

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060518\
 Data File : VX002338.D
 Acq On : 06 Jun 2018 06:37
 Operator : JC/MD
 Sample : J3309-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-07

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	93	rBV	1723381	2140056	46.12%	3.575%
2	1.404	130	134	140	rVB	570256	570586	12.30%	0.953%
3	1.794	193	198	212	rVB	1471693	2004289	43.19%	3.348%
4	2.001	227	232	245	rVB	497755	737585	15.89%	1.232%
5	2.270	271	276	289	rVB	199379	313800	6.76%	0.524%
6	2.843	361	370	373	rVV3	99450	316260	6.82%	0.528%
7	2.892	373	378	381	rVV	328702	680359	14.66%	1.137%
8	2.934	381	385	400	rVB	625564	1339403	28.86%	2.238%
9	3.172	415	424	438	rVB	248432	586867	12.65%	0.980%
10	3.995	552	559	567	rVB	44153	109779	2.37%	0.183%
11	4.403	614	626	638	rVB	662887	1922024	41.42%	3.211%
12	5.507	792	807	811	rBV	153990	414747	8.94%	0.693%
13	5.580	811	819	828	rVV	700432	2209699	47.62%	3.691%
14	5.665	828	833	859	rVB	239386	827845	17.84%	1.383%
15	5.940	864	878	890	rVV6	95830	334806	7.21%	0.559%
16	6.074	890	900	907	rVV	177516	457673	9.86%	0.765%
17	6.147	907	912	920	rVV2	50609	137129	2.96%	0.229%
18	6.306	921	938	943	rVV2	221654	932946	20.10%	1.559%
19	6.367	943	948	956	rVV3	130474	413554	8.91%	0.691%
20	6.458	956	963	979	rVB	165896	467192	10.07%	0.780%
21	6.866	1020	1030	1039	rBV	353712	839575	18.09%	1.403%
22	7.464	1114	1128	1141	rBV	807405	1905529	41.06%	3.183%
23	7.738	1166	1173	1182	rVB2	82655	158276	3.41%	0.264%
24	7.933	1198	1205	1206	rBV	68630	106974	2.31%	0.179%
25	7.958	1207	1209	1215	rVV2	82456	134131	2.89%	0.224%
26	8.214	1238	1251	1256	rBV	64248	135057	2.91%	0.226%
27	8.580	1305	1311	1319	rVB3	100927	202934	4.37%	0.339%
28	8.671	1319	1326	1329	rBV2	157016	299184	6.45%	0.500%
29	8.720	1329	1334	1341	rVV	871646	1415280	30.50%	2.364%
30	8.787	1341	1345	1350	rVV	88910	131792	2.84%	0.220%
31	8.842	1350	1354	1358	rVV	69710	101374	2.18%	0.169%
32	8.915	1363	1366	1373	rVB4	56926	100214	2.16%	0.167%
33	9.043	1382	1387	1393	rVB	150110	234286	5.05%	0.391%
34	9.165	1398	1407	1412	rVV	95605	182501	3.93%	0.305%

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Integrator: RTE
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35	9.226	1412	1417	1429	rVB	75232	134410	2.90%	0.225%
36	9.628	1478	1483	1487	rBV	135993	197418	4.25%	0.330%
37	10.116	1554	1563	1568	rVV	744959	1011398	21.80%	1.690%
38	10.189	1568	1575	1581	rVV	101363	196685	4.24%	0.329%
39	10.256	1581	1586	1591	rVV	125609	187198	4.03%	0.313%
40	10.360	1591	1603	1609	rVB	246754	478441	10.31%	0.799%
41	10.646	1642	1650	1656	rBV6	44488	139765	3.01%	0.233%
42	10.701	1656	1659	1669	rVB	72821	111772	2.41%	0.187%
43	10.841	1670	1682	1689	rBV2	82376	152476	3.29%	0.255%
44	10.914	1689	1694	1700	rBV6	50466	119047	2.57%	0.199%
45	11.024	1706	1712	1716	rBV	462296	607146	13.08%	1.014%
46	11.140	1726	1731	1736	rBV	794662	975323	21.02%	1.629%
47	11.366	1763	1768	1774	rVV	476326	608864	13.12%	1.017%
48	11.670	1813	1818	1827	rBV	229709	323078	6.96%	0.540%
49	11.774	1827	1835	1838	rBV	75286	110474	2.38%	0.185%
50	11.951	1861	1864	1870	rVV	127967	162504	3.50%	0.271%
51	12.079	1880	1885	1891	rBV	920715	1180934	25.45%	1.973%
52	12.237	1907	1911	1916	rVB	171970	208873	4.50%	0.349%
53	12.304	1916	1922	1927	rBV	2469082	2985833	64.34%	4.988%
54	12.372	1927	1933	1943	rVB2	233027	414754	8.94%	0.693%
55	12.463	1943	1948	1954	rBV	404050	504776	10.88%	0.843%
56	12.670	1974	1982	1985	rBV	882351	1069358	23.04%	1.786%
57	12.707	1985	1988	1992	rVV	350493	448178	9.66%	0.749%
58	12.762	1992	1997	2004	rVV2	1430776	2313846	49.86%	3.865%
59	12.817	2004	2006	2010	rVV	98000	127665	2.75%	0.213%
60	12.902	2014	2020	2025	rVB	267114	407718	8.79%	0.681%
61	12.981	2025	2033	2037	rBV	1028670	1327241	28.60%	2.217%
62	13.024	2037	2040	2046	rVB2	103213	149346	3.22%	0.249%
63	13.091	2046	2051	2056	rBV3	213978	387512	8.35%	0.647%
64	13.158	2058	2062	2065	rVV	111920	157232	3.39%	0.263%
65	13.201	2065	2069	2071	rVV2	335472	445706	9.60%	0.745%
66	13.231	2071	2074	2083	rVB2	579609	841989	18.14%	1.407%
67	13.316	2083	2088	2089	rBV2	152092	187869	4.05%	0.314%
68	13.353	2089	2094	2100	rVV2	3319356	4640438	100.00%	7.752%
69	13.408	2100	2103	2105	rVV2	130795	177835	3.83%	0.297%
70	13.432	2105	2107	2109	rVV	100528	110061	2.37%	0.184%
71	13.487	2111	2116	2122	rVB	588662	782446	16.86%	1.307%

