

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082318\
 Data File : VX004227.D
 Acq On : 23 Aug 2018 15:26
 Operator : JC/MD
 Sample : J4579-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0561

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\82X082218W.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	1.075	75	80	94	rBV	1786096	2305543	7.51%	0.914%
2	1.788	191	197	208	rVB	399779	556367	1.81%	0.221%
3	1.995	226	231	244	rVB	720057	1021028	3.33%	0.405%
4	2.392	288	296	302	rBV	274401	540563	1.76%	0.214%
5	2.843	362	370	372	rBV	271723	554943	1.81%	0.220%
6	2.885	372	377	400	rVB	940875	2416391	7.87%	0.958%
7	3.166	413	423	438	rVB	876774	1975631	6.44%	0.783%
8	3.513	471	480	495	rVB	709942	1594098	5.19%	0.632%
9	4.391	614	624	638	rVB	984422	2859327	9.31%	1.134%
10	5.507	790	807	811	rBV	213249	646750	2.11%	0.256%
11	5.574	811	818	828	rVV3	600709	2075036	6.76%	0.823%
12	5.659	828	832	838	rVB	191483	436687	1.42%	0.173%
13	5.922	862	875	889	rBV	438536	1326664	4.32%	0.526%
14	6.068	889	899	913	rVB	232026	590065	1.92%	0.234%
15	6.257	913	930	940	rBV3	122333	481587	1.57%	0.191%
16	6.452	954	962	974	rVB	197644	545451	1.78%	0.216%
17	6.702	992	1003	1012	rBV	323429	781328	2.55%	0.310%
18	6.860	1020	1029	1037	rBV	493496	1153722	3.76%	0.457%
19	7.458	1115	1127	1139	rBV	1924086	4429205	14.43%	1.756%
20	7.732	1165	1172	1182	rVB	264752	508861	1.66%	0.202%
21	8.665	1318	1325	1328	rBV2	414749	719399	2.34%	0.285%
22	8.714	1328	1333	1343	rVB	1265272	2153523	7.01%	0.854%
23	8.970	1370	1375	1381	rBV	220199	332754	1.08%	0.132%
24	9.037	1381	1386	1393	rVB	365482	600238	1.96%	0.238%
25	9.165	1398	1407	1413	rBV	192077	313185	1.02%	0.124%
26	9.622	1478	1482	1487	rVB	510281	701076	2.28%	0.278%
27	9.677	1487	1491	1496	rBV	211284	308137	1.00%	0.122%
28	10.116	1558	1563	1568	rVB	925243	1275568	4.15%	0.506%
29	10.256	1580	1586	1592	rBV	9573347	12421828	40.46%	4.925%
30	10.360	1594	1603	1609	rBV	12871600	17612018	57.37%	6.983%
31	10.701	1653	1659	1669	rVB	11113329	14016729	45.66%	5.557%
32	11.018	1706	1711	1725	rVB	2245675	2867724	9.34%	1.137%
33	11.140	1725	1731	1736	rBV	1049796	1380814	4.50%	0.547%
34	11.360	1762	1767	1774	rBV	3249232	3872060	12.61%	1.535%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	11.439	1774	1780	1787	rVV	15691478	29248034	95.27%	11.596%
36	11.506	1787	1791	1801	rVB	8268465	10404348	33.89%	4.125%
37	11.670	1812	1818	1827	rBV	14602235	17947800	58.46%	7.116%
38	11.811	1836	1841	1846	rVB	25194013	30699604	100.00%	12.171%
39	11.945	1860	1863	1870	rVB	559089	706963	2.30%	0.280%
40	12.030	1870	1877	1880	rBV	1340397	1695286	5.52%	0.672%
41	12.079	1880	1885	1890	rVB3	1636236	2647253	8.62%	1.050%
42	12.146	1890	1896	1901	rBV	15533525	18234439	59.40%	7.229%
43	12.231	1906	1910	1916	rVB	287676	352660	1.15%	0.140%
44	12.298	1916	1921	1924	rBV	4221592	5699189	18.56%	2.260%
45	12.372	1929	1933	1942	rVB3	2769258	4014628	13.08%	1.592%
46	12.463	1943	1948	1953	rBV	817292	1049560	3.42%	0.416%
47	12.518	1953	1957	1963	rVB	2099855	2493998	8.12%	0.989%
48	12.585	1963	1968	1970	rBV	1871321	2336564	7.61%	0.926%
49	12.609	1970	1972	1977	rVB	2197902	2544299	8.29%	1.009%
50	12.670	1977	1982	1985	rBV	2425185	2735753	8.91%	1.085%
51	12.701	1985	1987	1991	rVB	590804	643675	2.10%	0.255%
52	12.756	1991	1996	2004	rBV2	1721242	2968002	9.67%	1.177%
53	12.908	2015	2021	2026	rVB2	1554216	2013607	6.56%	0.798%
54	12.981	2026	2033	2036	rBV	1164523	1599151	5.21%	0.634%
55	13.018	2036	2039	2045	rVB	2481636	2923179	9.52%	1.159%
56	13.085	2045	2050	2055	rBV4	262169	558237	1.82%	0.221%
57	13.195	2065	2068	2070	rBV2	320766	395507	1.29%	0.157%
58	13.225	2070	2073	2077	rVB	715930	817151	2.66%	0.324%
59	13.310	2083	2087	2089	rBV2	362584	506270	1.65%	0.201%
60	13.347	2089	2093	2098	rVB2	4624845	6103490	19.88%	2.420%
61	13.481	2110	2115	2122	rVB	2939726	3620595	11.79%	1.435%
62	13.560	2123	2128	2132	rBV3	212388	344521	1.12%	0.137%
63	13.640	2137	2141	2147	rVB2	437614	797929	2.60%	0.316%
64	13.725	2152	2155	2160	rVB2	502032	743858	2.42%	0.295%
65	13.835	2165	2173	2180	rVB	2638382	3520446	11.47%	1.396%
66	13.914	2180	2186	2190	rVB	377268	502367	1.64%	0.199%
67	13.987	2190	2198	2202	rBV2	542313	789928	2.57%	0.313%
68	14.286	2240	2247	2252	rBV	775286	1309429	4.27%	0.519%
69	14.542	2283	2289	2298	rBV2	788843	1035535	3.37%	0.411%
70	14.694	2307	2314	2318	rBV	1136472	1558770	5.08%	0.618%
71	14.841	2331	2338	2343	rBV	1017946	1289972	4.20%	0.511%

