

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX033120\
 Data File : VX015450.D
 Acq On : 31 Mar 2020 13:49
 Operator : JC/SP
 Sample : L2050-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WATER-COMP

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\82X032720W.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	1.064	25	29	43	rBV	3934571	4425907	10.37%	1.086%
2	2.417	245	251	261	rBV	761412	1221430	2.86%	0.300%
3	5.472	739	752	762	rBV	670973	1912466	4.48%	0.469%
4	5.636	768	779	790	rBV	1143403	3178039	7.45%	0.780%
5	6.039	835	845	852	rBV	772563	2032282	4.76%	0.499%
6	6.118	852	858	870	rVB2	236192	638150	1.50%	0.157%
7	6.831	965	975	988	rBV	1822059	4394949	10.30%	1.079%
8	8.697	1275	1281	1287	rBV	3943005	6035259	14.15%	1.481%
9	8.770	1287	1293	1304	rVB	4963504	7878909	18.47%	1.934%
10	10.099	1505	1511	1521	rVB	4143637	5475614	12.83%	1.344%
11	10.239	1528	1534	1545	rVB	2055242	2775639	6.51%	0.681%
12	10.343	1545	1551	1559	rBV	16304729	21750912	50.98%	5.338%
13	10.684	1601	1607	1619	rBV	11104056	14635971	34.31%	3.592%
14	11.001	1653	1659	1662	rBV	614246	809146	1.90%	0.199%
15	11.123	1673	1679	1685	rVV	3649141	4580309	10.74%	1.124%
16	11.343	1710	1715	1722	rBV	2251544	2809449	6.59%	0.690%
17	11.422	1722	1728	1735	rVV	17388841	30136155	70.64%	7.396%
18	11.495	1735	1740	1750	rVB	9189774	11800242	27.66%	2.896%
19	11.654	1760	1766	1774	rVB	10083227	12455052	29.19%	3.057%
20	11.794	1783	1789	1794	rBV	33593361	42663509	100.00%	10.471%
21	11.885	1801	1804	1809	rBV2	274703	507656	1.19%	0.125%
22	11.934	1809	1812	1817	rVB	419330	523565	1.23%	0.128%
23	12.013	1817	1825	1828	rBV	1302803	1628744	3.82%	0.400%
24	12.062	1828	1833	1839	rVB2	4847215	6232230	14.61%	1.530%
25	12.129	1839	1844	1849	rBV	15815996	19049440	44.65%	4.675%
26	12.257	1861	1865	1867	rBV	12500114	14879410	34.88%	3.652%
27	12.288	1867	1870	1872	rVB	5127558	6473990	15.17%	1.589%
28	12.361	1877	1882	1891	rVB2	7735930	11499853	26.95%	2.822%
29	12.446	1892	1896	1901	rBV2	820174	1092662	2.56%	0.268%
30	12.501	1901	1905	1911	rVB	2584013	3161135	7.41%	0.776%
31	12.574	1911	1917	1919	rBV	5719697	8148911	19.10%	2.000%
32	12.592	1919	1920	1925	rVB	5187610	5018224	11.76%	1.232%
33	12.653	1925	1930	1934	rBV	10370405	12088730	28.34%	2.967%
34	12.684	1934	1935	1940	rVB	723748	652640	1.53%	0.160%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX033120\
 Data File : VX015450.D
 Acq On : 31 Mar 2020 13:49
 Operator : JC/SP
 Sample : L2050-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WATER-COMP

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\82X032720W.M
 Title : SW846 8260

35	12.745	1940	1945	1953	rBV4	2317398	4651589	10.90%	1.142%
36	12.836	1956	1960	1964	rVB	417878	445581	1.04%	0.109%
37	12.891	1964	1969	1974	rVB	3220334	4075966	9.55%	1.000%
38	12.964	1974	1981	1984	rBV	7329115	9425952	22.09%	2.313%
39	13.001	1984	1987	1993	rVB	11672749	14148458	33.16%	3.472%
40	13.062	1993	1997	1999	rBV	591702	772659	1.81%	0.190%
41	13.208	2014	2021	2025	rBV	5209877	6669514	15.63%	1.637%
42	13.251	2025	2028	2031	rVB2	583290	623573	1.46%	0.153%
43	13.293	2031	2035	2037	rBV2	925775	1256811	2.95%	0.308%
44	13.330	2037	2041	2047	rVB2	13202394	18342624	42.99%	4.502%
45	13.409	2052	2054	2058	rVV	701626	684521	1.60%	0.168%
46	13.464	2058	2063	2071	rVB	5488740	6778837	15.89%	1.664%
47	13.543	2071	2076	2079	rBV2	700481	992143	2.33%	0.243%
48	13.586	2079	2083	2085	rVV	1157970	1480827	3.47%	0.363%
49	13.623	2085	2089	2096	rVV3	2058888	4252466	9.97%	1.044%
50	13.684	2096	2099	2100	rVV	653988	820055	1.92%	0.201%
51	13.708	2100	2103	2108	rVB2	1586002	2575883	6.04%	0.632%
52	13.818	2114	2121	2127	rBV	8474719	10878085	25.50%	2.670%
53	13.897	2129	2134	2139	rVB2	2102077	2761556	6.47%	0.678%
54	13.964	2139	2145	2150	rBV2	2318023	3422420	8.02%	0.840%
55	14.013	2150	2153	2156	rVB	791779	852003	2.00%	0.209%
56	14.117	2166	2170	2174	rBV	2060509	2458909	5.76%	0.603%
57	14.269	2188	2195	2203	rBV	5752750	9227938	21.63%	2.265%
58	14.360	2207	2210	2215	rVV	881331	1023788	2.40%	0.251%
59	14.415	2215	2219	2223	rVB	1825799	2175881	5.10%	0.534%
60	14.525	2231	2237	2242	rBV2	3670432	5030433	11.79%	1.235%
61	14.659	2254	2259	2260	rBV2	1776079	2098798	4.92%	0.515%
62	14.677	2260	2262	2265	rVV	2471186	2833530	6.64%	0.695%
63	14.714	2265	2268	2273	rVB	1161149	1495202	3.50%	0.367%
64	14.769	2273	2277	2280	rBV2	370078	454102	1.06%	0.111%
65	14.824	2280	2286	2289	rVV	1158804	1540281	3.61%	0.378%
66	14.860	2289	2292	2295	rVV	673662	993585	2.33%	0.244%
67	14.891	2295	2297	2302	rVB2	515216	547900	1.28%	0.134%
68	14.946	2302	2306	2313	rBV2	677749	1211502	2.84%	0.297%
69	15.098	2324	2331	2336	rBV3	360642	620946	1.46%	0.152%
70	15.232	2348	2353	2360	rVV2	1506748	2813769	6.60%	0.691%
71	15.427	2380	2385	2388	rBV	368890	577414	1.35%	0.142%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

