

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120519\
 Data File : VX013773.D
 Acq On : 05 Dec 2019 13:50
 Operator : JC/SP
 Sample : K6174-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WATER-COMP

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\82X112619W.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	1.070	26	30	39	rBV	877574	1042666	6.58%	0.529%
2	1.783	142	147	157	rBV	2198946	2978304	18.80%	1.511%
3	1.991	176	181	194	rVB	2183138	3253477	20.54%	1.650%
4	2.259	220	225	235	rVB	205071	332646	2.10%	0.169%
5	2.838	313	320	322	rBV2	332109	591669	3.74%	0.300%
6	2.887	322	328	333	rBV	1638531	3085042	19.48%	1.565%
7	3.167	363	374	391	rBV	1407255	3180957	20.08%	1.613%
8	3.521	422	432	444	rVV	986103	2256954	14.25%	1.145%
9	3.814	471	480	490	rVB	214304	531190	3.35%	0.269%
10	4.399	565	576	589	rVV	3175648	9257432	58.45%	4.696%
11	5.295	714	723	734	rVB2	279631	856108	5.41%	0.434%
12	5.490	741	755	760	rBV	368587	1119080	7.07%	0.568%
13	5.575	760	769	778	rBV2	3083640	9728931	61.42%	4.935%
14	5.929	813	827	839	rBV5	917303	3030569	19.13%	1.537%
15	6.057	839	848	855	rVV	383409	998674	6.31%	0.507%
16	6.258	870	881	890	rVV3	797320	2879056	18.18%	1.460%
17	6.356	890	897	904	rVV	753587	2095105	13.23%	1.063%
18	6.447	904	912	925	rVB	1263382	3533496	22.31%	1.792%
19	6.697	942	953	965	rBV	542490	1346887	8.50%	0.683%
20	6.850	968	978	989	rBV2	980391	3058709	19.31%	1.551%
21	7.252	1036	1044	1052	rVB	222355	475586	3.00%	0.241%
22	7.459	1064	1078	1088	rBV	6927853	15838795	100.00%	8.034%
23	7.728	1115	1122	1129	rBV	811637	1568554	9.90%	0.796%
24	7.843	1135	1141	1148	rVB	332124	649533	4.10%	0.329%
25	7.929	1148	1155	1156	rBV2	267187	463816	2.93%	0.235%
26	7.947	1156	1158	1165	rVV	344189	619103	3.91%	0.314%
27	8.020	1165	1170	1176	rVV	479189	927337	5.85%	0.470%
28	8.209	1195	1201	1205	rVV	549393	957859	6.05%	0.486%
29	8.258	1205	1209	1217	rVV3	354398	713992	4.51%	0.362%
30	8.337	1217	1222	1225	rVV	379739	635313	4.01%	0.322%
31	8.380	1225	1229	1236	rVB4	311886	620402	3.92%	0.315%
32	8.496	1243	1248	1252	rBV	274017	435244	2.75%	0.221%
33	8.569	1254	1260	1268	rVB	1071749	1897882	11.98%	0.963%
34	8.660	1268	1275	1278	rBV2	1272287	2180154	13.76%	1.106%

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35	8.709	1278	1283	1291	rVB2	2448409	4284020	27.05%	2.173%
36	8.831	1298	1303	1308	rBV	520028	901854	5.69%	0.457%
37	8.904	1312	1315	1321	rVB	398014	657519	4.15%	0.334%
38	8.977	1321	1327	1331	rBV3	272500	486372	3.07%	0.247%
39	9.038	1331	1337	1342	rVV	903144	1492819	9.43%	0.757%
40	9.160	1349	1357	1362	rVB2	547120	944543	5.96%	0.479%
41	9.215	1362	1366	1372	rBV2	443110	711687	4.49%	0.361%
42	9.483	1406	1410	1413	rBV	280552	433108	2.73%	0.220%
43	9.514	1413	1415	1418	rVV	284587	358466	2.26%	0.182%
44	9.569	1418	1424	1428	rVV4	479336	972826	6.14%	0.493%
45	9.617	1428	1432	1436	rVV	1284888	1817094	11.47%	0.922%
46	9.672	1436	1441	1445	rVV2	691032	1157217	7.31%	0.587%
47	9.813	1459	1464	1470	rBV3	276413	448114	2.83%	0.227%
48	9.886	1470	1476	1479	rBV2	174975	341151	2.15%	0.173%
49	10.111	1508	1513	1517	rVV	1925130	2709606	17.11%	1.374%
50	10.178	1521	1524	1528	rVV	472543	778914	4.92%	0.395%
51	10.245	1531	1535	1541	rVV	1266069	1810984	11.43%	0.919%
52	10.331	1544	1549	1550	rVV2	309052	468772	2.96%	0.238%
53	10.355	1550	1553	1559	rVV2	407340	835645	5.28%	0.424%
54	10.636	1592	1599	1607	rVV3	260994	584462	3.69%	0.296%
55	10.831	1621	1631	1639	rBV3	447986	884167	5.58%	0.448%
56	10.910	1639	1644	1652	rVV4	364879	672982	4.25%	0.341%
57	11.014	1656	1661	1665	rVV	3399878	4264436	26.92%	2.163%
58	11.135	1675	1681	1685	rBV	1675374	2200594	13.89%	1.116%
59	11.355	1712	1717	1724	rBV	3569398	4283757	27.05%	2.173%
60	11.452	1727	1733	1736	rVV	678965	914160	5.77%	0.464%
61	11.666	1763	1768	1776	rBV	924048	1206353	7.62%	0.612%
62	11.916	1803	1809	1810	rBV	372929	448353	2.83%	0.227%
63	11.940	1810	1813	1818	rVB	835029	1036375	6.54%	0.526%
64	12.019	1818	1826	1829	rBV	304529	430096	2.72%	0.218%
65	12.074	1829	1835	1840	rVB2	2238180	3334867	21.06%	1.692%
66	12.135	1840	1845	1849	rBV	733482	919762	5.81%	0.467%
67	12.227	1856	1860	1864	rVB	404777	476304	3.01%	0.242%
68	12.294	1864	1871	1878	rBV	4934214	6301065	39.78%	3.196%
69	12.367	1878	1883	1890	rVB2	1339987	2417383	15.26%	1.226%
70	12.458	1893	1898	1902	rVB	753218	958472	6.05%	0.486%
71	12.513	1903	1907	1914	rBV	633121	743798	4.70%	0.377%

