

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091619\
 Data File : VX012419.D
 Acq On : 16 Sep 2019 20:39
 Operator : JC/SP
 Sample : K4830-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S13-0283

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\82X091619W.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	43	rBV	879173	1038441	71.08%	5.292%
2	2.600	274	281	292	rVB	90947	163814	11.21%	0.835%
3	2.838	311	320	322	rBV	24080	49521	3.39%	0.252%
4	2.881	322	327	330	rVV	40673	87166	5.97%	0.444%
5	2.923	330	334	343	rVV	52485	106356	7.28%	0.542%
6	3.161	364	373	383	rBV	71897	168334	11.52%	0.858%
7	4.393	565	575	588	rVB2	122651	359481	24.61%	1.832%
8	5.490	744	755	761	rBV	100682	291018	19.92%	1.483%
9	5.569	761	768	775	rVV2	132070	392308	26.85%	1.999%
10	5.648	775	781	790	rVB	113446	297785	20.38%	1.518%
11	5.923	811	826	838	rBV5	34542	121987	8.35%	0.622%
12	6.051	838	847	854	rVV	121937	328371	22.48%	1.674%
13	6.136	854	861	871	rVV	254933	672236	46.02%	3.426%
14	6.258	872	881	889	rVV4	29252	108457	7.42%	0.553%
15	6.349	890	896	904	rVV4	28699	81079	5.55%	0.413%
16	6.447	904	912	924	rVB2	33721	96970	6.64%	0.494%
17	6.849	968	978	989	rBV	199172	480347	32.88%	2.448%
18	7.453	1065	1077	1089	rBV	157511	382679	26.20%	1.950%
19	7.727	1116	1122	1133	rVB2	23735	47108	3.22%	0.240%
20	7.843	1133	1141	1150	rVB2	16718	34863	2.39%	0.178%
21	8.026	1165	1171	1180	rVB3	17984	41339	2.83%	0.211%
22	8.386	1225	1230	1236	rVB2	20878	40237	2.75%	0.205%
23	8.550	1252	1257	1268	rVB2	17208	49916	3.42%	0.254%
24	8.660	1268	1275	1278	rBV2	77095	138073	9.45%	0.704%
25	8.709	1278	1283	1291	rVV	560447	922846	63.17%	4.703%
26	8.837	1299	1304	1308	rVV	47154	85183	5.83%	0.434%
27	8.904	1312	1315	1322	rVB	26521	44764	3.06%	0.228%
28	9.038	1331	1337	1343	rVB	87507	148468	10.16%	0.757%
29	9.160	1349	1357	1363	rVB	52324	86094	5.89%	0.439%
30	9.568	1418	1424	1428	rBV3	44012	71149	4.87%	0.363%
31	9.617	1428	1432	1437	rVB	100183	140551	9.62%	0.716%
32	9.672	1437	1441	1446	rBV	64926	96266	6.59%	0.491%
33	10.111	1507	1513	1517	rVV	411449	567387	38.84%	2.892%
34	10.184	1521	1525	1529	rVV	53716	81818	5.60%	0.417%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	10.245	1529	1535	1541	rVV2	27061	49579	3.39%	0.253%
36	10.330	1545	1549	1551	rVV3	35163	54208	3.71%	0.276%
37	10.355	1551	1553	1559	rVB5	37937	63609	4.35%	0.324%
38	10.623	1591	1597	1607	rVB7	31017	98044	6.71%	0.500%
39	10.830	1619	1631	1640	rBV3	68084	136228	9.33%	0.694%
40	10.910	1640	1644	1649	rBV3	44926	78108	5.35%	0.398%
41	10.958	1649	1652	1657	rVB3	29186	41571	2.85%	0.212%
42	11.013	1657	1661	1667	rBV	186803	234456	16.05%	1.195%
43	11.135	1674	1681	1686	rBV	478688	638598	43.71%	3.255%
44	11.178	1686	1688	1692	rVV2	30955	38728	2.65%	0.197%
45	11.251	1695	1700	1707	rVV7	42108	104297	7.14%	0.532%
46	11.355	1713	1717	1723	rVV	150045	193193	13.22%	0.985%
47	11.452	1728	1733	1735	rVV2	22602	35099	2.40%	0.179%
48	11.483	1735	1738	1743	rVB	28198	34427	2.36%	0.175%
49	11.666	1764	1768	1770	rBV2	42039	56675	3.88%	0.289%
50	11.751	1777	1782	1786	rBV4	33034	62688	4.29%	0.319%
51	11.800	1786	1790	1795	rVV	1042745	1318497	90.25%	6.720%
52	11.916	1801	1809	1810	rBV5	20834	36579	2.50%	0.186%
53	11.940	1810	1813	1821	rVB	57646	86006	5.89%	0.438%
54	12.074	1830	1835	1840	rBV	577719	711355	48.69%	3.625%
55	12.141	1840	1846	1851	rVV	1141355	1460866	100.00%	7.445%
56	12.226	1856	1860	1864	rBV	34337	40464	2.77%	0.206%
57	12.294	1866	1871	1876	rVV	617592	759515	51.99%	3.871%
58	12.361	1876	1882	1891	rVB2	93047	176256	12.07%	0.898%
59	12.458	1893	1898	1903	rBV	68569	87100	5.96%	0.444%
60	12.513	1903	1907	1913	rBV	149779	180602	12.36%	0.920%
61	12.580	1913	1918	1921	rBV	98070	143567	9.83%	0.732%
62	12.604	1921	1922	1926	rVV	71488	58761	4.02%	0.299%
63	12.665	1927	1932	1935	rVV	130967	168868	11.56%	0.861%
64	12.696	1935	1937	1942	rVV	82306	109137	7.47%	0.556%
65	12.751	1942	1946	1954	rVV4	180166	405098	27.73%	2.065%
66	12.848	1958	1962	1965	rVV2	42359	55798	3.82%	0.284%
67	12.897	1965	1970	1975	rVB2	98613	153040	10.48%	0.780%
68	12.976	1975	1983	1986	rBV2	99715	188900	12.93%	0.963%
69	13.013	1986	1989	1995	rVB2	148187	183853	12.59%	0.937%
70	13.080	1995	2000	2005	rBV3	88449	165217	11.31%	0.842%
71	13.147	2008	2011	2015	rVV	39570	49877	3.41%	0.254%

