

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122618\
 Data File : VX006816.D
 Acq On : 26 Dec 2018 14:16
 Operator : JC/MD
 Sample : J6549-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TW-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\82X121818W.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	1.276	17	20	27	rBV2	31153	47359	1.59%	0.146%
2	1.623	66	77	81	rBV4	39441	114030	3.83%	0.350%
3	2.446	207	212	218	rBV	23870	38105	1.28%	0.117%
4	5.501	701	713	726	rBV	355345	1002081	33.62%	3.079%
5	5.659	729	739	757	rVB	643320	1806639	60.62%	5.551%
6	6.068	796	806	813	rBV	371818	957320	32.12%	2.942%
7	6.141	813	818	828	rVB3	89148	231602	7.77%	0.712%
8	6.860	924	936	950	rBV	925265	2169907	72.81%	6.668%
9	8.714	1234	1240	1248	rBV	1958909	2980235	100.00%	9.158%
10	8.781	1248	1251	1261	rVB	19034	35875	1.20%	0.110%
11	9.518	1367	1372	1377	rBV	47341	71043	2.38%	0.218%
12	10.110	1461	1469	1482	rBV	1924451	2646282	88.79%	8.132%
13	10.250	1486	1492	1503	rVV	1276618	1618364	54.30%	4.973%
14	10.360	1503	1510	1515	rVB	60069	96036	3.22%	0.295%
15	10.695	1561	1565	1571	rVB	48859	67868	2.28%	0.209%
16	10.939	1596	1605	1606	rBV8	15165	36819	1.24%	0.113%
17	11.018	1612	1618	1625	rBV	166824	230947	7.75%	0.710%
18	11.134	1631	1637	1649	rBV	1671399	2187145	73.39%	6.721%
19	11.359	1669	1674	1680	rVB	248835	310790	10.43%	0.955%
20	11.506	1693	1698	1703	rVB2	29648	44984	1.51%	0.138%
21	11.664	1720	1724	1732	rBV	132292	170099	5.71%	0.523%
22	11.768	1735	1741	1744	rVV	100770	135249	4.54%	0.416%
23	11.804	1744	1747	1751	rVB	51976	66055	2.22%	0.203%
24	11.920	1761	1766	1767	rBV	28555	33840	1.14%	0.104%
25	11.945	1767	1770	1777	rVB	62272	79496	2.67%	0.244%
26	12.079	1786	1792	1797	rBV	2107796	2687635	90.18%	8.259%
27	12.140	1797	1802	1808	rBV	429342	530371	17.80%	1.630%
28	12.231	1813	1817	1820	rVB	59671	72500	2.43%	0.223%
29	12.298	1820	1828	1836	rBV	1704444	2101093	70.50%	6.456%
30	12.371	1836	1840	1847	rVB3	79603	158790	5.33%	0.488%
31	12.463	1847	1855	1859	rBV3	94979	146909	4.93%	0.451%
32	12.518	1859	1864	1870	rVB4	41423	66967	2.25%	0.206%
33	12.585	1871	1875	1877	rBV	37855	47042	1.58%	0.145%
34	12.664	1884	1888	1892	rBV	637129	774265	25.98%	2.379%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122618\
 Data File : VX006816.D
 Acq On : 26 Dec 2018 14:16
 Operator : JC/MD
 Sample : J6549-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TW-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\82X121818W.M
 Title : SW846 8260

35	12.701	1892	1894	1899	rVV	98261	101469	3.40%	0.312%
36	12.756	1899	1903	1908	rVV2	446155	702437	23.57%	2.158%
37	12.810	1908	1912	1917	rVB2	349402	428649	14.38%	1.317%
38	12.896	1922	1926	1932	rVB2	75290	107345	3.60%	0.330%
39	12.975	1932	1939	1943	rBV	705765	922671	30.96%	2.835%
40	13.018	1943	1946	1951	rVB2	105364	138391	4.64%	0.425%
41	13.073	1951	1955	1961	rBV4	76204	152192	5.11%	0.468%
42	13.146	1965	1967	1971	rVB	42010	47960	1.61%	0.147%
43	13.194	1971	1975	1976	rBV3	27851	32784	1.10%	0.101%
44	13.225	1976	1980	1986	rVB	230331	283117	9.50%	0.870%
45	13.310	1989	1994	1995	rBV5	36013	49853	1.67%	0.153%
46	13.347	1995	2000	2005	rVV2	1177353	1681593	56.42%	5.167%
47	13.396	2005	2008	2011	rVB3	104063	116181	3.90%	0.357%
48	13.481	2017	2022	2026	rVB	311229	392714	13.18%	1.207%
49	13.530	2026	2030	2033	rBV2	118368	141415	4.75%	0.435%
50	13.603	2037	2042	2045	rBV	424645	548654	18.41%	1.686%
51	13.646	2045	2049	2054	rVB5	89400	171701	5.76%	0.528%
52	13.719	2054	2061	2067	rBV3	127197	231231	7.76%	0.711%
53	13.835	2076	2080	2088	rVB2	99565	196410	6.59%	0.604%
54	13.914	2088	2093	2094	rBV3	44237	59677	2.00%	0.183%
55	13.981	2100	2104	2108	rVV3	88805	115885	3.89%	0.356%
56	14.024	2108	2111	2123	rVB6	110143	186300	6.25%	0.572%
57	14.158	2129	2133	2138	rBV	92163	117831	3.95%	0.362%
58	14.213	2138	2142	2146	rVB	74103	102226	3.43%	0.314%
59	14.261	2146	2150	2159	rBV5	112640	177639	5.96%	0.546%
60	14.371	2165	2168	2173	rVB4	61600	96925	3.25%	0.298%
61	14.432	2173	2178	2187	rVB3	93850	192396	6.46%	0.591%
62	14.566	2193	2200	2210	rBV	263975	466858	15.67%	1.435%
63	14.700	2219	2222	2232	rVB	21344	41324	1.39%	0.127%
64	14.792	2234	2237	2241	rVV2	51344	79593	2.67%	0.245%
65	14.841	2241	2245	2253	rVB	90268	140569	4.72%	0.432%
66	14.956	2259	2264	2275	rVV3	183074	324929	10.90%	0.998%
67	15.048	2275	2279	2287	rVV5	55175	105555	3.54%	0.324%
68	15.151	2291	2296	2303	rVB	49177	92179	3.09%	0.283%
69	15.389	2331	2335	2339	rBV7	22997	33154	1.11%	0.102%

