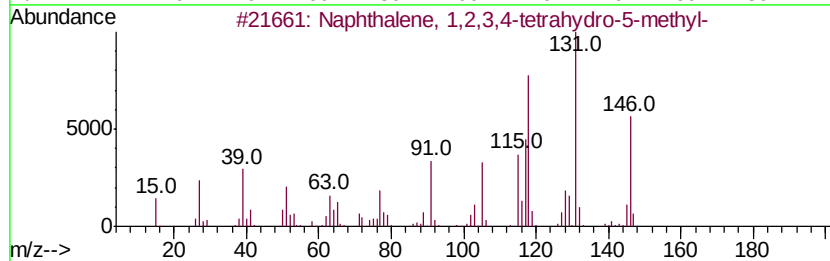
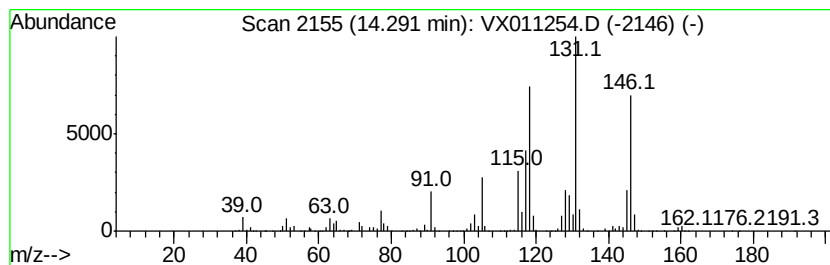
Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	66.03 ug/l	1116390	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 97
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 97
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 89
4		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 70
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
OWL-C1-GW

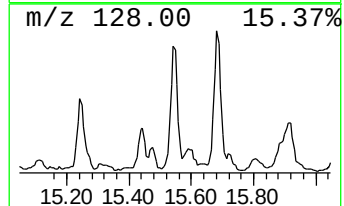
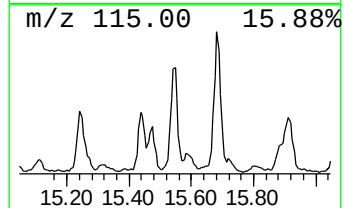
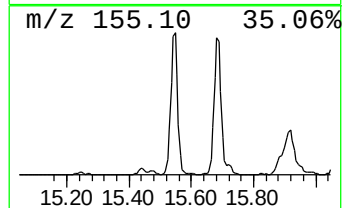
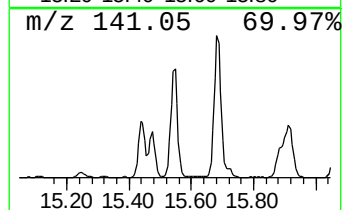
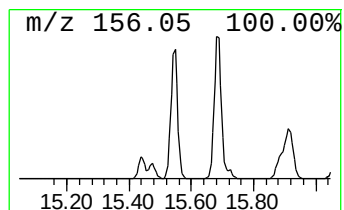
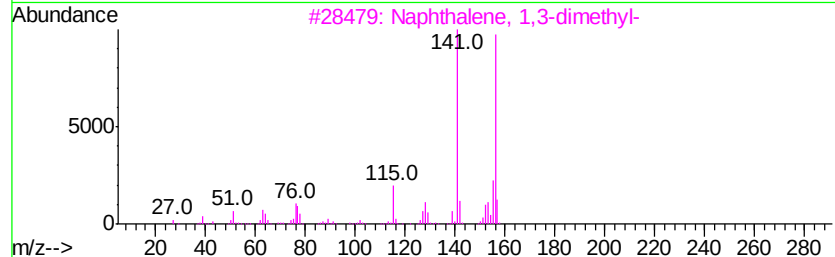
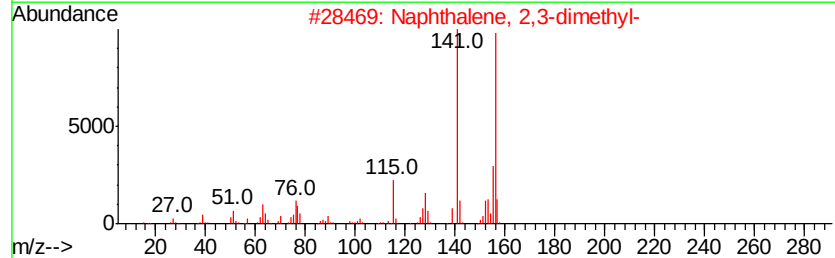
