

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060618\
 Data File : VX002364.D
 Acq On : 06 Jun 2018 20:22
 Operator : JC/MD
 Sample : J3312-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 W-04

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\82X060418W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.075	76	80	94	rBV	1409387	1735441	100.00%	10.811%
2	1.270	108	112	116	rBV	45203	62275	3.59%	0.388%
3	1.325	116	121	139	rVV	79261	418713	24.13%	2.608%
4	2.617	328	333	340	rBV	13368	26115	1.50%	0.163%
5	5.507	796	807	821	rBV	145443	416780	24.02%	2.596%
6	5.666	823	833	851	rVB	234954	659554	38.00%	4.109%
7	6.074	889	900	918	rBV	167496	438345	25.26%	2.731%
8	6.867	1019	1030	1045	rBV	339059	797627	45.96%	4.969%
9	8.720	1328	1334	1343	rVB	836374	1277130	73.59%	7.956%
10	9.043	1381	1387	1392	rBV2	16537	27010	1.56%	0.168%
11	9.573	1468	1474	1479	rBV4	9213	18040	1.04%	0.112%
12	9.683	1487	1492	1495	rBV	12943	19017	1.10%	0.118%
13	10.116	1557	1563	1570	rBV	701680	938182	54.06%	5.844%
14	10.189	1570	1575	1579	rVB2	14067	21755	1.25%	0.136%
15	10.342	1595	1600	1609	rBV8	12750	34971	2.02%	0.218%
16	10.652	1636	1651	1652	rBV5	15618	34482	1.99%	0.215%
17	10.841	1679	1682	1687	rVB3	13944	18325	1.06%	0.114%
18	10.915	1688	1694	1703	rBV5	13825	29986	1.73%	0.187%
19	11.024	1707	1712	1716	rBV	117474	154890	8.93%	0.965%
20	11.140	1726	1731	1736	rBV	693315	889181	51.24%	5.539%
21	11.189	1736	1739	1742	rVB	15216	18683	1.08%	0.116%
22	11.280	1751	1754	1760	rVB3	15778	22043	1.27%	0.137%
23	11.366	1762	1768	1776	rBV	75828	102818	5.92%	0.641%
24	11.677	1813	1819	1826	rBV5	14715	29437	1.70%	0.183%
25	11.774	1830	1835	1839	rBV	36141	46155	2.66%	0.288%
26	11.927	1854	1860	1861	rBV	65901	90298	5.20%	0.563%
27	11.951	1861	1864	1871	rVB	135637	167879	9.67%	1.046%
28	12.085	1880	1886	1891	rBV	784395	1006546	58.00%	6.270%
29	12.238	1906	1911	1916	rVB	56286	68395	3.94%	0.426%
30	12.305	1916	1922	1927	rVV	213981	259021	14.93%	1.614%
31	12.390	1929	1936	1943	rVB2	54195	106819	6.16%	0.665%
32	12.463	1943	1948	1955	rBV	156804	201985	11.64%	1.258%
33	12.670	1972	1982	1985	rBV3	48080	85003	4.90%	0.530%
34	12.707	1985	1988	1993	rVV	144459	187405	10.80%	1.167%

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Integrator: RTE
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35	12.768	1993	1998	2004	rVV4	187195	363240	20.93%	2.263%
36	12.823	2004	2007	2011	rVB	29186	36326	2.09%	0.226%
37	12.902	2011	2020	2024	rBV	95432	157309	9.06%	0.980%
38	12.981	2025	2033	2038	rBV	305690	444258	25.60%	2.767%
39	13.030	2038	2041	2046	rVB	22045	30747	1.77%	0.192%
40	13.085	2046	2050	2057	rVB2	55035	89030	5.13%	0.555%
41	13.158	2057	2062	2065	rBV	35355	52059	3.00%	0.324%
42	13.201	2065	2069	2072	rBV	119127	149147	8.59%	0.929%
43	13.231	2072	2074	2076	rVV	45436	44357	2.56%	0.276%
44	13.256	2076	2078	2083	rVB2	37554	41215	2.37%	0.257%
45	13.310	2083	2087	2090	rBV3	19002	29796	1.72%	0.186%
46	13.353	2090	2094	2099	rVB2	256977	335705	19.34%	2.091%
47	13.408	2099	2103	2106	rBV2	28137	43460	2.50%	0.271%
48	13.487	2113	2116	2121	rVB2	22163	29426	1.70%	0.183%
49	13.573	2126	2130	2132	rVV2	42091	59405	3.42%	0.370%
50	13.603	2132	2135	2138	rVV	62582	78878	4.55%	0.491%
51	13.646	2138	2142	2145	rVV2	126279	182101	10.49%	1.134%
52	13.701	2145	2151	2153	rVV3	107780	217341	12.52%	1.354%
53	13.725	2153	2155	2164	rVB2	142166	196678	11.33%	1.225%
54	13.810	2164	2169	2173	rBV	52489	74676	4.30%	0.465%
55	13.859	2173	2177	2182	rVB4	28551	44913	2.59%	0.280%
56	13.914	2182	2186	2191	rVB	146203	186269	10.73%	1.160%
57	13.987	2191	2198	2202	rBV	173660	255964	14.75%	1.595%
58	14.036	2202	2206	2210	rBV2	41376	53259	3.07%	0.332%
59	14.134	2218	2222	2225	rBV2	47769	66382	3.83%	0.414%
60	14.170	2225	2228	2232	rVB	46403	52985	3.05%	0.330%
61	14.280	2241	2246	2255	rVB2	89572	155527	8.96%	0.969%
62	14.383	2259	2263	2267	rVB	62679	76340	4.40%	0.476%
63	14.438	2267	2272	2276	rVB2	55169	71838	4.14%	0.448%
64	14.481	2276	2279	2284	rVB	30154	37829	2.18%	0.236%
65	14.542	2284	2289	2295	rBV2	196303	282598	16.28%	1.760%
66	14.676	2308	2311	2317	rVB4	43611	68775	3.96%	0.428%
67	14.737	2317	2321	2325	rVB2	55998	70931	4.09%	0.442%
68	14.786	2326	2329	2333	rVB2	18076	23259	1.34%	0.145%
69	14.847	2333	2339	2343	rBV2	439232	597860	34.45%	3.724%
70	14.914	2347	2350	2353	rVB5	28007	36287	2.09%	0.226%
71	14.987	2353	2362	2366	rVB8	20355	47433	2.73%	0.295%

