

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042220\
 Data File : VX015872.D
 Acq On : 23 Apr 2020 04:20
 Operator : JC/SP
 Sample : L2380-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY001MW031-200421

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\82X042220W.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	1.064	24	29	54	rBV2	18771005	61417634	100.00%	19.195%
2	2.826	310	318	321	rBV	755756	1540948	2.51%	0.482%
3	2.875	321	326	330	rVV	1342663	2928890	4.77%	0.915%
4	2.911	330	332	343	rVV2	671044	1318340	2.15%	0.412%
5	3.161	362	373	392	rBV3	5061601	12956121	21.10%	4.049%
6	3.503	421	429	441	rVB	285141	661479	1.08%	0.207%
7	3.795	468	477	486	rBV	482363	1164859	1.90%	0.364%
8	3.905	486	495	501	rBV	284766	724613	1.18%	0.226%
9	4.173	525	539	547	rVV3	248131	750228	1.22%	0.234%
10	4.283	547	557	564	rVV	485329	1439274	2.34%	0.450%
11	4.381	564	573	585	rVV	2536097	7294246	11.88%	2.280%
12	5.472	739	752	758	rBV	587722	1779907	2.90%	0.556%
13	5.563	758	767	774	rVV5	1518804	5871612	9.56%	1.835%
14	5.630	776	778	783	rVV	1009159	2212231	3.60%	0.691%
15	5.703	783	790	808	rVV	1902741	6341469	10.33%	1.982%
16	5.905	812	823	836	rVV2	1250237	3841205	6.25%	1.200%
17	6.039	836	845	850	rVV	557395	1427192	2.32%	0.446%
18	6.118	850	858	867	rVV	1697818	4534029	7.38%	1.417%
19	6.240	869	878	884	rVV	1049864	3582980	5.83%	1.120%
20	6.325	884	892	901	rVV3	3033368	9813253	15.98%	3.067%
21	6.429	901	909	921	rVB4	959226	3152130	5.13%	0.985%
22	6.837	963	976	981	rBV	1474945	3827597	6.23%	1.196%
23	6.923	986	990	1000	rVB4	972989	2336196	3.80%	0.730%
24	7.234	1032	1041	1049	rBV	761905	1662951	2.71%	0.520%
25	7.435	1062	1074	1086	rBV4	1972288	5810412	9.46%	1.816%
26	7.618	1099	1104	1109	rBV	378470	718958	1.17%	0.225%
27	7.715	1115	1120	1128	rVB	438865	842932	1.37%	0.263%
28	7.947	1145	1158	1164	rBV7	393343	1336203	2.18%	0.418%
29	8.038	1164	1173	1185	rVB	1249373	2894718	4.71%	0.905%
30	8.160	1185	1193	1197	rBV	1403546	2954507	4.81%	0.923%
31	8.240	1202	1206	1216	rVB3	416728	856861	1.40%	0.268%
32	8.362	1216	1226	1232	rBV4	320155	730942	1.19%	0.228%
33	8.550	1250	1257	1267	rVB4	646109	1481664	2.41%	0.463%
34	8.654	1267	1274	1276	rBV2	444271	764593	1.24%	0.239%

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35	8.697	1276	1281	1287	rVB	3063511	4876822	7.94%	1.524%
36	8.770	1287	1293	1298	rBV	685941	1080435	1.76%	0.338%
37	8.886	1305	1312	1320	rVB3	373815	830697	1.35%	0.260%
38	9.148	1346	1355	1360	rBV3	317819	635360	1.03%	0.199%
39	9.209	1360	1365	1374	rVB	712745	1137384	1.85%	0.355%
40	9.502	1404	1413	1416	rBV3	339953	806393	1.31%	0.252%
41	10.056	1499	1504	1507	rBV2	590416	1010098	1.64%	0.316%
42	10.099	1507	1511	1515	rVB	3113057	3961149	6.45%	1.238%
43	10.184	1521	1525	1528	rVB2	535790	640699	1.04%	0.200%
44	10.239	1528	1534	1539	rBV	25980404	35484925	57.78%	11.090%
45	10.343	1546	1551	1557	rVB	4817385	6521317	10.62%	2.038%
46	10.416	1557	1563	1566	rVV	456324	643111	1.05%	0.201%
47	10.684	1602	1607	1616	rVB	3570165	4481528	7.30%	1.401%
48	10.831	1627	1631	1637	rBV2	400500	647714	1.05%	0.202%
49	11.001	1652	1659	1664	rVV	4741106	6279729	10.22%	1.963%
50	11.056	1664	1668	1671	rVV	519270	653713	1.06%	0.204%
51	11.123	1675	1679	1684	rVV	3183104	4122512	6.71%	1.288%
52	11.202	1688	1692	1700	rVB2	429872	839670	1.37%	0.262%
53	11.343	1711	1715	1724	rVV	3726411	4956807	8.07%	1.549%
54	11.440	1724	1731	1734	rVV	743992	1104561	1.80%	0.345%
55	11.489	1734	1739	1747	rVB2	901496	1429416	2.33%	0.447%
56	11.654	1761	1766	1775	rVB	4778878	5956443	9.70%	1.862%
57	11.794	1785	1789	1793	rVB	1631123	2021206	3.29%	0.632%
58	12.062	1828	1833	1839	rVB	3964602	4846261	7.89%	1.515%
59	12.129	1839	1844	1848	rBV	1012551	1323677	2.16%	0.414%
60	12.221	1854	1859	1864	rVB	2360654	2886696	4.70%	0.902%
61	12.288	1864	1870	1875	rBV	25044108	33113125	53.91%	10.349%
62	12.355	1877	1881	1888	rVB	721168	968499	1.58%	0.303%
63	12.452	1893	1897	1902	rVV	599234	792829	1.29%	0.248%
64	12.684	1932	1935	1939	rVV	1031393	1244166	2.03%	0.389%
65	12.739	1939	1944	1952	rVB	3960269	5171493	8.42%	1.616%
66	12.964	1975	1981	1985	rBV	2009259	2658174	4.33%	0.831%
67	13.208	2017	2021	2031	rVB	1459498	1915490	3.12%	0.599%
68	13.336	2037	2042	2046	rVB	4056895	4968379	8.09%	1.553%
69	13.385	2046	2050	2056	rVV2	1035122	1235601	2.01%	0.386%
70	13.464	2059	2063	2069	rVB	974612	1207184	1.97%	0.377%
71	13.623	2086	2089	2094	rVV5	457073	729713	1.19%	0.228%

