

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX063018\
 Data File : VX003116.D
 Acq On : 01 Jul 2018 09:05
 Operator : JC/MD
 Sample : J3799-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-6

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\82X062518W.M
 Title : SW846 8260

Signal : TIC: VX003115.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.075	76	80	94	rBV	1961235	2328314	100.00%	13.405%
2	1.276	108	113	127	rBV	162540	395822	17.00%	2.279%
3	5.507	793	807	822	rBV	181703	527256	22.65%	3.036%
4	5.666	822	833	849	rVB	272534	768779	33.02%	4.426%
5	6.074	889	900	922	rBV	190862	504973	21.69%	2.907%
6	6.867	1020	1030	1044	rBV	390775	909737	39.07%	5.238%
7	8.720	1327	1334	1349	rVB	929972	1457331	62.59%	8.390%
8	10.116	1551	1563	1570	rBV	837781	1105583	47.48%	6.365%
9	10.360	1595	1603	1610	rVB2	22187	43371	1.86%	0.250%
10	11.018	1706	1711	1716	rBV	63952	89441	3.84%	0.515%
11	11.140	1725	1731	1737	rBV	880330	1093433	46.96%	6.295%
12	11.280	1751	1754	1762	rVB3	17391	29370	1.26%	0.169%
13	11.360	1762	1767	1772	rBV	61873	87616	3.76%	0.504%
14	11.671	1814	1818	1821	rBV	25534	36305	1.56%	0.209%
15	11.774	1826	1835	1838	rVV	29318	49731	2.14%	0.286%
16	11.811	1838	1841	1847	rVB	22469	33779	1.45%	0.194%
17	11.921	1853	1859	1861	rBV	46238	66504	2.86%	0.383%
18	11.945	1861	1863	1871	rVV	81691	119288	5.12%	0.687%
19	12.079	1880	1885	1891	rBV	1014707	1281788	55.05%	7.380%
20	12.238	1907	1911	1915	rVB	67440	82965	3.56%	0.478%
21	12.305	1917	1922	1926	rBV	104352	123333	5.30%	0.710%
22	12.372	1930	1933	1941	rVB4	36882	82865	3.56%	0.477%
23	12.463	1941	1948	1953	rBV	164549	217637	9.35%	1.253%
24	12.524	1954	1958	1964	rVB	57475	77325	3.32%	0.445%
25	12.670	1975	1982	1985	rBV	98541	124762	5.36%	0.718%
26	12.707	1985	1988	1993	rVV4	28470	56560	2.43%	0.326%
27	12.762	1993	1997	2004	rVV4	188735	352415	15.14%	2.029%
28	12.817	2004	2006	2009	rVV	29629	31759	1.36%	0.183%
29	12.896	2013	2019	2024	rVB2	104892	178504	7.67%	1.028%
30	12.981	2025	2033	2037	rBV	325166	483132	20.75%	2.782%
31	13.024	2037	2040	2046	rVB2	101575	142992	6.14%	0.823%
32	13.091	2046	2051	2056	rVB2	80853	128193	5.51%	0.738%
33	13.152	2056	2061	2065	rBV	47087	65868	2.83%	0.379%
34	13.201	2065	2069	2072	rBV	149804	186364	8.00%	1.073%

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35	13.256	2075	2078	2083	rVV	41001	51468	2.21%	0.296%
36	13.311	2083	2087	2090	rVV2	80328	110223	4.73%	0.635%
37	13.353	2090	2094	2098	rVV2	150911	229139	9.84%	1.319%
38	13.402	2098	2102	2105	rVV4	29465	58135	2.50%	0.335%
39	13.481	2109	2115	2121	rVB4	39224	85437	3.67%	0.492%
40	13.560	2126	2128	2131	rBV3	16894	23541	1.01%	0.136%
41	13.603	2131	2135	2138	rBV2	46879	52409	2.25%	0.302%
42	13.640	2138	2141	2145	rBV3	112894	165690	7.12%	0.954%
43	13.701	2145	2151	2154	rBV2	109476	153007	6.57%	0.881%
44	13.810	2165	2169	2173	rBV	80119	101796	4.37%	0.586%
45	13.859	2173	2177	2181	rVB3	41454	64483	2.77%	0.371%
46	13.914	2181	2186	2190	rBV	180138	232496	9.99%	1.339%
47	13.987	2191	2198	2202	rBV	215273	298802	12.83%	1.720%
48	14.036	2202	2206	2210	rBV2	51641	76682	3.29%	0.441%
49	14.134	2218	2222	2225	rBV3	66389	96584	4.15%	0.556%
50	14.164	2225	2227	2231	rVB3	61366	68246	2.93%	0.393%
51	14.274	2241	2245	2256	rVB2	94336	205975	8.85%	1.186%
52	14.377	2257	2262	2267	rVB	106012	151672	6.51%	0.873%
53	14.432	2267	2271	2276	rBV2	86841	126491	5.43%	0.728%
54	14.481	2276	2279	2285	rVB	36884	51176	2.20%	0.295%
55	14.548	2285	2290	2294	rBV2	112630	171028	7.35%	0.985%
56	14.640	2302	2305	2308	rBV2	18990	26604	1.14%	0.153%
57	14.676	2308	2311	2313	rVV3	63297	80089	3.44%	0.461%
58	14.737	2317	2321	2324	rVV2	86146	120687	5.18%	0.695%
59	14.786	2325	2329	2334	rVV	45068	72047	3.09%	0.415%
60	14.853	2335	2340	2344	rVV	77677	146202	6.28%	0.842%
61	14.908	2347	2349	2353	rVB2	46206	52680	2.26%	0.303%
62	14.950	2353	2356	2359	rBV2	28081	52274	2.25%	0.301%
63	14.981	2359	2361	2366	rVB4	32735	43098	1.85%	0.248%
64	15.042	2367	2371	2375	rVB7	17948	32862	1.41%	0.189%
65	15.121	2381	2384	2388	rVB5	23763	31085	1.34%	0.179%
66	15.249	2399	2405	2413	rBV2	78540	165157	7.09%	0.951%
67	15.396	2425	2429	2434	rBV2	46479	81704	3.51%	0.470%
68	15.481	2440	2443	2447	rVB2	29392	38194	1.64%	0.220%
69	15.554	2451	2455	2459	rBV3	26065	38091	1.64%	0.219%
70	15.597	2460	2462	2470	rVB4	22906	45908	1.97%	0.264%
71	15.688	2471	2477	2482	rBV3	56490	110491	4.75%	0.636%

