

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091619\
 Data File : VX012420.D
 Acq On : 16 Sep 2019 21:02
 Operator : JC/SP
 Sample : K4830-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S13-0280

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	42	rBV	746809	880651	6.00%	0.759%
2	1.777	141	146	157	rVB	764682	1090167	7.43%	0.939%
3	1.990	175	181	191	rVB	738671	1023335	6.97%	0.882%
4	2.381	237	245	258	rVB	190741	387926	2.64%	0.334%
5	2.832	310	319	322	rBV	260554	533872	3.64%	0.460%
6	2.881	322	327	330	rVV	808831	1637287	11.15%	1.411%
7	2.923	330	334	357	rVB	922551	1893661	12.90%	1.632%
8	3.161	362	373	393	rBV	865751	1994153	13.58%	1.718%
9	3.515	420	431	443	rBV	313317	707111	4.82%	0.609%
10	4.392	565	575	589	rVB	1406613	4124247	28.10%	3.554%
11	5.228	698	712	735	rVB3	49458	217933	1.48%	0.188%
12	5.490	741	755	759	rBV	98243	277722	1.89%	0.239%
13	5.575	759	769	779	rBV2	1071083	3465406	23.61%	2.986%
14	5.715	787	792	811	rVB	159341	487344	3.32%	0.420%
15	5.923	814	826	839	rBV5	319839	1039881	7.08%	0.896%
16	6.051	839	847	855	rVV	119422	310495	2.12%	0.268%
17	6.130	855	860	870	rVV2	67423	186365	1.27%	0.161%
18	6.258	870	881	890	rVV3	177521	652215	4.44%	0.562%
19	6.349	890	896	904	rVV	172573	477237	3.25%	0.411%
20	6.447	904	912	927	rVB	282865	781800	5.33%	0.674%
21	6.849	968	978	987	rBV	200749	486741	3.32%	0.419%
22	7.459	1065	1078	1090	rBV	1703068	3942925	26.86%	3.397%
23	7.727	1115	1122	1133	rVB	198186	387838	2.64%	0.334%
24	7.843	1133	1141	1149	rVB3	84888	174442	1.19%	0.150%
25	8.026	1163	1171	1187	rVB2	83142	192754	1.31%	0.166%
26	8.380	1225	1229	1236	rVB3	82034	157387	1.07%	0.136%
27	8.660	1267	1275	1279	rBV2	340510	632015	4.31%	0.545%
28	8.709	1279	1283	1290	rVB	582563	1002400	6.83%	0.864%
29	8.837	1298	1304	1308	rBV	160616	269069	1.83%	0.232%
30	8.904	1312	1315	1321	rVB2	105006	170221	1.16%	0.147%
31	9.038	1331	1337	1343	rVB	321819	526414	3.59%	0.454%
32	9.160	1348	1357	1363	rBV	177861	284380	1.94%	0.245%
33	9.568	1417	1424	1428	rBV3	144438	242600	1.65%	0.209%
34	9.617	1428	1432	1437	rVB	335541	465692	3.17%	0.401%

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35	9.672	1437	1441	1446	rBV	229871	344140	2.34%	0.297%
36	10.111	1508	1513	1518	rVB	420041	546382	3.72%	0.471%
37	10.184	1518	1525	1529	rBV	190462	285516	1.94%	0.246%
38	10.251	1529	1536	1545	rBV	11084305	14679575	100.00%	12.649%
39	10.355	1545	1553	1559	rVB	2116390	3029775	20.64%	2.611%
40	10.641	1591	1600	1605	rBV3	75547	199675	1.36%	0.172%
41	10.696	1605	1609	1619	rVB	415910	574410	3.91%	0.495%
42	10.830	1620	1631	1638	rBV2	145869	251088	1.71%	0.216%
43	10.910	1639	1644	1650	rBV3	93283	157601	1.07%	0.136%
44	11.013	1656	1661	1674	rVB	1512812	1943768	13.24%	1.675%
45	11.135	1674	1681	1685	rBV	465118	611297	4.16%	0.527%
46	11.355	1712	1717	1724	rBV	1792881	2185807	14.89%	1.883%
47	11.428	1725	1729	1735	rVV	244536	431687	2.94%	0.372%
48	11.501	1735	1741	1750	rVB	966195	1233673	8.40%	1.063%
49	11.666	1758	1768	1777	rBV	3843177	4930934	33.59%	4.249%
50	11.806	1786	1791	1796	rVB	11066146	13423768	91.45%	11.567%
51	11.916	1801	1809	1811	rBV2	88079	155189	1.06%	0.134%
52	11.940	1811	1813	1820	rVB	242961	307401	2.09%	0.265%
53	12.074	1829	1835	1840	rVB2	639462	1044723	7.12%	0.900%
54	12.141	1840	1846	1851	rBV	5398489	6701846	45.65%	5.775%
55	12.226	1856	1860	1866	rBV	127316	163000	1.11%	0.140%
56	12.293	1866	1871	1878	rBV	3166072	4018841	27.38%	3.463%
57	12.367	1878	1883	1893	rVB3	663032	1127611	7.68%	0.972%
58	12.458	1893	1898	1903	rBV2	296139	402037	2.74%	0.346%
59	12.513	1903	1907	1913	rBV	766616	897606	6.11%	0.773%
60	12.580	1913	1918	1920	rBV	812753	993063	6.76%	0.856%
61	12.604	1920	1922	1927	rVV	837310	942433	6.42%	0.812%
62	12.665	1927	1932	1935	rVV	1122905	1312536	8.94%	1.131%
63	12.696	1935	1937	1942	rVV	355145	394965	2.69%	0.340%
64	12.751	1942	1946	1954	rVB2	901562	1463779	9.97%	1.261%
65	12.897	1965	1970	1975	rVB2	615103	824907	5.62%	0.711%
66	12.976	1975	1983	1986	rBV	681203	966904	6.59%	0.833%
67	13.013	1986	1989	1995	rVB	1167732	1373530	9.36%	1.184%
68	13.080	1995	2000	2005	rBV3	160689	292255	1.99%	0.252%
69	13.190	2015	2018	2020	rBV2	158894	179813	1.22%	0.155%
70	13.220	2020	2023	2027	rVV	505308	565057	3.85%	0.487%
71	13.305	2033	2037	2039	rBV2	169242	238820	1.63%	0.206%

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72	13.342	2039	2043	2048	rVB2	2186572	2944875	20.06%	2.538%
73	13.476	2060	2065	2072	rVB	1140875	1416882	9.65%	1.221%
74	13.555	2073	2078	2081	rBV3	115059	162388	1.11%	0.140%
75	13.598	2081	2085	2087	rVV	156107	221985	1.51%	0.191%
76	13.635	2087	2091	2096	rVV4	303285	623615	4.25%	0.537%
77	13.696	2097	2101	2102	rVV	170454	249838	1.70%	0.215%
78	13.720	2102	2105	2111	rVB2	264869	430671	2.93%	0.371%
79	13.830	2115	2123	2131	rVB	3711202	4747289	32.34%	4.091%
80	13.909	2131	2136	2140	rVB	156184	220787	1.50%	0.190%
81	13.982	2140	2148	2152	rBV2	232016	362360	2.47%	0.312%
82	14.281	2190	2197	2202	rVB2	313308	555686	3.79%	0.479%
83	14.537	2234	2239	2247	rBV2	279768	395802	2.70%	0.341%
84	14.689	2257	2264	2268	rBV	1222995	1589451	10.83%	1.370%
85	14.836	2282	2288	2293	rBV	928271	1204601	8.21%	1.038%
86	15.543	2398	2404	2409	rBV	133824	199546	1.36%	0.172%
87	15.683	2420	2427	2430	rBV	162524	262413	1.79%	0.226%
88	15.720	2430	2433	2440	rVB	112380	166823	1.14%	0.144%

