

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100318\
 Data File : VX004953.D
 Acq On : 03 Oct 2018 07:59
 Operator : JC/MD
 Sample : J5084-01 10X
 Misc : 5.73g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-8(11-11.5)

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\82X092618W.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	1.404	37	41	47	rBV	337931	322859	4.20%	0.291%
2	1.763	95	100	109	rBV	2302104	3342871	43.44%	3.016%
3	1.983	131	136	151	rVB	1215752	1815986	23.60%	1.639%
4	2.105	151	156	164	rVV	184738	259935	3.38%	0.235%
5	2.257	176	181	193	rVB	219065	344958	4.48%	0.311%
6	2.385	193	202	217	rBV	105299	213163	2.77%	0.192%
7	2.836	269	276	278	rVV	489069	890543	11.57%	0.804%
8	2.879	278	283	295	rVB	1498768	3216416	41.80%	2.902%
9	3.159	320	329	347	rVB	955420	2119768	27.55%	1.913%
10	3.385	355	366	378	rBV	118570	285463	3.71%	0.258%
11	3.513	378	387	401	rBV	699850	1577417	20.50%	1.423%
12	3.659	402	411	417	rBV3	89096	227840	2.96%	0.206%
13	3.739	417	424	429	rVV	122151	279880	3.64%	0.253%
14	3.806	429	435	444	rVB	183602	429541	5.58%	0.388%
15	3.915	444	453	459	rBV3	98794	249830	3.25%	0.225%
16	4.184	490	497	506	rVB3	140117	362305	4.71%	0.327%
17	4.299	506	516	523	rVV2	346125	1051848	13.67%	0.949%
18	4.391	523	531	544	rVV2	547543	1644333	21.37%	1.484%
19	5.299	672	680	697	rVB2	135648	437697	5.69%	0.395%
20	5.494	697	712	719	rBV3	188692	621393	8.08%	0.561%
21	5.598	719	729	736	rBV5	540341	1655667	21.52%	1.494%
22	5.720	742	749	770	rVB3	752356	2488012	32.33%	2.245%
23	5.921	770	782	797	rBV2	661414	1914129	24.88%	1.727%
24	6.067	797	806	815	rBV2	182296	444338	5.77%	0.401%
25	6.256	824	837	841	rBV2	204446	616045	8.01%	0.556%
26	6.348	842	852	862	rVV2	1619199	4610553	59.92%	4.160%
27	6.451	862	869	880	rVB2	191743	555957	7.23%	0.502%
28	6.695	899	909	918	rBV2	398752	999246	12.99%	0.902%
29	6.860	928	936	942	rBV2	353725	901567	11.72%	0.813%
30	6.933	943	948	956	rVB3	212554	602856	7.83%	0.544%
31	7.256	991	1001	1008	rBV3	118916	283181	3.68%	0.256%
32	7.457	1022	1034	1045	rBV4	515885	1386150	18.01%	1.251%
33	7.573	1045	1053	1058	rBV	242586	514028	6.68%	0.464%
34	7.634	1058	1063	1068	rBV	268107	495607	6.44%	0.447%

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35	7.732	1074	1079	1089	rVB	154381	309527	4.02%	0.279%
36	7.848	1090	1098	1105	rVB3	122198	254796	3.31%	0.230%
37	8.055	1122	1132	1144	rVB	776432	1646963	21.40%	1.486%
38	8.177	1144	1152	1161	rBV2	769789	1831752	23.80%	1.653%
39	8.256	1161	1165	1174	rVB2	311423	580151	7.54%	0.523%
40	8.341	1174	1179	1182	rBV	407745	662998	8.62%	0.598%
41	8.378	1182	1185	1192	rVB	233366	382407	4.97%	0.345%
42	8.500	1200	1205	1214	rBV	444832	852925	11.08%	0.770%
43	8.573	1214	1217	1227	rVB4	136405	279053	3.63%	0.252%
44	8.677	1227	1234	1237	rBV2	446725	920370	11.96%	0.830%
45	8.713	1237	1240	1247	rVB	837430	1316155	17.10%	1.188%
46	8.786	1247	1252	1257	rBV	430561	685796	8.91%	0.619%
47	8.975	1277	1283	1290	rBV2	422566	801233	10.41%	0.723%
48	9.042	1290	1294	1304	rVV4	127486	295604	3.84%	0.267%
49	9.164	1304	1314	1320	rVV3	126548	278971	3.63%	0.252%
50	9.561	1374	1379	1385	rBV3	187270	411420	5.35%	0.371%
51	9.963	1441	1445	1450	rVB	455073	635347	8.26%	0.573%
52	10.067	1457	1462	1466	rBV	387002	514495	6.69%	0.464%
53	10.115	1466	1470	1474	rVV	733681	965015	12.54%	0.871%
54	10.256	1487	1493	1499	rBV	2133319	2970033	38.60%	2.680%
55	10.359	1504	1510	1516	rVV	5494476	7694851	100.00%	6.943%
56	10.414	1516	1519	1531	rVB	586079	962137	12.50%	0.868%
57	10.640	1551	1556	1560	rBV	169148	238379	3.10%	0.215%
58	10.701	1560	1566	1575	rVV	1592855	2215185	28.79%	1.999%
59	10.835	1580	1588	1592	rBV	195638	315236	4.10%	0.284%
60	10.914	1592	1601	1605	rBV3	130908	235201	3.06%	0.212%
61	11.018	1612	1618	1625	rBV	330852	493635	6.42%	0.445%
62	11.140	1632	1638	1641	rBV	972009	1389072	18.05%	1.253%
63	11.243	1649	1655	1659	rBV	252386	328886	4.27%	0.297%
64	11.359	1663	1674	1681	rBV	1235730	1642592	21.35%	1.482%
65	11.438	1681	1687	1689	rBV	2656254	4049660	52.63%	3.654%
66	11.505	1694	1698	1710	rVB2	1693878	2851569	37.06%	2.573%
67	11.670	1720	1725	1732	rBV	1508469	1955482	25.41%	1.764%
68	11.810	1743	1748	1755	rVB	6369628	7627237	99.12%	6.882%
69	11.944	1768	1770	1774	rVB	174701	198263	2.58%	0.179%
70	12.024	1780	1783	1787	rVB	229843	262186	3.41%	0.237%
71	12.079	1787	1792	1797	rVB	1017905	1384238	17.99%	1.249%

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Stop Thrs : 0 Peak Location: TOP