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72	13.702	2100	2102	2107	rVB	547888	654261	1.07%	0.204%
73	13.818	2113	2121	2129	rVB	1953735	2488082	4.05%	0.778%
74	13.909	2129	2136	2141	rBV	1442180	1768457	2.88%	0.553%
75	14.824	2281	2286	2291	rBV2	603589	908657	1.48%	0.284%

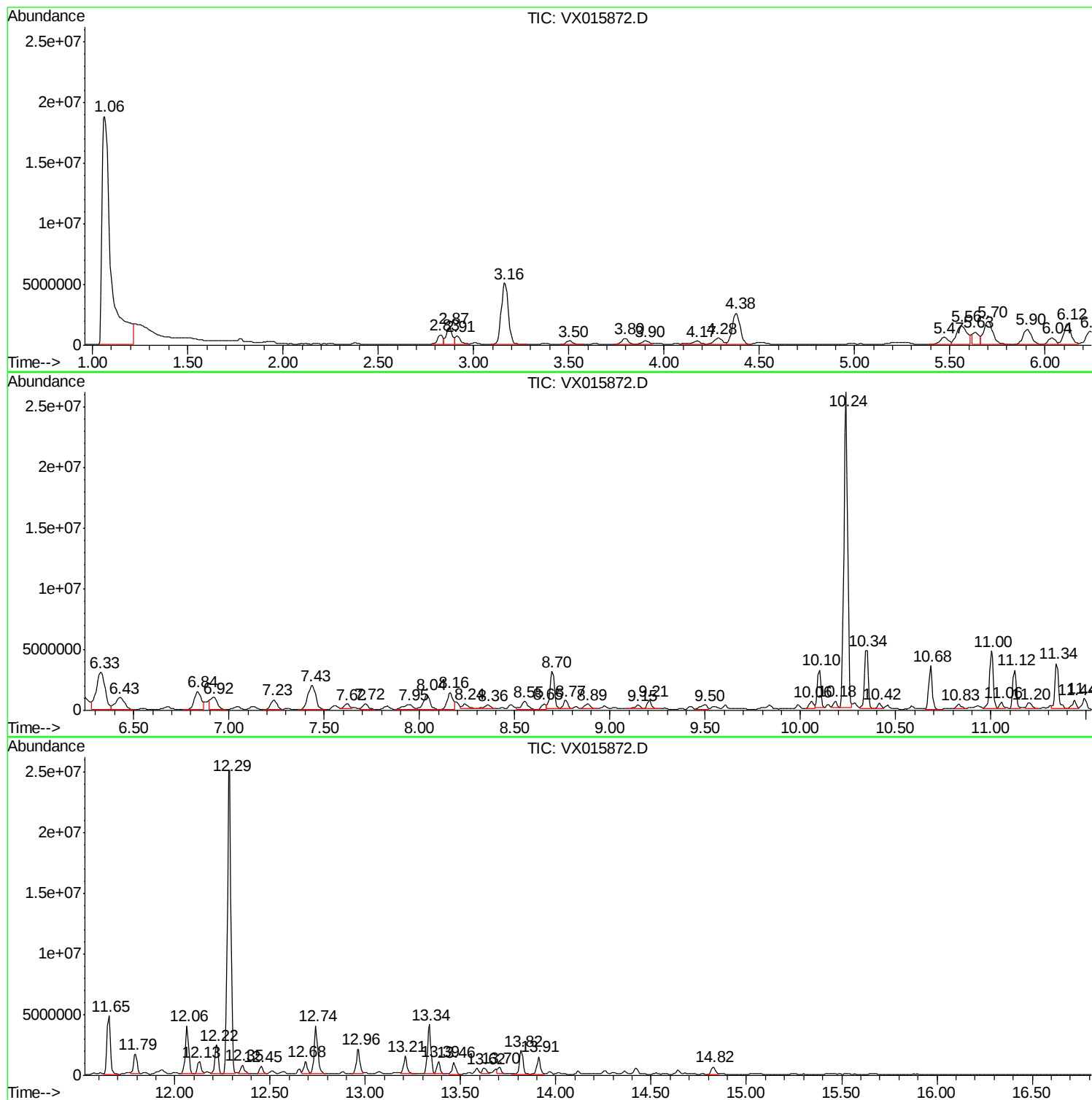
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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



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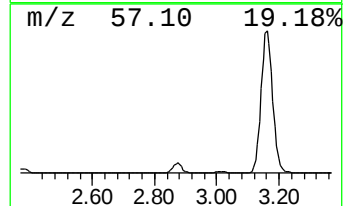
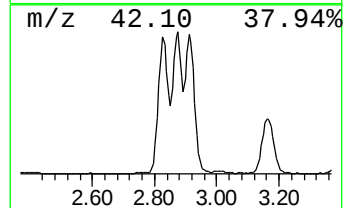
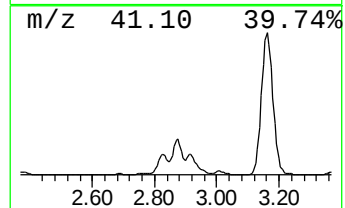
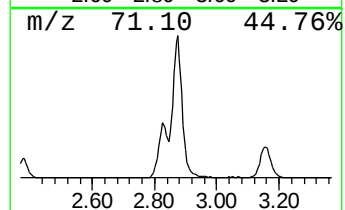
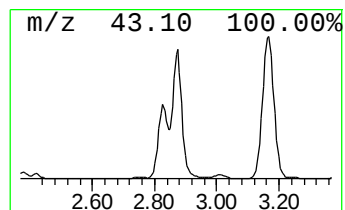
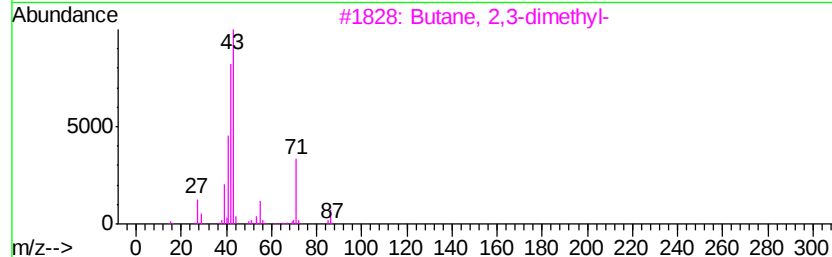
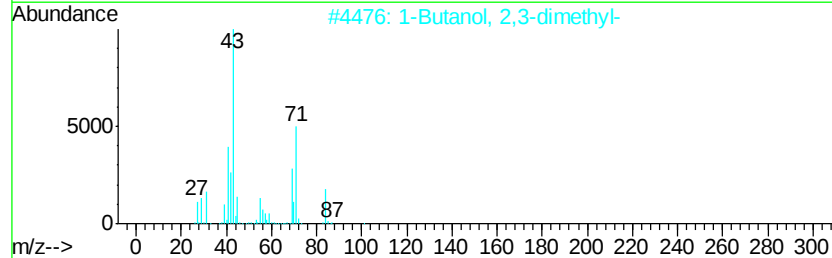
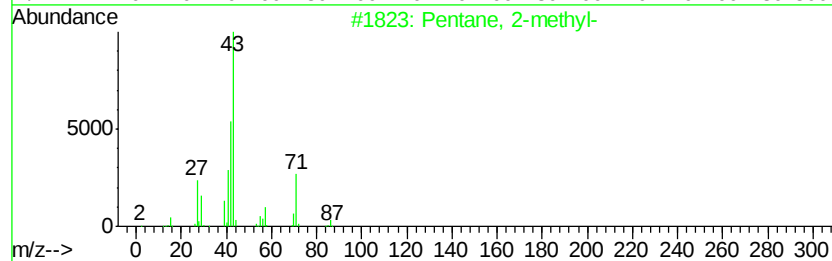
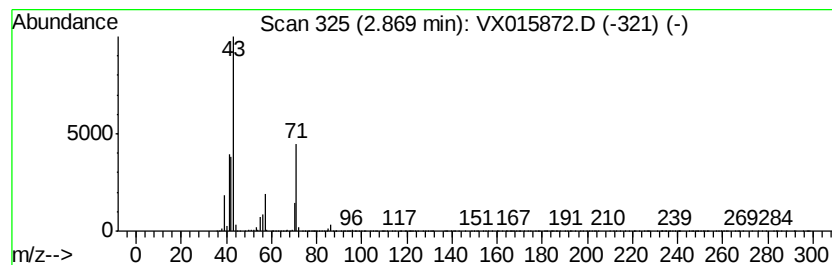
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	66.20 ug/l	2928890	Pentafluorobenzene	5.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	64
2			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	56
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	36
5			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	30



