

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021119\
 Data File : VX007464.D
 Acq On : 11 Feb 2019 13:33
 Operator : JC/SP
 Sample : K1433-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY001MW038-190207

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\82X020119W.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	2.843	266	277	281	rBV2	141164	370249	1.24%	0.187%
2	2.891	281	285	290	rVV	250383	439621	1.47%	0.222%
3	3.202	322	336	350	rVB3	761185	2625101	8.79%	1.325%
4	3.519	381	388	403	rVB	168582	403060	1.35%	0.203%
5	3.812	428	436	446	rBV	180715	439028	1.47%	0.222%
6	3.922	446	454	461	rBV2	126498	335505	1.12%	0.169%
7	4.306	507	517	523	rVV	140447	407599	1.36%	0.206%
8	4.397	523	532	545	rVV	1215768	3547944	11.87%	1.791%
9	5.306	655	681	694	rBV3	314147	1194426	4.00%	0.603%
10	5.507	698	714	718	rBV	328980	985025	3.30%	0.497%
11	5.580	718	726	735	rVV3	848223	2713920	9.08%	1.370%
12	5.665	735	740	745	rVV	405097	1081960	3.62%	0.546%
13	5.726	745	750	771	rVB2	636851	2053136	6.87%	1.037%
14	5.927	771	783	798	rBV4	393870	1221796	4.09%	0.617%
15	6.068	798	806	810	rBV	299998	731692	2.45%	0.369%
16	6.147	810	819	830	rVB	6254659	15832092	52.99%	7.993%
17	6.263	830	838	841	rBV	308106	805224	2.69%	0.407%
18	6.360	846	854	863	rVV5	1223973	4403265	14.74%	2.223%
19	6.458	863	870	882	rVB2	348825	1003693	3.36%	0.507%
20	6.702	897	910	927	rVB2	160414	411465	1.38%	0.208%
21	6.860	927	936	943	rBV2	796888	2246199	7.52%	1.134%
22	6.939	946	949	959	rVB4	480832	1134504	3.80%	0.573%
23	7.256	991	1001	1009	rBV	404419	887141	2.97%	0.448%
24	7.458	1021	1034	1047	rBV3	1473537	3843491	12.86%	1.940%
25	7.732	1073	1079	1088	rVB	301484	574891	1.92%	0.290%
26	7.927	1104	1111	1112	rBV	227089	320580	1.07%	0.162%
27	7.957	1113	1116	1124	rVV3	278318	599907	2.01%	0.303%
28	8.055	1124	1132	1144	rVB	367068	845509	2.83%	0.427%
29	8.177	1144	1152	1154	rBV	435658	815853	2.73%	0.412%
30	8.214	1154	1158	1163	rVV	1057714	1866119	6.25%	0.942%
31	8.256	1163	1165	1175	rVB3	176645	316730	1.06%	0.160%
32	8.500	1200	1205	1210	rBV	219769	365538	1.22%	0.185%
33	8.573	1211	1217	1226	rVB2	977379	1737979	5.82%	0.877%
34	8.671	1226	1233	1235	rBV2	171020	312073	1.04%	0.158%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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 Title : SW846 8260

35	8.713	1235	1240	1247	rVB	1651466	2610040	8.74%	1.318%
36	8.909	1264	1272	1279	rVB	385427	710430	2.38%	0.359%
37	8.988	1279	1285	1289	rBV	215334	331386	1.11%	0.167%
38	9.165	1305	1314	1318	rBV3	159134	339605	1.14%	0.171%
39	9.219	1318	1323	1335	rVB	508111	846073	2.83%	0.427%
40	9.628	1385	1390	1393	rBV2	308409	467117	1.56%	0.236%
41	10.085	1459	1465	1466	rBV	407344	567371	1.90%	0.286%
42	10.110	1466	1469	1474	rVB	1580572	2123755	7.11%	1.072%
43	10.201	1480	1484	1488	rVB3	245593	353740	1.18%	0.179%
44	10.250	1488	1492	1497	rVV	2905878	3693987	12.36%	1.865%
45	10.305	1497	1501	1505	rVV	238275	440347	1.47%	0.222%
46	10.359	1505	1510	1515	rVB	2170725	2724120	9.12%	1.375%
47	10.847	1585	1590	1596	rVB	213333	328533	1.10%	0.166%
48	11.018	1611	1618	1625	rVB	6874795	8457317	28.30%	4.270%
49	11.140	1633	1638	1642	rVB	1436274	1823092	6.10%	0.920%
50	11.225	1647	1652	1659	rVB	638189	775566	2.60%	0.392%
51	11.359	1670	1674	1684	rVB	7380797	8630068	28.88%	4.357%
52	11.451	1684	1689	1693	rVV	1092516	1363703	4.56%	0.688%
53	11.506	1693	1698	1705	rVB	946679	1408969	4.72%	0.711%
54	11.664	1719	1724	1733	rBV	2249531	2833835	9.48%	1.431%
55	11.804	1743	1747	1751	rBV	1703476	1954960	6.54%	0.987%
56	11.847	1751	1754	1761	rVB	365926	464574	1.55%	0.235%
57	11.920	1761	1766	1768	rBV2	446352	655173	2.19%	0.331%
58	11.945	1768	1770	1777	rVB	586066	638347	2.14%	0.322%
59	12.079	1786	1792	1798	rBV	1846113	2381246	7.97%	1.202%
60	12.140	1798	1802	1807	rBV	1154370	1426916	4.78%	0.720%
61	12.298	1822	1828	1835	rBV	24563999	29880077	100.00%	15.085%
62	12.371	1835	1840	1850	rVB3	1202374	2055477	6.88%	1.038%
63	12.463	1850	1855	1860	rVB2	344759	459881	1.54%	0.232%
64	12.518	1860	1864	1868	rVB	293907	341800	1.14%	0.173%
65	12.585	1870	1875	1878	rBV	324504	425685	1.42%	0.215%
66	12.670	1883	1889	1892	rBV	4373957	5379643	18.00%	2.716%
67	12.701	1892	1894	1898	rVV	1321455	1353431	4.53%	0.683%
68	12.755	1898	1903	1910	rVB	4812475	6047172	20.24%	3.053%
69	12.975	1932	1939	1943	rBV	2690162	3217913	10.77%	1.625%
70	13.018	1943	1946	1951	rVB	685898	767520	2.57%	0.387%
71	13.225	1976	1980	1989	rVB	3182001	3825524	12.80%	1.931%

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Integration Parameters: RTEINT.P

Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
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72	13.347	1995	2000	2005	rVB2	6285951	8130847	27.21%	4.105%
73	13.402	2005	2009	2017	rVB3	475332	726175	2.43%	0.367%
74	13.481	2017	2022	2028	rVB	947655	1181904	3.96%	0.597%
75	13.603	2038	2042	2045	rVV	444596	604746	2.02%	0.305%
76	13.639	2045	2048	2053	rVV2	451496	654921	2.19%	0.331%
77	13.700	2053	2058	2059	rVV3	261633	365173	1.22%	0.184%
78	13.719	2059	2061	2068	rVB2	548342	728623	2.44%	0.368%
79	13.834	2072	2080	2089	rBV	17737173	21201873	70.96%	10.704%
80	13.926	2089	2095	2099	rBV	219777	299923	1.00%	0.151%
81	14.133	2124	2129	2136	rBV	295841	368453	1.23%	0.186%
82	14.273	2147	2152	2157	rBV	228911	379177	1.27%	0.191%
83	14.694	2216	2221	2231	rVB	2248928	2717248	9.09%	1.372%
84	14.840	2239	2245	2253	rBV	1363165	1740978	5.83%	0.879%