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72	12.146	1797	1803	1807	rBV	1461395	1894178	24.62%	1.709%
73	12.304	1823	1829	1831	rBV	1470978	2144623	27.87%	1.935%
74	12.371	1836	1840	1847	rVB3	1635901	2561449	33.29%	2.311%
75	12.475	1847	1857	1861	rBV3	386522	620800	8.07%	0.560%
76	12.524	1861	1865	1871	rVB2	630927	985571	12.81%	0.889%
77	12.585	1871	1875	1877	rBV	737896	923033	12.00%	0.833%
78	12.609	1877	1879	1884	rVV	723829	866144	11.26%	0.782%
79	12.670	1884	1889	1892	rVV	1386423	1627403	21.15%	1.468%
80	12.700	1892	1894	1899	rVV	147046	195197	2.54%	0.176%
81	12.755	1899	1903	1914	rVB2	404896	912116	11.85%	0.823%
82	12.908	1924	1928	1931	rVB	255925	316005	4.11%	0.285%
83	12.981	1931	1940	1943	rBV	629341	928295	12.06%	0.838%
84	13.017	1943	1946	1952	rVB	914463	1120499	14.56%	1.011%
85	13.078	1952	1956	1958	rBV	188409	260964	3.39%	0.235%
86	13.103	1958	1960	1962	rVV2	237128	263938	3.43%	0.238%
87	13.225	1972	1980	1984	rVB3	588232	918408	11.94%	0.829%
88	13.316	1990	1995	1997	rBV4	239196	375782	4.88%	0.339%
89	13.347	1997	2000	2006	rVB2	922381	1287613	16.73%	1.162%
90	13.426	2011	2013	2018	rVB	199623	221238	2.88%	0.200%
91	13.481	2018	2022	2031	rVB2	204366	302888	3.94%	0.273%
92	13.603	2038	2042	2044	rBV	163587	192714	2.50%	0.174%
93	13.639	2044	2048	2054	rVB2	256495	472150	6.14%	0.426%
94	13.725	2059	2062	2067	rVB3	174942	288026	3.74%	0.260%
95	13.834	2075	2080	2086	rVB	892159	1099659	14.29%	0.992%
96	13.987	2098	2105	2109	rBV4	129912	244407	3.18%	0.221%
97	14.133	2125	2129	2133	rBV	190649	230109	2.99%	0.208%
98	14.273	2148	2152	2158	rBV	177000	279303	3.63%	0.252%
99	14.694	2213	2221	2227	rBV	873120	1146863	14.90%	1.035%
100	14.840	2238	2245	2254	rVB	424471	549750	7.14%	0.496%

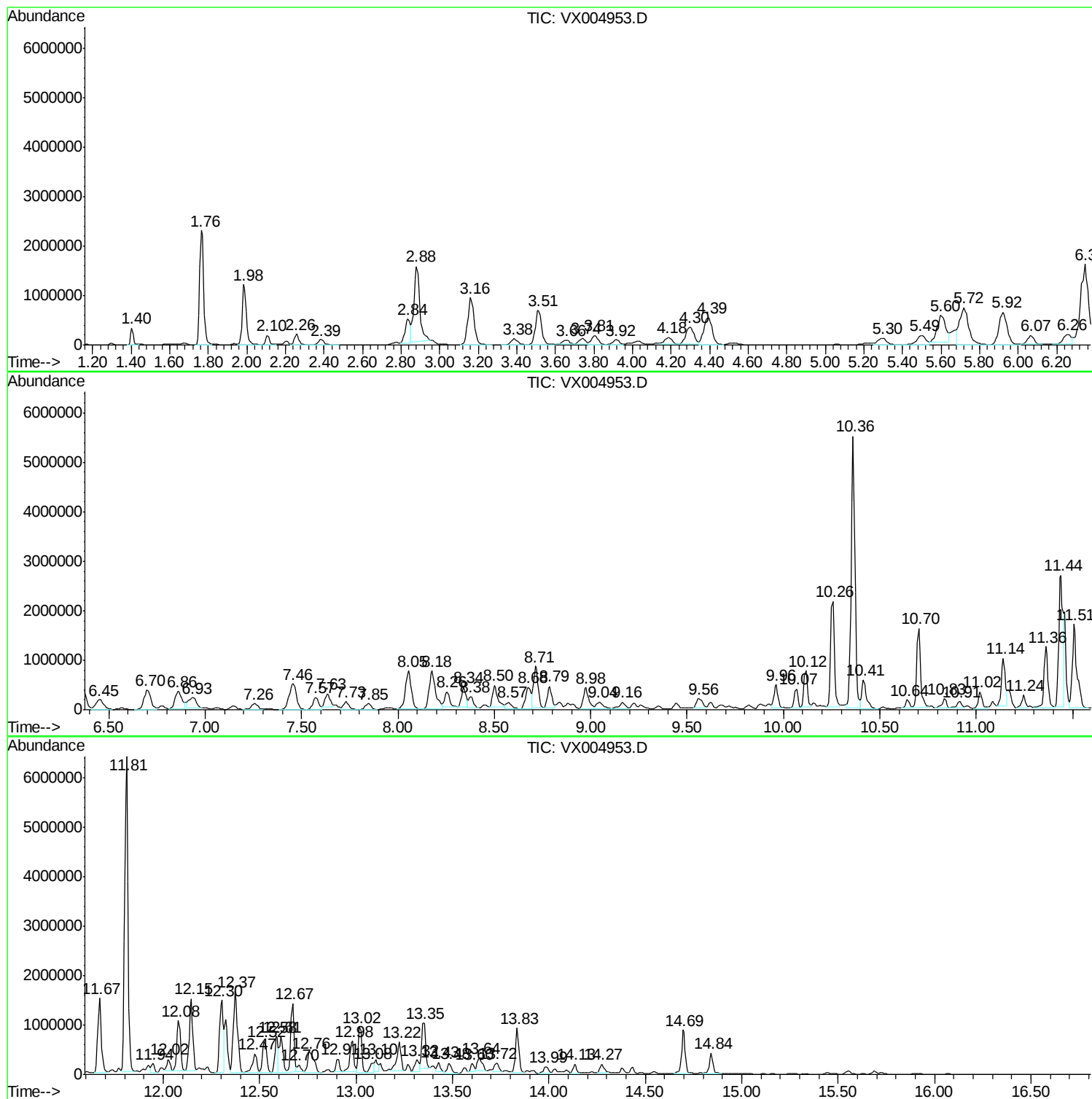
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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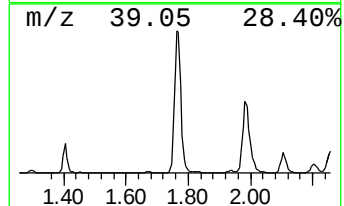
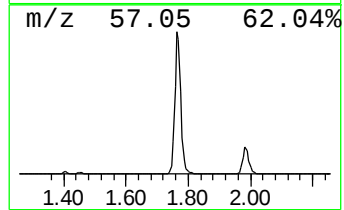
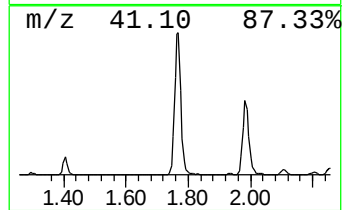
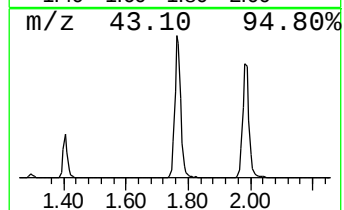
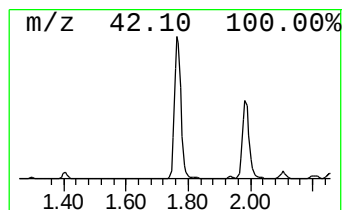
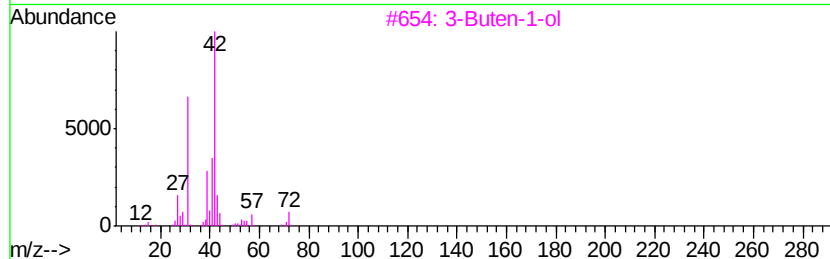
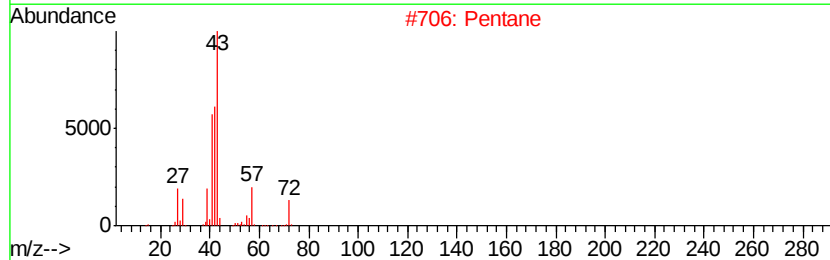
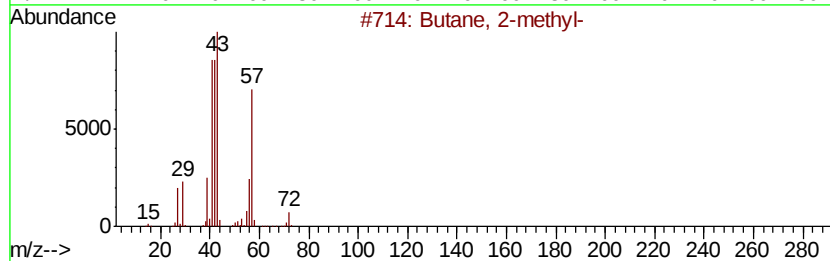
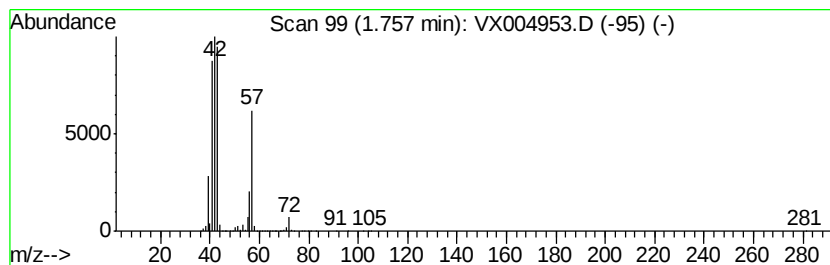
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	100.95 ug/l	3342870	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2		Pentane	72	C5H12	000109-66-0	50
3		3-Buten-1-ol	72	C4H8O	000627-27-0	12
4		1-Butene	56	C4H8	000106-98-9	9
5		3-Buten-2-ol	72	C4H8O	000598-32-3	9