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72	15.255	2401	2406	2413	rVB2	43455	79728	4.59%	0.497%
73	15.328	2414	2418	2424	rVB3	20509	29551	1.70%	0.184%
74	15.450	2434	2438	2440	rBV2	19523	27286	1.57%	0.170%
75	15.481	2440	2443	2448	rVV	37056	52923	3.05%	0.330%
76	15.554	2448	2455	2460	rVV2	71711	123521	7.12%	0.769%
77	15.603	2460	2463	2468	rVV4	15261	24807	1.43%	0.155%
78	15.694	2470	2478	2482	rVV2	98081	197862	11.40%	1.233%
79	15.731	2482	2484	2491	rVB	28486	45097	2.60%	0.281%
80	15.816	2491	2498	2505	rBV2	25153	55208	3.18%	0.344%
81	15.895	2506	2511	2513	rVV	31236	49564	2.86%	0.309%
82	15.926	2513	2516	2527	rVB2	48988	93755	5.40%	0.584%
83	16.078	2535	2541	2552	rVB2	32876	61685	3.55%	0.384%
84	16.286	2572	2575	2581	rVB6	10279	17815	1.03%	0.111%

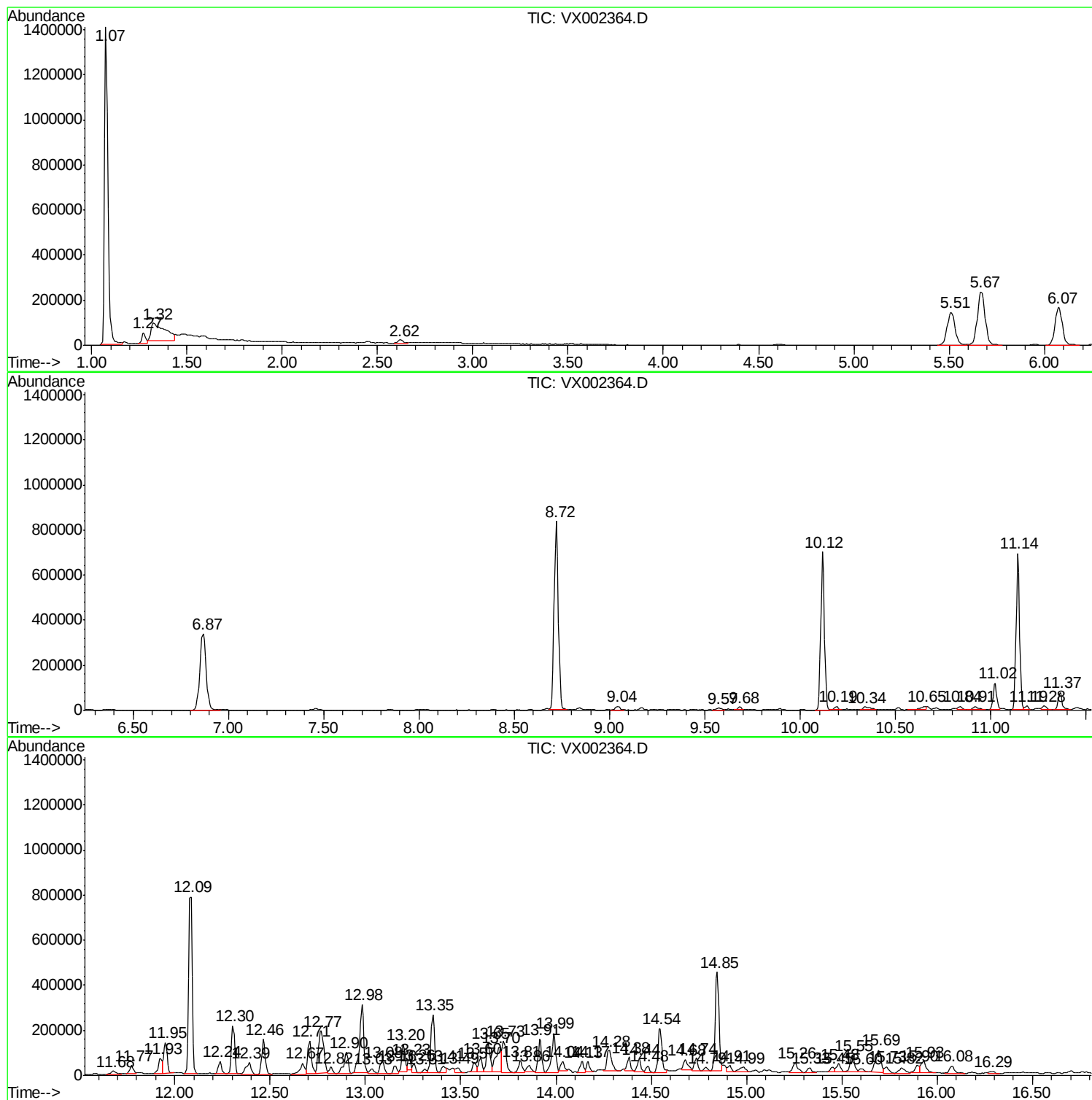
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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