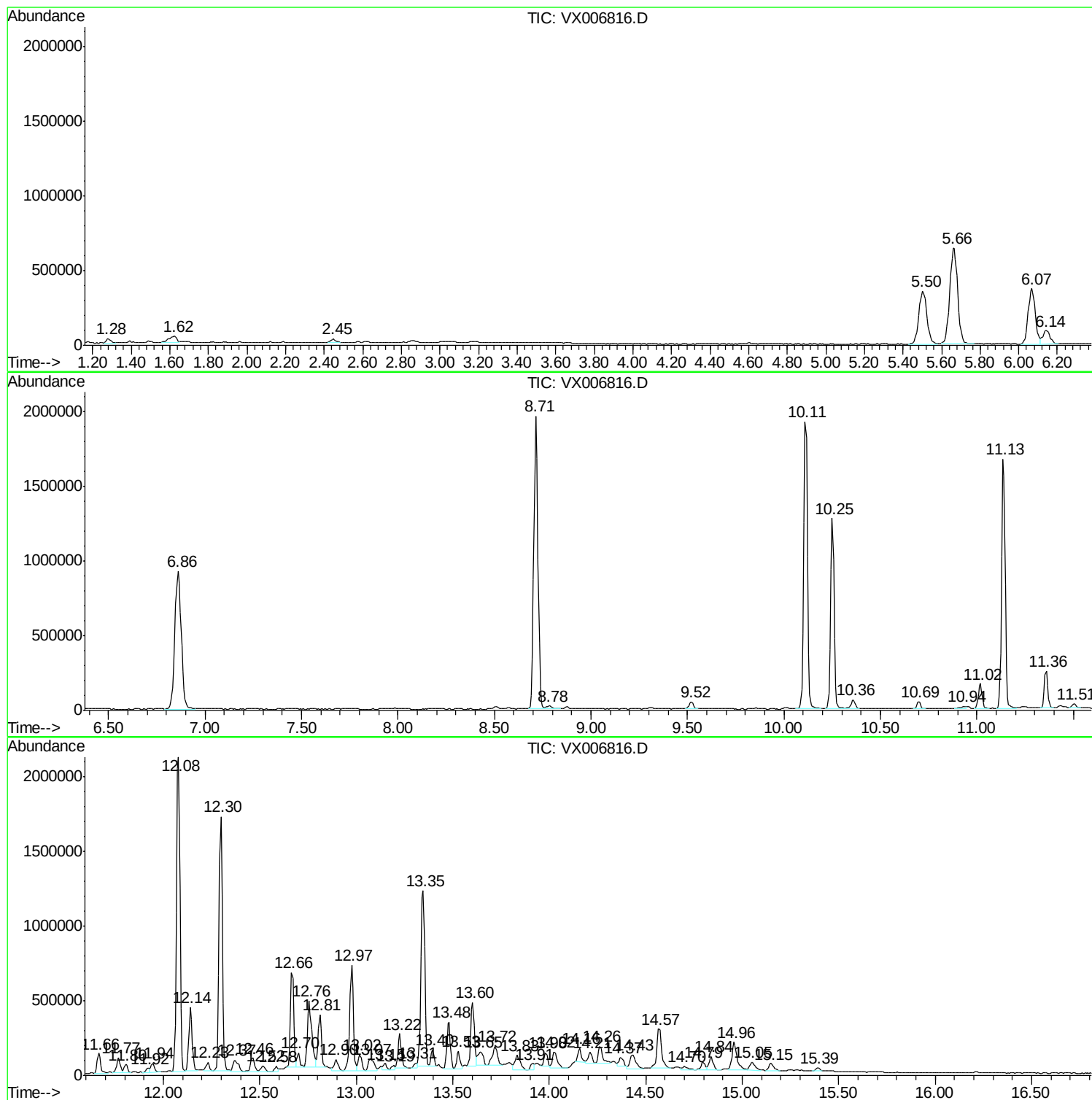
Sum of corrected areas: 32543549

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TW-1

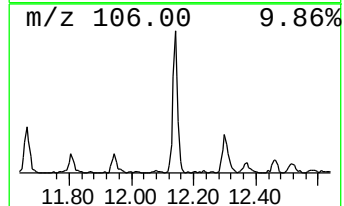
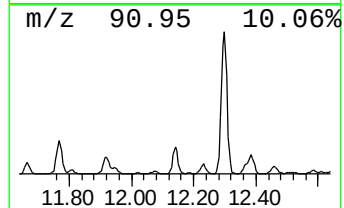
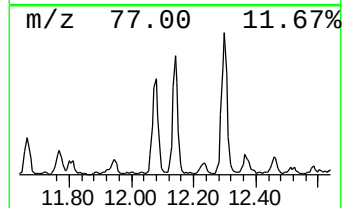
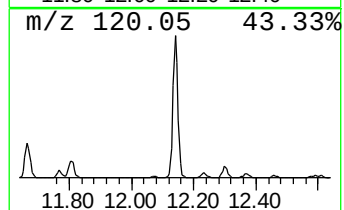
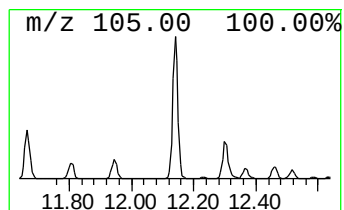
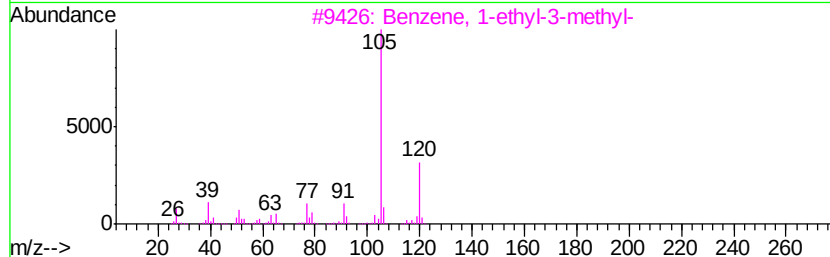
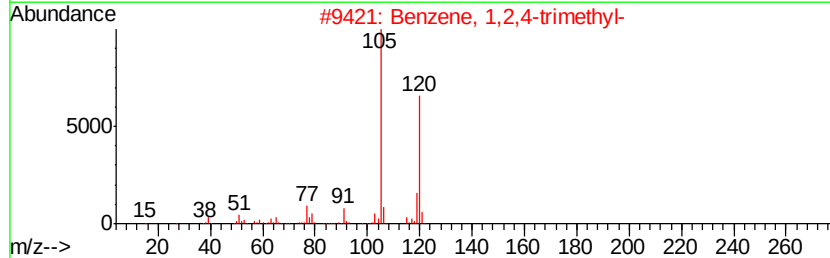
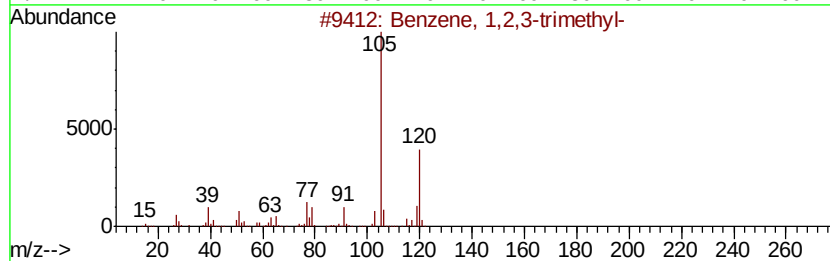
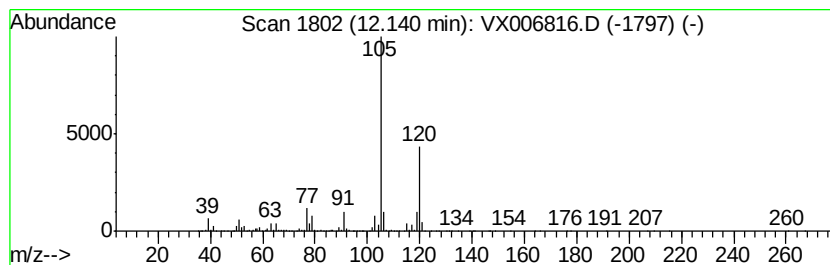
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	9.87 ug/l	530371	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