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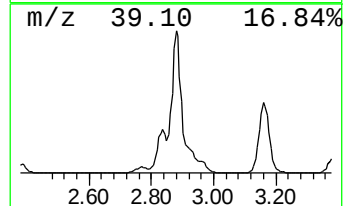
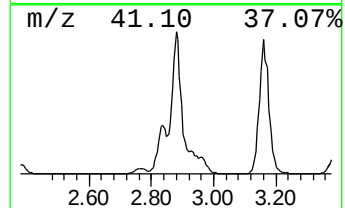
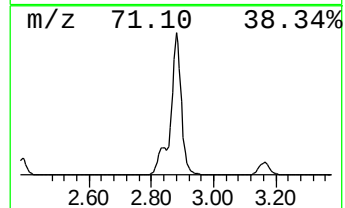
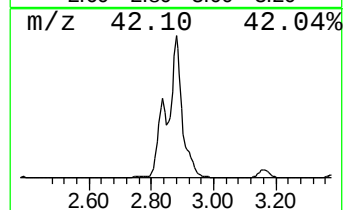
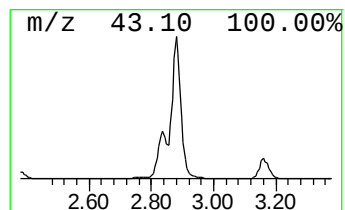
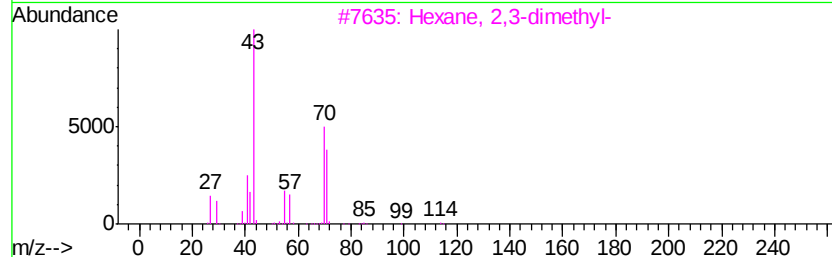
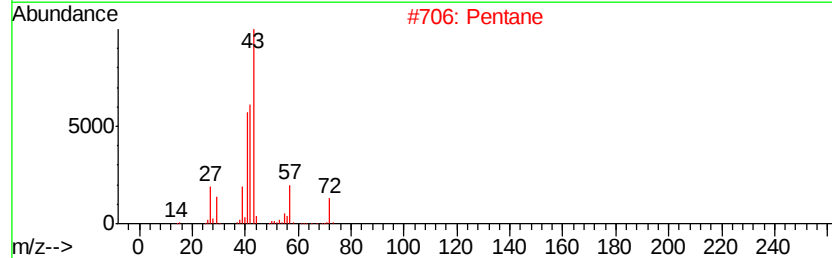
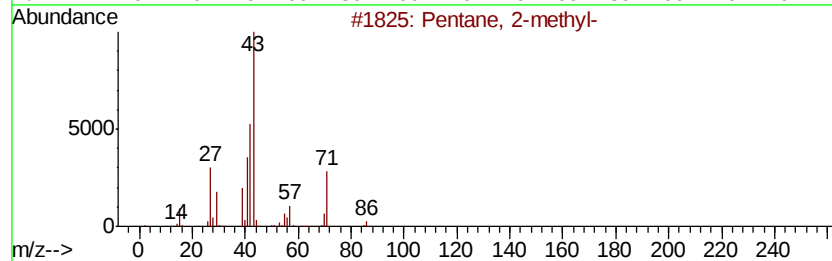
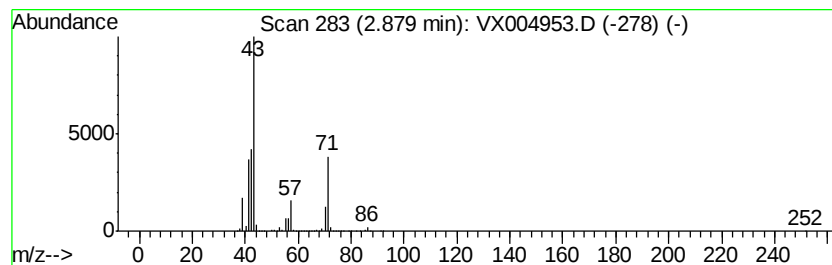
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	97.13 ug/l	3216420	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2		Pentane	72	C5H12	000109-66-0	38
3		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
4		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38
5		Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	33



