

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082218\
 Data File : VX004211.D
 Acq On : 22 Aug 2018 20:50
 Operator : JC/MD
 Sample : J4579-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0546

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	95	rBV	1644103	2128258	11.67%	1.285%
2	2.837	361	369	372	rBV	366377	736318	4.04%	0.445%
3	2.885	372	377	396	rVB	972249	2083726	11.43%	1.258%
4	3.166	413	423	440	rVB	826023	1867859	10.24%	1.128%
5	3.513	471	480	500	rVB	1082471	2480965	13.60%	1.498%
6	4.300	598	609	615	rBV	97914	294745	1.62%	0.178%
7	4.397	615	625	640	rVB	881152	2577990	14.14%	1.557%
8	5.501	791	806	810	rBV	207440	563157	3.09%	0.340%
9	5.580	810	819	828	rVV4	883863	3231459	17.72%	1.951%
10	5.665	828	833	837	rVV	319968	850718	4.66%	0.514%
11	5.720	838	842	862	rVB	332141	913079	5.01%	0.551%
12	5.928	862	876	890	rBV6	580560	2032140	11.14%	1.227%
13	6.068	890	899	909	rBV	221396	547517	3.00%	0.331%
14	6.257	918	930	938	rBV2	525739	1564130	8.58%	0.945%
15	6.354	938	946	954	rVV2	674843	1963341	10.77%	1.186%
16	6.452	954	962	976	rVB	932580	2566319	14.07%	1.550%
17	6.702	993	1003	1019	rVB	220566	527464	2.89%	0.319%
18	6.860	1019	1029	1038	rBV	471280	1111042	6.09%	0.671%
19	7.464	1114	1128	1141	rBV	8366686	18236782	100.00%	11.013%
20	7.732	1165	1172	1183	rVB	660298	1255495	6.88%	0.758%
21	7.848	1183	1191	1199	rVB	147957	298101	1.63%	0.180%
22	8.177	1236	1245	1254	rBV2	324039	738507	4.05%	0.446%
23	8.384	1275	1279	1286	rVB2	155991	298929	1.64%	0.181%
24	8.665	1318	1325	1329	rBV2	786542	1488319	8.16%	0.899%
25	8.714	1329	1333	1341	rVB	1254861	2144201	11.76%	1.295%
26	8.842	1348	1354	1358	rBV	360299	616461	3.38%	0.372%
27	8.909	1362	1365	1371	rVB	200300	312349	1.71%	0.189%
28	9.037	1381	1386	1398	rVB	681922	1123199	6.16%	0.678%
29	9.165	1398	1407	1412	rBV	339040	547198	3.00%	0.330%
30	9.573	1468	1474	1478	rBV2	258435	414481	2.27%	0.250%
31	9.622	1478	1482	1487	rVV	785041	1084827	5.95%	0.655%
32	9.677	1487	1491	1496	rVV	388379	552770	3.03%	0.334%
33	10.116	1558	1563	1568	rBV	951623	1233860	6.77%	0.745%
34	10.189	1568	1575	1580	rVV	317747	466006	2.56%	0.281%

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35	10.250	1580	1585	1591	rVB	611199	835118	4.58%	0.504%
36	10.360	1591	1603	1609	rBV	1473609	2252665	12.35%	1.360%
37	10.646	1642	1650	1655	rBV2	93560	239886	1.32%	0.145%
38	10.701	1655	1659	1670	rVB	524912	699535	3.84%	0.422%
39	10.835	1670	1681	1689	rBV3	173573	302372	1.66%	0.183%
40	10.914	1689	1694	1702	rBV4	110600	185924	1.02%	0.112%
41	11.018	1706	1711	1725	rVB	1932164	2395995	13.14%	1.447%
42	11.140	1725	1731	1735	rBV	1053042	1286778	7.06%	0.777%
43	11.360	1762	1767	1774	rBV	5049206	5970784	32.74%	3.606%
44	11.439	1774	1780	1782	rBV	8726247	12889539	70.68%	7.784%
45	11.506	1787	1791	1800	rVB	7501141	9426436	51.69%	5.693%
46	11.670	1812	1818	1826	rBV	3834199	4662757	25.57%	2.816%
47	11.811	1836	1841	1846	rVB	14957068	17257553	94.63%	10.422%
48	11.920	1852	1859	1861	rBV	217326	310961	1.71%	0.188%
49	11.945	1861	1863	1869	rVB	565596	665233	3.65%	0.402%
50	12.024	1869	1876	1880	rBV	1221530	1526574	8.37%	0.922%
51	12.079	1880	1885	1890	rVB3	1475905	2439320	13.38%	1.473%
52	12.146	1890	1896	1901	rBV	3840010	4652577	25.51%	2.810%
53	12.231	1905	1910	1915	rVB	248908	299805	1.64%	0.181%
54	12.304	1915	1922	1923	rBV2	975502	1185206	6.50%	0.716%
55	12.323	1923	1925	1929	rVV	1247179	1402690	7.69%	0.847%
56	12.371	1929	1933	1942	rVB3	1827294	2701757	14.81%	1.632%
57	12.463	1942	1948	1953	rBV	395155	480720	2.64%	0.290%
58	12.518	1953	1957	1964	rBV	971281	1188750	6.52%	0.718%
59	12.591	1964	1969	1970	rBV	965220	1215997	6.67%	0.734%
60	12.609	1970	1972	1977	rVV	1249081	1434922	7.87%	0.867%
61	12.670	1977	1982	1985	rVV	1299852	1504206	8.25%	0.908%
62	12.701	1985	1987	1992	rVV	515154	590258	3.24%	0.356%
63	12.756	1992	1996	2004	rVB2	1278136	2138753	11.73%	1.292%
64	12.902	2015	2020	2025	rVB2	783331	1044195	5.73%	0.631%
65	12.981	2025	2033	2036	rBV	675135	906770	4.97%	0.548%
66	13.018	2036	2039	2045	rVB	1359190	1627967	8.93%	0.983%
67	13.085	2045	2050	2055	rBV3	125395	238961	1.31%	0.144%
68	13.225	2070	2073	2077	rVB	372375	433807	2.38%	0.262%
69	13.347	2088	2093	2099	rVB2	2687081	3599081	19.74%	2.173%
70	13.481	2110	2115	2122	rVB	1082492	1282642	7.03%	0.775%
71	13.646	2137	2142	2146	rBV3	214404	342047	1.88%	0.207%