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72	12.580	1914	1918	1925	rVB	553944	687196	4.34%	0.349%
73	12.666	1925	1932	1935	rBV	4370470	5399479	34.09%	2.739%
74	12.696	1935	1937	1941	rVB	990380	1021117	6.45%	0.518%
75	12.751	1941	1946	1954	rBV2	4788522	7078759	44.69%	3.591%
76	12.897	1965	1970	1975	rVB3	614376	1020741	6.44%	0.518%
77	12.970	1975	1982	1987	rBV	2594174	3396693	21.45%	1.723%
78	13.019	1987	1990	1995	rVB3	294885	410399	2.59%	0.208%
79	13.080	1995	2000	2005	rBV4	598288	1218016	7.69%	0.618%
80	13.190	2015	2018	2020	rBV2	402430	485440	3.06%	0.246%
81	13.220	2020	2023	2027	rVB	2392622	2650871	16.74%	1.345%
82	13.342	2038	2043	2049	rVB2	8364420	11330187	71.53%	5.747%
83	13.477	2061	2065	2072	rVB	1861422	2344441	14.80%	1.189%
84	13.556	2073	2078	2081	rBV2	478117	670873	4.24%	0.340%
85	13.598	2081	2085	2088	rVV	1213446	1676062	10.58%	0.850%
86	13.635	2088	2091	2096	rVV4	1133863	1802209	11.38%	0.914%
87	13.696	2096	2101	2102	rVV3	666342	1027962	6.49%	0.521%
88	13.714	2102	2104	2110	rVB2	1452120	2062989	13.02%	1.046%
89	13.830	2115	2123	2132	rVB	3144422	4502898	28.43%	2.284%
90	13.983	2140	2148	2152	rBV3	560792	944928	5.97%	0.479%
91	14.025	2152	2155	2159	rVB	389587	432584	2.73%	0.219%
92	14.129	2167	2172	2175	rBV	536944	669665	4.23%	0.340%
93	14.269	2190	2195	2205	rBV2	677598	1350652	8.53%	0.685%
94	14.373	2208	2212	2218	rVV	304217	513818	3.24%	0.261%
95	14.428	2218	2221	2225	rVB	410969	484858	3.06%	0.246%
96	14.537	2234	2239	2243	rBV2	309203	430563	2.72%	0.218%
97	14.690	2260	2264	2274	rVB	2540203	3318814	20.95%	1.683%
98	14.836	2282	2288	2297	rBV	1937004	2558727	16.15%	1.298%
99	15.543	2397	2404	2416	rVV	226697	387192	2.44%	0.196%
100	15.683	2418	2427	2430	rVV	270936	440856	2.78%	0.224%

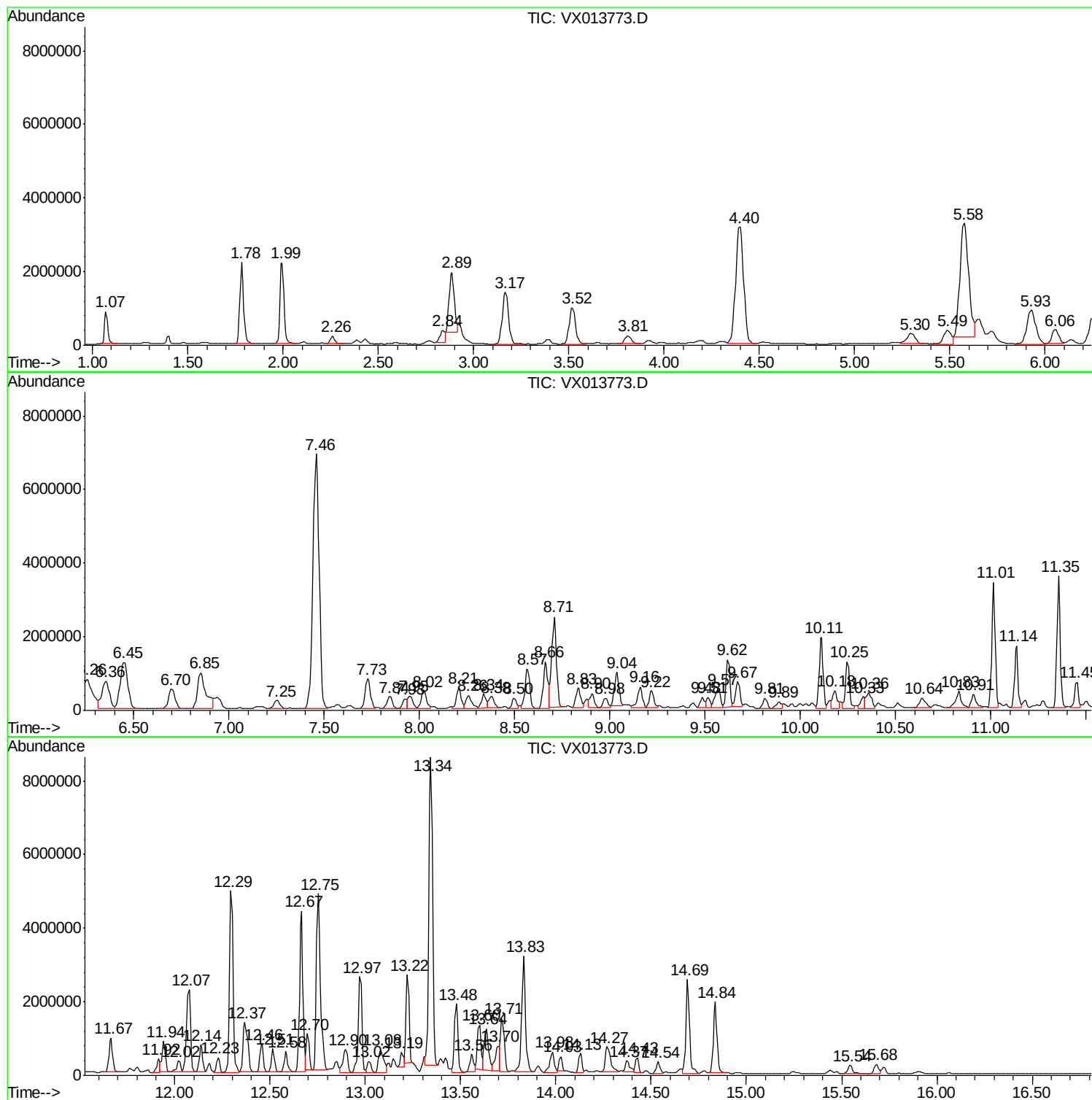
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Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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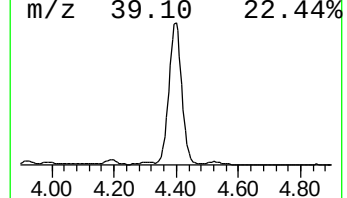
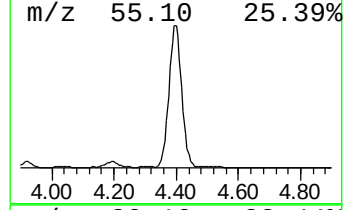
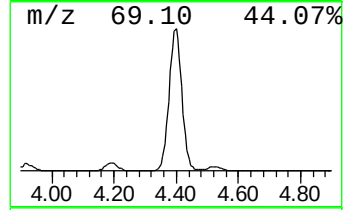
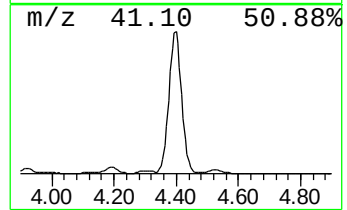
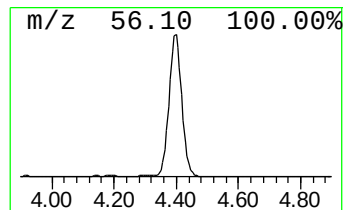
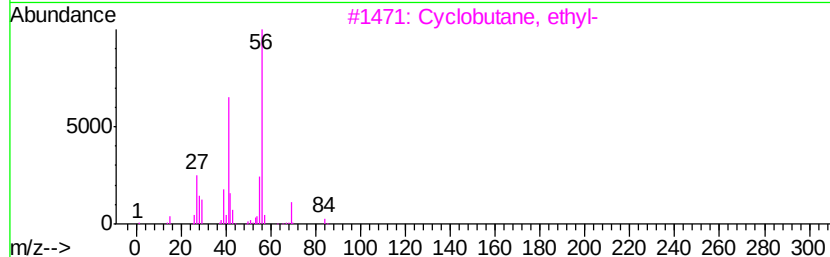
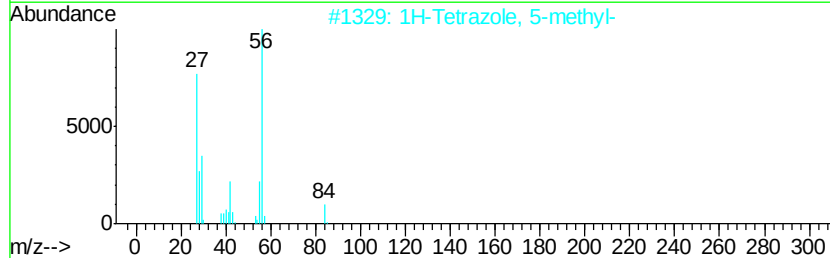
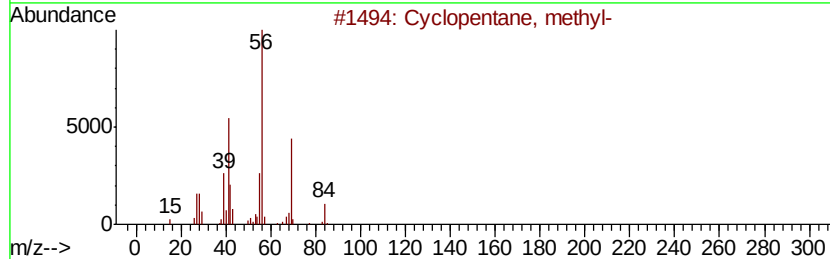
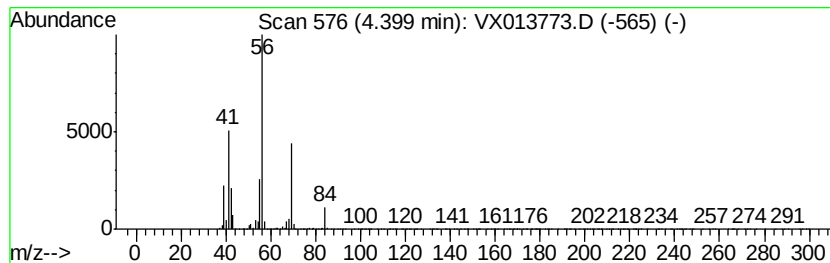
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	47.58 ug/l	9257430	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		Cyclohexane	84	C6H12	000110-82-7	56
5		Cyclopropane, propyl-	84	C6H12	002415-72-7	53