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72	13.190	2015	2018	2021	rVV2	82676	110112	7.54%	0.561%
73	13.220	2021	2023	2026	rVV	51575	58140	3.98%	0.296%
74	13.263	2026	2030	2033	rVB2	41007	59305	4.06%	0.302%
75	13.306	2033	2037	2039	rBV2	64144	86356	5.91%	0.440%
76	13.342	2039	2043	2048	rVV2	358524	542902	37.16%	2.767%
77	13.385	2048	2050	2053	rVV	49787	53482	3.66%	0.273%
78	13.421	2053	2056	2059	rVB	35408	35939	2.46%	0.183%
79	13.476	2060	2065	2071	rVB	160255	213094	14.59%	1.086%
80	13.556	2071	2078	2081	rBV5	42894	71199	4.87%	0.363%
81	13.598	2081	2085	2087	rVV	44575	55528	3.80%	0.283%
82	13.635	2087	2091	2096	rVV4	107732	218699	14.97%	1.115%
83	13.696	2097	2101	2102	rVV	54690	81207	5.56%	0.414%
84	13.720	2102	2105	2111	rVB3	86460	153118	10.48%	0.780%
85	13.805	2115	2119	2120	rBV	43207	49611	3.40%	0.253%
86	13.830	2120	2123	2130	rVB2	202000	302472	20.70%	1.542%
87	13.909	2131	2136	2142	rBV3	50882	86822	5.94%	0.442%
88	13.982	2143	2148	2152	rVB2	72118	96312	6.59%	0.491%
89	14.025	2152	2155	2159	rBV3	35166	48756	3.34%	0.248%
90	14.129	2167	2172	2174	rBV2	24469	38682	2.65%	0.197%
91	14.263	2189	2194	2202	rBV4	62658	138250	9.46%	0.705%
92	14.372	2208	2212	2216	rVB	43262	57253	3.92%	0.292%
93	14.427	2216	2221	2225	rVB2	53818	75027	5.14%	0.382%
94	14.537	2234	2239	2251	rBV2	82132	143482	9.82%	0.731%
95	14.689	2257	2264	2267	rVV3	86722	152185	10.42%	0.776%
96	14.726	2267	2270	2274	rVV2	43107	55204	3.78%	0.281%
97	14.836	2283	2288	2292	rVV2	112220	165439	11.32%	0.843%
98	14.958	2302	2308	2315	rVB8	24508	60481	4.14%	0.308%
99	15.244	2348	2355	2363	rBV2	30983	69528	4.76%	0.354%
100	15.384	2373	2378	2384	rBV2	32568	61147	4.19%	0.312%

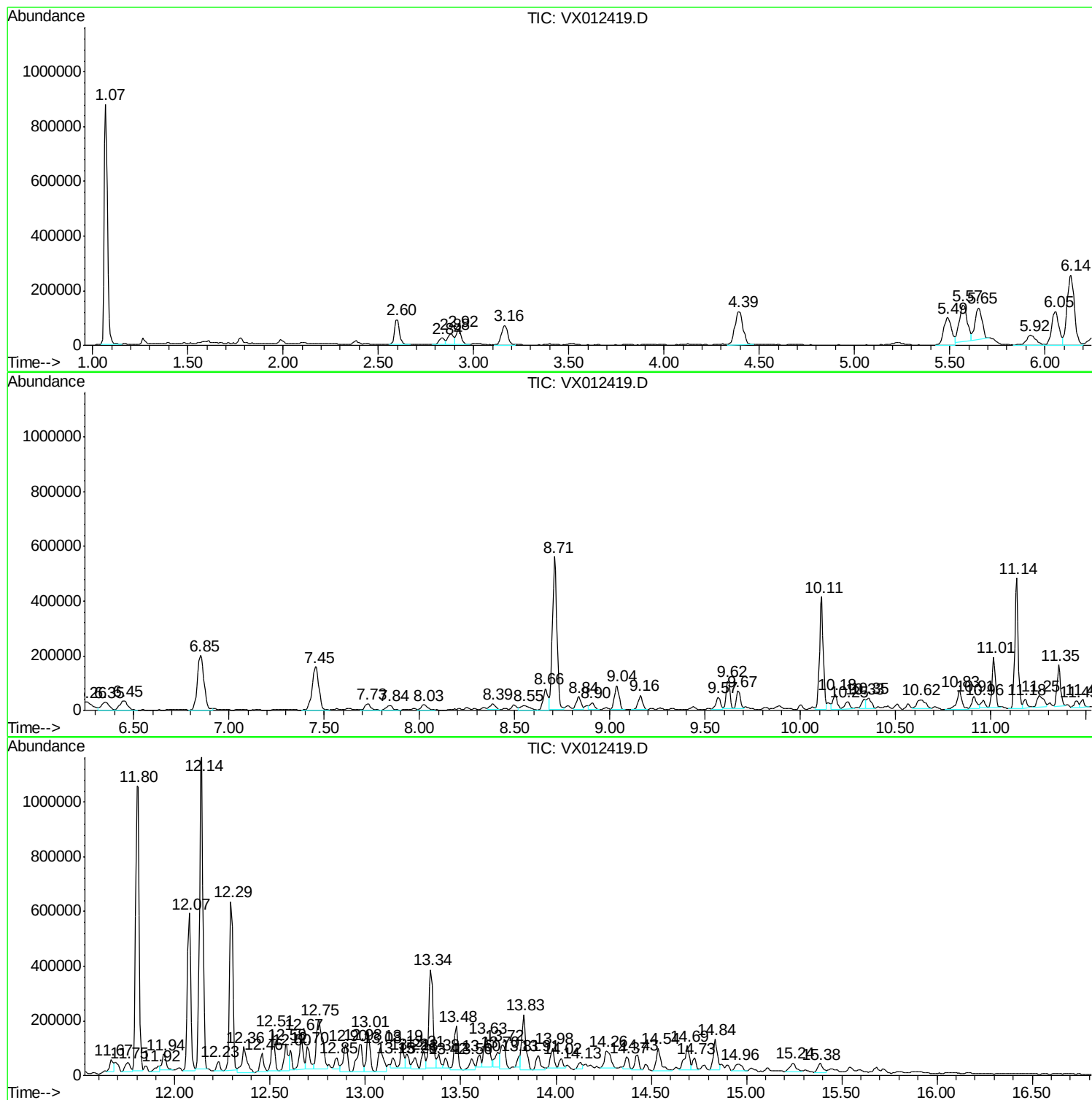
Sum of corrected areas: 19621008

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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



