

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071619\
 Data File : VX010916.D
 Acq On : 16 Jul 2019 19:02
 Operator : JC/SP
 Sample : K3791-16
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0610

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\82X071619W.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	2.001	134	139	146	rVB	68535	94819	1.01%	0.112%
2	2.848	268	278	281	rBV	180134	401614	4.28%	0.474%
3	2.891	281	285	301	rVB	452263	950221	10.12%	1.121%
4	3.177	322	332	345	rBV	390086	904517	9.63%	1.067%
5	3.525	378	389	402	rBV	412064	957155	10.19%	1.129%
6	4.305	507	517	524	rBV2	48716	149959	1.60%	0.177%
7	4.403	524	533	548	rVB	464224	1359893	14.48%	1.604%
8	5.500	698	713	718	rBV	131178	373652	3.98%	0.441%
9	5.586	718	727	735	rVV2	493630	1737081	18.50%	2.049%
10	5.659	735	739	745	rVV	235371	636527	6.78%	0.751%
11	5.732	745	751	766	rVB2	165969	504788	5.38%	0.596%
12	5.939	772	785	797	rBV5	295636	1019761	10.86%	1.203%
13	6.061	797	805	818	rVB	127345	327242	3.48%	0.386%
14	6.262	825	838	846	rBV2	269824	810512	8.63%	0.956%
15	6.360	846	854	862	rVV2	342339	1005557	10.71%	1.186%
16	6.457	862	870	884	rVB	492141	1331562	14.18%	1.571%
17	6.707	900	911	923	rBV2	49376	121428	1.29%	0.143%
18	6.860	926	936	946	rBV	339062	812915	8.66%	0.959%
19	7.469	1020	1036	1046	rBV	4112923	9390583	100.00%	11.079%
20	7.732	1073	1079	1089	rVB	334766	653033	6.95%	0.770%
21	7.847	1091	1098	1108	rVB	76977	154925	1.65%	0.183%
22	8.055	1121	1132	1144	rVB3	114289	374866	3.99%	0.442%
23	8.177	1144	1152	1161	rBV2	170290	412265	4.39%	0.486%
24	8.390	1182	1187	1193	rVB3	81972	157342	1.68%	0.186%
25	8.664	1225	1232	1235	rBV2	422497	730175	7.78%	0.861%
26	8.713	1235	1240	1248	rVB	774536	1365248	14.54%	1.611%
27	8.841	1255	1261	1265	rBV	203117	338625	3.61%	0.400%
28	8.908	1269	1272	1279	rVB	107151	168908	1.80%	0.199%
29	9.036	1288	1293	1304	rVB	362062	612821	6.53%	0.723%
30	9.164	1304	1314	1319	rBV	190500	303448	3.23%	0.358%
31	9.573	1375	1381	1385	rBV	136199	219589	2.34%	0.259%
32	9.621	1385	1389	1394	rVV	430150	578675	6.16%	0.683%
33	9.676	1394	1398	1403	rVV	209259	297029	3.16%	0.350%
34	10.109	1464	1469	1475	rBV	712964	955188	10.17%	1.127%

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Integrator: RTE
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35	10.182	1475	1481	1486	rVV	175225	267852	2.85%	0.316%
36	10.249	1486	1492	1498	rVB	244700	332501	3.54%	0.392%
37	10.359	1501	1510	1516	rBV	891339	1376055	14.65%	1.624%
38	10.646	1547	1557	1562	rBV3	55826	152462	1.62%	0.180%
39	10.694	1562	1565	1577	rVB	266853	370059	3.94%	0.437%
40	10.835	1577	1588	1593	rBV2	99143	172333	1.84%	0.203%
41	11.018	1613	1618	1632	rVB	443470	578681	6.16%	0.683%
42	11.133	1632	1637	1642	rBV	641741	819068	8.72%	0.966%
43	11.359	1668	1674	1681	rBV	538676	664997	7.08%	0.785%
44	11.432	1681	1686	1688	rBV	4482556	5896048	62.79%	6.956%
45	11.505	1694	1698	1707	rVB	4236830	4957746	52.79%	5.849%
46	11.664	1719	1724	1732	rBV	2017487	2570804	27.38%	3.033%
47	11.804	1742	1747	1753	rVB	6499964	8074636	85.99%	9.527%
48	11.944	1767	1770	1775	rVB	126452	158037	1.68%	0.186%
49	12.023	1775	1783	1786	rBV	770111	897021	9.55%	1.058%
50	12.072	1786	1791	1797	rVB3	943777	1551372	16.52%	1.830%
51	12.139	1797	1802	1808	rBV	2105556	2562170	27.28%	3.023%
52	12.231	1812	1817	1822	rVB	150273	187973	2.00%	0.222%
53	12.298	1823	1828	1830	rBV2	598036	812229	8.65%	0.958%
54	12.322	1830	1832	1835	rVV	722012	740090	7.88%	0.873%
55	12.371	1835	1840	1848	rVB2	1148944	1577837	16.80%	1.862%
56	12.456	1849	1854	1859	rBV	248892	307421	3.27%	0.363%
57	12.517	1859	1864	1870	rBV	657904	793596	8.45%	0.936%
58	12.584	1870	1875	1877	rBV	687355	883208	9.41%	1.042%
59	12.609	1877	1879	1884	rVB	844114	854328	9.10%	1.008%
60	12.664	1884	1888	1891	rBV	800179	963979	10.27%	1.137%
61	12.700	1891	1894	1898	rVV	341842	394923	4.21%	0.466%
62	12.755	1898	1903	1911	rVB2	855208	1401744	14.93%	1.654%
63	12.901	1922	1927	1932	rVB2	545361	707989	7.54%	0.835%
64	12.974	1932	1939	1942	rBV	475252	599952	6.39%	0.708%
65	13.017	1942	1946	1952	rVB	934154	1110124	11.82%	1.310%
66	13.084	1952	1957	1961	rBV2	90610	170396	1.81%	0.201%
67	13.194	1971	1975	1977	rBV2	96162	127653	1.36%	0.151%
68	13.224	1977	1980	1983	rVB	243160	278300	2.96%	0.328%
69	13.304	1990	1993	1995	rBV2	81332	102367	1.09%	0.121%
70	13.340	1995	1999	2005	rVB2	1693924	2382213	25.37%	2.811%
71	13.474	2017	2021	2027	rVB	628792	817889	8.71%	0.965%