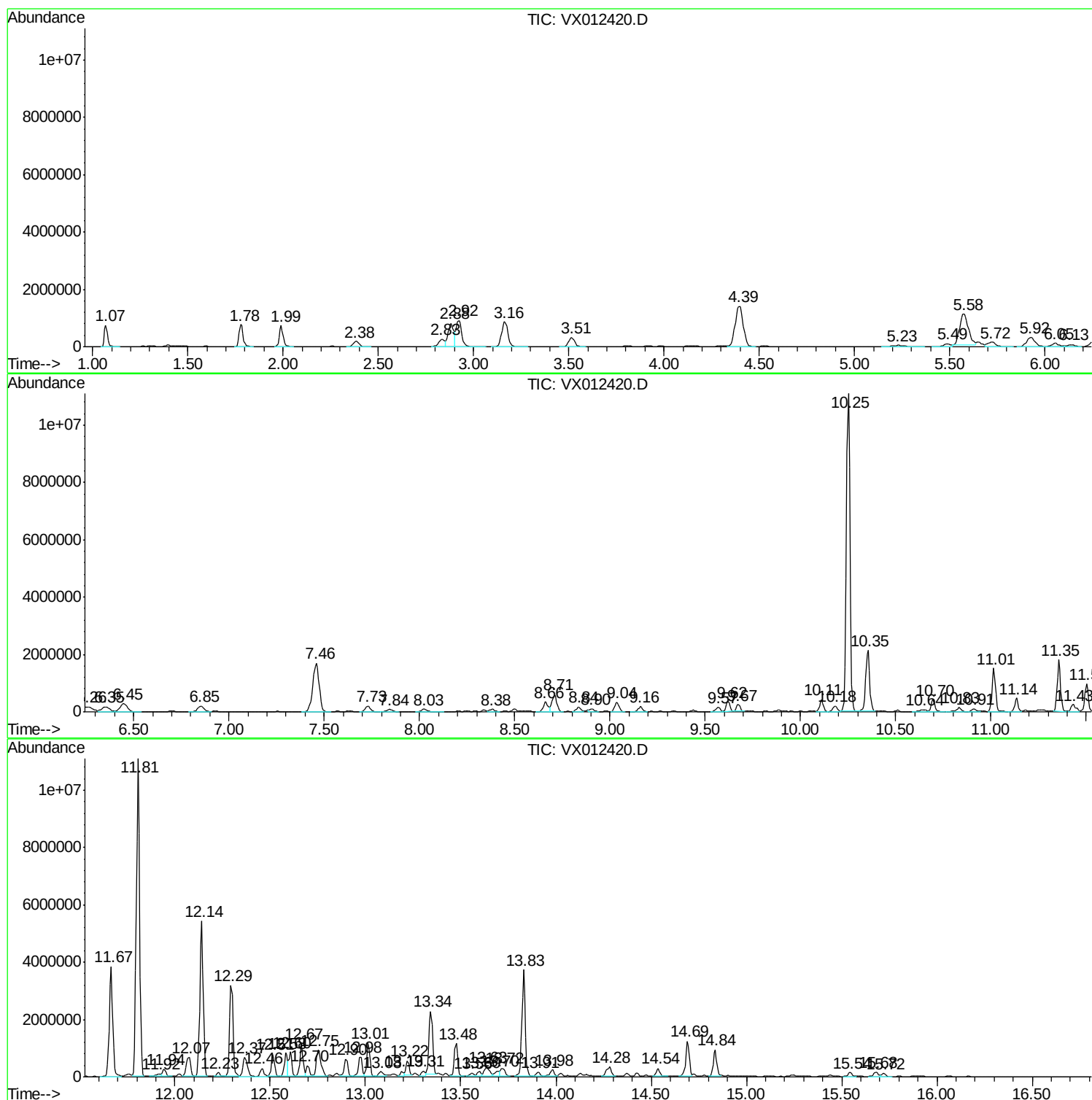
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampled :
S13-0280

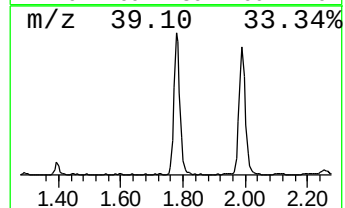
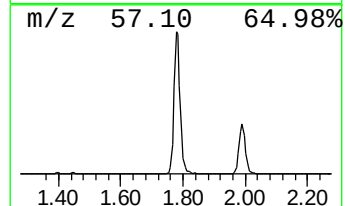
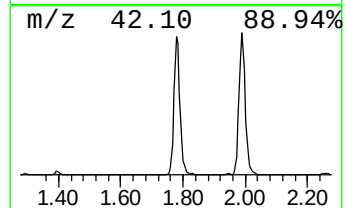
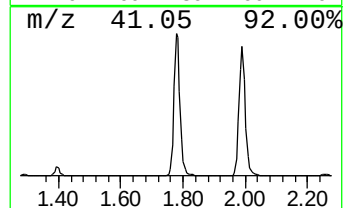
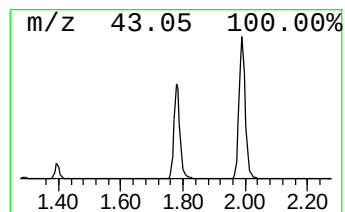
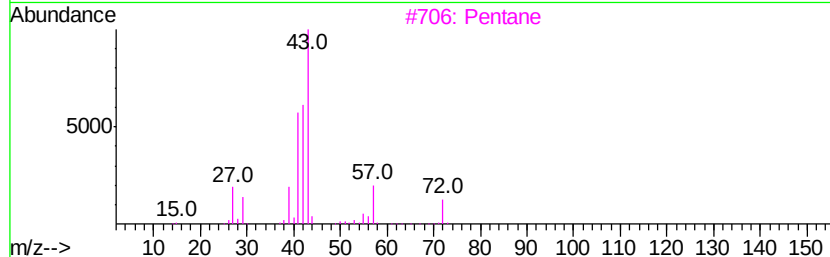
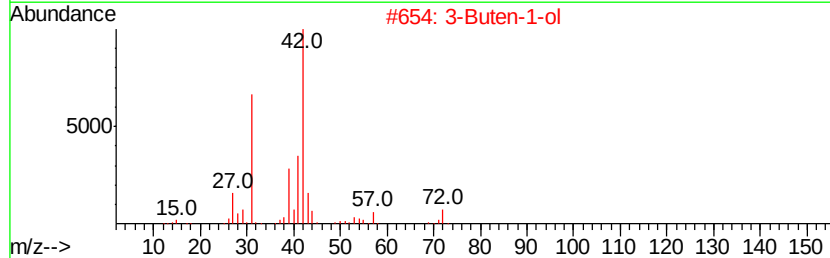
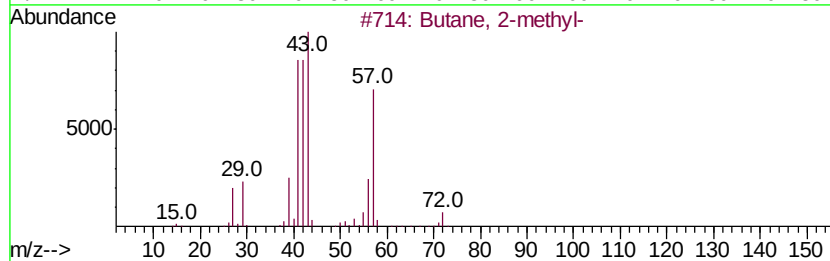
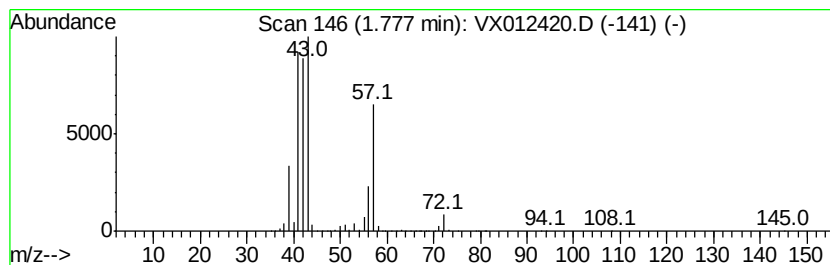
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.78	111.85 ug/l	1090170	Pentafluorobenzene	5.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	94
2		3-Buten-1-ol	72	C4H8O	000627-27-0	47
3		Pentane	72	C5H12	000109-66-0	36
4		1-Propene, 2-methyl-	56	C4H8	000115-11-7	9
5		Azetidine	57	C3H7N	000503-29-7	9



