

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062818\
 Data File : VX003020.D
 Acq On : 29 Jun 2018 08:01
 Operator : JC/MD
 Sample : J3720-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-23

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

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	92	rBV	1411488	1704917	100.00%	8.516%
2	1.276	108	113	129	rBV	209498	848531	49.77%	4.238%
3	2.617	328	333	339	rBV	22618	40183	2.36%	0.201%
4	5.507	796	807	820	rBV	189666	537787	31.54%	2.686%
5	5.665	822	833	852	rVB	279820	798722	46.85%	3.990%
6	6.074	888	900	920	rBV	200332	529783	31.07%	2.646%
7	6.866	1018	1030	1047	rBV	404065	948179	55.61%	4.736%
8	8.720	1327	1334	1348	rVB	975646	1523348	89.35%	7.609%
9	9.043	1381	1387	1397	rVB	21544	36446	2.14%	0.182%
10	9.573	1468	1474	1479	rBV5	10160	19574	1.15%	0.098%
11	10.116	1557	1563	1571	rBV	875377	1218848	71.49%	6.088%
12	10.189	1571	1575	1579	rVB	15036	23004	1.35%	0.115%
13	10.335	1595	1599	1603	rBV3	15962	29451	1.73%	0.147%
14	10.512	1624	1628	1635	rVB2	11697	18348	1.08%	0.092%
15	10.646	1638	1650	1657	rBV6	21379	65593	3.85%	0.328%
16	10.841	1678	1682	1686	rVB	27303	39421	2.31%	0.197%
17	10.914	1689	1694	1706	rBV7	20667	46502	2.73%	0.232%
18	11.018	1706	1711	1716	rBV	67639	93100	5.46%	0.465%
19	11.140	1725	1731	1736	rBV	912377	1132755	66.44%	5.658%
20	11.189	1736	1739	1742	rVB	19393	24628	1.44%	0.123%
21	11.280	1750	1754	1760	rVB4	23094	34000	1.99%	0.170%
22	11.366	1763	1768	1771	rBV	27482	37678	2.21%	0.188%
23	11.445	1777	1781	1786	rBV4	11901	19863	1.17%	0.099%
24	11.676	1812	1819	1825	rBV2	21527	41002	2.40%	0.205%
25	11.774	1831	1835	1838	rBV	32278	38873	2.28%	0.194%
26	11.920	1852	1859	1861	rBV	59155	87013	5.10%	0.435%
27	11.951	1861	1864	1869	rVB	79288	102577	6.02%	0.512%
28	12.079	1880	1885	1891	rBV	1115364	1366940	80.18%	6.828%
29	12.176	1897	1901	1906	rBV6	7834	17864	1.05%	0.089%
30	12.237	1906	1911	1915	rVV	73356	94811	5.56%	0.474%
31	12.304	1917	1922	1927	rVV	257034	319445	18.74%	1.596%
32	12.371	1929	1933	1934	rVV2	49916	60131	3.53%	0.300%
33	12.390	1934	1936	1941	rVB	54870	67321	3.95%	0.336%
34	12.463	1943	1948	1955	rBV	116665	148467	8.71%	0.742%

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35	12.670	1974	1982	1985	rBV	228890	294910	17.30%	1.473%
36	12.707	1985	1988	1993	rVV2	74824	120708	7.08%	0.603%
37	12.762	1993	1997	2004	rVV4	177379	359369	21.08%	1.795%
38	12.817	2004	2006	2012	rVB2	32495	39580	2.32%	0.198%
39	12.877	2012	2016	2017	rBV2	50463	63881	3.75%	0.319%
40	12.896	2017	2019	2024	rVB	104148	127470	7.48%	0.637%
41	12.981	2024	2033	2038	rBV2	321285	505686	29.66%	2.526%
42	13.030	2038	2041	2046	rBV2	30769	49078	2.88%	0.245%
43	13.085	2046	2050	2056	rVB2	83816	132953	7.80%	0.664%
44	13.152	2056	2061	2065	rBV	64989	92490	5.42%	0.462%
45	13.195	2065	2068	2072	rBV2	119292	147255	8.64%	0.736%
46	13.225	2072	2073	2076	rVV	24124	22720	1.33%	0.113%
47	13.255	2076	2078	2082	rVB	33688	36693	2.15%	0.183%
48	13.316	2082	2088	2090	rBV2	50302	80901	4.75%	0.404%
49	13.347	2090	2093	2100	rVB4	84277	141285	8.29%	0.706%
50	13.408	2100	2103	2104	rBV2	29370	32118	1.88%	0.160%
51	13.426	2104	2106	2109	rVV	46170	49977	2.93%	0.250%
52	13.457	2109	2111	2113	rVV3	25755	32953	1.93%	0.165%
53	13.481	2113	2115	2122	rVB2	55600	75979	4.46%	0.380%
54	13.566	2125	2129	2132	rBV3	46268	73623	4.32%	0.368%
55	13.603	2132	2135	2138	rVV2	61429	72709	4.26%	0.363%
56	13.640	2138	2141	2145	rVV3	142604	184260	10.81%	0.920%
57	13.701	2145	2151	2153	rVV2	103858	223395	13.10%	1.116%
58	13.725	2153	2155	2164	rVB3	142523	222078	13.03%	1.109%
59	13.810	2164	2169	2173	rBV	86212	115509	6.78%	0.577%
60	13.859	2173	2177	2181	rVB2	49765	64420	3.78%	0.322%
61	13.914	2181	2186	2190	rVB2	144315	201283	11.81%	1.005%
62	13.987	2190	2198	2202	rBV2	189935	306428	17.97%	1.531%
63	14.036	2202	2206	2209	rBV2	42612	60590	3.55%	0.303%
64	14.133	2218	2222	2225	rBV3	61734	84192	4.94%	0.421%
65	14.164	2225	2227	2231	rVB	39728	40945	2.40%	0.205%
66	14.225	2231	2237	2240	rBV3	26861	47592	2.79%	0.238%
67	14.274	2240	2245	2254	rVB3	133410	236258	13.86%	1.180%
68	14.383	2255	2263	2267	rBV	117163	174054	10.21%	0.869%
69	14.432	2267	2271	2276	rVB2	82789	111869	6.56%	0.559%
70	14.481	2276	2279	2283	rVB	60966	74180	4.35%	0.371%
71	14.542	2284	2289	2293	rBV2	161295	244663	14.35%	1.222%