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Instrument :
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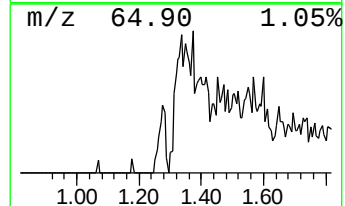
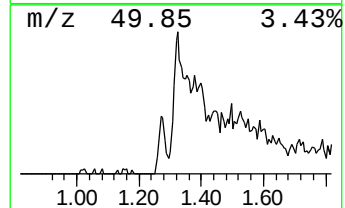
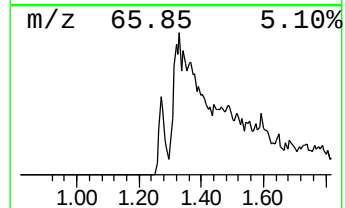
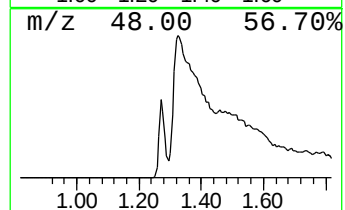
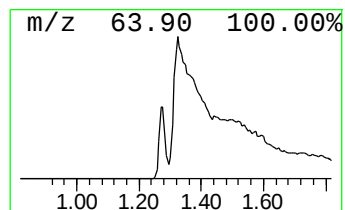
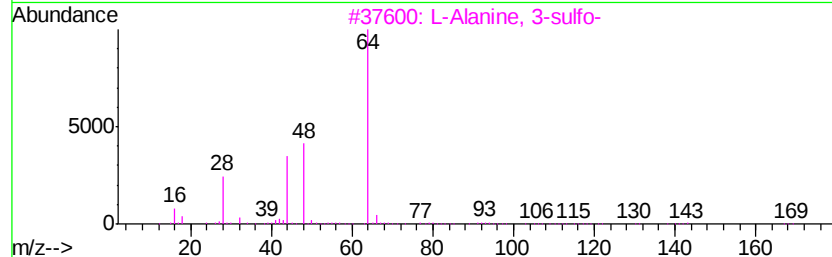
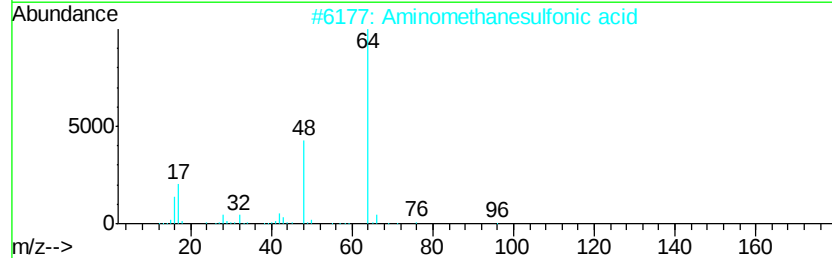
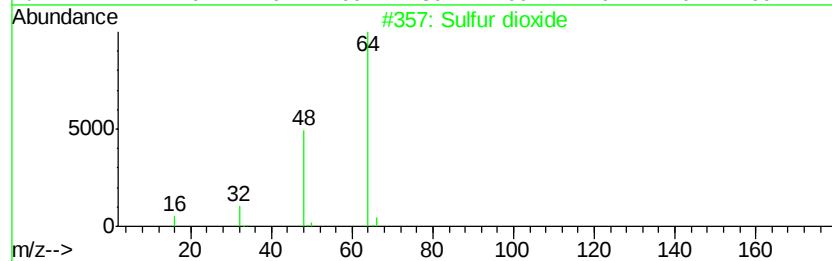
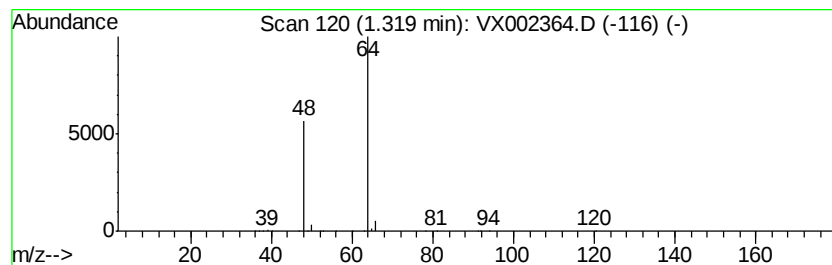
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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.
1.32	31.74 ug/l	418713	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	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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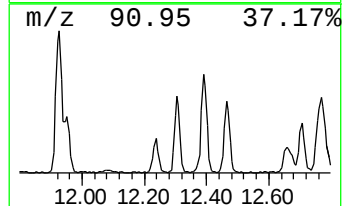
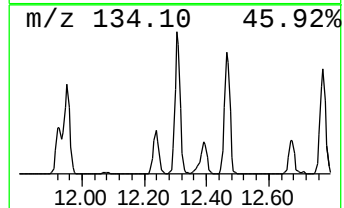
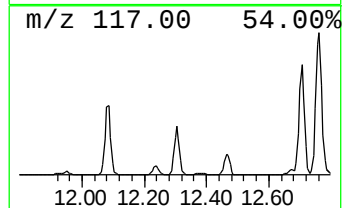
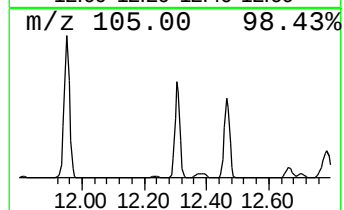
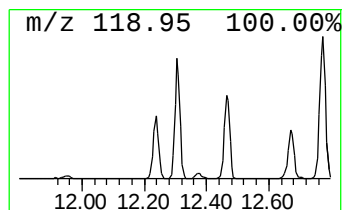
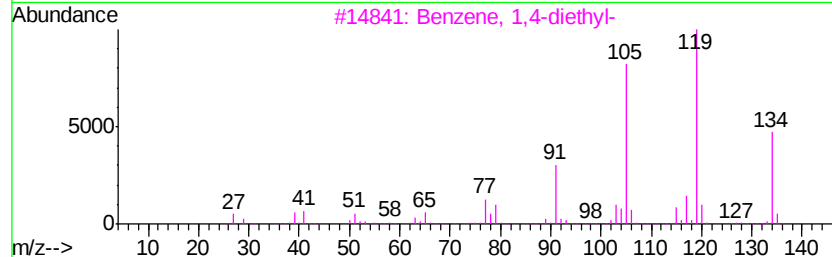
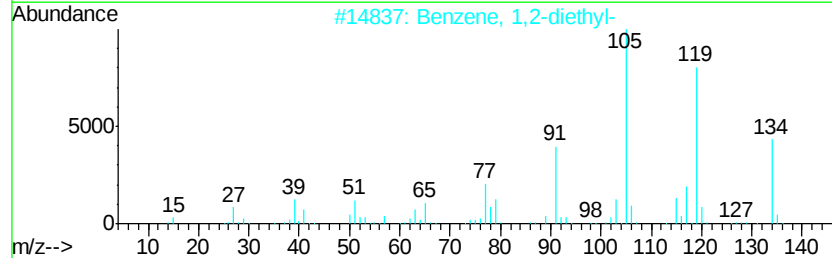
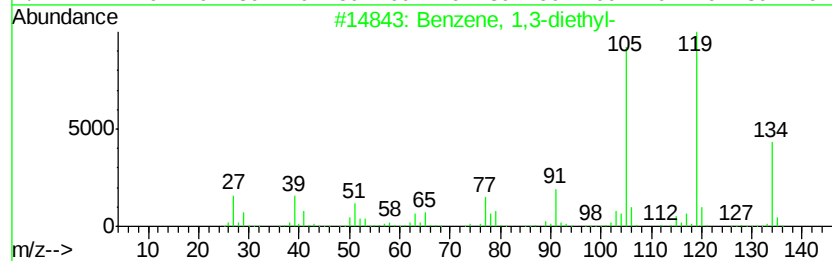
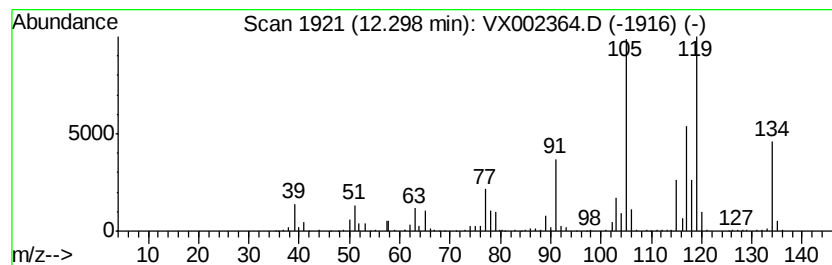
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	12.87 ug/l	259021	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 93
2		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 90
3		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 76
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 58
5		Benzene, tert-butyl-	134 C10H14	000098-06-6 49