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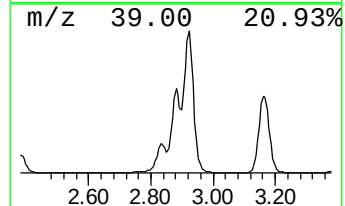
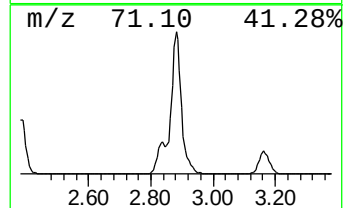
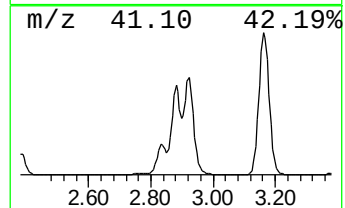
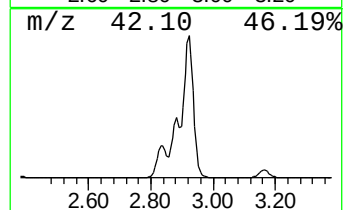
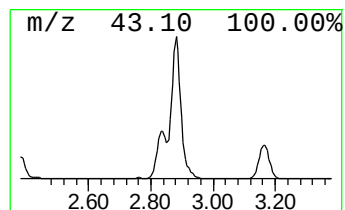
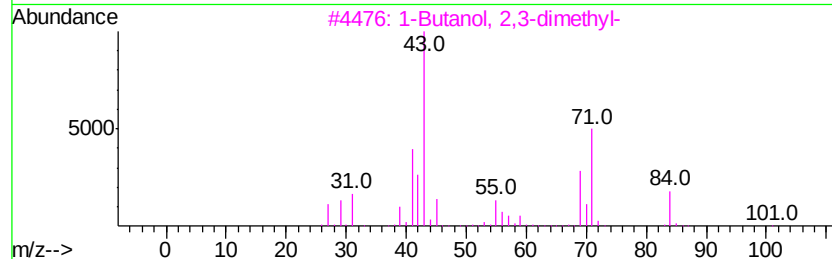
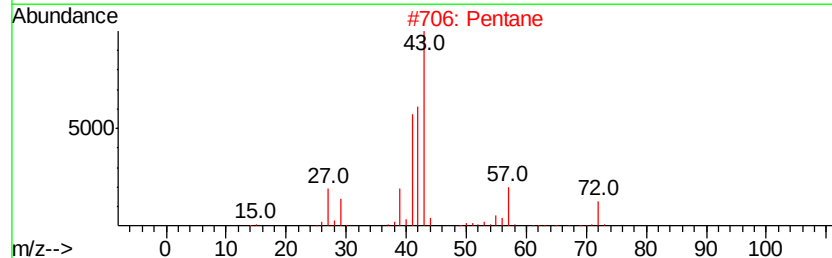
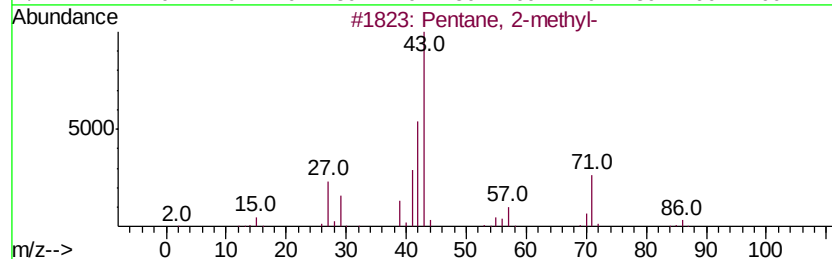
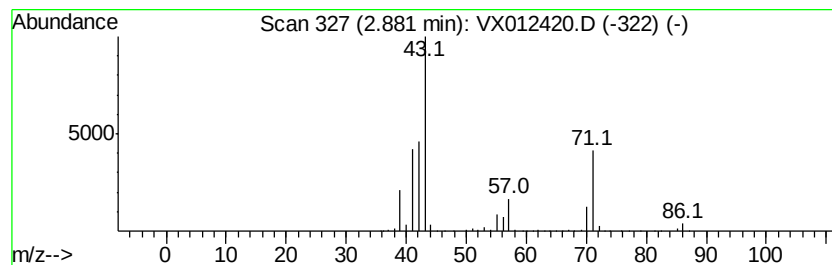
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	167.98 ug/l	1637290	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2		Pentane	72	C5H12	000109-66-0	46
3		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
4		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28