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72	13.566	2122	2129	2132	rBV3	179059	279685	6.03%	0.467%
73	13.603	2132	2135	2138	rVV	375444	525106	11.32%	0.877%
74	13.646	2138	2142	2146	rVV3	637495	1055775	22.75%	1.764%
75	13.701	2146	2151	2152	rVV2	412088	582825	12.56%	0.974%
76	13.725	2152	2155	2161	rVV2	722612	1140248	24.57%	1.905%
77	13.810	2165	2169	2172	rVV	148623	213665	4.60%	0.357%
78	13.853	2172	2176	2182	rVB3	224833	392335	8.45%	0.655%
79	13.914	2182	2186	2191	rBV	361234	453804	9.78%	0.758%
80	13.987	2191	2198	2202	rBV2	532092	741888	15.99%	1.239%
81	14.036	2202	2206	2209	rVV	214984	285802	6.16%	0.477%
82	14.140	2218	2223	2225	rBV2	213005	284288	6.13%	0.475%
83	14.170	2225	2228	2231	rVB	95286	105047	2.26%	0.175%
84	14.280	2241	2246	2253	rBV2	400826	742650	16.00%	1.241%
85	14.383	2259	2263	2267	rBV	188274	214683	4.63%	0.359%
86	14.438	2267	2272	2276	rVB	240045	312724	6.74%	0.522%
87	14.481	2276	2279	2284	rVB	88556	112530	2.42%	0.188%
88	14.542	2284	2289	2294	rBV2	737930	1002967	21.61%	1.676%
89	14.682	2307	2312	2313	rBV3	117546	162669	3.51%	0.272%
90	14.700	2313	2315	2318	rVV	126636	140162	3.02%	0.234%
91	14.737	2318	2321	2325	rVB	125616	150393	3.24%	0.251%
92	14.847	2333	2339	2343	rBV	1298145	1732122	37.33%	2.894%
93	14.987	2353	2362	2367	rVB4	52365	133630	2.88%	0.223%
94	15.255	2400	2406	2414	rBV2	128133	223511	4.82%	0.373%
95	15.554	2450	2455	2460	rVV	249453	401385	8.65%	0.671%
96	15.694	2471	2478	2482	rVV	341051	558936	12.04%	0.934%
97	15.731	2482	2484	2491	rVB	161367	219788	4.74%	0.367%
98	15.901	2506	2512	2513	rVV2	73387	124504	2.68%	0.208%
99	15.926	2513	2516	2525	rVB	117957	211157	4.55%	0.353%
100	16.078	2536	2541	2553	rVB	54986	107017	2.31%	0.179%

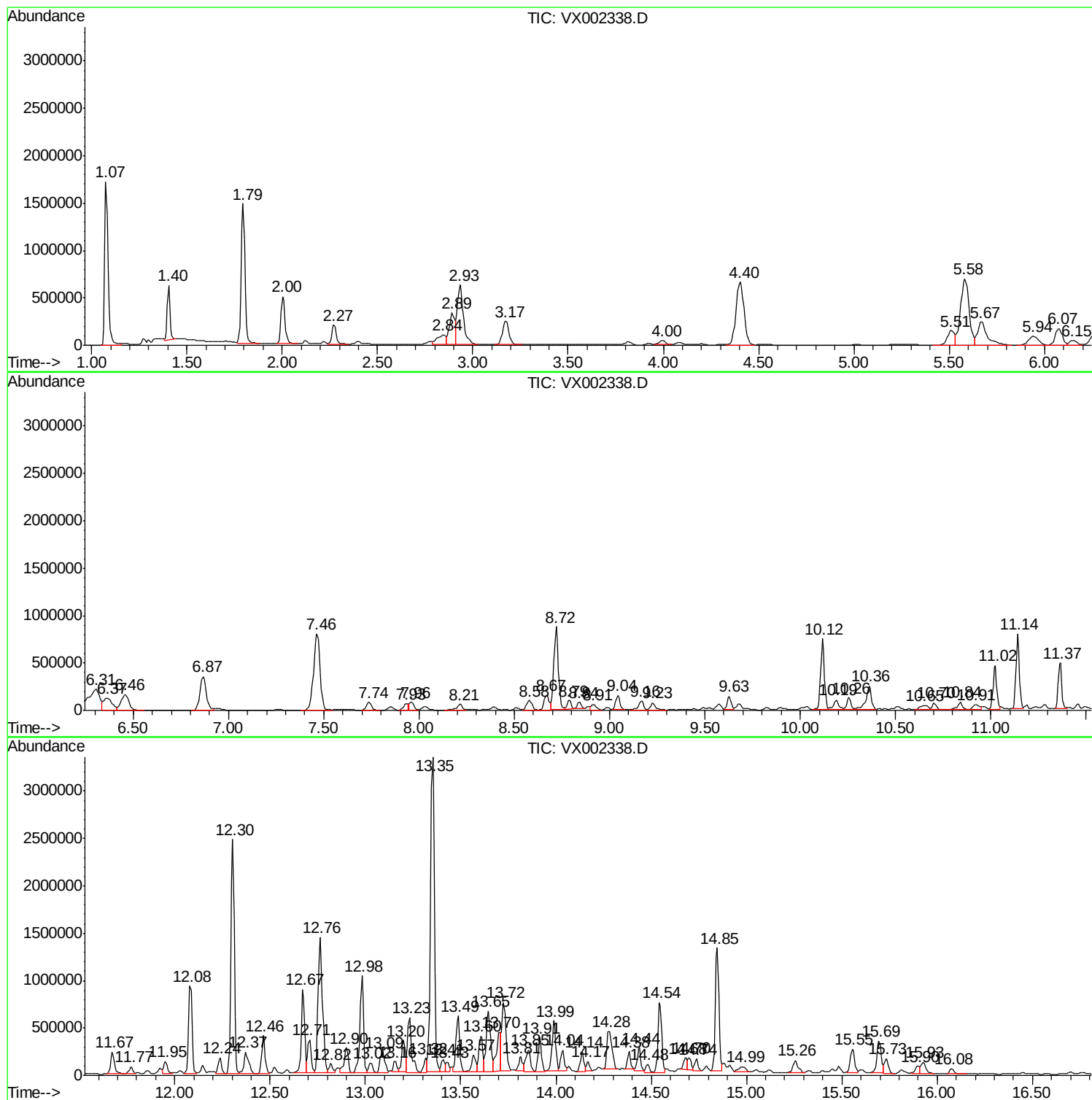
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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