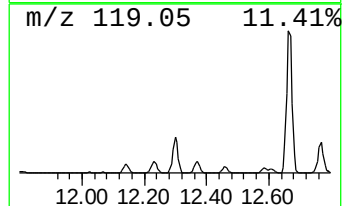
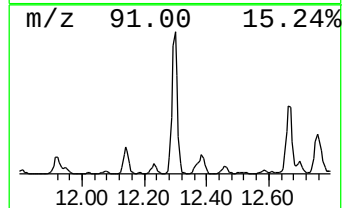
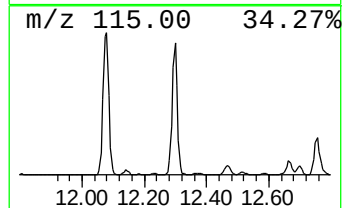
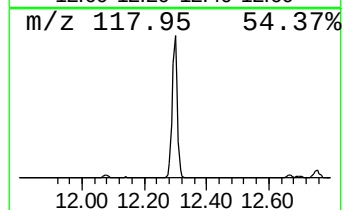
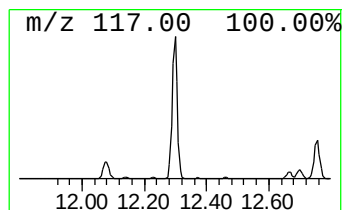
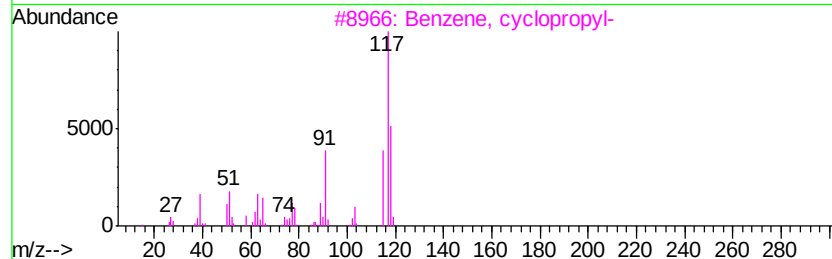
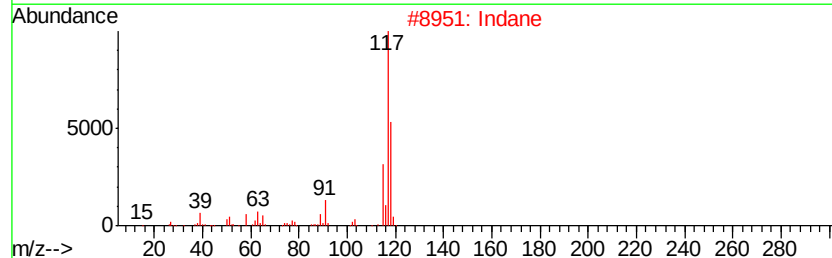
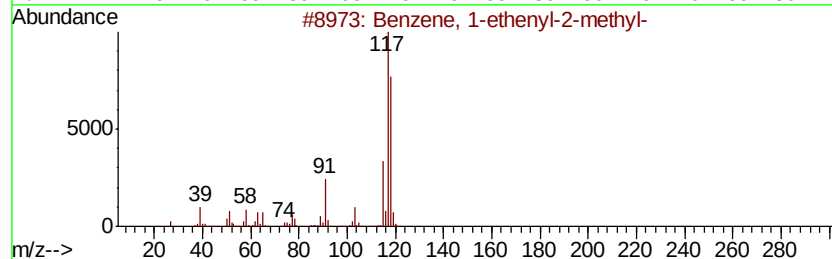
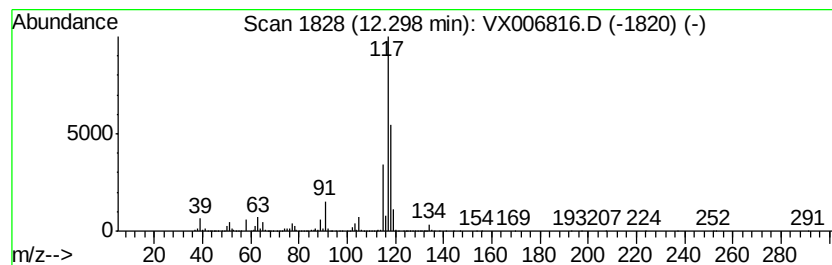
Instrument :
MSVOA_X
ClientSampleId :
TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	39.09 ug/l	2101090	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 81
2		Indane	118 C9H10	000496-11-7 81
3		Benzene, cyclopropyl-	118 C9H10	000873-49-4 81
4		Benzene, 2-propenyl-	118 C9H10	000300-57-2 74
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

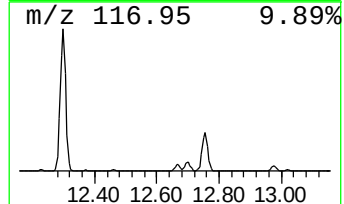
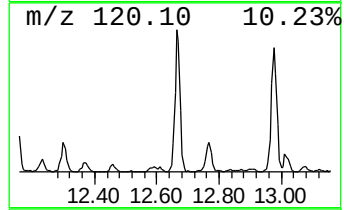
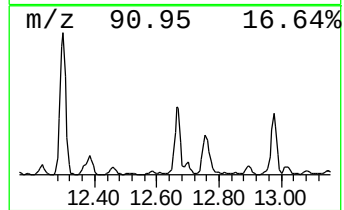
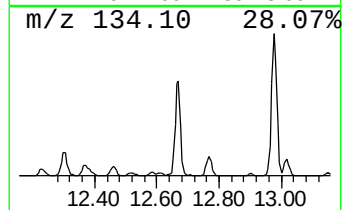
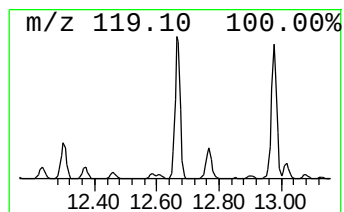
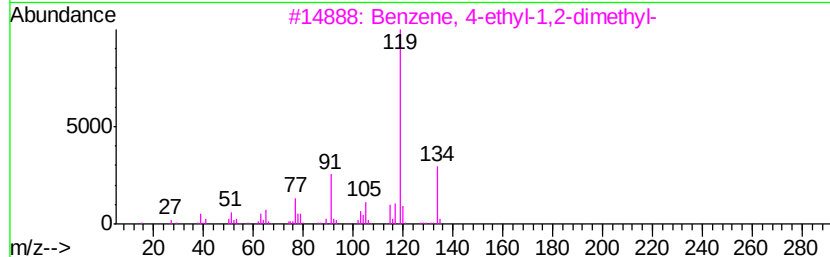
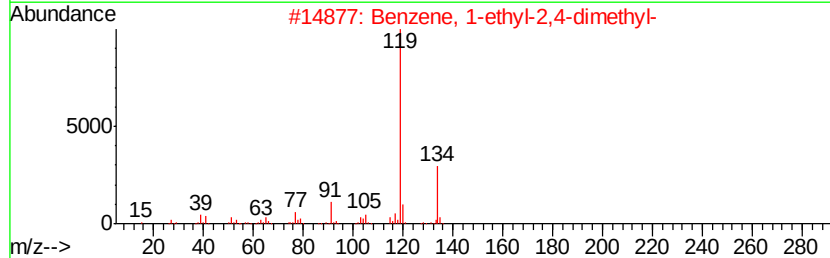
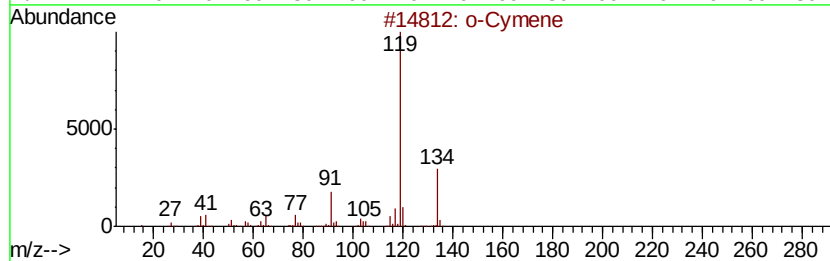
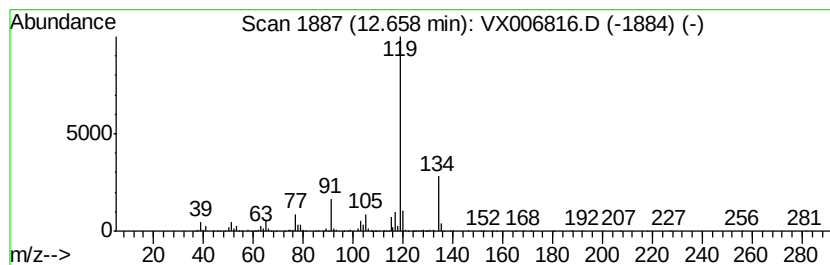
Instrument :
MSVOA_X
ClientSampleId :
TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 3 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	14.40 ug/l	774265	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 95
2		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 95
3		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 95
4		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 94
5		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-1