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72	15.810	2493	2497	2503	rBV2	29298	60747	2.61%	0.350%
73	15.895	2507	2511	2513	rVV3	42401	76462	3.28%	0.440%
74	15.926	2513	2516	2525	rVB2	54031	98177	4.22%	0.565%
75	16.072	2536	2540	2553	rVB	35101	77735	3.34%	0.448%
76	16.261	2568	2571	2580	rVB7	14996	31469	1.35%	0.181%
77	16.420	2594	2597	2604	rVB2	14869	25628	1.10%	0.148%
78	16.566	2618	2621	2628	rVB7	12252	24054	1.03%	0.138%

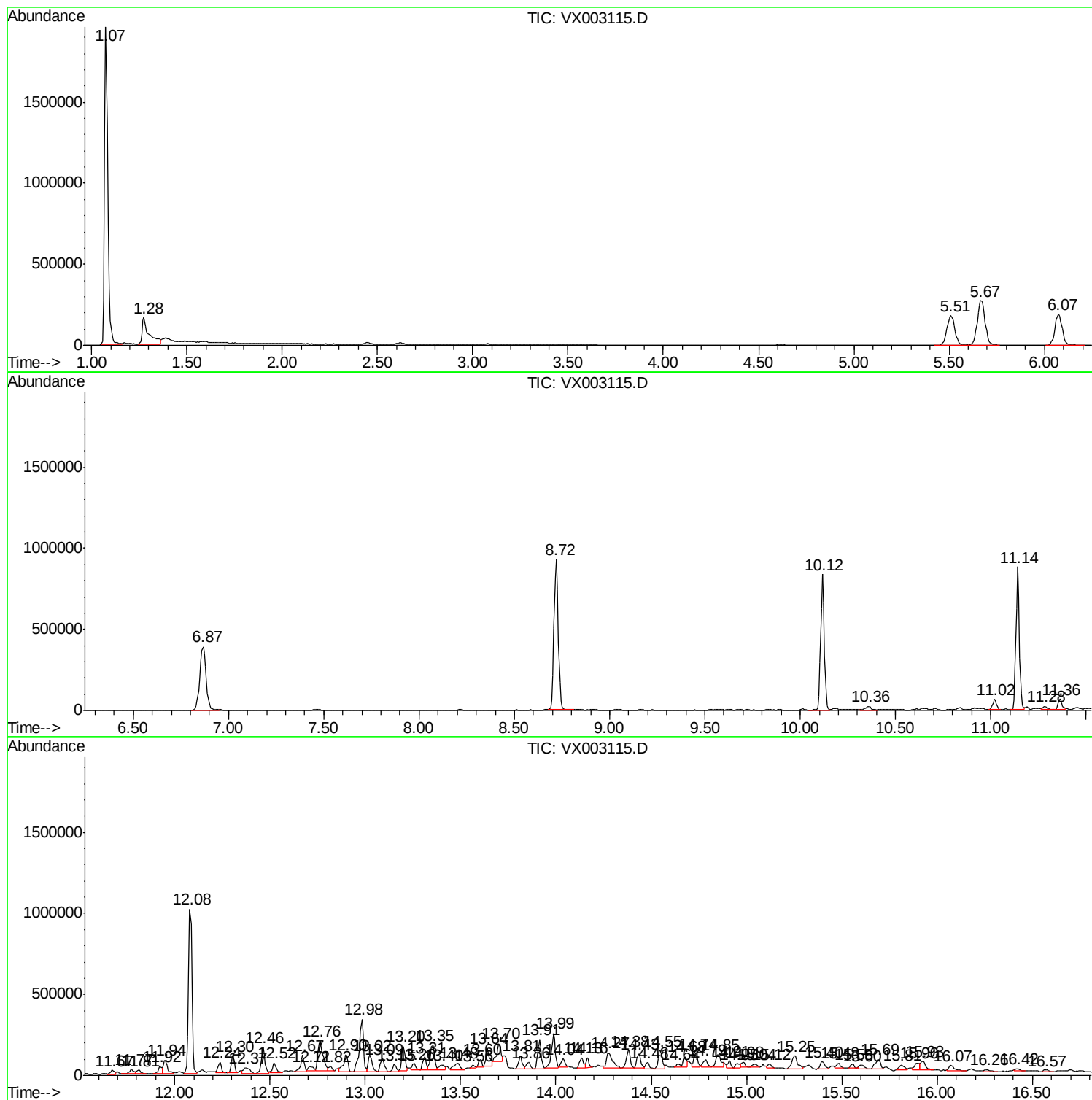
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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