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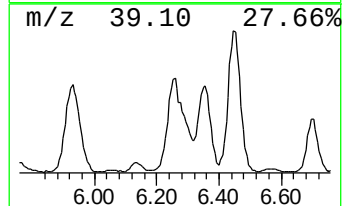
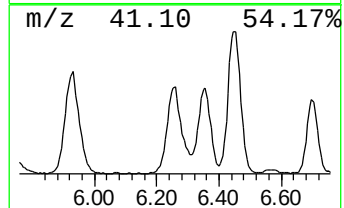
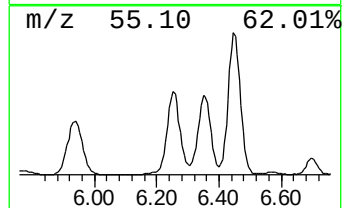
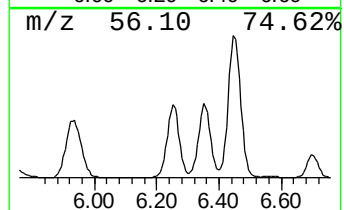
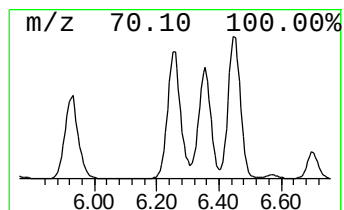
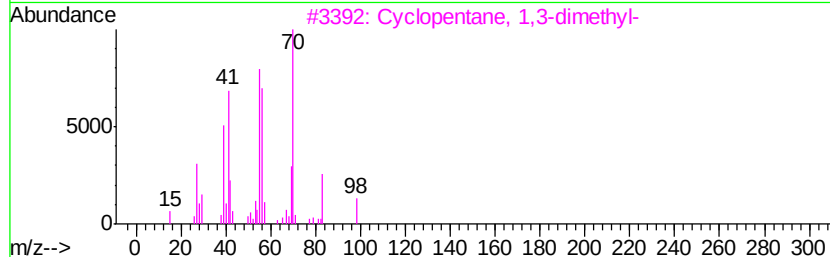
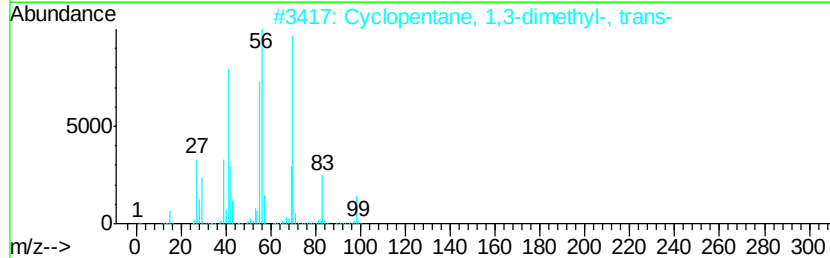
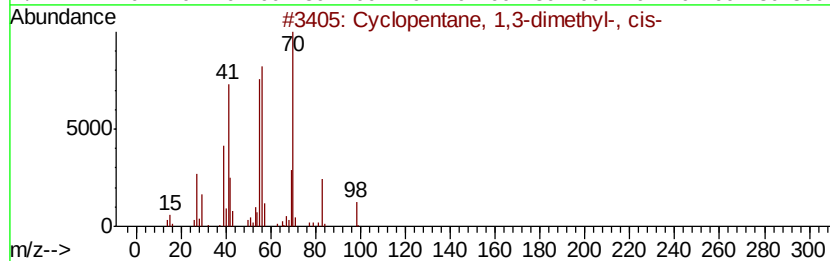
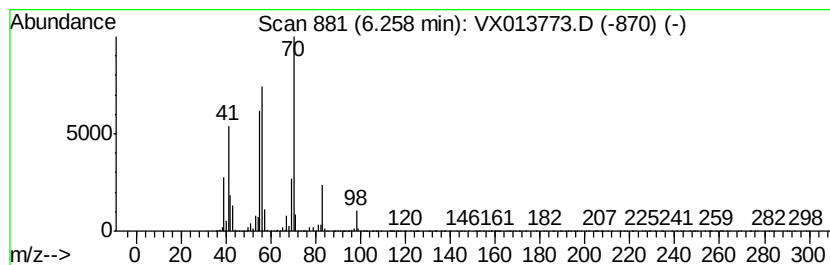
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Peak Number 2 Cyclopentane, 1,3-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.26	47.06 ug/l	2879060	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	94
2		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
3		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	87
4		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	76
5		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	49