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72	13.560	2030	2035	2038	rVV2	68951	98610	1.05%	0.116%
73	13.596	2038	2041	2044	rVV	96682	135075	1.44%	0.159%
74	13.639	2044	2048	2053	rVV3	176640	348596	3.71%	0.411%
75	13.694	2053	2057	2059	rVV2	105042	178850	1.90%	0.211%
76	13.718	2059	2061	2067	rVV2	165465	229774	2.45%	0.271%
77	13.834	2071	2080	2088	rVB2	369635	598856	6.38%	0.707%
78	13.986	2097	2105	2108	rBV2	85832	134091	1.43%	0.158%
79	14.273	2145	2152	2160	rBV3	127331	259492	2.76%	0.306%
80	14.535	2190	2195	2207	rBV2	88697	137929	1.47%	0.163%
81	14.694	2216	2221	2231	rVB	2373302	2887298	30.75%	3.407%
82	14.834	2239	2244	2253	rBV	1429799	1901492	20.25%	2.243%
83	15.267	2306	2315	2320	rBV2	63039	110064	1.17%	0.130%
84	15.444	2334	2344	2347	rBV2	59047	118214	1.26%	0.139%
85	15.547	2353	2361	2365	rBV	154341	243119	2.59%	0.287%
86	15.687	2378	2384	2387	rBV	158747	254363	2.71%	0.300%
87	15.724	2387	2390	2398	rVB	122365	180179	1.92%	0.213%
88	15.913	2411	2421	2432	rVB2	42801	112952	1.20%	0.133%

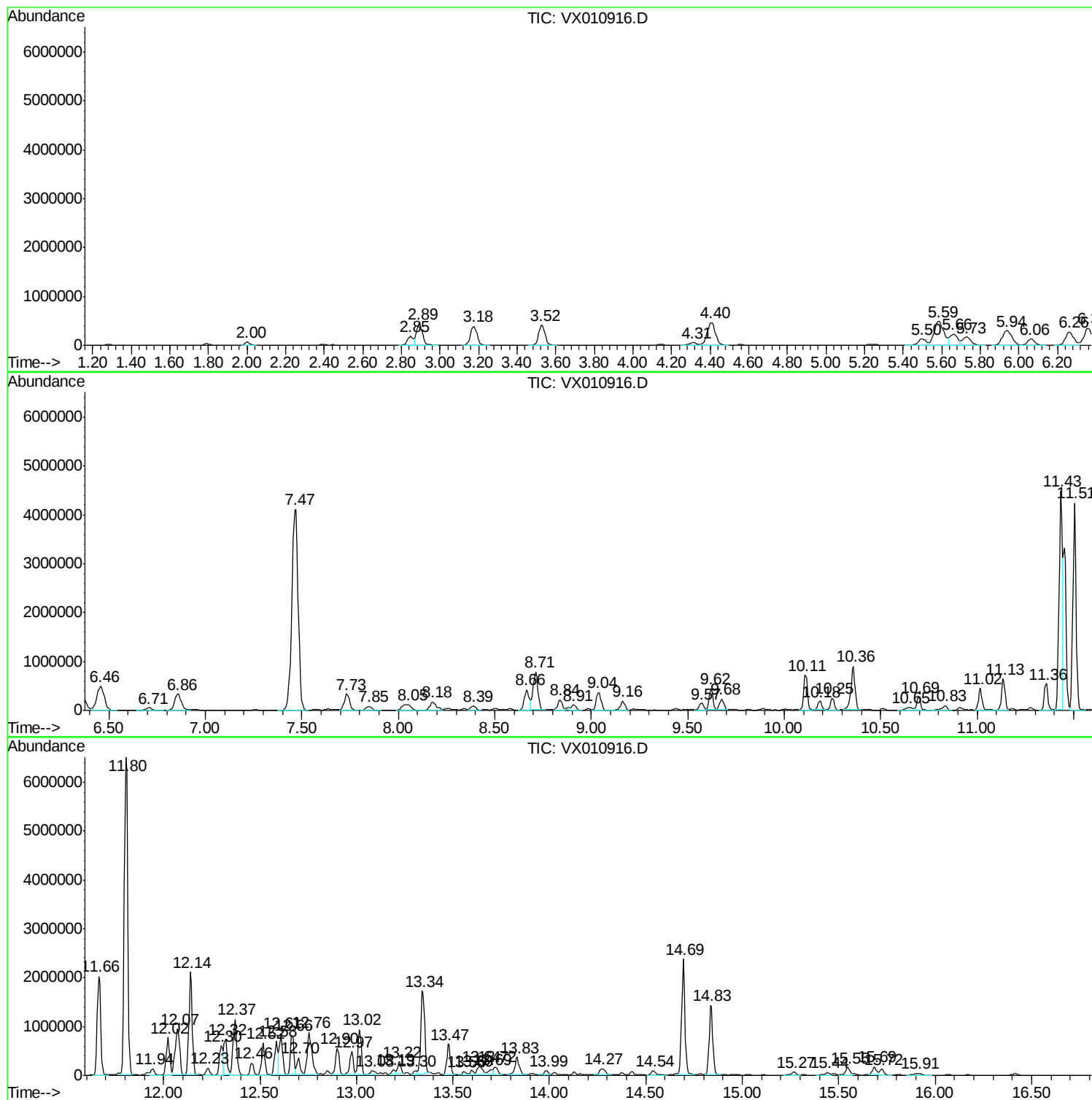
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampled :
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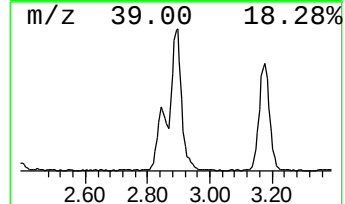
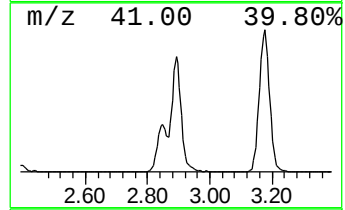
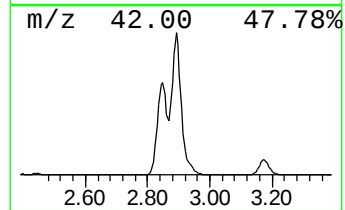
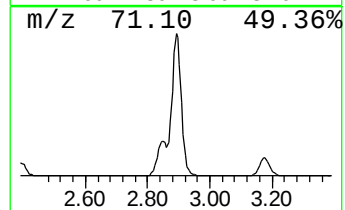
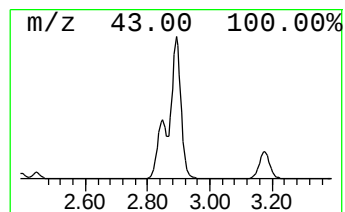
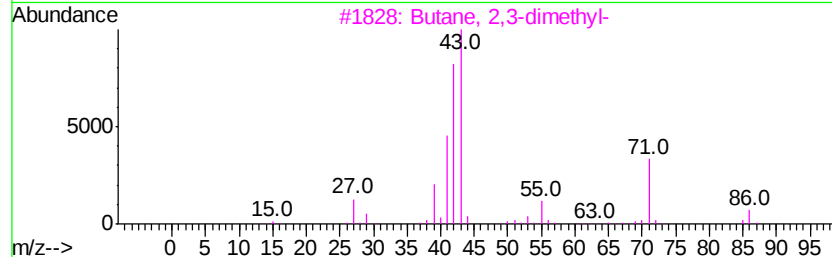
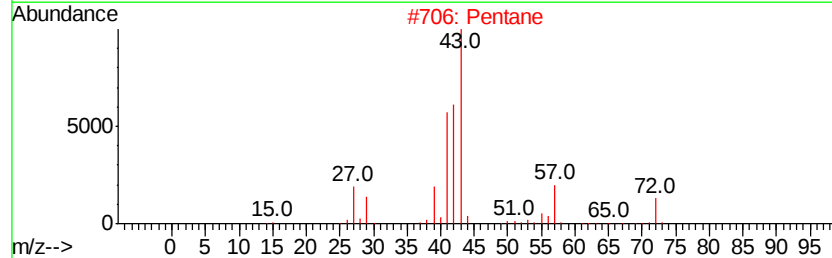
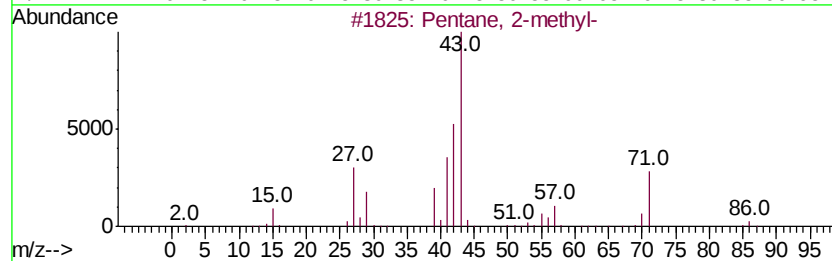
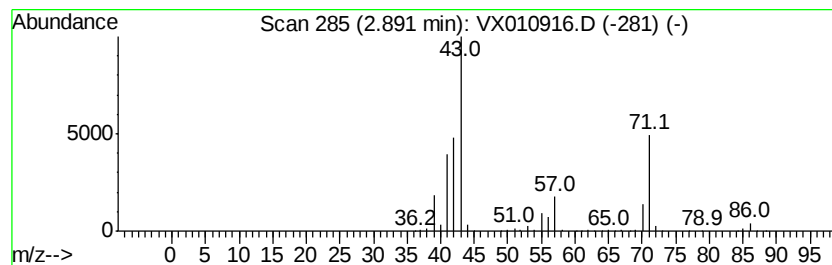
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	74.64 ug/l	950221	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	86
2		Pentane	72	C5H12	000109-66-0	49
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
4		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	42
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40