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\82X032720W.M
Title : SW846 8260

72	15.531	2394	2402	2406	rBV	801364	1325720	3.11%	0.325%
73	15.574	2406	2409	2414	rVB3	277709	473805	1.11%	0.116%
74	15.665	2419	2424	2428	rVV	834597	1464502	3.43%	0.359%
75	15.702	2428	2430	2437	rVB2	399613	610006	1.43%	0.150%

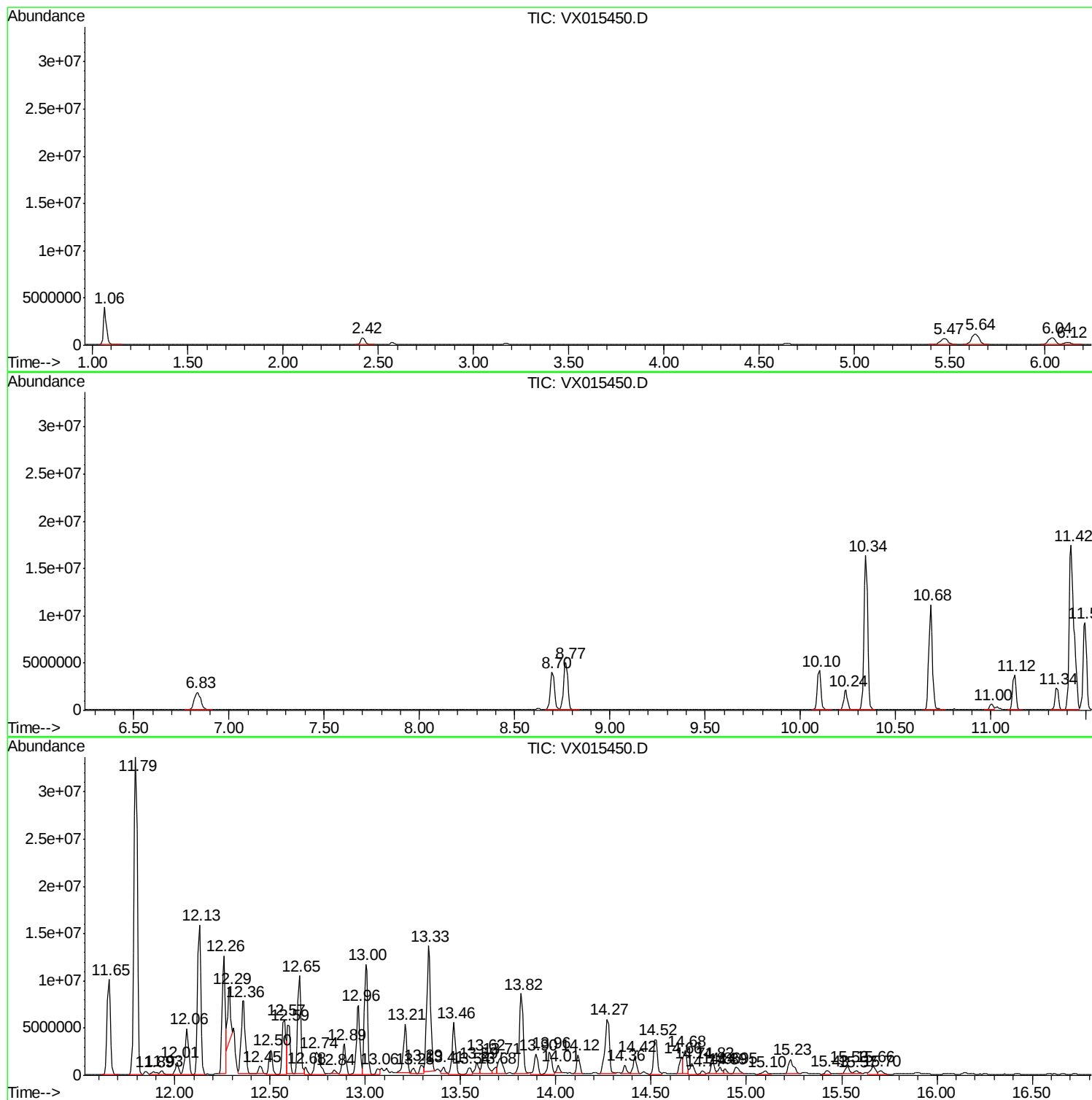
Sum of corrected areas: 407456113

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

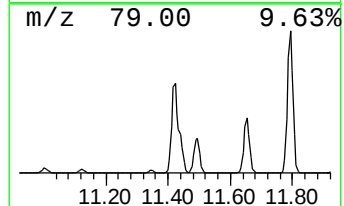
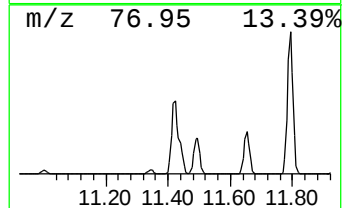
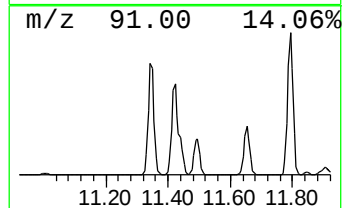
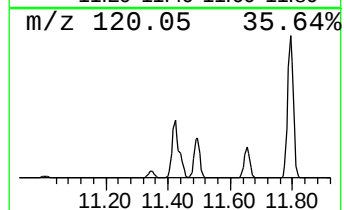
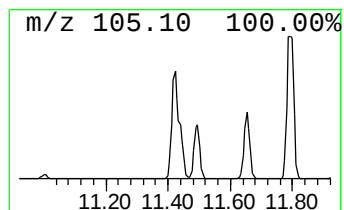
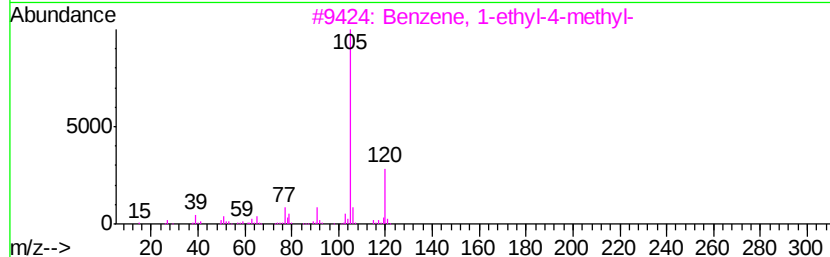
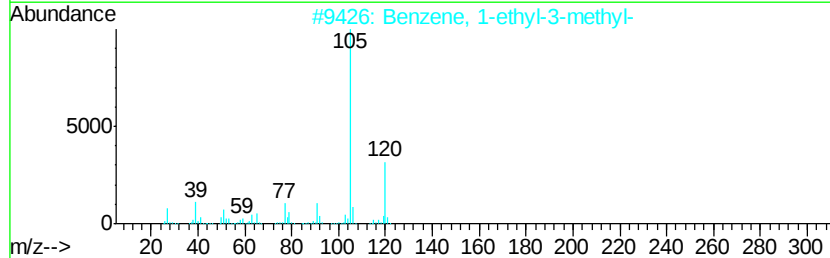
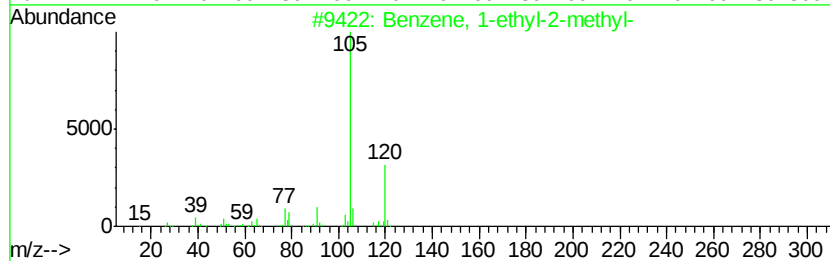