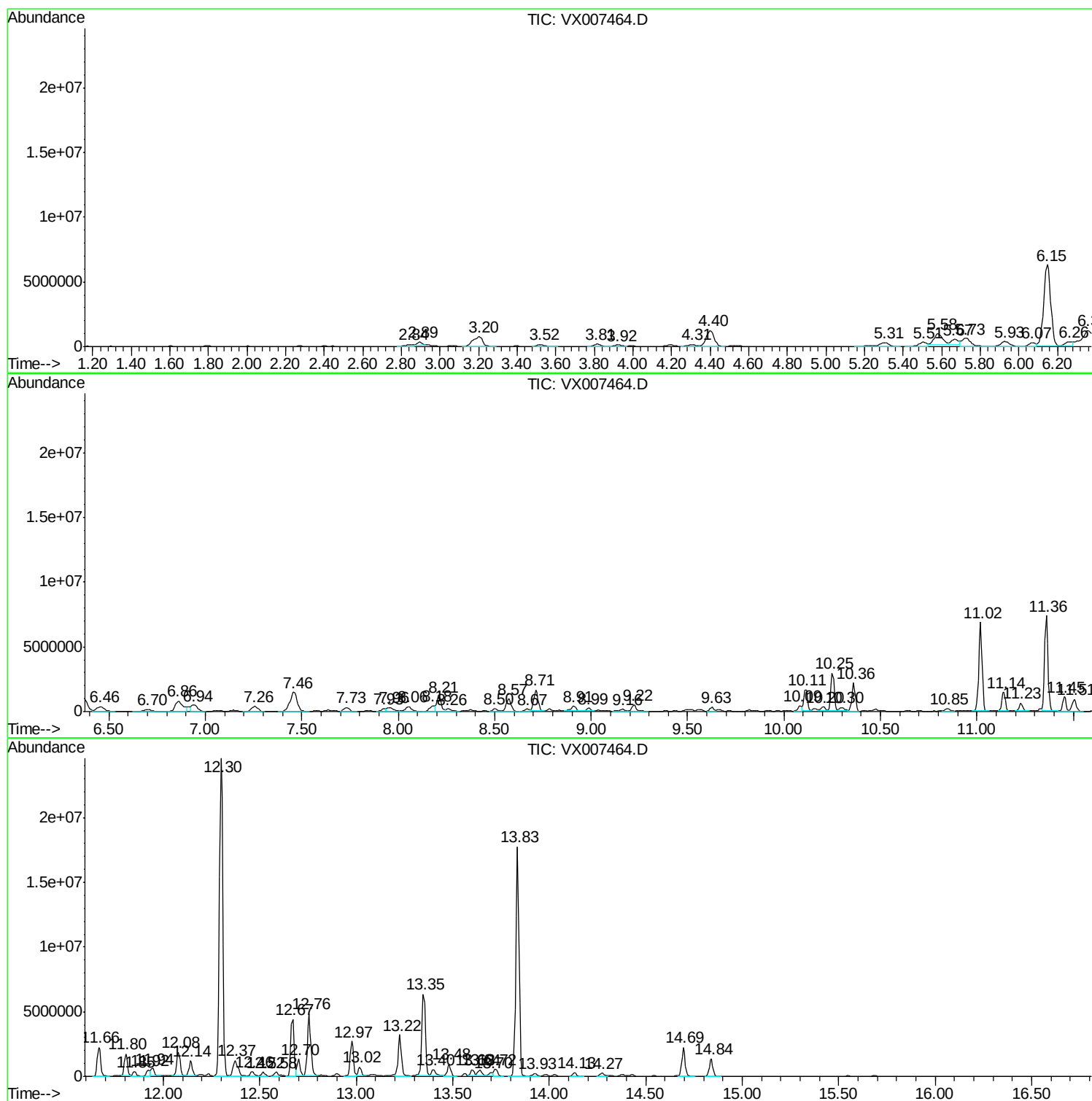
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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



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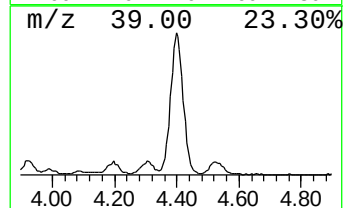
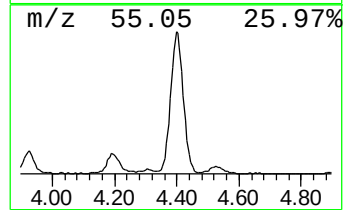
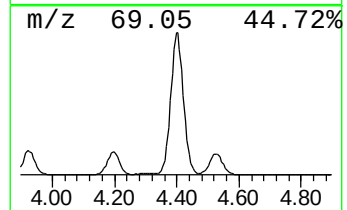
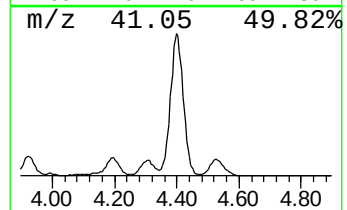
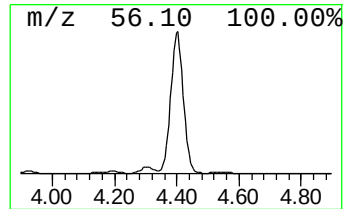
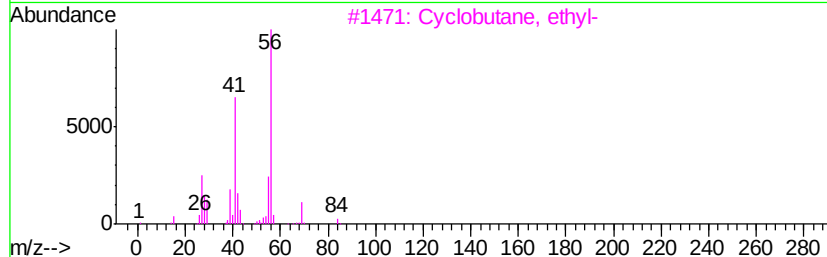
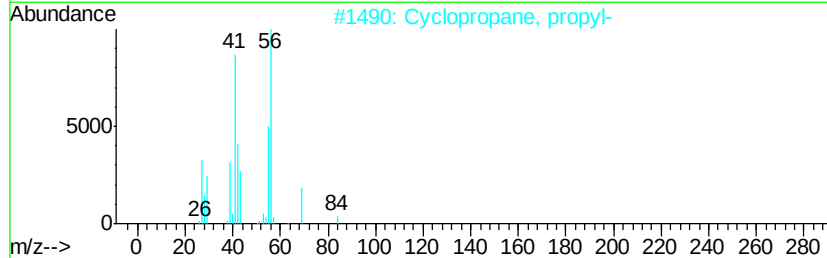
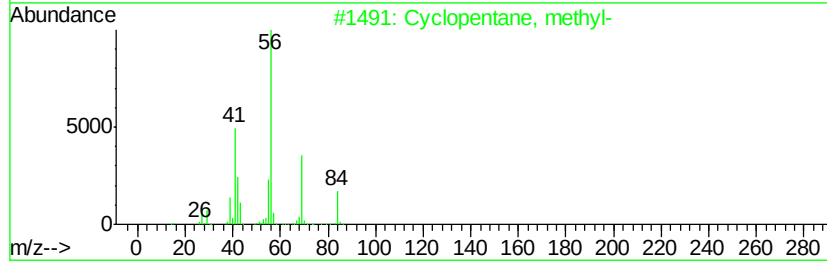
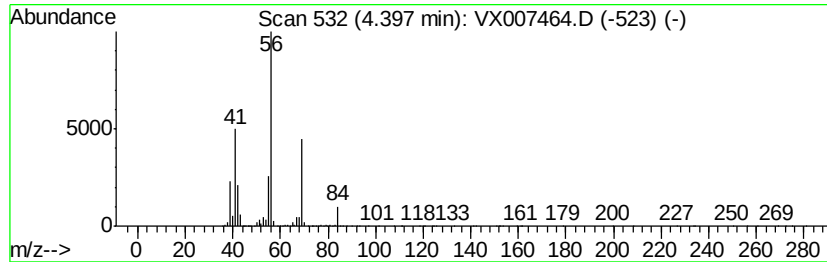
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	163.96 ug/l	3547940	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	80
2		Cyclopropane, propyl-	84	C6H12	002415-72-7	72
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50