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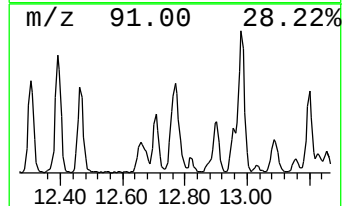
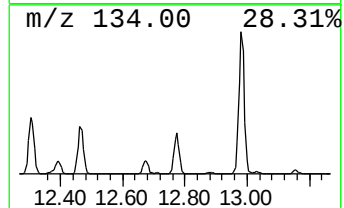
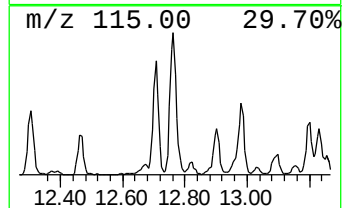
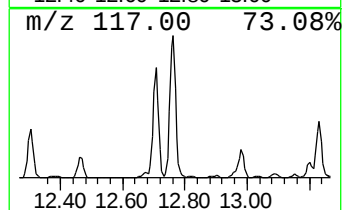
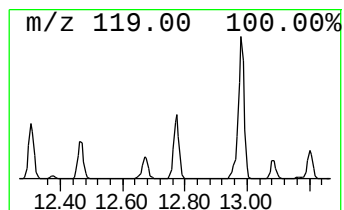
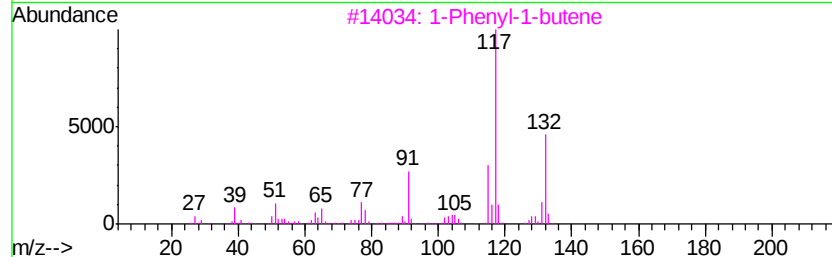
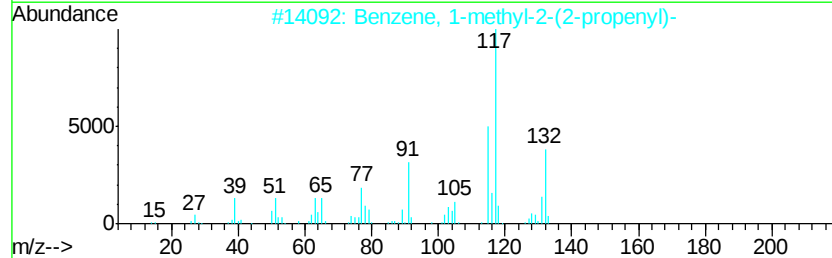
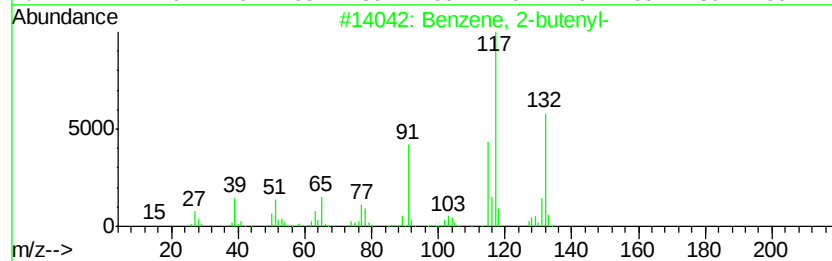
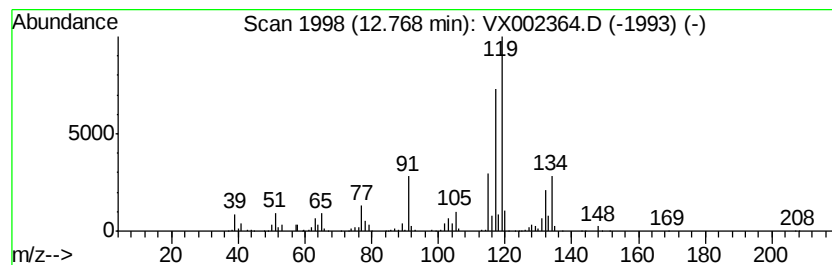
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-butenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	18.04 ug/l	363240	1,4-Dichlorobenzene-d4	12.09

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	84
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	84
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	60
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	60