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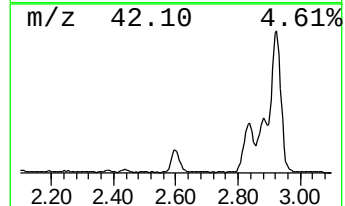
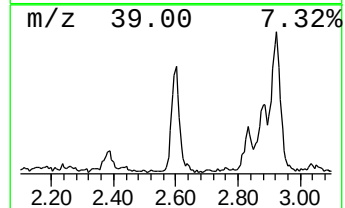
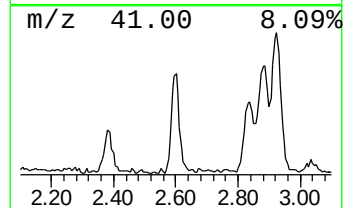
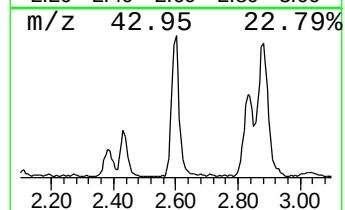
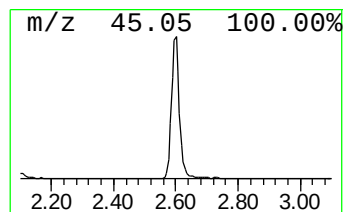
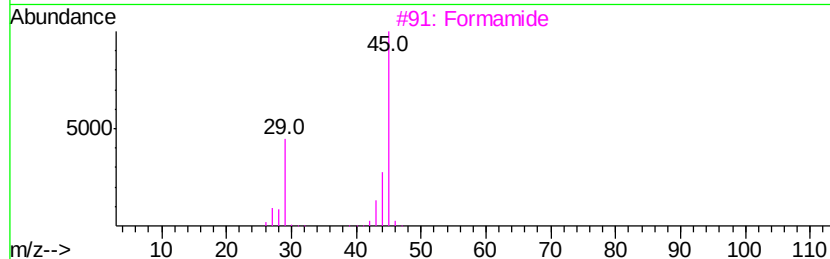
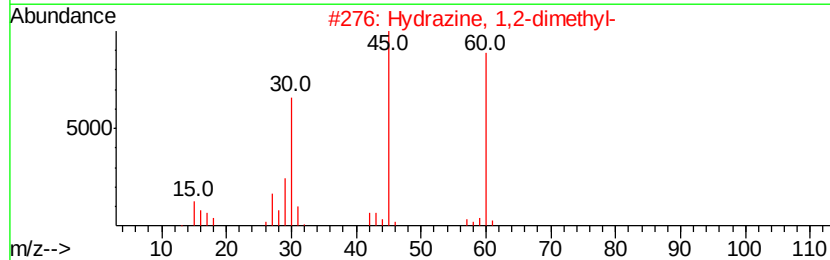
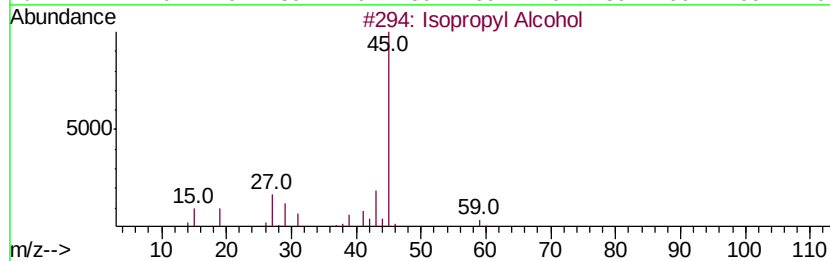
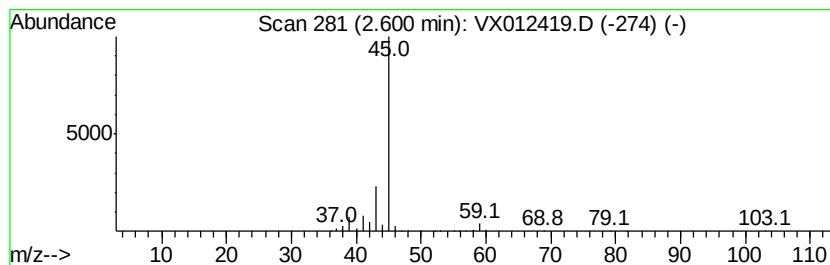
Instrument :
MSVOA_X
ClientSampleId :
S13-0283

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 2 Isopropyl Alcohol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
2.60	27.51 ug/l	163814	Pentafluorobenzene	5.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isopropyl Alcohol	60	C3H8O	000067-63-0	83
2		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3		Formamide	45	CH3NO	000075-12-7	5
4		2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	4
5		Oxirane, (methoxymethyl)-	88	C4H8O2	000930-37-0	4



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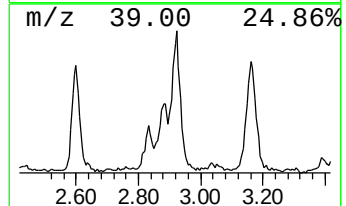
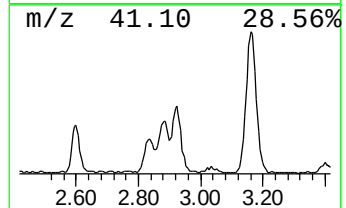
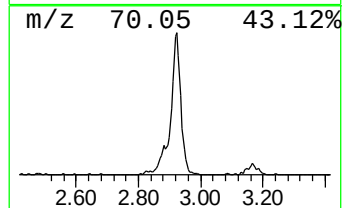
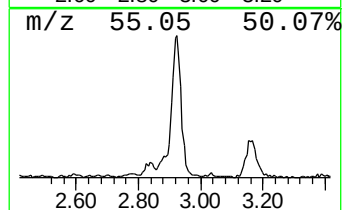
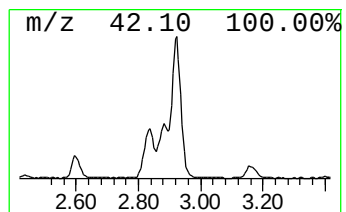
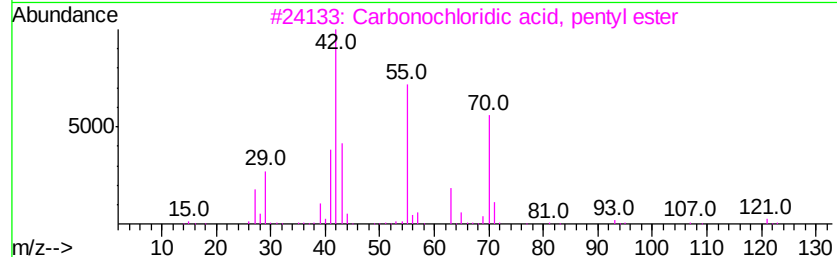
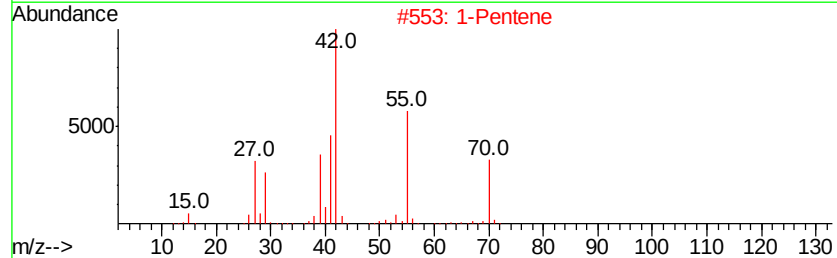
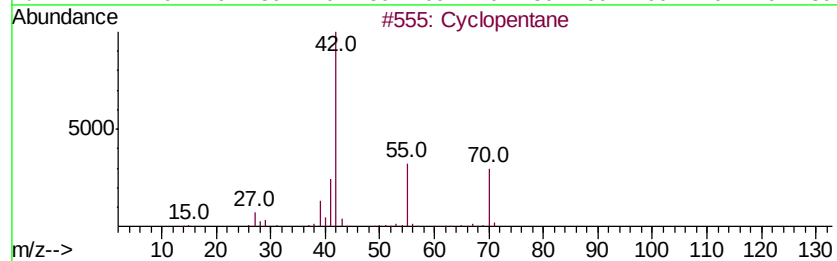
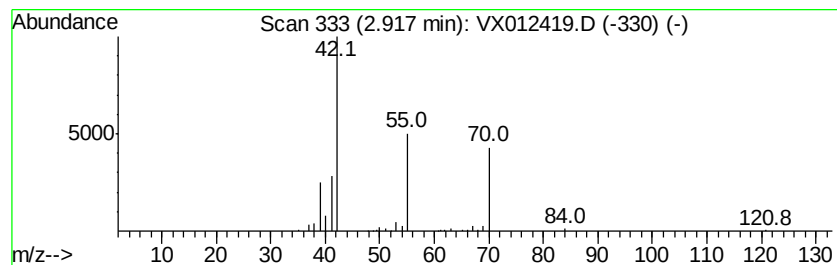
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	17.86 ug/l	106356	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	72
2		1-Pentene	70	C5H10	000109-67-1	59
3		Carbonochloridic acid, pentyl ester	150	C6H11ClO2	000638-41-5	56
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	38
5		2-Pentene	70	C5H10	000109-68-2	27