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72	14.639	2302	2305	2308	rBV3	18182	24362	1.43%	0.122%
73	14.676	2308	2311	2317	rVV3	88947	143307	8.41%	0.716%
74	14.731	2317	2320	2324	rVB2	116656	141888	8.32%	0.709%
75	14.786	2326	2329	2333	rVV	45445	59091	3.47%	0.295%
76	14.853	2335	2340	2343	rVV	126536	197510	11.58%	0.987%
77	14.908	2347	2349	2353	rVB3	45615	57552	3.38%	0.287%
78	14.981	2353	2361	2366	rBV8	27317	74685	4.38%	0.373%
79	15.115	2376	2383	2387	rBV7	21052	47630	2.79%	0.238%
80	15.255	2400	2406	2413	rVB2	78442	152207	8.93%	0.760%
81	15.328	2413	2418	2424	rVB7	32203	61447	3.60%	0.307%
82	15.450	2434	2438	2440	rBV2	50649	66106	3.88%	0.330%
83	15.481	2440	2443	2447	rVB	70447	95848	5.62%	0.479%
84	15.554	2448	2455	2460	rBV	212381	333731	19.57%	1.667%
85	15.694	2470	2478	2483	rBV	333647	541182	31.74%	2.703%
86	15.810	2492	2497	2503	rBV2	55525	111317	6.53%	0.556%
87	15.895	2505	2511	2512	rVV	100823	147225	8.64%	0.735%
88	15.926	2512	2516	2525	rVB	149385	298057	17.48%	1.489%
89	16.078	2535	2541	2552	rVB	90316	172900	10.14%	0.864%
90	16.176	2554	2557	2563	rVB5	14871	23495	1.38%	0.117%
91	16.261	2566	2571	2572	rBV3	10798	17392	1.02%	0.087%
92	16.572	2617	2622	2627	rBV3	15821	31450	1.84%	0.157%
93	16.700	2633	2643	2648	rBV	40671	87658	5.14%	0.438%
94	16.755	2648	2652	2657	rBV	41258	74573	4.37%	0.372%

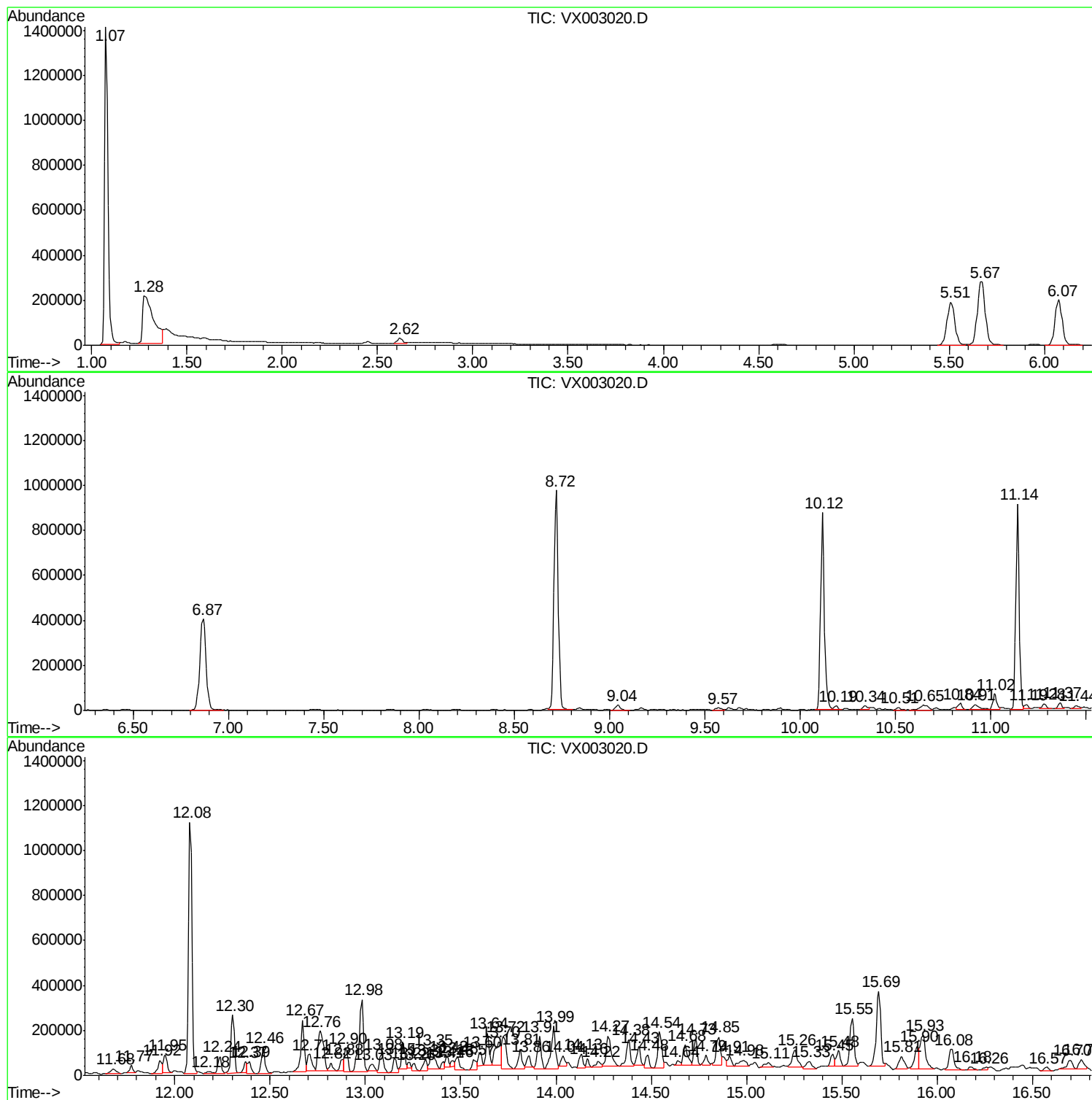
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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