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Integration Parameters: RTEINT.P

Integrator: RTE
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Stop Thrs : 0 Peak Location: TOP

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72	13.725	2152	2155	2160	rVB2	253683	380396	2.09%	0.230%
73	13.835	2165	2173	2180	rVB	717136	1049363	5.75%	0.634%
74	13.987	2191	2198	2202	rBV3	142778	235843	1.29%	0.142%
75	14.274	2240	2245	2255	rVB3	224174	449194	2.46%	0.271%
76	14.542	2284	2289	2298	rBV2	157789	234689	1.29%	0.142%
77	14.700	2310	2315	2325	rVB	5136910	6391216	35.05%	3.860%
78	14.841	2333	2338	2348	rBV	3379428	4015133	22.02%	2.425%
79	15.273	2400	2409	2414	rBV	169191	266950	1.46%	0.161%
80	15.444	2428	2437	2441	rBV2	125809	240371	1.32%	0.145%
81	15.554	2449	2455	2459	rBV	338507	540386	2.96%	0.326%
82	15.688	2467	2477	2481	rBV	387184	671286	3.68%	0.405%
83	15.725	2481	2483	2491	rVB	285946	399656	2.19%	0.241%
84	15.914	2504	2514	2527	rVB	90495	246216	1.35%	0.149%

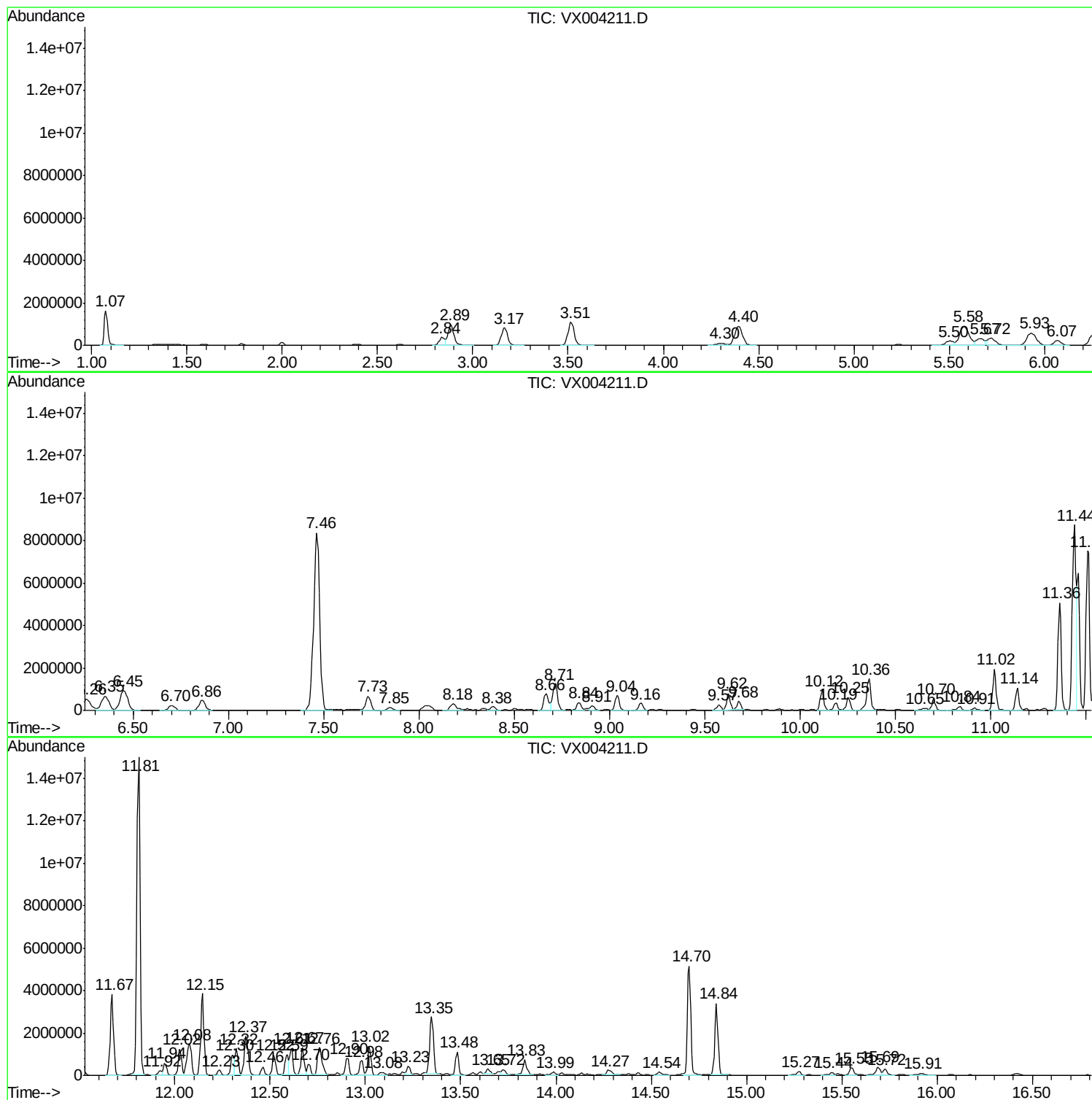
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260