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Instrument :
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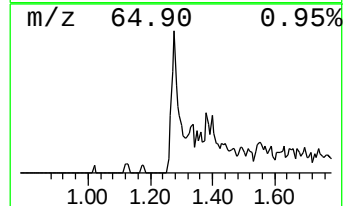
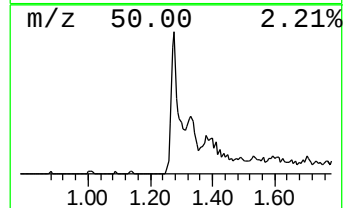
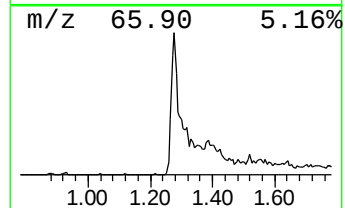
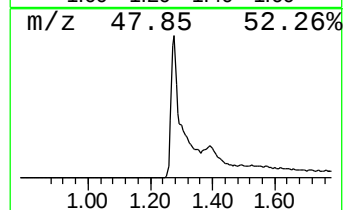
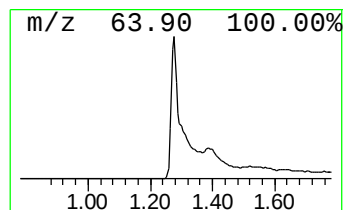
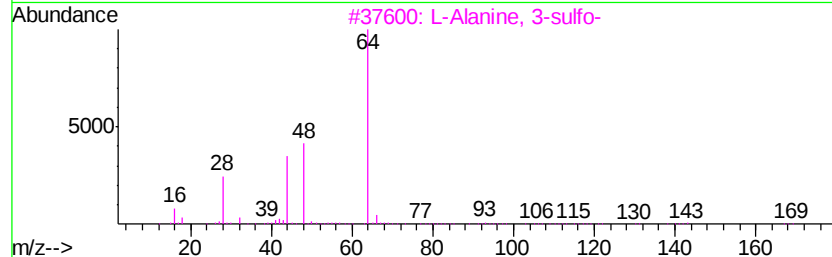
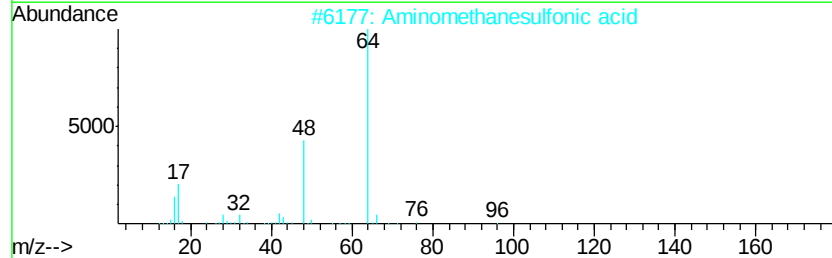
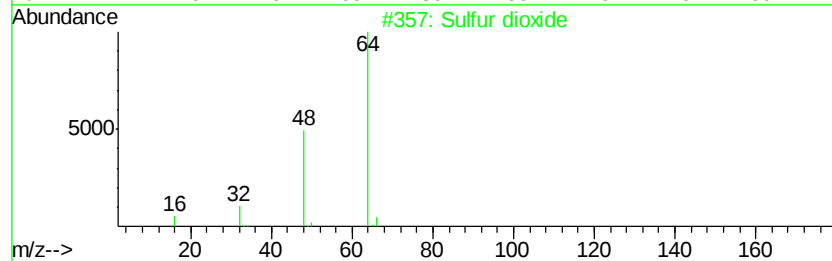
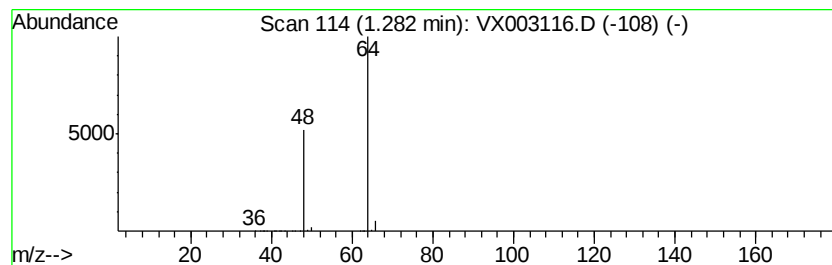
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 2 Sulfur dioxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.28	25.74 ug/l	395822	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfur dioxide	64	O2S	007446-09-5	83
2		Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3		L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4		Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4
5		Ethyl Chloride	64	C2H5Cl	000075-00-3	3



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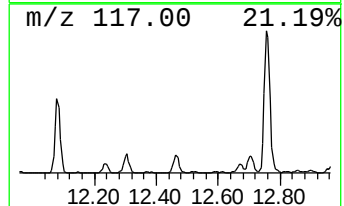
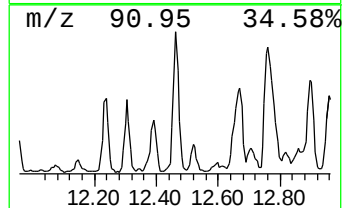
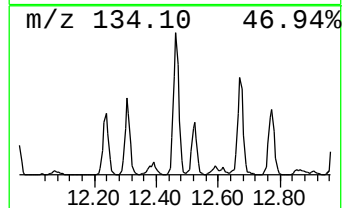
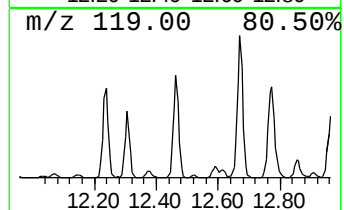
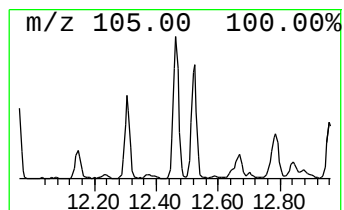
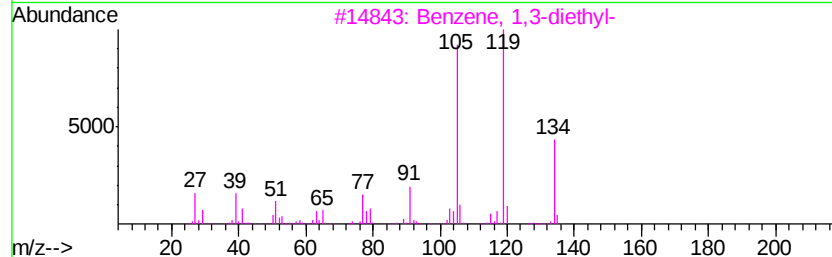
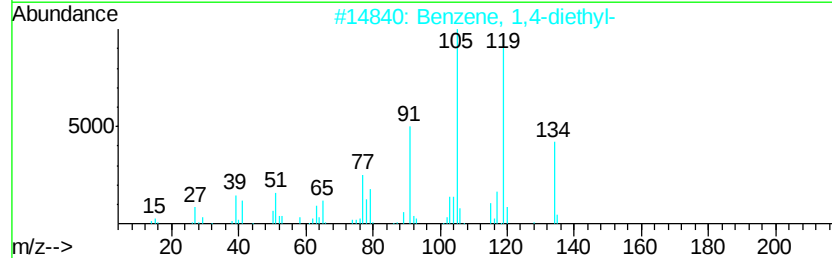
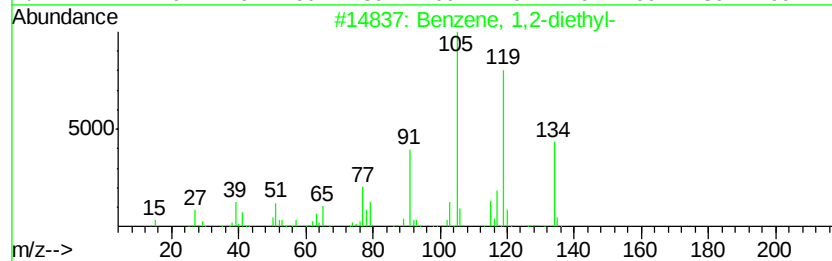
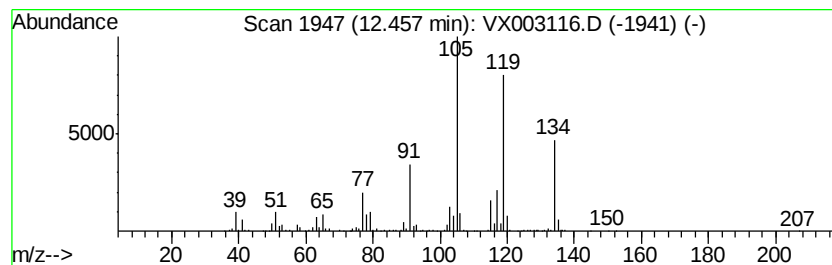
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	8.49 ug/l	217637	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 96
2		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 93
3		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 87
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 74
5		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 62