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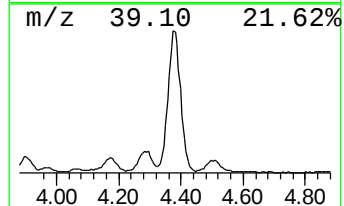
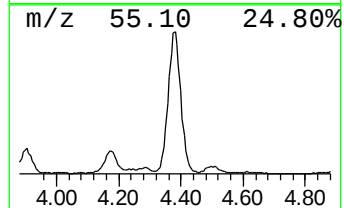
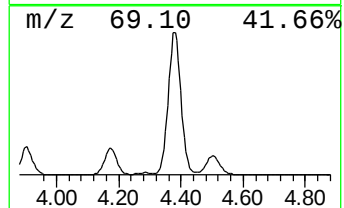
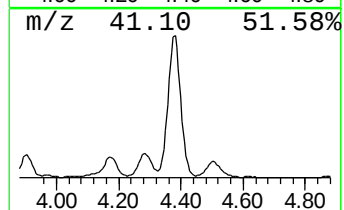
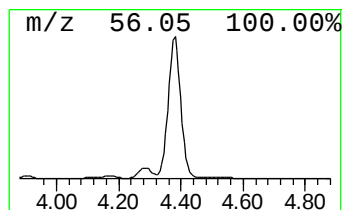
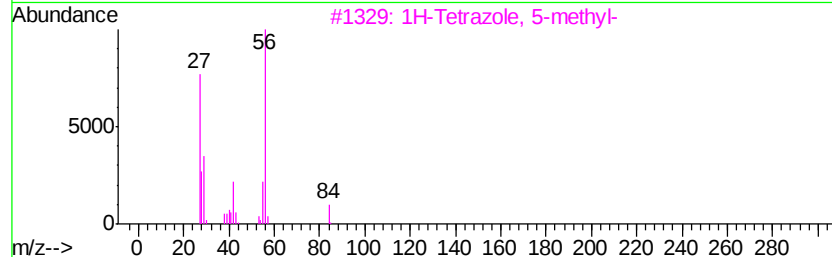
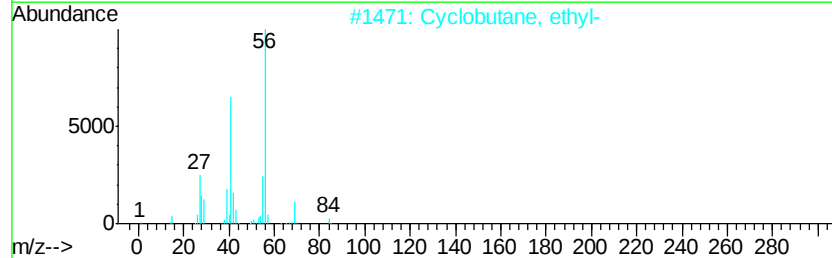
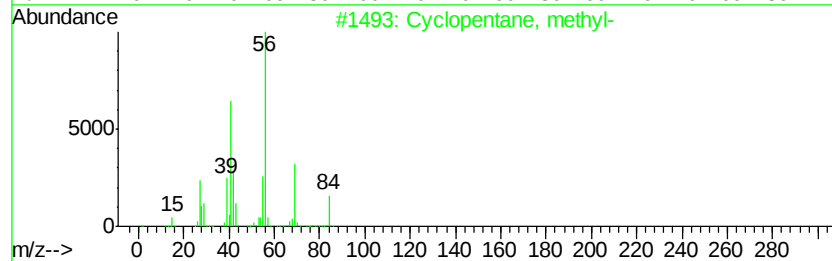
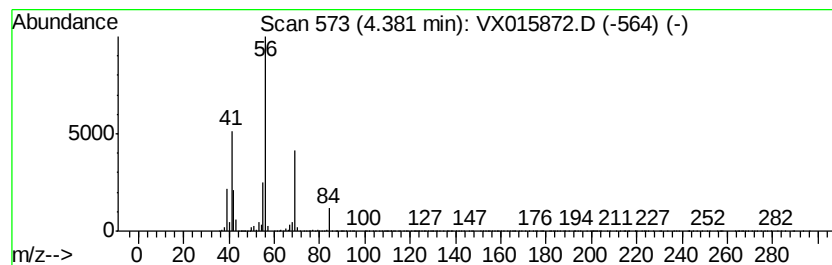
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	164.86 ug/l	7294250	Pentafluorobenzene	5.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	46
5		Cyclopropane, 1-ethyl-2-methyl-,...	84	C6H12	019781-68-1	38