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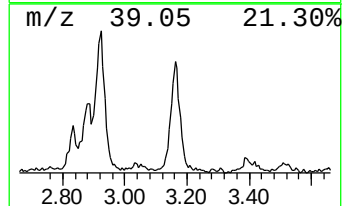
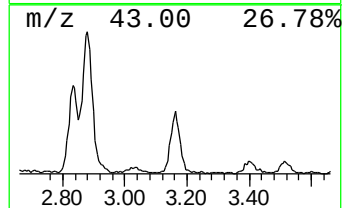
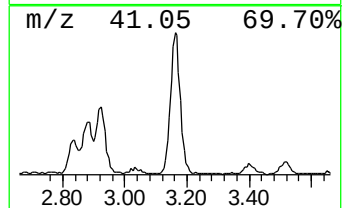
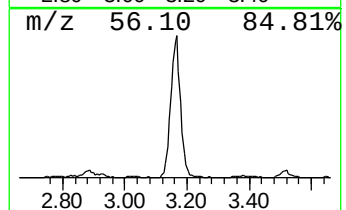
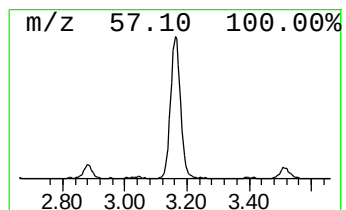
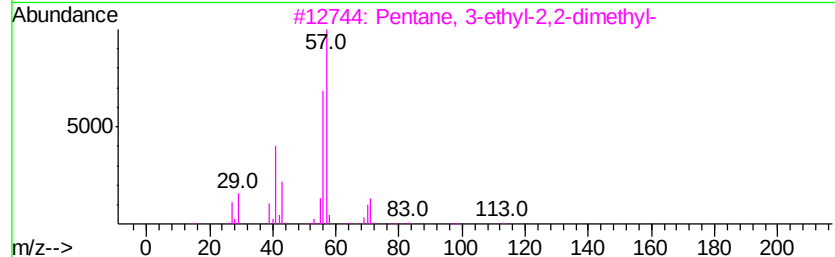
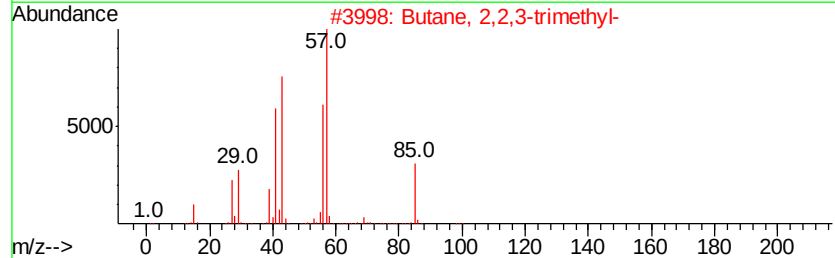
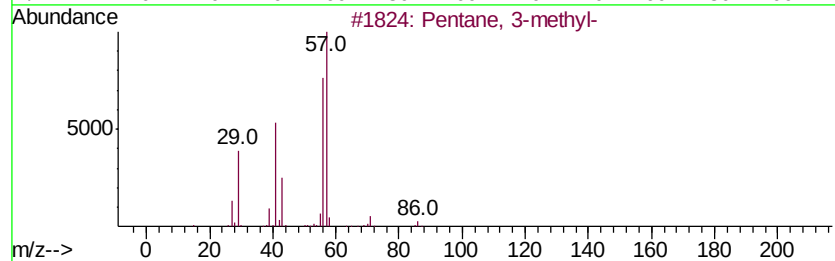
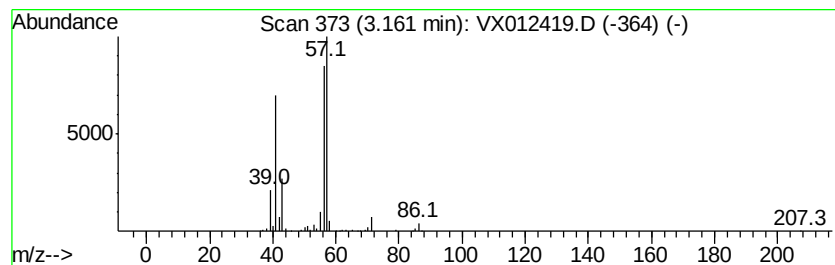
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	28.26 ug/l	168334	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
3		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56
5		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0283

