

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
 Data File : VX010703.D
 Acq On : 08 Jul 2019 18:09
 Operator : JC/SP
 Sample : K3664-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0579

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\82X061919W.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	1.788	100	104	113	rVB	157391	228766	2.93%	0.429%
2	2.001	133	139	149	rVB	120446	172297	2.20%	0.323%
3	2.398	196	204	209	rBV	50035	102361	1.31%	0.192%
4	2.843	269	277	281	rBV	84126	190393	2.44%	0.357%
5	2.891	281	285	289	rVV	151178	323632	4.14%	0.607%
6	2.934	289	292	302	rVB	125359	229198	2.93%	0.430%
7	3.172	321	331	343	rBV	305692	709804	9.08%	1.332%
8	3.525	380	389	403	rVB	57682	130576	1.67%	0.245%
9	4.409	523	534	548	rVB	518036	1527780	19.55%	2.866%
10	5.245	658	671	685	rVB2	22622	80117	1.03%	0.150%
11	5.501	700	713	718	rBV	126942	360651	4.61%	0.677%
12	5.580	718	726	734	rVV	553819	1737344	22.23%	3.259%
13	5.665	734	740	746	rVV	222987	642229	8.22%	1.205%
14	5.726	747	750	767	rVV	80542	207826	2.66%	0.390%
15	5.940	770	785	797	rBV4	88066	306472	3.92%	0.575%
16	6.062	797	805	812	rVV	124558	324842	4.16%	0.609%
17	6.141	813	818	827	rVV	44498	117154	1.50%	0.220%
18	6.263	828	838	847	rVV2	109280	360661	4.61%	0.677%
19	6.360	847	854	861	rVV	108826	298148	3.81%	0.559%
20	6.458	861	870	883	rVB	177910	492778	6.31%	0.924%
21	6.860	926	936	947	rBV	326016	767563	9.82%	1.440%
22	7.464	1023	1035	1047	rBV	1358212	3060425	39.16%	5.741%
23	7.732	1072	1079	1089	rVB	96897	188487	2.41%	0.354%
24	7.848	1089	1098	1106	rBV2	39999	81365	1.04%	0.153%
25	8.031	1121	1128	1139	rVB2	46585	104024	1.33%	0.195%
26	8.665	1224	1232	1235	rBV2	174263	307030	3.93%	0.576%
27	8.713	1235	1240	1248	rVB	733617	1187206	15.19%	2.227%
28	8.842	1254	1261	1265	rBV	91404	158106	2.02%	0.297%
29	8.909	1268	1272	1279	rVB	48415	86908	1.11%	0.163%
30	9.037	1288	1293	1300	rVB	180729	290823	3.72%	0.546%
31	9.165	1305	1314	1320	rBV	86446	141226	1.81%	0.265%
32	9.573	1375	1381	1384	rBV	53288	82368	1.05%	0.155%
33	9.622	1384	1389	1394	rVB	182432	252218	3.23%	0.473%
34	9.677	1394	1398	1402	rBV	69893	99619	1.27%	0.187%

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35	10.110	1464	1469	1475	rBV	685743	916879	11.73%	1.720%
36	10.183	1476	1481	1486	rVB	91947	135162	1.73%	0.254%
37	10.250	1486	1492	1501	rBV	1217955	1555756	19.91%	2.918%
38	10.360	1501	1510	1516	rVB	970089	1282069	16.40%	2.405%
39	10.835	1576	1588	1596	rBV2	51217	91531	1.17%	0.172%
40	11.018	1612	1618	1630	rVB	887768	1131077	14.47%	2.122%
41	11.134	1630	1637	1642	rBV	616687	779827	9.98%	1.463%
42	11.359	1669	1674	1681	rVB	1346291	1602536	20.51%	3.006%
43	11.451	1683	1689	1693	rVV	335874	412735	5.28%	0.774%
44	11.506	1693	1698	1706	rVB	79069	113808	1.46%	0.213%
45	11.664	1713	1724	1734	rBV	570486	743398	9.51%	1.395%
46	11.804	1742	1747	1753	rVB	6316063	7815317	100.00%	14.661%
47	11.920	1758	1766	1767	rBV	78263	110154	1.41%	0.207%
48	11.945	1767	1770	1780	rVB	229481	283924	3.63%	0.533%
49	12.079	1786	1792	1797	rVB2	778753	1094614	14.01%	2.053%
50	12.140	1797	1802	1808	rBV	3285498	3945916	50.49%	7.402%
51	12.231	1812	1817	1821	rVB	99658	122786	1.57%	0.230%
52	12.298	1821	1828	1834	rBV	750893	917275	11.74%	1.721%
53	12.365	1834	1839	1849	rVB2	348428	595640	7.62%	1.117%
54	12.457	1849	1854	1859	rBV	242483	289504	3.70%	0.543%
55	12.518	1859	1864	1870	rBV	605678	735052	9.41%	1.379%
56	12.585	1870	1875	1877	rBV	605090	764149	9.78%	1.433%
57	12.609	1877	1879	1883	rVV	553207	569577	7.29%	1.068%
58	12.664	1883	1888	1892	rVV	745929	918097	11.75%	1.722%
59	12.701	1892	1894	1898	rVV	197056	211000	2.70%	0.396%
60	12.755	1898	1903	1910	rVB2	548002	868072	11.11%	1.628%
61	12.902	1921	1927	1932	rVB2	504590	656547	8.40%	1.232%
62	12.975	1932	1939	1942	rBV	493357	616855	7.89%	1.157%
63	13.018	1942	1946	1952	rVB	820389	1003430	12.84%	1.882%
64	13.079	1952	1956	1962	rVB3	87540	120856	1.55%	0.227%
65	13.194	1971	1975	1977	rBV	92776	108827	1.39%	0.204%
66	13.225	1977	1980	1983	rVB	167308	176789	2.26%	0.332%
67	13.304	1989	1993	1995	rBV	97317	120412	1.54%	0.226%
68	13.341	1995	1999	2005	rVB2	1384667	1900370	24.32%	3.565%
69	13.475	2017	2021	2027	rVB	614603	788063	10.08%	1.478%
70	13.560	2027	2035	2038	rBV3	61372	98403	1.26%	0.185%
71	13.597	2038	2041	2044	rVV	86486	124056	1.59%	0.233%

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72	13.639	2044	2048	2053	rVV3	169799	341444	4.37%	0.641%
73	13.694	2053	2057	2059	rVV3	92563	157551	2.02%	0.296%
74	13.719	2059	2061	2069	rVB2	152377	216995	2.78%	0.407%
75	13.835	2069	2080	2088	rBV	567143	864014	11.06%	1.621%
76	13.908	2088	2092	2097	rVB	72030	91958	1.18%	0.173%
77	13.981	2097	2104	2108	rBV2	115860	182827	2.34%	0.343%
78	14.286	2146	2154	2159	rVB3	160638	297259	3.80%	0.558%
79	14.536	2190	2195	2208	rBV2	158455	214250	2.74%	0.402%
80	14.694	2213	2221	2225	rBV	725893	887975	11.36%	1.666%
81	14.834	2239	2244	2250	rBV	501394	656241	8.40%	1.231%
82	15.548	2354	2361	2366	rBV	67039	108366	1.39%	0.203%
83	15.688	2375	2384	2387	rBV	77466	136769	1.75%	0.257%
84	15.724	2387	2390	2397	rVB	54617	80453	1.03%	0.151%