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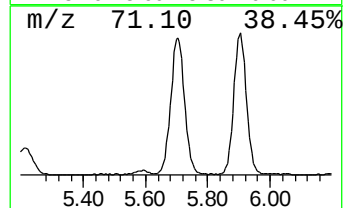
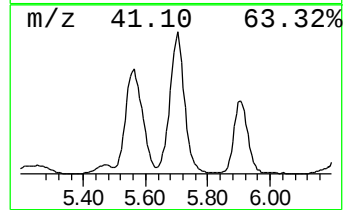
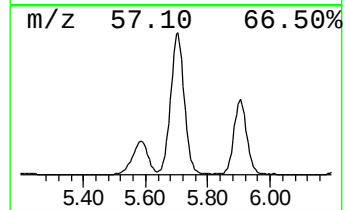
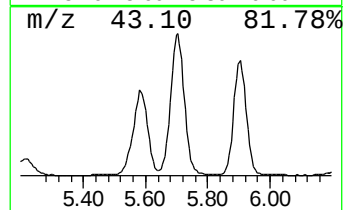
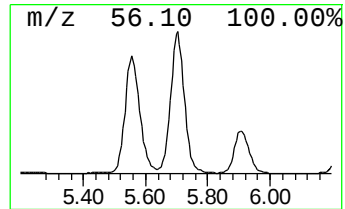
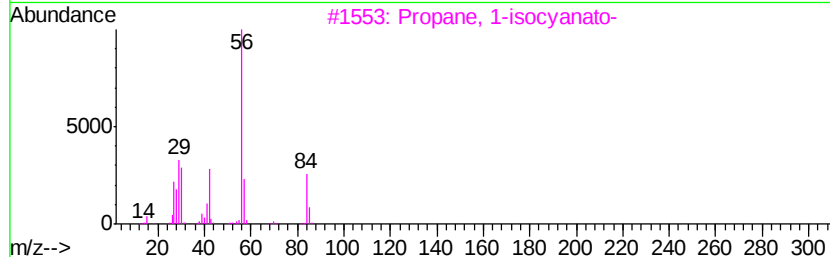
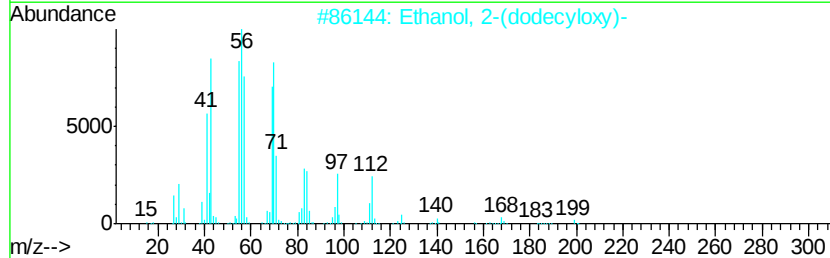
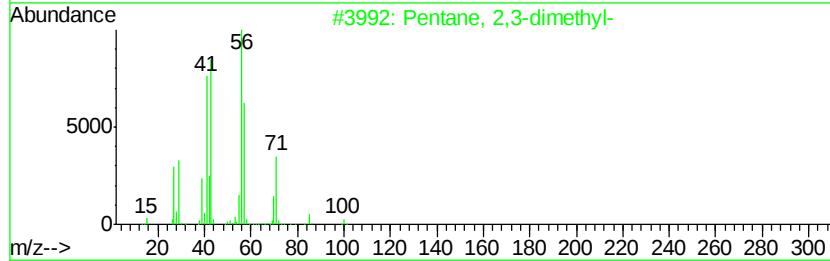
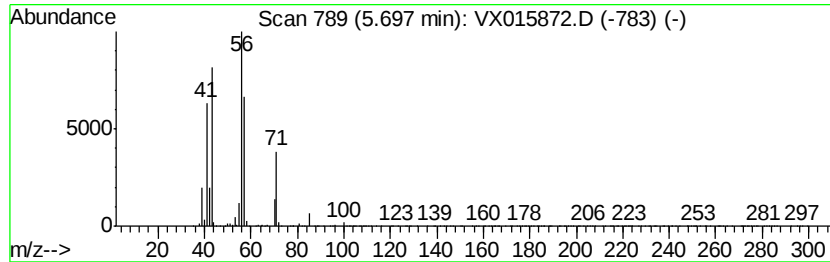
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	143.33 ug/l	6341470	Pentafluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2		Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	56
3		Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	50
4		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	40
5		Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	37



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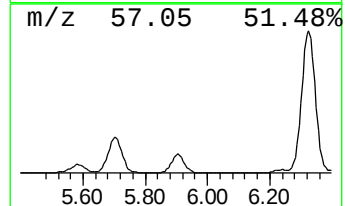
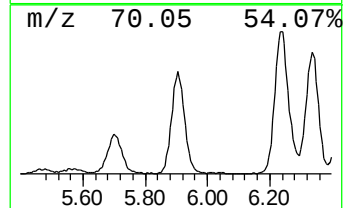
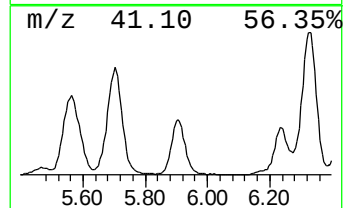
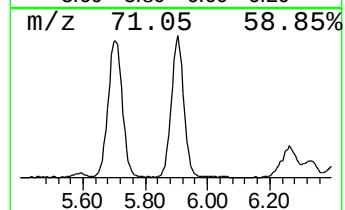
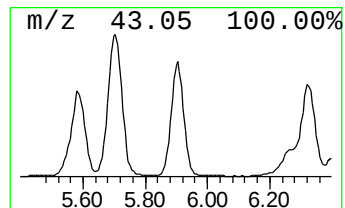
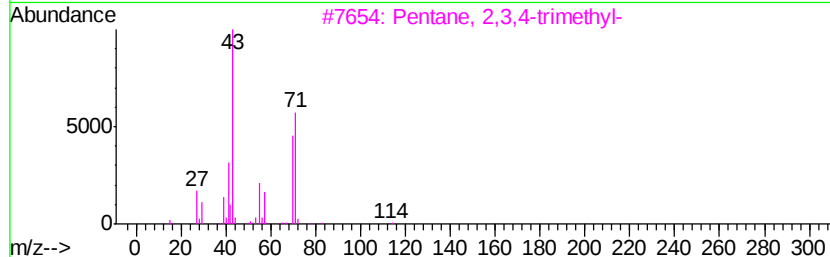
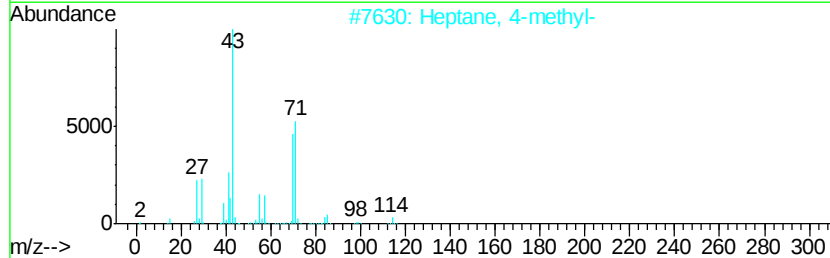
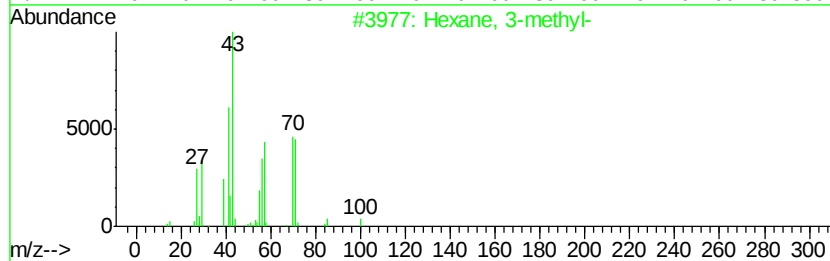
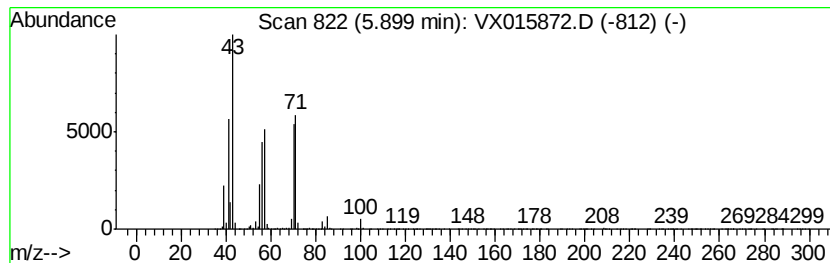
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Quant Title : SW846 8260