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



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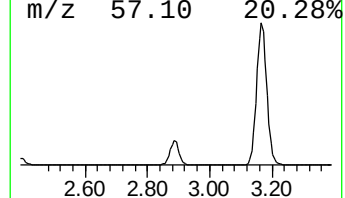
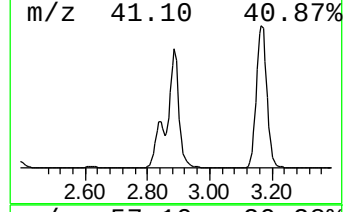
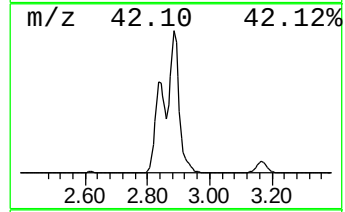
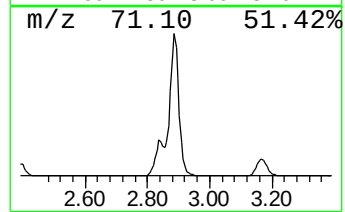
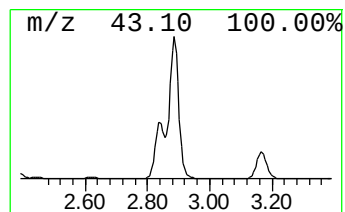
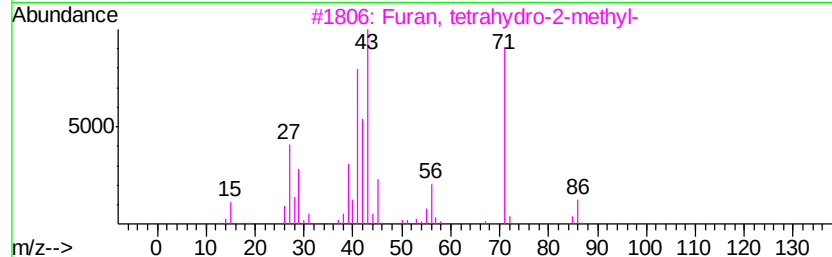
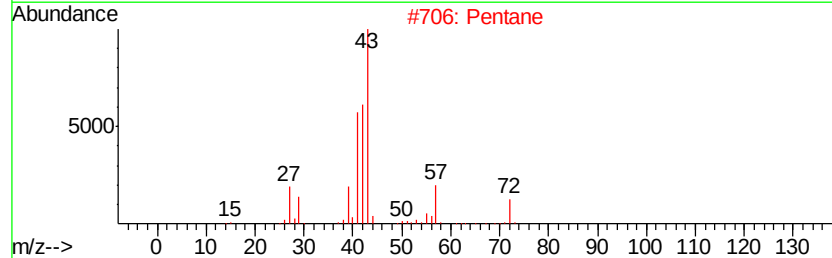
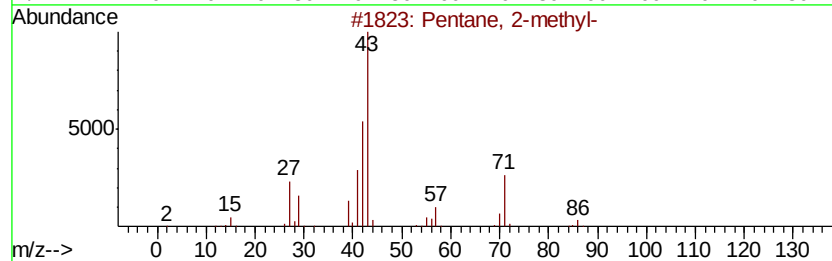
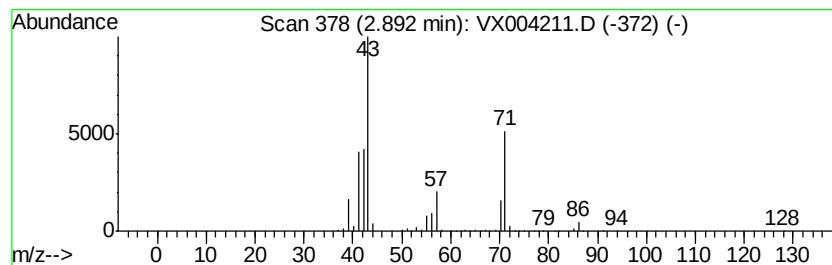
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	122.47 ug/l	2083730	Pentafluorobenzene	5.67

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	46
3		Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	43
4		Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
5		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38



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ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampled :
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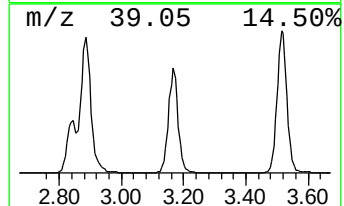
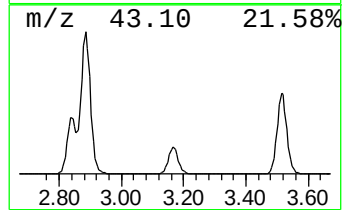
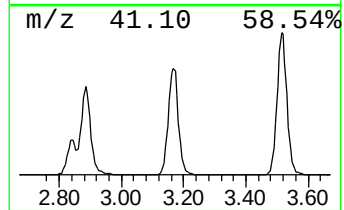
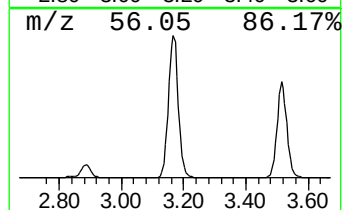
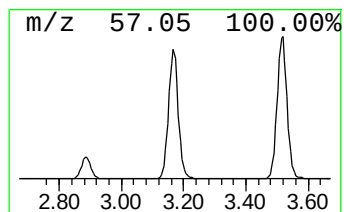
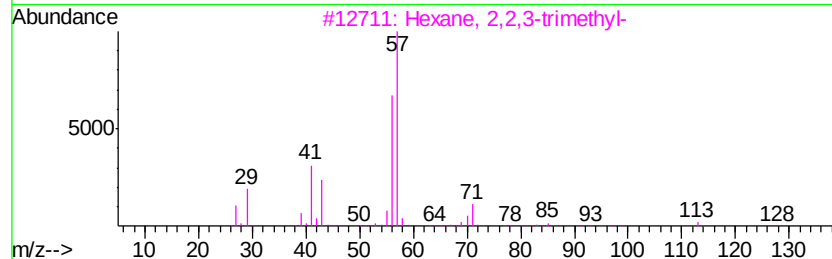
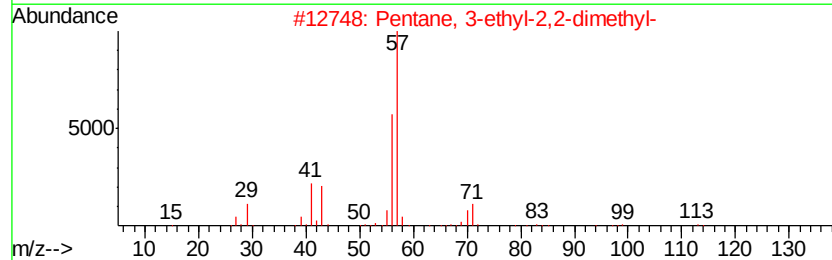
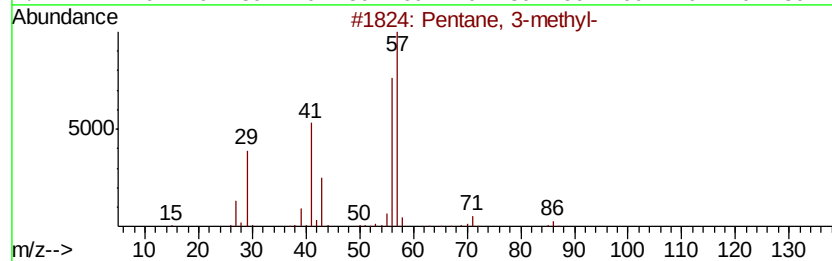
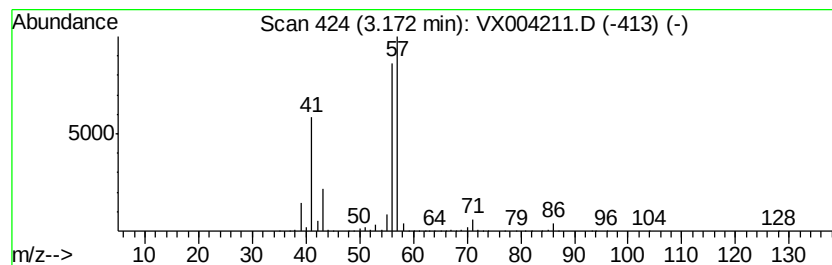
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Pentane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	109.78 ug/l	1867860	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	78
3		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	74
4		Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	64
5		Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	39