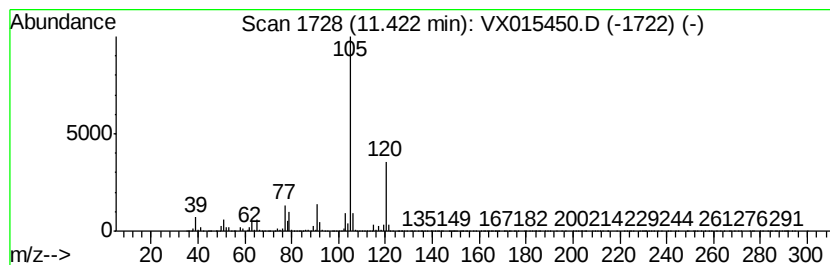
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	241.78 ug/l	30136200	1,4-Dichlorobenzene-d4	12.06

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	94
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
WATER-COMP

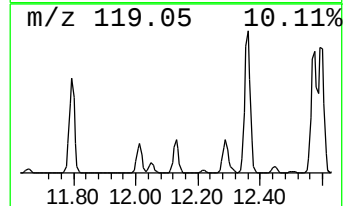
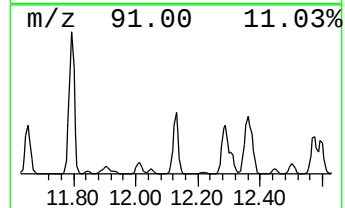
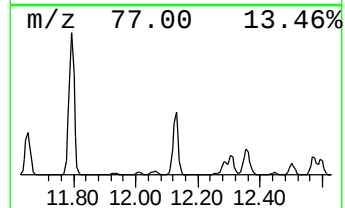
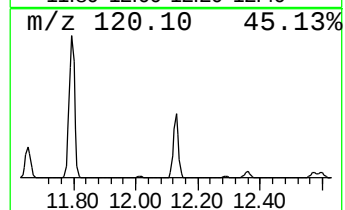
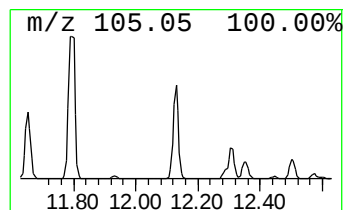
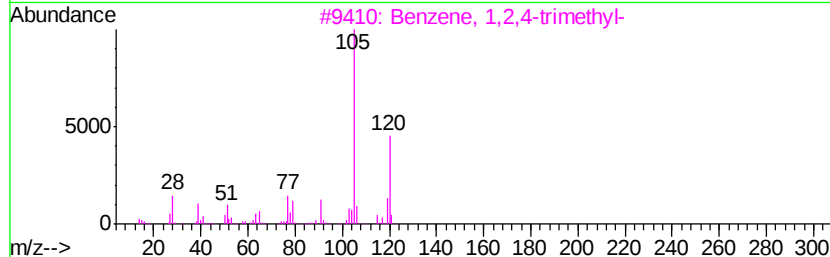
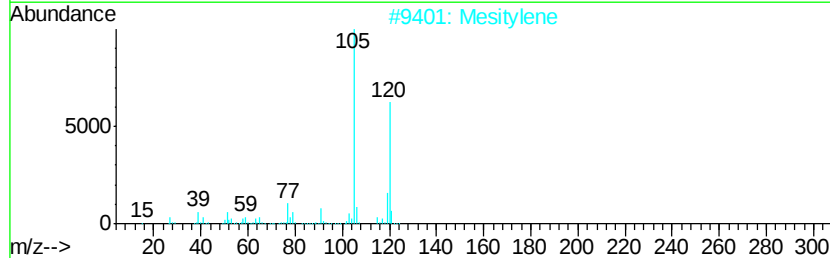
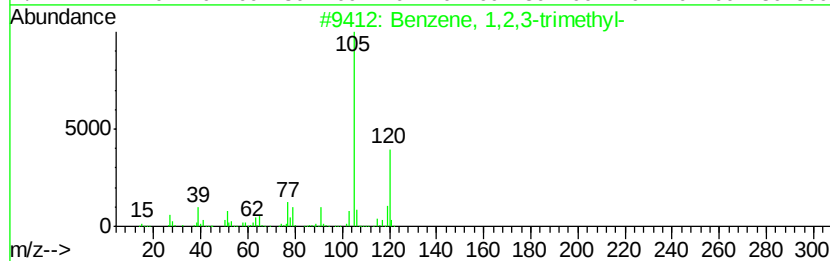
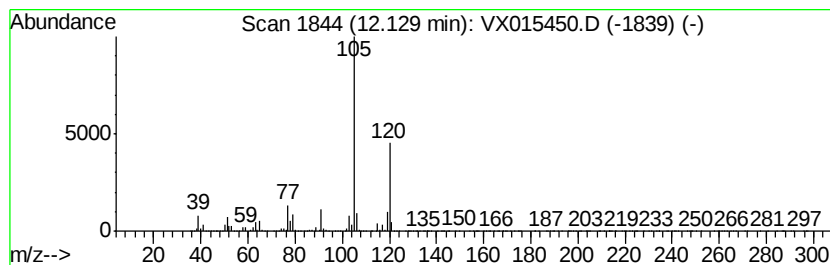
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	152.83 ug/l	19049400	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