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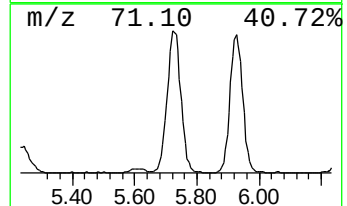
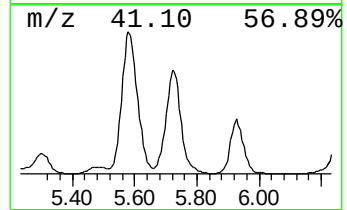
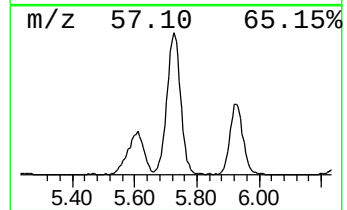
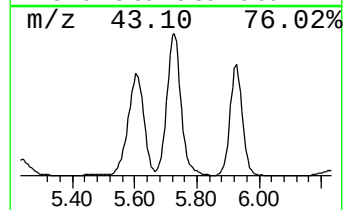
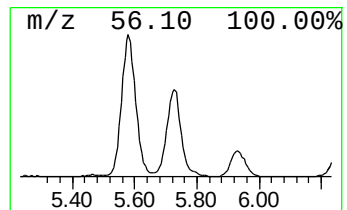
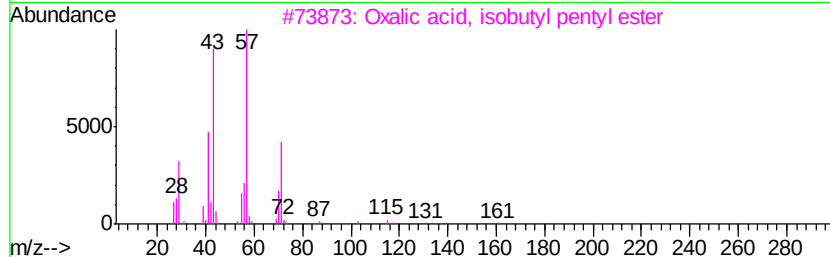
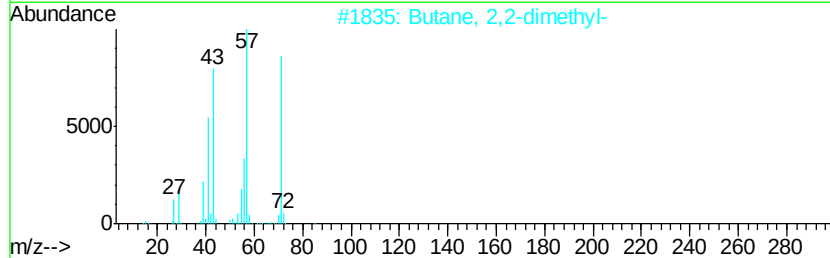
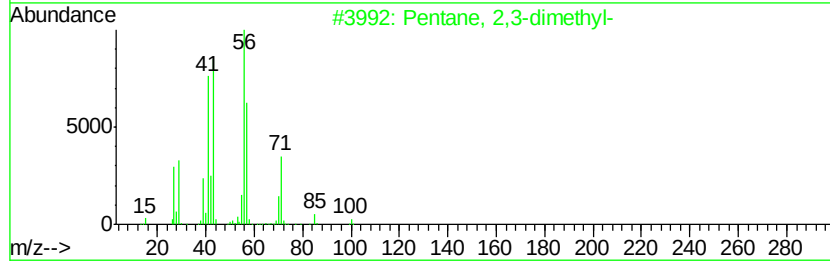
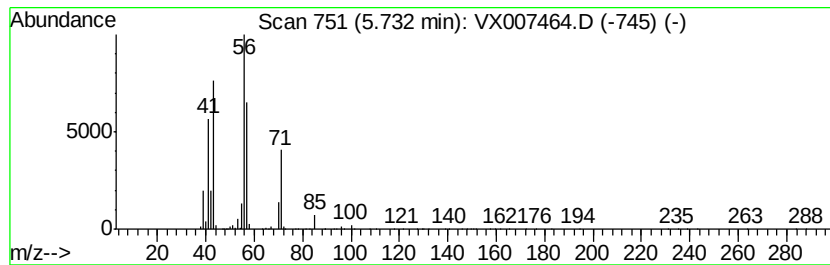
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.73	94.88 ug/l	2053140	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	94
2		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	38
3		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	38
4		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	38
5		Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	38