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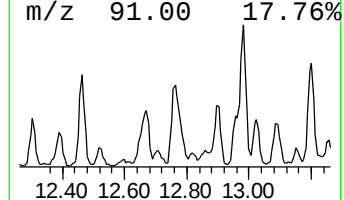
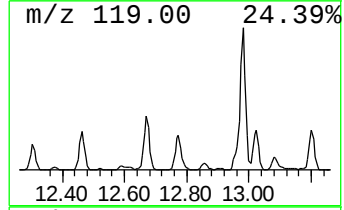
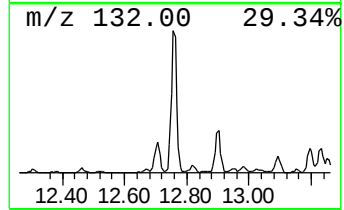
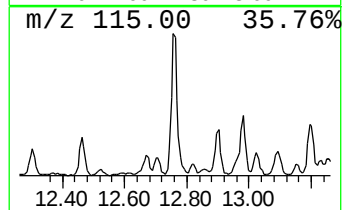
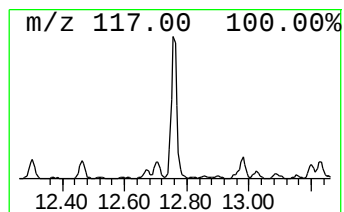
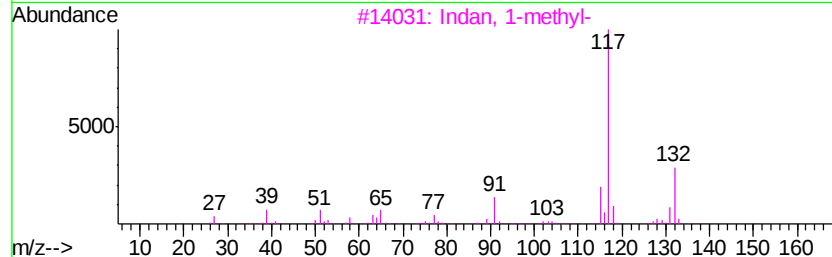
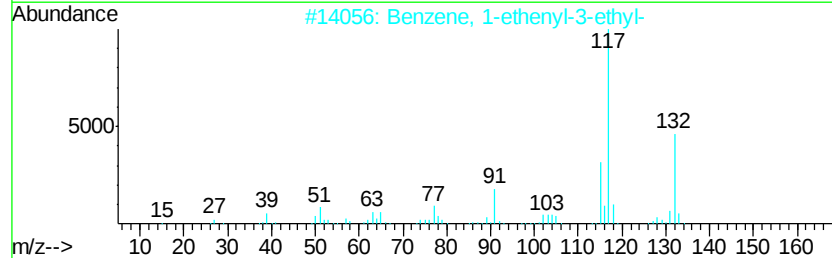
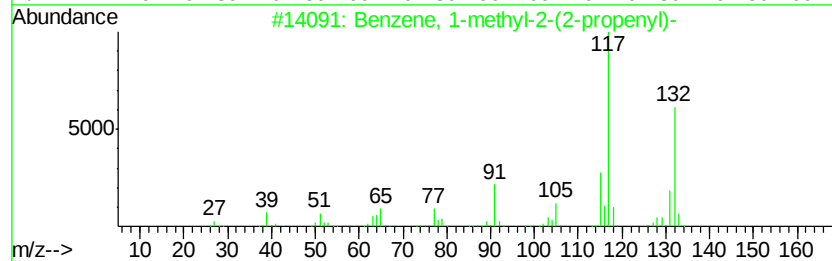
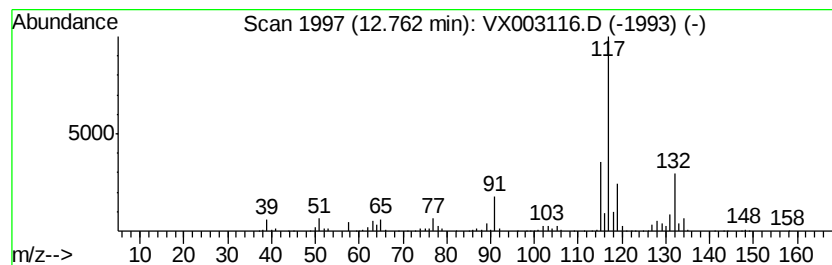
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-2-(2-prop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	13.75 ug/l	352415	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
3		Indan, 1-methyl-	132	C10H12	000767-58-8	81
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	80