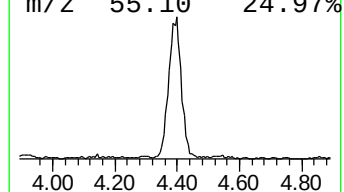
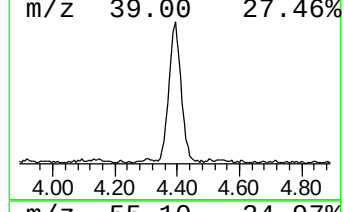
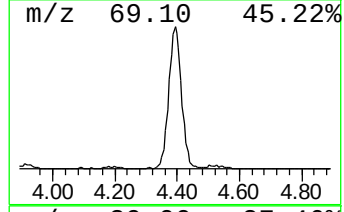
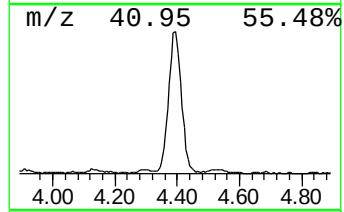
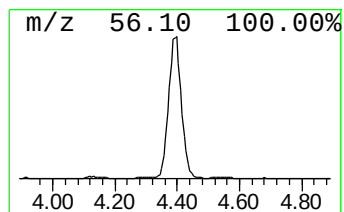
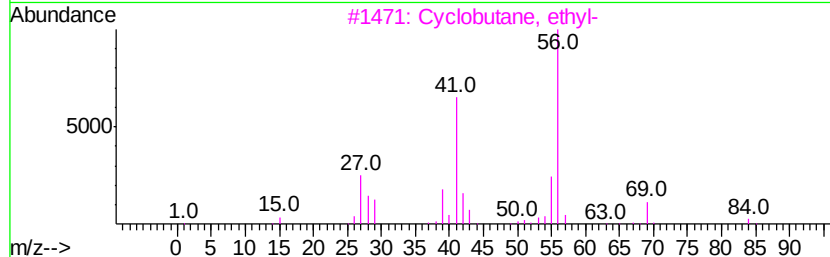
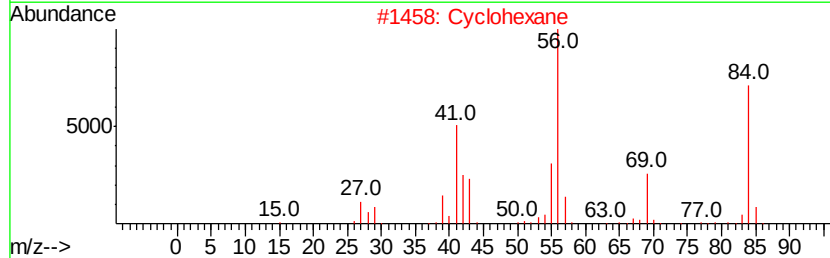
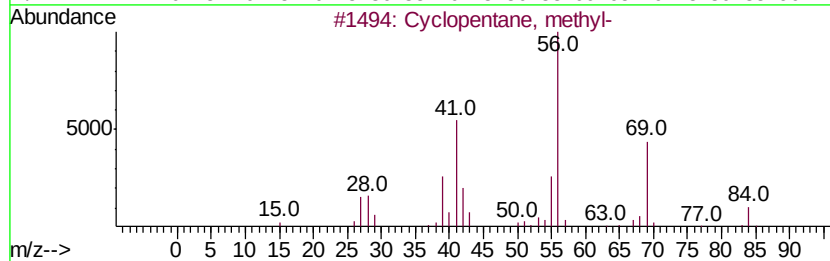
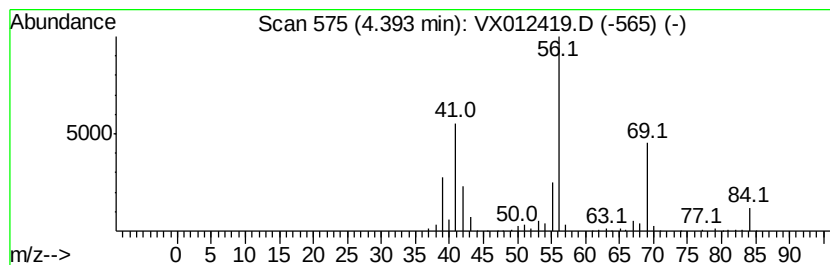
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	60.36 ug/l	359481	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		Cyclohexane	84	C6H12	000110-82-7	78
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59
5		Cyclobutane	56	C4H8	000287-23-0	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0283

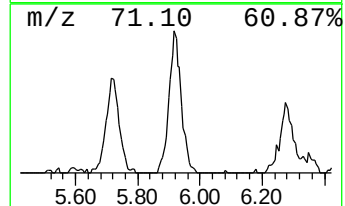
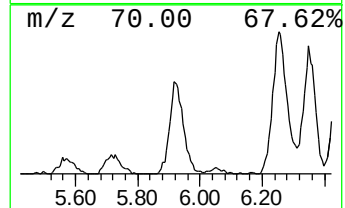
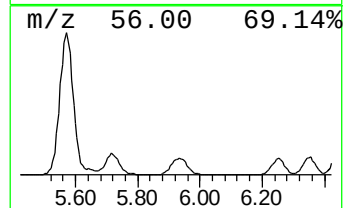
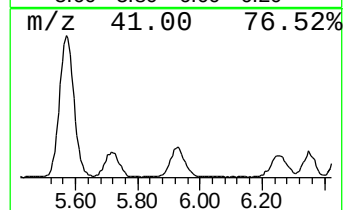
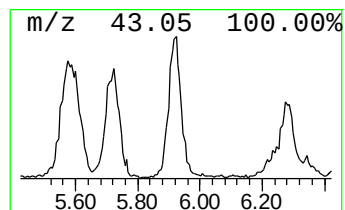
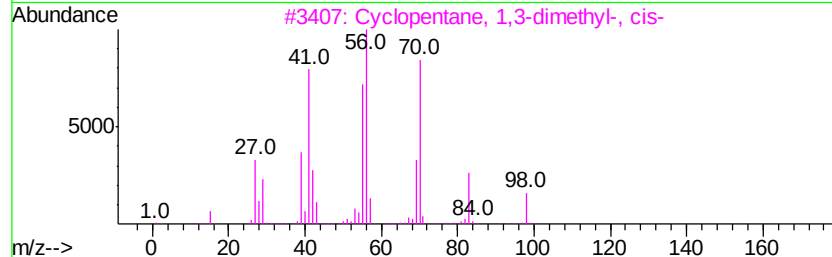
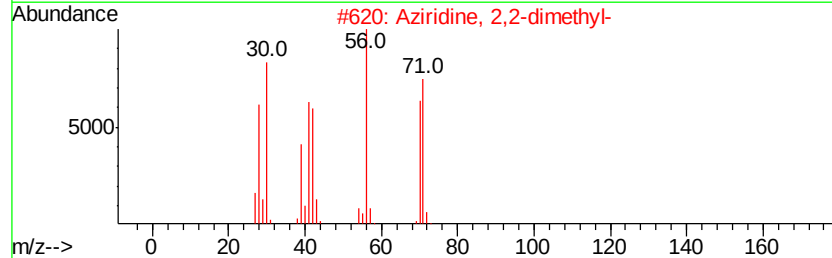
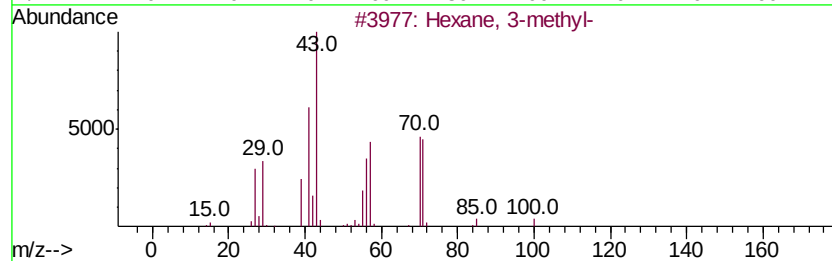
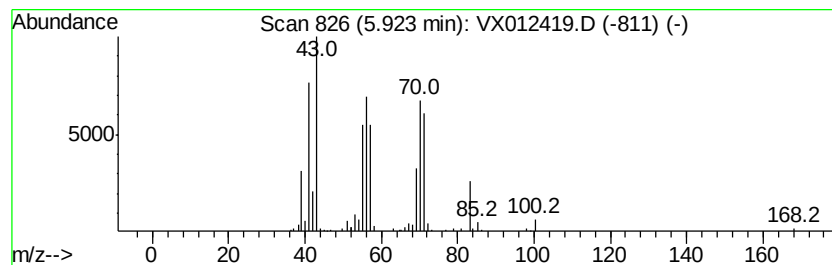
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	20.48 ug/l	121987	Pentafluorobenzene	5.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16		000589-34-4	87
2	Aziridine, 2,2-dimethyl-	71	C4H9N		002658-24-4	53
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14		002532-58-3	50
4	Propanoic acid, 2-methyl-, 3-met...	158	C9H18O2		002050-01-3	47
5	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14		001192-18-3	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0283