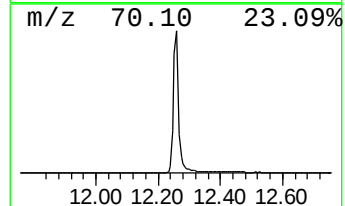
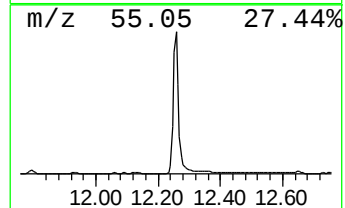
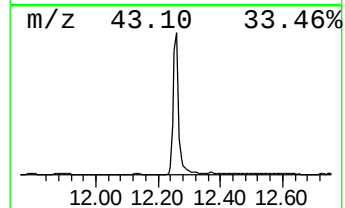
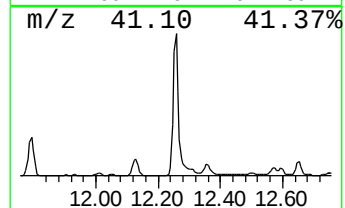
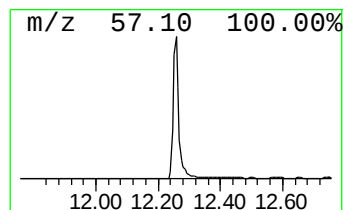
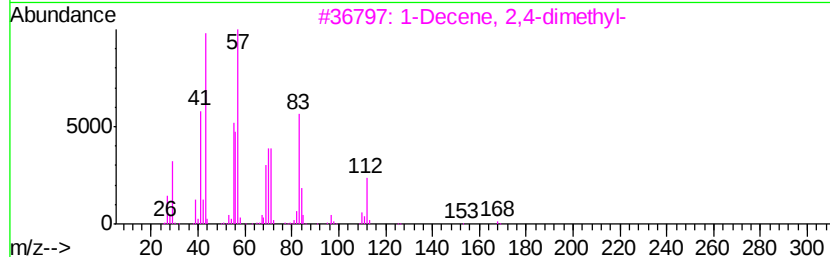
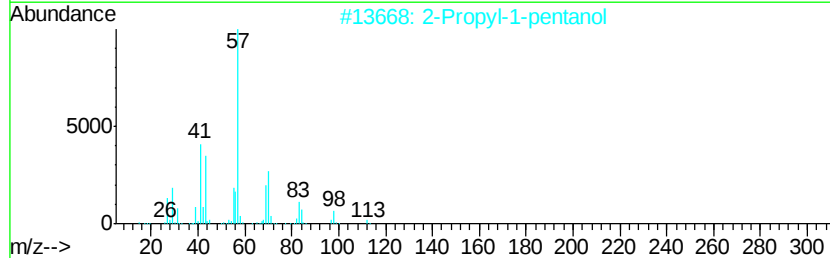
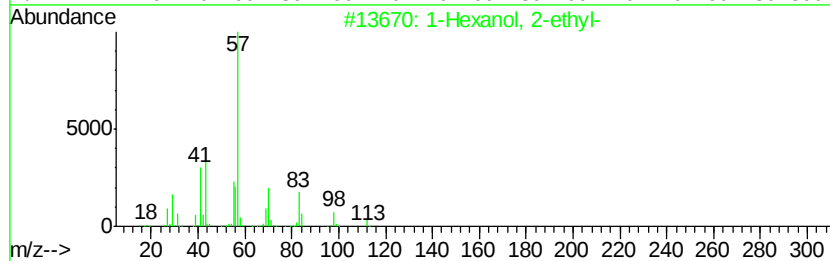
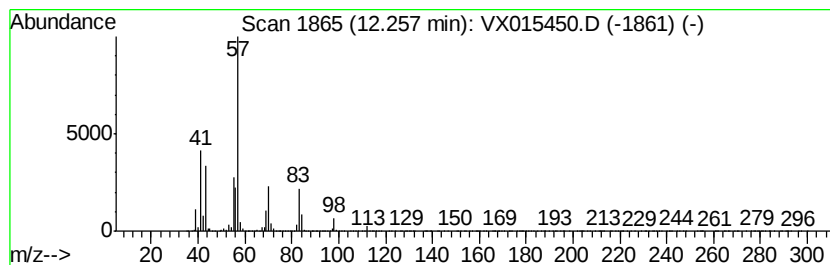
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	119.38 ug/l	14879400	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
3		1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	47
4		2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	47
5		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