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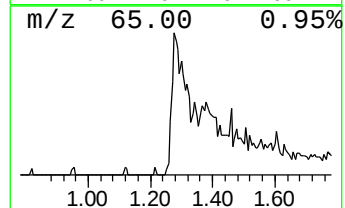
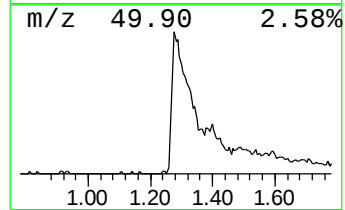
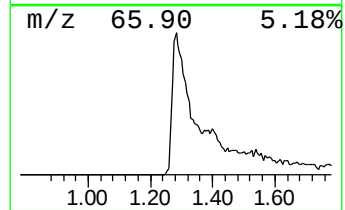
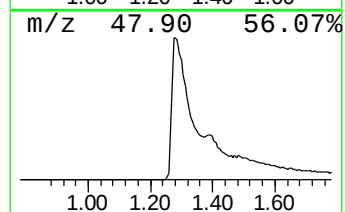
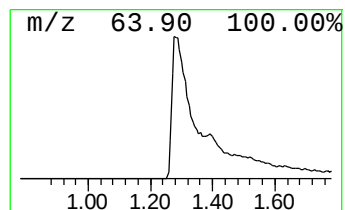
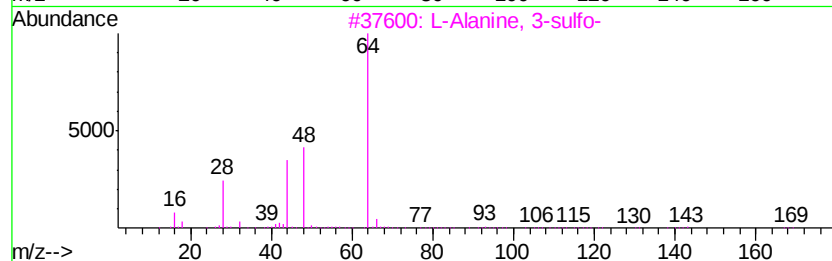
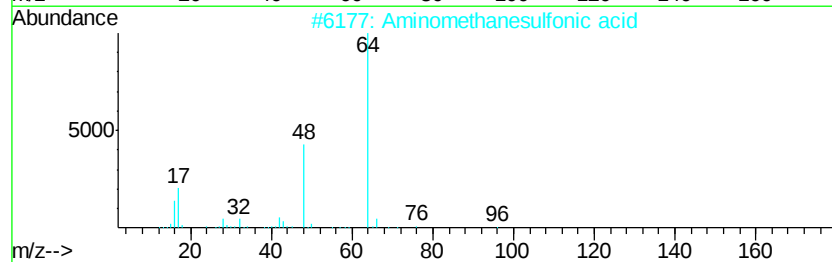
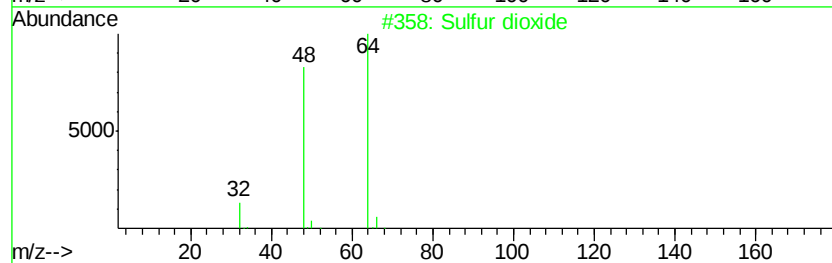
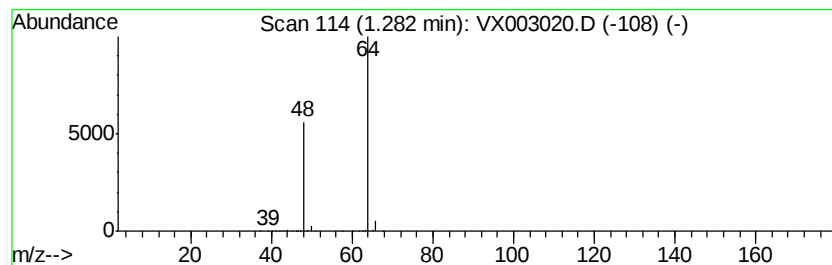
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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	53.12 ug/l	848531	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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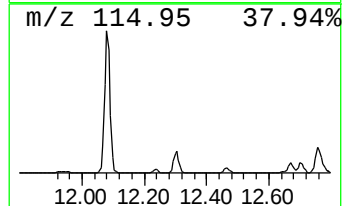
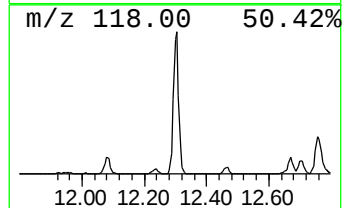
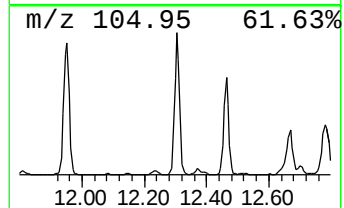
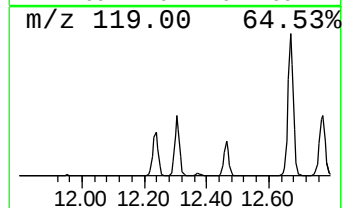
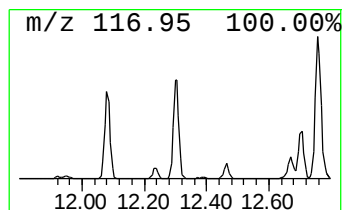
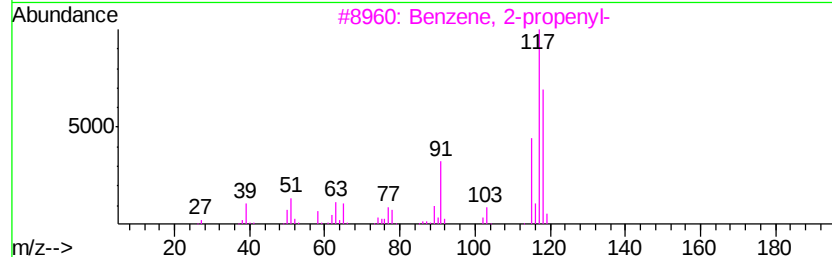
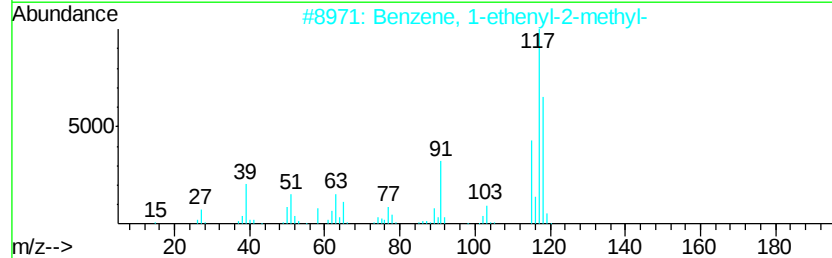
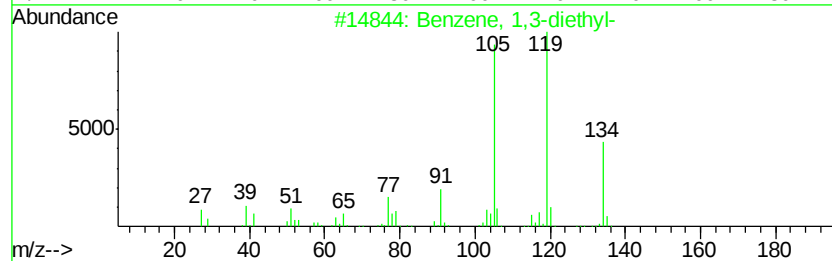
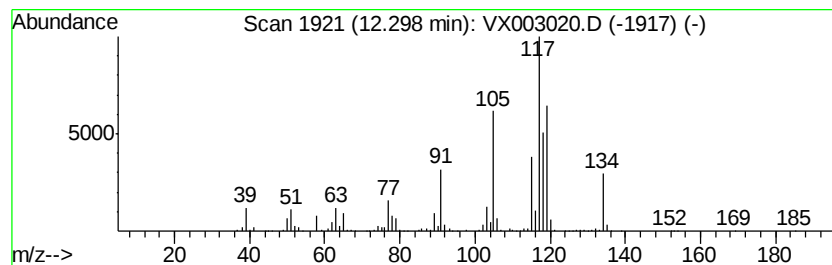
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,3-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	11.68 ug/l	319445	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 87
2		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 86
3		Benzene, 2-propenyl-	118 C9H10	000300-57-2 80
4		Benzene, 1-ethenyl-4-methyl-	118 C9H10	000622-97-9 60
5		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 55