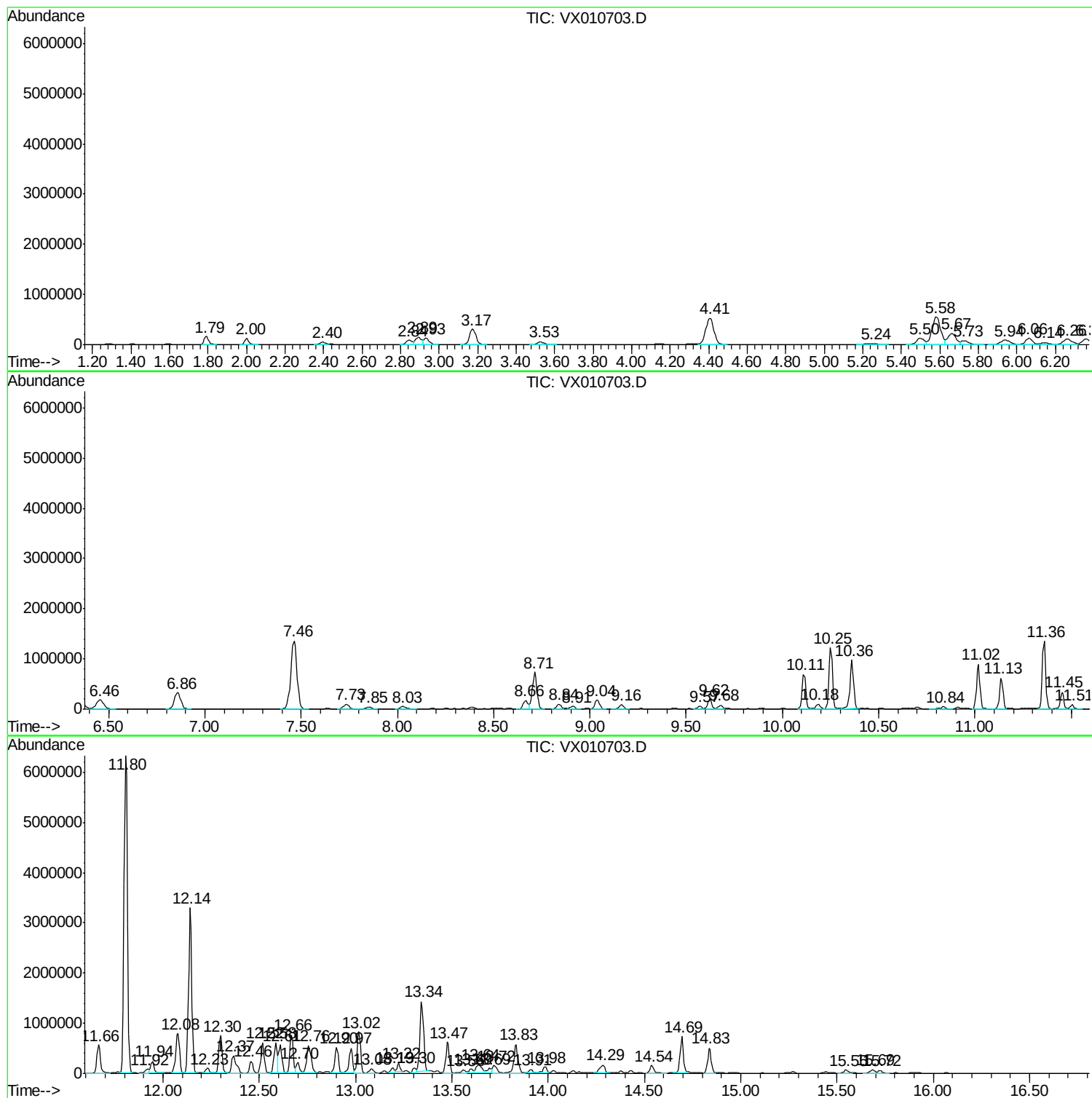
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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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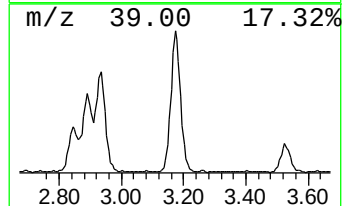
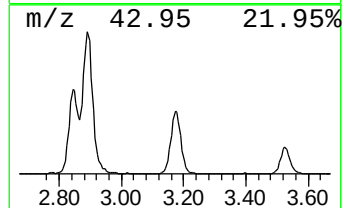
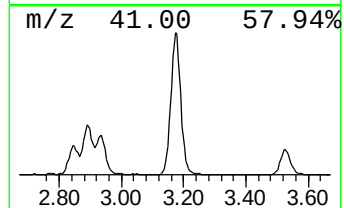
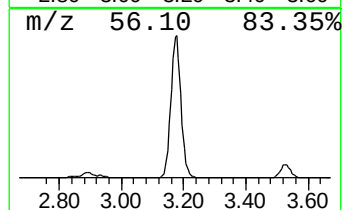
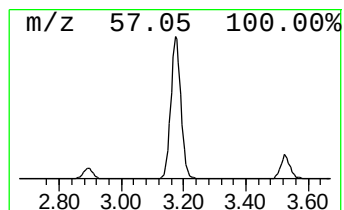
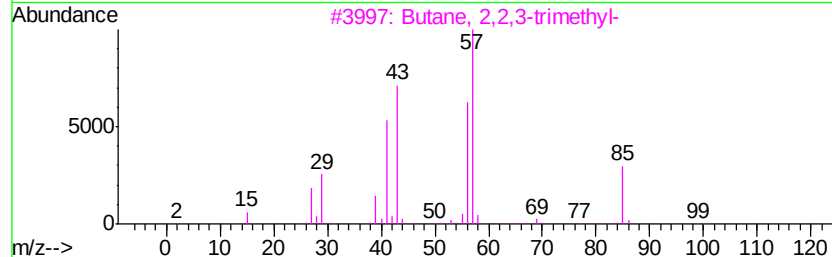
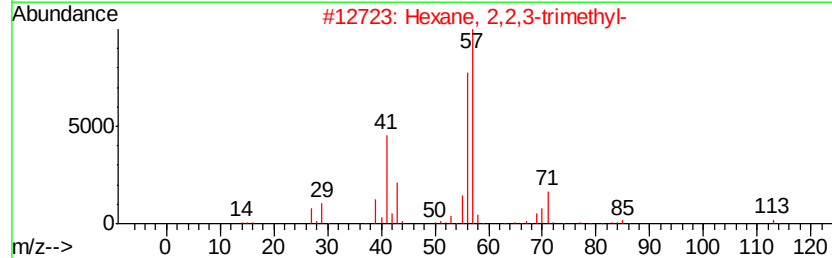
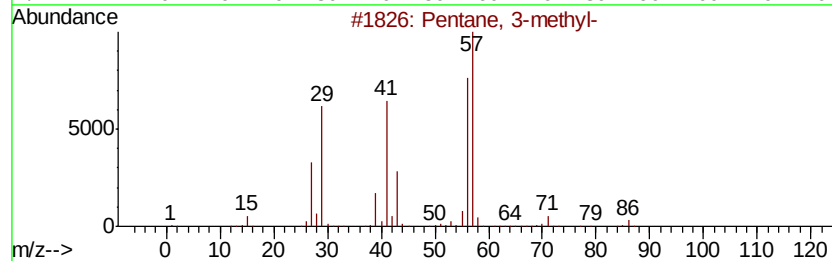
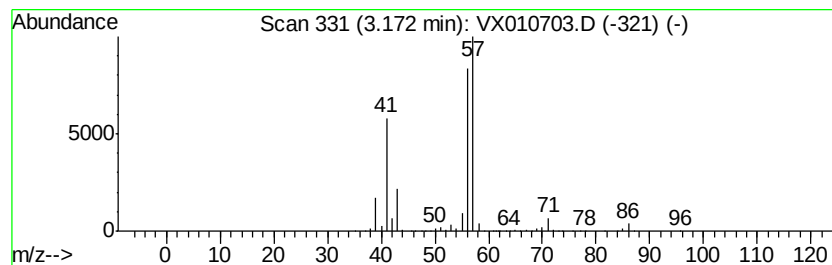
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	55.26 ug/l	709804	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
4		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
5		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	42