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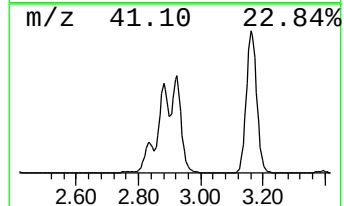
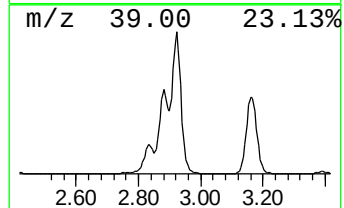
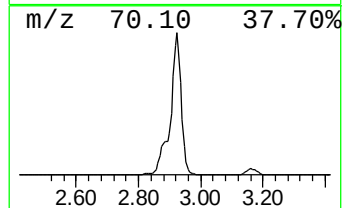
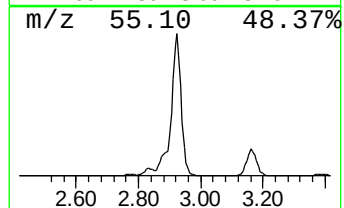
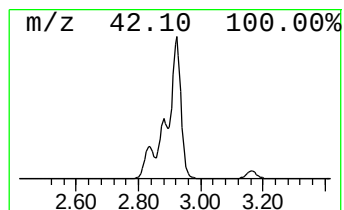
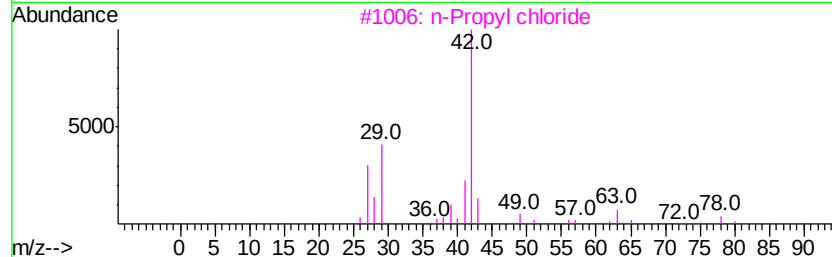
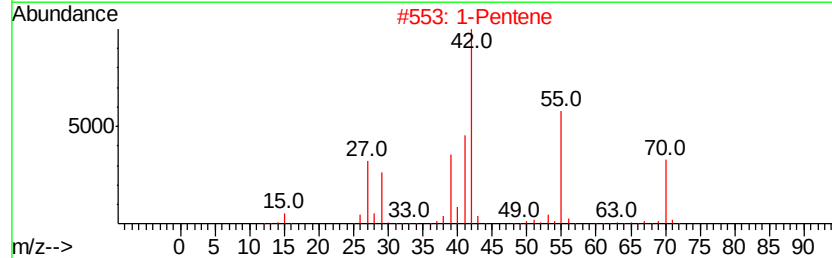
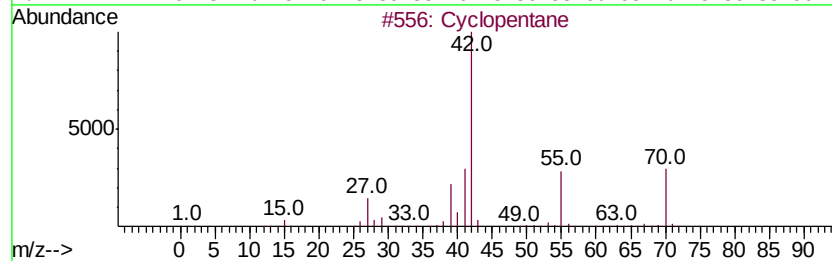
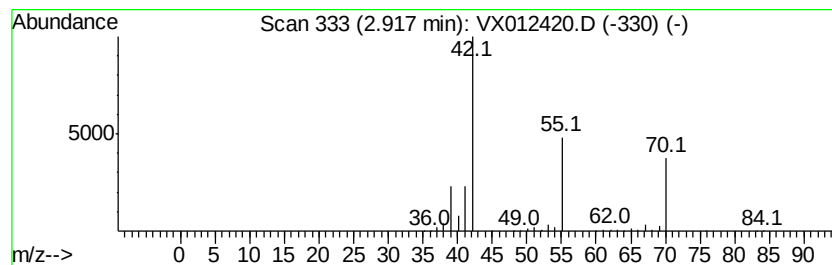
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	194.28 ug/l	1893660	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	86
2		1-Pentene	70	C5H10	000109-67-1	74
3		n-Propyl chloride	78	C3H7Cl	000540-54-5	25
4		Cyclobutanone	70	C4H6O	001191-95-3	9
5		Cyclopropane, ethyl-	70	C5H10	001191-96-4	9



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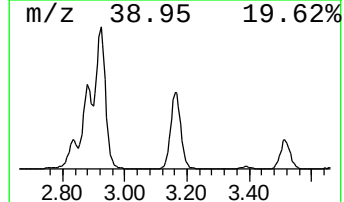
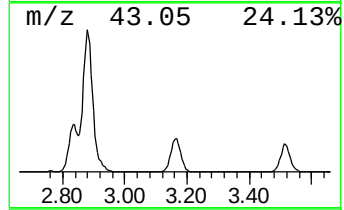
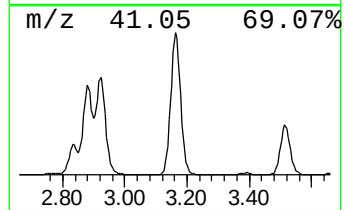
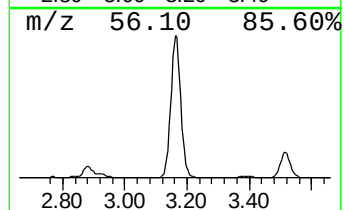
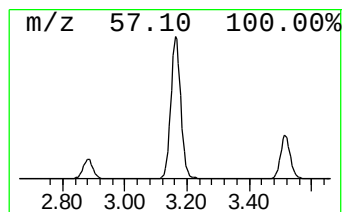
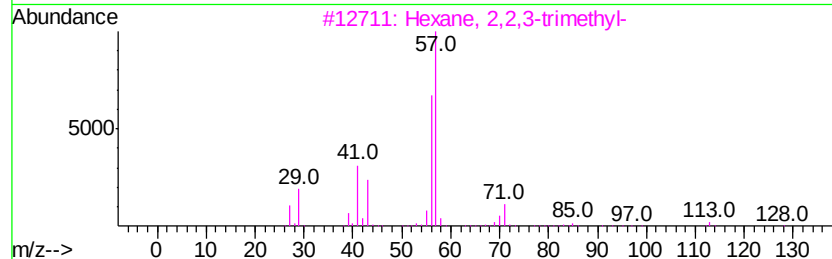
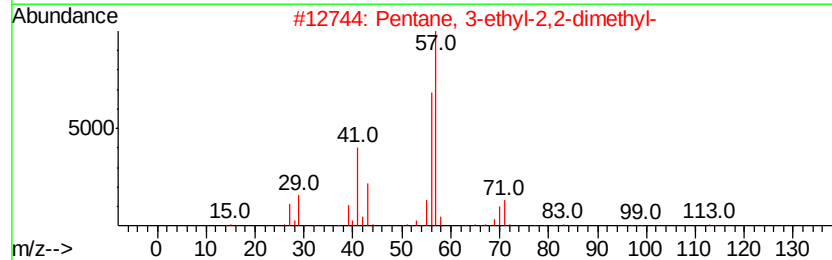
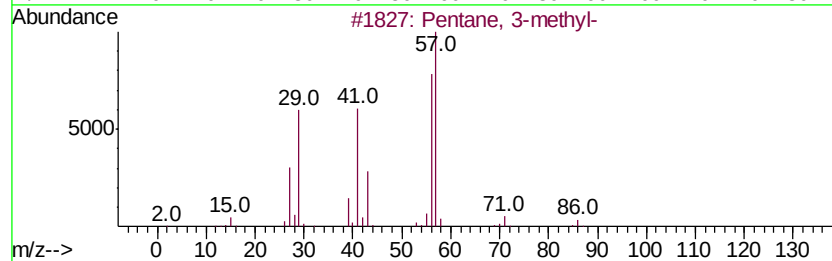
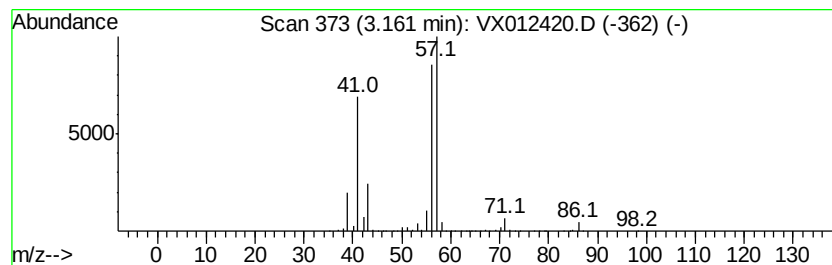
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	204.59 ug/l	1994150	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
3		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	43
4		n-Hexane	86	C6H14	000110-54-3	40
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0280