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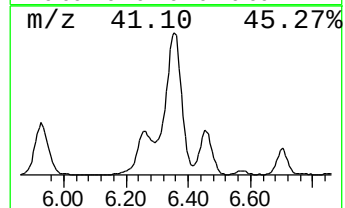
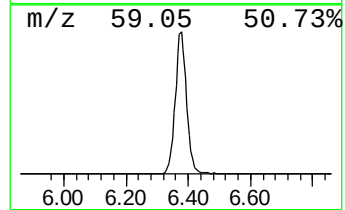
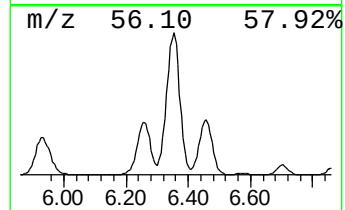
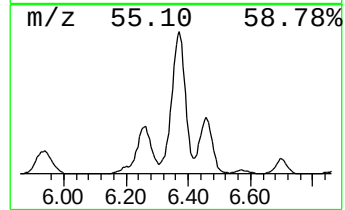
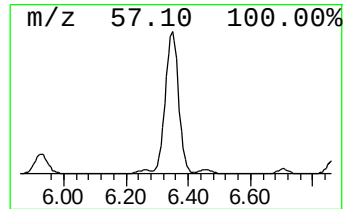
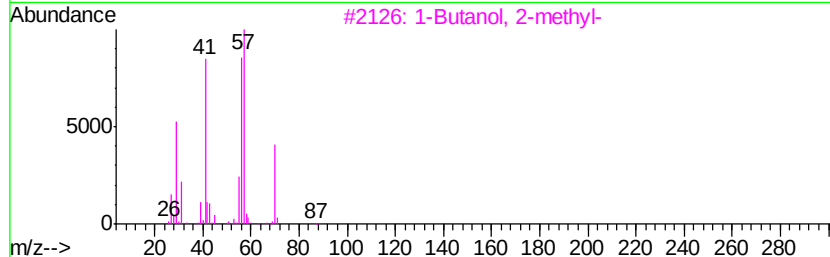
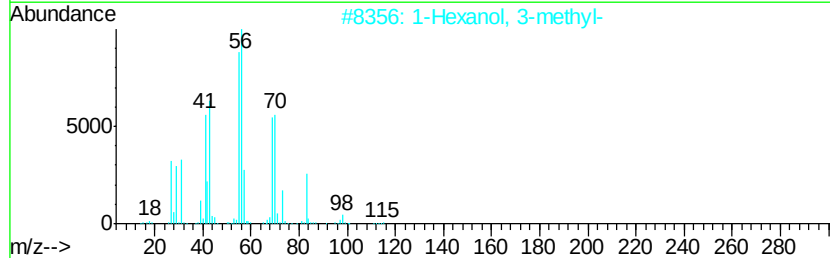
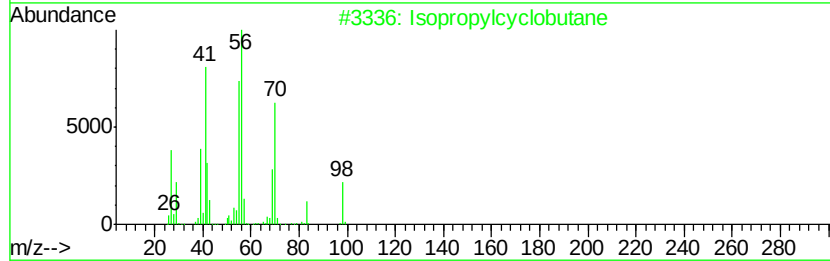
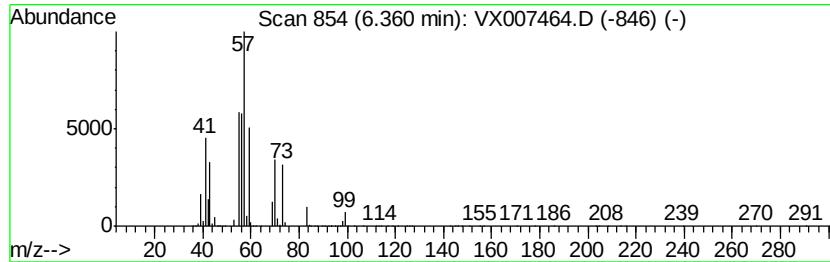
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Quant Title : SW846 8260

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

Peak Number 3 unknown6.36 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.36	98.02 ug/l	4403270	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropylcyclobutane	98	C7H14	000872-56-0	37
2		1-Hexanol, 3-methyl-	116	C7H16O	013231-81-7	37
3		1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	35
4		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	32
5		Acetamide, N-2-propenyl-	99	C5H9NO	000692-33-1	27



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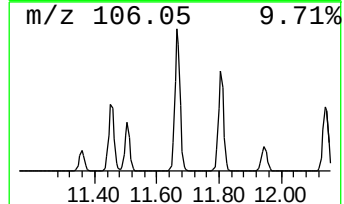
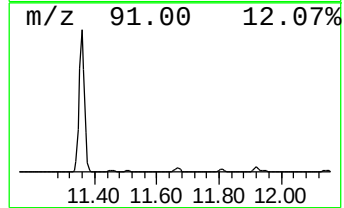
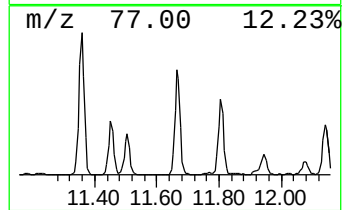
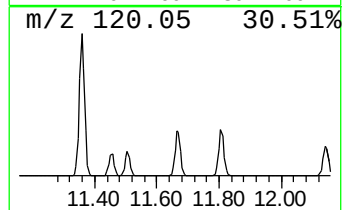
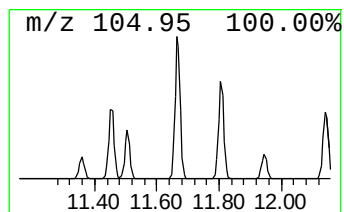
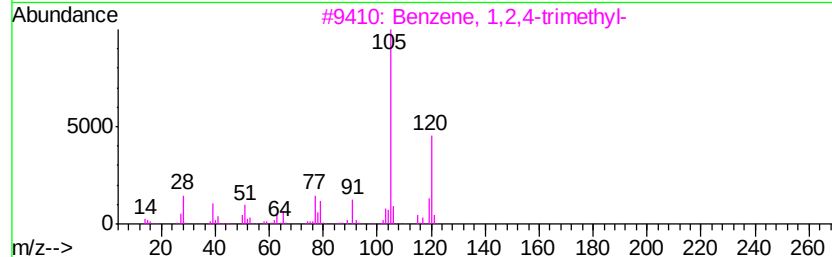
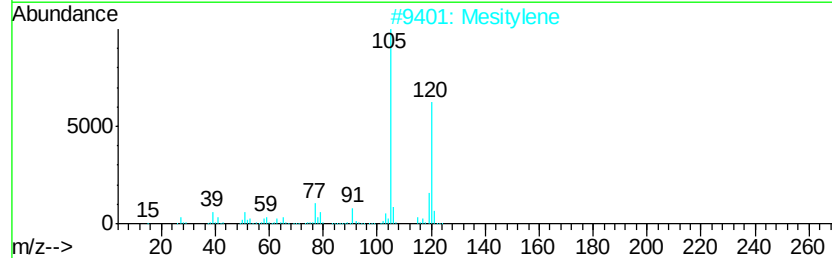
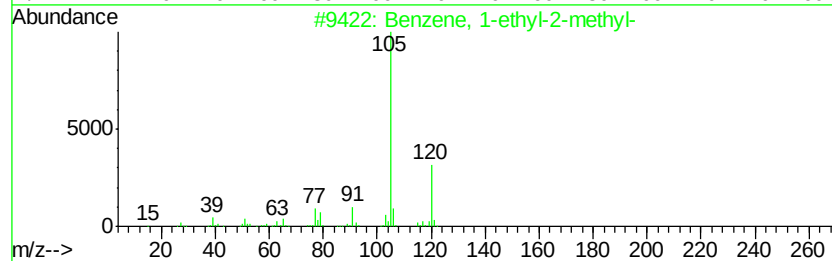
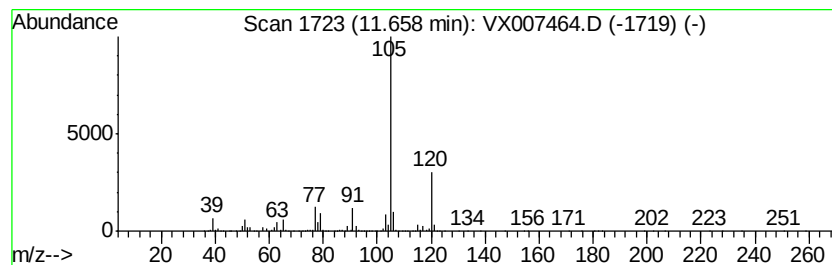
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	59.50 ug/l	2833840	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93	
2	Mesitylene	120	C9H12	000108-67-8	90	
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90	
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90	
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW038-190207