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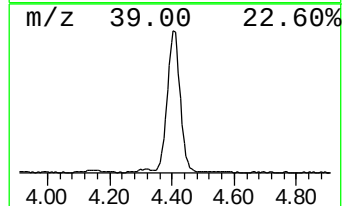
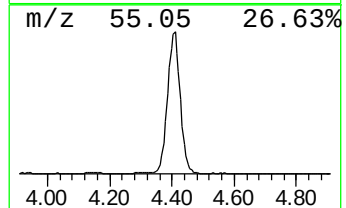
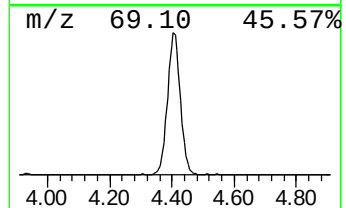
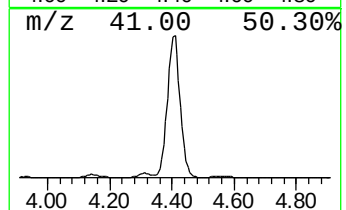
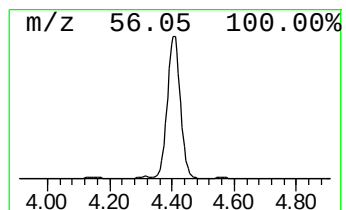
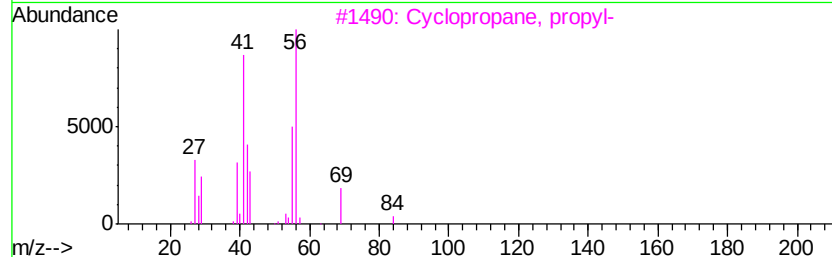
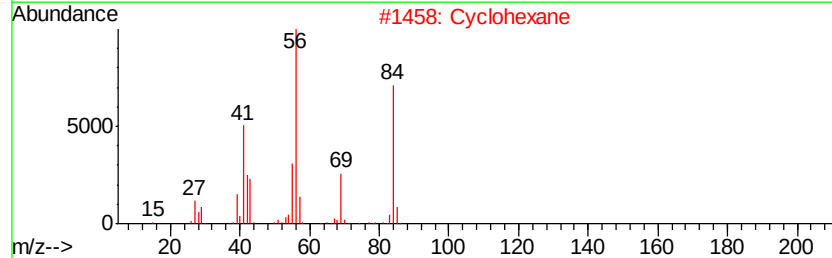
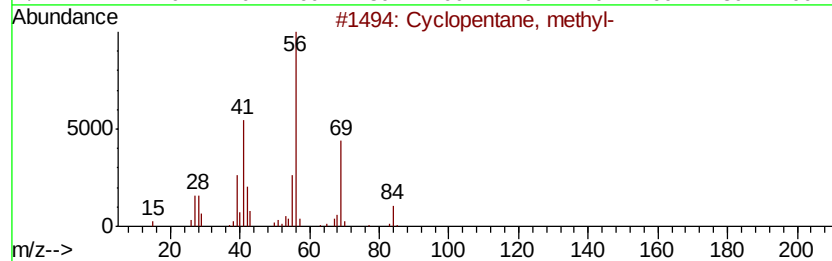
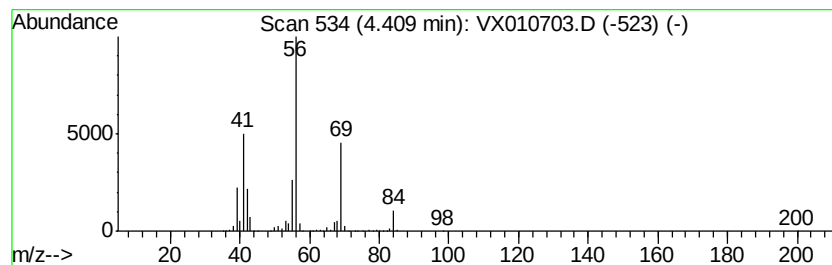
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	118.94 ug/l	1527780	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclohexane	84	C6H12	000110-82-7	78
3		Cyclopropane, propyl-	84	C6H12	002415-72-7	78
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64