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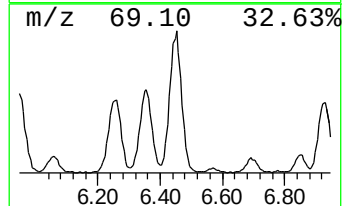
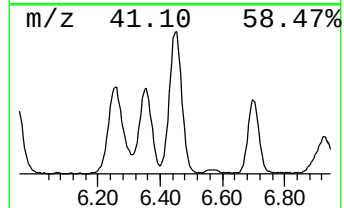
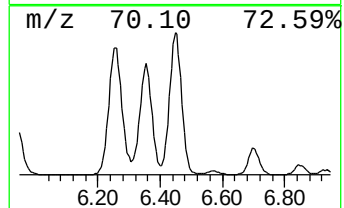
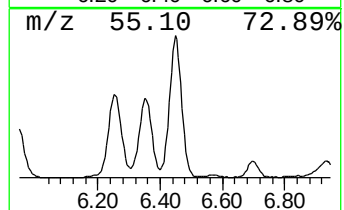
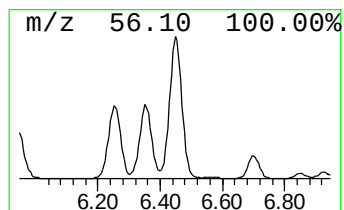
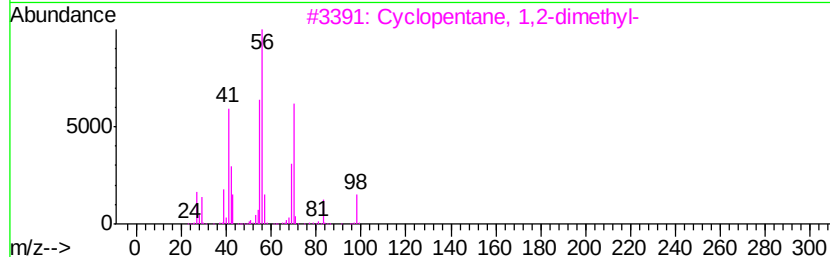
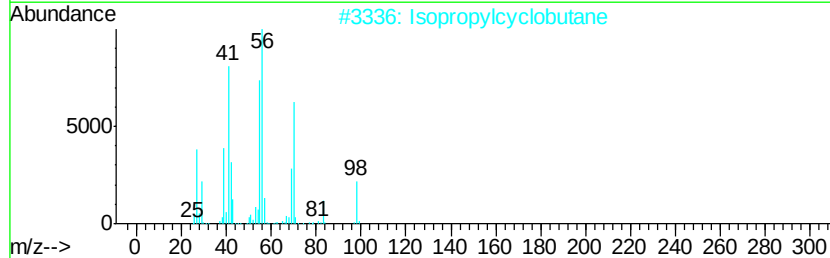
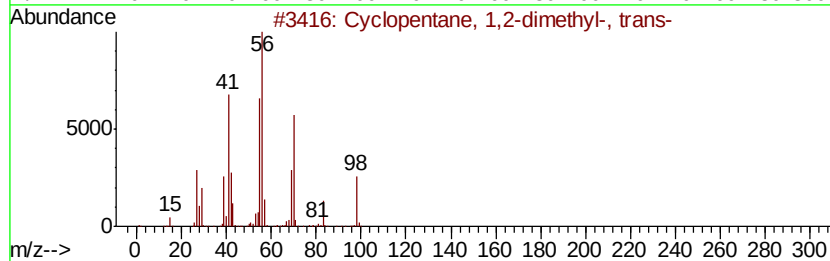
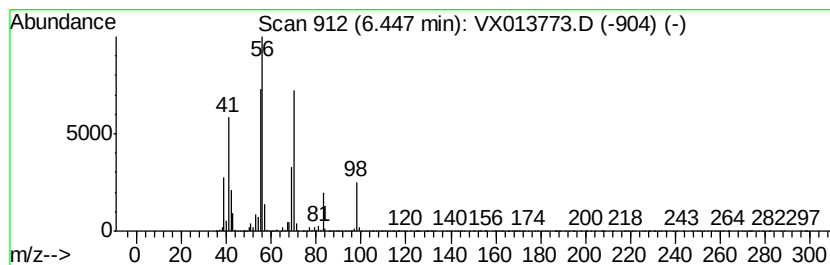
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentane, 1,2-dimethyl-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	57.76 ug/l	3533500	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	94
2		Isopropylcyclobutane	98	C7H14	000872-56-0	94
3		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
4		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	86
5		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

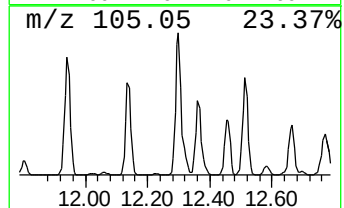
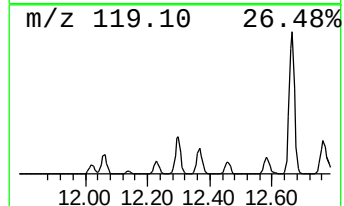
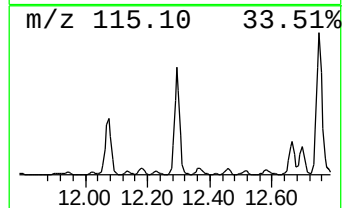
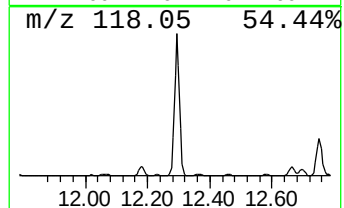
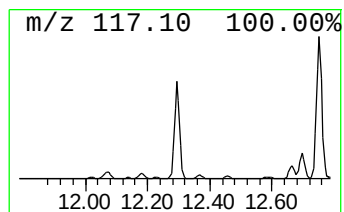
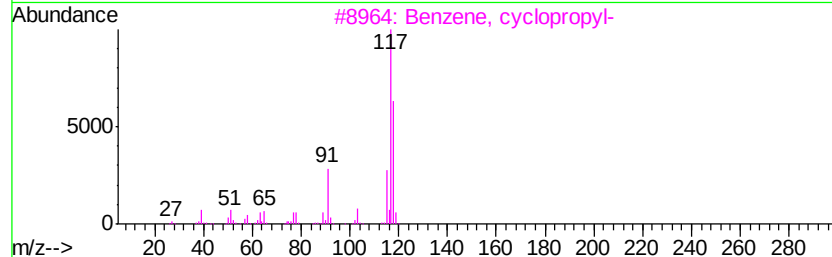
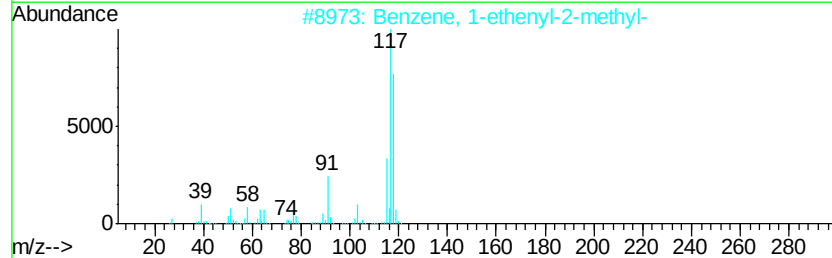
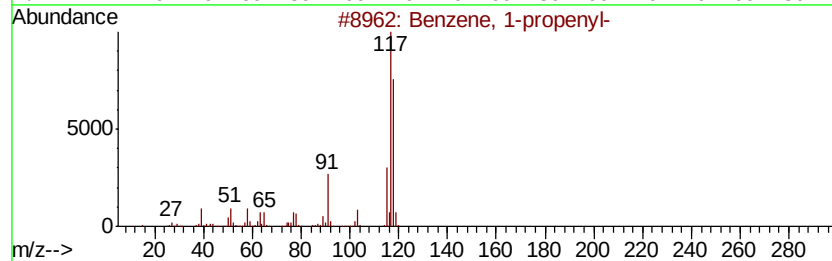
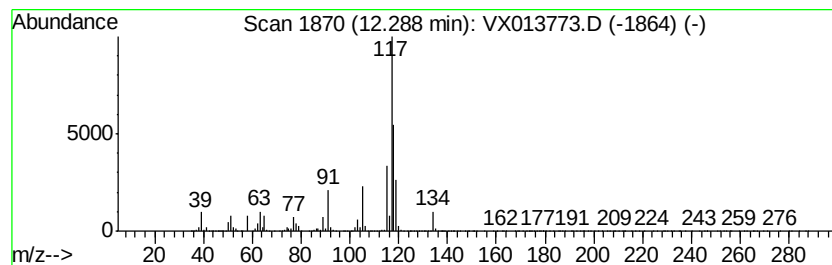
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-propenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	94.47 ug/l	6301070	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	90
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	87
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
5		Indane	118	C9H10	000496-11-7	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