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Integration Parameters: RTEINT.P
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Sampling : 1 Min Area: 3 % of largest Peak
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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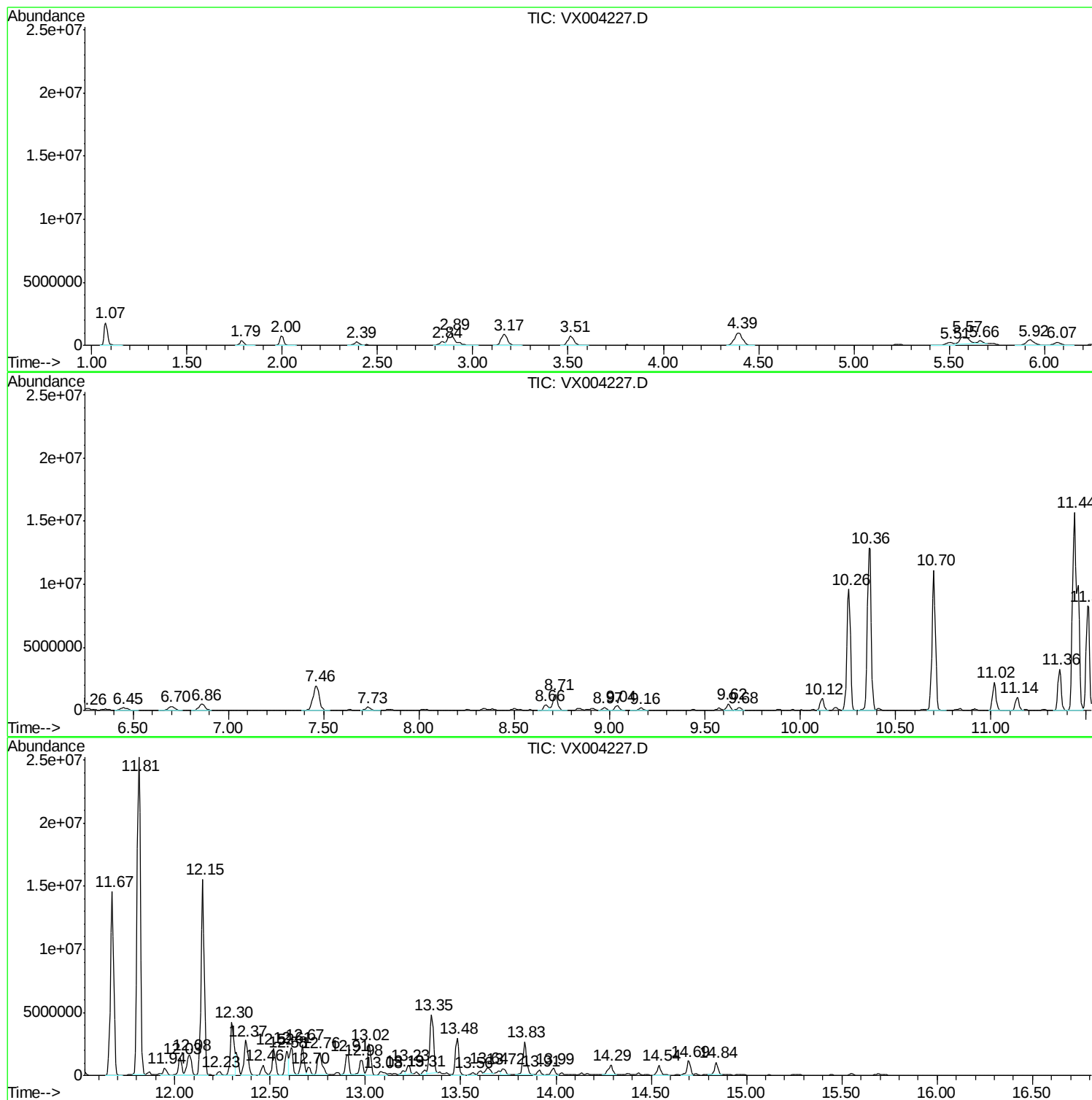
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA X/WATER
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Instrument :
MSVOA_X
ClientSampleId :
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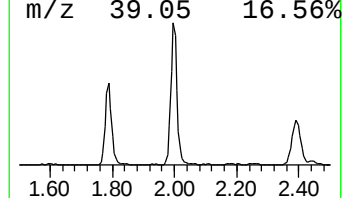
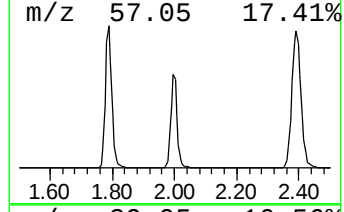
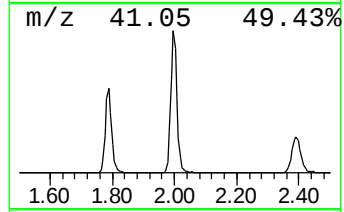
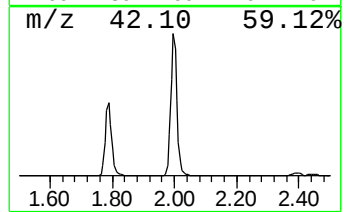
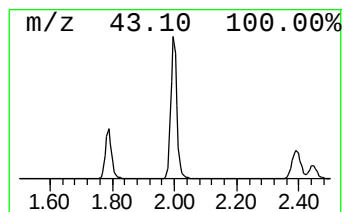
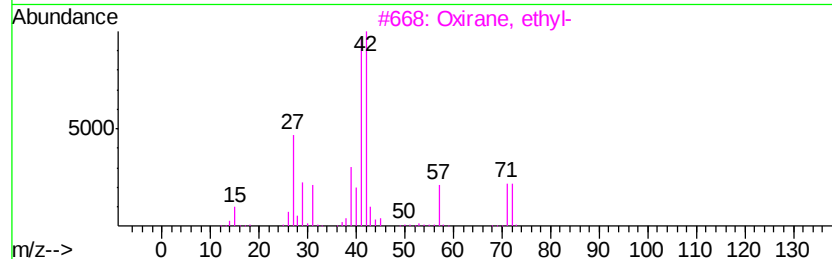
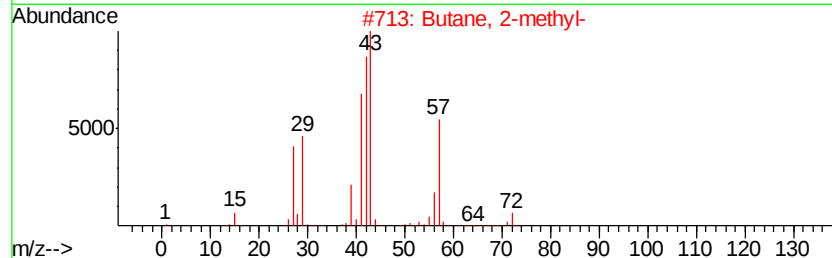
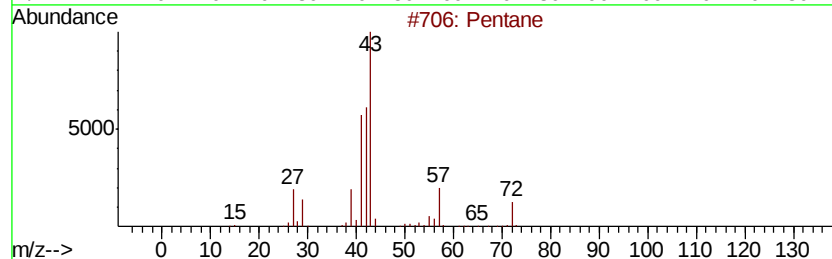
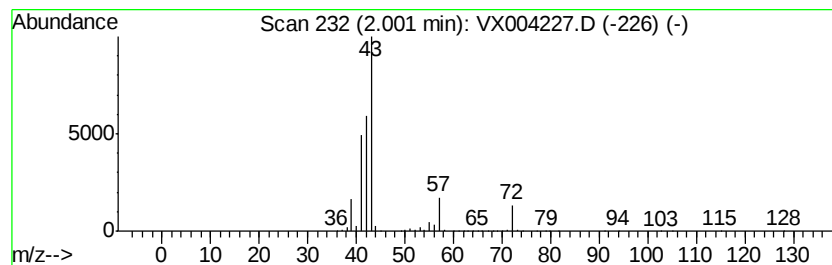
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	116.91 ug/l	1021030	Pentafluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	91
2			Butane, 2-methyl-	72	C5H12	000078-78-4	38
3			Oxirane, ethyl-	72	C4H8O	000106-88-7	33
4			Propanal, 2-methyl-	72	C4H8O	000078-84-2	9
5			2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	9