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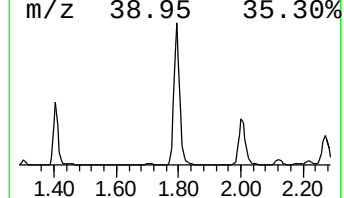
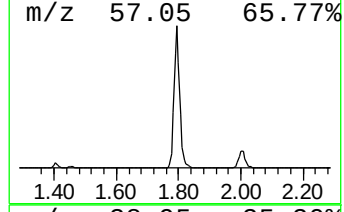
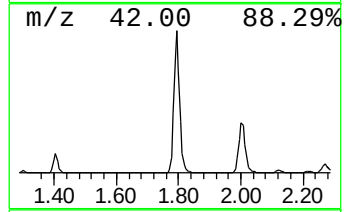
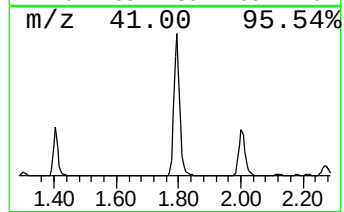
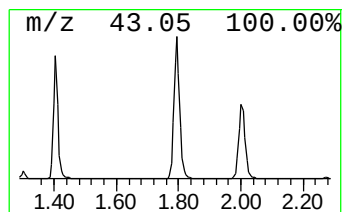
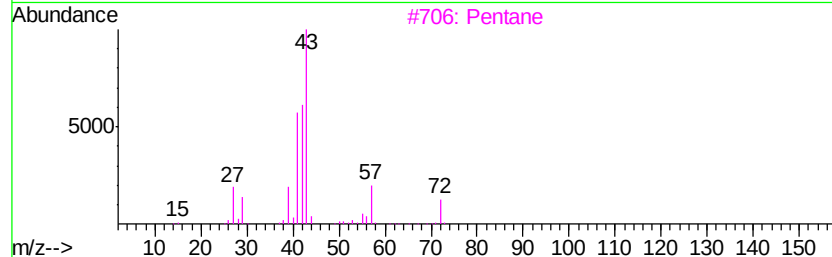
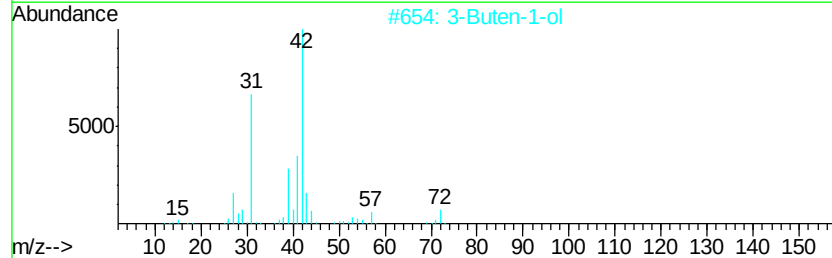
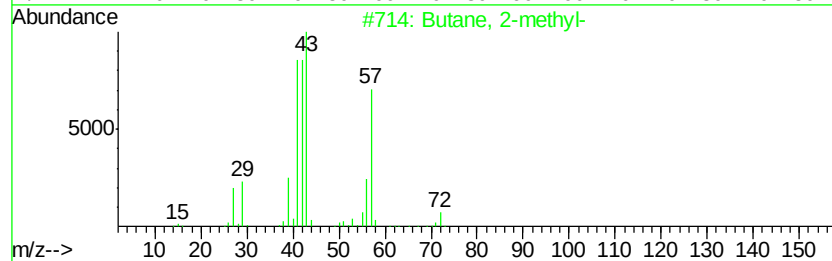
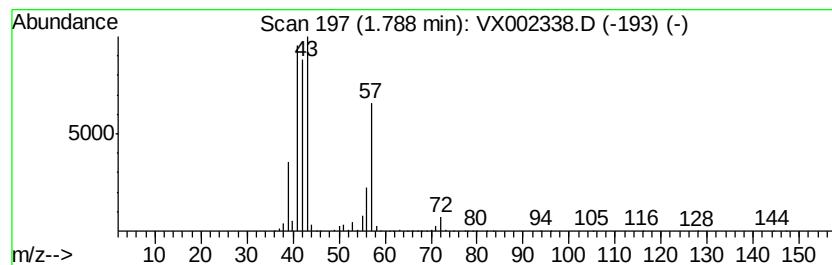
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.79	121.06 ug/l	2004290	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		3-Buten-1-ol	72	C4H8O	000627-27-0	47
3		Pentane	72	C5H12	000109-66-0	36
4		1-Propene, 2-methyl-	56	C4H8	000115-11-7	9
5		1-Butene	56	C4H8	000106-98-9	9