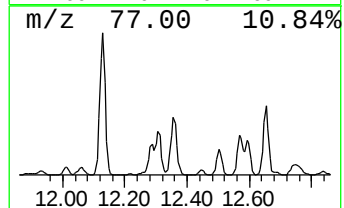
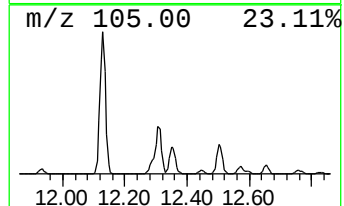
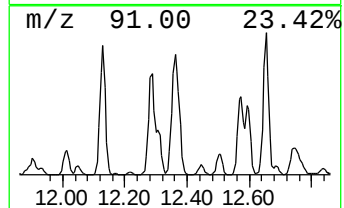
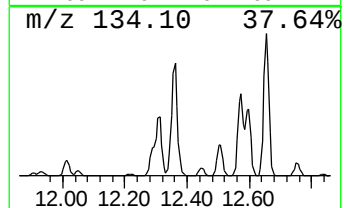
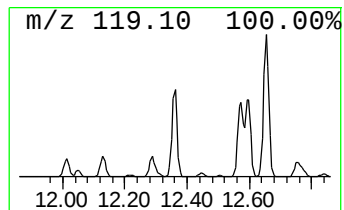
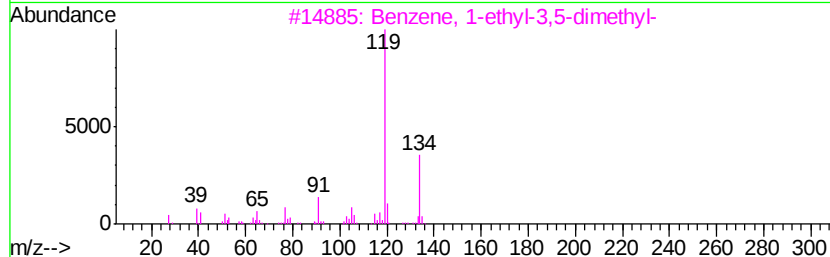
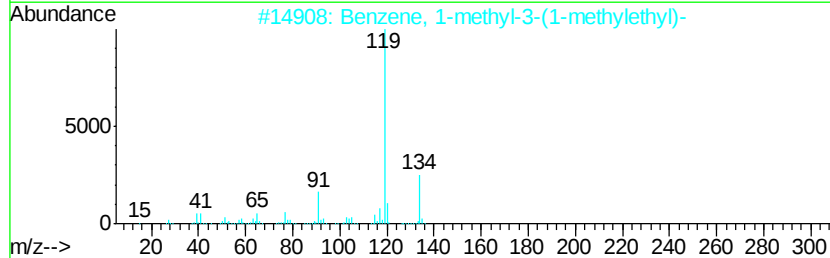
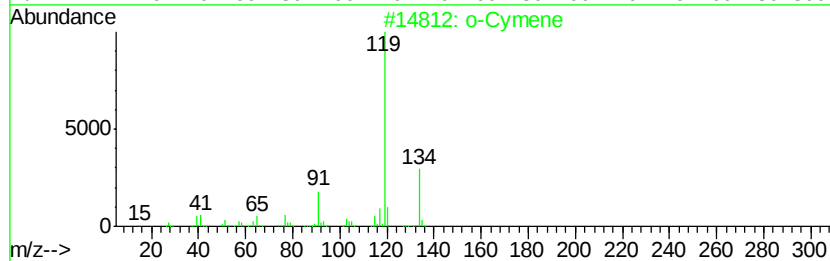
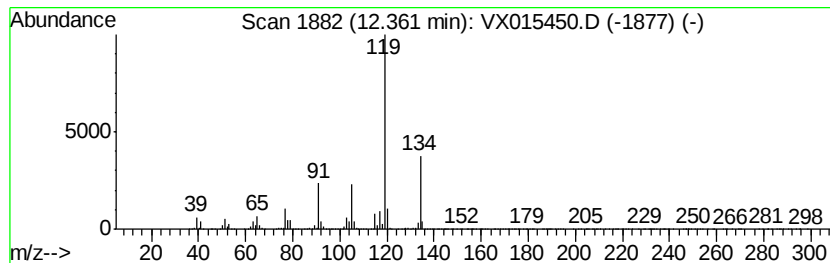
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	92.26 ug/l	11499900	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	95
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

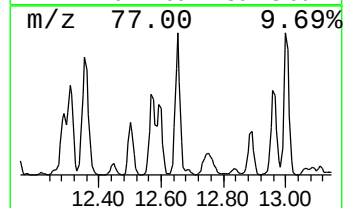
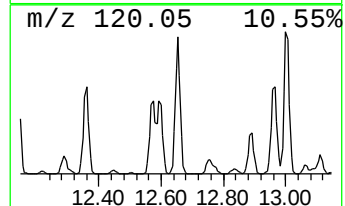
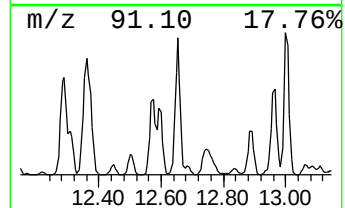
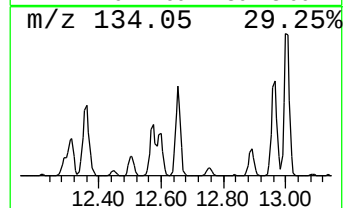
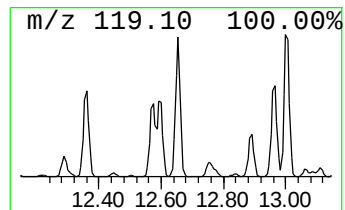
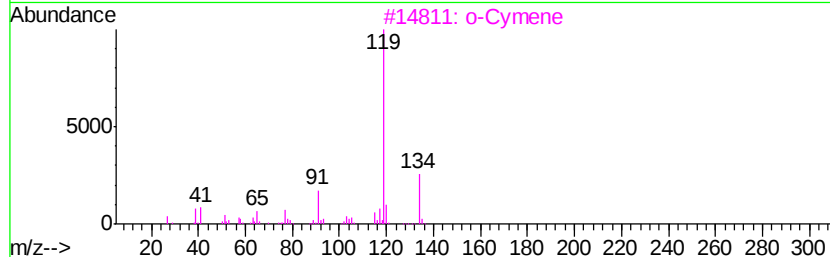
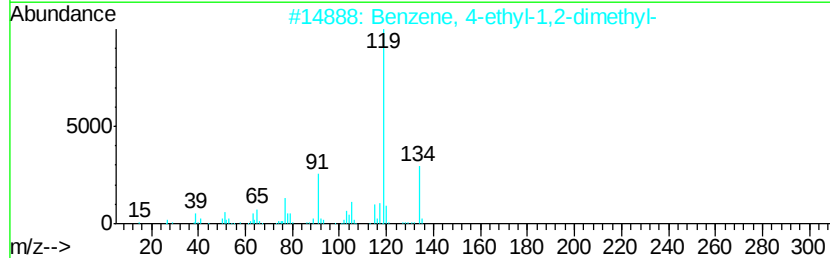
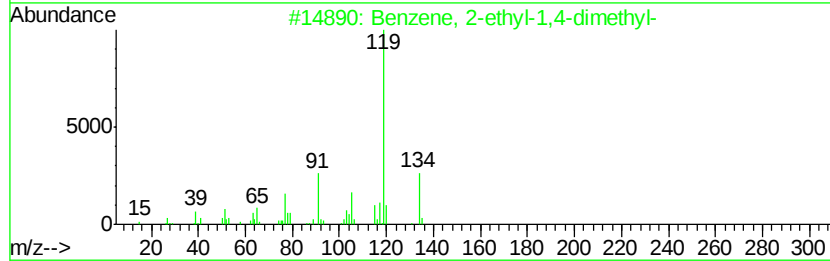
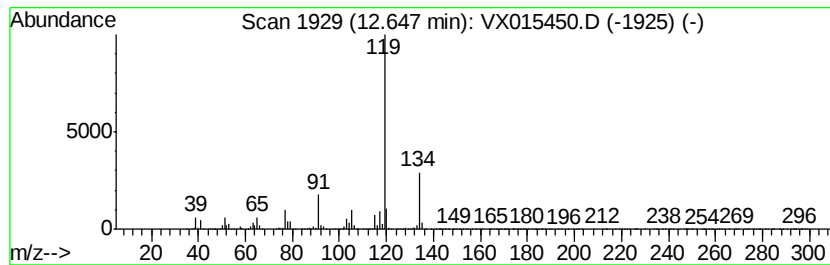
Instrument :
MSVOA_X
ClientSampled :
WATER-COMP