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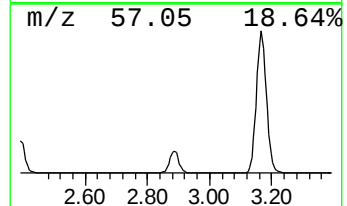
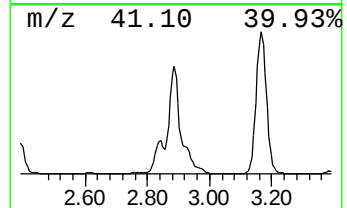
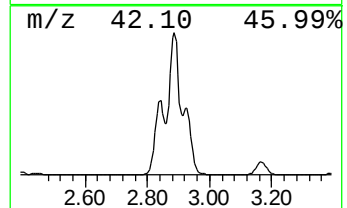
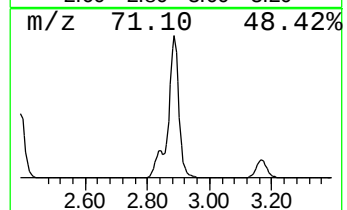
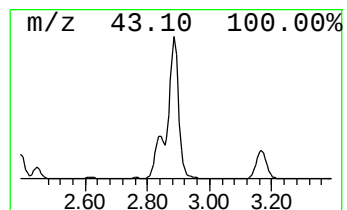
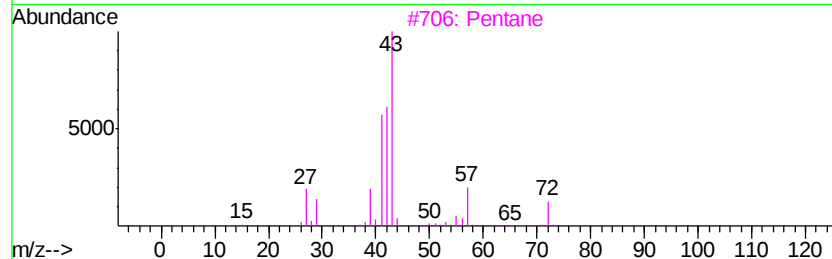
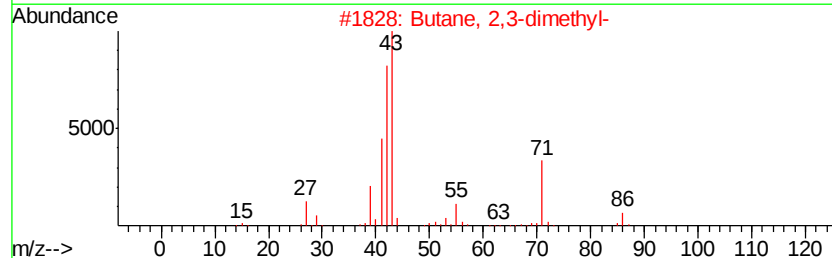
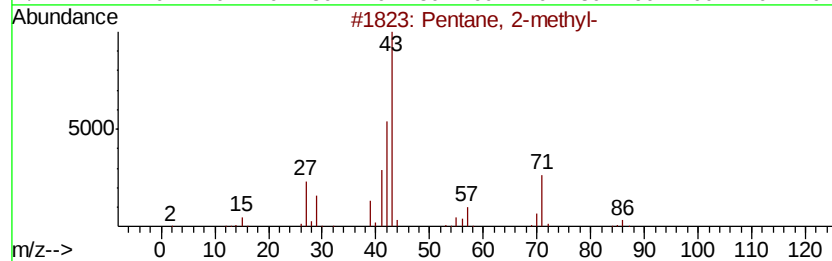
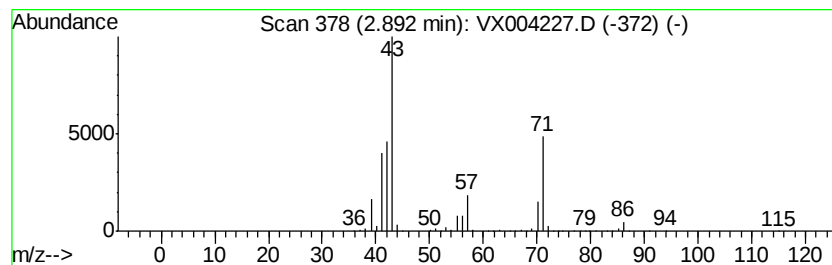
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	276.67 ug/l	2416390	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
3		Pentane	72	C5H12	000109-66-0	46
4		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28