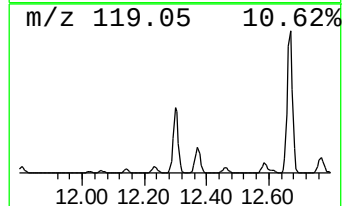
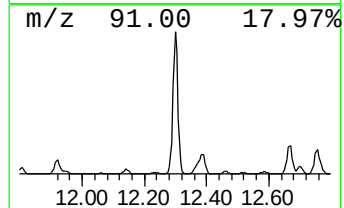
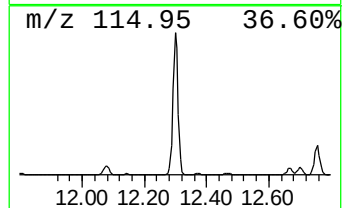
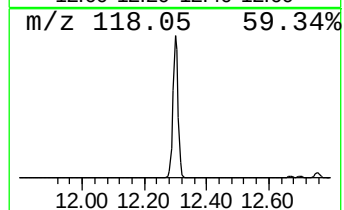
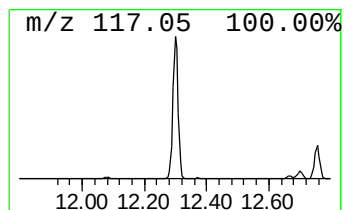
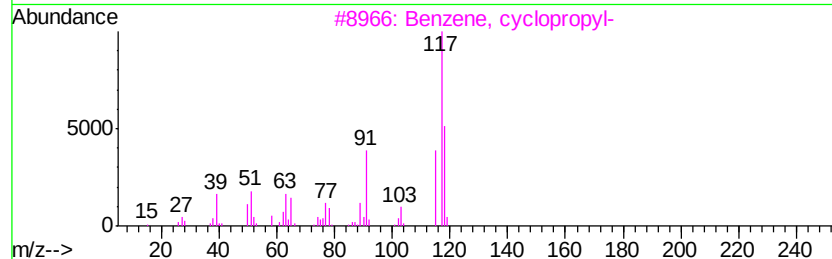
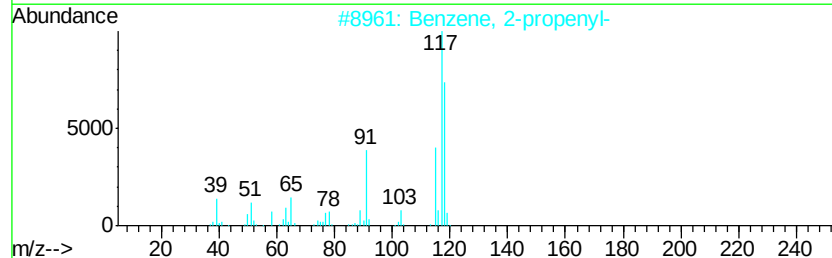
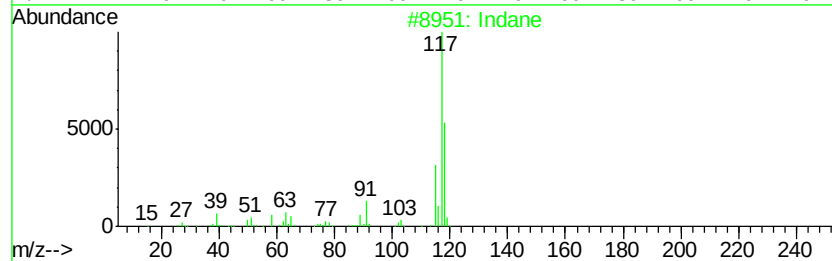
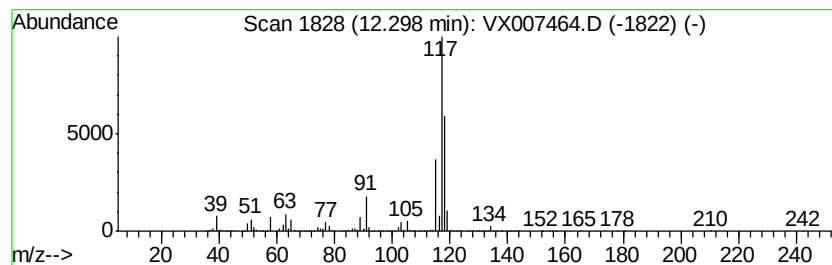
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Quant Title : SW846 8260

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

Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	627.40 ug/l	29880100	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	90
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	87
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
5		Benzeneethanol, .beta.-ethenyl-....	162	C11H14O	040596-14-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW038-190207