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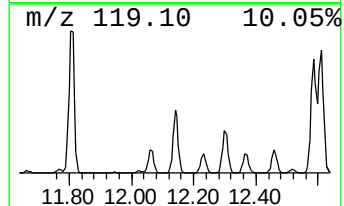
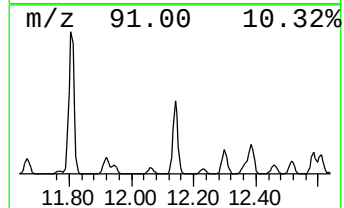
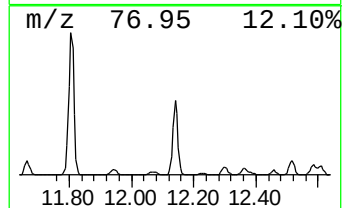
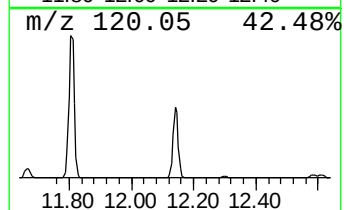
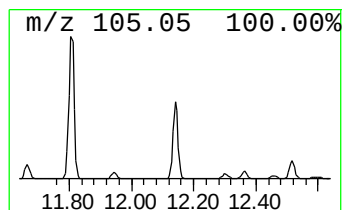
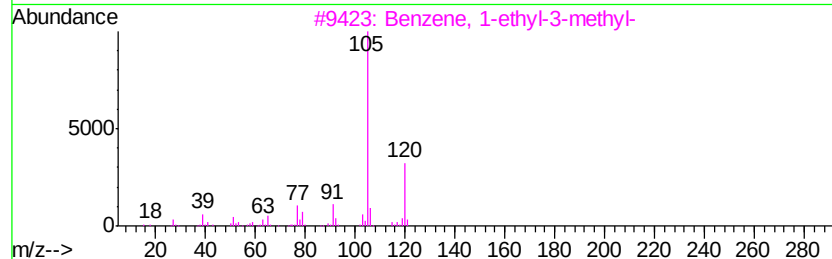
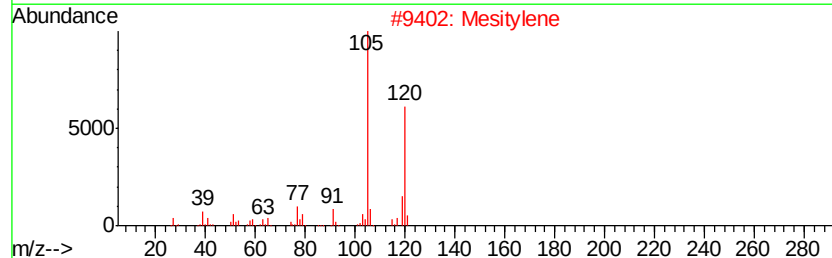
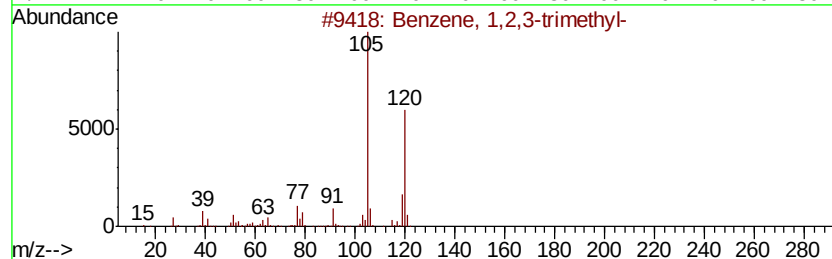
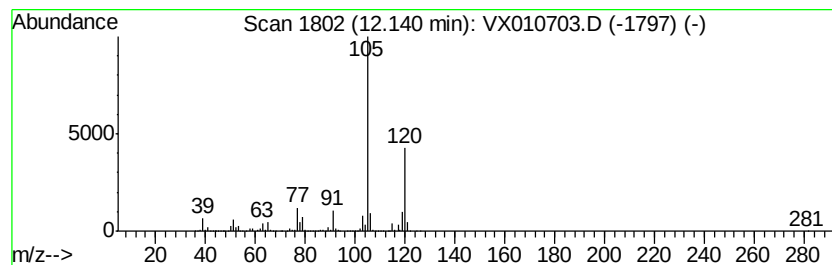
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	180.24 ug/l	3945920	1,4-Dichlorobenzene-d4	12.08

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	91
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-4-methyl-	120	C9H12	000622-96-8	90



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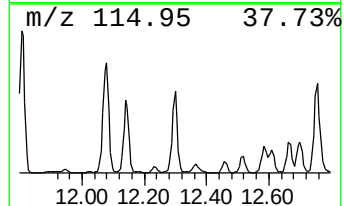
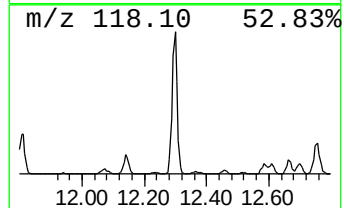
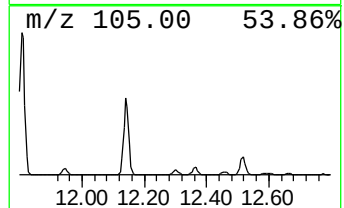
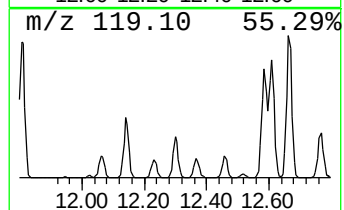
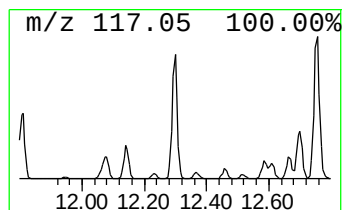
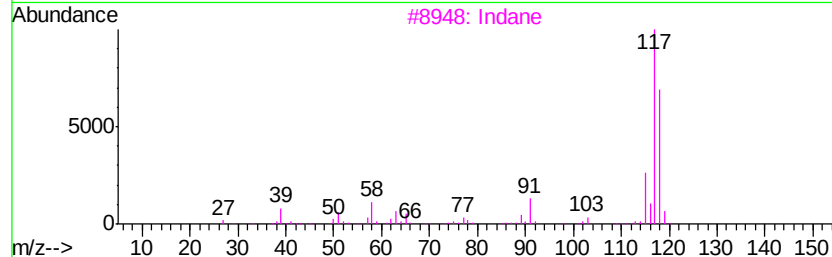
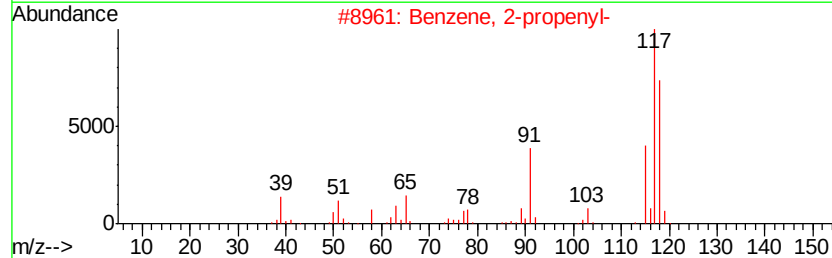
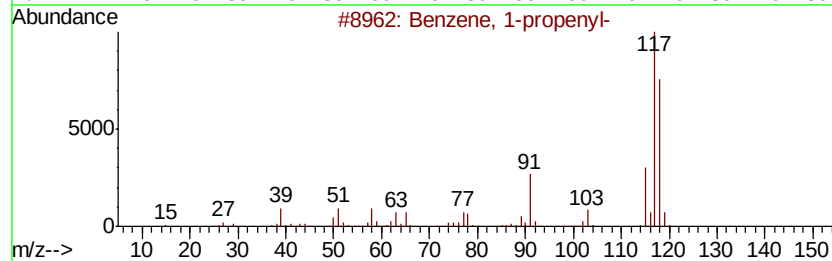
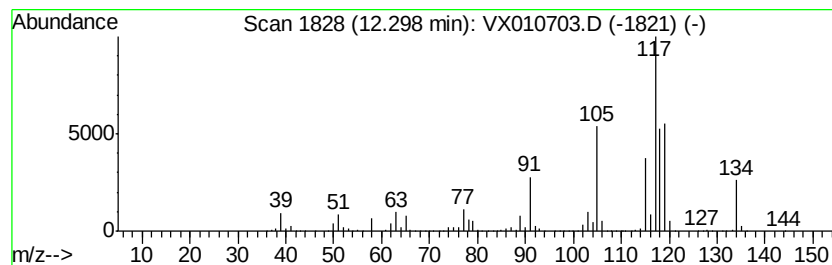
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-propenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	41.90 ug/l	917275	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	90
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
3		Indane	118	C9H10	000496-11-7	60
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	60
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