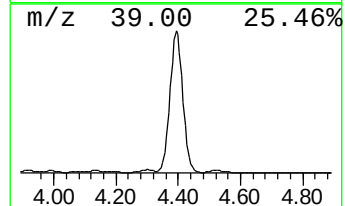
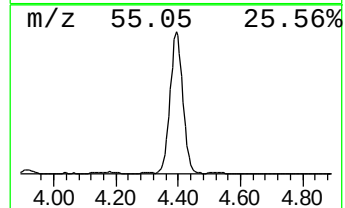
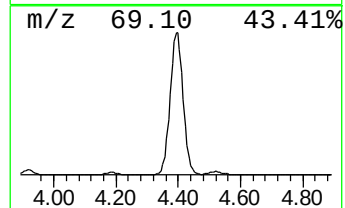
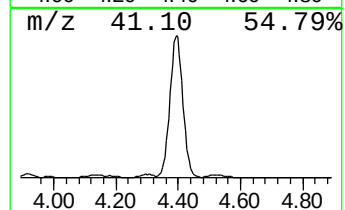
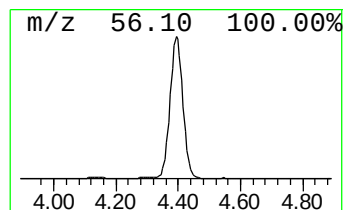
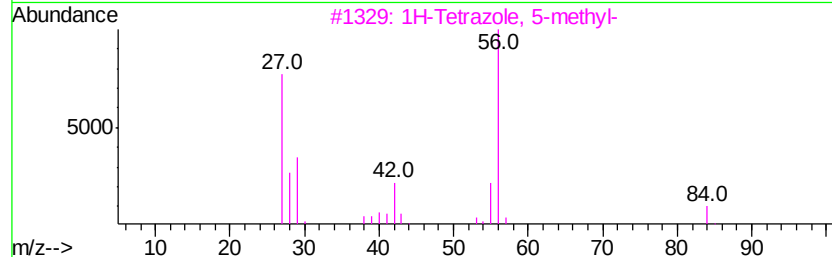
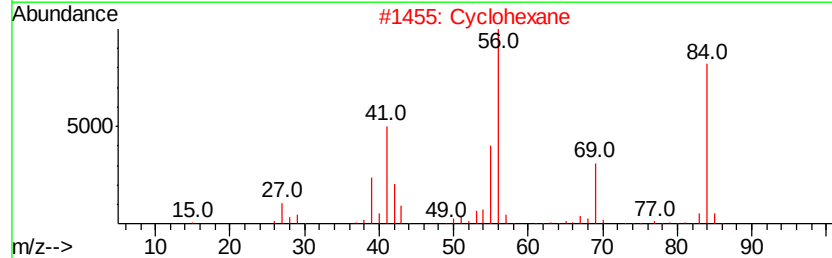
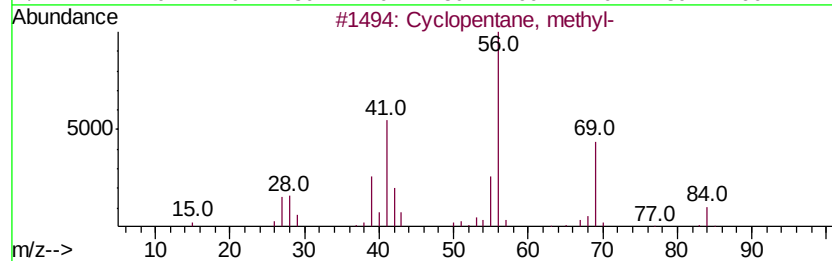
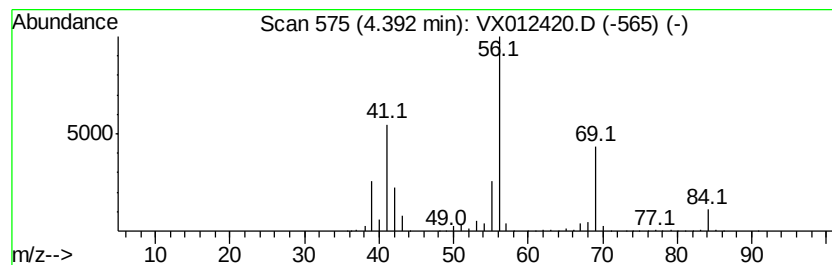
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	423.14 ug/l	4124250	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclohexane	84	C6H12	000110-82-7	86
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		Cyclobutane	56	C4H8	000287-23-0	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0280