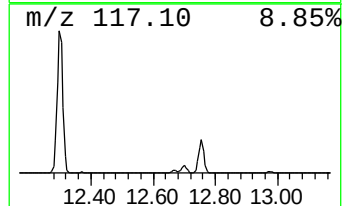
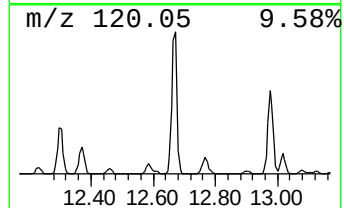
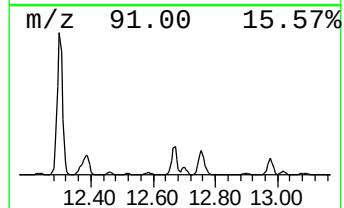
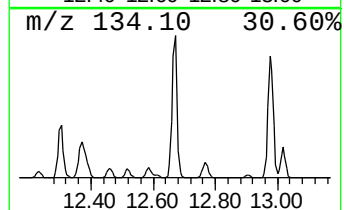
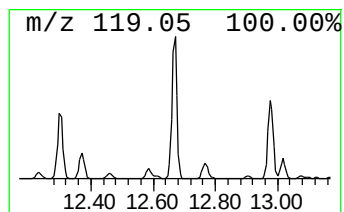
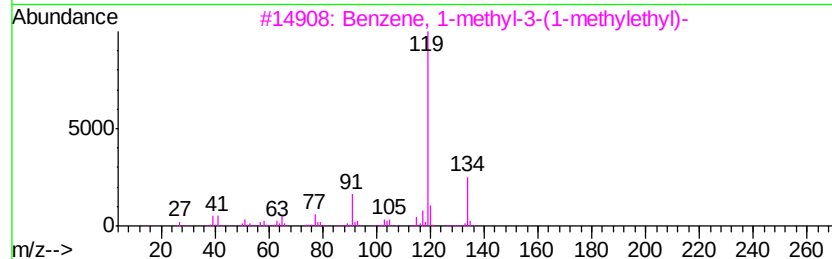
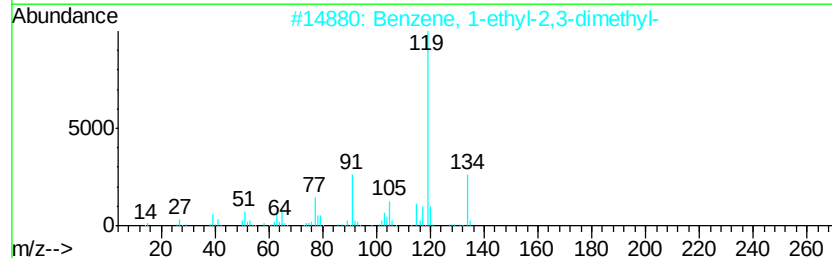
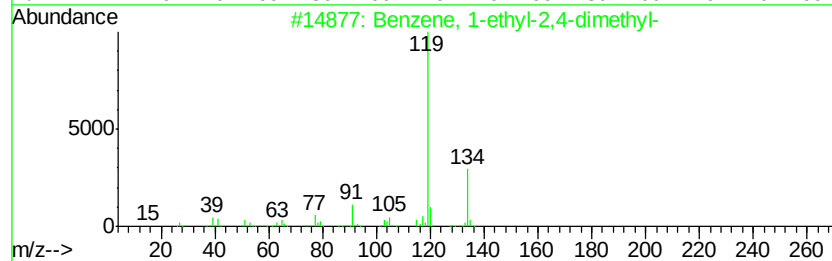
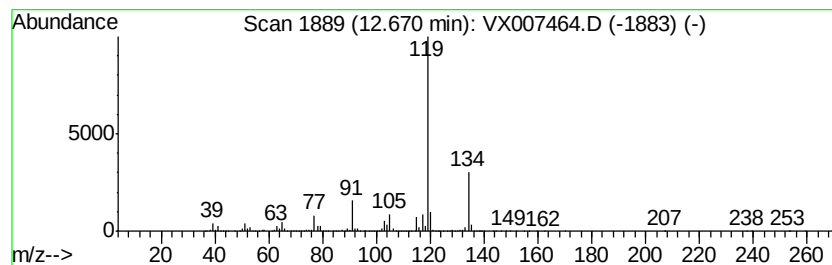
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	112.96 ug/l	5379640	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW038-190207

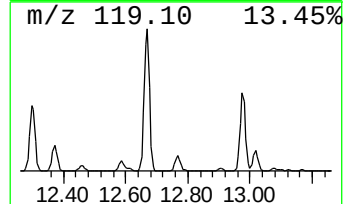
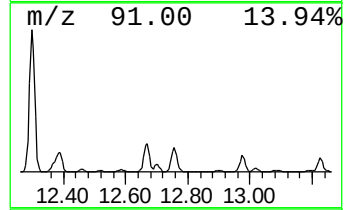
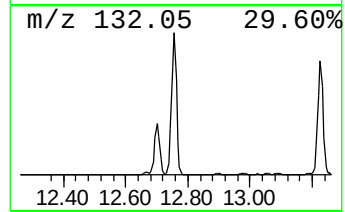
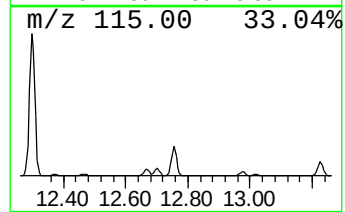
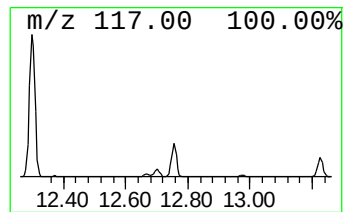
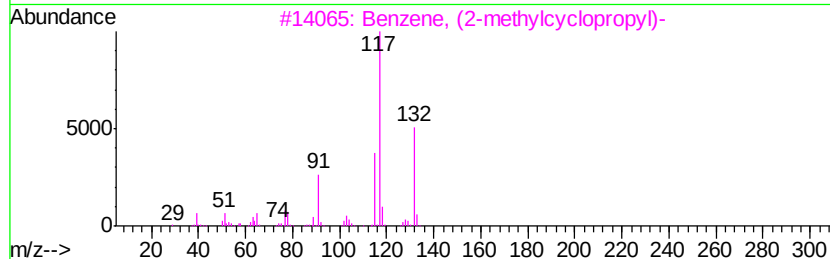
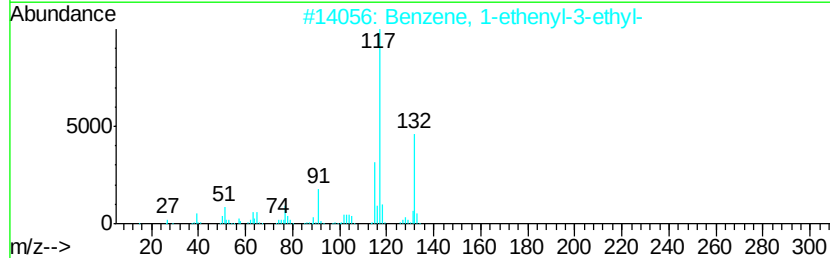
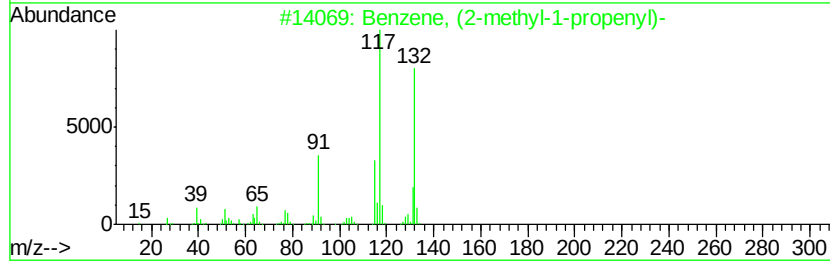
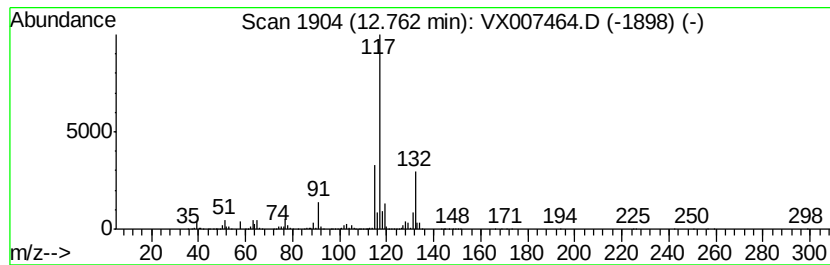
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-propen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	126.98 ug/l	6047170	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-methylcyclopropyl)-	132	C10H12	1000327-39-0	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\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW038-190207