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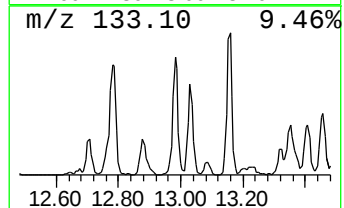
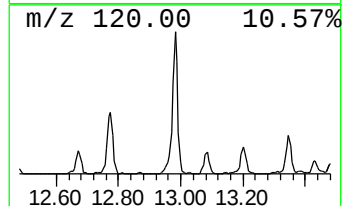
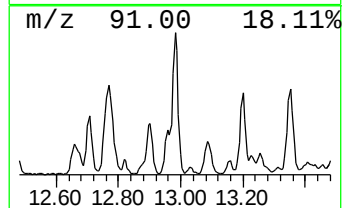
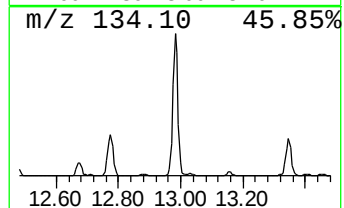
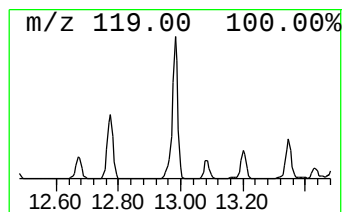
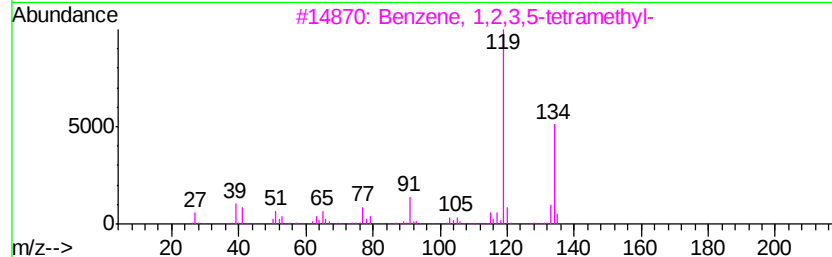
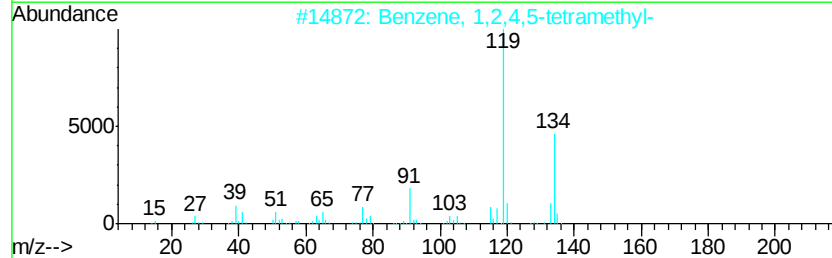
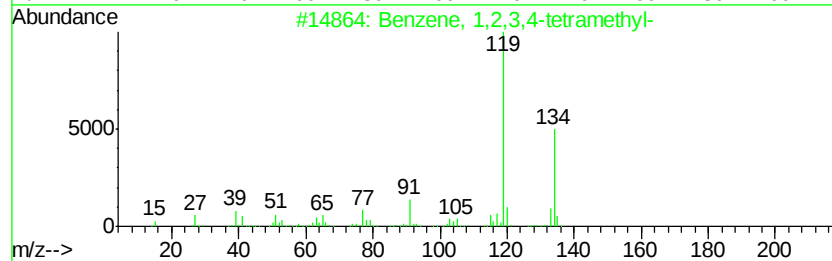
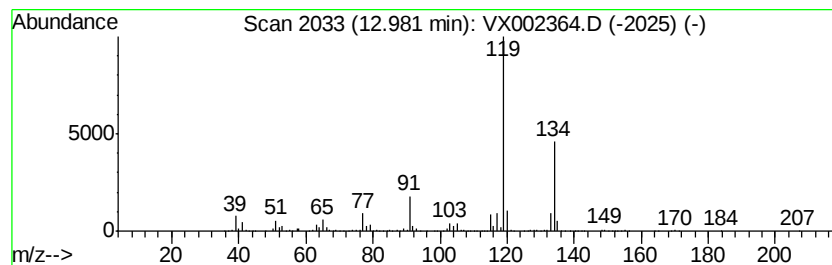
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	22.07 ug/l	444258	1,4-Dichlorobenzene-d4	12.09

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002364.D
Acq On : 06 Jun 2018 20:22
Operator : JC/MD
Sample : J3312-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-04

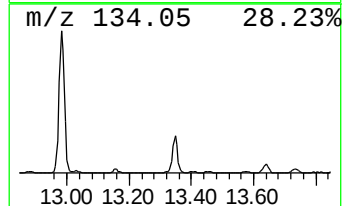
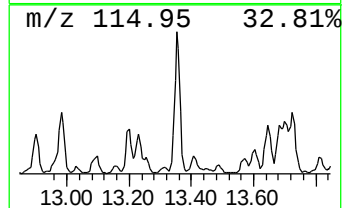
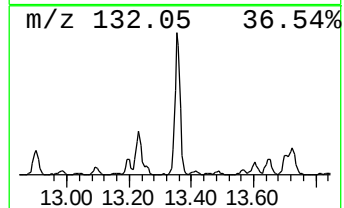
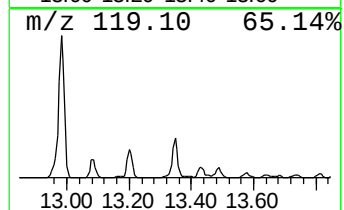
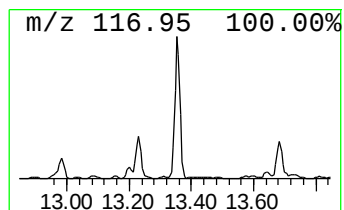
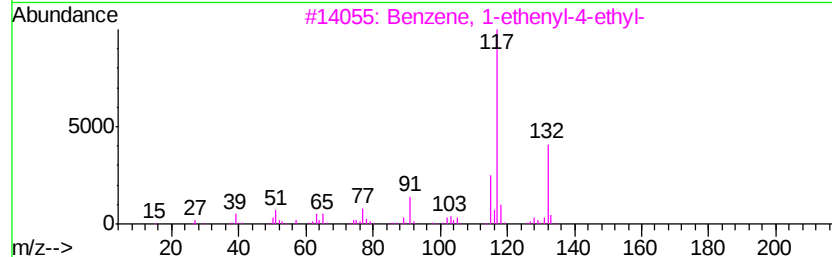
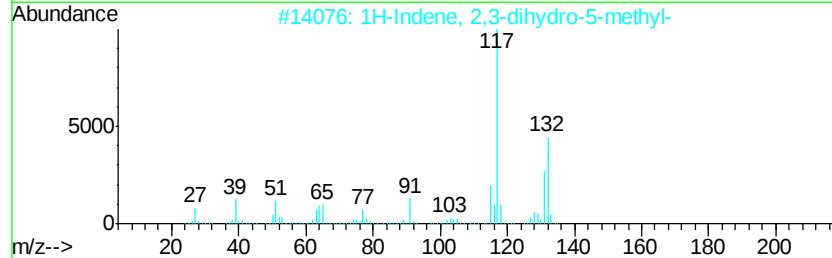
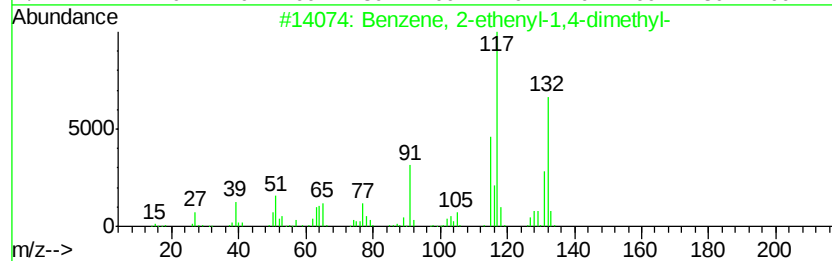
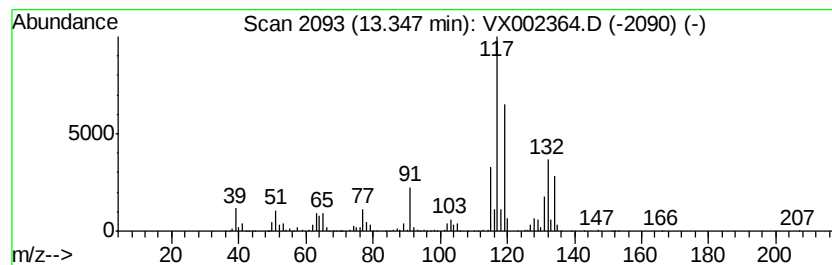
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	16.68 ug/l	335705	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002364.D
Acq On : 06 Jun 2018 20:22
Operator : JC/MD
Sample : J3312-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-04