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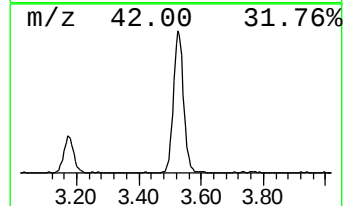
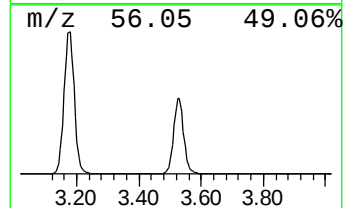
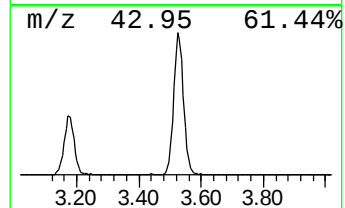
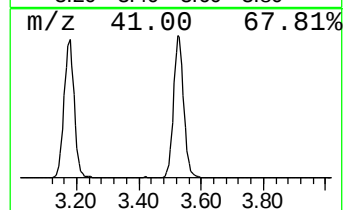
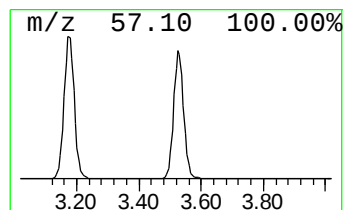
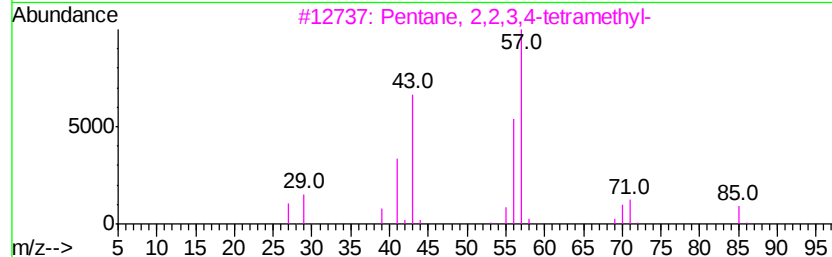
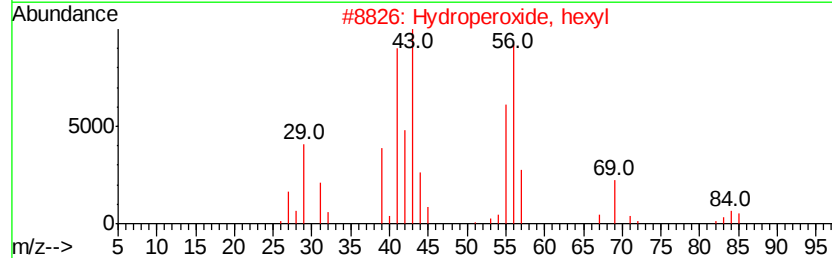
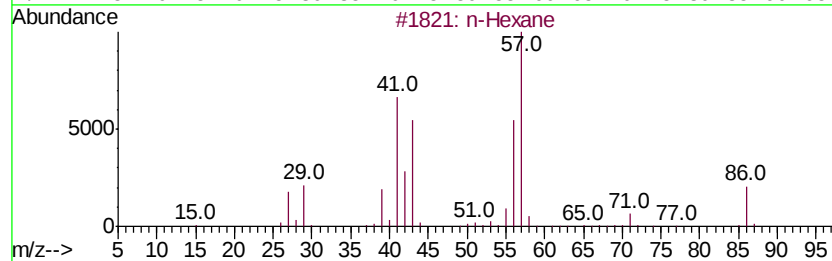
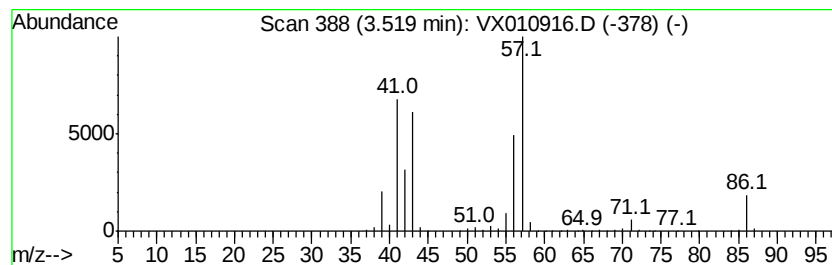
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	75.19 ug/l	957155	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	94
2		Hydroperoxide, hexyl	118	C6H14O2	004312-76-9	47
3		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	42
4		Methane, isocyanato-	57	C2H3NO	000624-83-9	37
5		Butanal, 2-methyl-	86	C5H10O	000096-17-3	35