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.65	96.99 ug/l	12088700	1,4-Dichlorobenzene-d4	12.06		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		o-Cymene	134	C10H14	000527-84-4	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:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

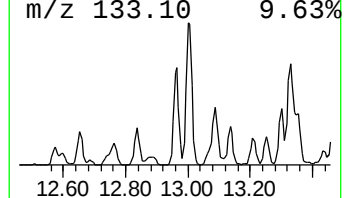
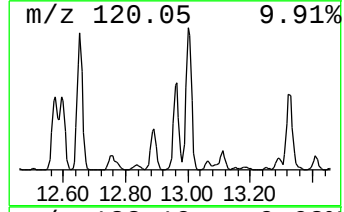
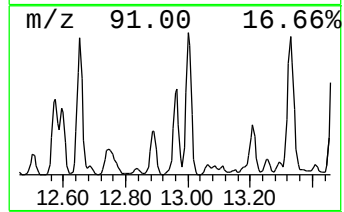
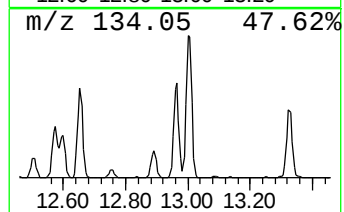
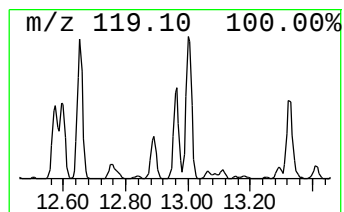
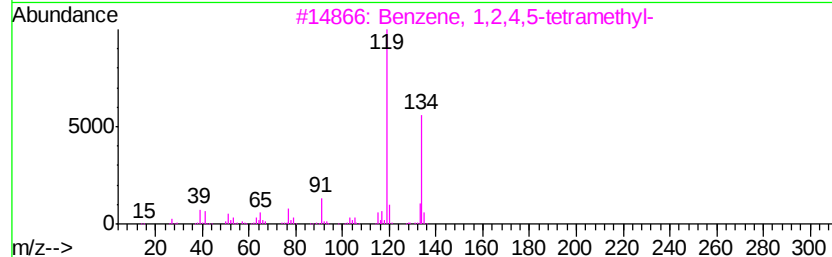
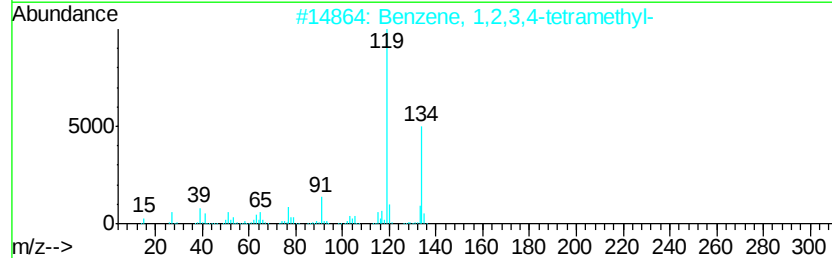
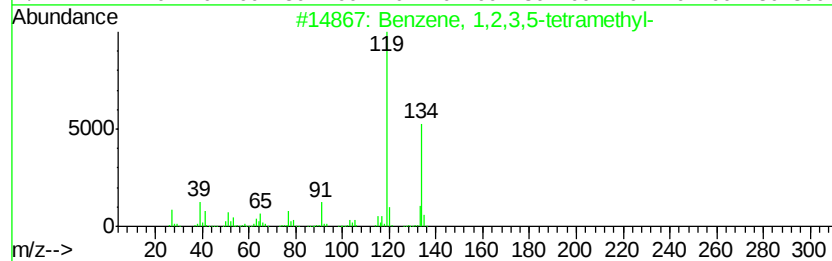
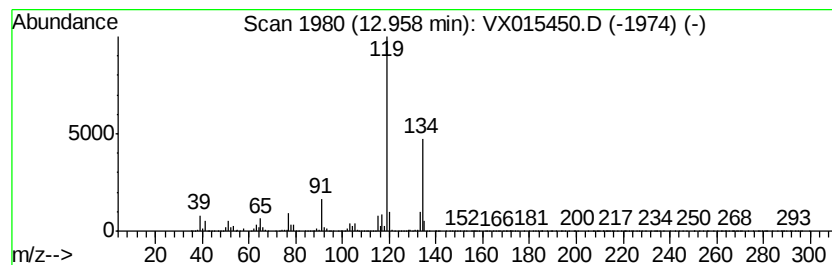
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	75.62 ug/l	9425950	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