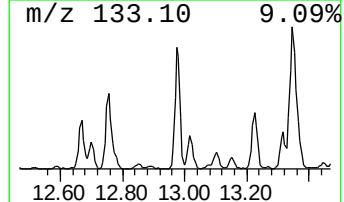
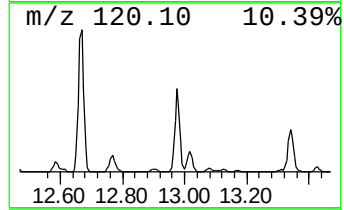
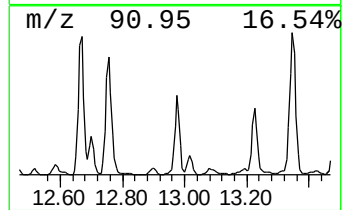
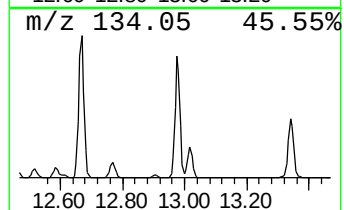
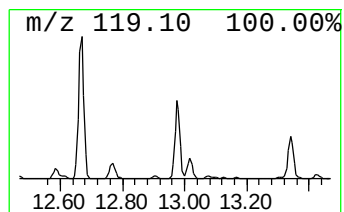
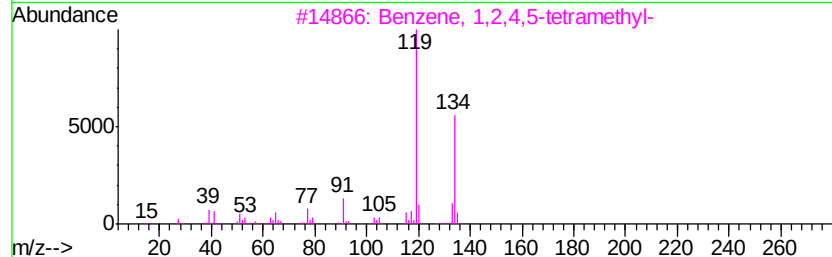
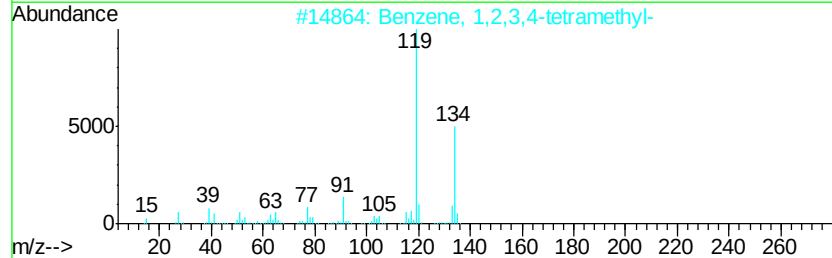
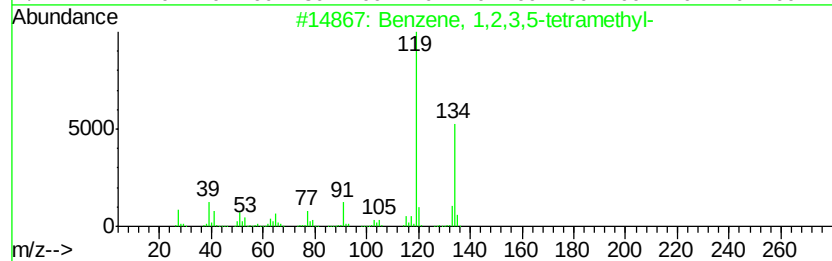
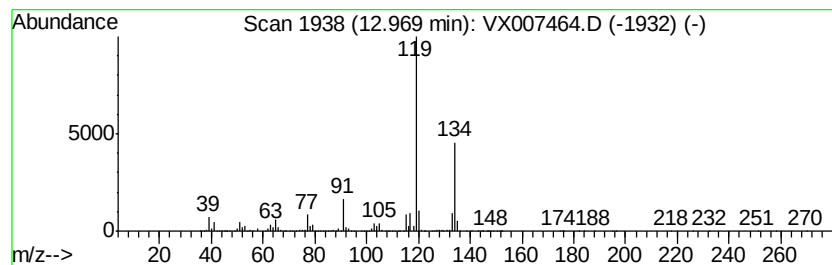
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	67.57 ug/l	3217910	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

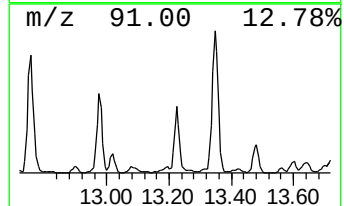
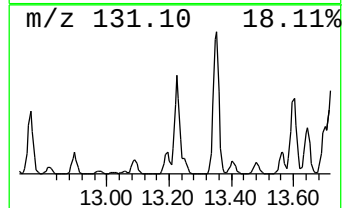
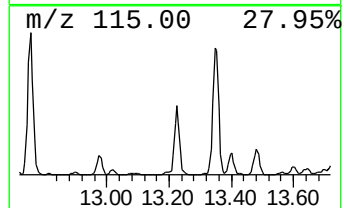
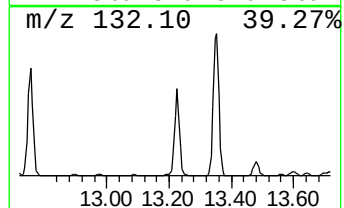
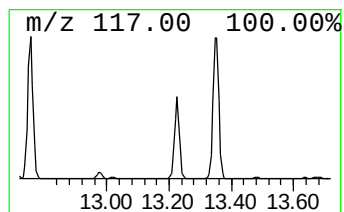
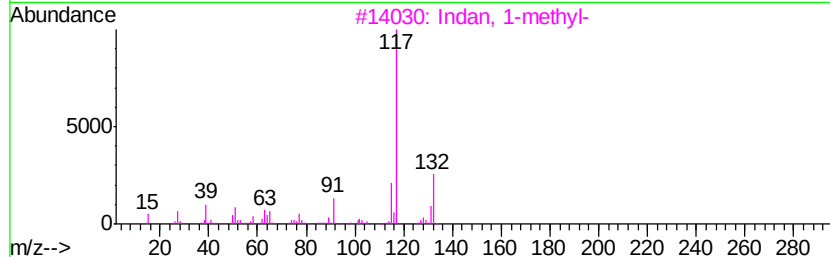
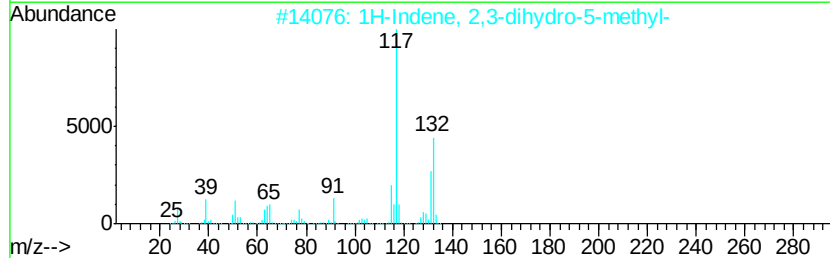
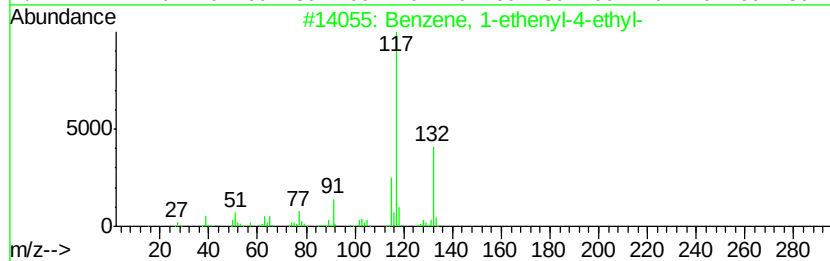
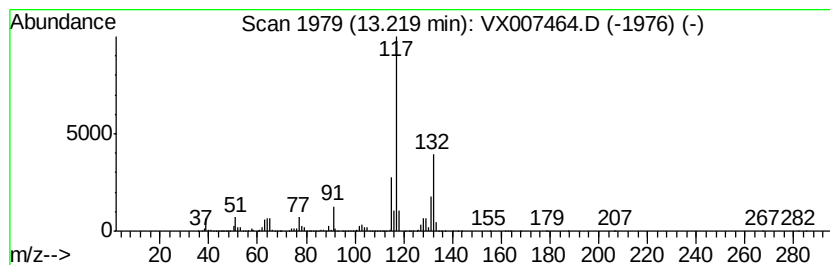
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MSVOA_X
ClientSampleId :
KY001MW038-190207