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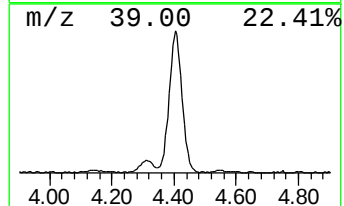
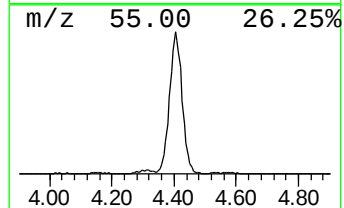
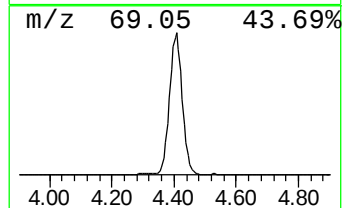
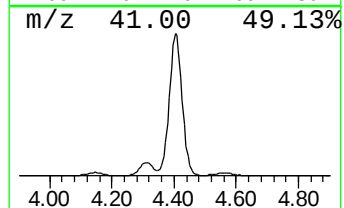
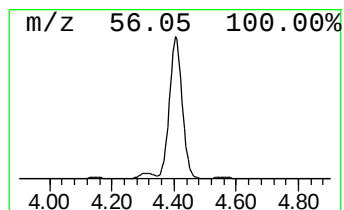
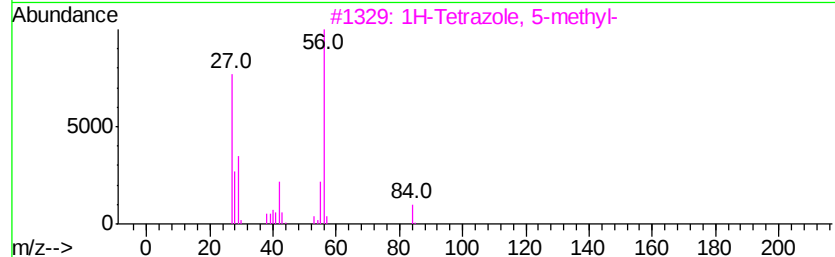
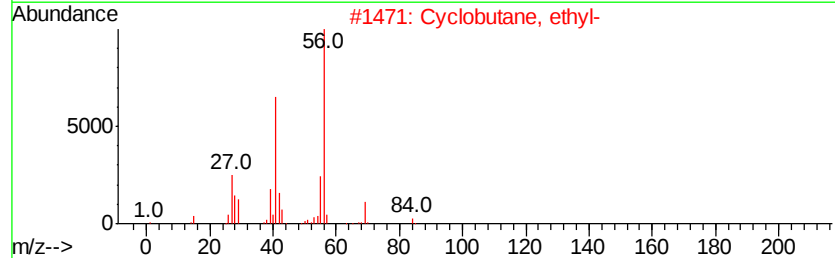
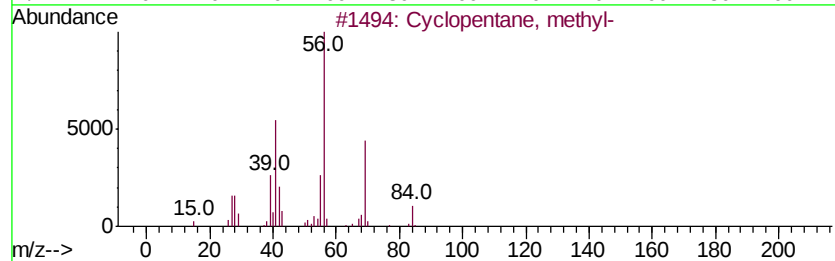
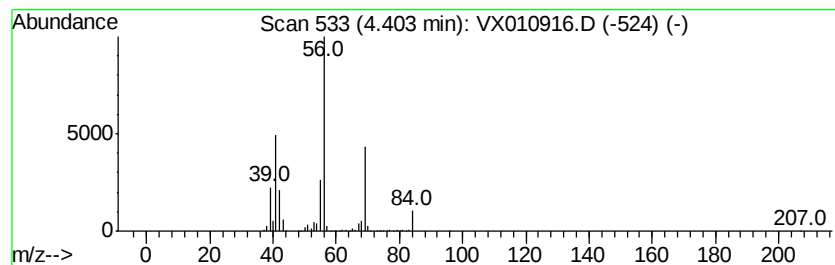
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	106.82 ug/l	1359890	Pentafluorobenzene	5.66

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	72
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4		Cyclopropane, propyl-	84	C6H12	002415-72-7	50
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	46