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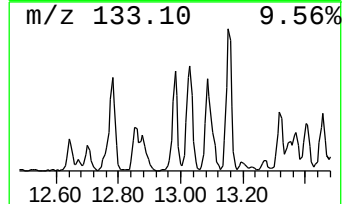
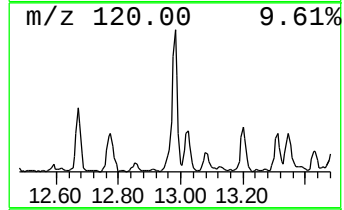
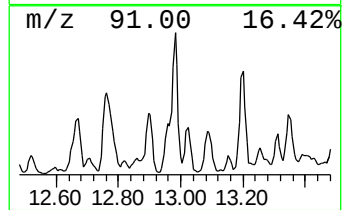
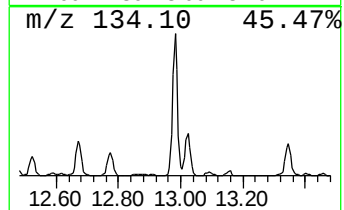
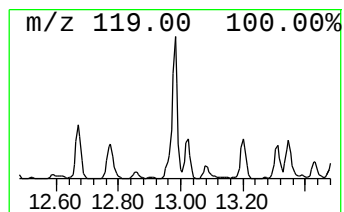
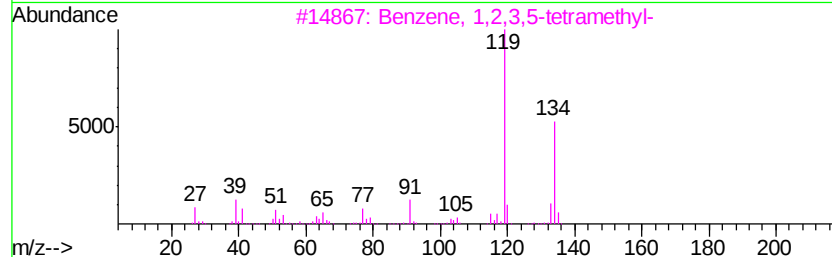
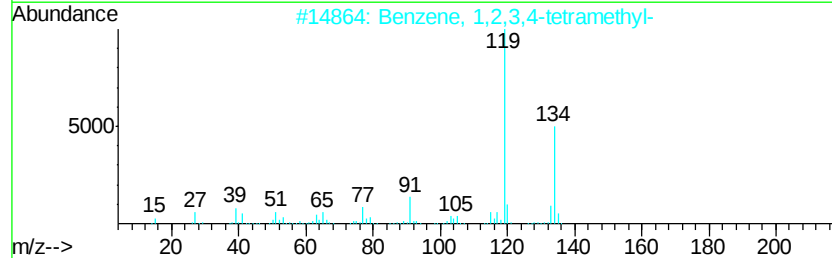
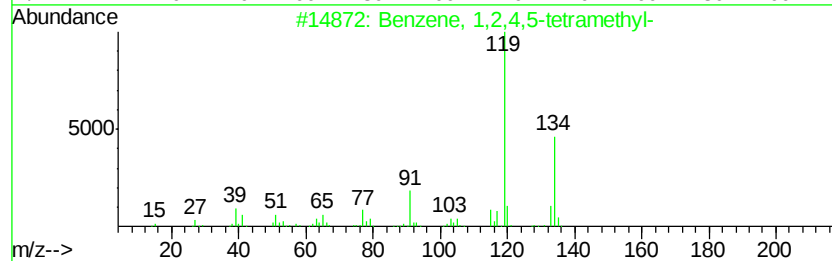
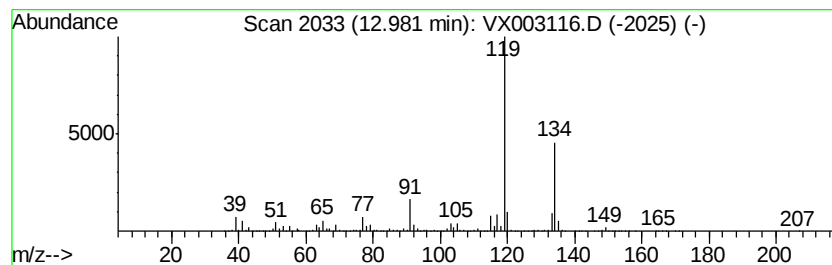
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	18.85 ug/l	483132	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX063018\
Data File : VX003116.D
Acq On : 01 Jul 2018 09:05
Operator : JC/MD
Sample : J3799-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

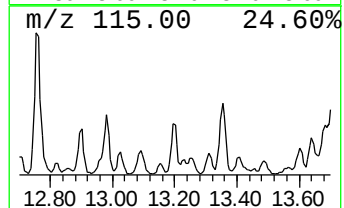
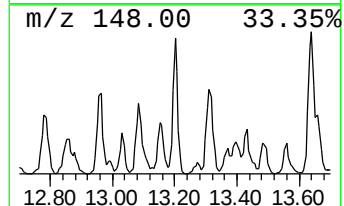
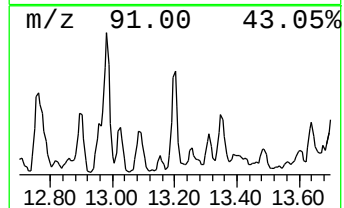
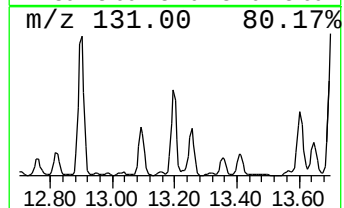
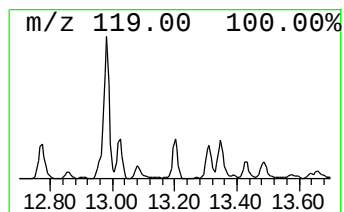
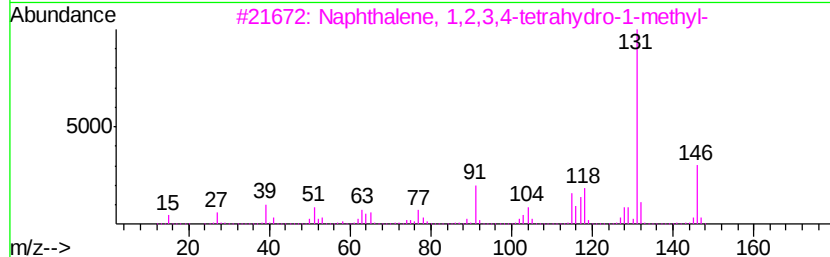
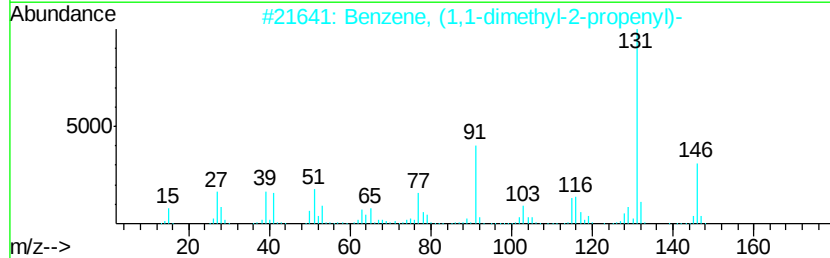
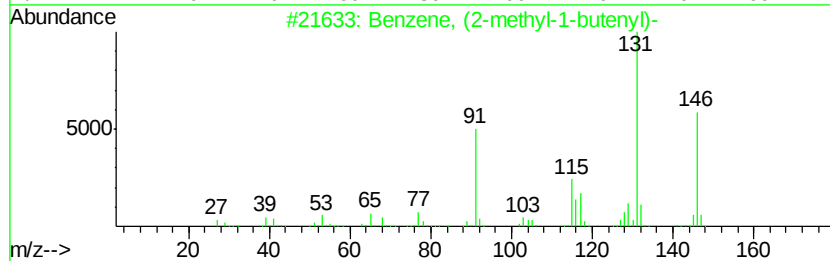
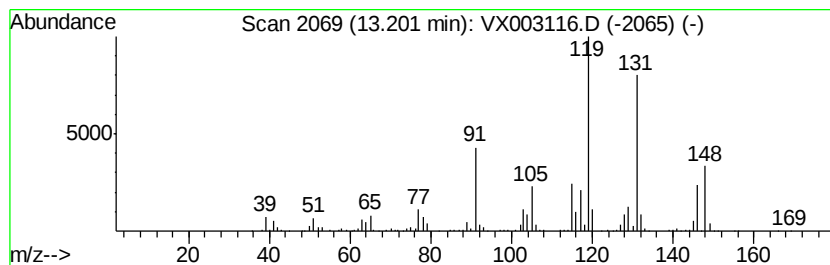
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	7.27 ug/l	186364	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	50
2		Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	46
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	46
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	46
5		1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX063018\
Data File : VX003116.D
Acq On : 01 Jul 2018 09:05
Operator : JC/MD
Sample : J3799-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