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

Peak Number 5 Hexane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.90	86.82 ug/l	3841210	Pentafluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	53
3		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
4		Heptane	100	C7H16	000142-82-5	53
5		Octane, 4,5-dimethyl-	142	C10H22	015869-96-2	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015872.D
Acq On : 23 Apr 2020 04:20
Operator : JC/SP
Sample : L2380-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

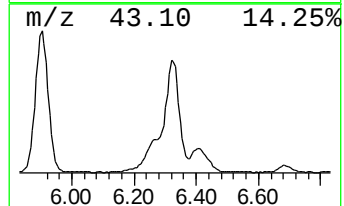
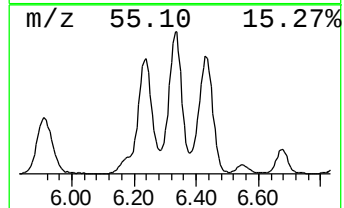
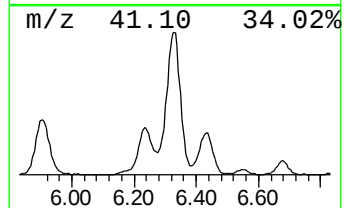
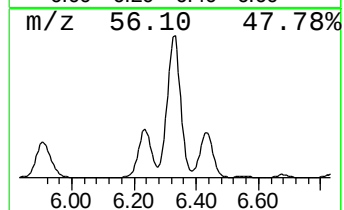
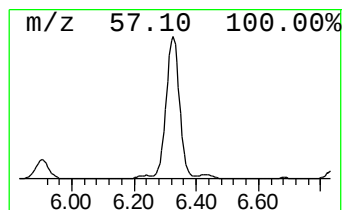
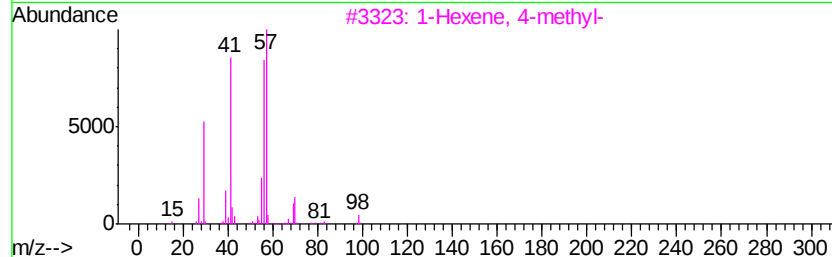
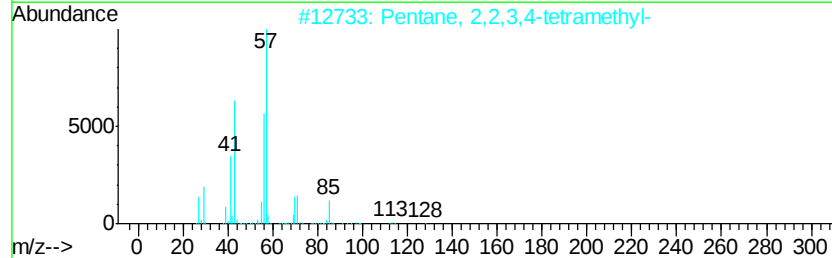
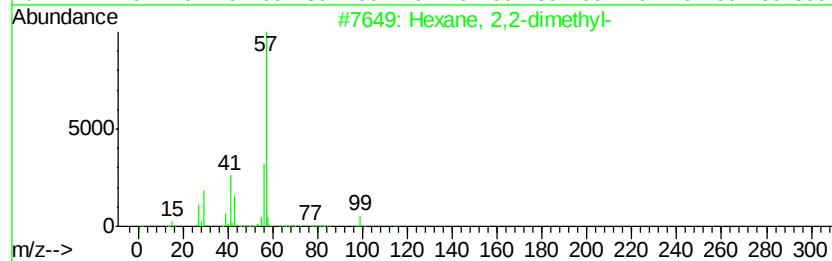
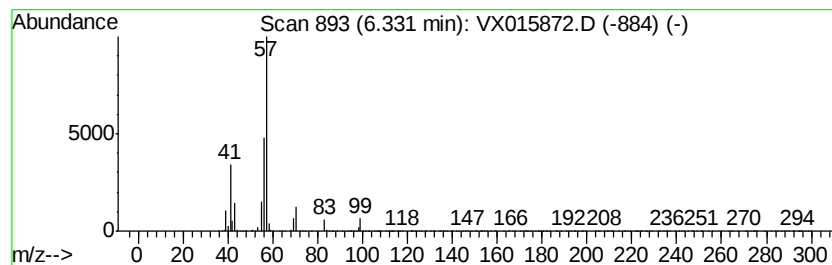
Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-200421