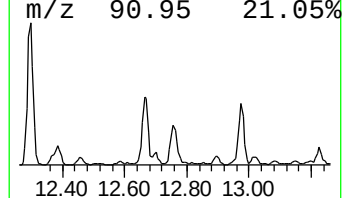
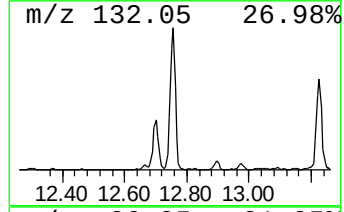
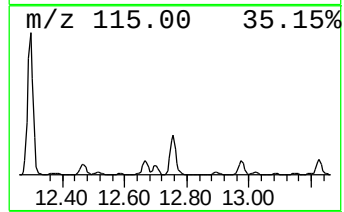
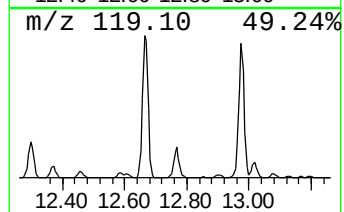
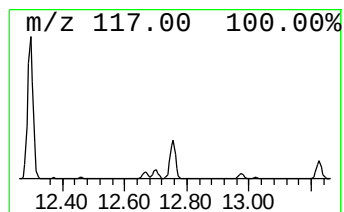
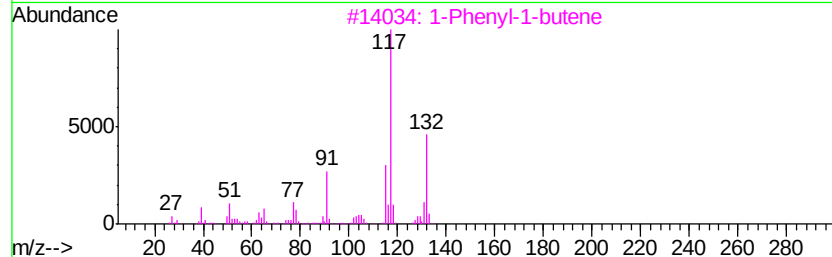
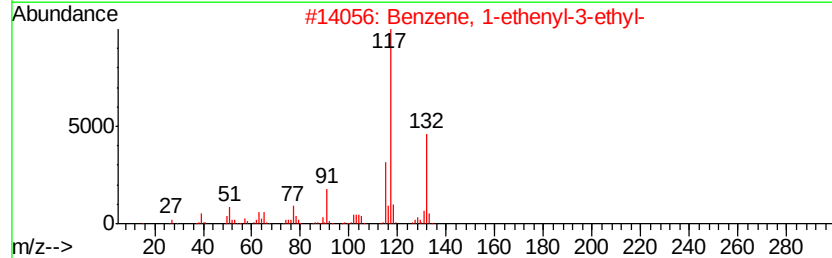
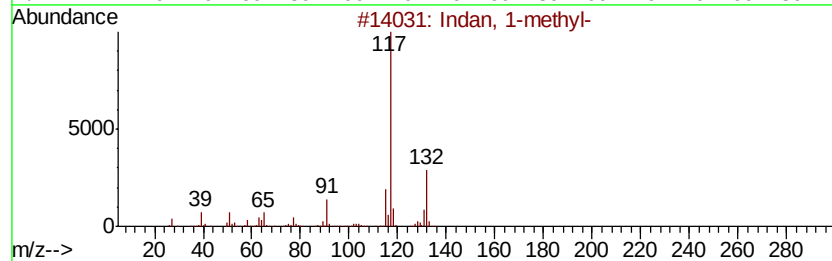
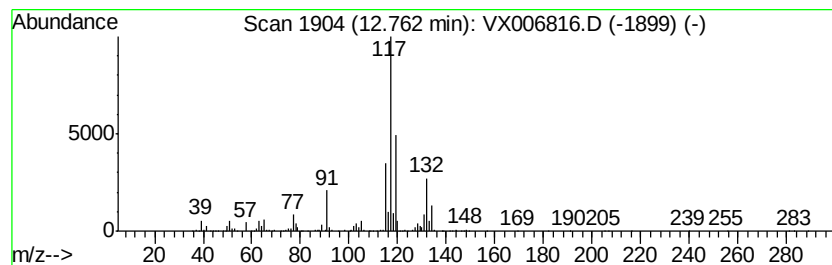
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	13.07 ug/l	702437	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-		132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	87
3	1-Phenyl-1-butene		132	C10H12	000824-90-8	81
4	Benzene, 2-butenyl-		132	C10H12	001560-06-1	74
5	Benzene, (2-methyl-1-propenyl)-		132	C10H12	000768-49-0	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-1