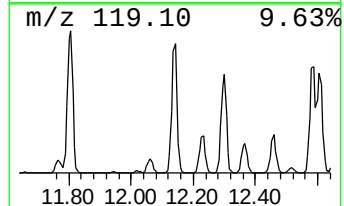
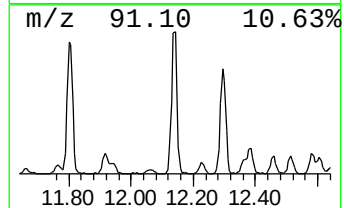
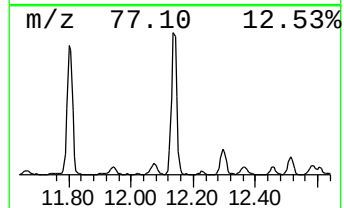
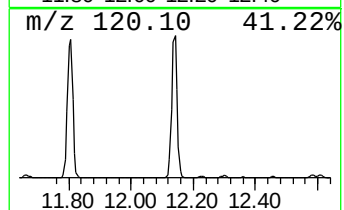
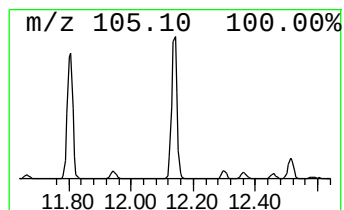
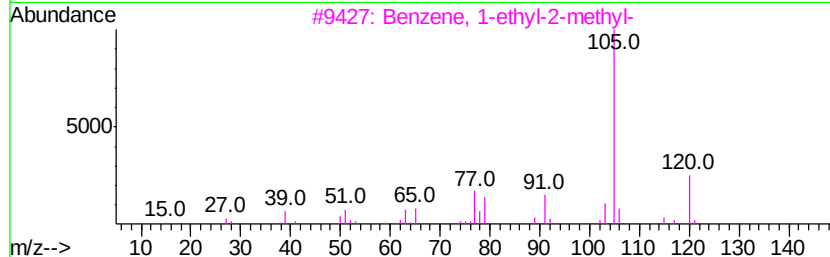
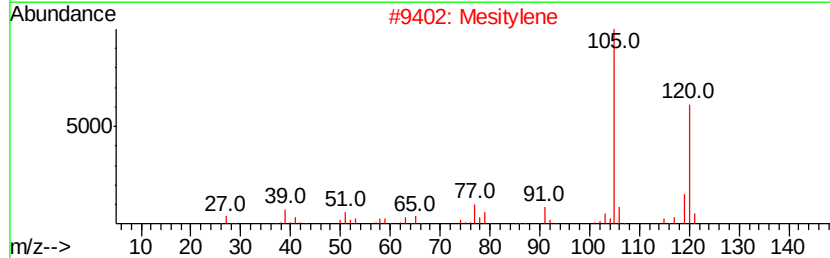
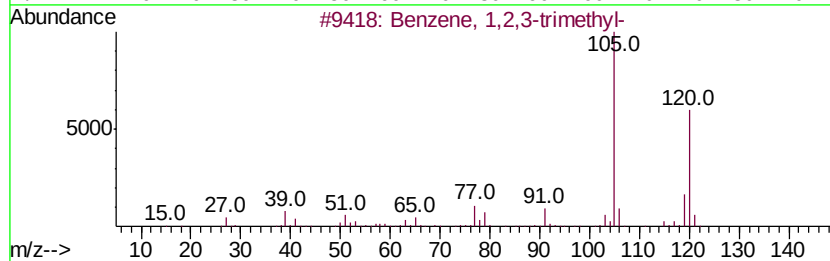
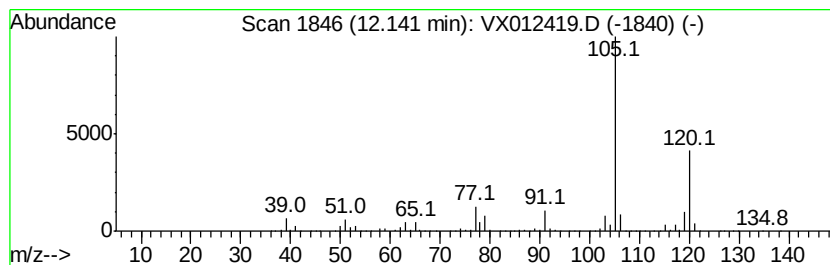
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	102.68 ug/l	1460870	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	93
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0283