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.M
Quant Title : SW846 8260

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

Peak Number 6 Hexane, 2,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.33	128.19 ug/l	9813250	1,4-Difluorobenzene	6.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
2	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	56
3	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	53
4	2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	45
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015872.D
Acq On : 23 Apr 2020 04:20
Operator : JC/SP
Sample : L2380-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-200421

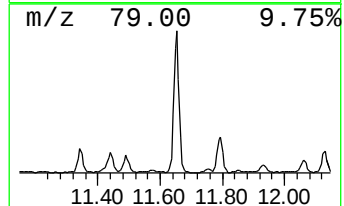
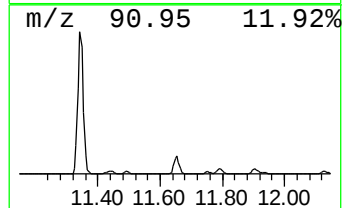
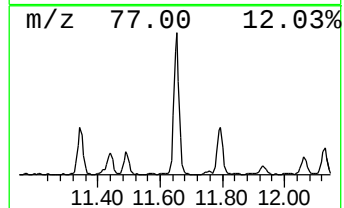
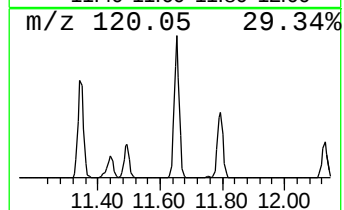
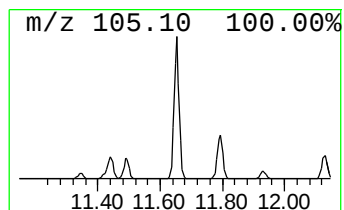
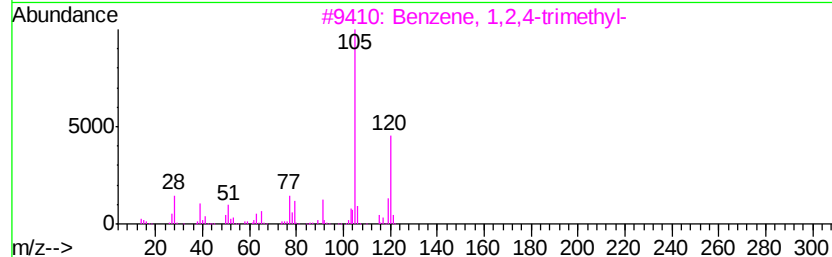
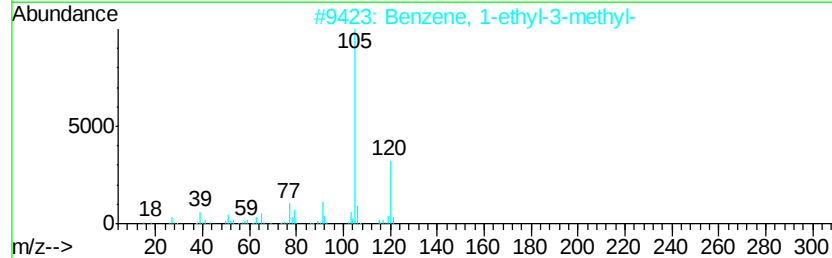
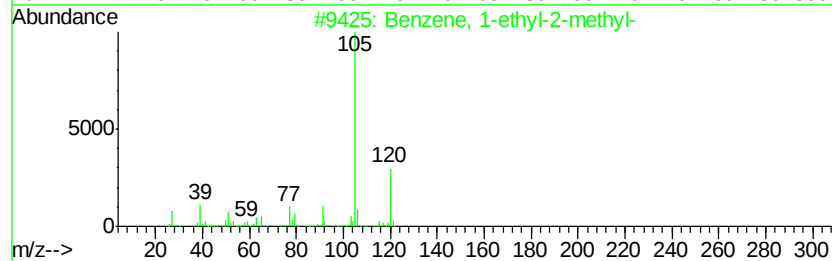
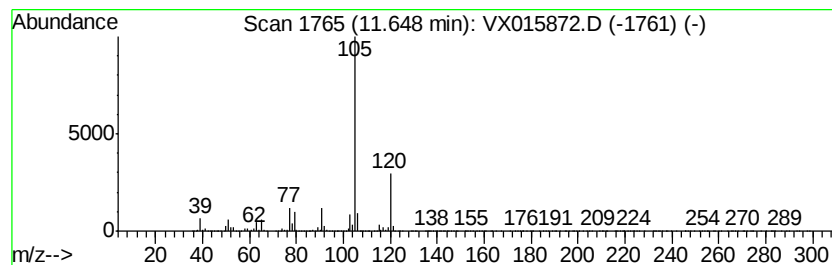
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	61.45 ug/l	5956440	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX042220\
 Data File : VX015872.D
 Acq On : 23 Apr 2020 04:20
 Operator : JC/SP
 Sample : L2380-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
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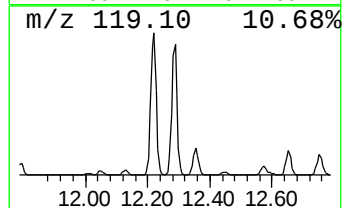
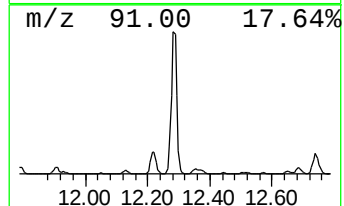
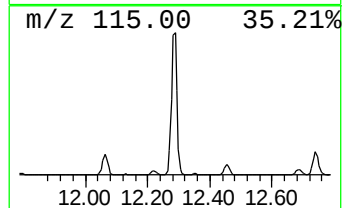
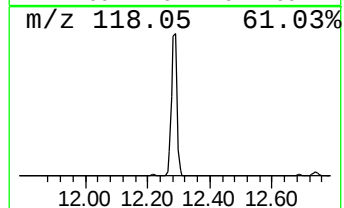
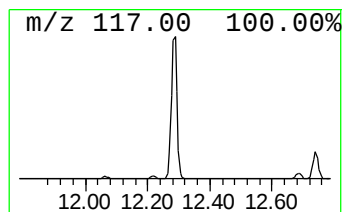
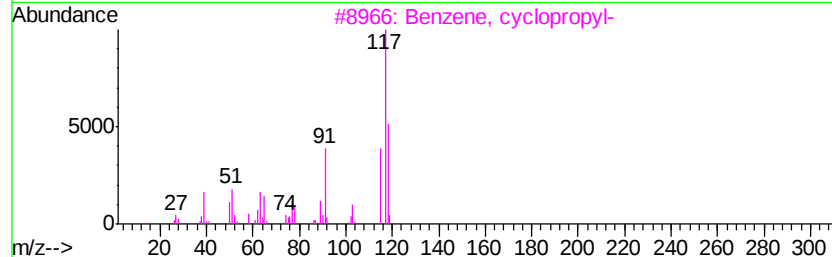
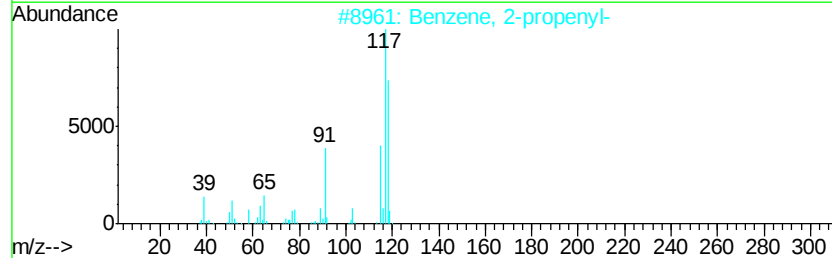
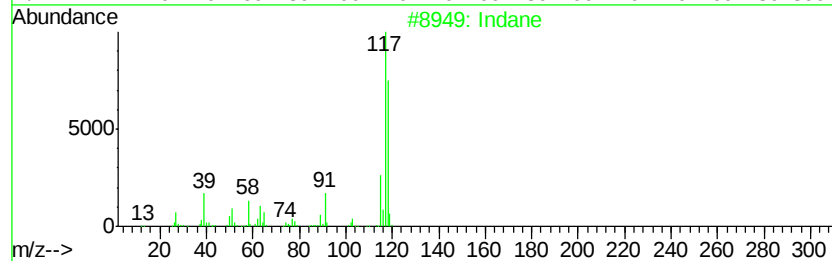
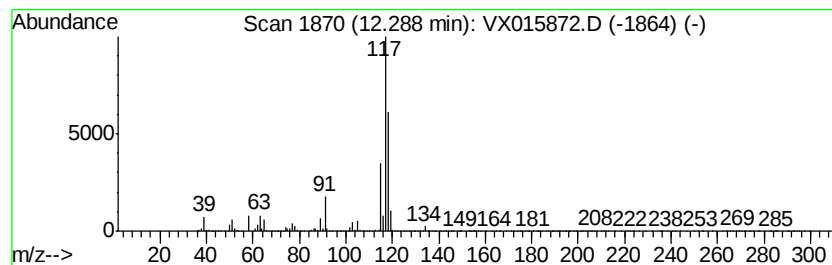
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 Quant Title : SW846 8260