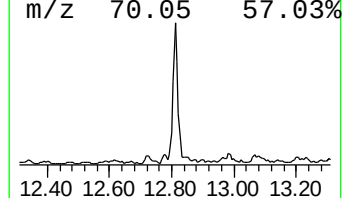
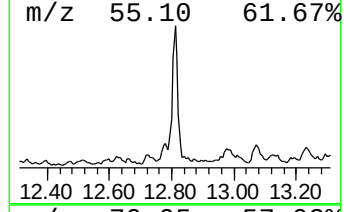
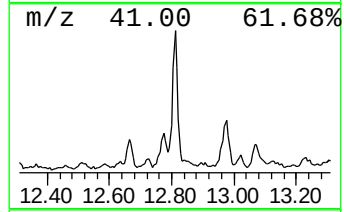
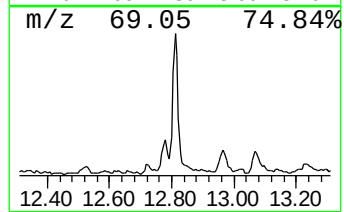
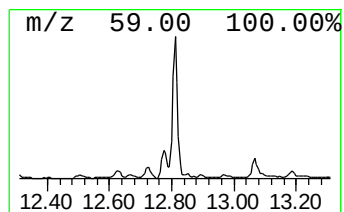
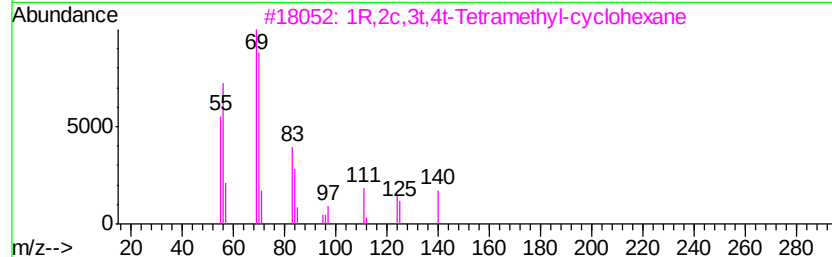
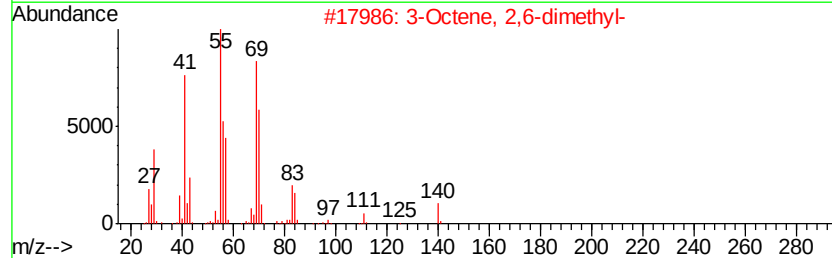
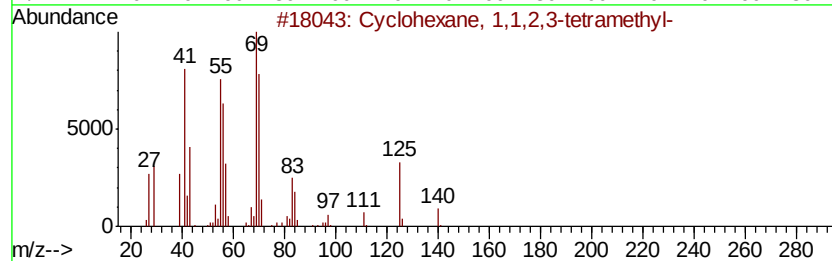
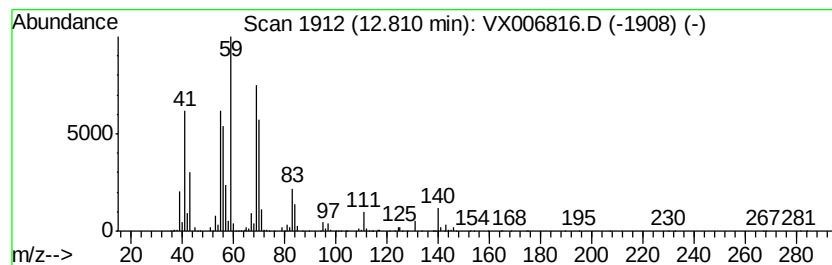
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1,1,2,3-tetram... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.81	7.97 ug/l	428649	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	93
2		3-Octene, 2,6-dimethyl-	140	C10H20	006874-28-8	62
3		1R,2c,3t,4t-Tetramethyl-cyclohexane	140	C10H20	1000144-07-3	62
4		3-Decene	140	C10H20	019398-37-9	58
5		4-Octene, 2,6-dimethyl-, [S-(Z)]-	140	C10H20	062960-77-4	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TW-1

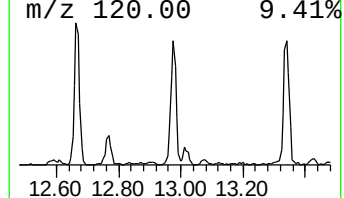
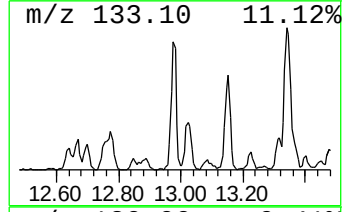
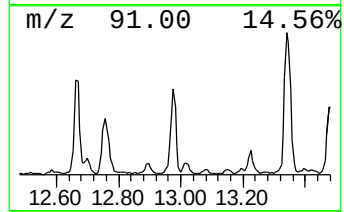
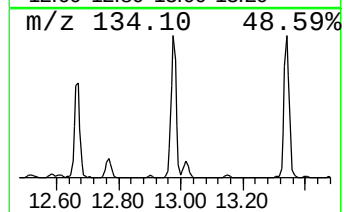
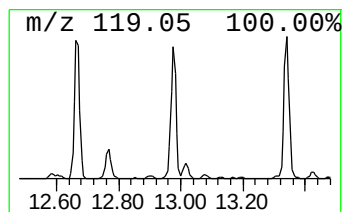
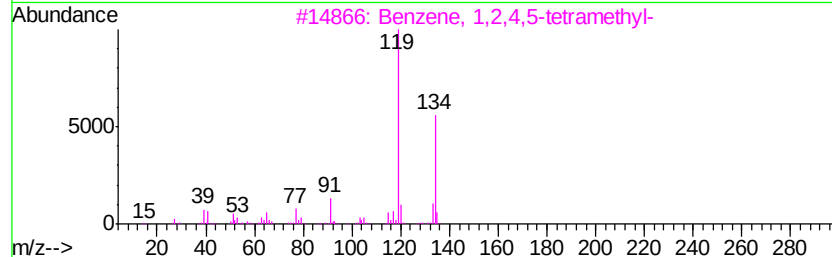
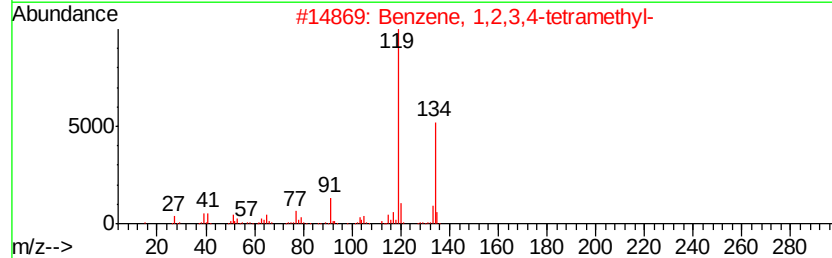
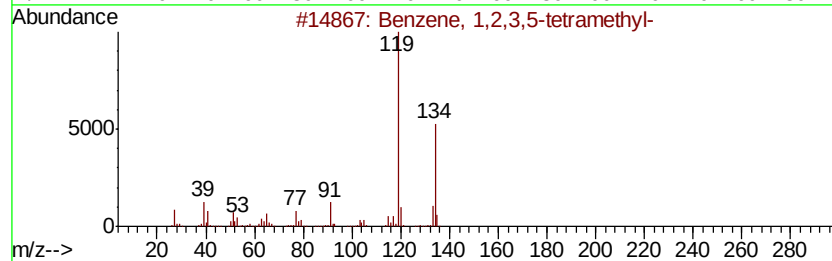
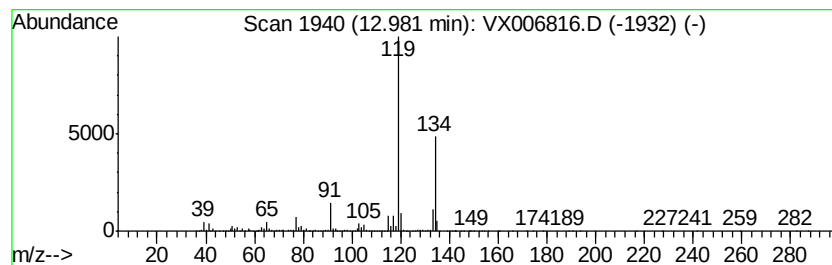
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	17.17 ug/l	922671	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TW-1