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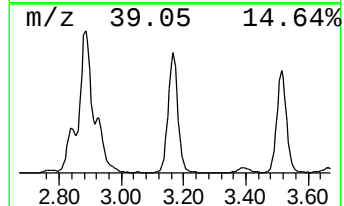
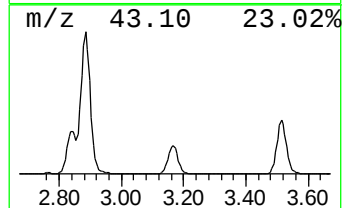
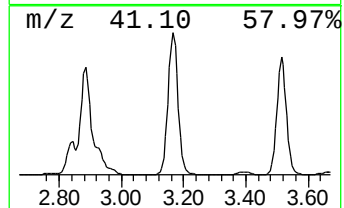
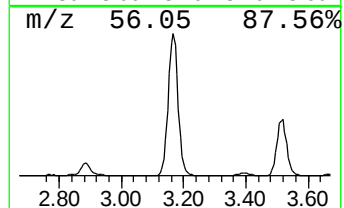
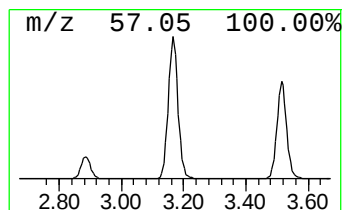
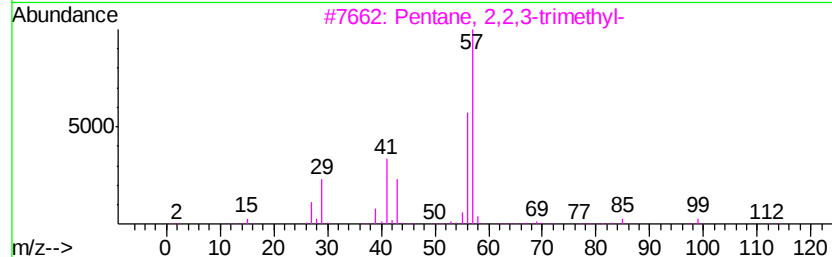
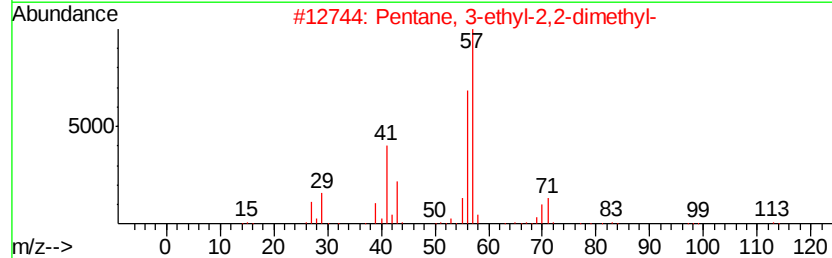
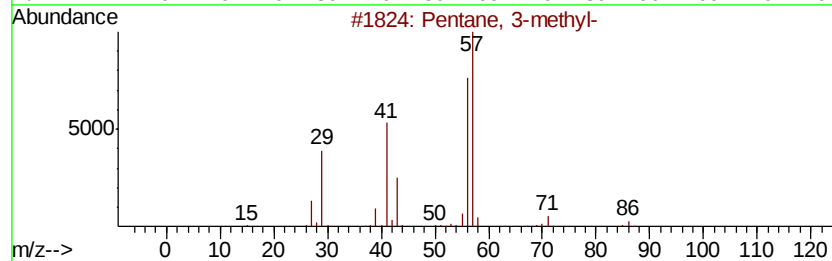
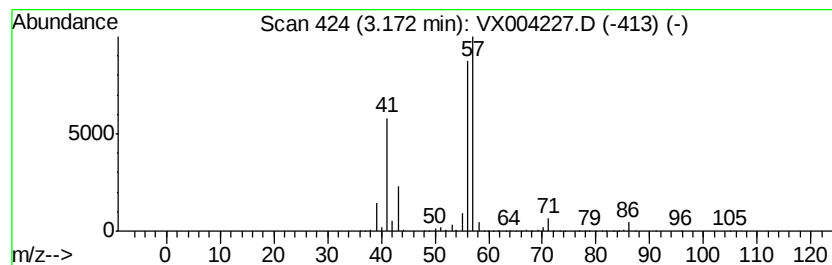
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	226.21 ug/l	1975630	Pentafluorobenzene	5.67

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	56
3		Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	39
4		Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	39
5		Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	38



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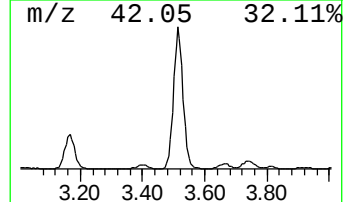
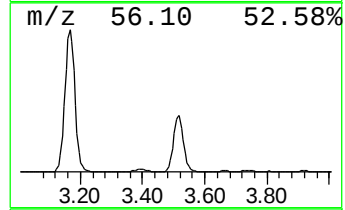
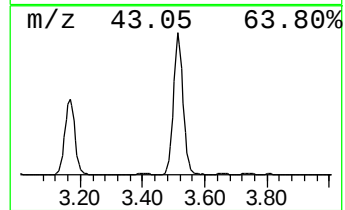
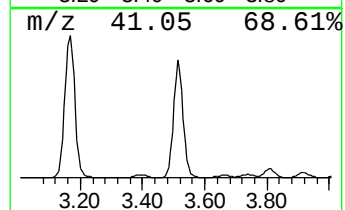
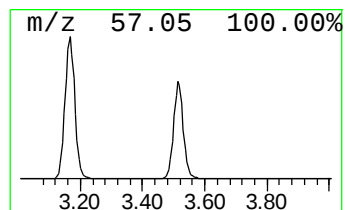
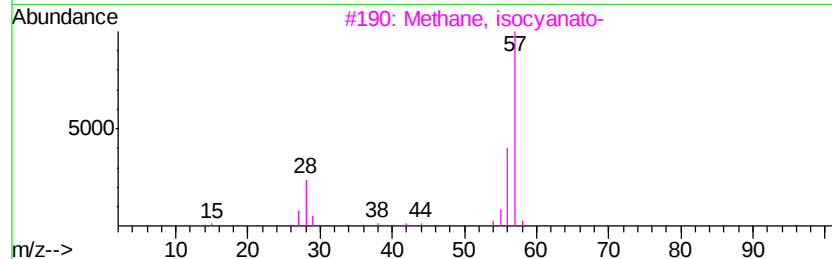
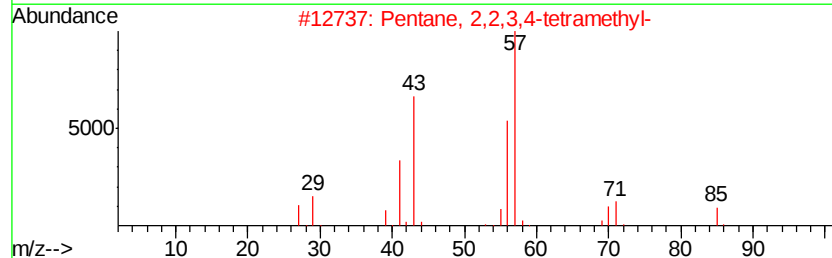
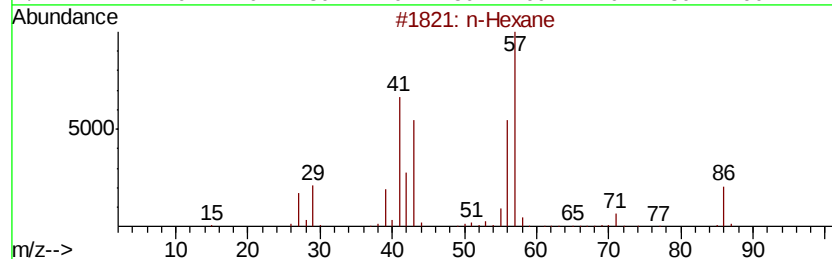
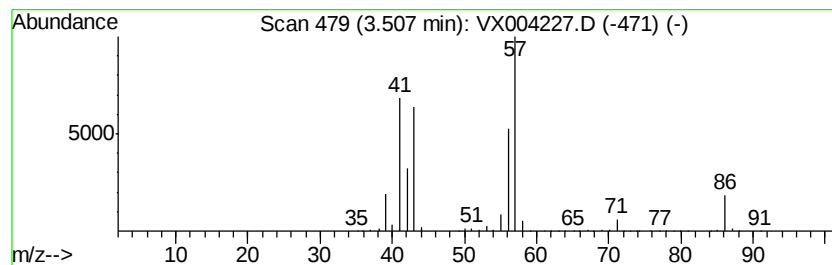
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Quant Title : SW846 8260