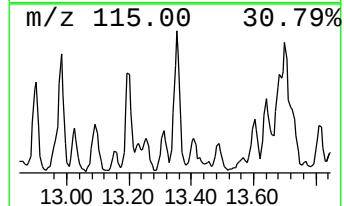
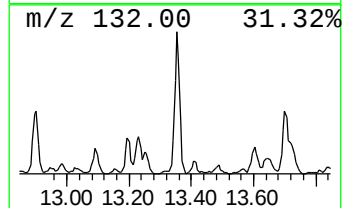
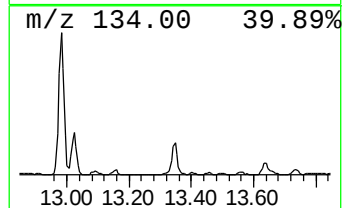
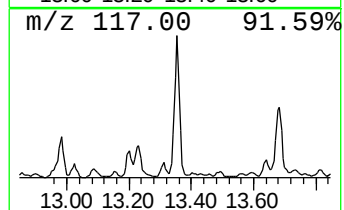
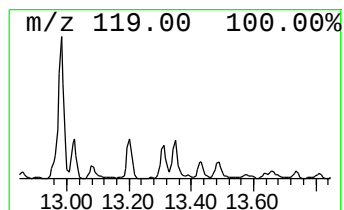
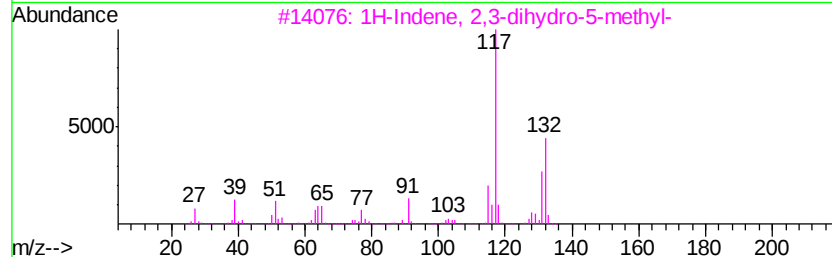
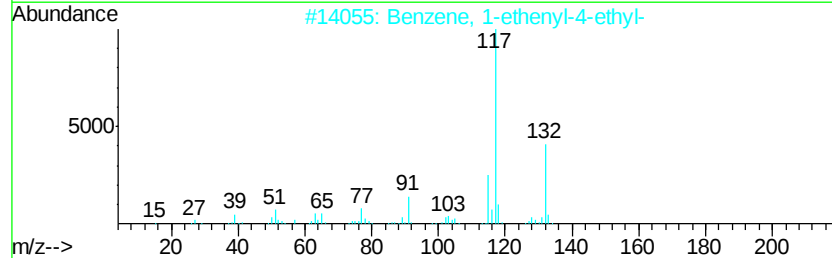
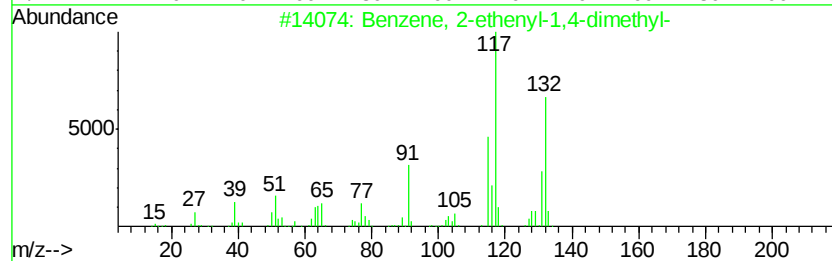
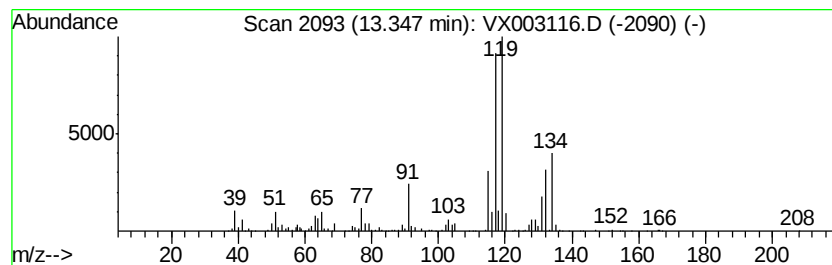
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	8.94 ug/l	229139	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	91
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX063018\
Data File : VX003116.D
Acq On : 01 Jul 2018 09:05
Operator : JC/MD
Sample : J3799-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-6