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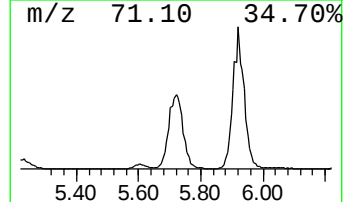
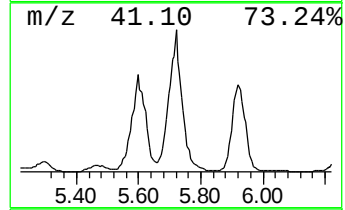
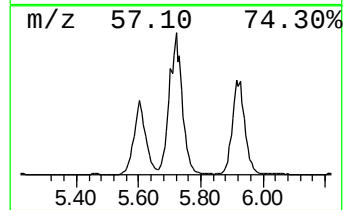
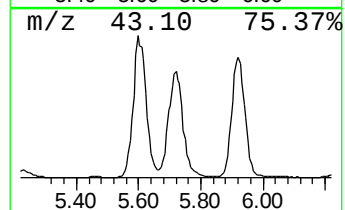
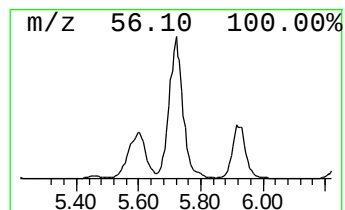
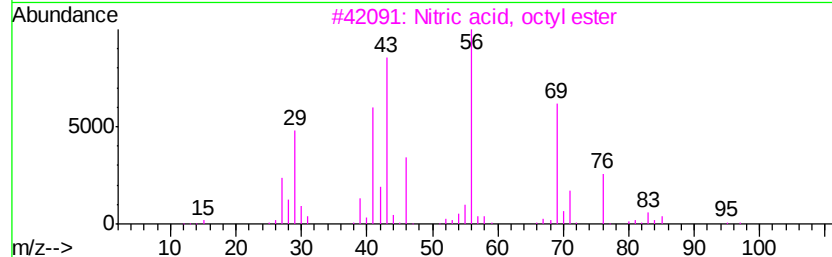
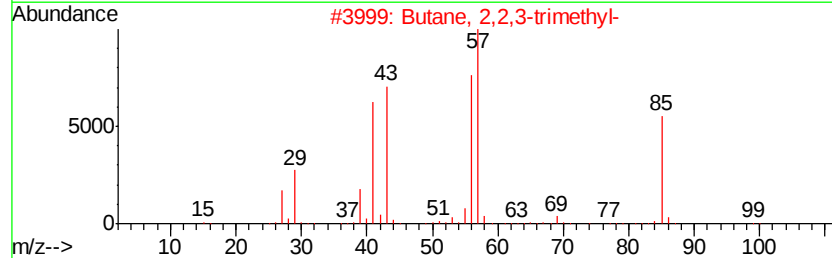
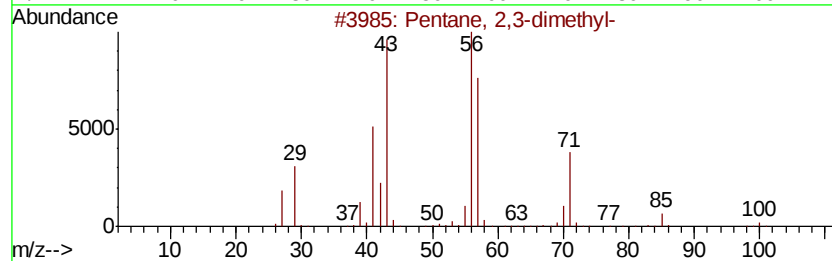
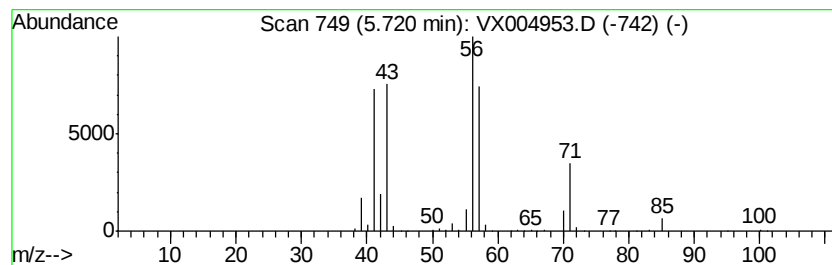
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.72	75.14 ug/l	2488010	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	83
2		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	47
3		Nitric acid, octyl ester	175	C8H17NO3	000629-39-0	45
4		Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	45
5		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73q/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-8(11-11.5)