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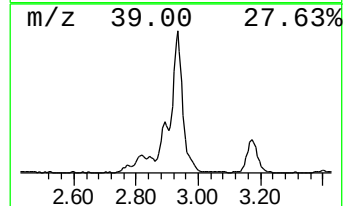
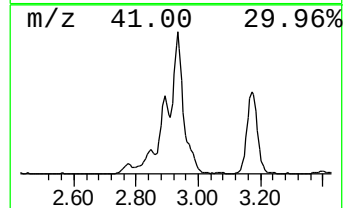
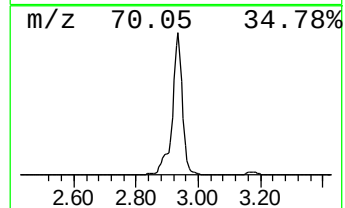
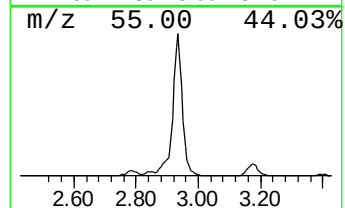
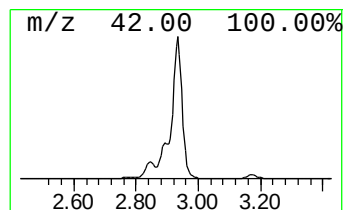
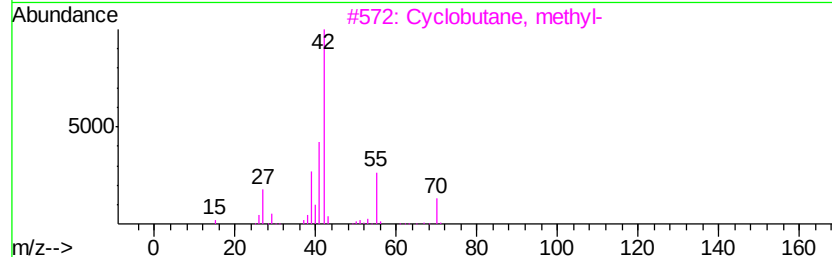
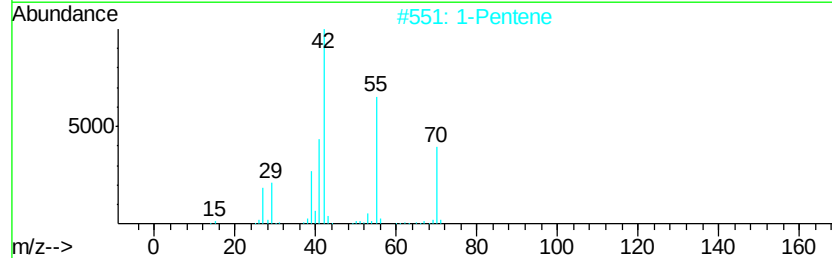
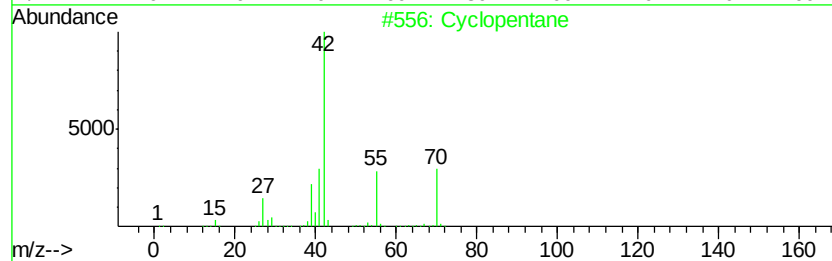
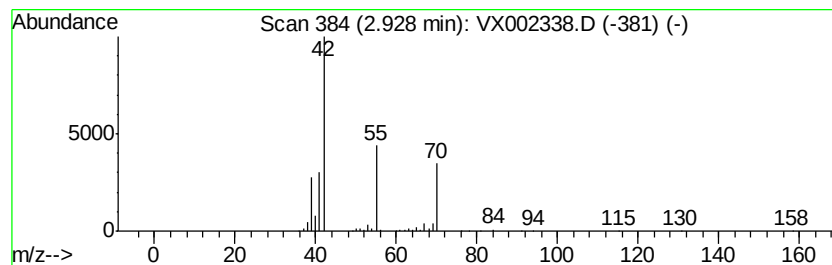
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	80.90 ug/l	1339400	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	86
2		1-Pentene	70	C5H10	000109-67-1	72
3		Cyclobutane, methyl-	70	C5H10	000598-61-8	56
4		3-Buten-1-ol	72	C4H8O	000627-27-0	36
5		Cyclobutanone	70	C4H6O	001191-95-3	9



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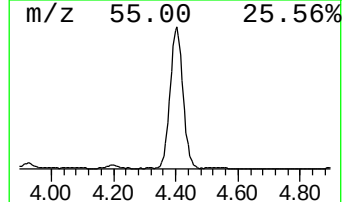
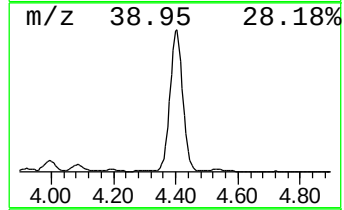
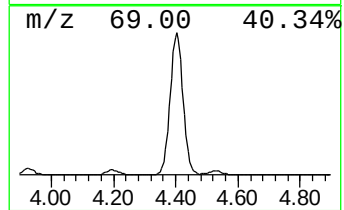
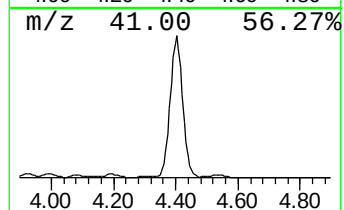
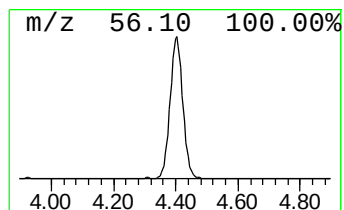
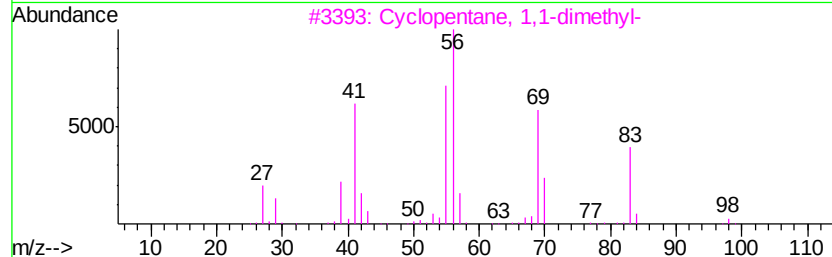
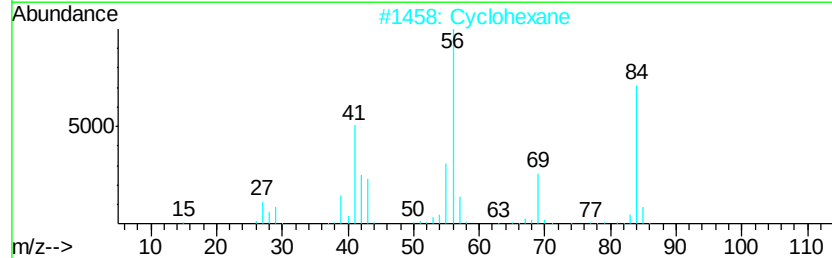
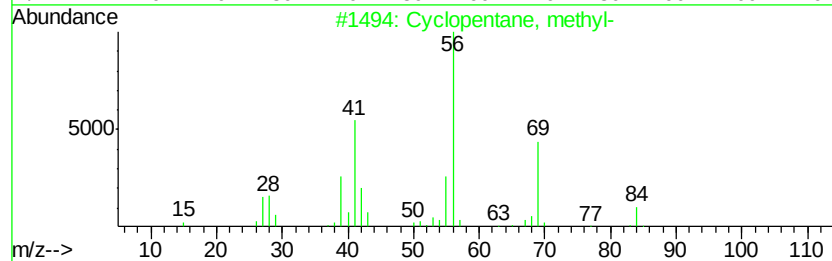
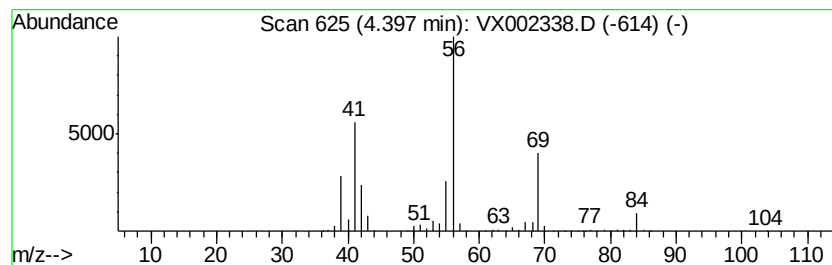
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	116.09 ug/l	1922020	Pentafluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclohexane	84	C6H12	000110-82-7	78
3			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	72
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5			Cyclobutane, ethyl-	84	C6H12	004806-61-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060518\
Data File : VX002338.D
Acq On : 06 Jun 2018 06:37
Operator : JC/MD
Sample : J3309-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