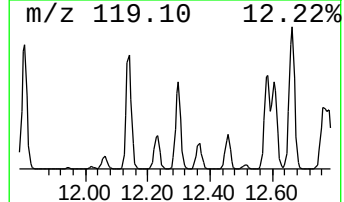
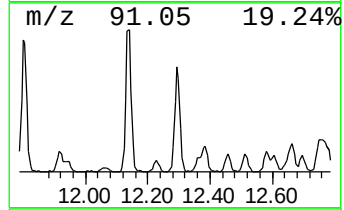
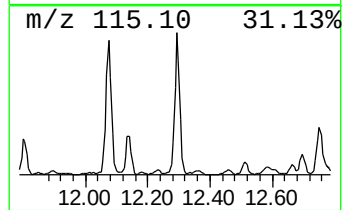
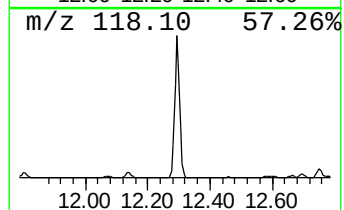
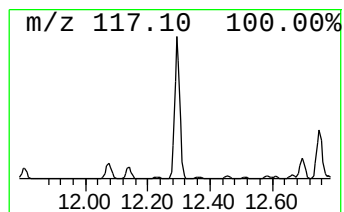
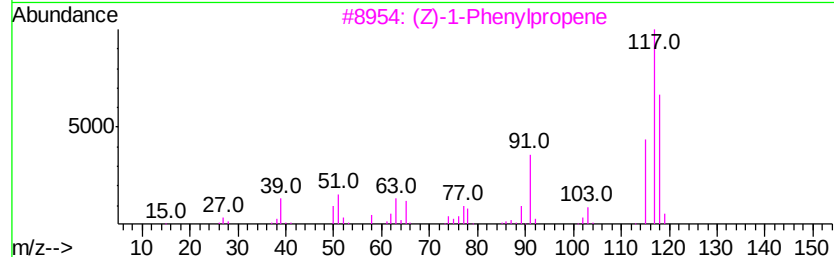
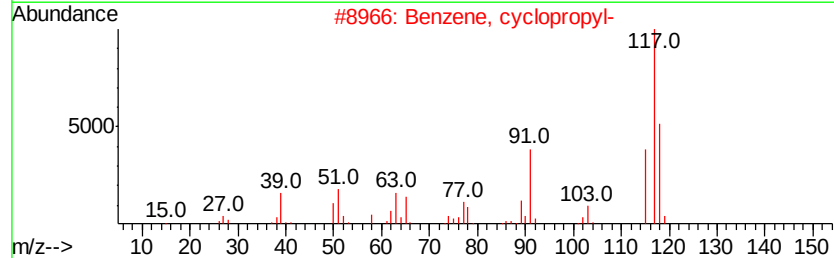
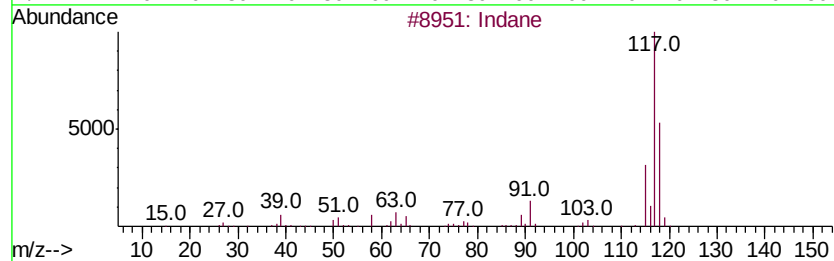
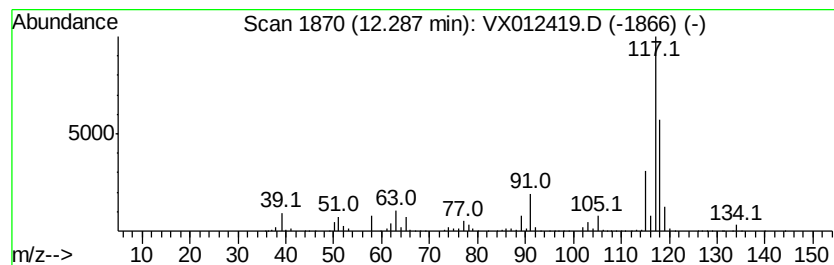
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Quant Title : SW846 8260

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

Peak Number 8 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	53.39 ug/l	759515	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	76
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3		(Z)-1-Phenylpropene	118	C9H10	000766-90-5	76
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

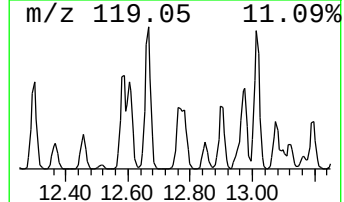
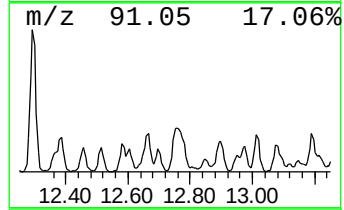
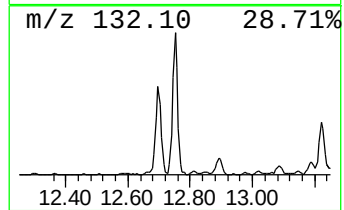
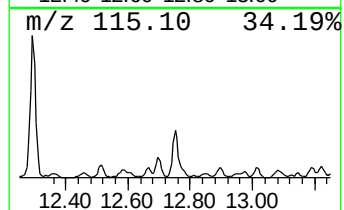
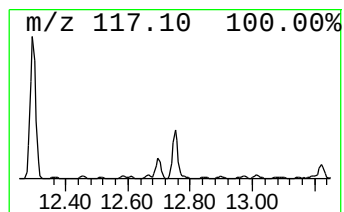
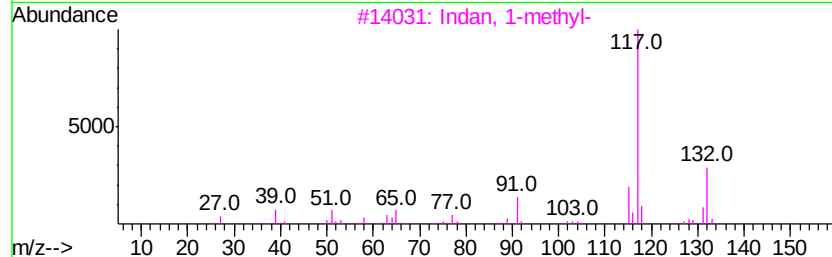
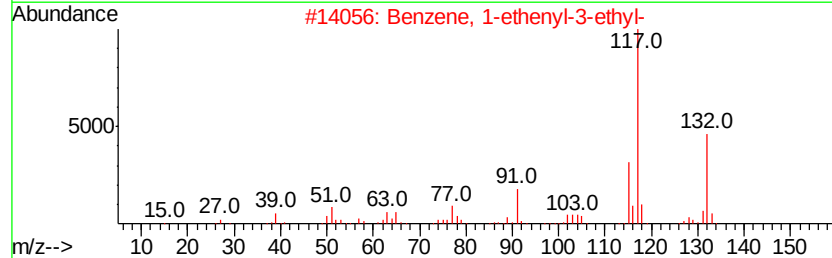
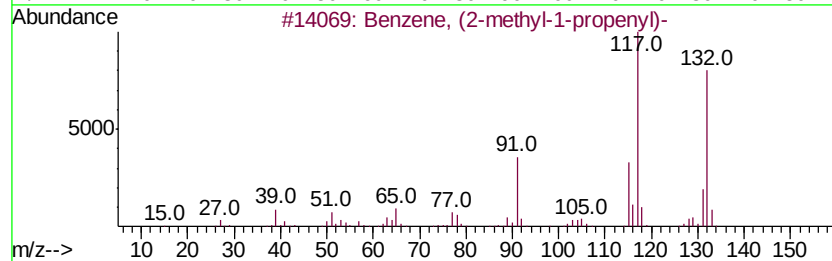
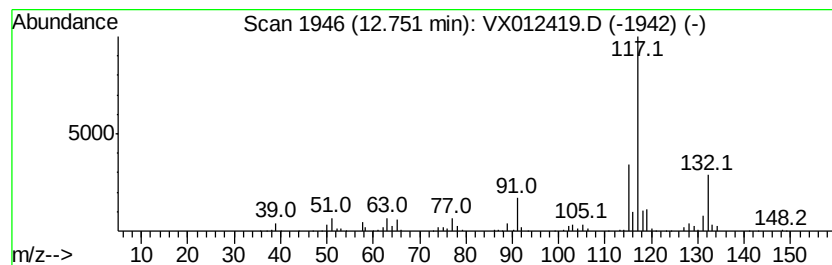
Instrument :
MSVOA_X
ClientSampled :
S13-0283