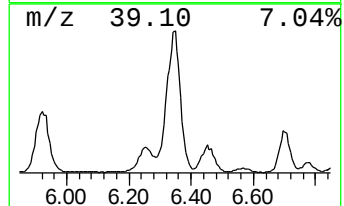
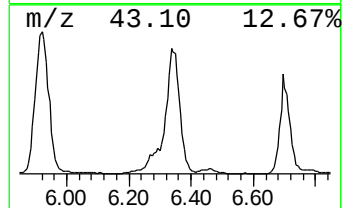
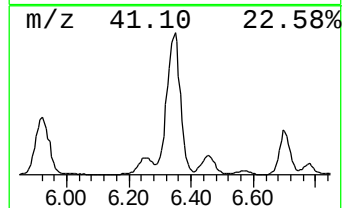
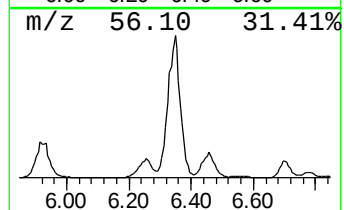
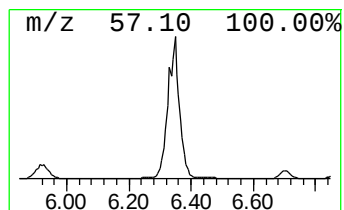
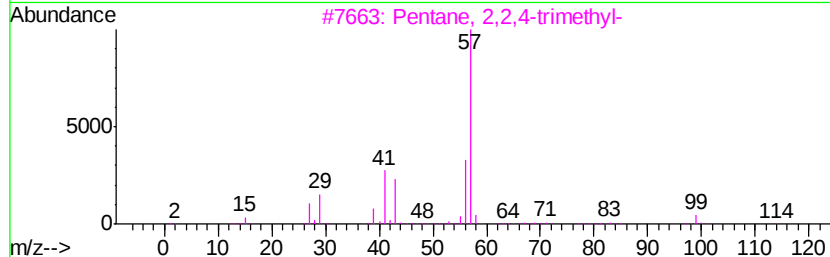
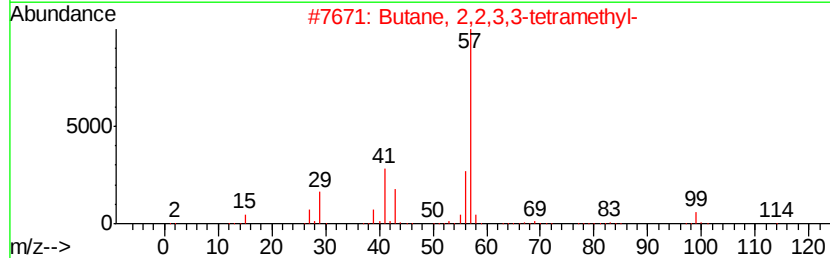
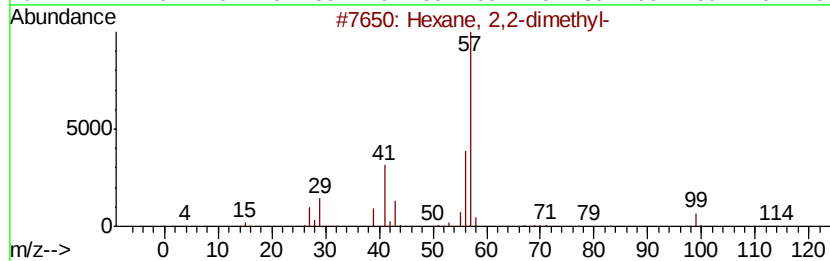
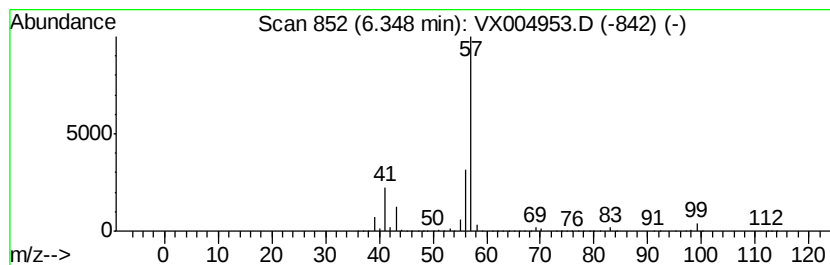
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Quant Title : SW846 8260

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

Peak Number 4 Hexane, 2,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	255.70 ug/l	4610550	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
4		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	56



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73q/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-8(11-11.5)

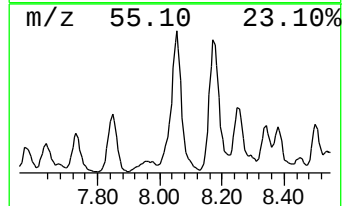
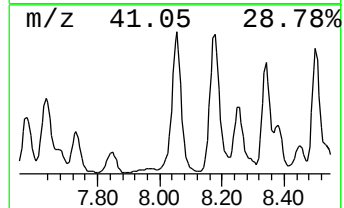
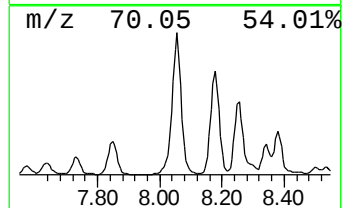
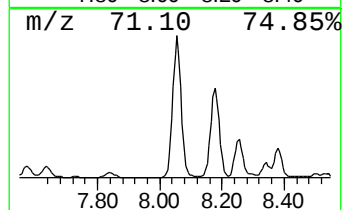
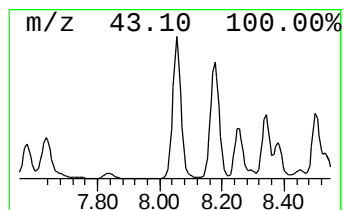
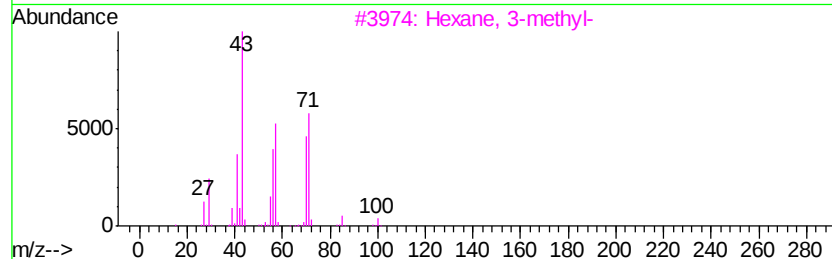
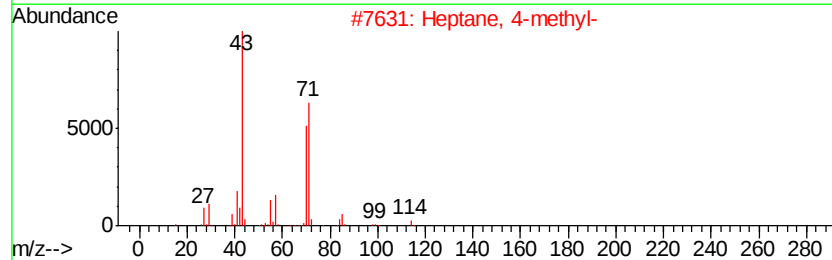
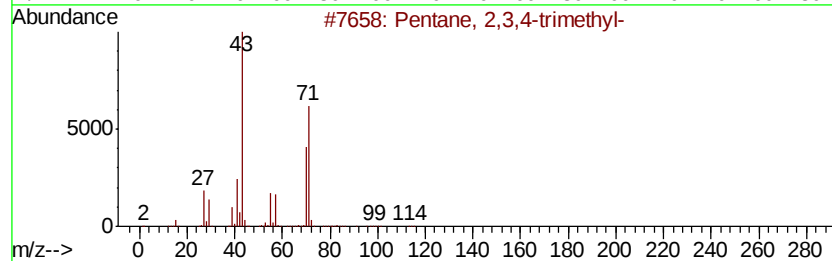
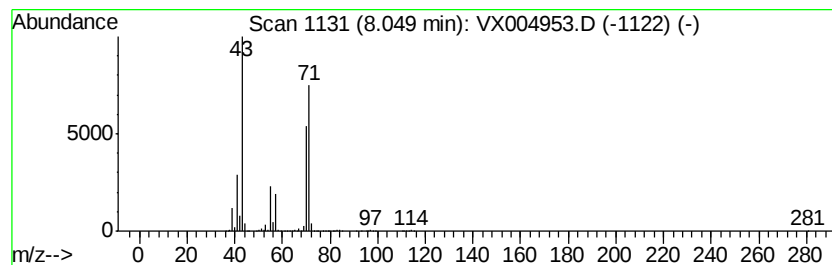
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,4-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	91.34 ug/l	1646960	1,4-Difluorobenzene	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91	
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	83	
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	78	
4	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74	
5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	74	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73q/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-8(11-11.5)