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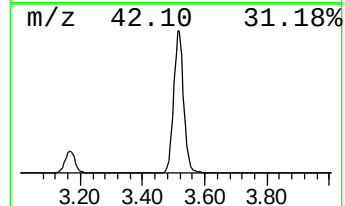
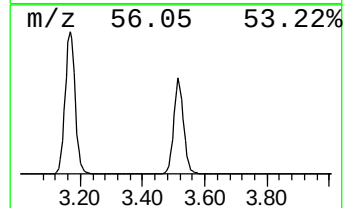
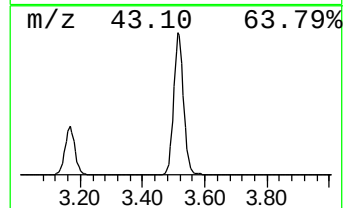
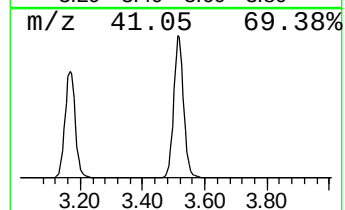
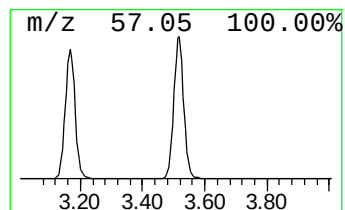
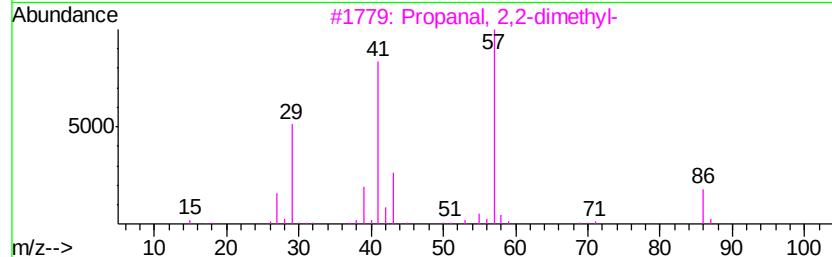
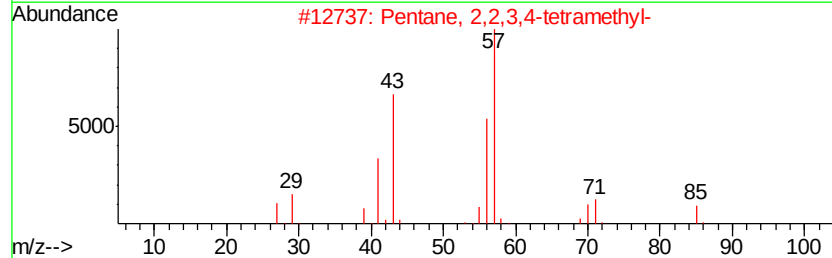
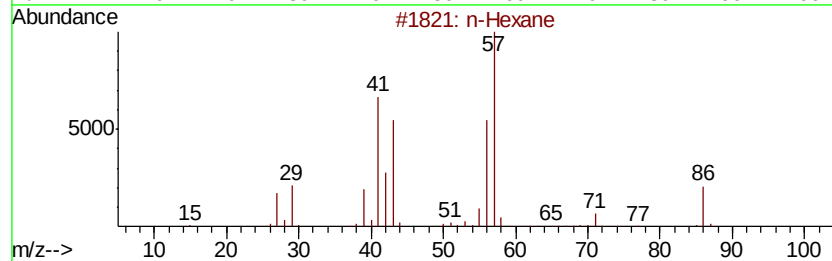
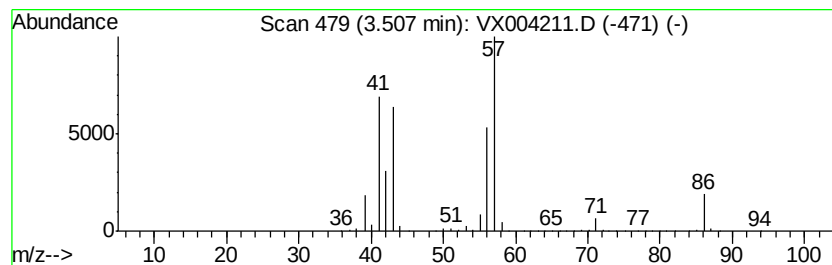
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Quant Title : SW846 8260