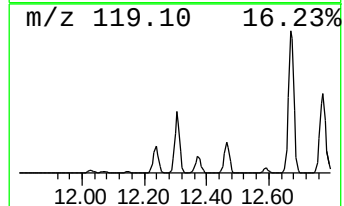
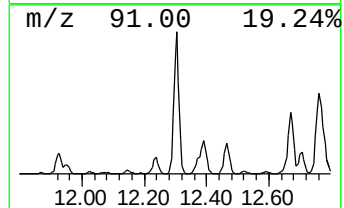
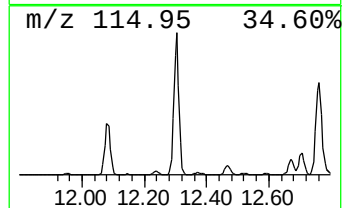
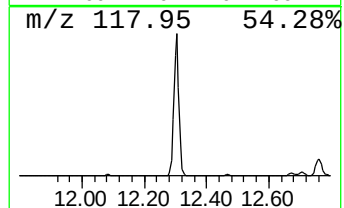
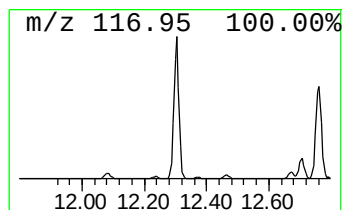
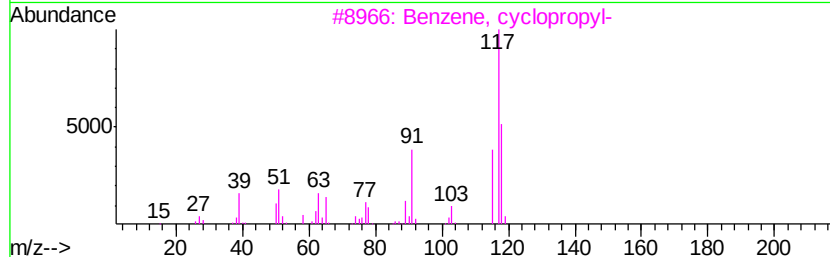
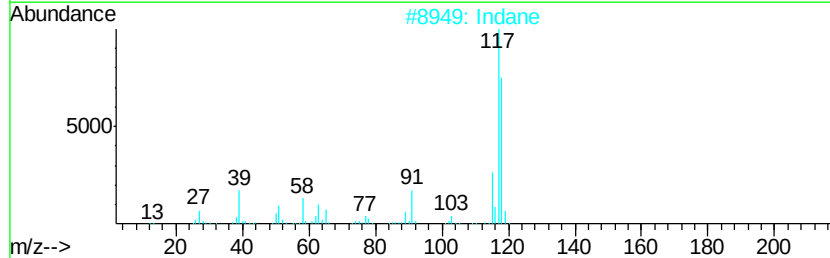
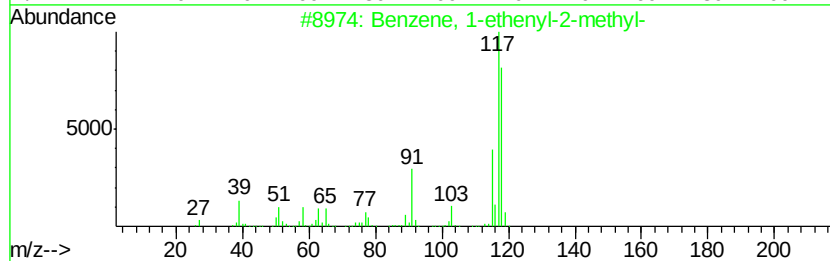
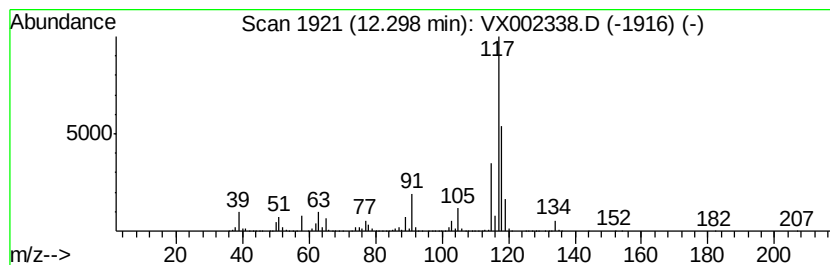
Instrument :
MSVOA_X
ClientSampleId :
MW-07

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 6 Benzene, 1-ethenyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	126.42 ug/l	2985830	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 86
2		Indane	118 C9H10	000496-11-7 76
3		Benzene, cyclopropyl-	118 C9H10	000873-49-4 70
4		Benzene, 2-propenyl-	118 C9H10	000300-57-2 70
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060518\
Data File : VX002338.D
Acq On : 06 Jun 2018 06:37
Operator : JC/MD
Sample : J3309-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-07