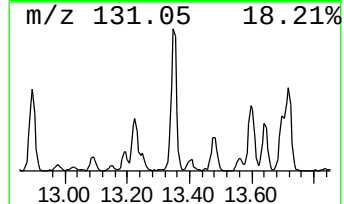
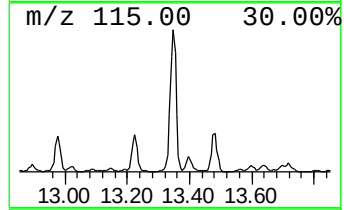
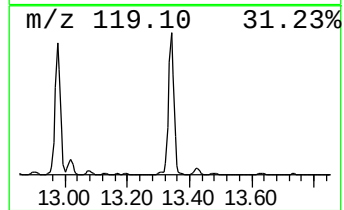
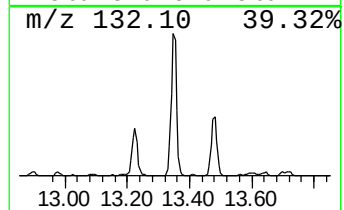
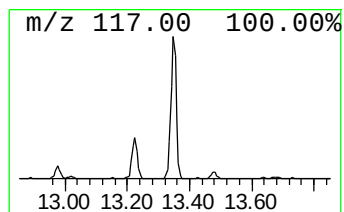
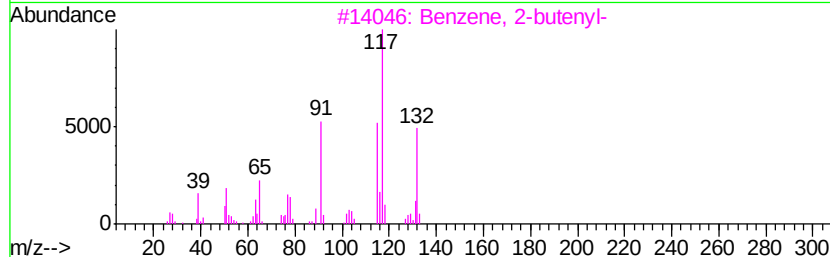
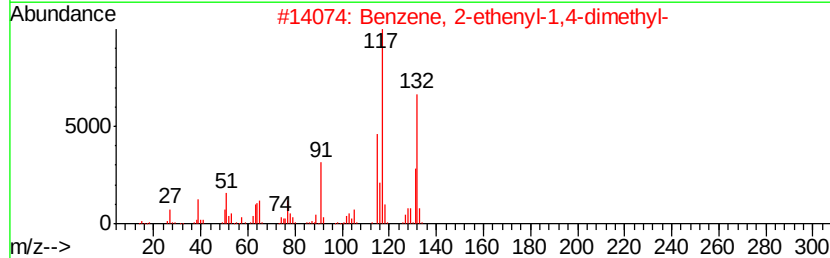
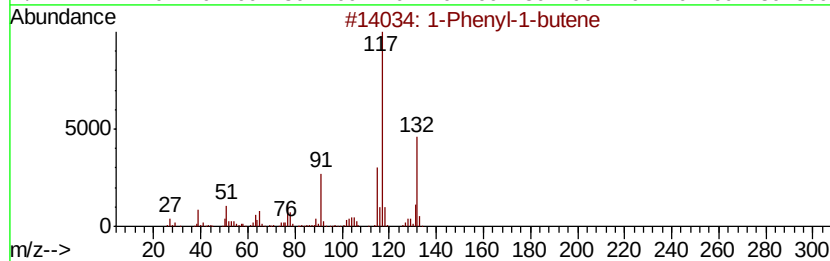
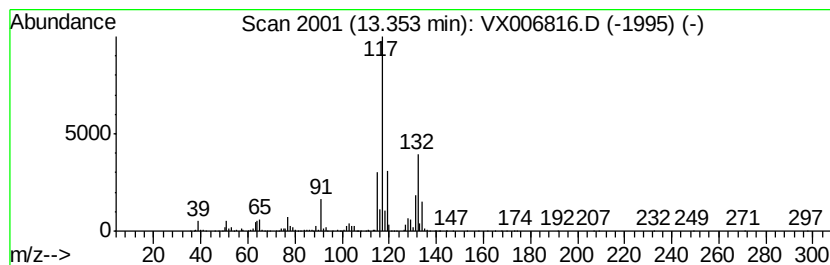
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	31.28 ug/l	1681590	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	84
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	78
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