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

Peak Number 5 n-Hexane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.51	182.52 ug/l	1594100	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	94
2		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	42
3		Methane, isocyanato-	57	C2H3NO	000624-83-9	37
4		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	35
5		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	28



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082318\
Data File : VX004227.D
Acq On : 23 Aug 2018 15:26
Operator : JC/MD
Sample : J4579-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0561

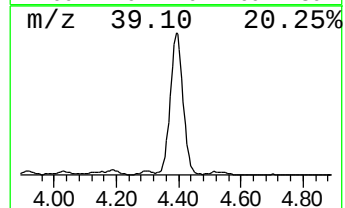
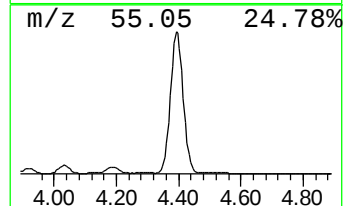
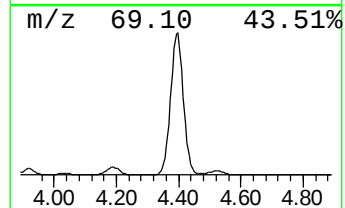
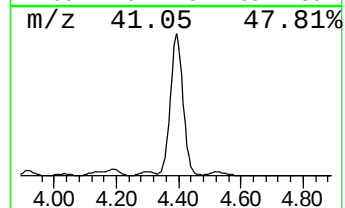
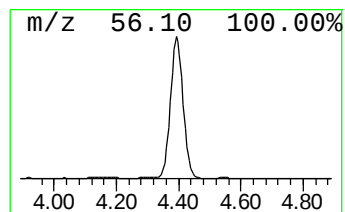
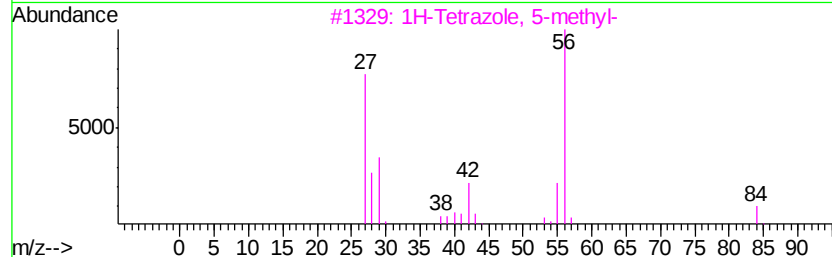
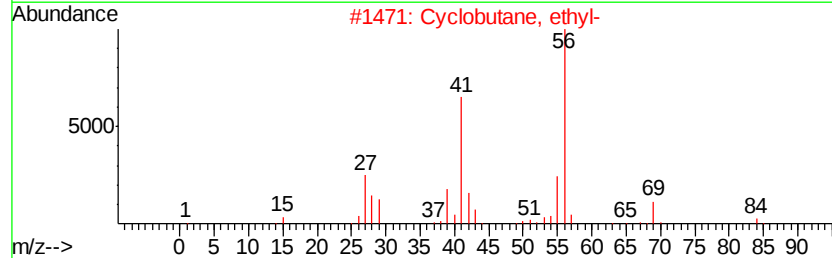
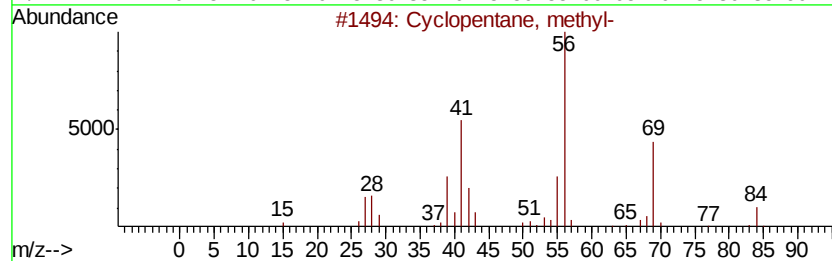
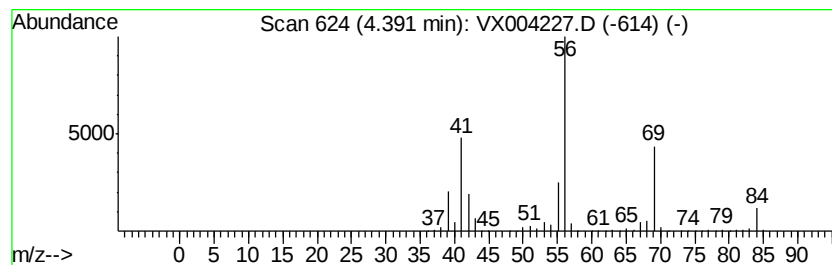
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	327.39 ug/l	2859330	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclohexane	84	C6H12	000110-82-7	64
5		Cyclobutane, butyl-	112	C8H16	013152-44-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082318\
Data File : VX004227.D
Acq On : 23 Aug 2018 15:26
Operator : JC/MD
Sample : J4579-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0561