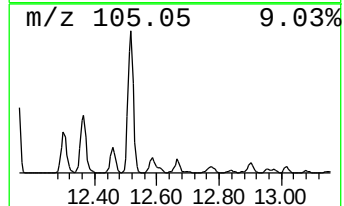
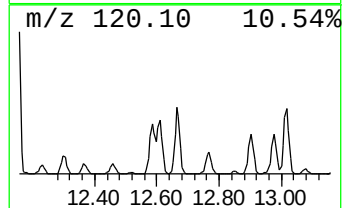
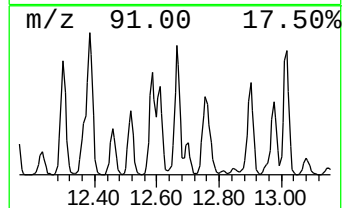
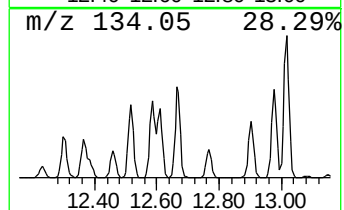
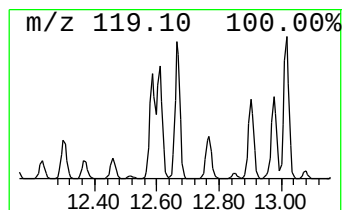
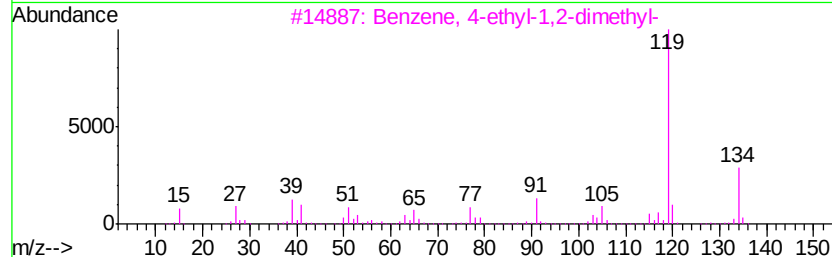
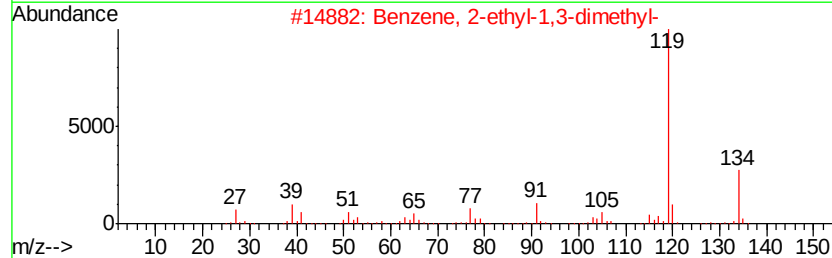
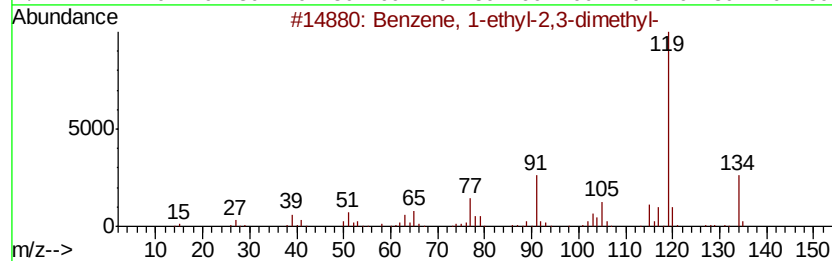
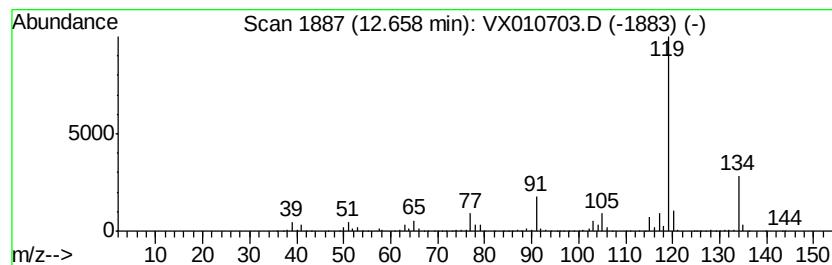
Instrument :
MSVOA_X
ClientSampleId :
FL-0579

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.66	41.94 ug/l	918097	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5	o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0579