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FL-0610

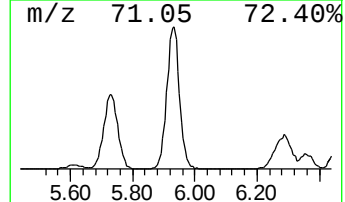
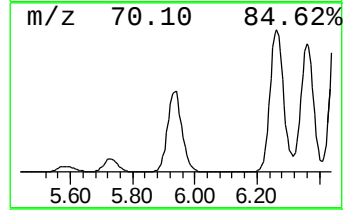
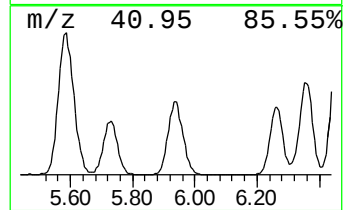
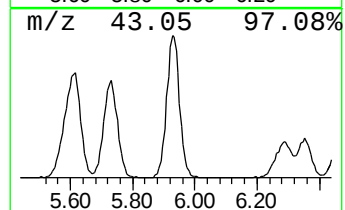
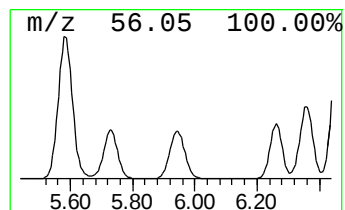
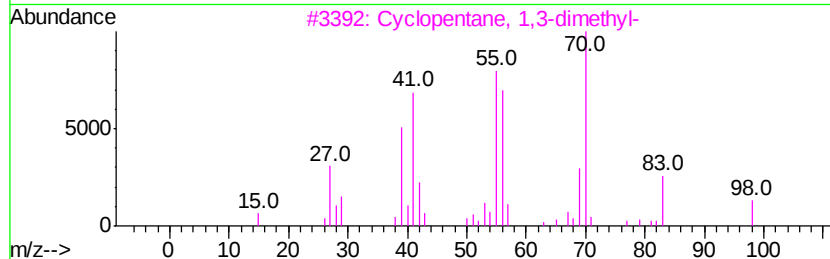
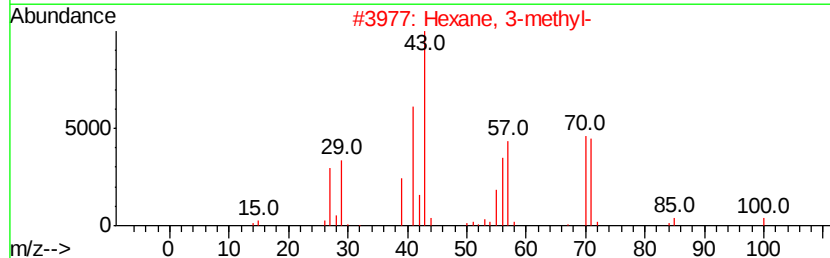
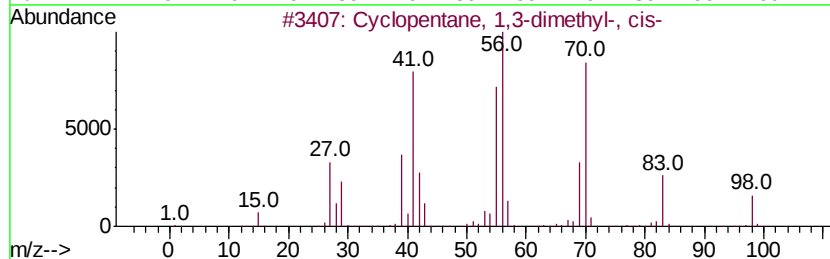
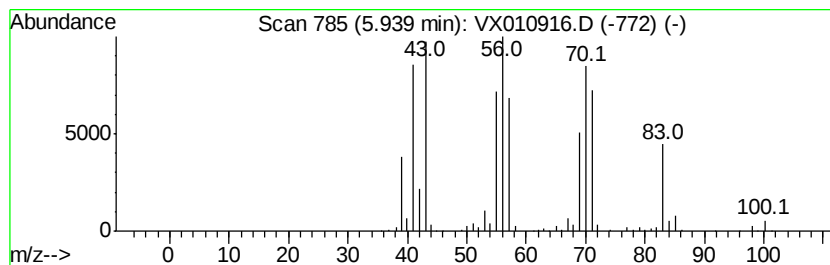
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Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, 1,3-dimethyl-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.94	80.10 ug/l	1019760	Pentafluorobenzene	5.66

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	50
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	49
3		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	47
4		1-Heptene, 4-methyl-	112	C8H16	013151-05-8	47
5		Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071619\
Data File : VX010916.D
Acq On : 16 Jul 2019 19:02
Operator : JC/SP
Sample : K3791-16
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0610

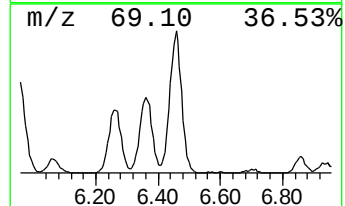
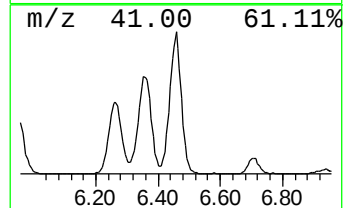
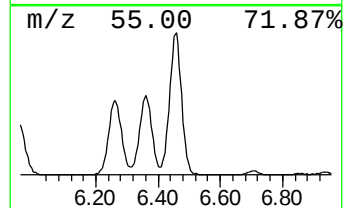
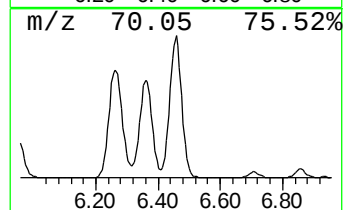
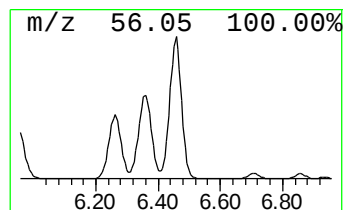
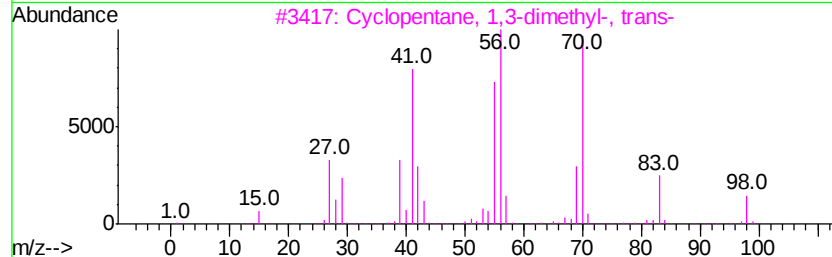
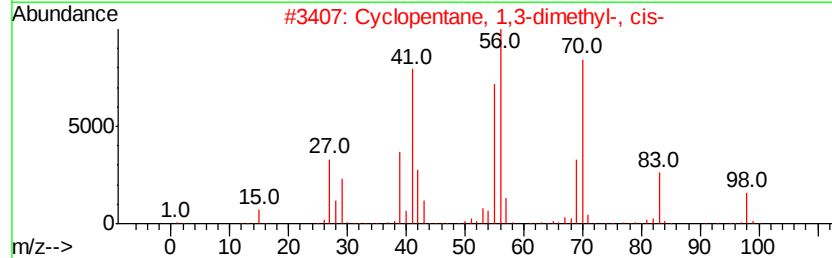
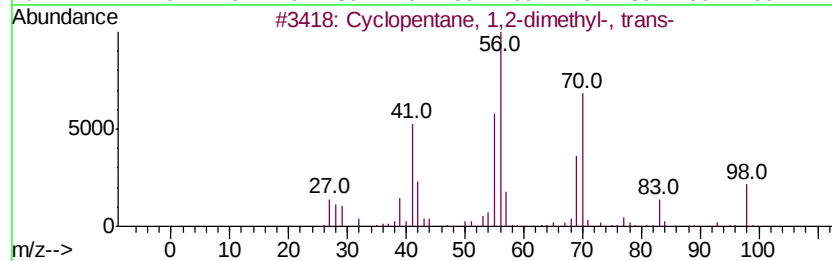
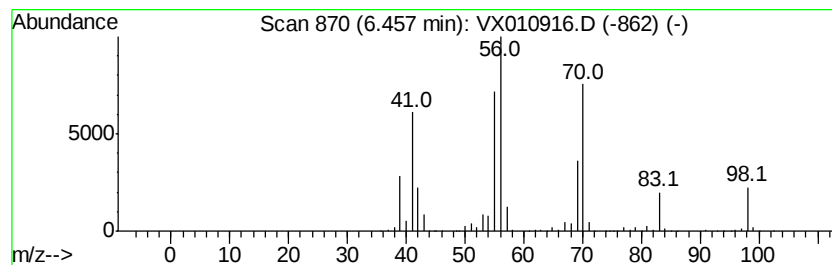
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, 1,2-dimethyl-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	81.90 ug/l	1331560	1,4-Difluorobenzene	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	90
2		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90
3		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	90
4		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	83
5		Isopropylcyclobutane	98	C7H14	000872-56-0	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071619\
Data File : VX010916.D
Acq On : 16 Jul 2019 19:02
Operator : JC/SP
Sample : K3791-16
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
FL-0610