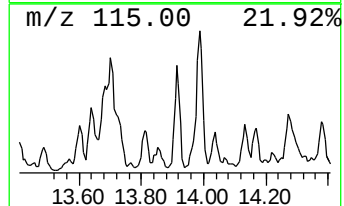
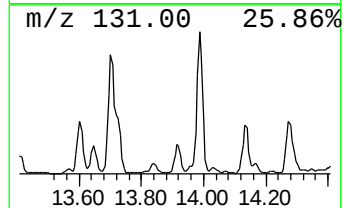
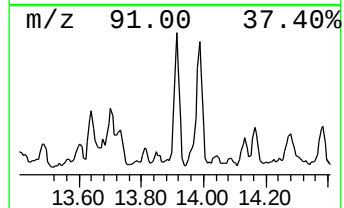
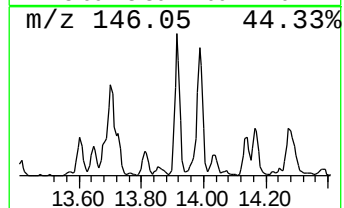
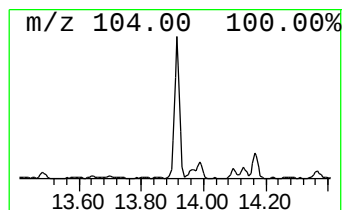
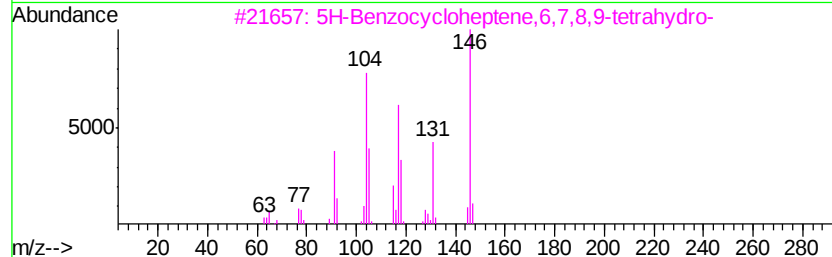
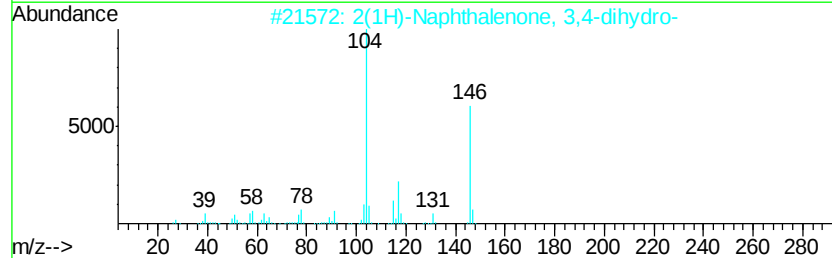
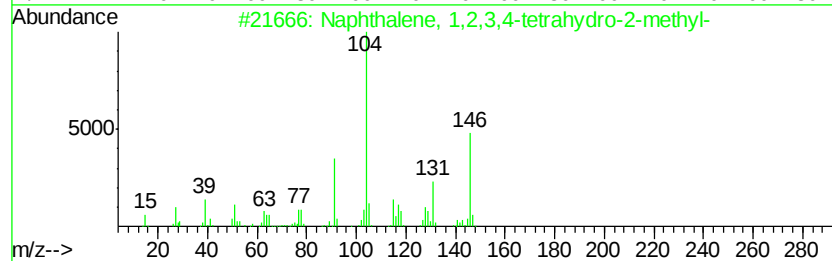
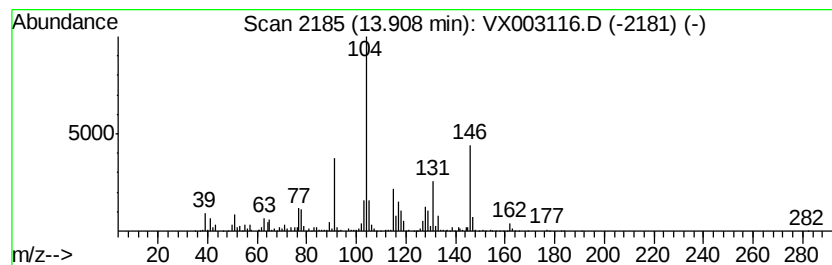
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	9.07 ug/l	232496	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	68
3		5H-Benzocycloheptene, 6,7,8,9-tet...	146	C11H14	001075-16-7	59
4		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	58
5		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX063018\
Data File : VX003116.D
Acq On : 01 Jul 2018 09:05
Operator : JC/MD
Sample : J3799-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

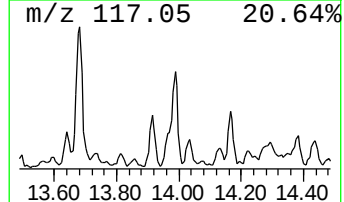
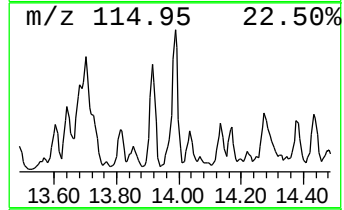
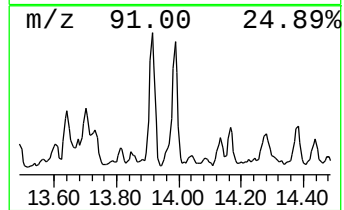
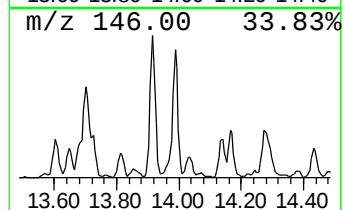
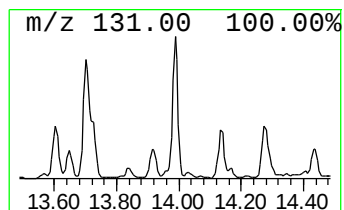
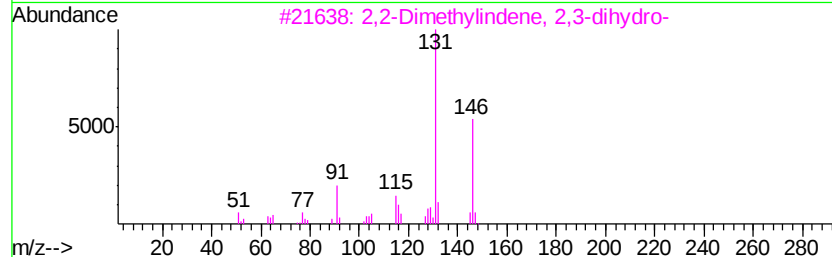
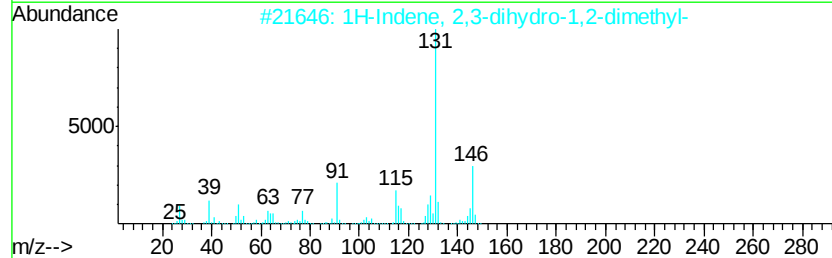
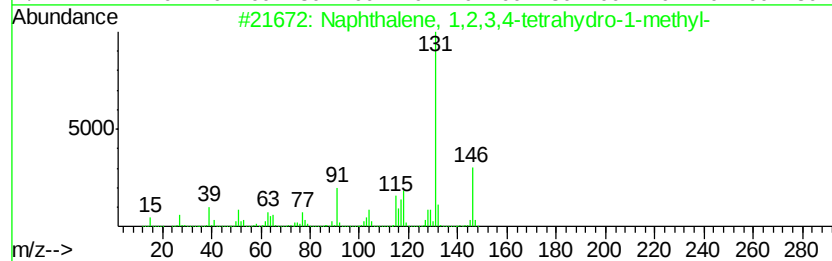
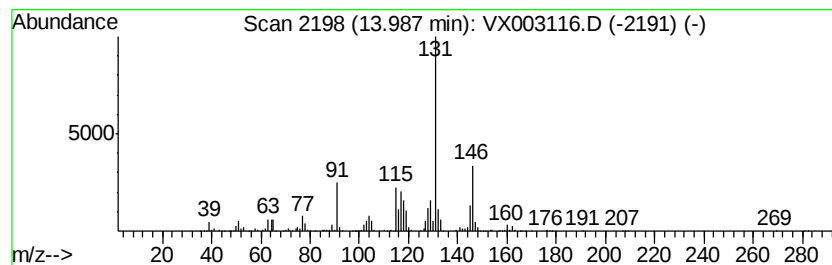
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	11.66 ug/l	298802	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	89
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX063018\
Data File : VX003116.D
Acq On : 01 Jul 2018 09:05
Operator : JC/MD
Sample : J3799-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