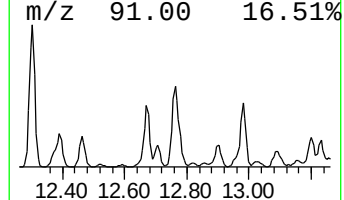
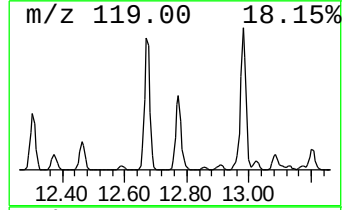
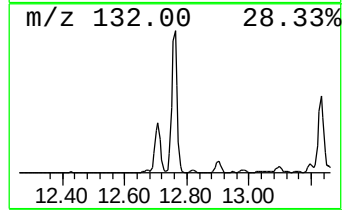
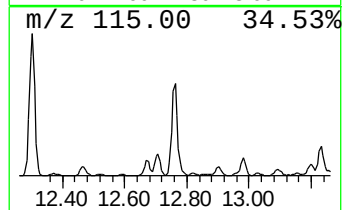
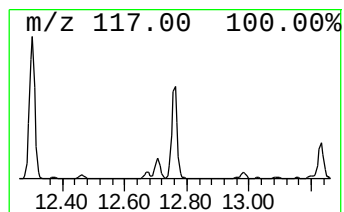
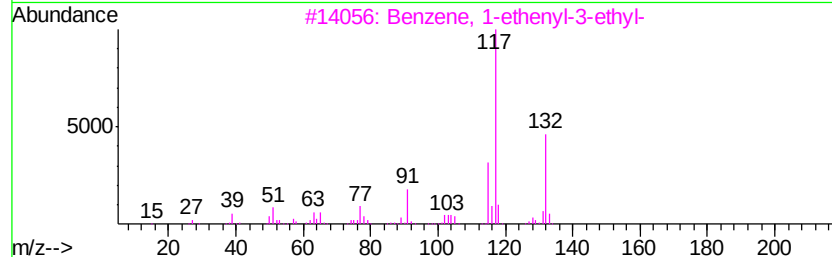
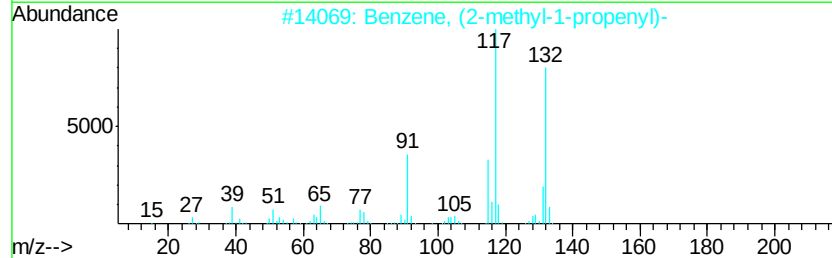
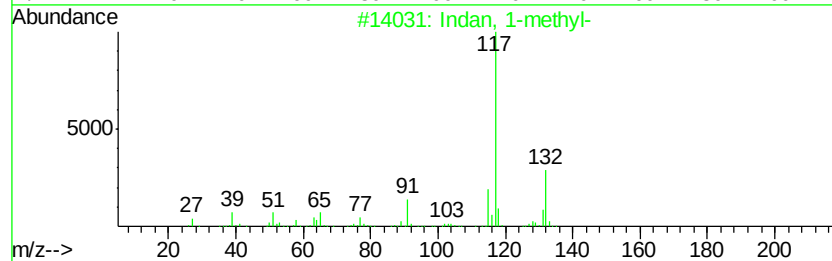
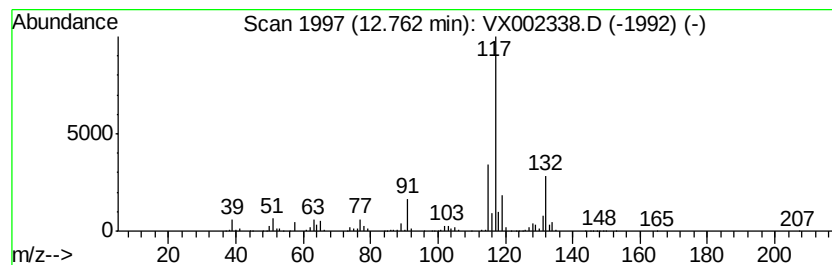
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	97.97 ug/l	2313850	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060518\
Data File : VX002338.D
Acq On : 06 Jun 2018 06:37
Operator : JC/MD
Sample : J3309-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-07