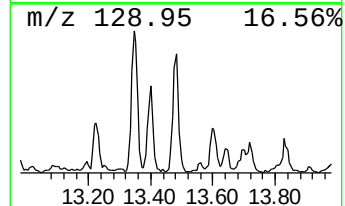
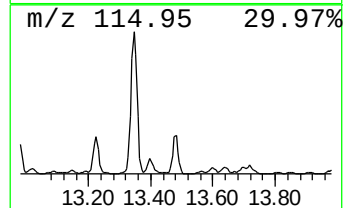
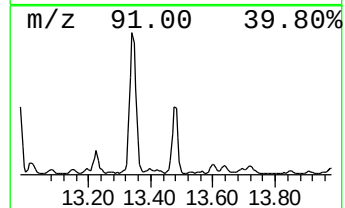
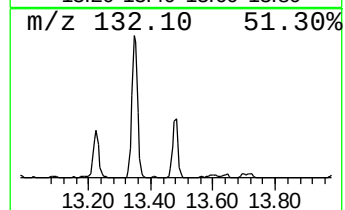
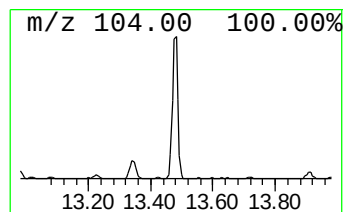
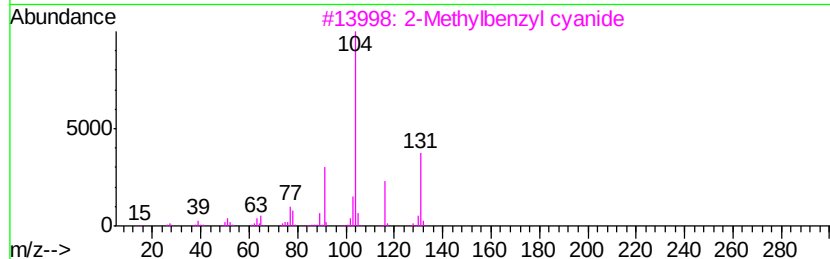
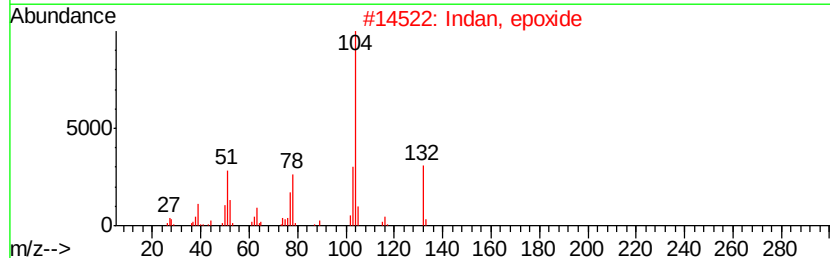
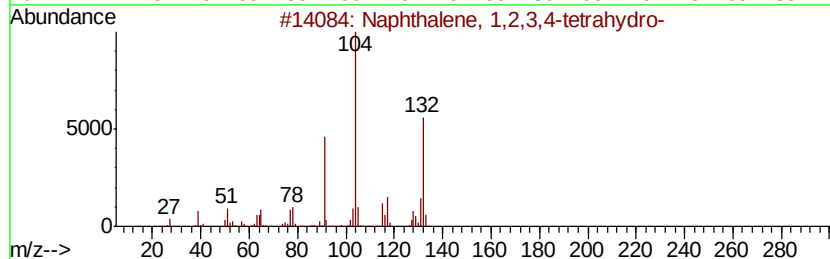
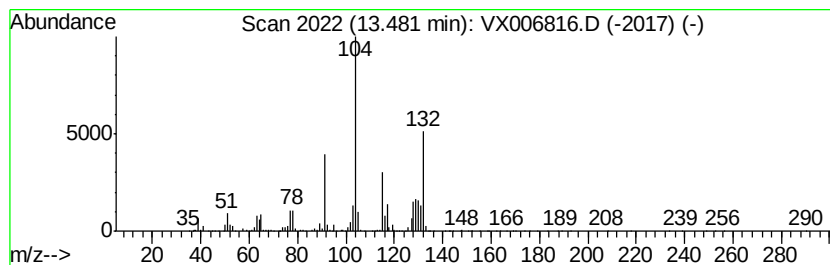
Instrument :
MSVOA_X
ClientSampleId :
TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	7.31 ug/l	392714	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 90
2		Indan, epoxide	132 C9H8O	000768-22-9 49
3		2-Methylbenzyl cyanide	131 C9H9N	022364-68-7 43
4		N-Ethylbenzimidoyl bromide	211 C9H10BrN	041182-87-0 43
5		2-Furanone, 4-phenyltetrahydro	162 C10H10O2	1000197-38-4 35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

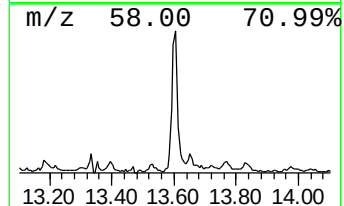
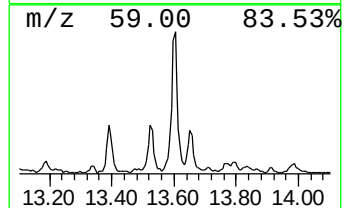
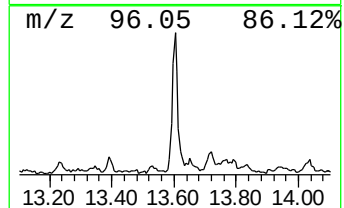
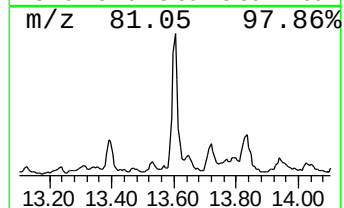
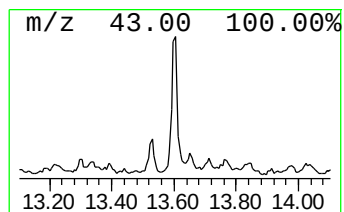
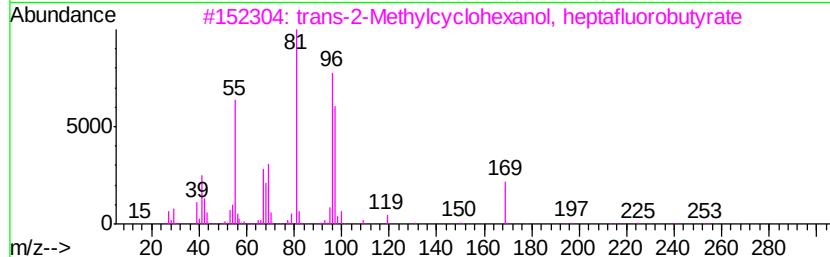
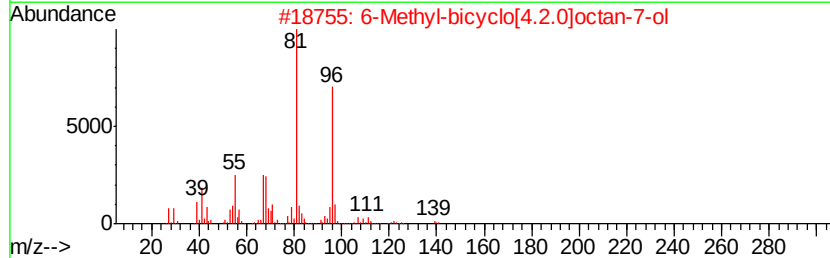
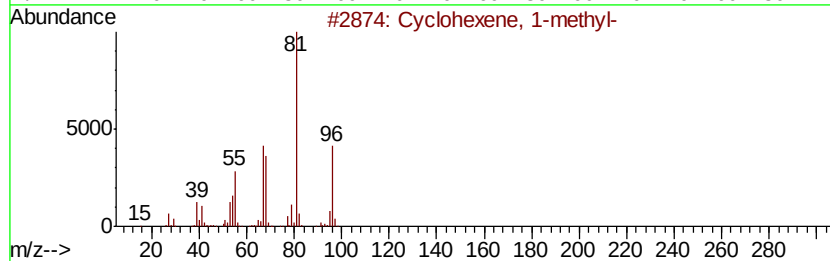
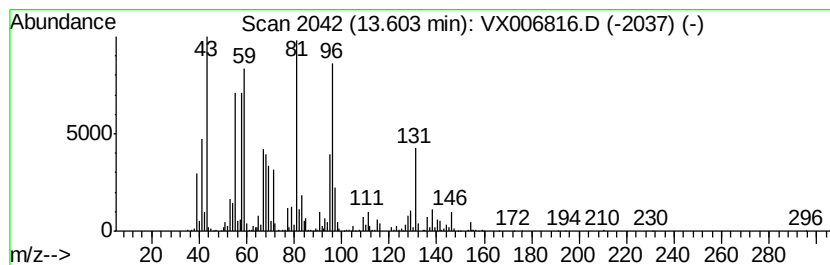
Instrument :
MSVOA_X
ClientSampleId :
TW-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 9 Cyclohexene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	10.21 ug/l	548654	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexene, 1-methyl-	96 C7H12	000591-49-1 55
2		6-Methyl-bicyclo[4.2.0]octan-7-ol	140 C9H16O	1000193-87-3 43
3		trans-2-Methylcyclohexanol, hept...	310 C11H13F7O2	1000364-13-5 35
4		Cyclohexene, 3-methyl-	96 C7H12	000591-48-0 35
5		cis-2-Methylcyclohexanol, heptaf...	310 C11H13F7O2	1000364-12-5 35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-1