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

Peak Number 4 n-Hexane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.51	145.82 ug/l	2480970	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	40
3		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	38
4		Methane, isocyanato-	57	C2H3NO	000624-83-9	32
5		3-Pentanone	86	C5H10O	000096-22-0	30



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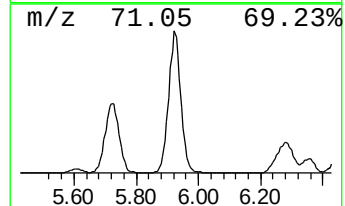
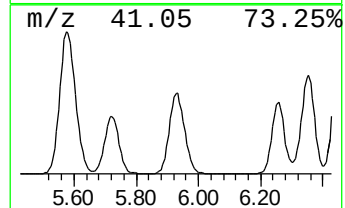
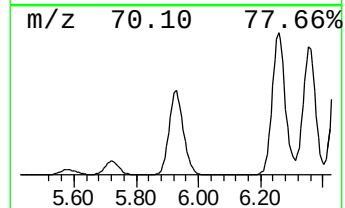
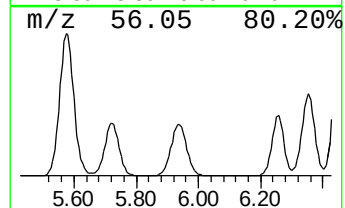
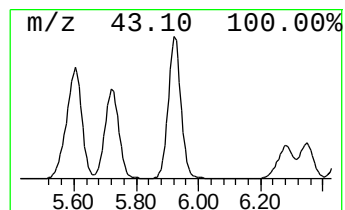
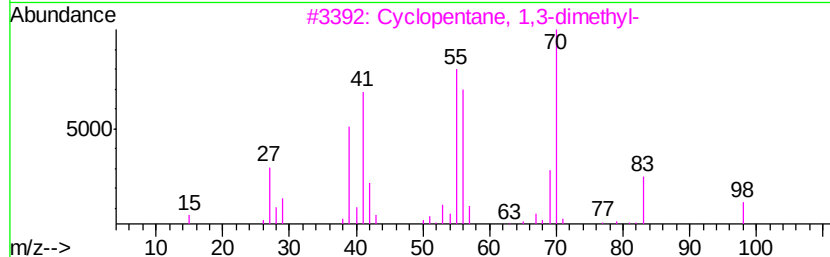
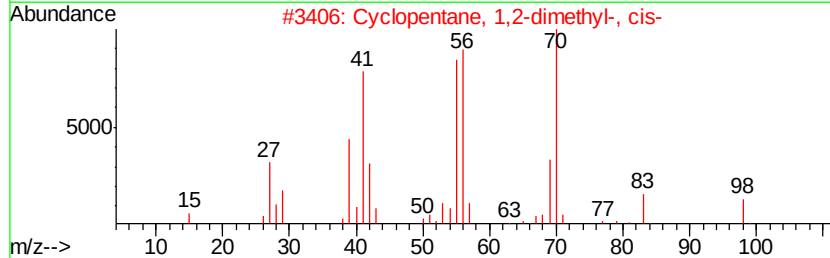
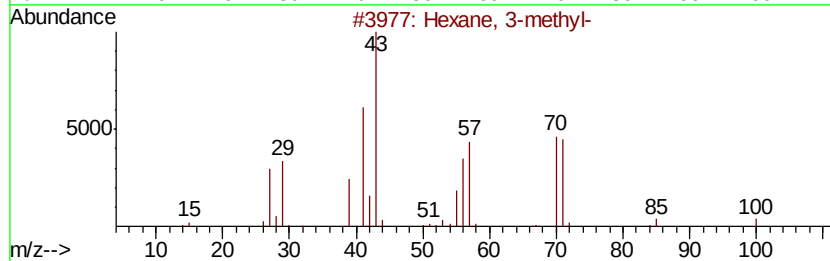
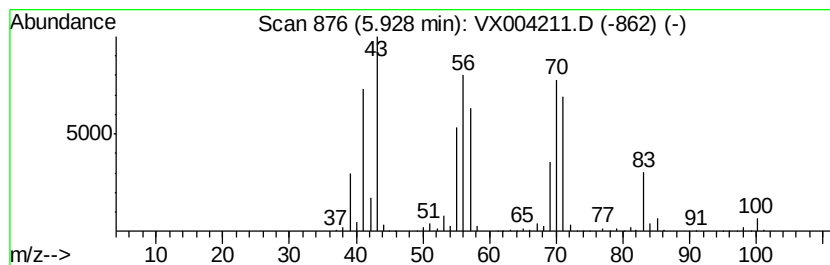
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	119.44 ug/l	2032140	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-		100	C7H16	000589-34-4	58
2	Cyclopentane, 1,2-dimethyl-, cis-		98	C7H14	001192-18-3	53
3	Cyclopentane, 1,3-dimethyl-		98	C7H14	002453-00-1	43
4	Hexane, 3,3,4-trimethyl-		128	C9H20	016747-31-2	43
5	Cyclopentane, 1,1-dimethyl-		98	C7H14	001638-26-2	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082218\
Data File : VX004211.D
Acq On : 22 Aug 2018 20:50
Operator : JC/MD
Sample : J4579-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0546