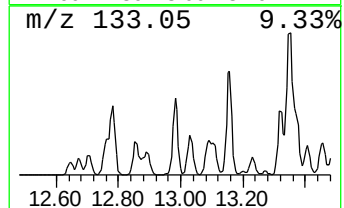
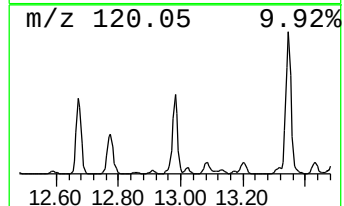
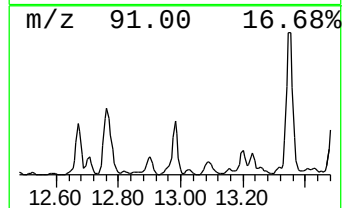
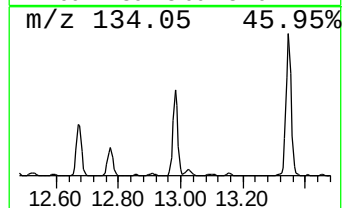
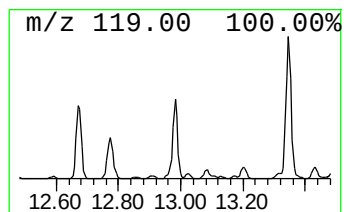
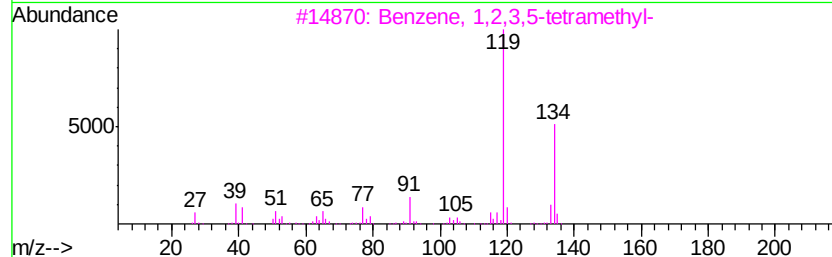
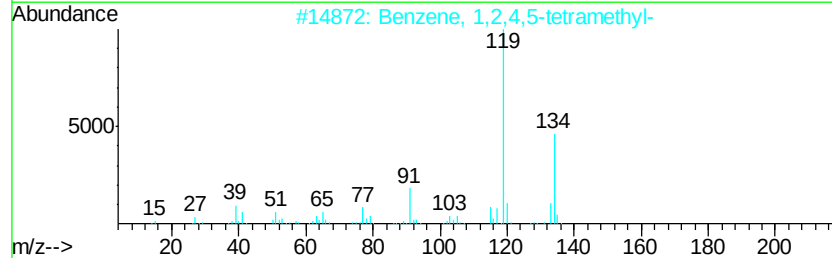
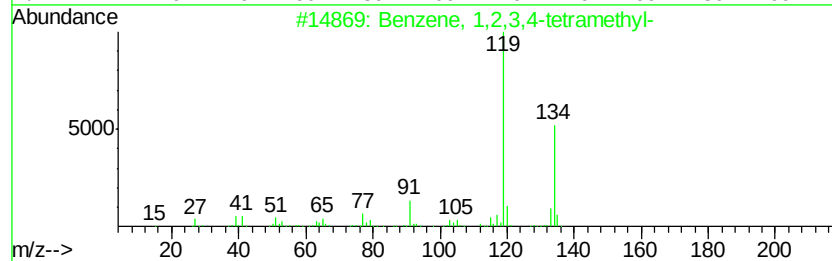
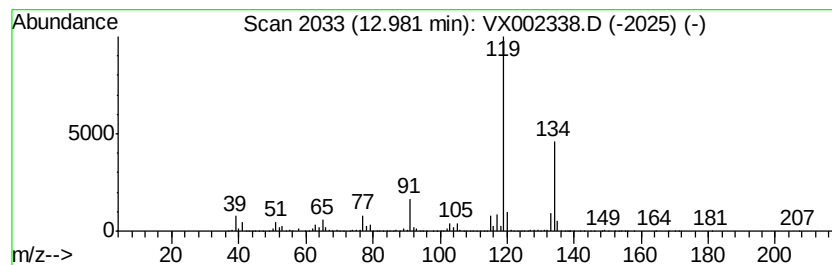
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	56.19 ug/l	1327240	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-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	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060518\
Data File : VX002338.D
Acq On : 06 Jun 2018 06:37
Operator : JC/MD
Sample : J3309-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-07

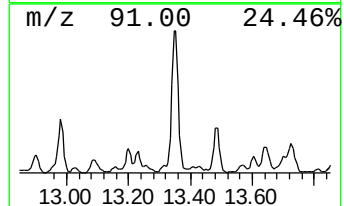
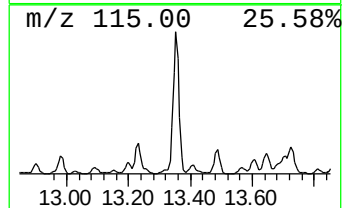
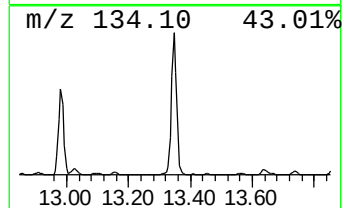
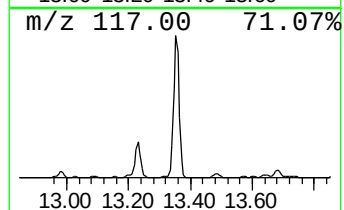
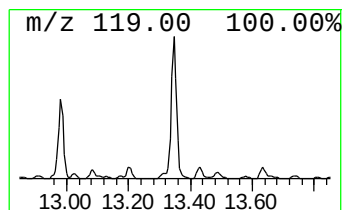
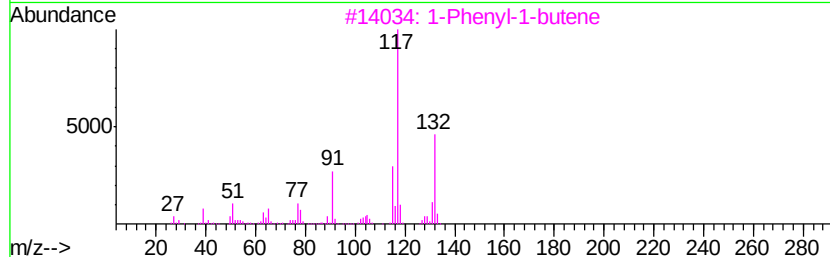
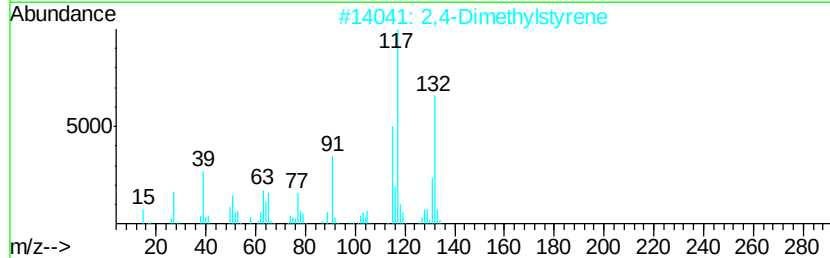
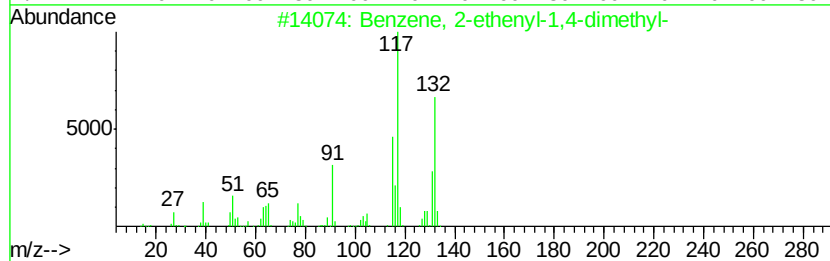
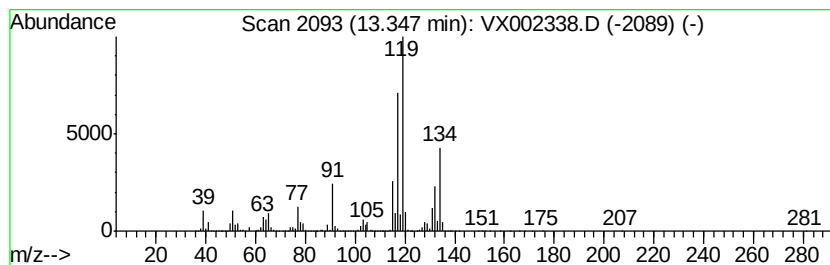
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	196.47 ug/l	4640440	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	95
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	92
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060518\
Data File : VX002338.D
Acq On : 06 Jun 2018 06:37
Operator : JC/MD
Sample : J3309-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-07