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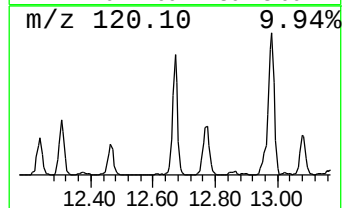
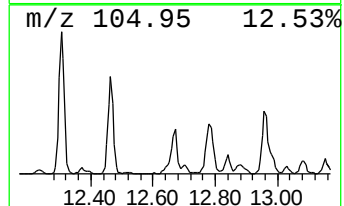
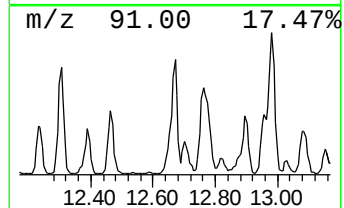
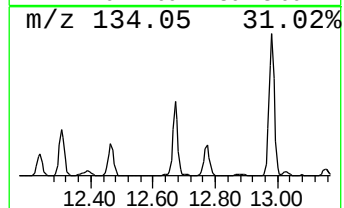
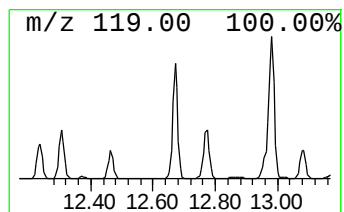
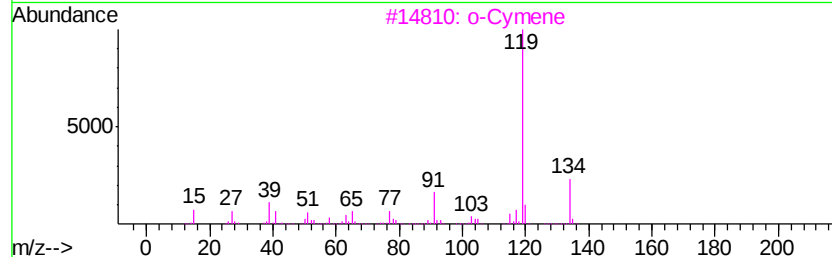
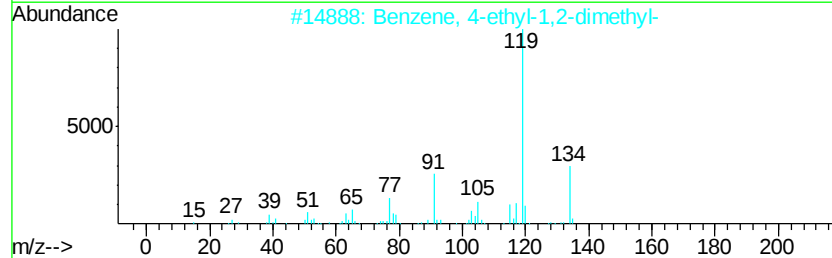
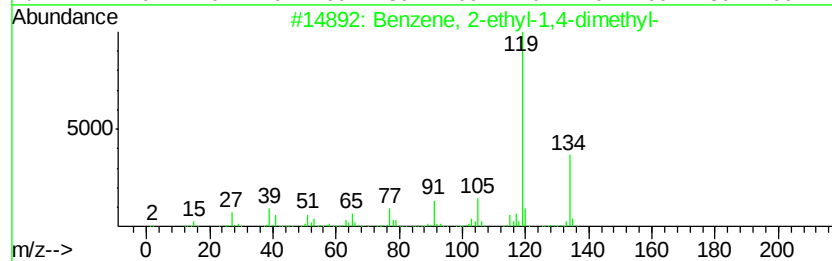
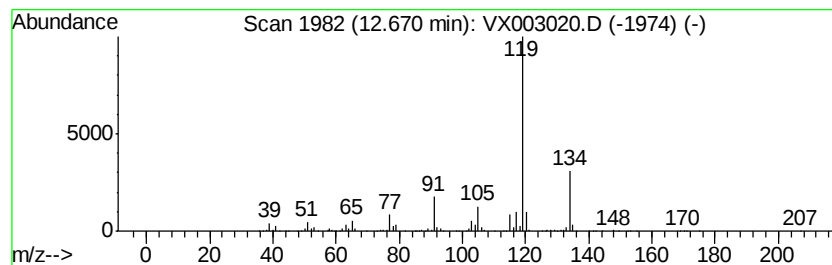
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	10.79 ug/l	294910	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		o-Cymene	134	C10H14	000527-84-4	94
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5		p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062818\
Data File : VX003020.D
Acq On : 29 Jun 2018 08:01
Operator : JC/MD
Sample : J3720-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-23