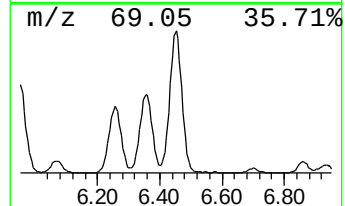
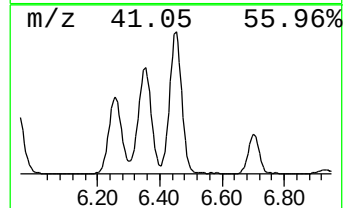
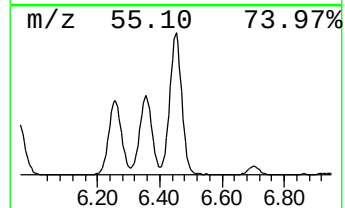
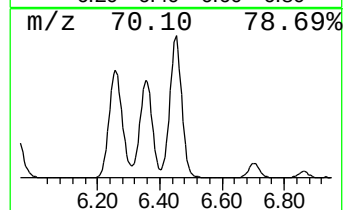
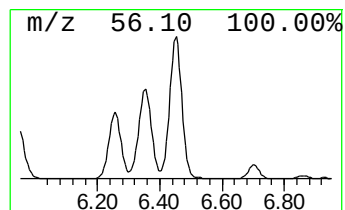
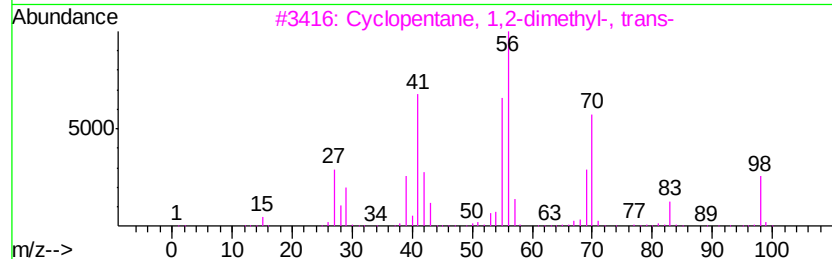
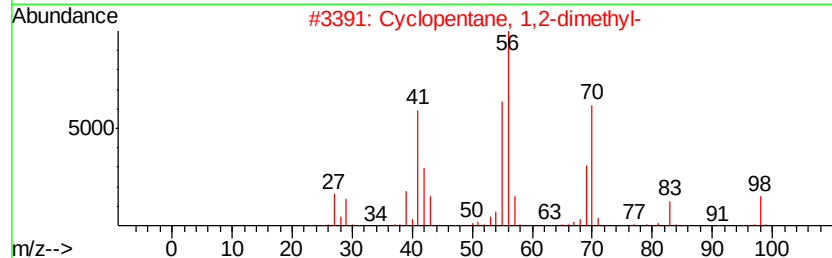
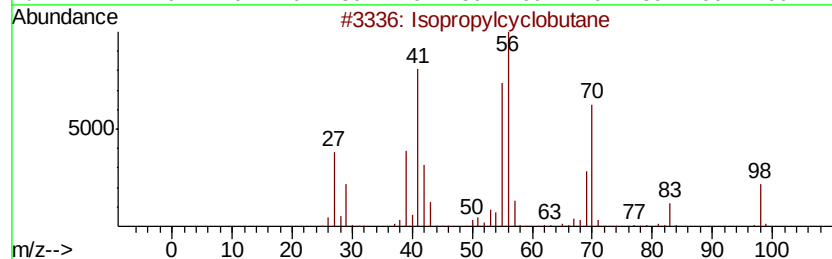
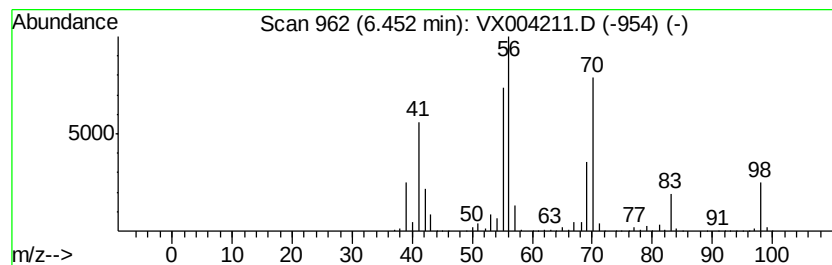
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Quant Title : SW846 8260

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

Peak Number 7 Isopropylcyclobutane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	115.49 ug/l	2566320	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropylcvclobutane	98	C7H14	000872-56-0	94
2		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
3		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
4		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	90
5		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082218\
Data File : VX004211.D
Acq On : 22 Aug 2018 20:50
Operator : JC/MD
Sample : J4579-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

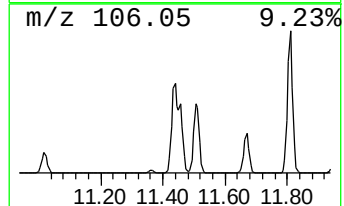
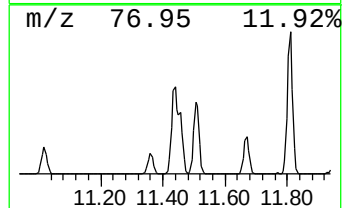
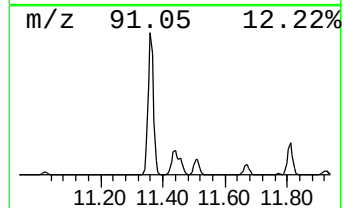
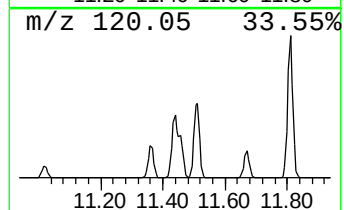
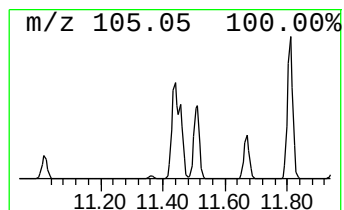
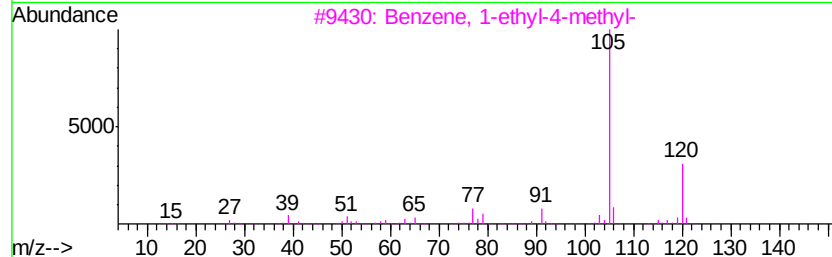
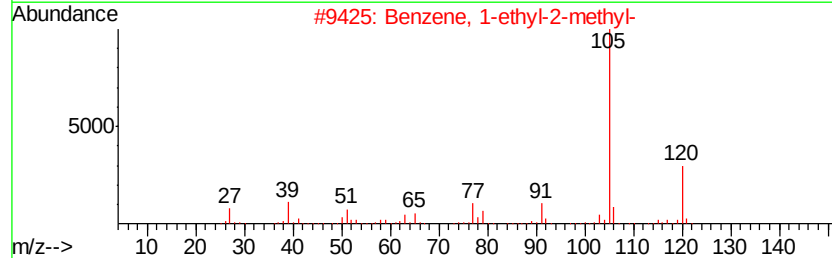
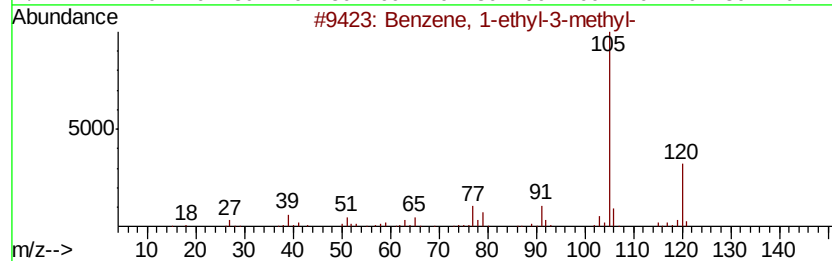
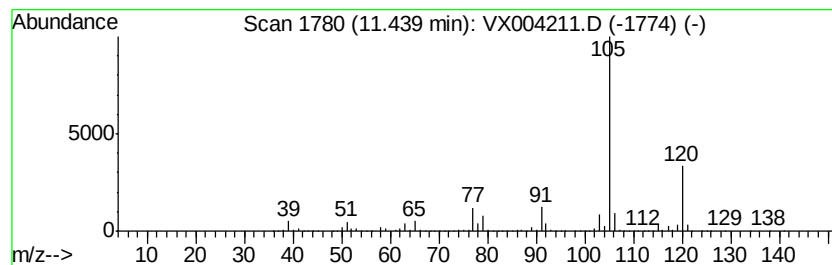
Instrument :
MSVOA_X
ClientSampled :
FL-0546