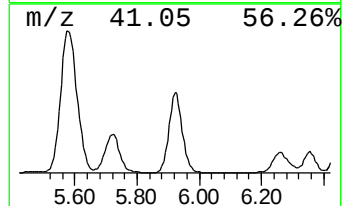
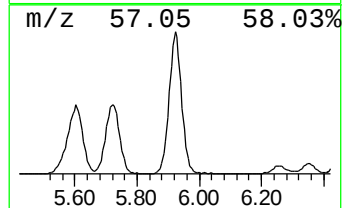
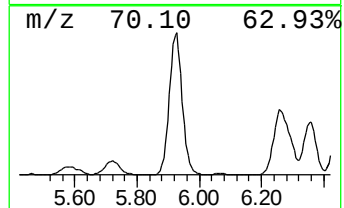
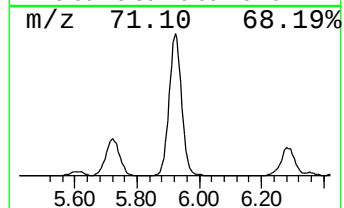
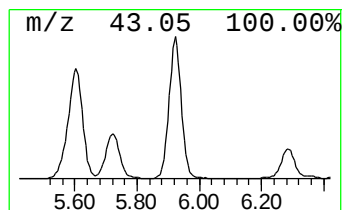
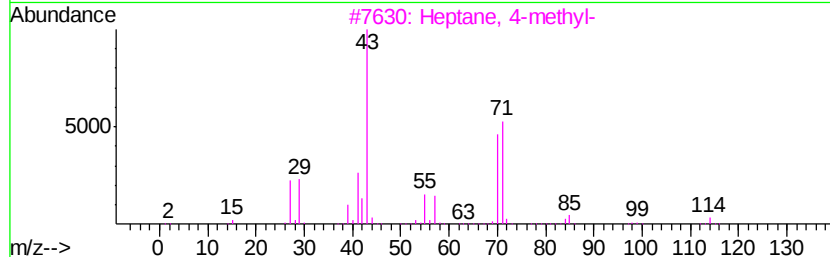
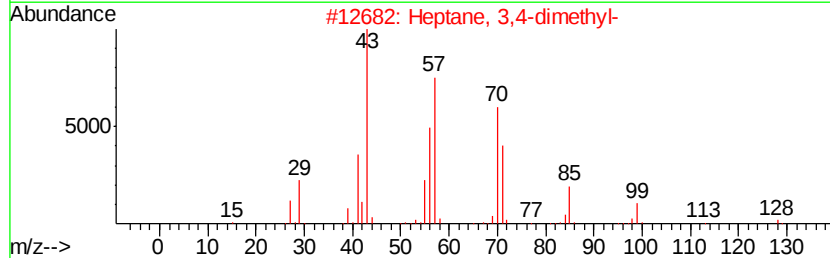
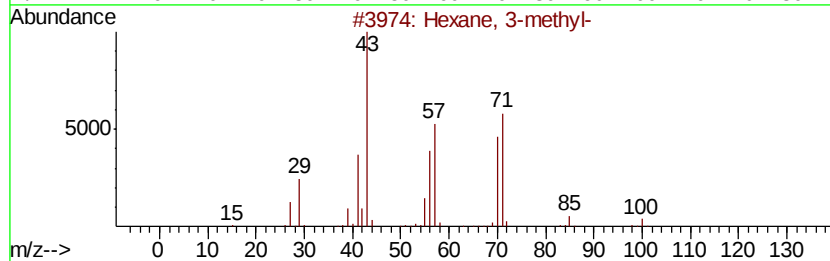
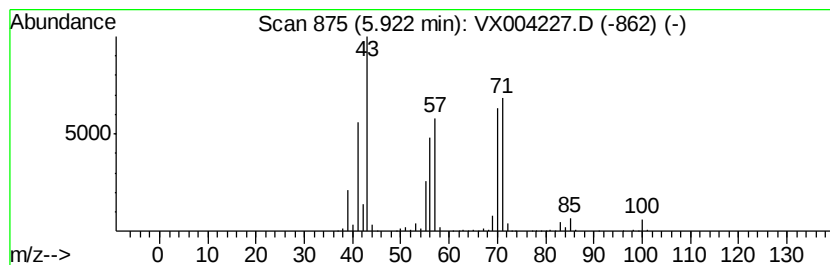
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Quant Title : SW846 8260

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

Peak Number 7 Hexane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	151.90 ug/l	1326660	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2		Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	72
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4		Heptane	100	C7H16	000142-82-5	58
5		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082318\
Data File : VX004227.D
Acq On : 23 Aug 2018 15:26
Operator : JC/MD
Sample : J4579-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