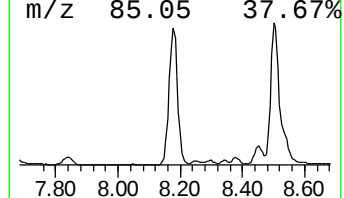
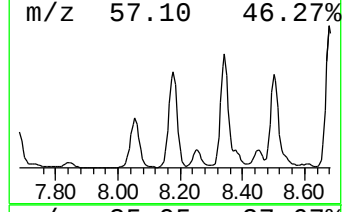
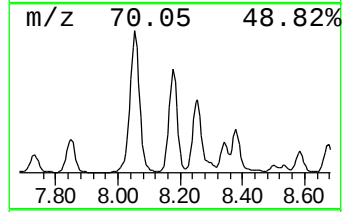
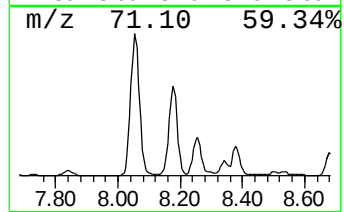
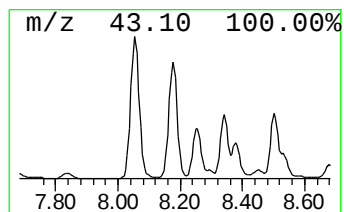
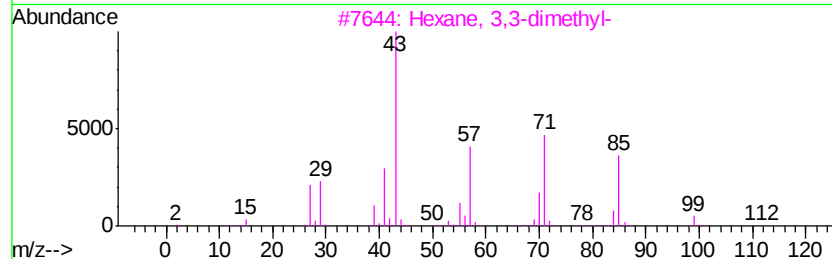
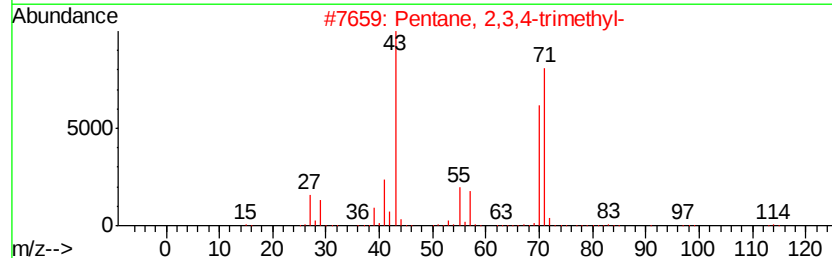
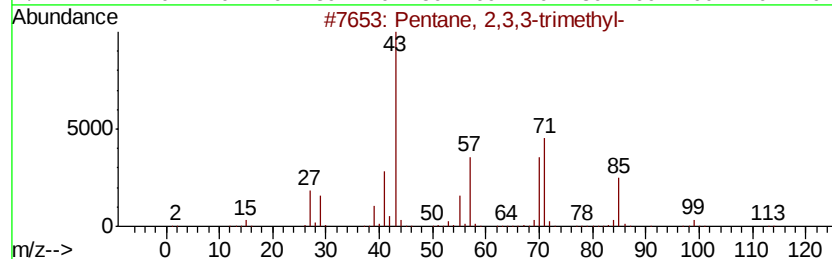
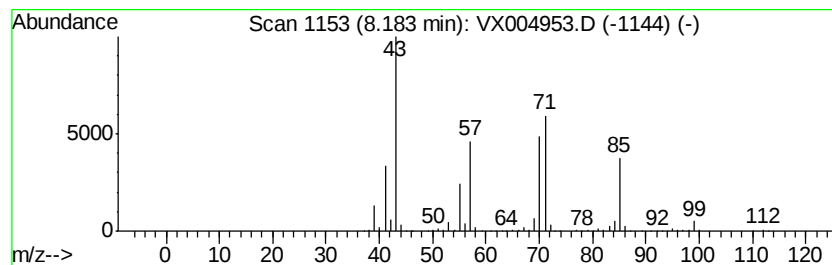
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

Peak Number 6 Pentane, 2,3,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	101.59 ug/l	1831750	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59
5		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73q/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