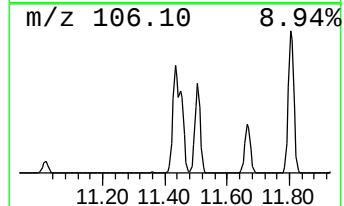
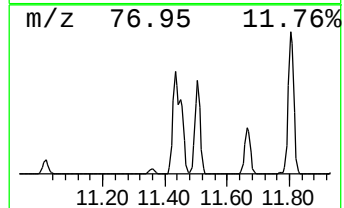
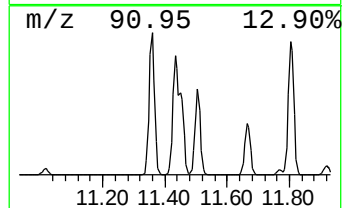
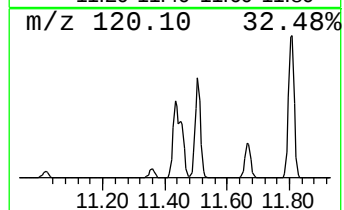
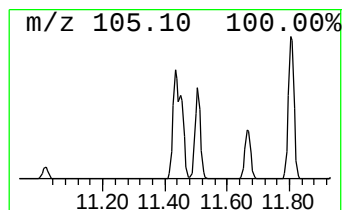
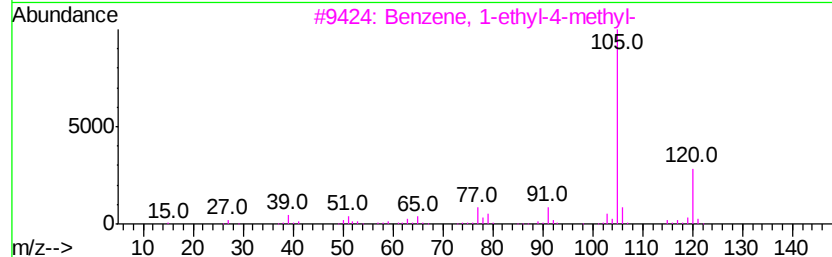
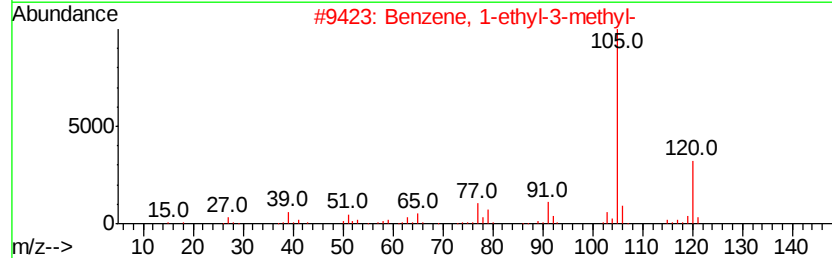
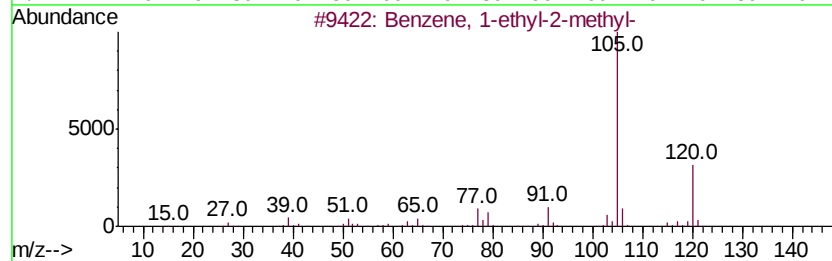
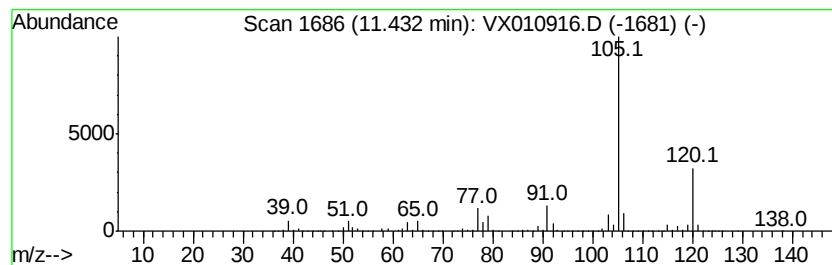
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	190.03 ug/l	5896050	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	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:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071619\
Data File : VX010916.D
Acq On : 16 Jul 2019 19:02
Operator : JC/SP
Sample : K3791-16
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
FL-0610