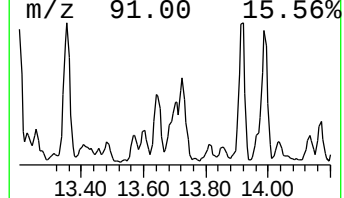
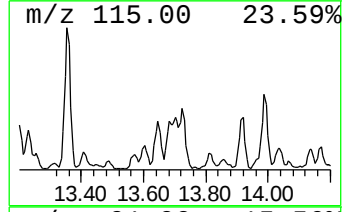
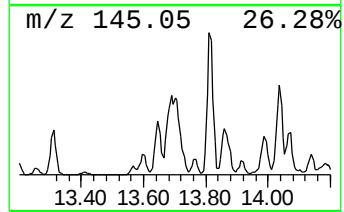
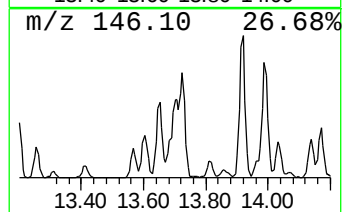
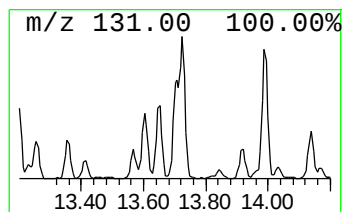
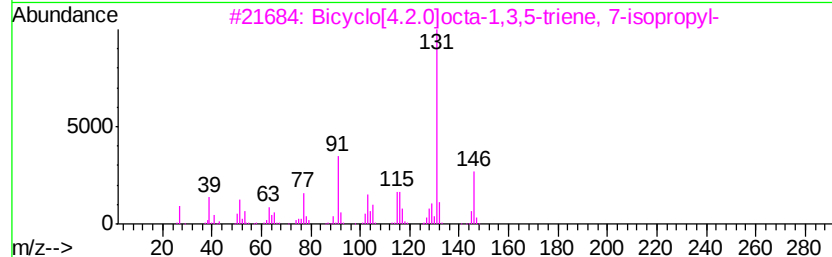
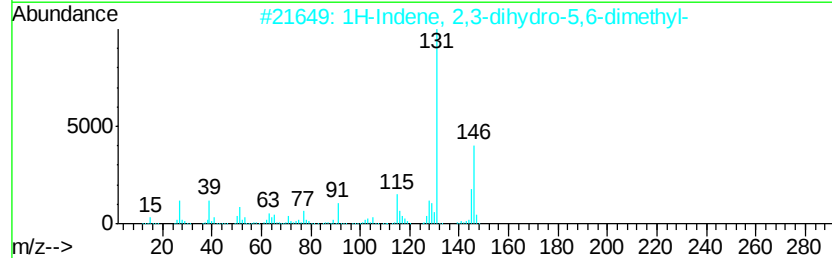
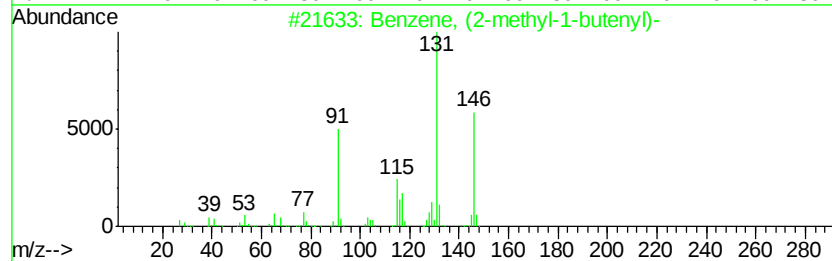
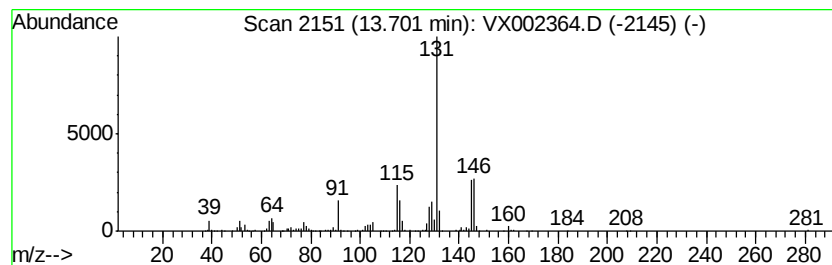
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	10.80 ug/l	217341	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90
2		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	90
3		Bicyclo[4.2.0]octa-1,3,5-triene...	146	C11H14	027087-54-3	87
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002364.D
Acq On : 06 Jun 2018 20:22
Operator : JC/MD
Sample : J3312-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-04