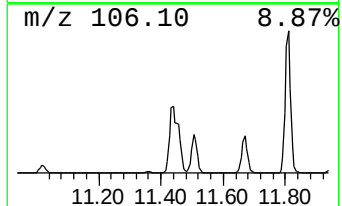
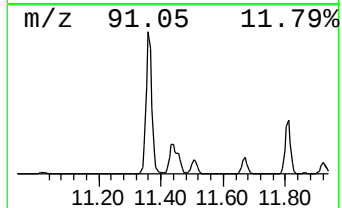
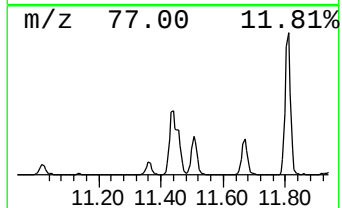
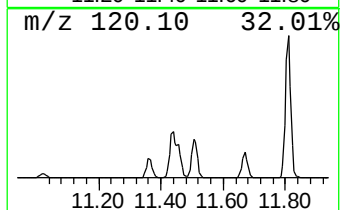
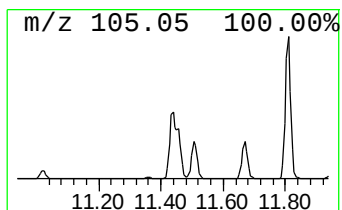
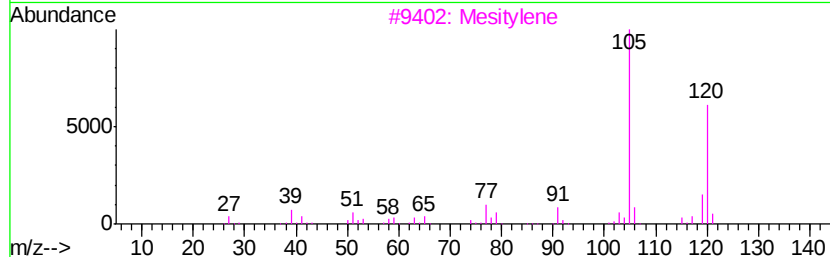
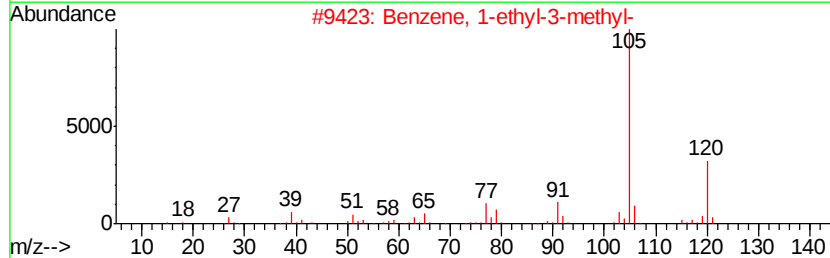
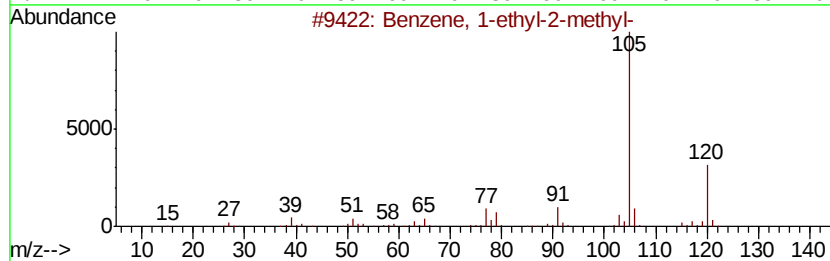
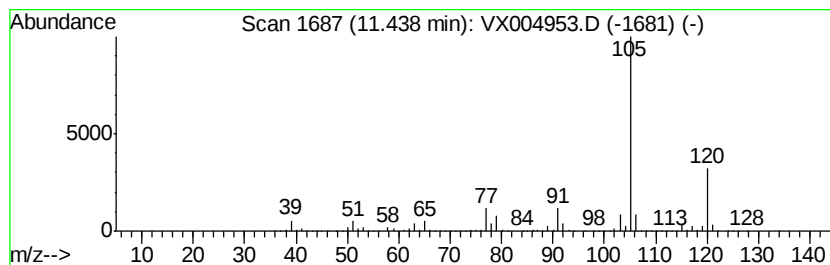
Instrument :
MSVOA_X
ClientSampleId :
B-8(11-11.5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
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 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.44	146.28 ug/l	4049660	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	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73q/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
B-8(11-11.5)

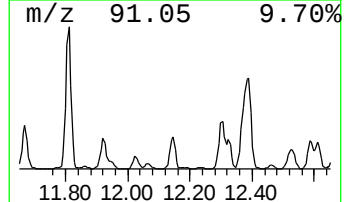
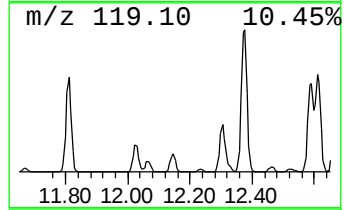
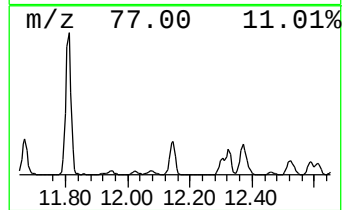
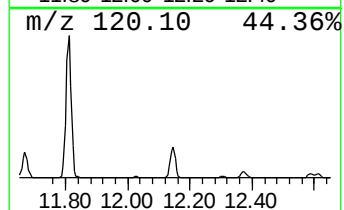
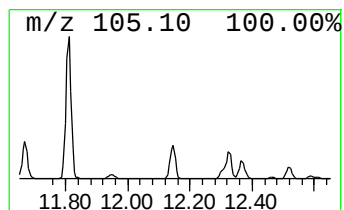
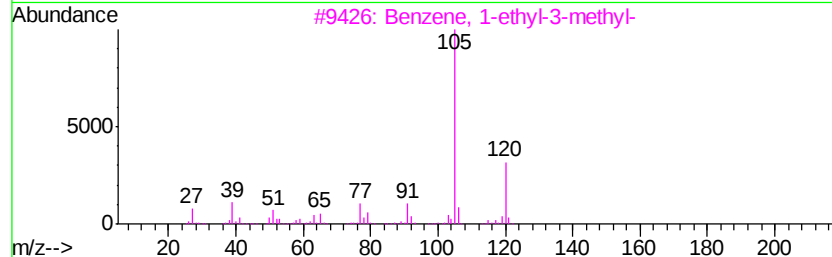
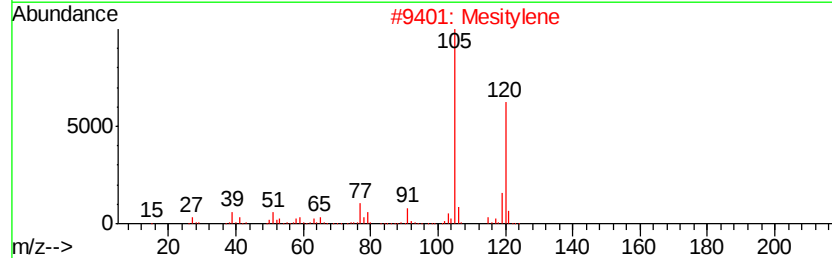
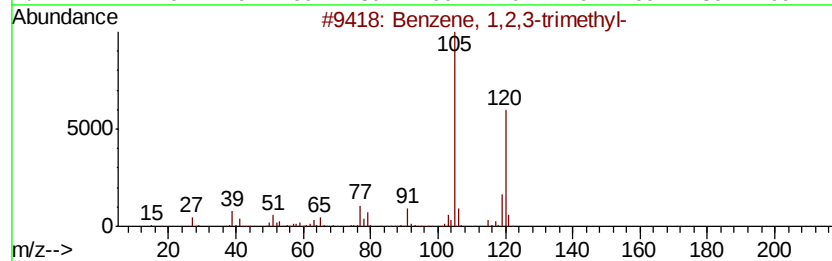
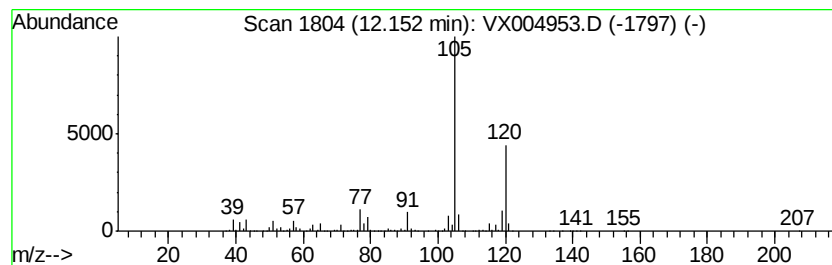
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	68.42 ug/l	1894180	1,4-Dichlorobenzene-d4	12.08