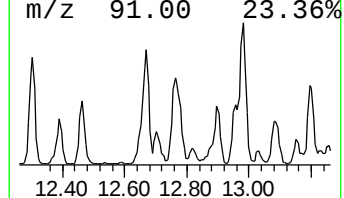
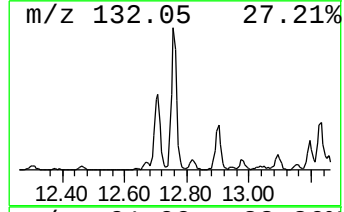
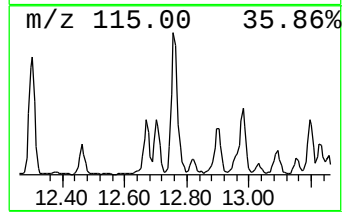
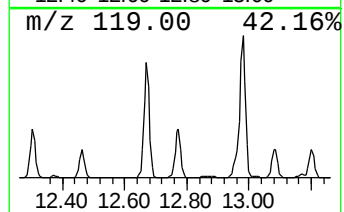
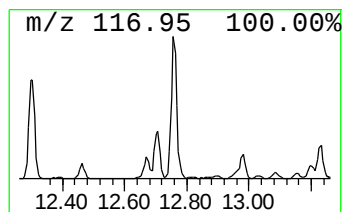
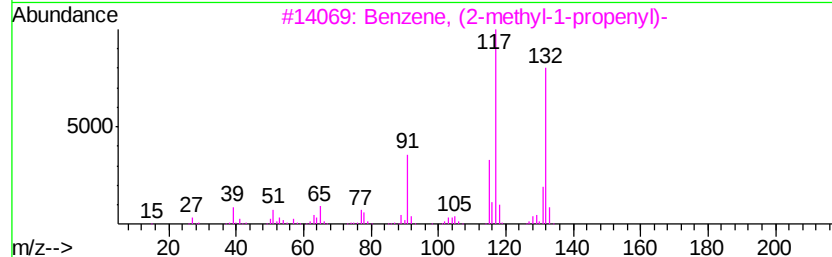
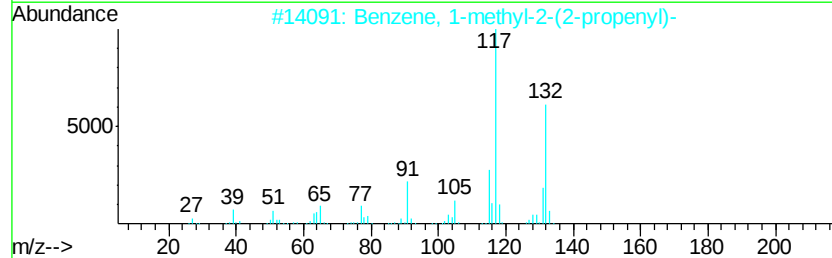
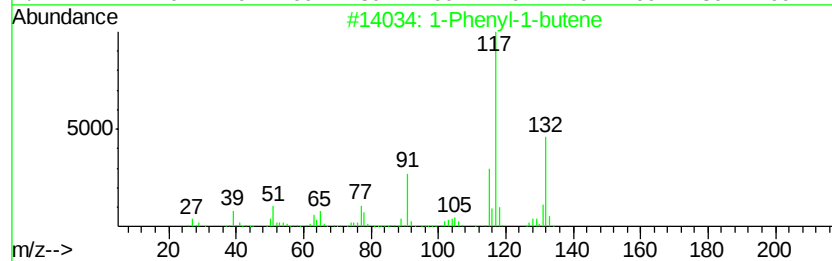
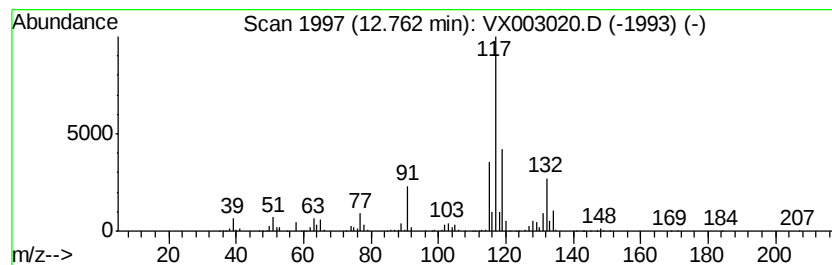
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 5 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	13.15 ug/l	359369	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	76
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	62
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062818\
Data File : VX003020.D
Acq On : 29 Jun 2018 08:01
Operator : JC/MD
Sample : J3720-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-23