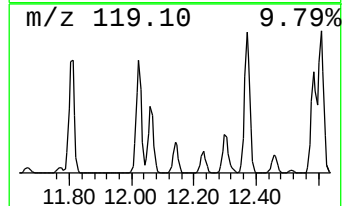
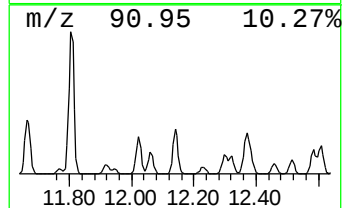
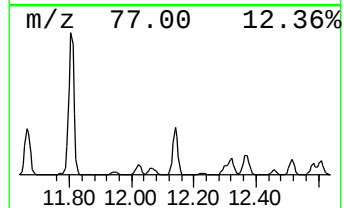
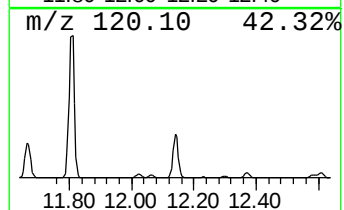
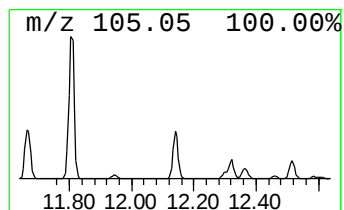
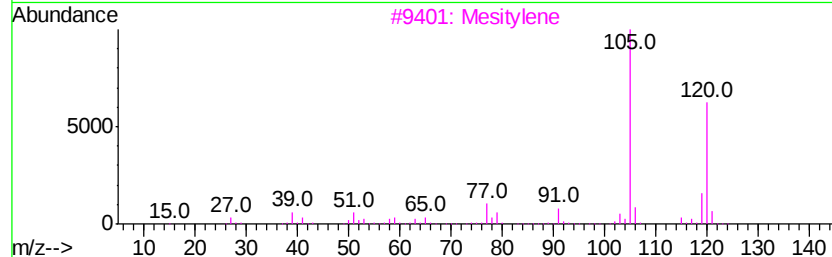
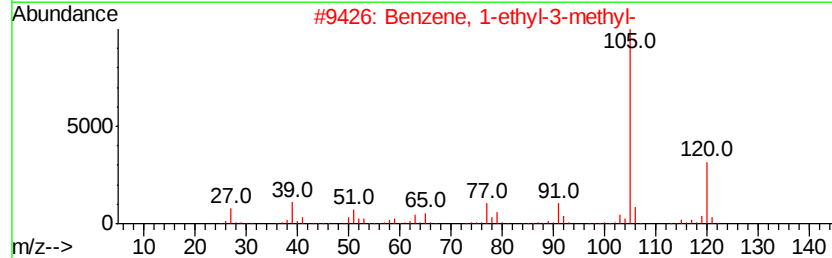
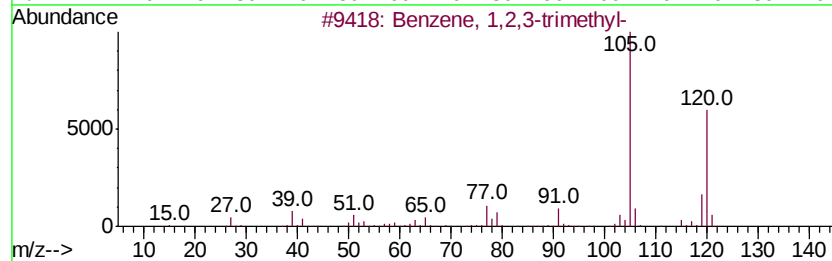
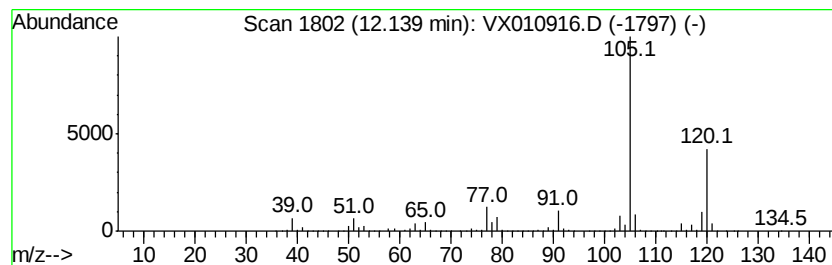
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	82.58 ug/l	2562170	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-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071619\
Data File : VX010916.D
Acq On : 16 Jul 2019 19:02
Operator : JC/SP
Sample : K3791-16
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0610

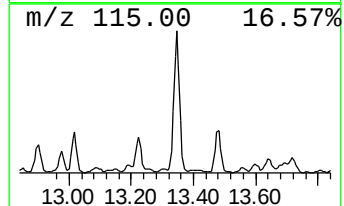
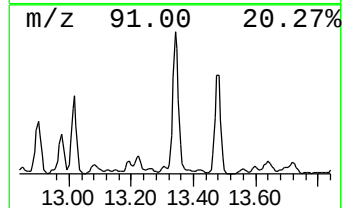
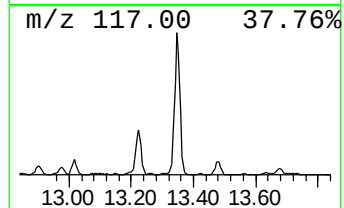
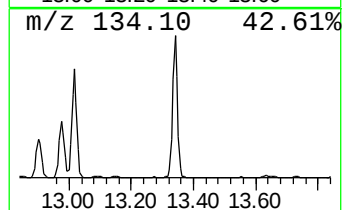
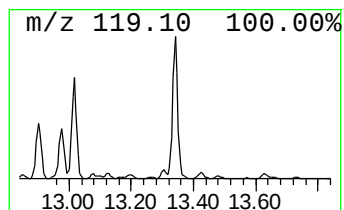
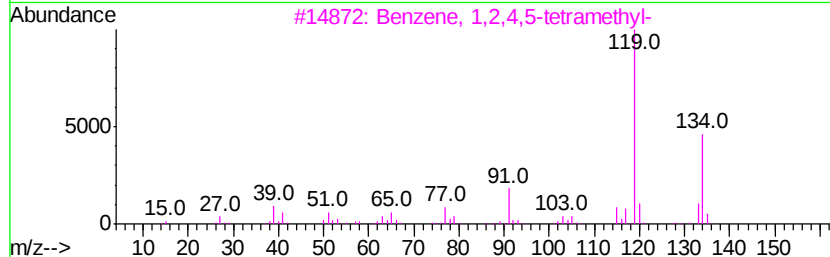
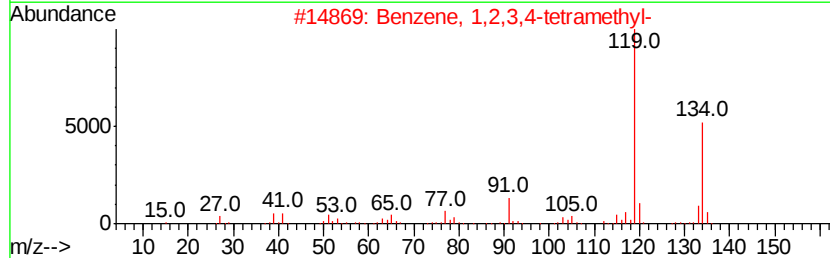
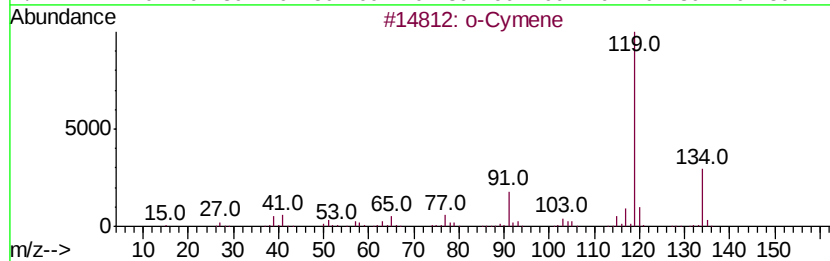
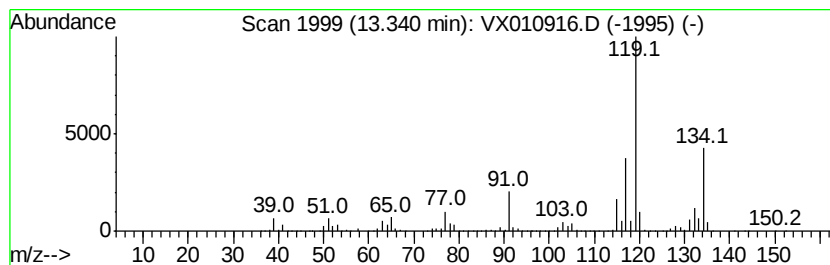
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Quant Title : SW846 8260