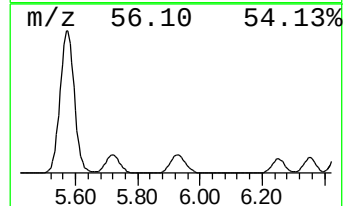
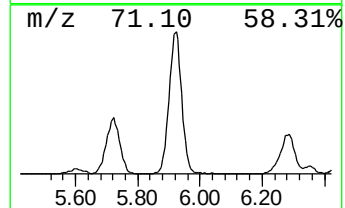
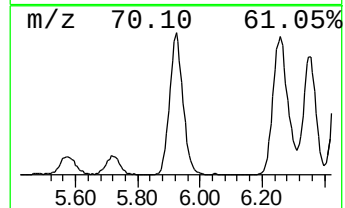
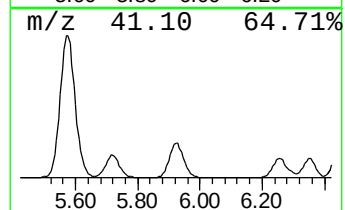
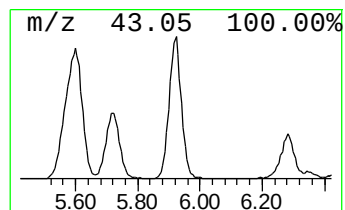
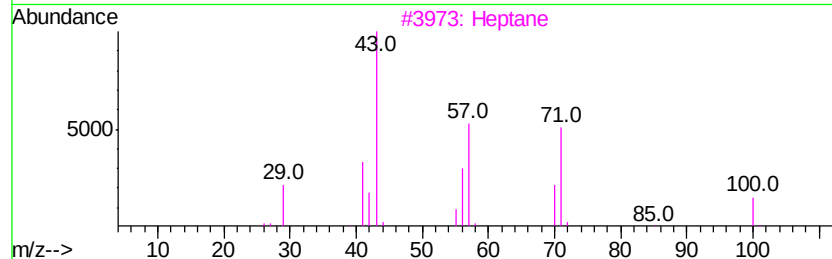
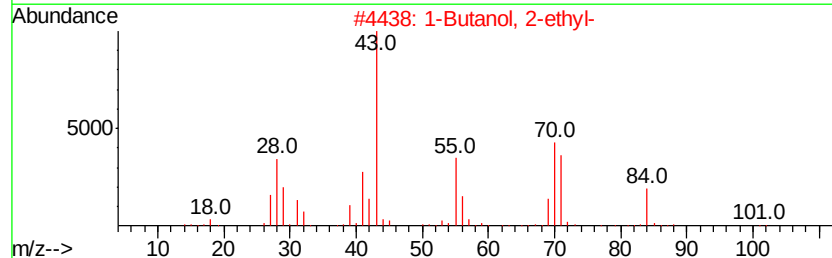
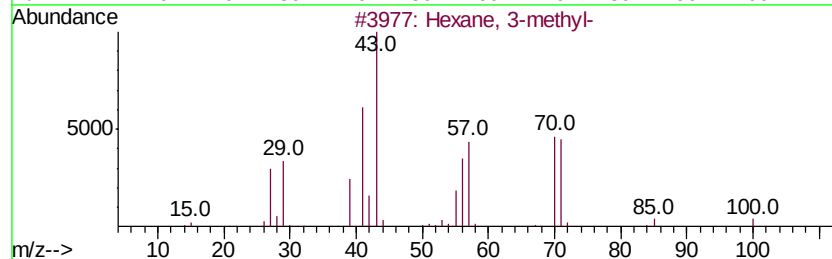
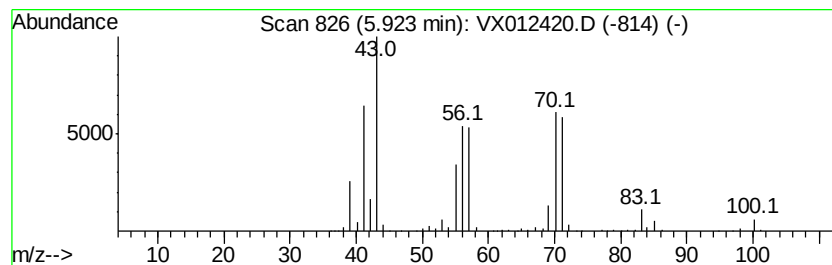
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	106.69 ug/l	1039880	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	93
2		1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53
3		Heptane	100	C7H16	000142-82-5	52
4		Heptane, 4-methyl-	114	C8H18	000589-53-7	50
5		3,3-Dimethyl-2-pentanol	116	C7H16O	019781-24-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
S13-0280

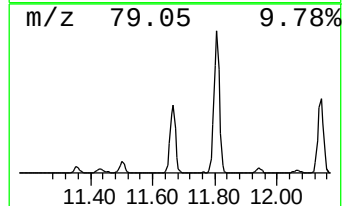
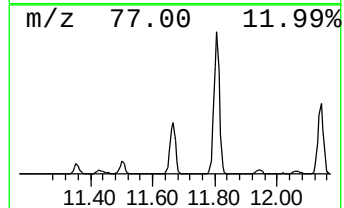
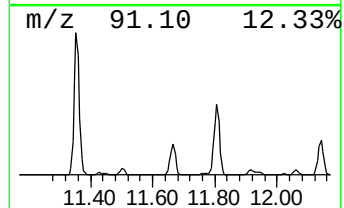
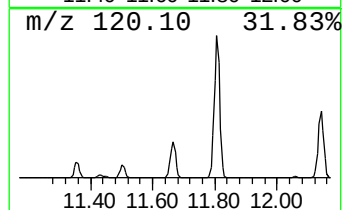
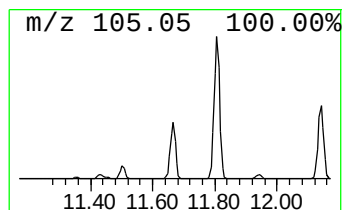
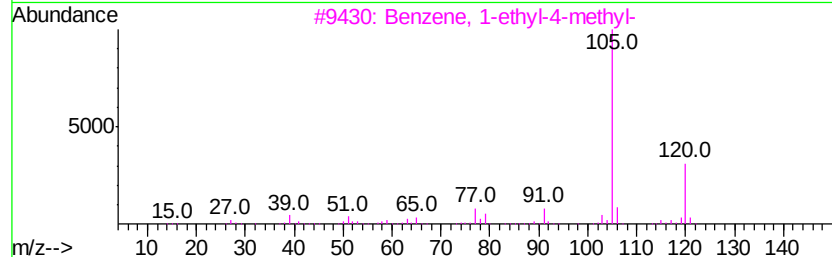
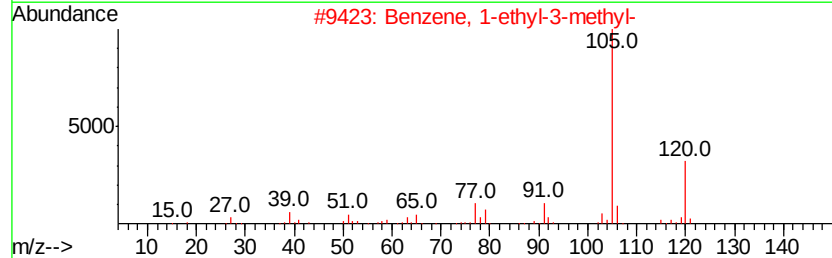
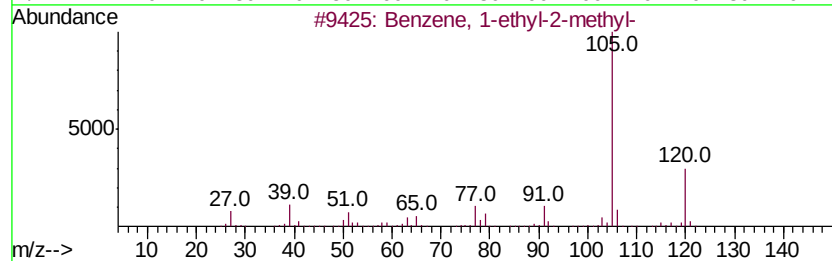
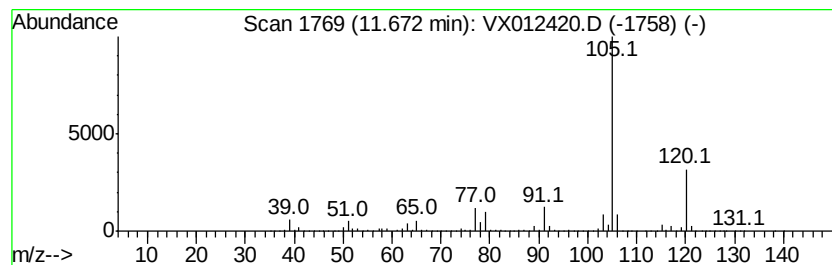
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	235.99 ug/l	4930930	1,4-Dichlorobenzene-d4	12.07

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-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0280