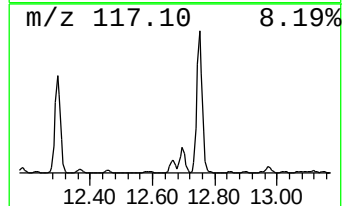
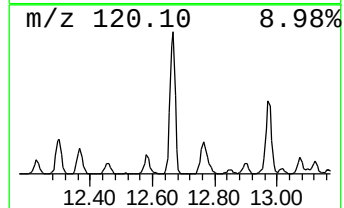
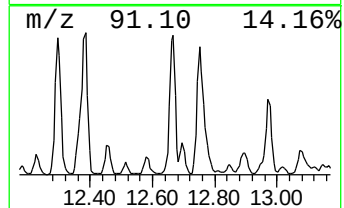
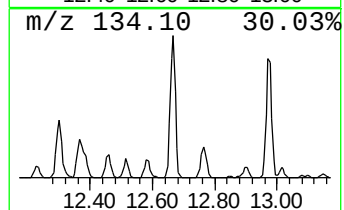
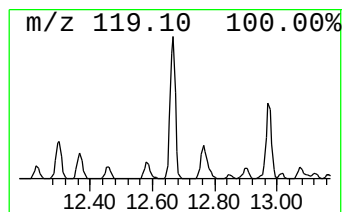
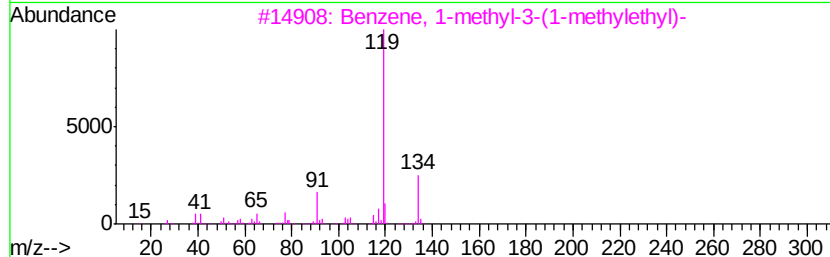
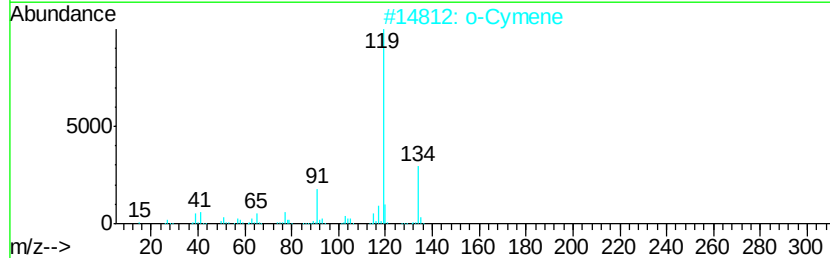
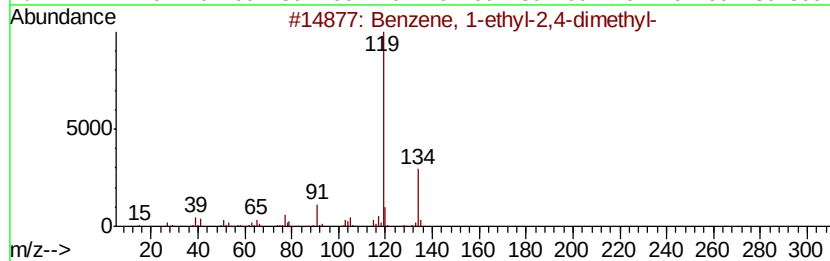
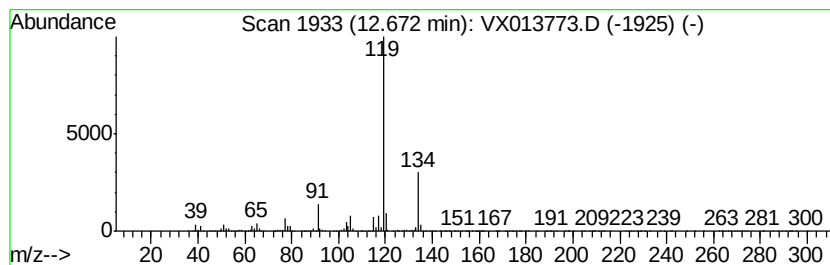
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	80.95 ug/l	5399480	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