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	264.20 ug/l	12889500	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 95
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, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082218\
Data File : VX004211.D
Acq On : 22 Aug 2018 20:50
Operator : JC/MD
Sample : J4579-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
FL-0546

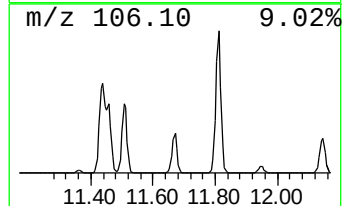
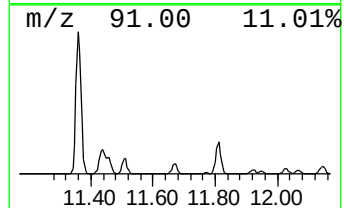
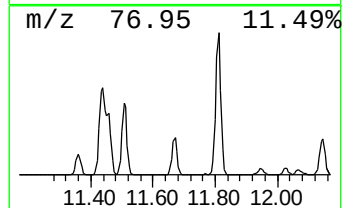
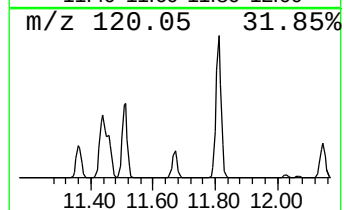
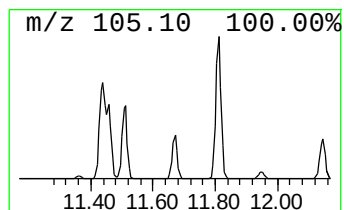
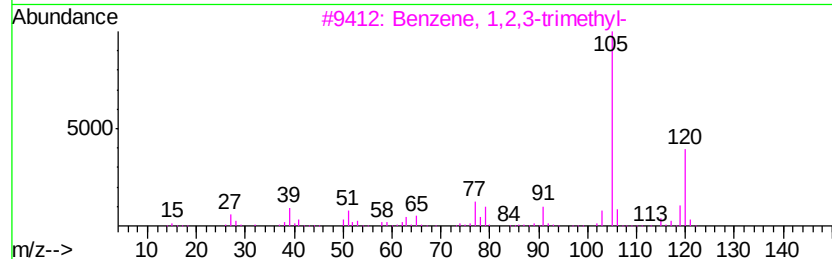
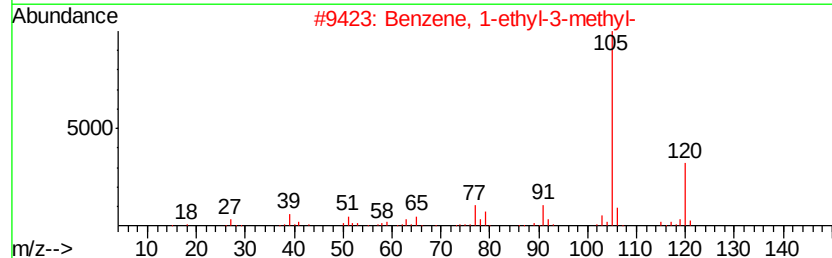
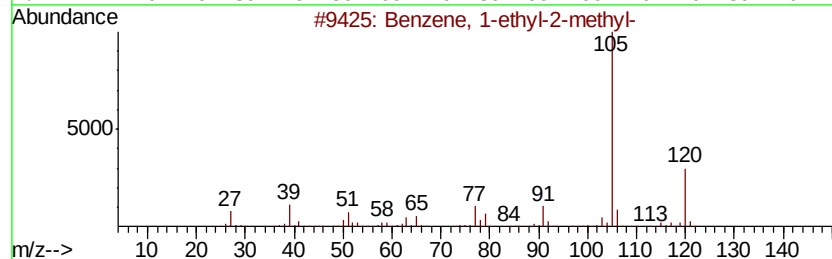
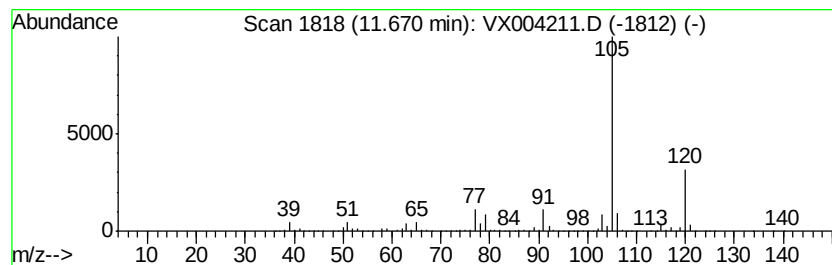
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	95.57 ug/l	4662760	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	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
4		Mesitylene	120	C9H12	000108-67-8	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082218\
Data File : VX004211.D
Acq On : 22 Aug 2018 20:50
Operator : JC/MD
Sample : J4579-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0546