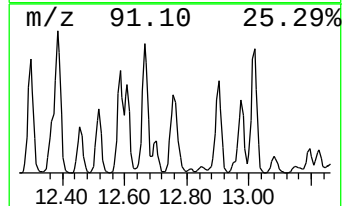
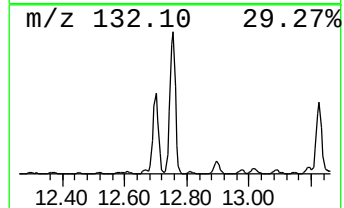
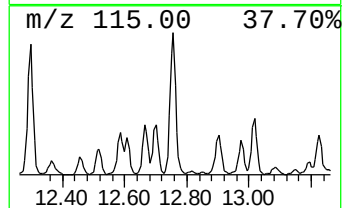
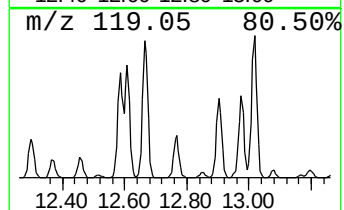
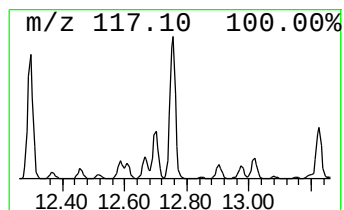
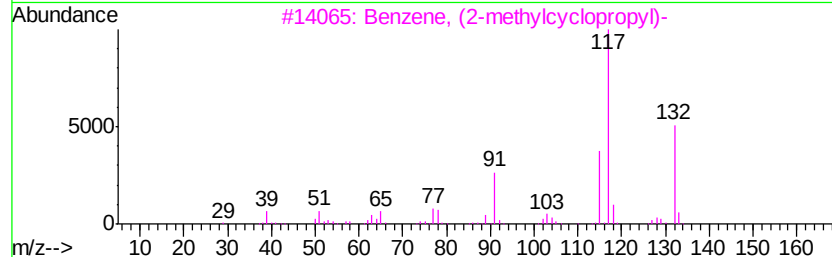
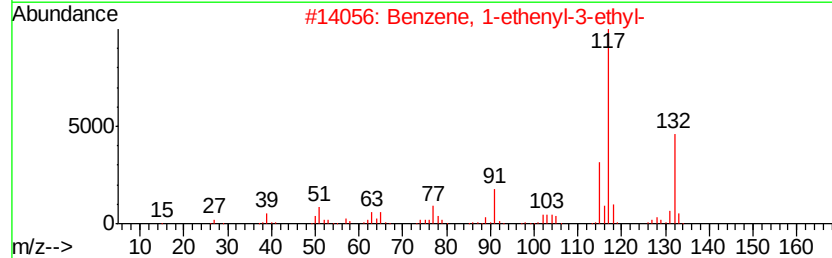
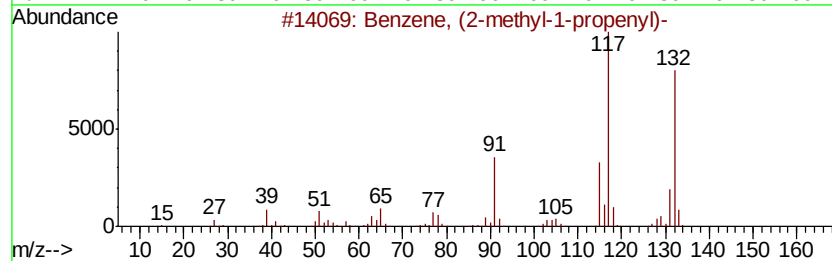
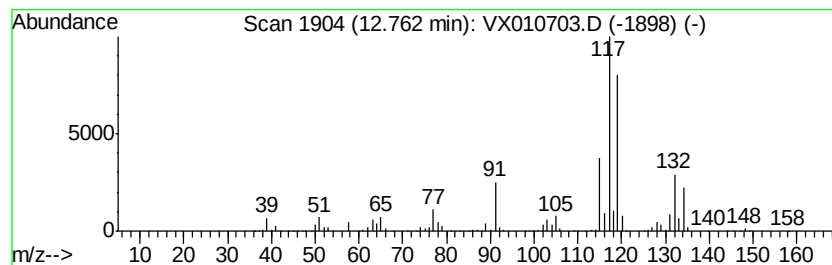
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	39.65 ug/l	868072	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		Indan, 1-methyl-	132	C10H12	000767-58-8	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

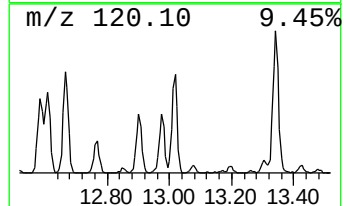
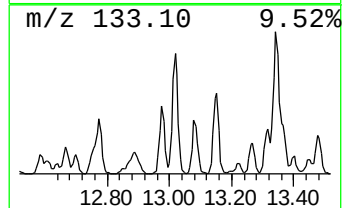
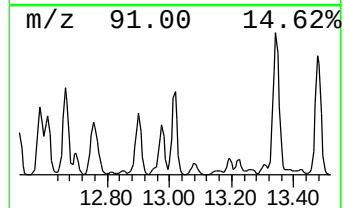
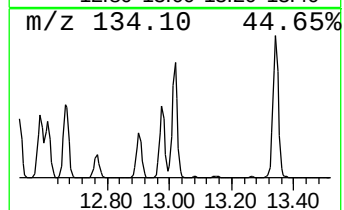
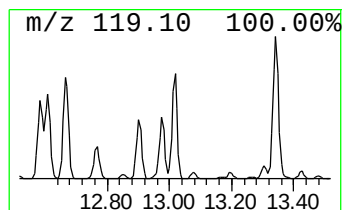
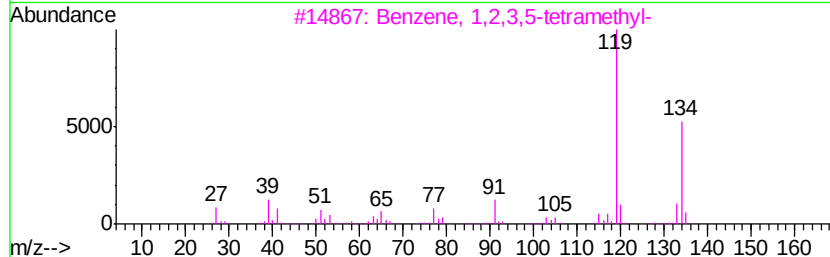
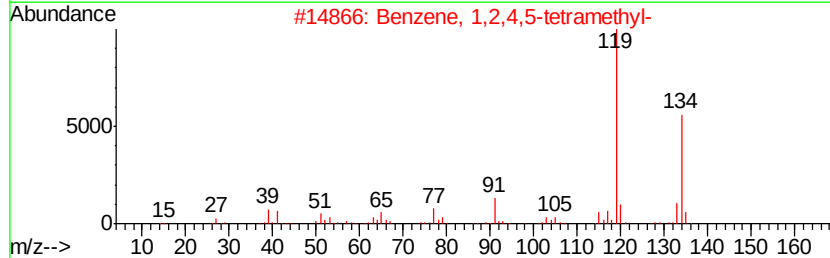
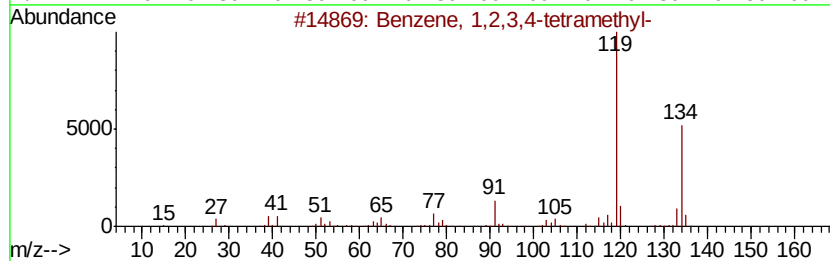
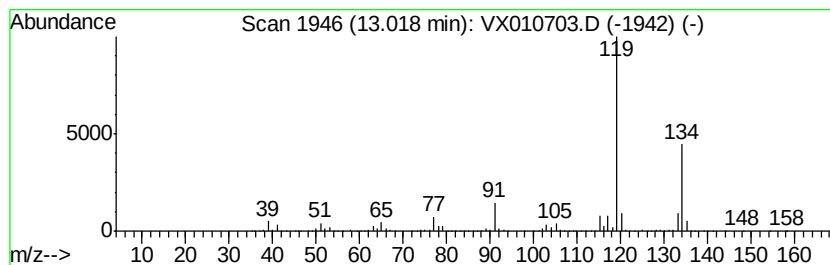
Instrument :
MSVOA_X
ClientSampleId :
FL-0579