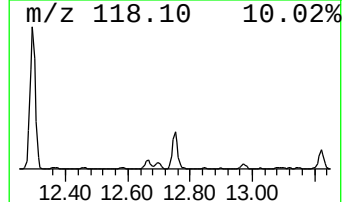
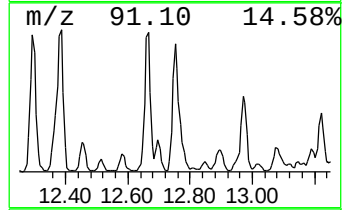
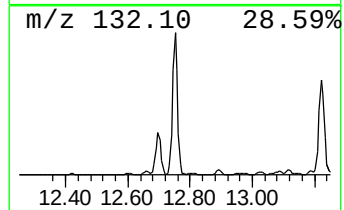
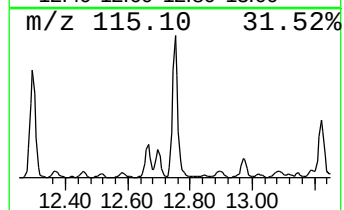
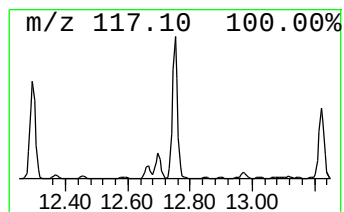
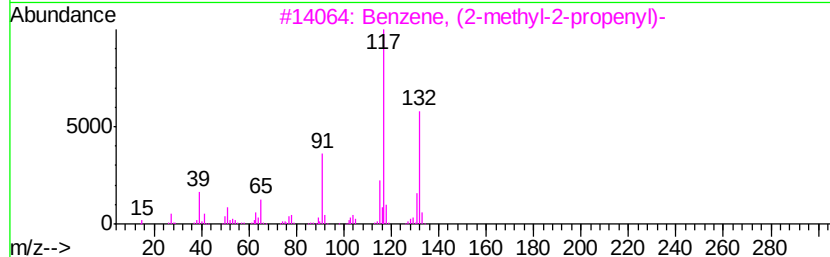
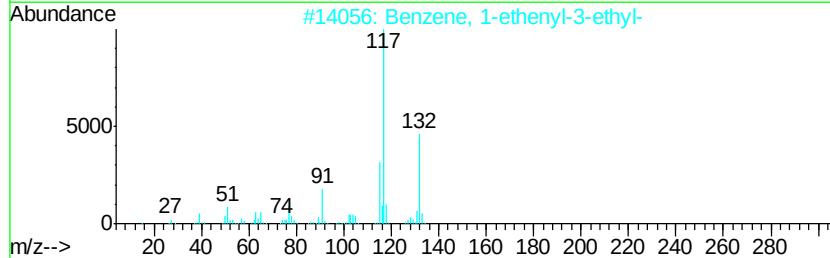
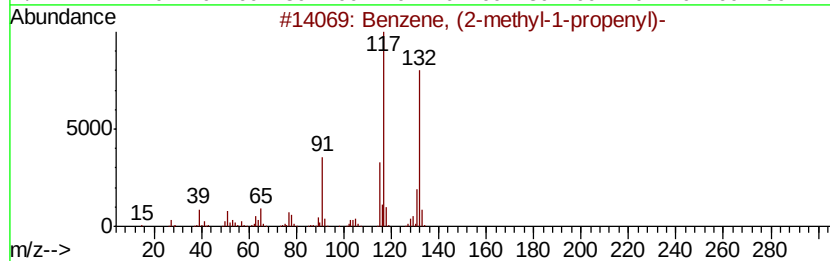
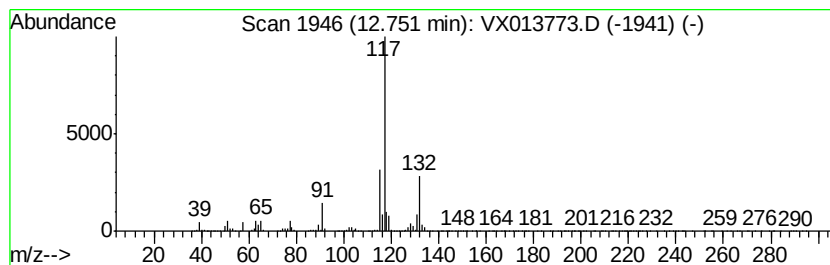
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-propen... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	106.13 ug/l	7078760	1,4-Dichlorobenzene-d4	12.07

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5		Indan, 1-methyl-	132	C10H12	000767-58-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

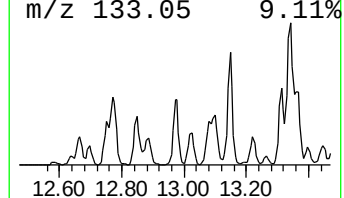
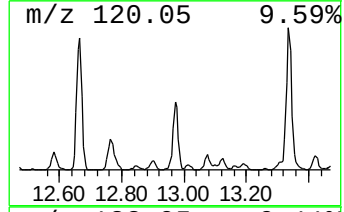
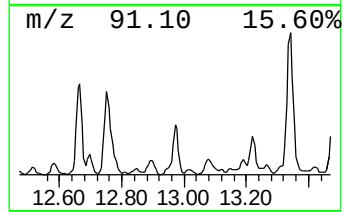
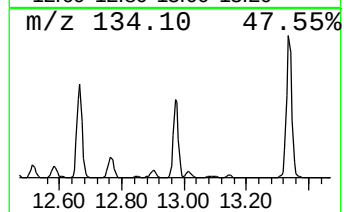
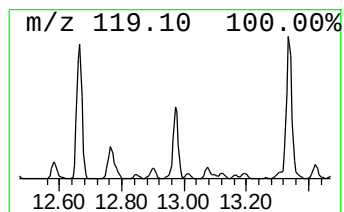
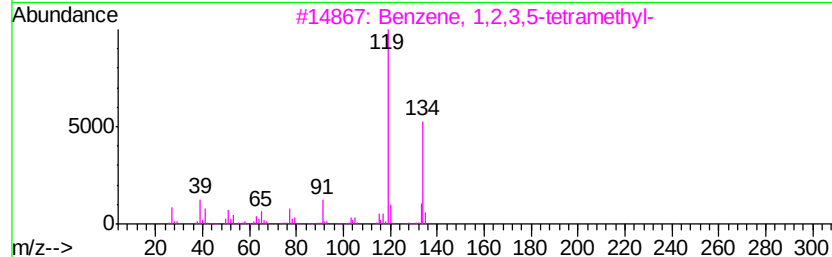
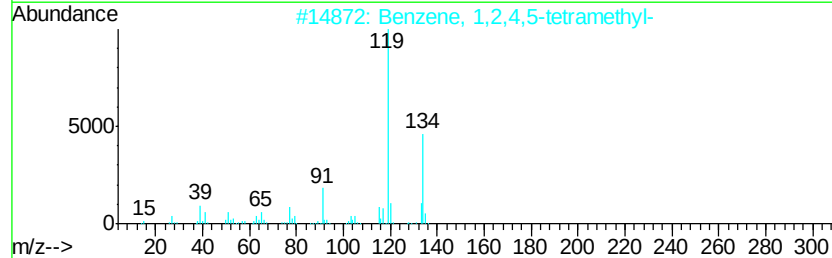
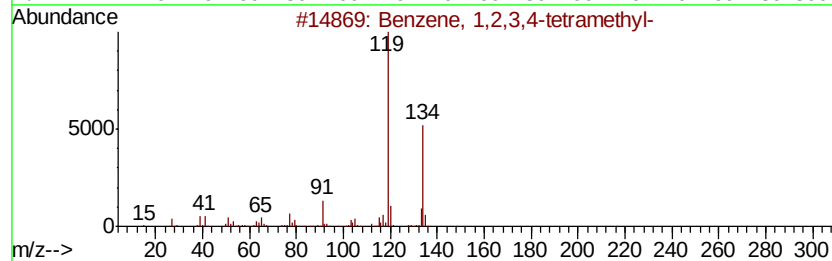
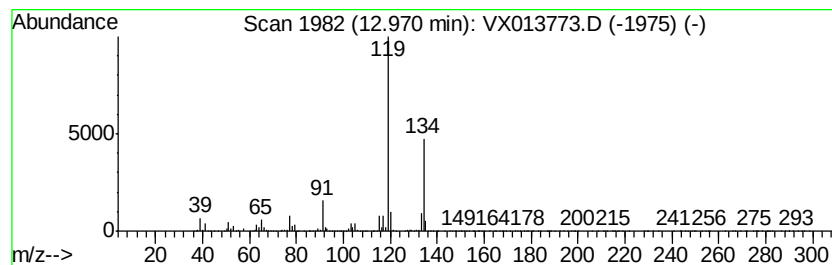
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	50.93 ug/l	3396690	1,4-Dichlorobenzene-d4	12.07