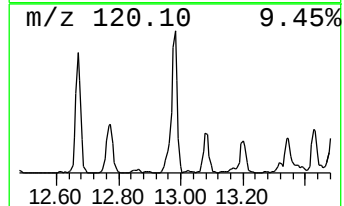
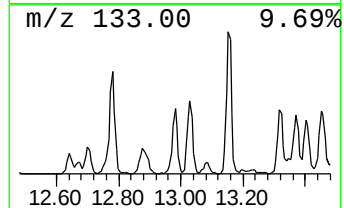
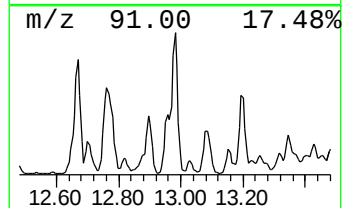
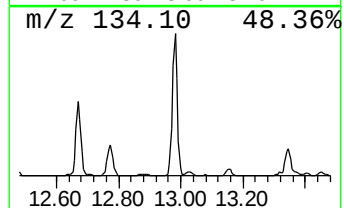
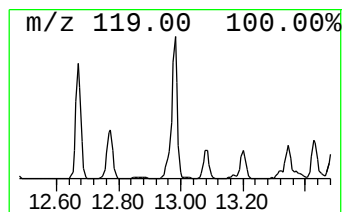
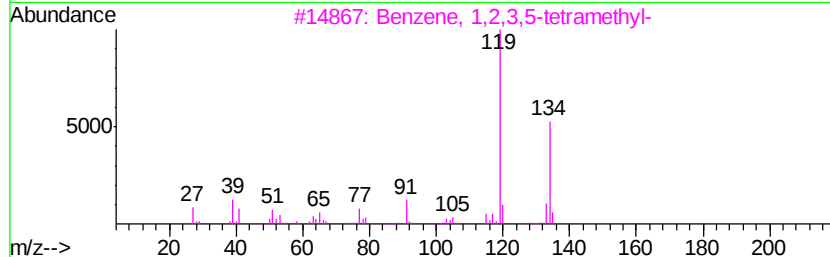
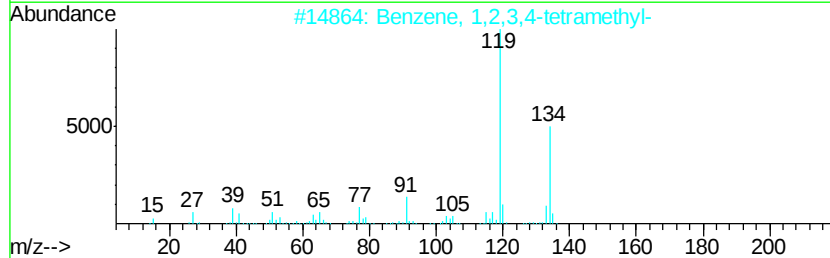
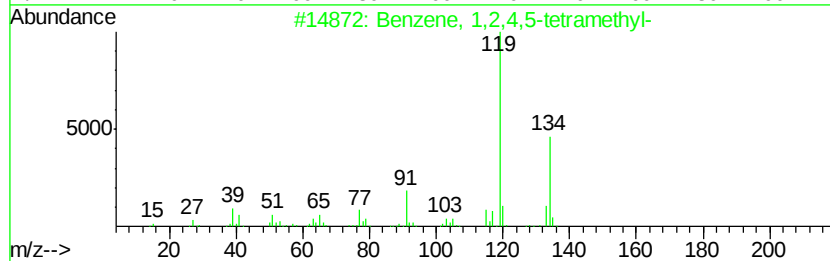
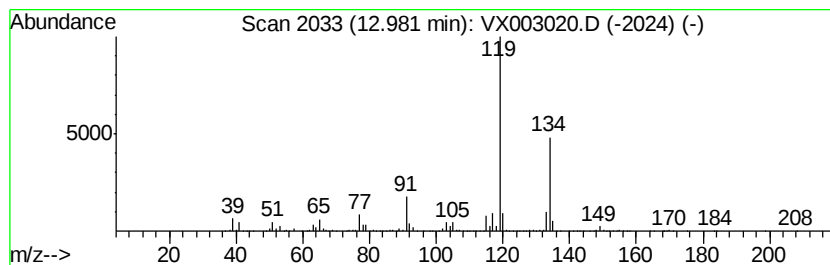
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	18.50 ug/l	505686	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	97
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-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062818\
Data File : VX003020.D
Acq On : 29 Jun 2018 08:01
Operator : JC/MD
Sample : J3720-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-23