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	45.83 ug/l	1003430	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 97
2		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 97
3		Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7 97
4		o-Cymene	134 C10H14	000527-84-4 96
5		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0579

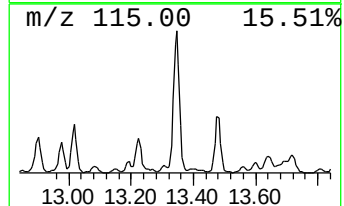
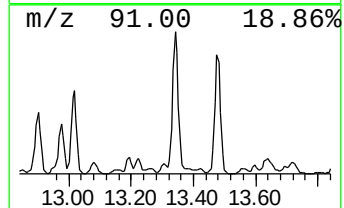
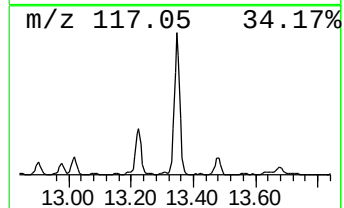
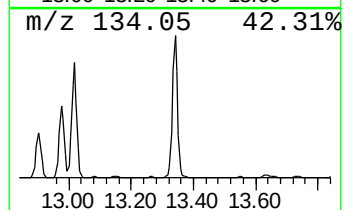
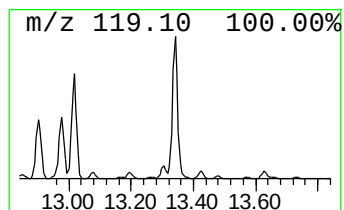
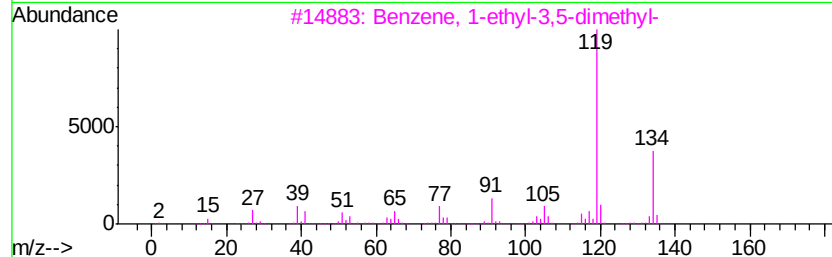
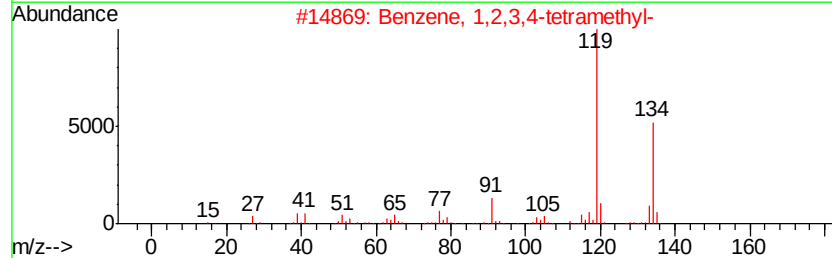
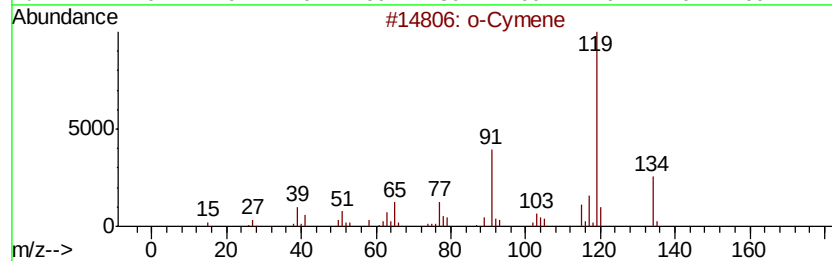
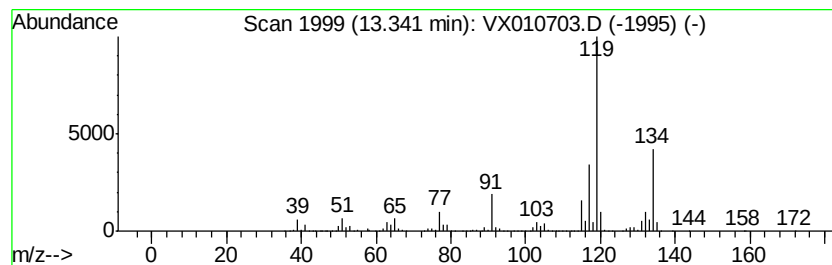
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Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	86.81 ug/l	1900370	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-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	81
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	76
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0579