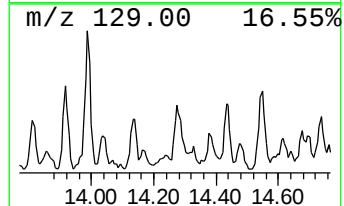
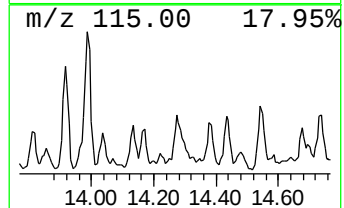
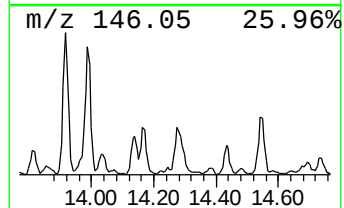
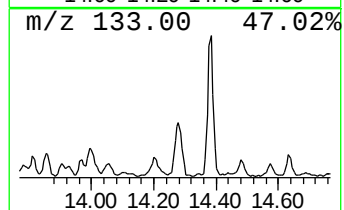
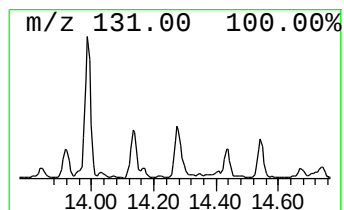
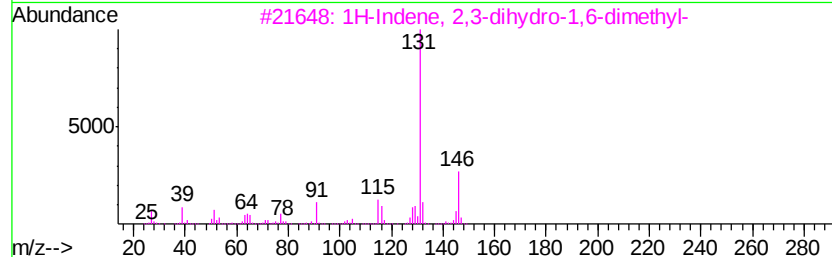
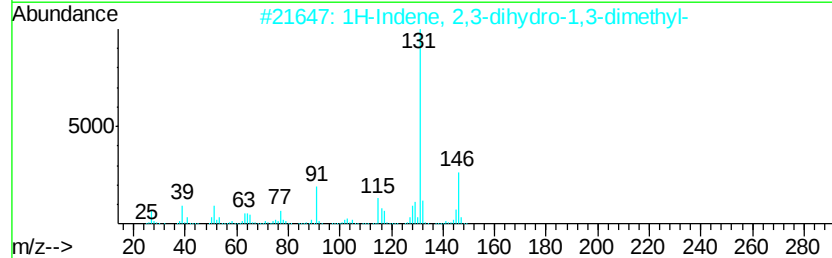
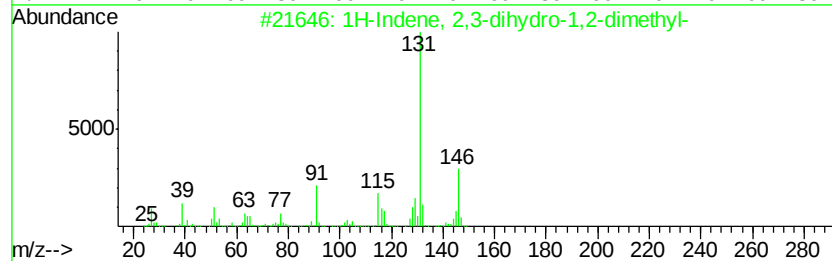
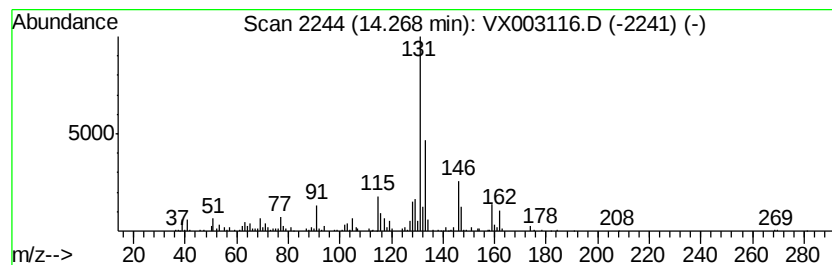
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	8.03 ug/l	205975	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	86
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
4		Benzene, (3-methyl-2-butenvl)-	146	C11H14	004489-84-3	64
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX063018\
Data File : VX003116.D
Acq On : 01 Jul 2018 09:05
Operator : JC/MD
Sample : J3799-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.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
Sulfur dioxide	1.28	25.7	ug/l	395822	1	5.67	768779	50.0
Benzene, 1,2-diet...	12.46	8.5	ug/l	217637	4	12.08	1281790	50.0
Benzene, 1-methyl...	12.76	13.8	ug/l	352415	4	12.08	1281790	50.0
Benzene, 1,2,4,5-...	12.98	18.9	ug/l	483132	4	12.08	1281790	50.0
Benzene, (2-methy...	13.20	7.3	ug/l	186364	4	12.08	1281790	50.0
Benzene, 2-etheny...	13.35	8.9	ug/l	229139	4	12.08	1281790	50.0
Naphthalene, 1,2,...	13.91	9.1	ug/l	232496	4	12.08	1281790	50.0
Naphthalene, 1,2,...	13.99	11.7	ug/l	298802	4	12.08	1281790	50.0
1H-Indene, 2,3-di...	14.27	8.0	ug/l	205975	4	12.08	1281790	50.0