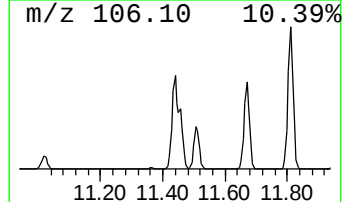
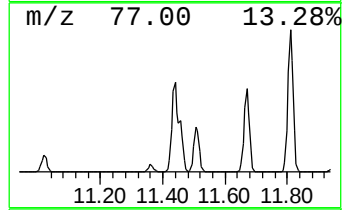
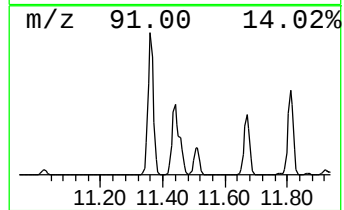
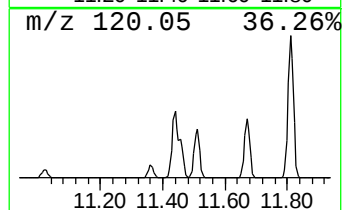
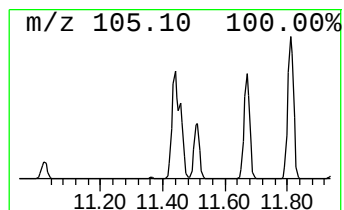
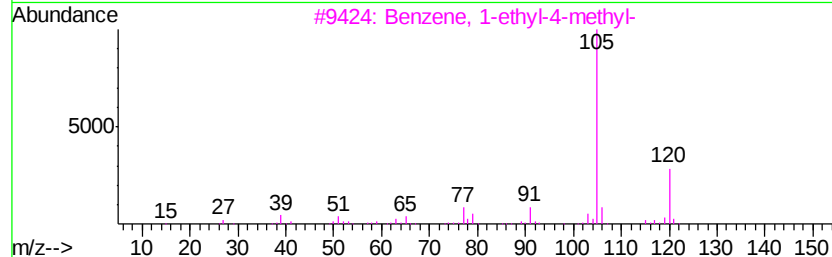
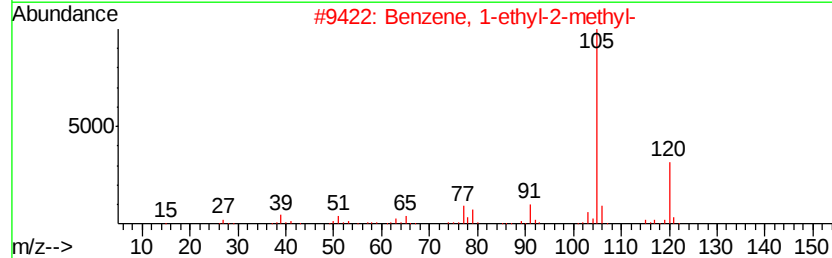
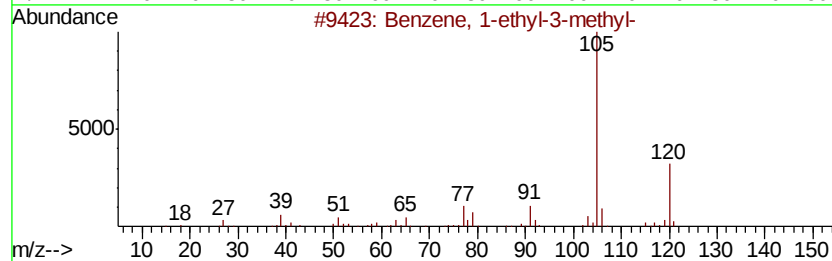
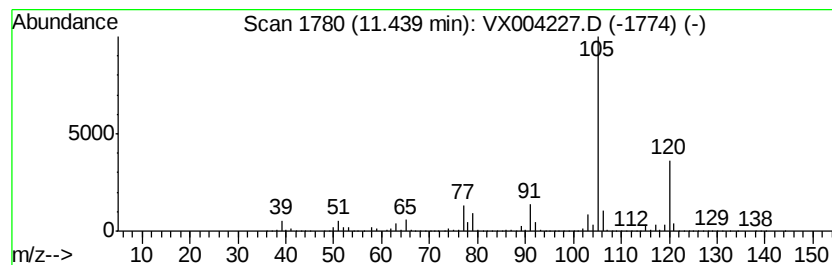
Instrument :
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ClientSampled :
FL-0561

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	552.42 ug/l	29248000	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 97
2		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 95
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91
4		Mesitylene	120 C9H12	000108-67-8 91
5		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082318\
Data File : VX004227.D
Acq On : 23 Aug 2018 15:26
Operator : JC/MD
Sample : J4579-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0561

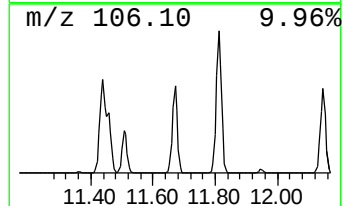
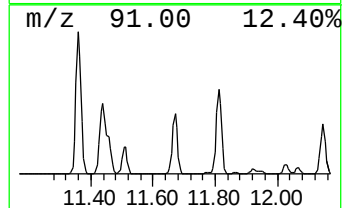
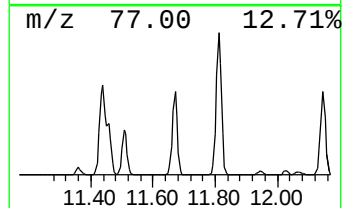
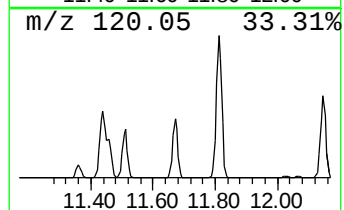
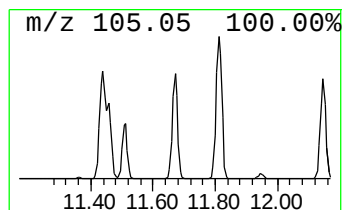
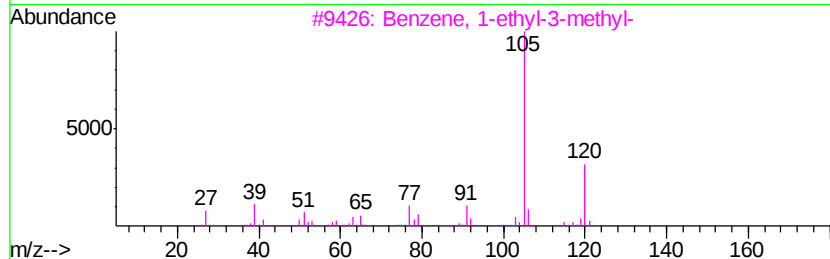
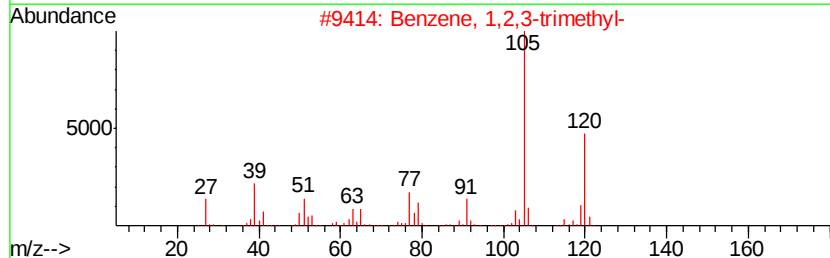
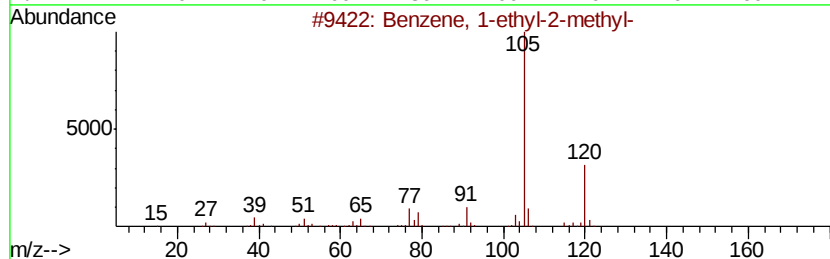
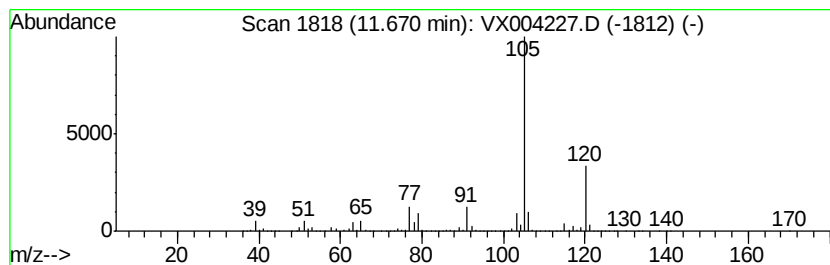
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	338.99 ug/l	17947800	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082318\
Data File : VX004227.D
Acq On : 23 Aug 2018 15:26
Operator : JC/MD
Sample : J4579-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0561