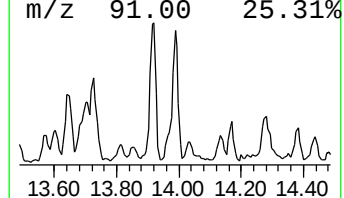
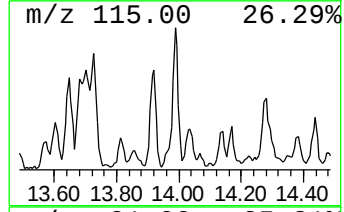
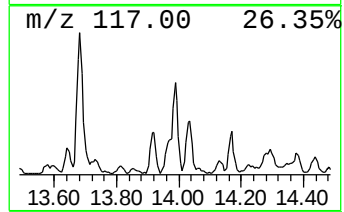
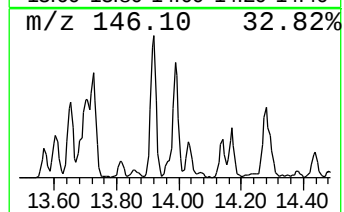
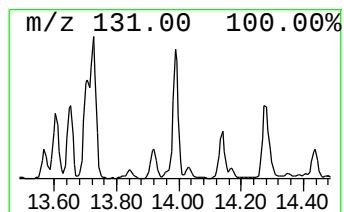
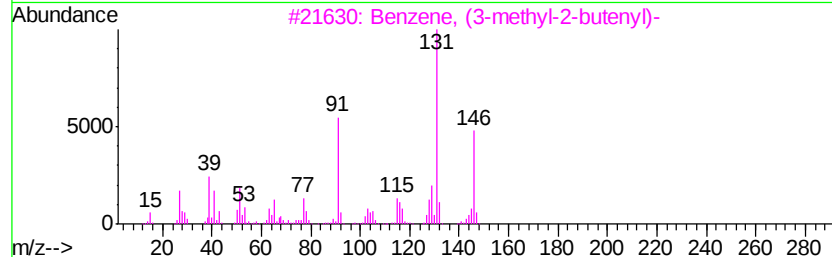
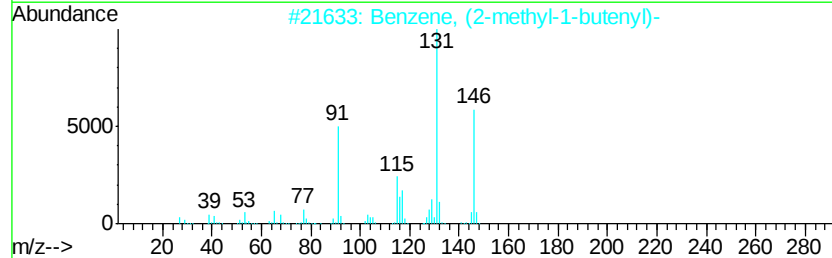
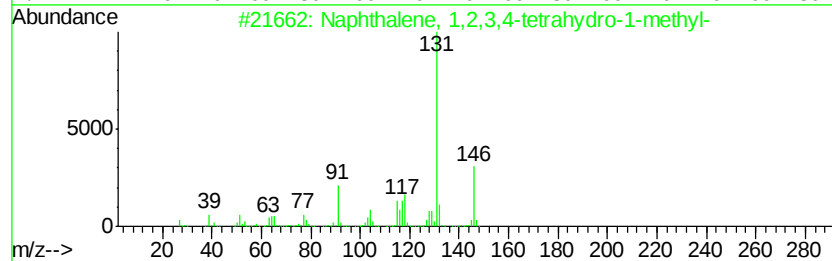
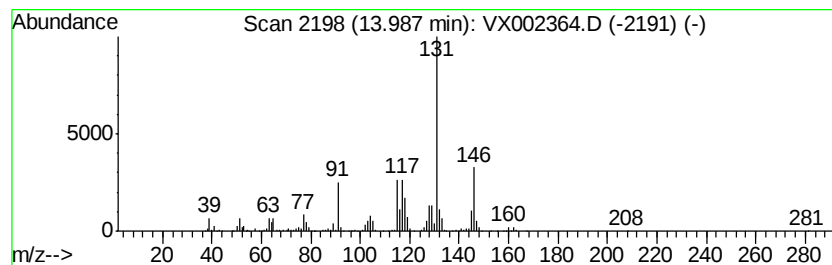
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	12.72 ug/l	255964	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	90
5		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002364.D
Acq On : 06 Jun 2018 20:22
Operator : JC/MD
Sample : J3312-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

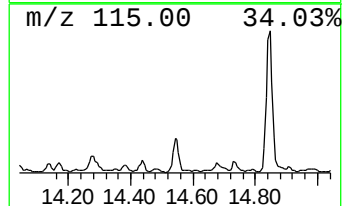
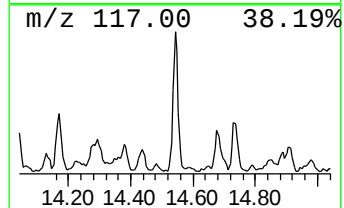
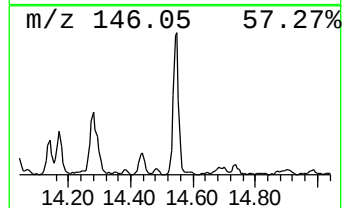
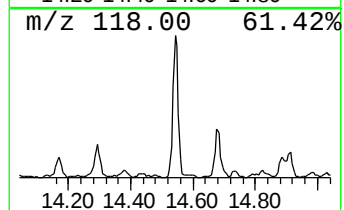
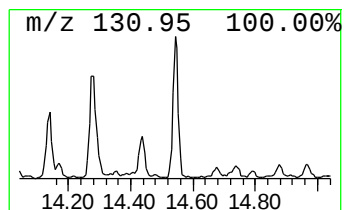
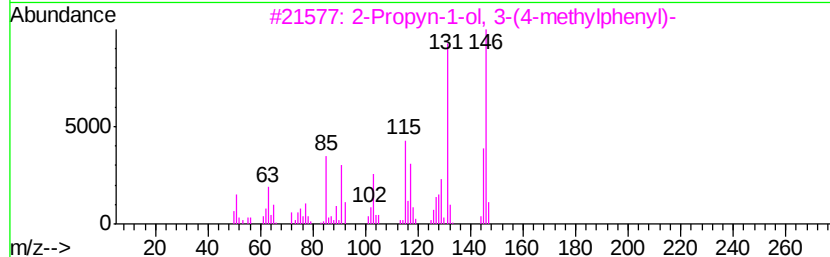
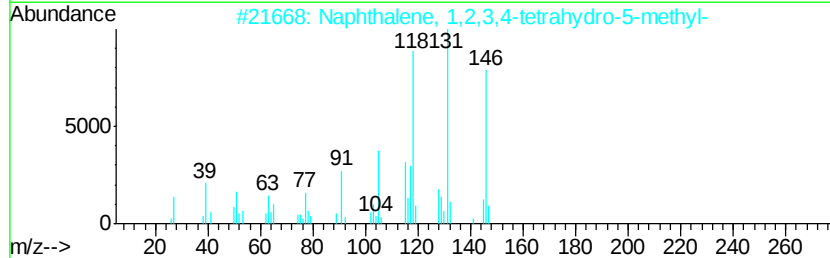
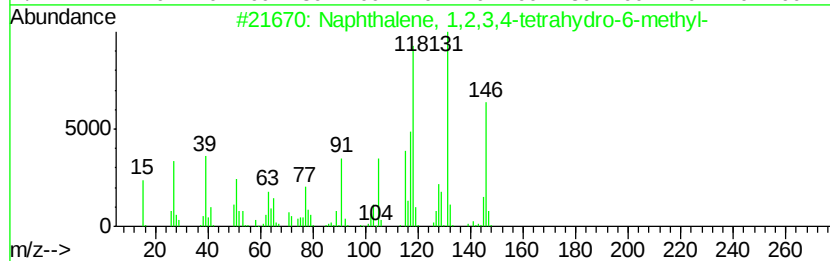
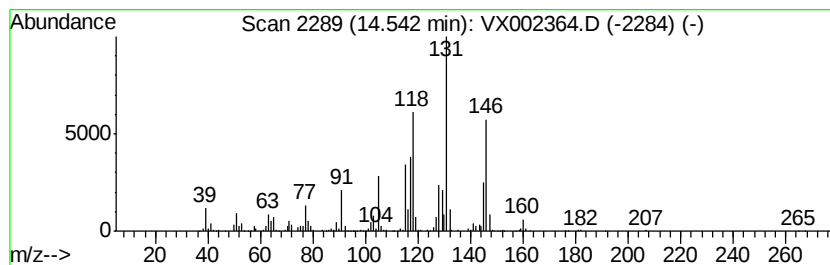
Instrument :
MSVOA_X
ClientSampled :
W-04