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

 Peak Number 8 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	341.64 ug/l	33113100	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	91
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	Benzene, 1-propenyl-		118	C9H10	000637-50-3	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015872.D
Acq On : 23 Apr 2020 04:20
Operator : JC/SP
Sample : L2380-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-200421

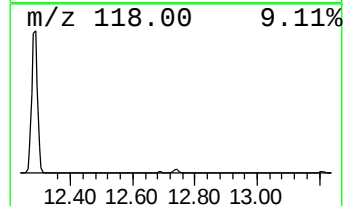
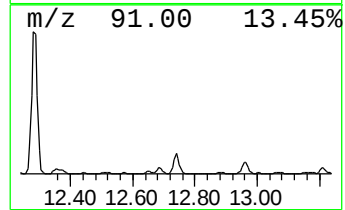
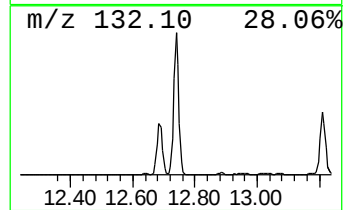
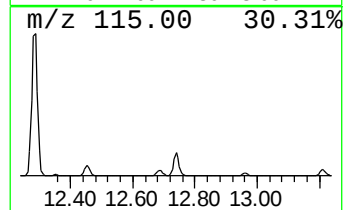
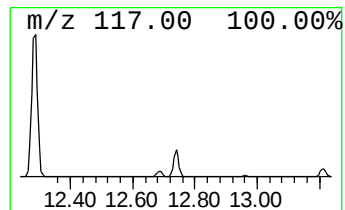
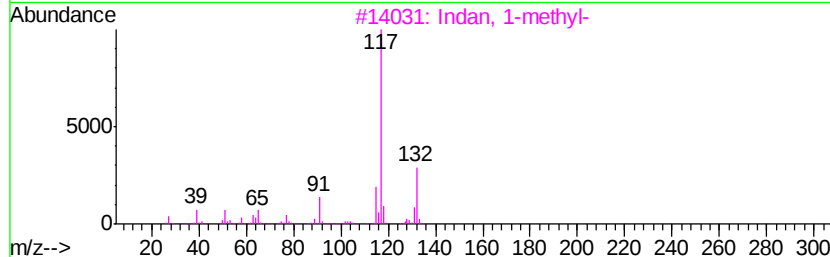
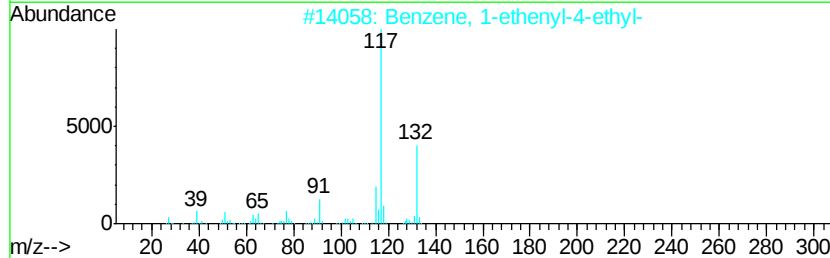
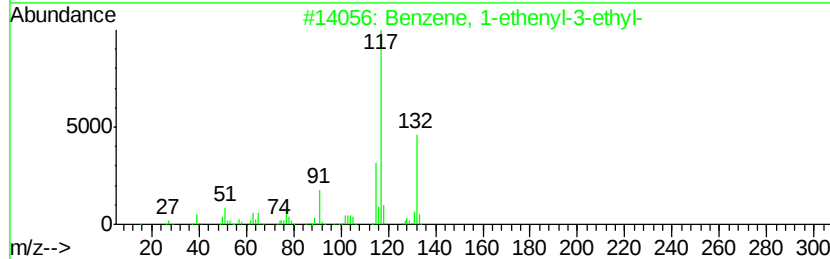
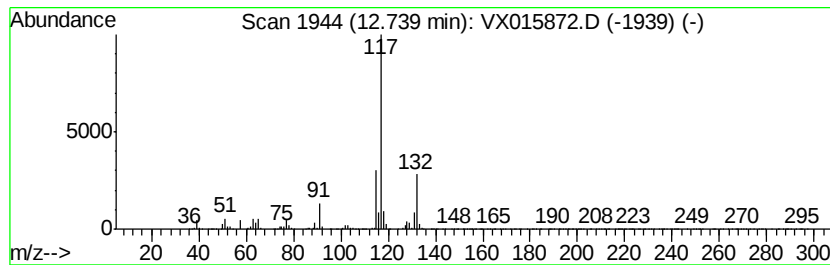
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	53.36 ug/l	5171490	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
3		Indan, 1-methyl-	132	C10H12	000767-58-8	90
4		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX042220\
Data File : VX015872.D
Acq On : 23 Apr 2020 04:20
Operator : JC/SP
Sample : L2380-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-200421