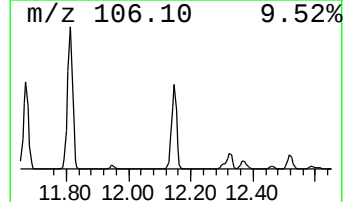
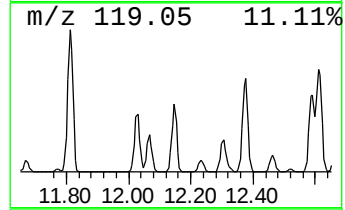
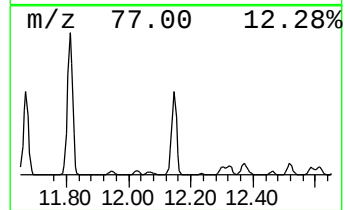
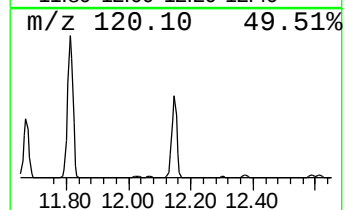
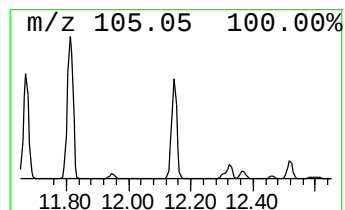
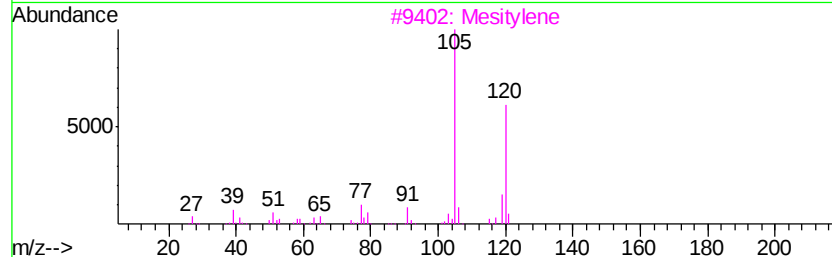
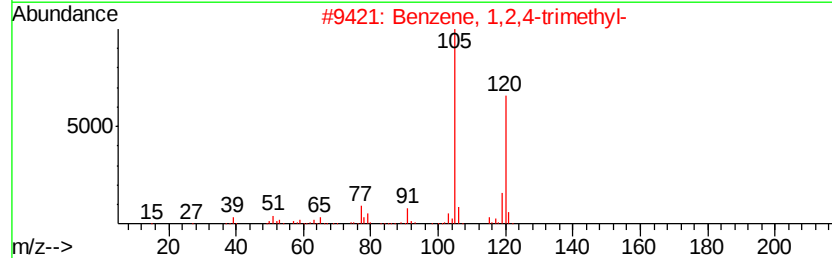
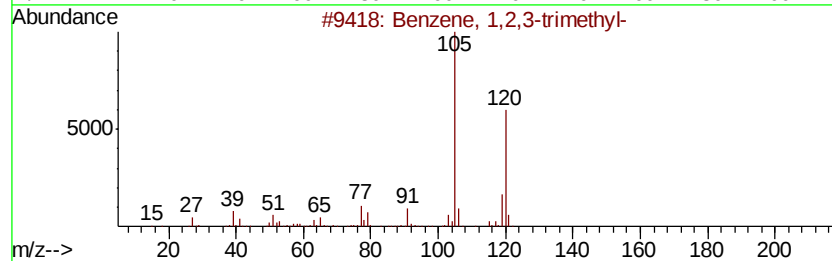
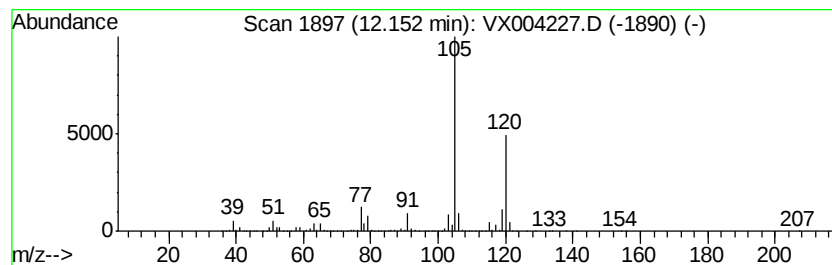
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	344.40 ug/l	18234400	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		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX082318\
Data File : VX004227.D
Acq On : 23 Aug 2018 15:26
Operator : JC/MD
Sample : J4579-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0561

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.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
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Pentane, 2-methyl-	2.89	276.7	ug/l	2416390	1	5.67	436687	50.0
Pentane, 3-methyl-	3.17	226.2	ug/l	1975630	1	5.67	436687	50.0
n-Hexane	3.51	182.5	ug/l	1594100	1	5.67	436687	50.0
Cyclopentane, met...	4.39	327.4	ug/l	2859330	1	5.67	436687	50.0
Hexane, 3-methyl-	5.92	151.9	ug/l	1326660	1	5.67	436687	50.0
Benzene, 1-ethyl-...	11.44	552.4	ug/l	29248000	4	12.08	2647250	50.0
Benzene, 1-ethyl-...	11.67	339.0	ug/l	17947800	4	12.08	2647250	50.0
Benzene, 1,2,3-tr...	12.15	344.4	ug/l	18234400	4	12.08	2647250	50.0