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	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

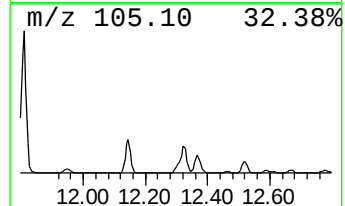
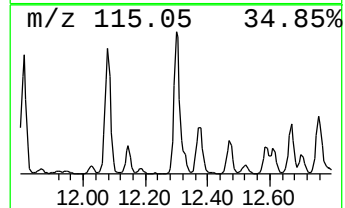
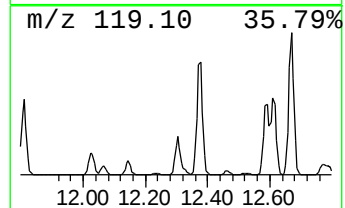
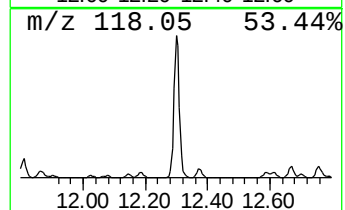
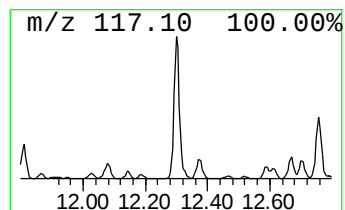
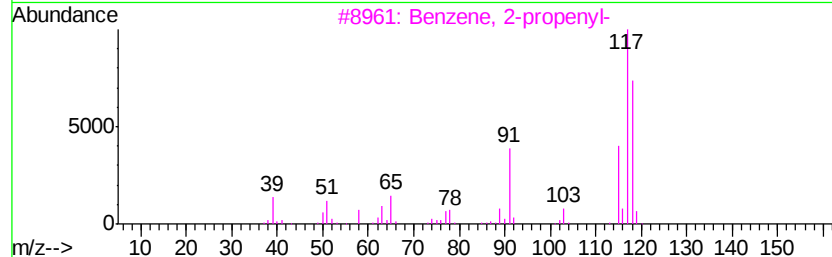
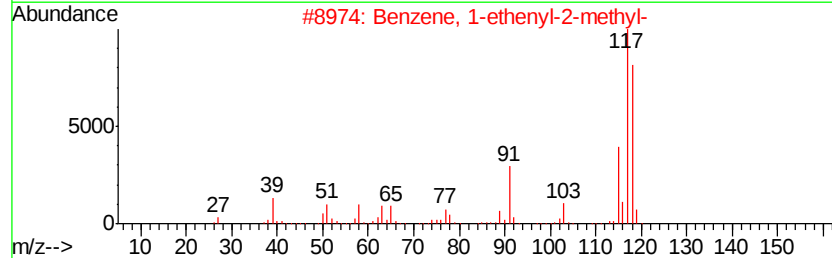
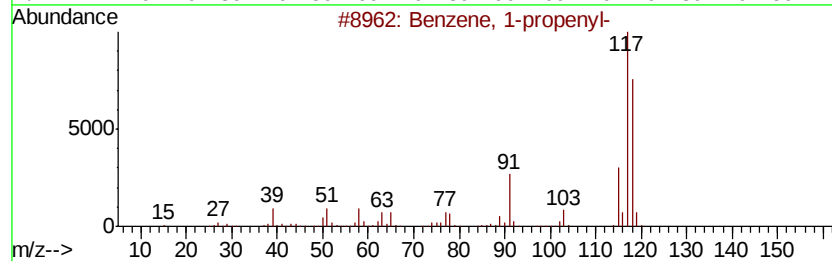
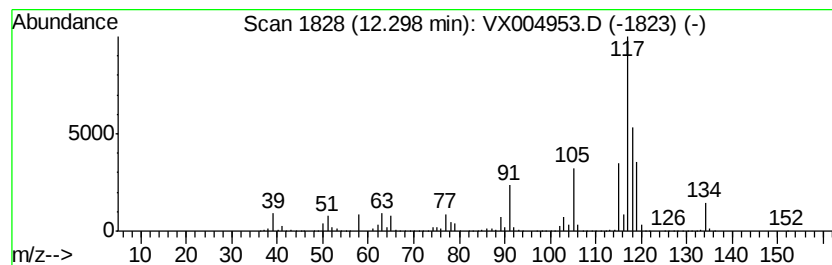
Instrument :
MSVOA_X
ClientSampled :
B-8(11-11.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-propenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.30	77.47 ug/l	2144620	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	55



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100318\
Data File : VX004953.D
Acq On : 03 Oct 2018 07:59
Operator : JC/MD
Sample : J5084-01 10X
Misc : 5.73g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-8(11-11.5)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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-	1.76	101.0	ug/l	3342870	1	5.66	1655670	50.0
Pentane, 2-methyl-	2.88	97.1	ug/l	3216420	1	5.66	1655670	50.0
Pentane, 2,3-dime...	5.72	75.1	ug/l	2488010	1	5.66	1655670	50.0
Hexane, 2,2-dimet...	6.35	255.7	ug/l	4610550	2	6.86	901567	50.0
Pentane, 2,3,4-tr...	8.05	91.3	ug/l	1646960	2	6.86	901567	50.0
Pentane, 2,3,3-tr...	8.18	101.6	ug/l	1831750	2	6.86	901567	50.0
Benzene, 1-ethyl-...	11.44	146.3	ug/l	4049660	4	12.08	1384240	50.0
Benzene, 1,2,3-tr...	12.15	68.4	ug/l	1894180	4	12.08	1384240	50.0
Benzene, 1-propenyl-	12.30	77.5	ug/l	2144620	4	12.08	1384240	50.0