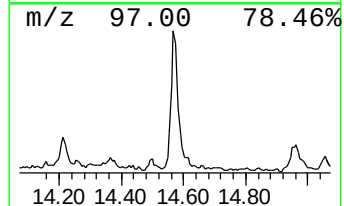
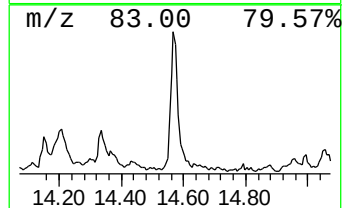
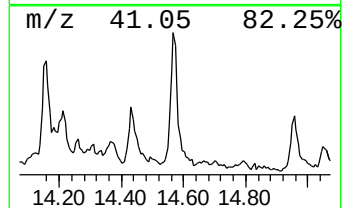
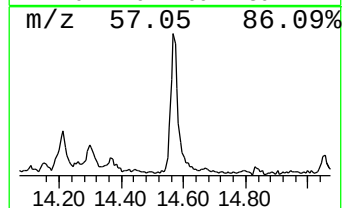
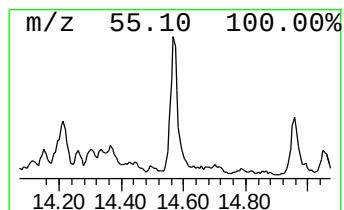
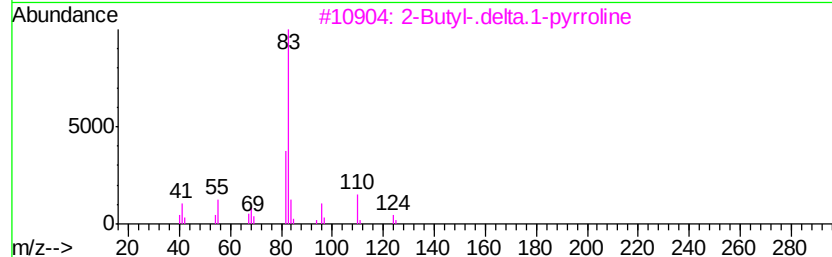
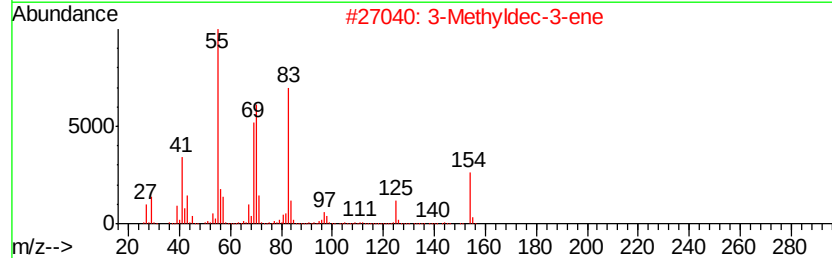
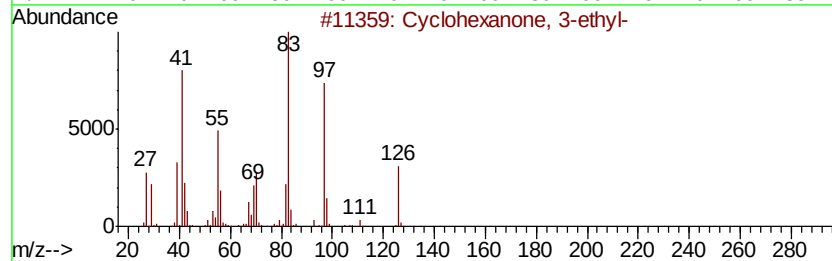
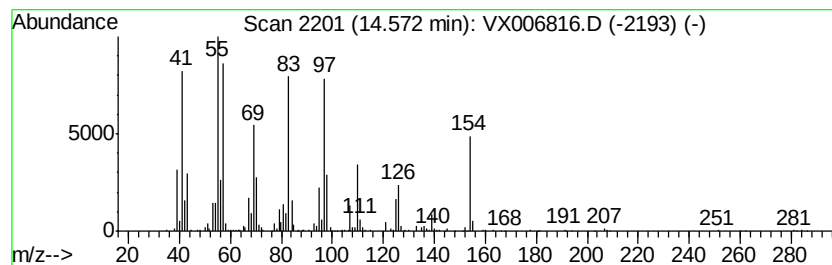
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 10 unknown14.57 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	8.69 ug/l	466858	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 3-ethyl-	126	C8H14O	022461-89-8	43
2		3-Methyldec-3-ene	154	C11H22	036969-75-2	43
3		2-Butyl-.delta.1-pyrroline	125	C8H15N	064319-86-4	42
4		2,3-Dimethyl-3-heptene	126	C9H18	1000113-49-3	42
5		trans-7-Oxabicyclo[4.3.0]nonane	126	C8H14O	027345-70-6	30



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX122618\
Data File : VX006816.D
Acq On : 26 Dec 2018 14:16
Operator : JC/MD
Sample : J6549-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.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, 1,2,3-tr...	12.14	9.9	ug/l	530371	4	12.08	2687640	50.0
Benzene, 1-etheny...	12.30	39.1	ug/l	2101090	4	12.08	2687640	50.0
o-Cymene	12.66	14.4	ug/l	774265	4	12.08	2687640	50.0
Indan, 1-methyl-	12.76	13.1	ug/l	702437	4	12.08	2687640	50.0
Cyclohexane, 1,1,...	12.81	8.0	ug/l	428649	4	12.08	2687640	50.0
Benzene, 1,2,3,5-...	12.98	17.2	ug/l	922671	4	12.08	2687640	50.0
1-Phenyl-1-butene	13.35	31.3	ug/l	1681590	4	12.08	2687640	50.0
Naphthalene, 1,2,...	13.48	7.3	ug/l	392714	4	12.08	2687640	50.0
Cyclohexene, 1-me...	13.60	10.2	ug/l	548654	4	12.08	2687640	50.0
unknown14.57	14.57	8.7	ug/l	466858	4	12.08	2687640	50.0