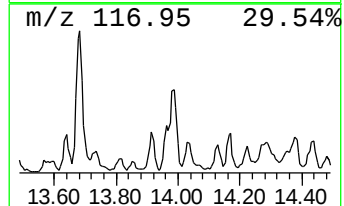
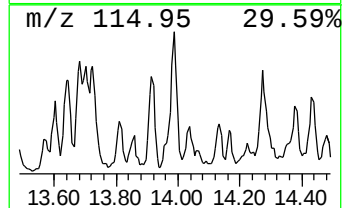
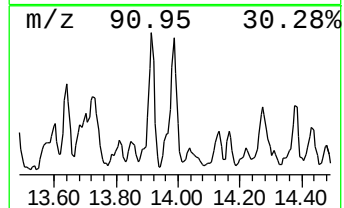
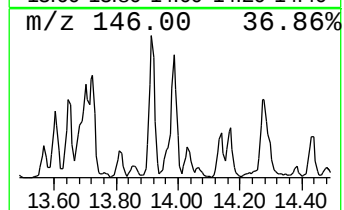
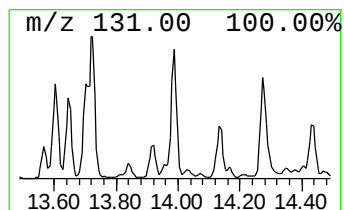
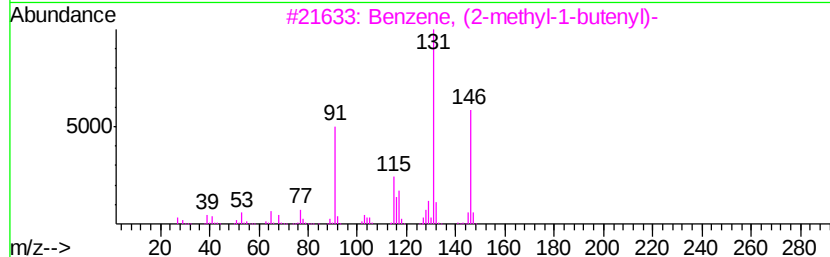
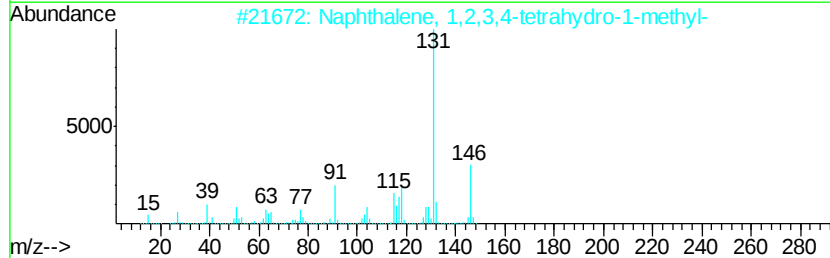
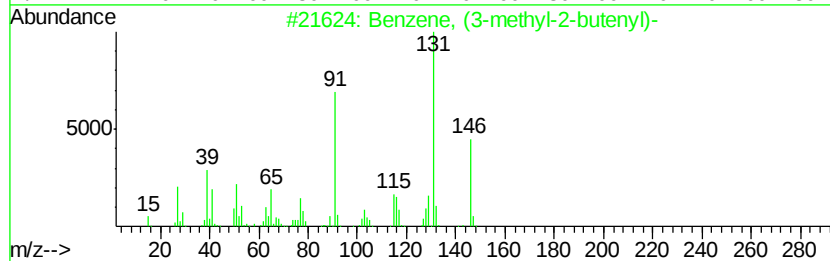
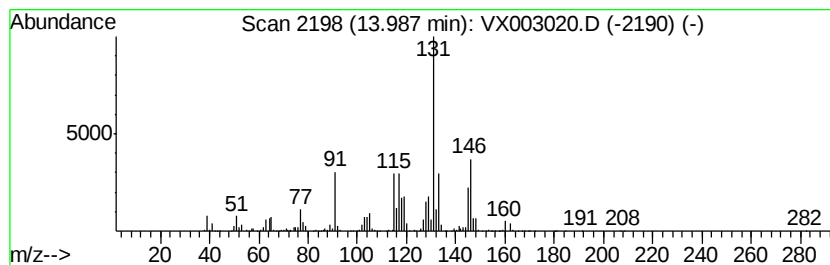
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (3-methyl-2-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	11.21 ug/l	306428	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70
3		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70
4		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70
5		2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062818\
Data File : VX003020.D
Acq On : 29 Jun 2018 08:01
Operator : JC/MD
Sample : J3720-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 47 Sample Multiplier: 1

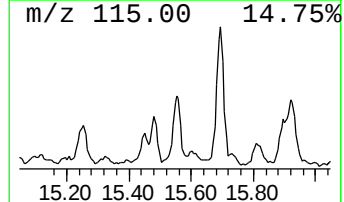
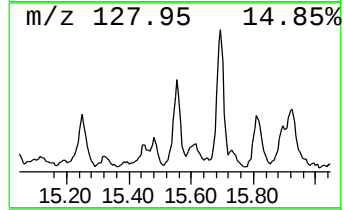
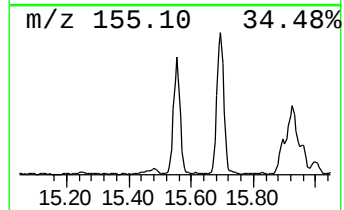
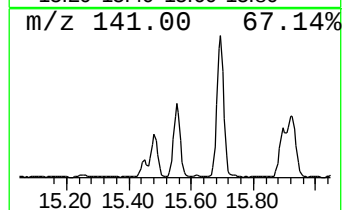
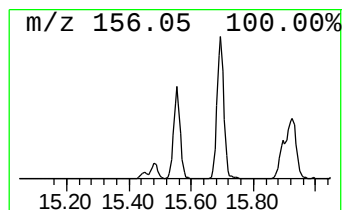
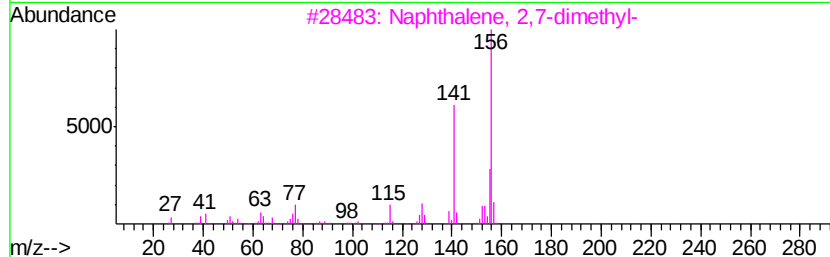
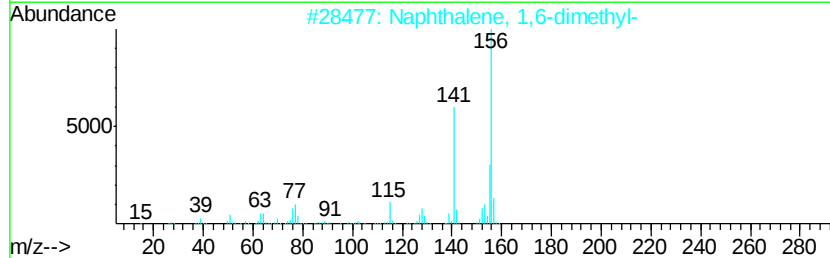
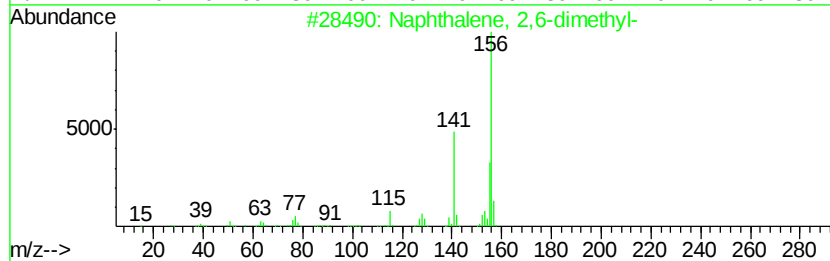
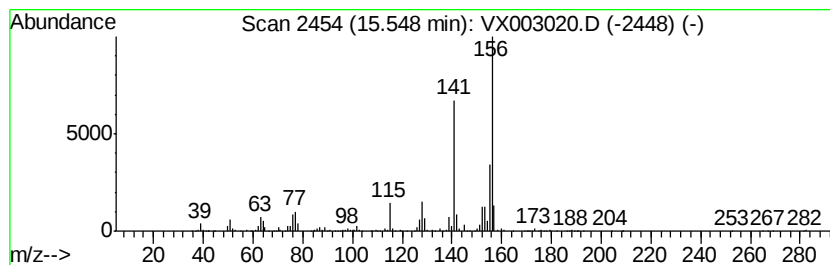
Instrument :
MSVOA_X
ClientSampleId :
MW-23