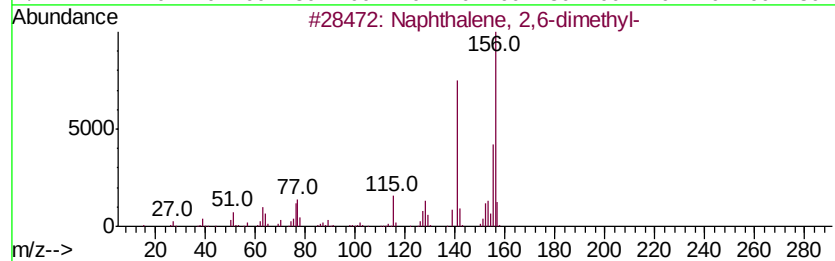
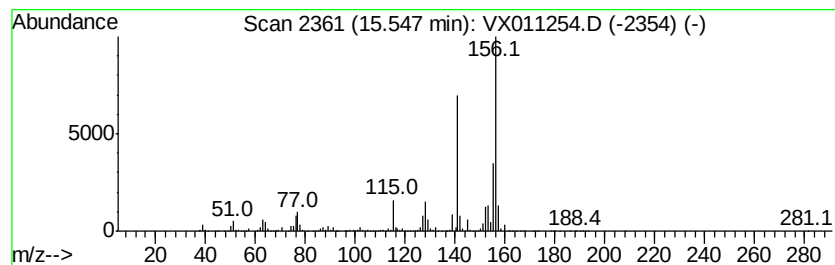
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	44.30 ug/l	749016	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

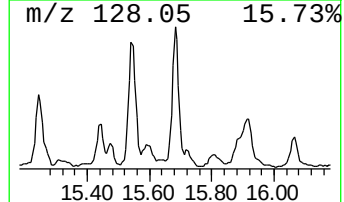
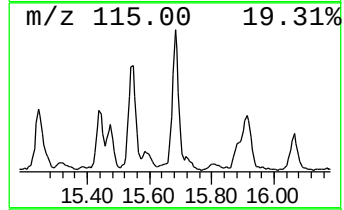
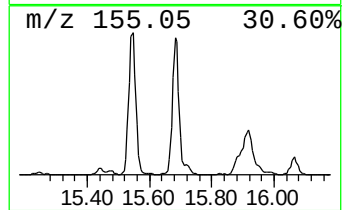
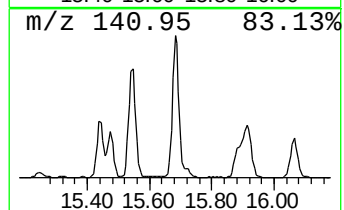
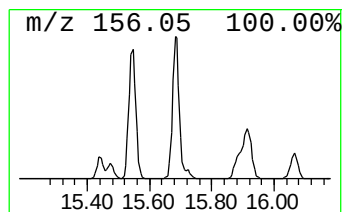
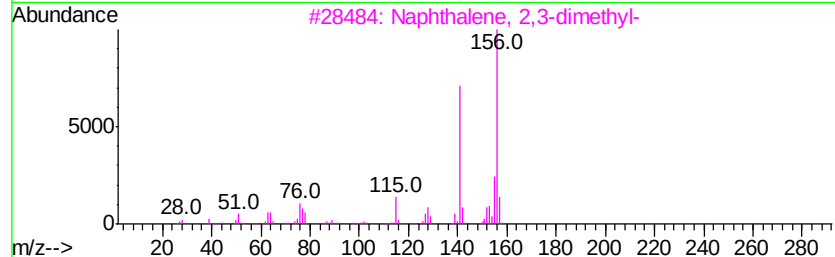
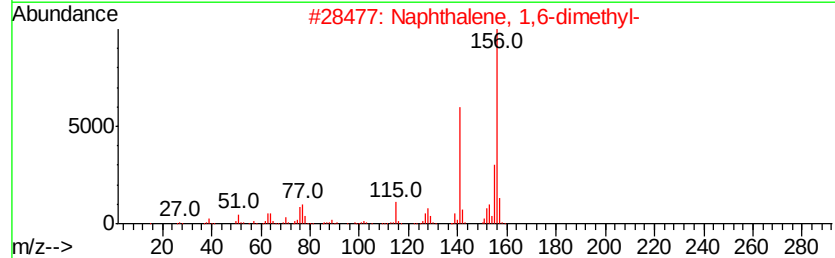
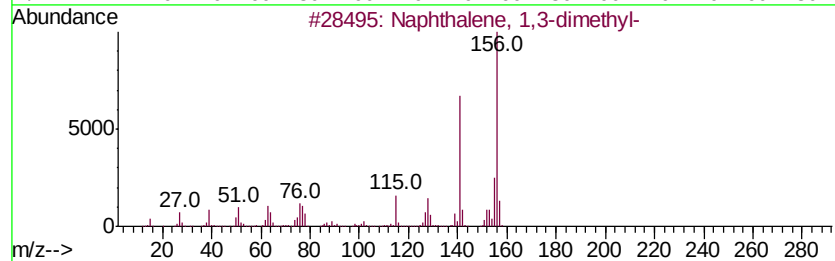
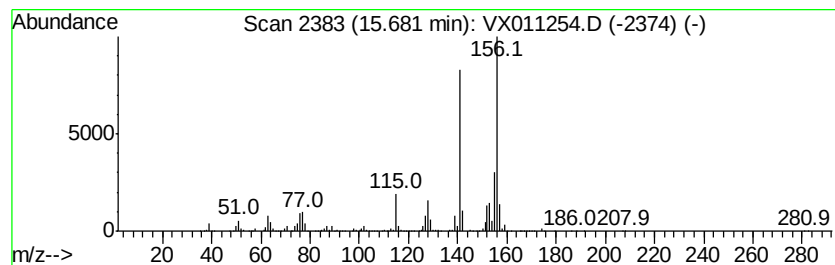
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	50.46 ug/l	853116	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073119\
Data File : VX011254.D
Acq On : 31 Jul 2019 19:37
Operator : JC/SP
Sample : K4048-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C1-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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, cyclopro...	12.30	49.0	ug/l	827682	4	12.08	845380	50.0
Benzene, 4-ethyl-...	12.66	44.8	ug/l	757845	4	12.08	845380	50.0
Benzene, (2-methy...	12.76	70.0	ug/l	1183530	4	12.08	845380	50.0
Benzene, 1,2,3,4-...	12.97	41.9	ug/l	708048	4	12.08	845380	50.0
1H-Indene, 2,3-di...	13.22	38.5	ug/l	650356	4	12.08	845380	50.0
Benzene, 2-etheny...	13.35	142.5	ug/l	2409780	4	12.08	845380	50.0
Naphthalene, 1,2,...	14.29	66.0	ug/l	1116390	4	12.08	845380	50.0
Naphthalene, 2,6-...	15.55	44.3	ug/l	749016	4	12.08	845380	50.0
Naphthalene, 1,3-...	15.68	50.5	ug/l	853116	4	12.08	845380	50.0