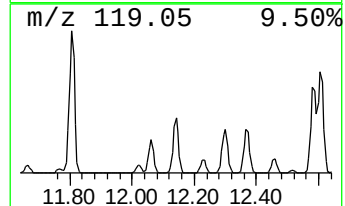
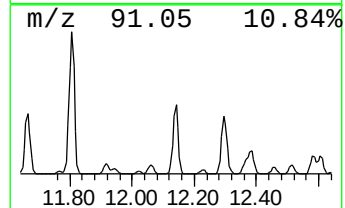
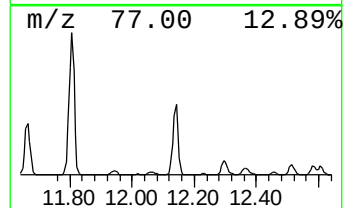
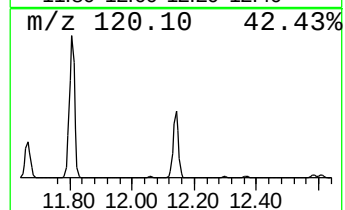
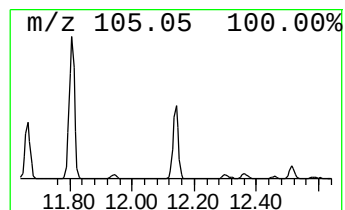
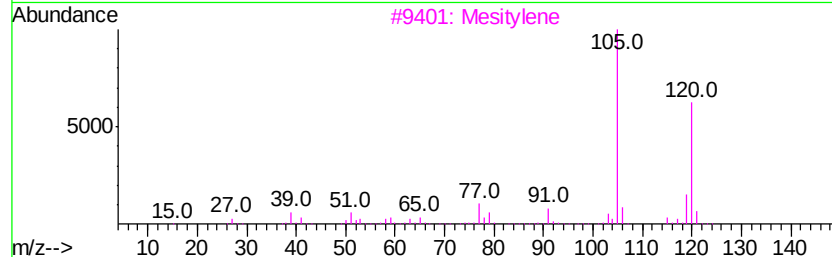
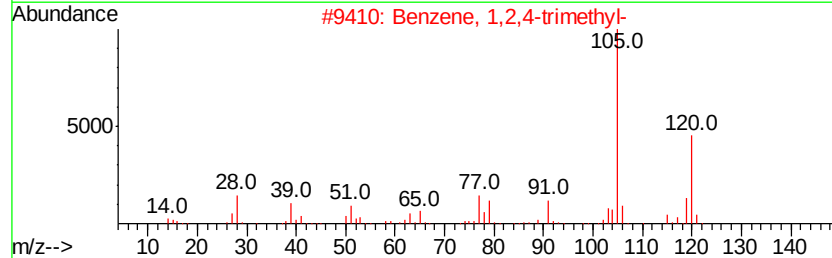
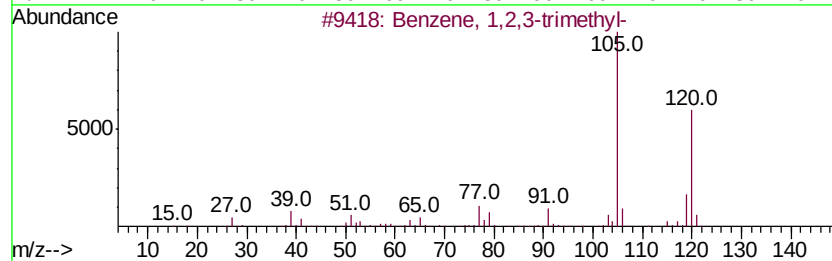
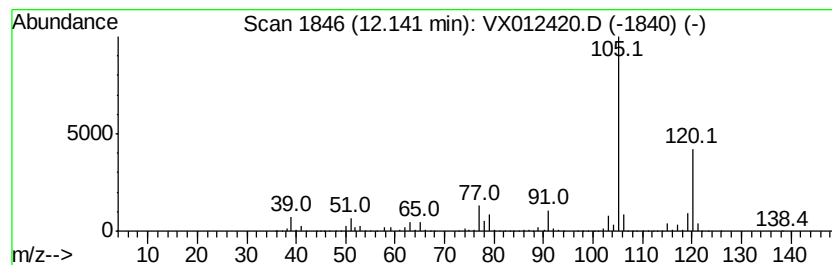
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	320.75 ug/l	6701850	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		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

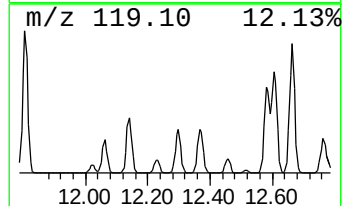
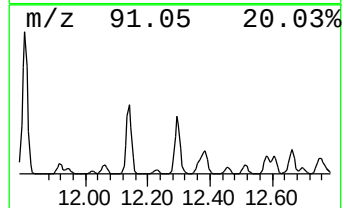
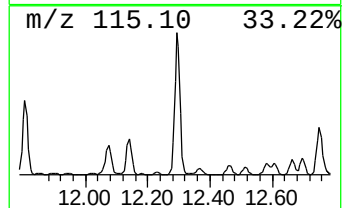
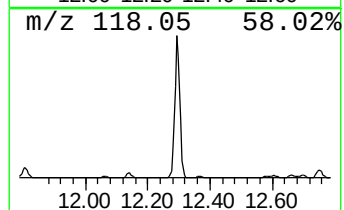
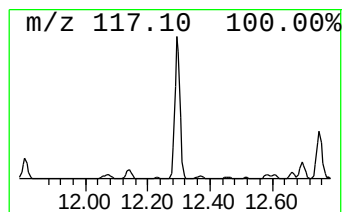
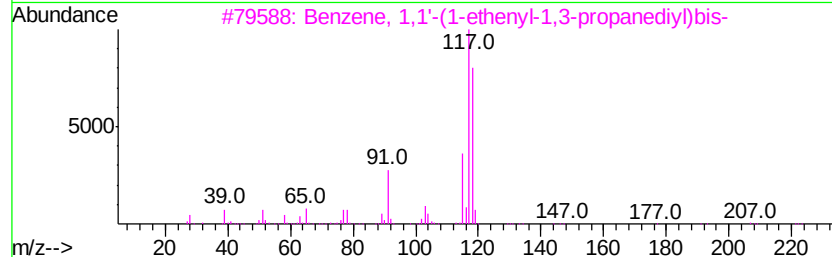
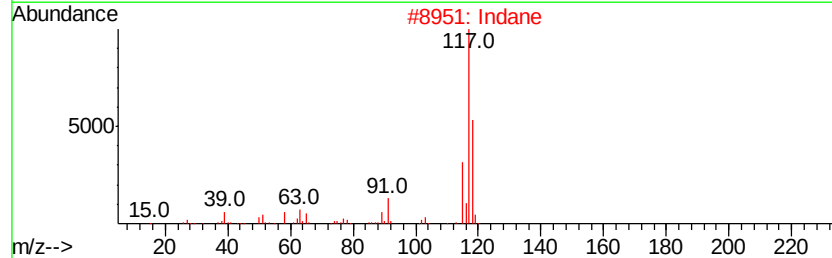
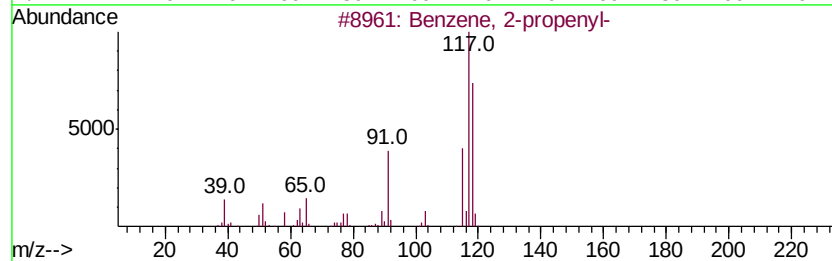
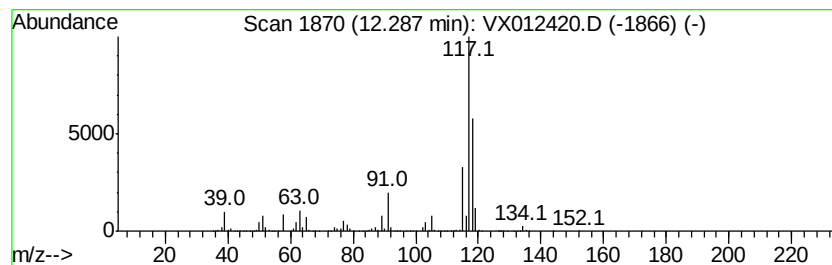
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ClientSampleId :
S13-0280