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	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
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\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

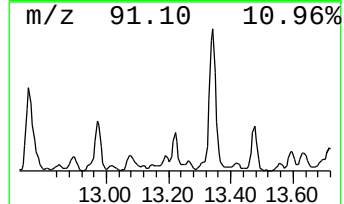
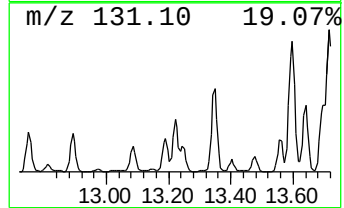
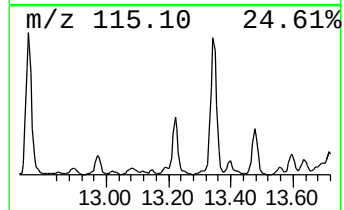
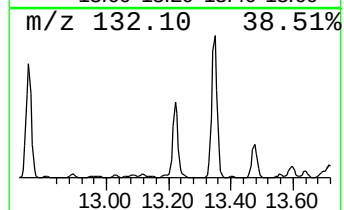
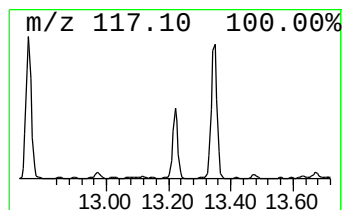
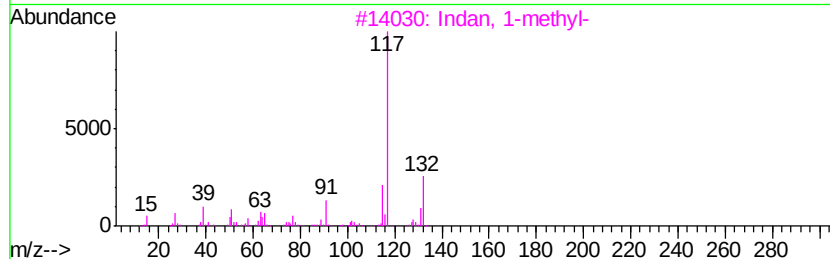
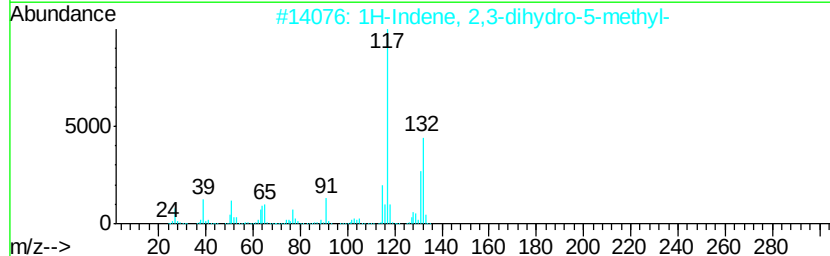
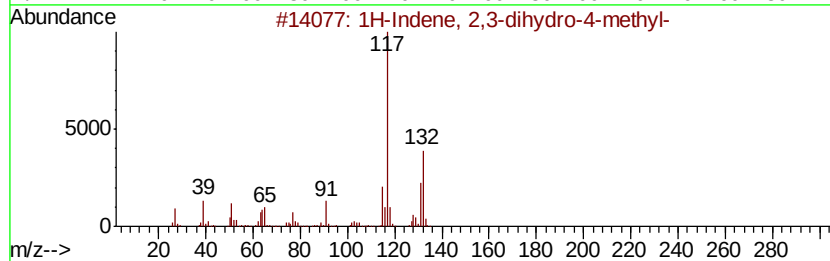
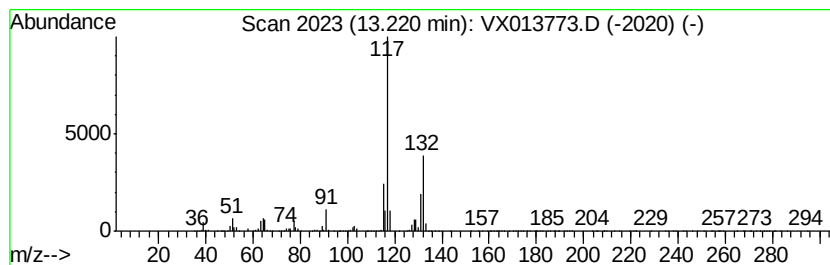
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	39.74 ug/l	2650870	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3		Indan, 1-methyl-	132	C10H12	000767-58-8	90
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

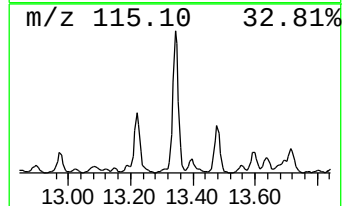
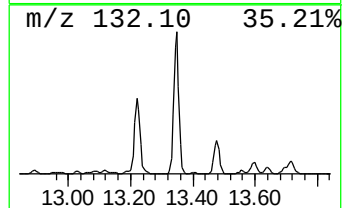
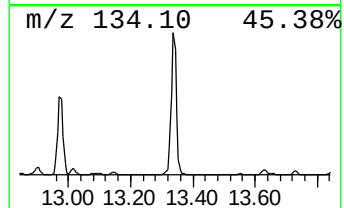
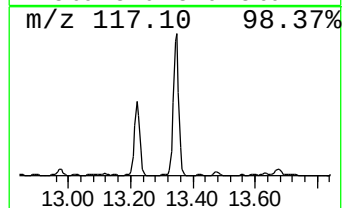
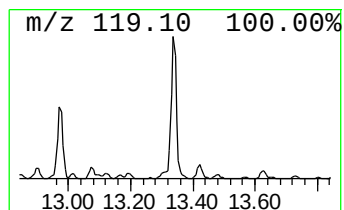
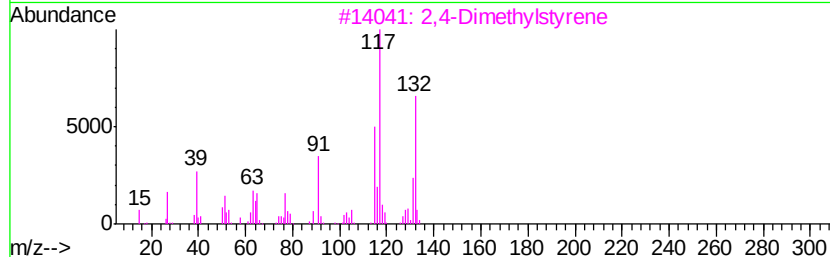
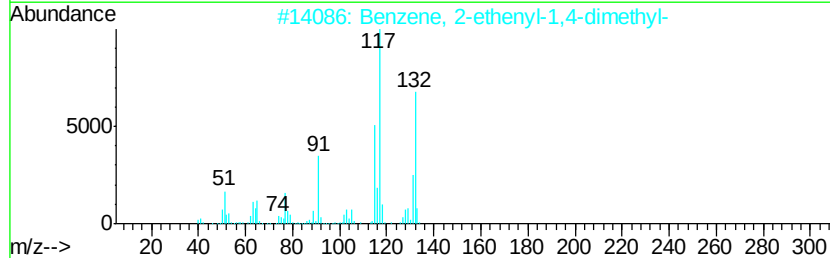
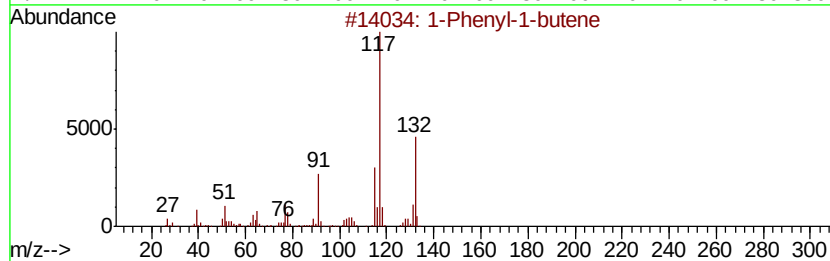
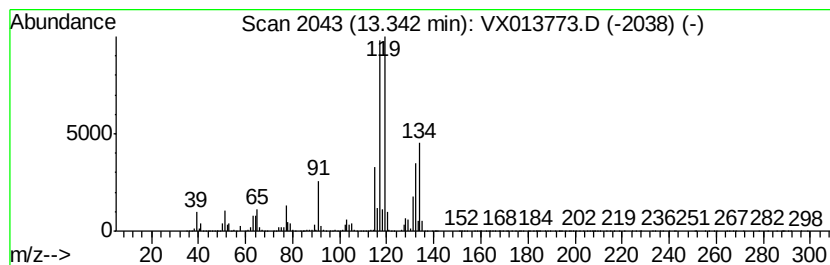
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Quant Title : SW846 8260