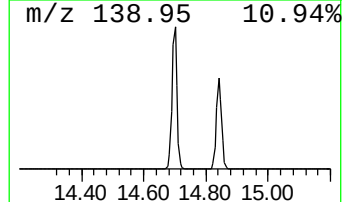
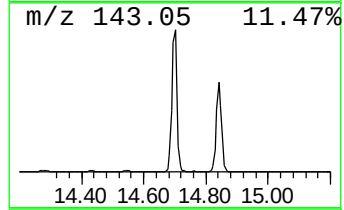
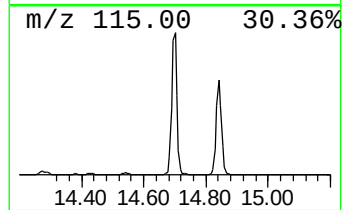
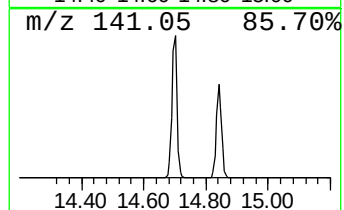
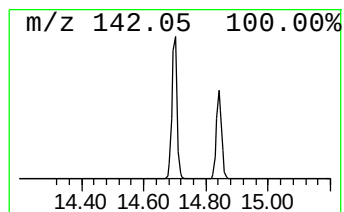
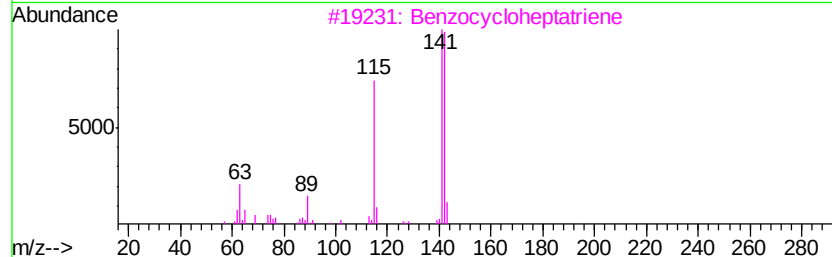
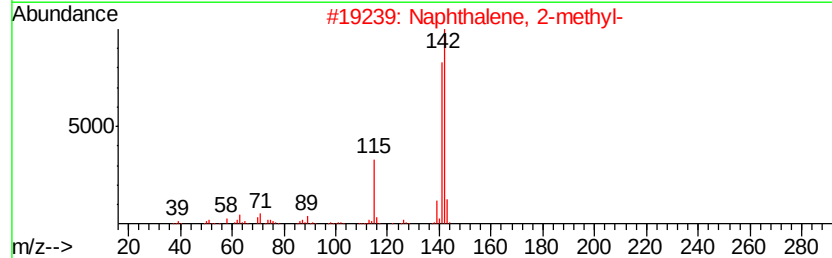
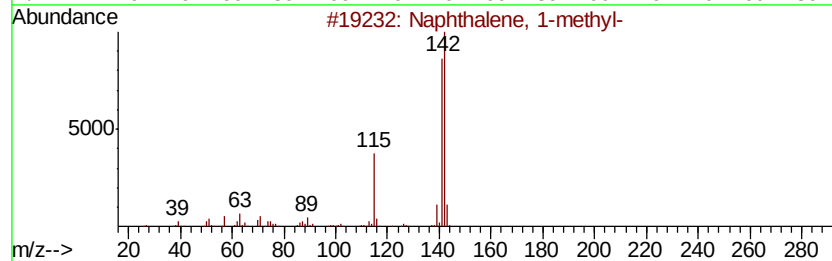
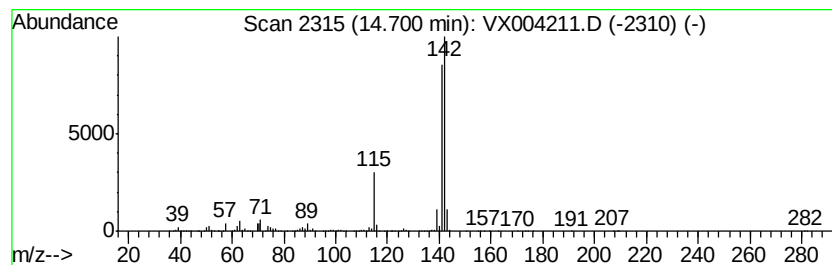
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	131.00 ug/l	6391220	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		3,4-Difluorobenzaldehyde	142	C7H4F2O	034036-07-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX082218\
Data File : VX004211.D
Acq On : 22 Aug 2018 20:50
Operator : JC/MD
Sample : J4579-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0546

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
Pentane, 2-methyl-	2.89	122.5	ug/l	2083730	1	5.67	850718	50.0
Pentane, 3-methyl-	3.17	109.8	ug/l	1867860	1	5.67	850718	50.0
n-Hexane	3.51	145.8	ug/l	2480970	1	5.67	850718	50.0
Hexane, 3-methyl-	5.93	119.4	ug/l	2032140	1	5.67	850718	50.0
Isopropylcyclobutane	6.45	115.5	ug/l	2566320	2	6.86	1111040	50.0
Benzene, 1-ethyl-...	11.44	264.2	ug/l	12889500	4	12.08	2439320	50.0
Benzene, 1-ethyl-...	11.67	95.6	ug/l	4662760	4	12.08	2439320	50.0
Naphthalene, 1-me...	14.70	131.0	ug/l	6391220	4	12.08	2439320	50.0