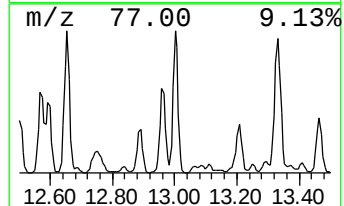
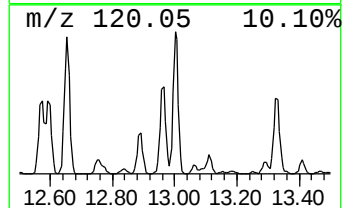
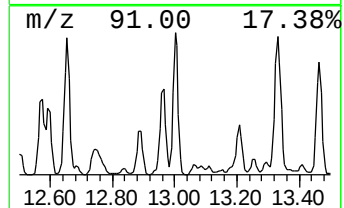
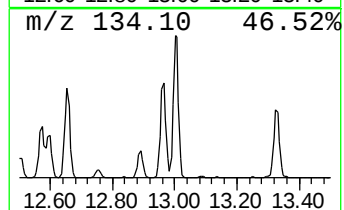
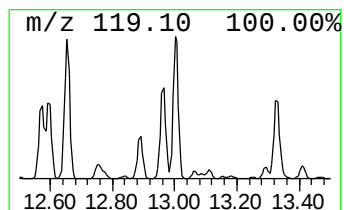
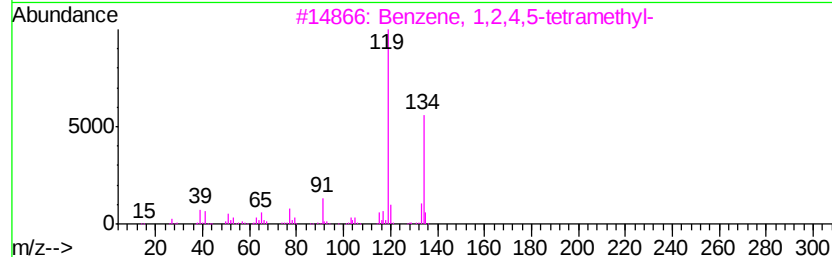
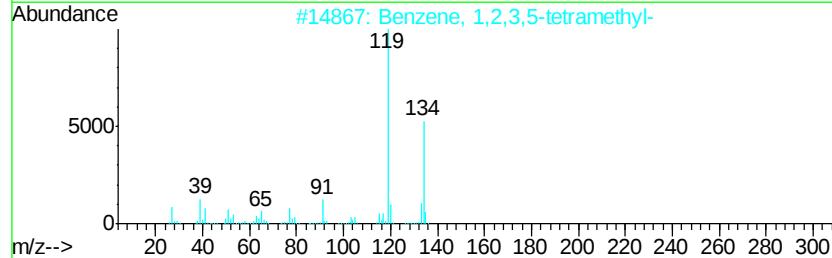
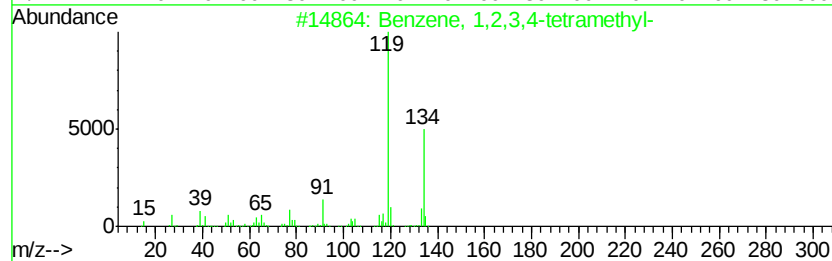
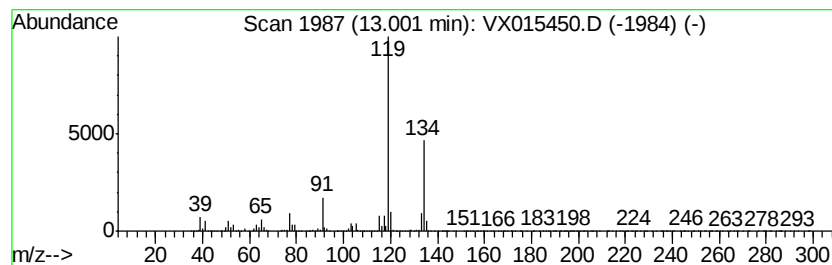
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	113.51 ug/l	14148500	1,4-Dichlorobenzene-d4	12.06

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	96
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

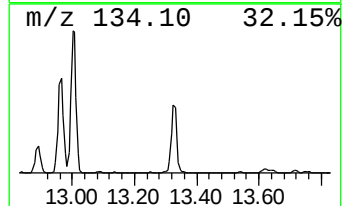
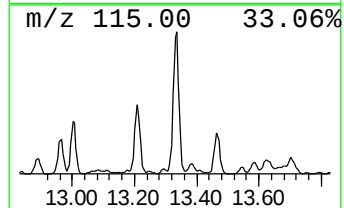
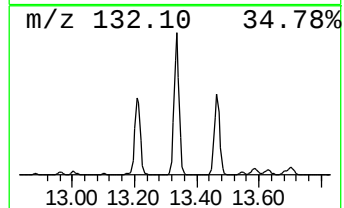
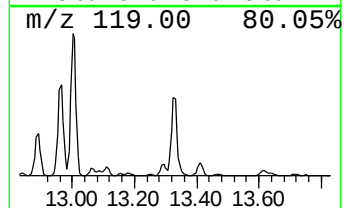
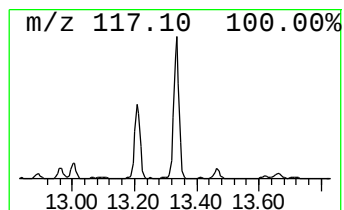
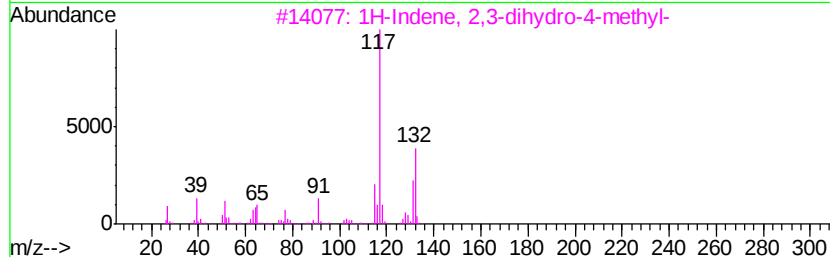
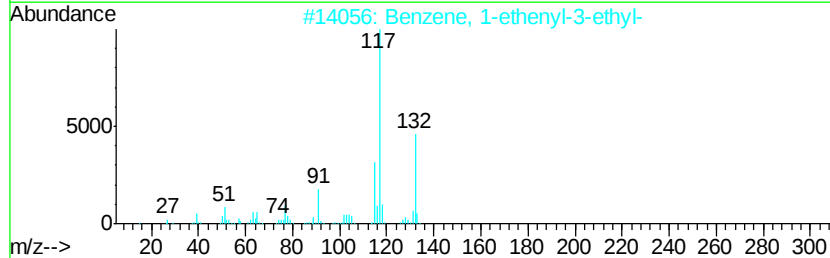
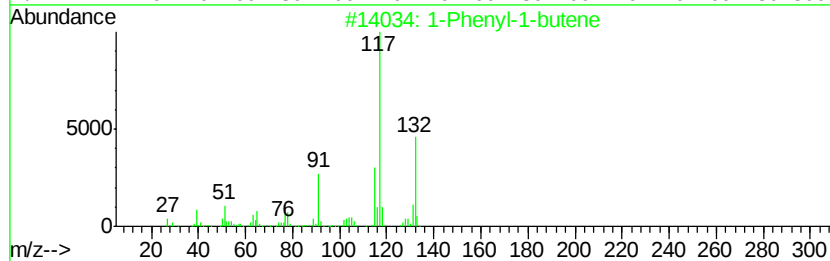
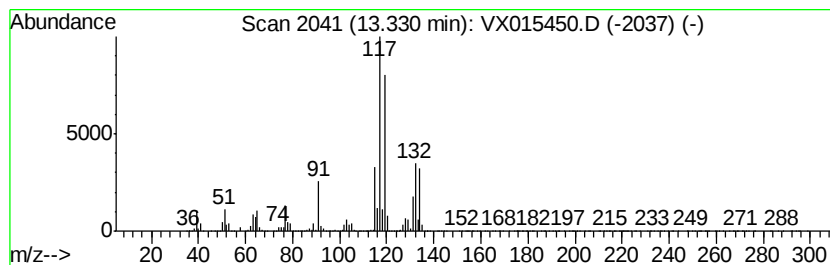
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	147.16 ug/l	18342600	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	86
5		2,4-Dimethylstyrene	132	C10H12	002234-20-0	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