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

Peak Number 9 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	76.78 ug/l	2382210	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	81
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	76



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071619\
Data File : VX010916.D
Acq On : 16 Jul 2019 19:02
Operator : JC/SP
Sample : K3791-16
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0610

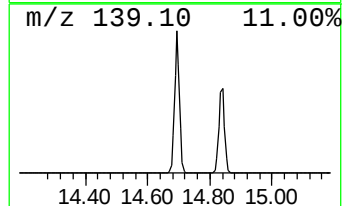
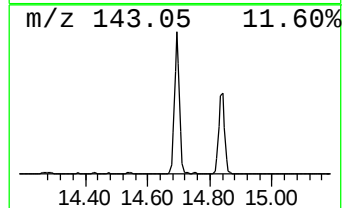
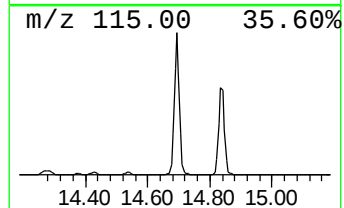
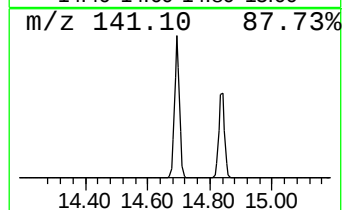
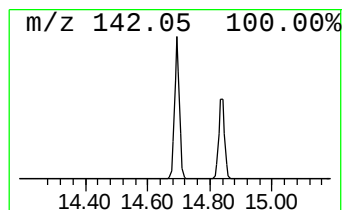
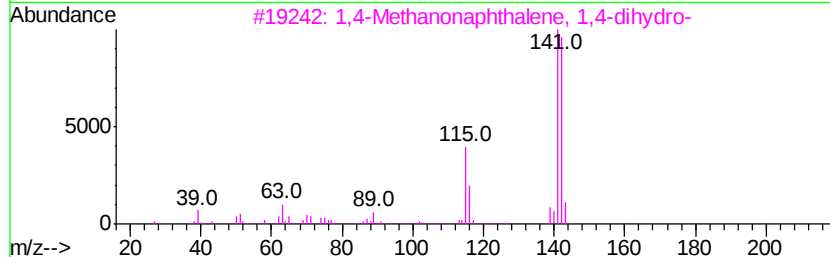
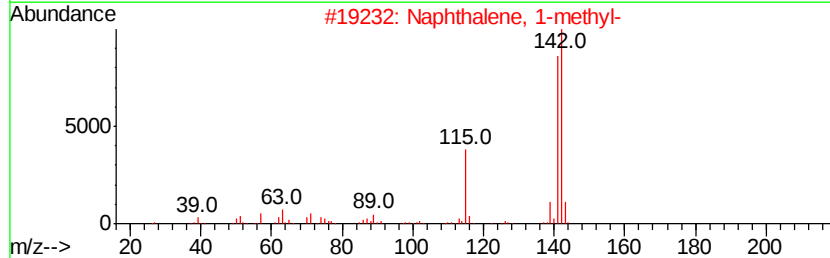
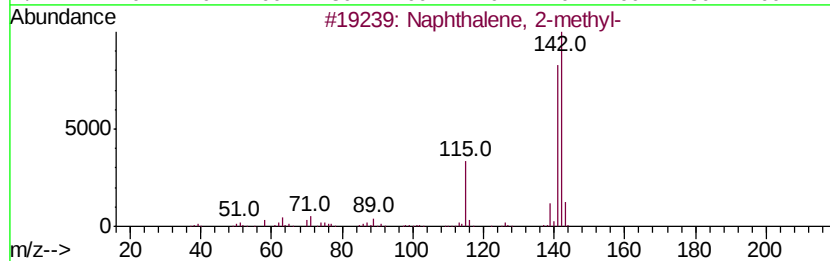
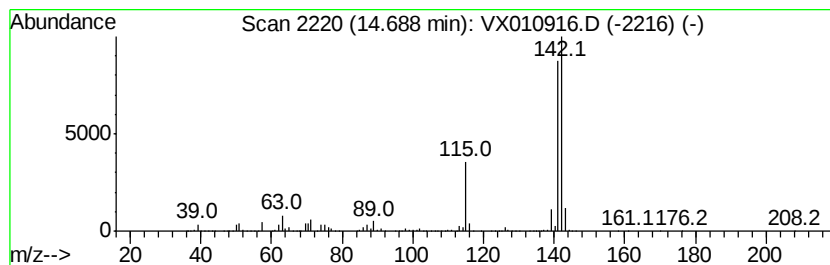
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	93.06 ug/l	2887300	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071619\
Data File : VX010916.D
Acq On : 16 Jul 2019 19:02
Operator : JC/SP
Sample : K3791-16
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0610

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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	74.6	ug/l	950221	1	5.66	636527	50.0
n-Hexane	3.52	75.2	ug/l	957155	1	5.66	636527	50.0
Cyclopentane, met...	4.40	106.8	ug/l	1359890	1	5.66	636527	50.0
Cyclopentane, 1,3...	5.94	80.1	ug/l	1019760	1	5.66	636527	50.0
Cyclopentane, 1,2...	6.46	81.9	ug/l	1331560	2	6.86	812915	50.0
Benzene, 1-ethyl-...	11.43	190.0	ug/l	5896050	4	12.08	1551370	50.0
Benzene, 1,2,3-tr...	12.14	82.6	ug/l	2562170	4	12.08	1551370	50.0
o-Cymene	13.34	76.8	ug/l	2382210	4	12.08	1551370	50.0
Naphthalene, 1-me...	14.69	93.1	ug/l	2887300	4	12.08	1551370	50.0