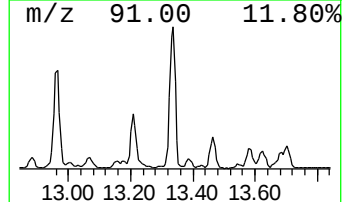
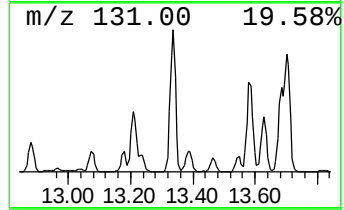
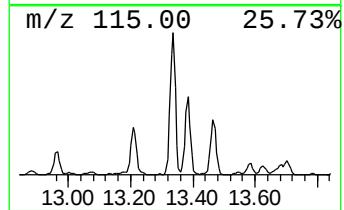
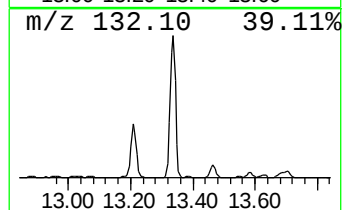
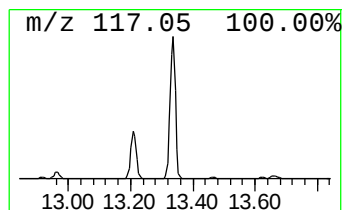
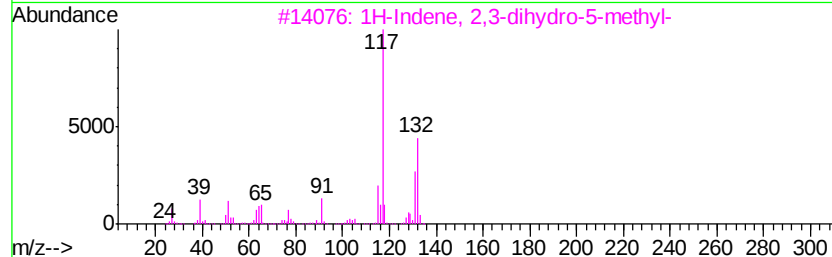
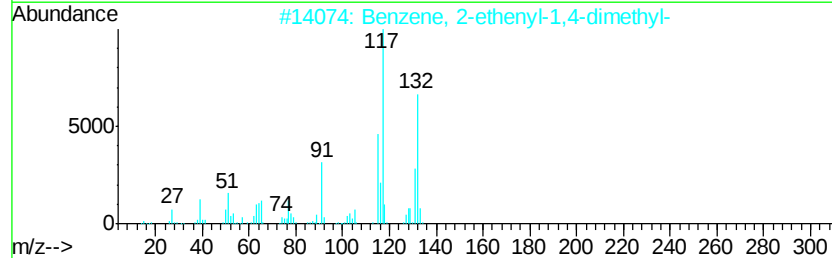
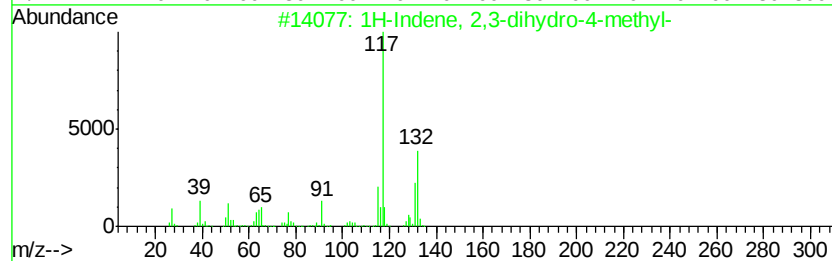
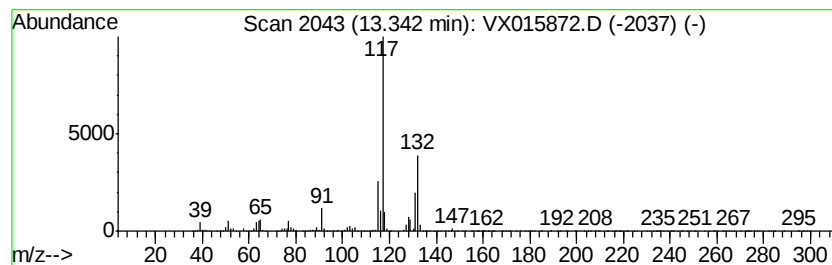
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	51.26 ug/l	4968380	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
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:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX042220\
Data File : VX015872.D
Acq On : 23 Apr 2020 04:20
Operator : JC/SP
Sample : L2380-14
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY001MW031-200421

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042220W.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
Pentane, 2-methyl-	2.87	66.2	ug/l	2928890	1	5.64	2212230	50.0
Cyclopentane, met...	4.38	164.9	ug/l	7294250	1	5.64	2212230	50.0
Pentane, 2,3-dime...	5.70	143.3	ug/l	6341470	1	5.64	2212230	50.0
Hexane, 3-methyl-	5.90	86.8	ug/l	3841210	1	5.64	2212230	50.0
Hexane, 2,2-dimet...	6.33	128.2	ug/l	9813250	2	6.83	3827600	50.0
Benzene, 1-ethyl-...	11.65	61.5	ug/l	5956440	4	12.06	4846260	50.0
Indane	12.29	341.6	ug/l	33113100	4	12.06	4846260	50.0
Benzene, 1-etheny...	12.74	53.4	ug/l	5171490	4	12.06	4846260	50.0
1H-Indene, 2,3-di...	13.34	51.3	ug/l	4968380	4	12.06	4846260	50.0