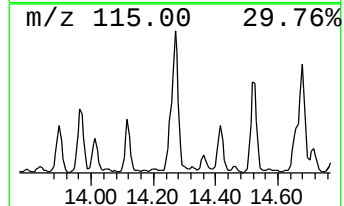
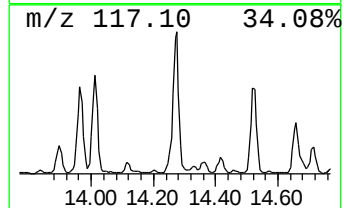
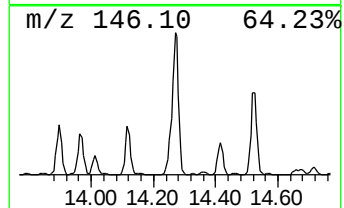
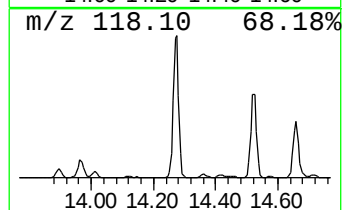
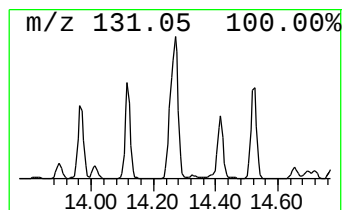
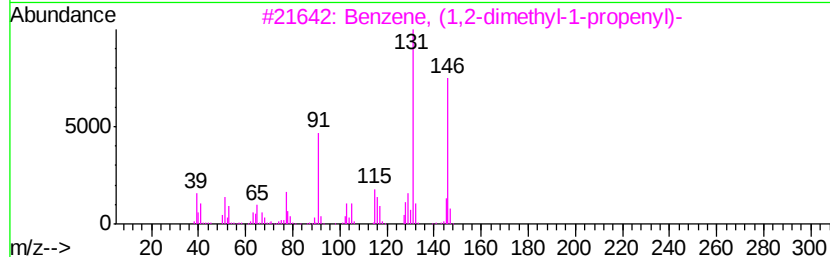
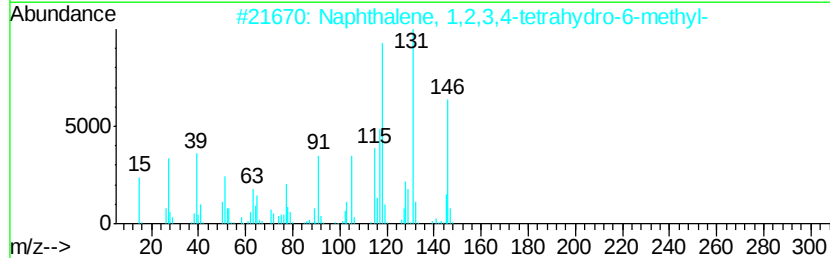
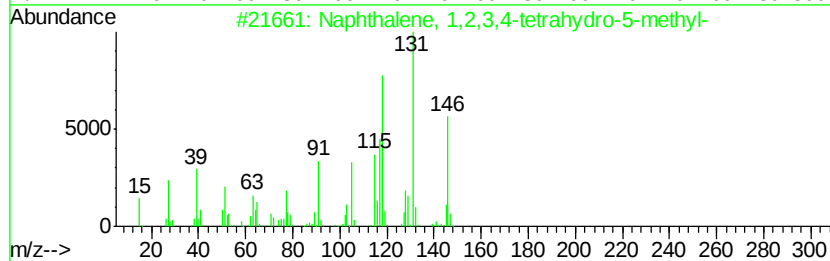
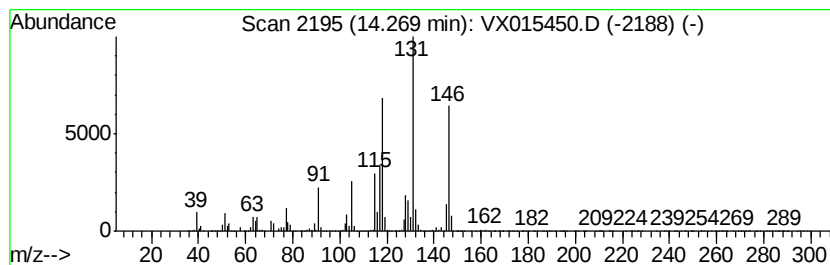
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	74.03 ug/l	9227940	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	97
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	92
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
5		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX033120\
Data File : VX015450.D
Acq On : 31 Mar 2020 13:49
Operator : JC/SP
Sample : L2050-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.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-...	11.42	241.8	ug/l	30136200	4	12.06	6232230	50.0
Benzene, 1,2,3-tr...	12.13	152.8	ug/l	19049400	4	12.06	6232230	50.0
1-Hexanol, 2-ethyl-	12.26	119.4	ug/l	14879400	4	12.06	6232230	50.0
o-Cymene	12.36	92.3	ug/l	11499900	4	12.06	6232230	50.0
Benzene, 2-ethyl-...	12.65	97.0	ug/l	12088700	4	12.06	6232230	50.0
Benzene, 1,2,3,5-...	12.96	75.6	ug/l	9425950	4	12.06	6232230	50.0
Benzene, 1,2,3,4-...	13.00	113.5	ug/l	14148500	4	12.06	6232230	50.0
1-Phenyl-1-butene	13.33	147.2	ug/l	18342600	4	12.06	6232230	50.0
Naphthalene, 1,2,...	14.27	74.0	ug/l	9227940	4	12.06	6232230	50.0