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	169.88 ug/l	11330200	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	92
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

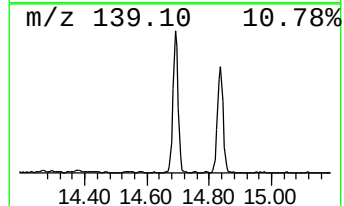
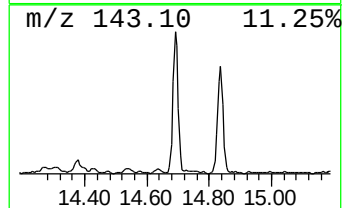
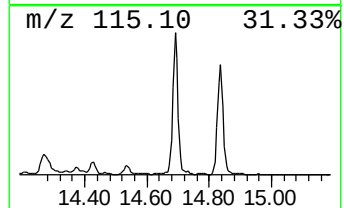
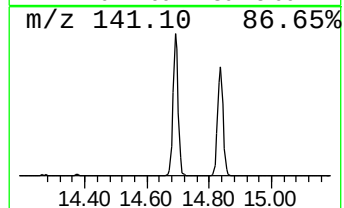
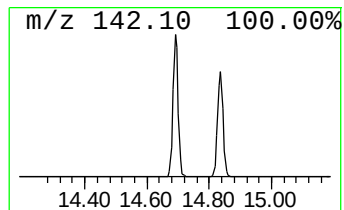
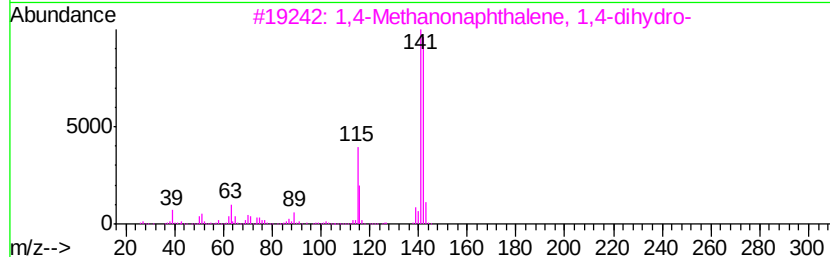
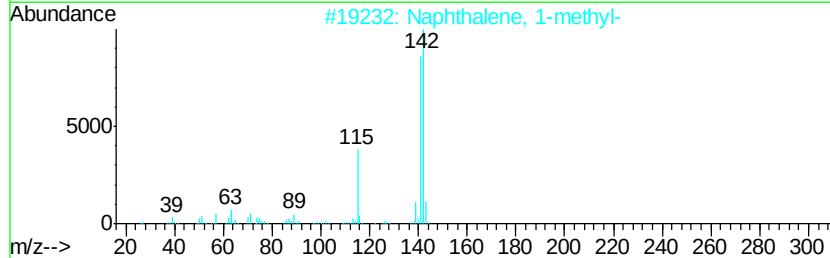
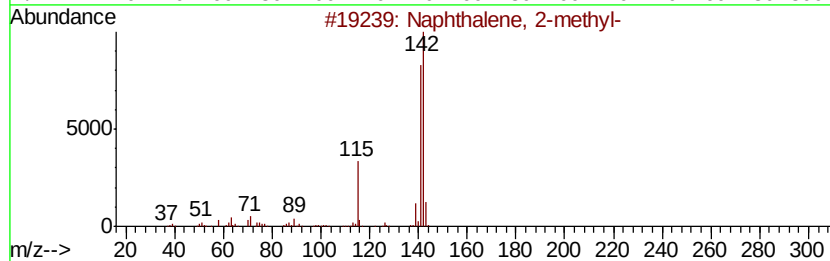
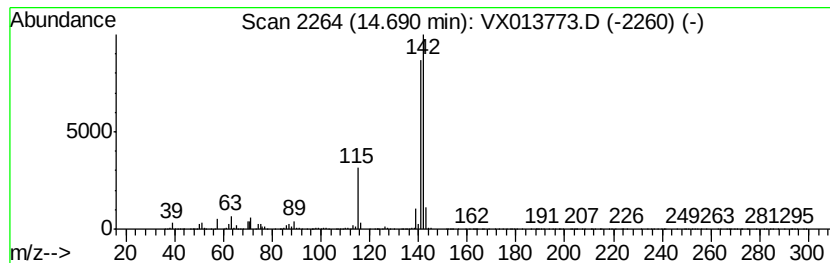
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	49.76 ug/l	3318810	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX120519\
Data File : VX013773.D
Acq On : 05 Dec 2019 13:50
Operator : JC/SP
Sample : K6174-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WATER-COMP

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.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
Cyclopentane, met...	4.40	47.6	ug/l	9257430	1	5.65	9728930	50.0
Cyclopentane, 1,3...	6.26	47.1	ug/l	2879060	2	6.85	3058710	50.0
Cyclopentane, 1,2...	6.45	57.8	ug/l	3533500	2	6.85	3058710	50.0
Benzene, 1-propenyl-	12.29	94.5	ug/l	6301070	4	12.07	3334870	50.0
Benzene, 1-ethyl-...	12.67	81.0	ug/l	5399480	4	12.07	3334870	50.0
Benzene, (2-methy...	12.75	106.1	ug/l	7078760	4	12.07	3334870	50.0
Benzene, 1,2,3,4-...	12.97	50.9	ug/l	3396690	4	12.07	3334870	50.0
1H-Indene, 2,3-di...	13.22	39.7	ug/l	2650870	4	12.07	3334870	50.0
1-Phenyl-1-butene	13.34	169.9	ug/l	11330200	4	12.07	3334870	50.0
Naphthalene, 1-me...	14.69	49.8	ug/l	3318810	4	12.07	3334870	50.0