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	14.04 ug/l	282598	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 93
3		2-Propyn-1-ol, 3-(4-methylphenyl)-	146 C10H10O	016017-24-6 68
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 64
5		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002364.D
Acq On : 06 Jun 2018 20:22
Operator : JC/MD
Sample : J3312-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
W-04

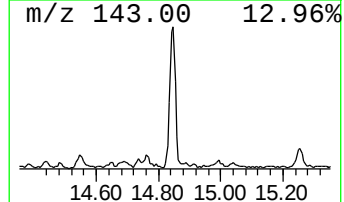
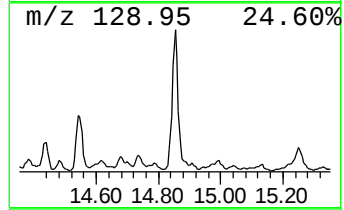
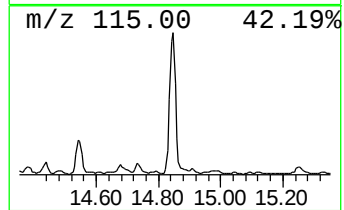
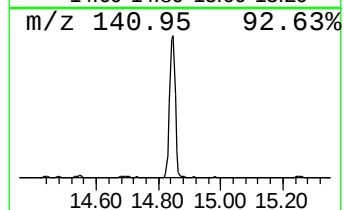
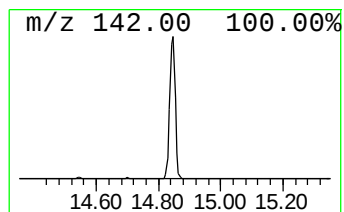
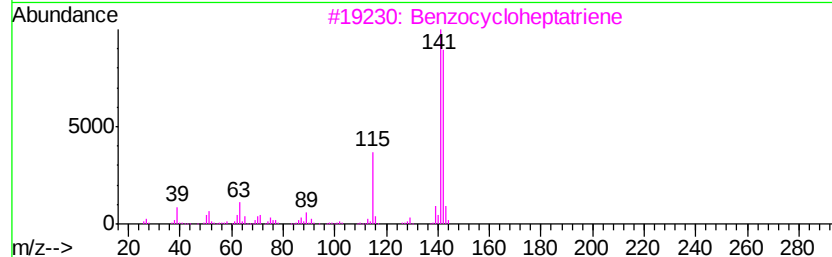
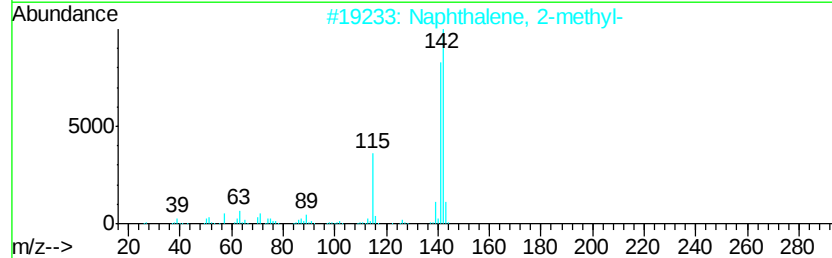
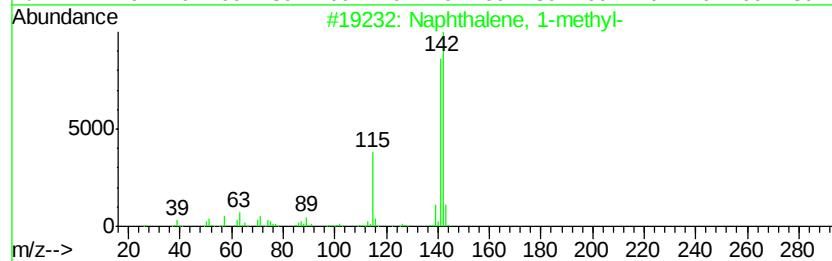
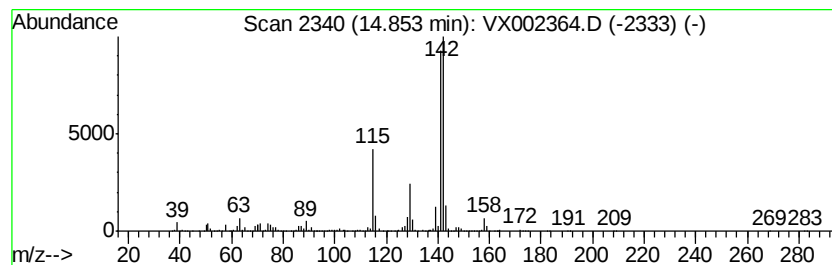
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	29.70 ug/l	597860	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	80



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002364.D
Acq On : 06 Jun 2018 20:22
Operator : JC/MD
Sample : J3312-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-04

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.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	1.32	31.7	ug/l	418713	1	5.67	659554	50.0
Benzene, 1,3-diet...	12.30	12.9	ug/l	259021	4	12.09	1006550	50.0
Benzene, 2-butenyl-	12.77	18.0	ug/l	363240	4	12.09	1006550	50.0
Benzene, 1,2,3,4-...	12.98	22.1	ug/l	444258	4	12.09	1006550	50.0
Benzene, 2-etheny...	13.35	16.7	ug/l	335705	4	12.09	1006550	50.0
Benzene, (2-methy...	13.70	10.8	ug/l	217341	4	12.09	1006550	50.0
Naphthalene, 1,2,...	13.99	12.7	ug/l	255964	4	12.09	1006550	50.0
Naphthalene, 1,2,...	14.54	14.0	ug/l	282598	4	12.09	1006550	50.0
Naphthalene, 1-me...	14.85	29.7	ug/l	597860	4	12.09	1006550	50.0