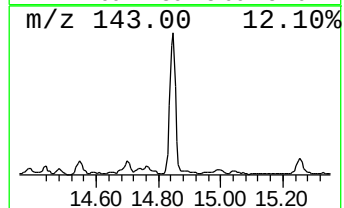
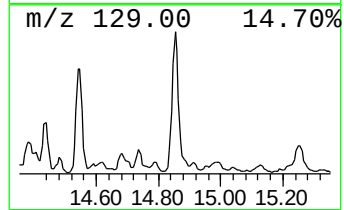
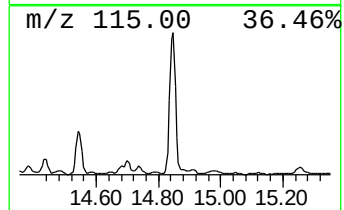
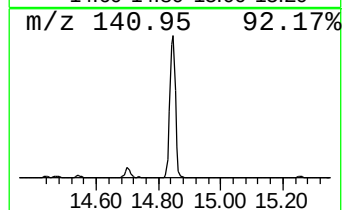
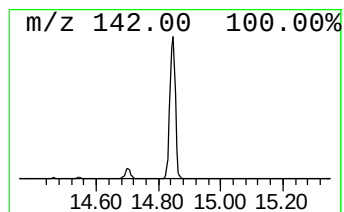
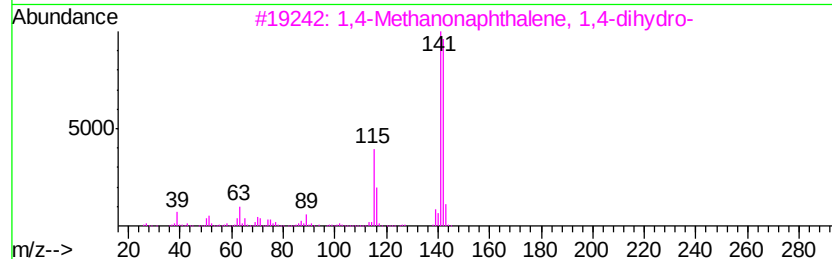
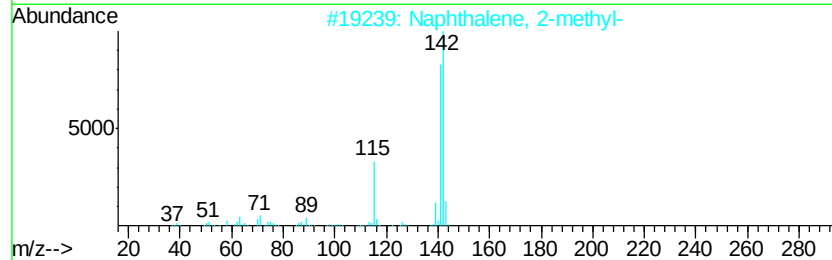
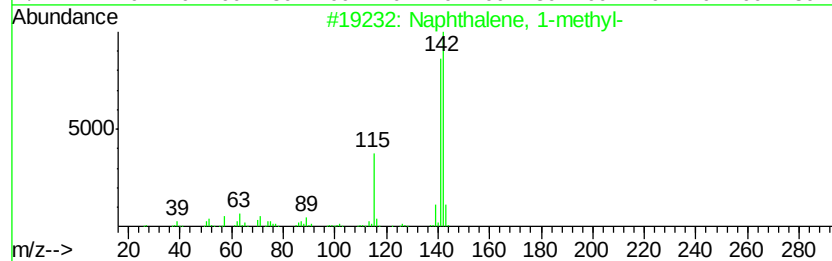
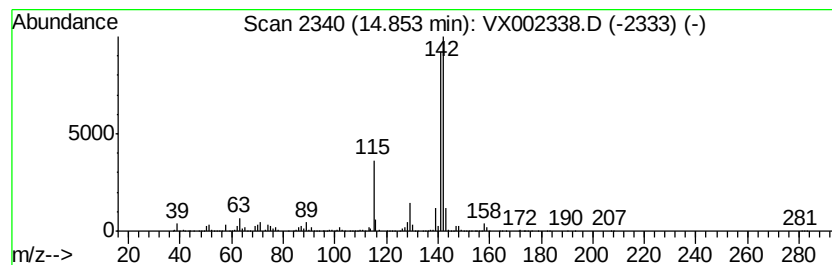
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	73.34 ug/l	1732120	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		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060518\
Data File : VX002338.D
Acq On : 06 Jun 2018 06:37
Operator : JC/MD
Sample : J3309-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-07

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
Butane, 2-methyl-	1.79	121.1	ug/l	2004290	1	5.67	827845	50.0
Cyclopentane	2.93	80.9	ug/l	1339400	1	5.67	827845	50.0
Cyclopentane, met...	4.40	116.1	ug/l	1922020	1	5.67	827845	50.0
Benzene, 1-etheny...	12.30	126.4	ug/l	2985830	4	12.09	1180930	50.0
Indan, 1-methyl-	12.76	98.0	ug/l	2313850	4	12.09	1180930	50.0
Benzene, 1,2,3,4-...	12.98	56.2	ug/l	1327240	4	12.09	1180930	50.0
Benzene, 2-etheny...	13.35	196.5	ug/l	4640440	4	12.09	1180930	50.0
Naphthalene, 1-me...	14.85	73.3	ug/l	1732120	4	12.09	1180930	50.0