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (2-methyl-1-propen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	28.47 ug/l	405098	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 91
2		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 90
3		Indan, 1-methyl-	132 C10H12	000767-58-8 87
4		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 80
5		Benzene, 2-butenyl-	132 C10H12	001560-06-1 74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

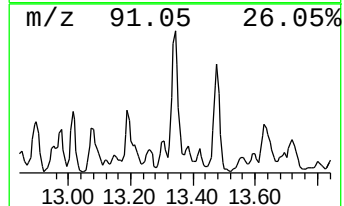
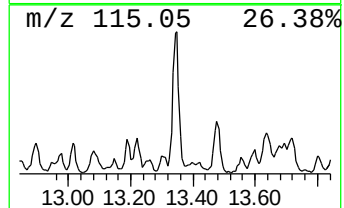
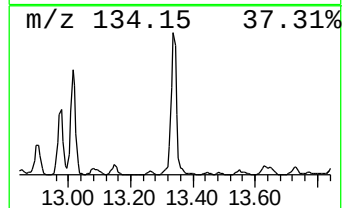
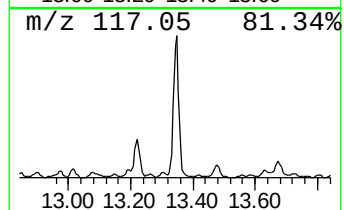
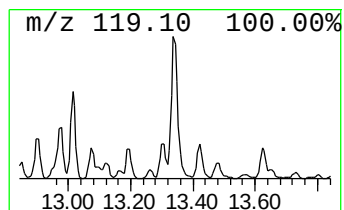
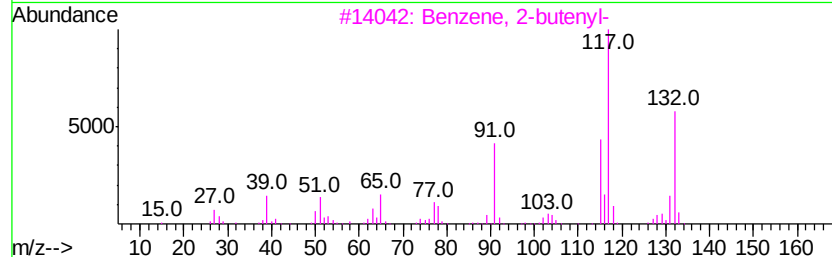
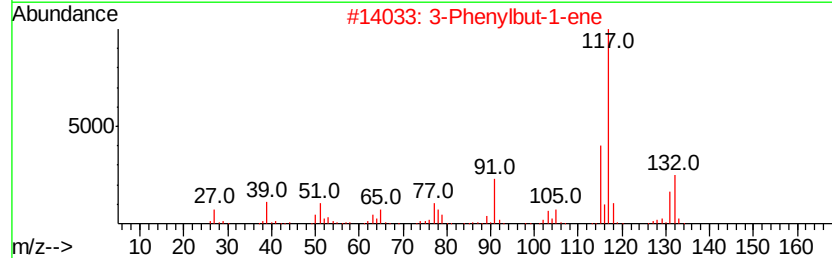
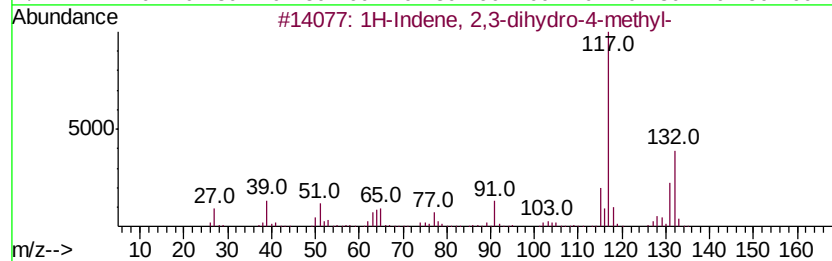
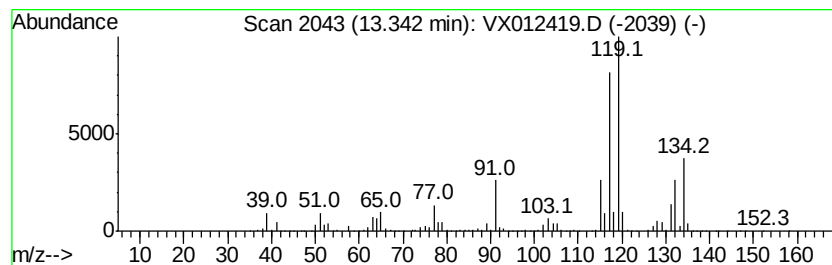
Instrument :
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ClientSampleId :
S13-0283

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
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 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.34	38.16 ug/l	542902	1,4-Dichlorobenzene-d4	12.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX091619\
Data File : VX012419.D
Acq On : 16 Sep 2019 20:39
Operator : JC/SP
Sample : K4830-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0283

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.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
Isopropyl Alcohol	2.60	27.5	ug/l	163814	1	5.65	297785	50.0
Cyclopentane	2.92	17.9	ug/l	106356	1	5.65	297785	50.0
Pentane, 3-methyl-	3.16	28.3	ug/l	168334	1	5.65	297785	50.0
Cyclopentane, met...	4.39	60.4	ug/l	359481	1	5.65	297785	50.0
Hexane, 3-methyl-	5.92	20.5	ug/l	121987	1	5.65	297785	50.0
Benzene, 1,2,3-tr...	12.14	102.7	ug/l	1460870	4	12.07	711355	50.0
Indane	12.29	53.4	ug/l	759515	4	12.07	711355	50.0
Benzene, (2-methy...	12.75	28.5	ug/l	405098	4	12.07	711355	50.0
1H-Indene, 2,3-di...	13.34	38.2	ug/l	542902	4	12.07	711355	50.0