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	80.33 ug/l	3825520	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 94
2		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 93
3		Indan, 1-methyl-	132 C10H12	000767-58-8 90
4		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 90
5		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW038-190207

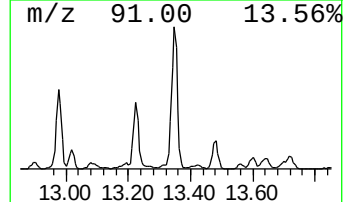
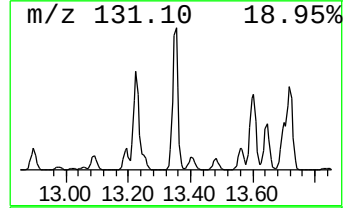
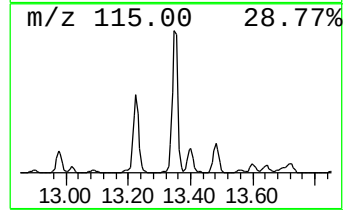
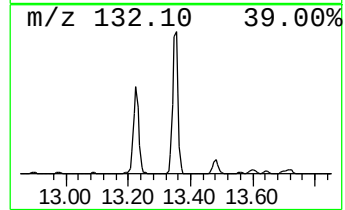
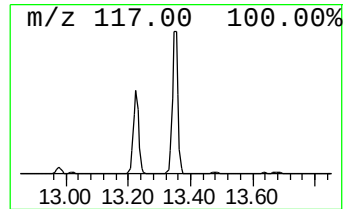
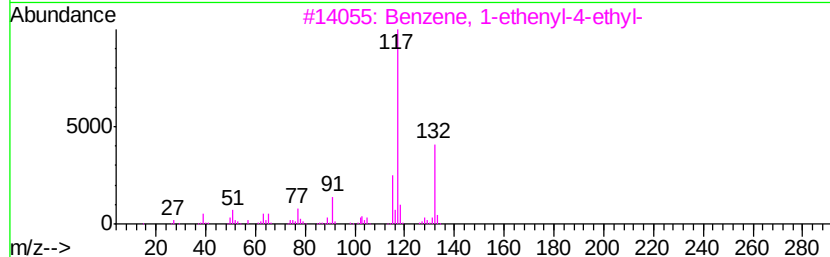
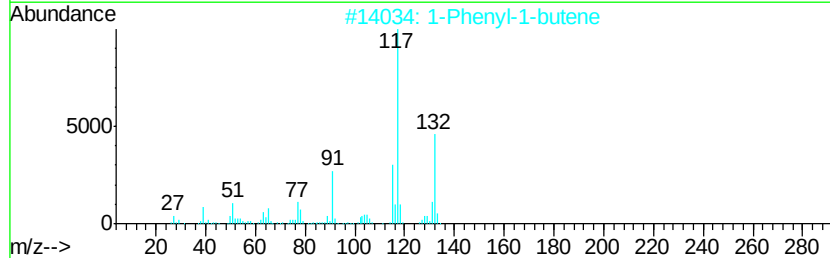
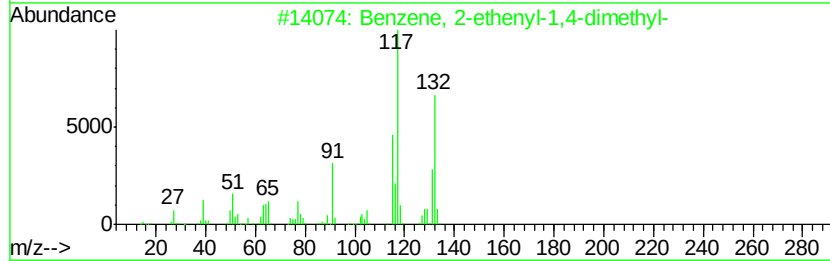
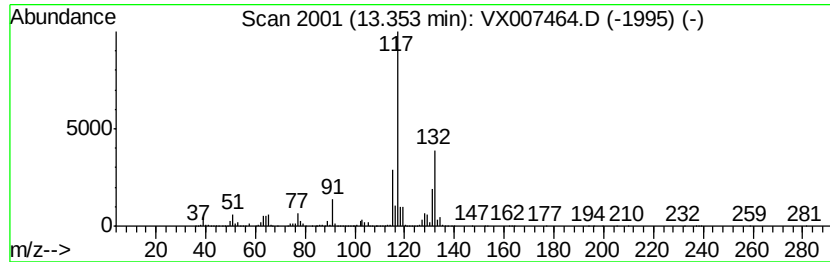
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	170.73 ug/l	8130850	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	91
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX021119\
Data File : VX007464.D
Acq On : 11 Feb 2019 13:33
Operator : JC/SP
Sample : K1433-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW038-190207

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.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	164.0	ug/l	3547940	1	5.67	1081960	50.0
Pentane, 2,3-dime...	5.73	94.9	ug/l	2053140	1	5.67	1081960	50.0
unknown6.36	6.36	98.0	ug/l	4403270	2	6.86	2246200	50.0
Benzene, 1-ethyl-...	11.66	59.5	ug/l	2833840	4	12.08	2381250	50.0
Indane	12.30	627.4	ug/l	29880100	4	12.08	2381250	50.0
Benzene, 1-ethyl-...	12.67	113.0	ug/l	5379640	4	12.08	2381250	50.0
Benzene, (2-methy...	12.76	127.0	ug/l	6047170	4	12.08	2381250	50.0
Benzene, 1,2,3,5-...	12.97	67.6	ug/l	3217910	4	12.08	2381250	50.0
Benzene, 1-etheny...	13.22	80.3	ug/l	3825520	4	12.08	2381250	50.0
Benzene, 2-etheny...	13.35	170.7	ug/l	8130850	4	12.08	2381250	50.0