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-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	192.34 ug/l	4018840	1,4-Dichlorobenzene-d4	12.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-propenyl-	118 C9H10	000300-57-2 76
2		Indane	118 C9H10	000496-11-7 76
3		Benzene, 1,1'-(1-ethenyl-1,3-pro...	222 C17H18	061141-97-7 64
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
5		(Z)-1-Phenylpropene	118 C9H10	000766-90-5 55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0280

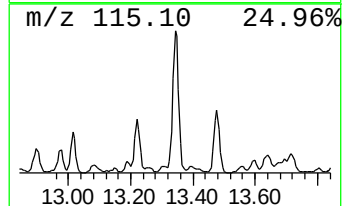
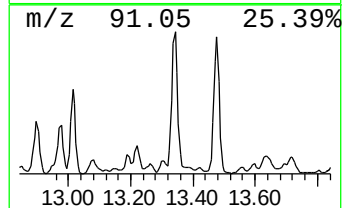
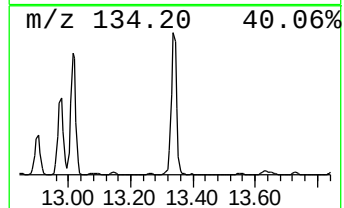
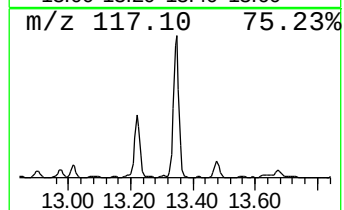
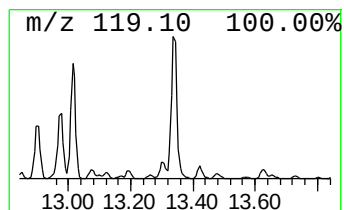
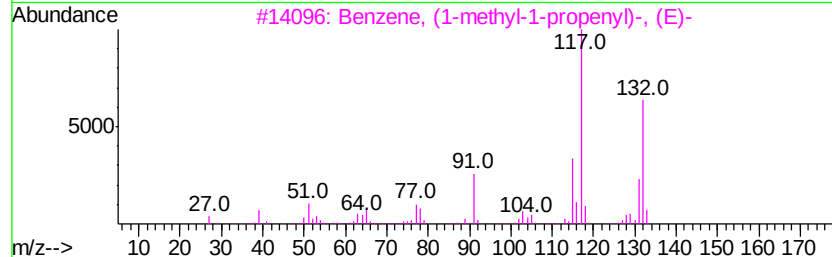
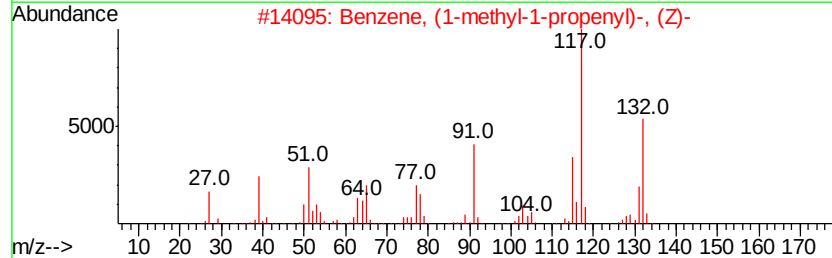
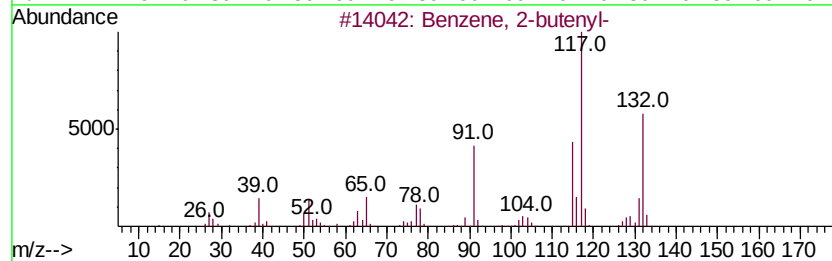
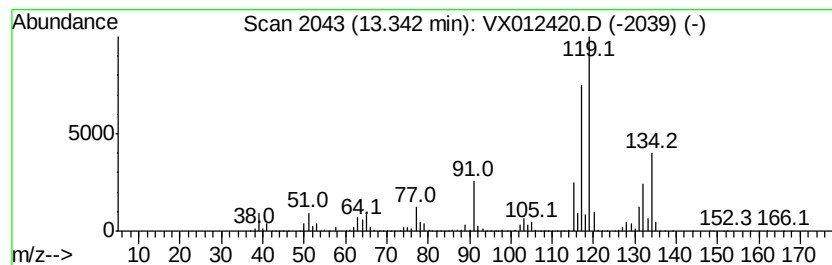
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-butenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	140.94 ug/l	2944880	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
5		o-Cymene	134	C10H14	000527-84-4	55



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091619\
Data File : VX012420.D
Acq On : 16 Sep 2019 21:02
Operator : JC/SP
Sample : K4830-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0280

Quant Method : Z:\VOASRV\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
Butane, 2-methyl-	1.78	111.8	ug/l	1090170	1	5.65	487344	50.0
Pentane, 2-methyl-	2.88	168.0	ug/l	1637290	1	5.65	487344	50.0
Cyclopentane	2.92	194.3	ug/l	1893660	1	5.65	487344	50.0
Pentane, 3-methyl-	3.16	204.6	ug/l	1994150	1	5.65	487344	50.0
Cyclopentane, met...	4.39	423.1	ug/l	4124250	1	5.65	487344	50.0
Hexane, 3-methyl-	5.92	106.7	ug/l	1039880	1	5.65	487344	50.0
Benzene, 1-ethyl-...	11.67	236.0	ug/l	4930930	4	12.07	1044720	50.0
Benzene, 1,2,3-tr...	12.14	320.8	ug/l	6701850	4	12.07	1044720	50.0
Benzene, 2-propenyl-	12.29	192.3	ug/l	4018840	4	12.07	1044720	50.0
Benzene, 2-butenyl-	13.34	140.9	ug/l	2944880	4	12.07	1044720	50.0