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, 2,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	12.21 ug/l	333731	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
2		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
3		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
5		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 96



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062818\
Data File : VX003020.D
Acq On : 29 Jun 2018 08:01
Operator : JC/MD
Sample : J3720-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-23

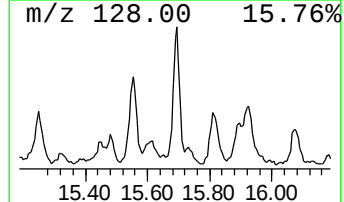
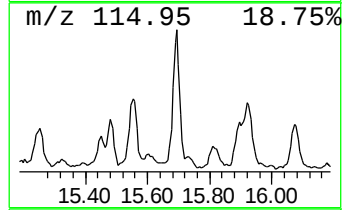
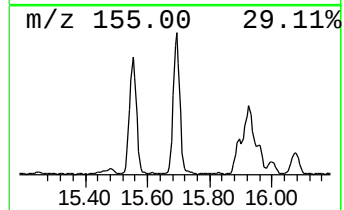
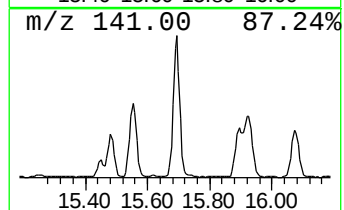
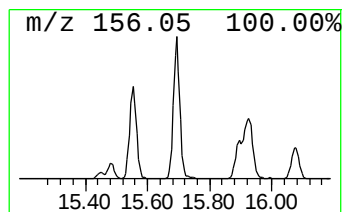
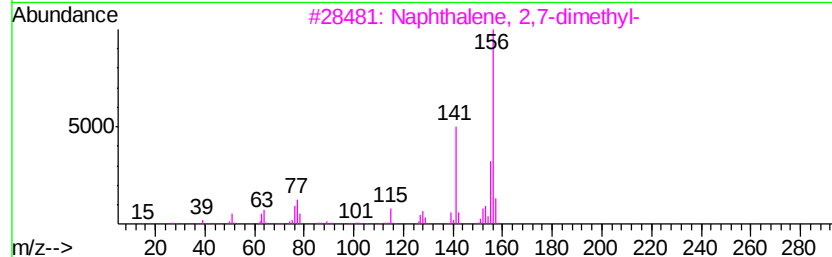
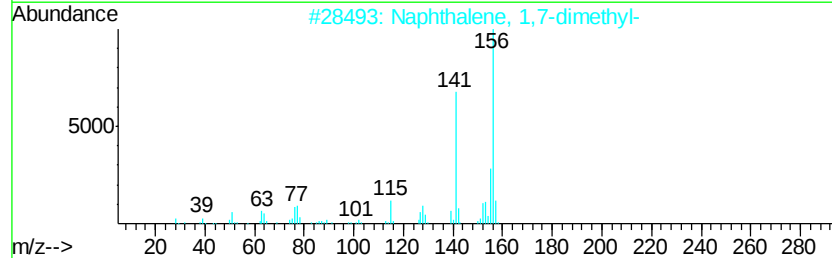
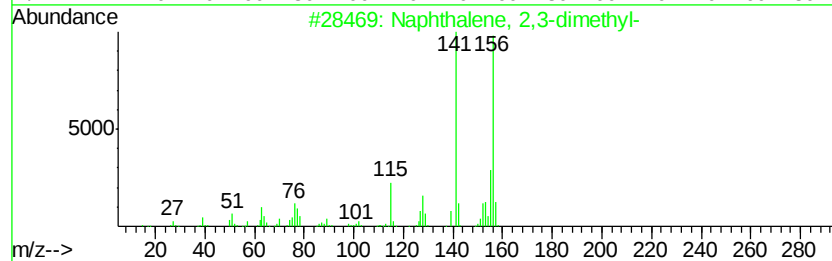
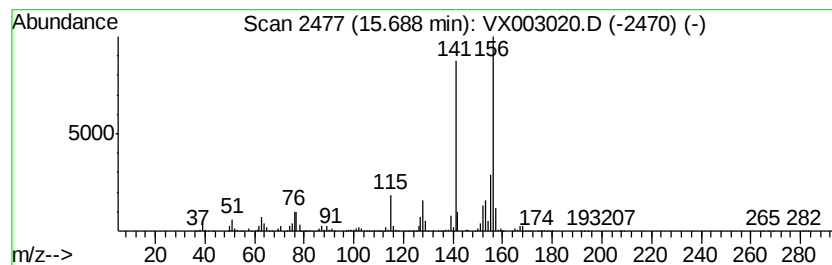
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	19.80 ug/l	541182	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062818\
Data File : VX003020.D
Acq On : 29 Jun 2018 08:01
Operator : JC/MD
Sample : J3720-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-23

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
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Benzene, 1,3-diet...	12.30	11.7	ug/l	319445	4	12.08	1366940	50.0
Benzene, 2-ethyl-...	12.67	10.8	ug/l	294910	4	12.08	1366940	50.0
1-Phenyl-1-butene	12.76	13.2	ug/l	359369	4	12.08	1366940	50.0
Benzene, 1,2,4,5-...	12.98	18.5	ug/l	505686	4	12.08	1366940	50.0
Benzene, (3-methy...	13.99	11.2	ug/l	306428	4	12.08	1366940	50.0
Naphthalene, 2,6-...	15.55	12.2	ug/l	333731	4	12.08	1366940	50.0
Naphthalene, 2,3-...	15.69	19.8	ug/l	541182	4	12.08	1366940	50.0