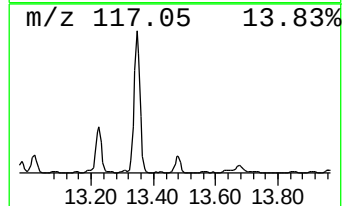
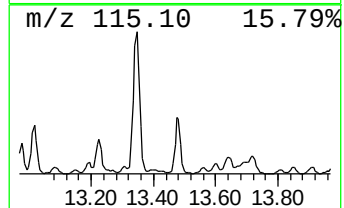
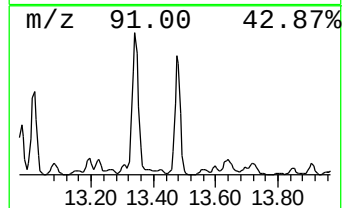
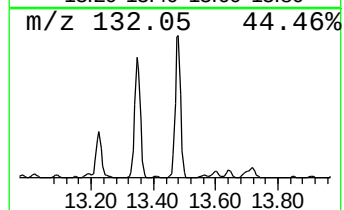
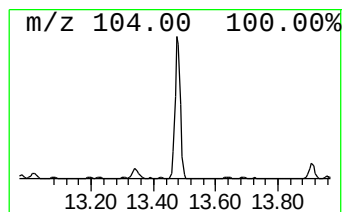
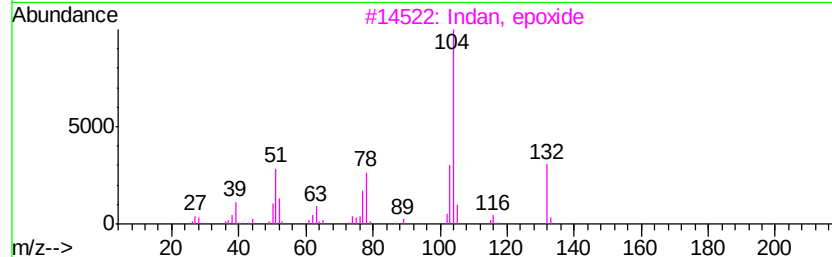
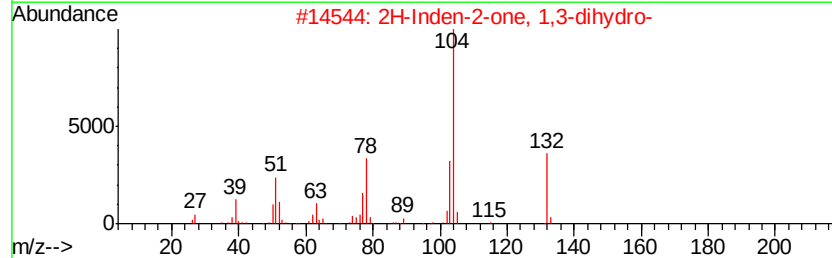
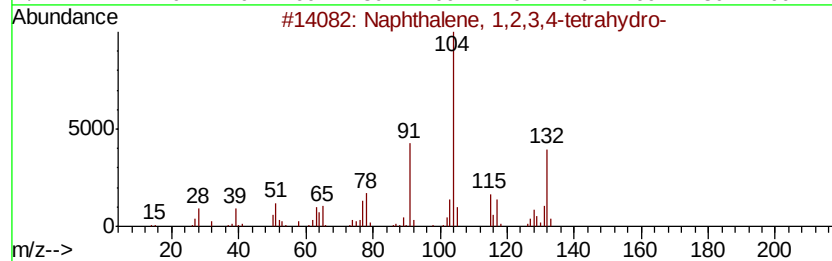
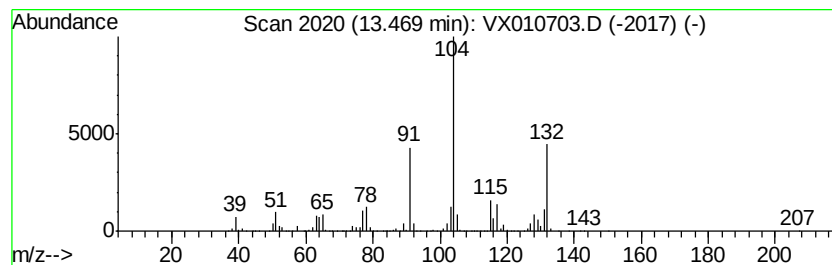
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	36.00 ug/l	788063	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
3		Indan, epoxide	132	C9H8O	000768-22-9	52
4		Benzene, cyclobutyl-	132	C10H12	004392-30-7	45
5		3-Phenylthiolane,S,S-dioxide	196	C10H12O2S	016766-64-6	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0579

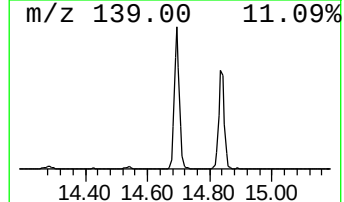
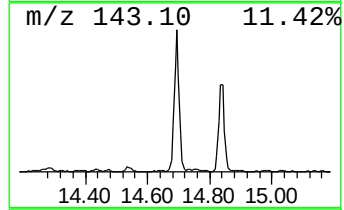
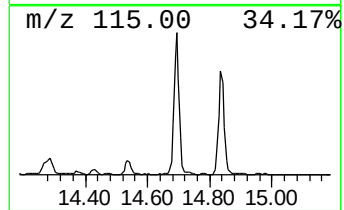
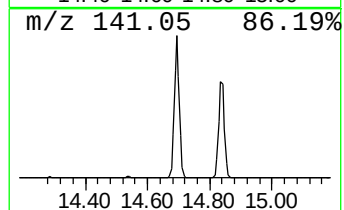
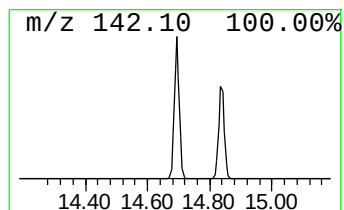
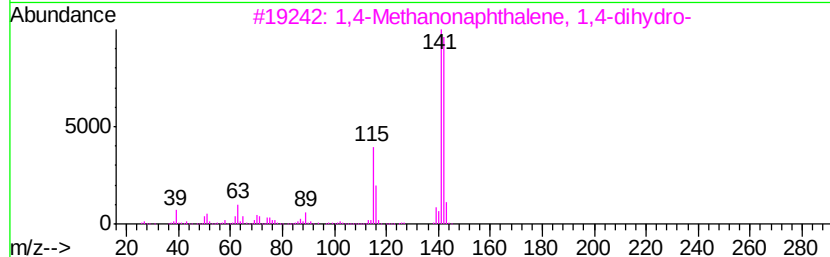
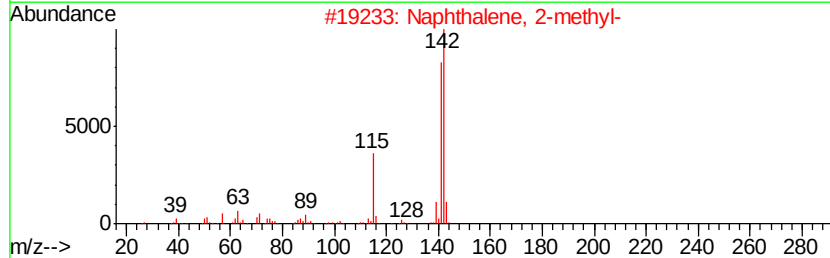
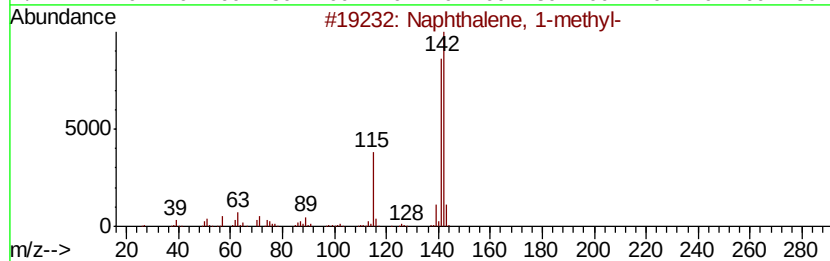
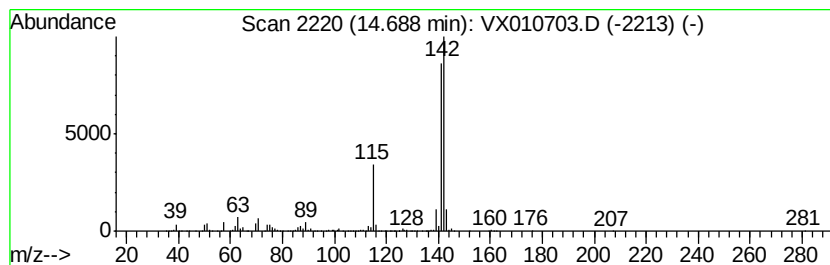
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	40.56 ug/l	887975	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		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010703.D
Acq On : 08 Jul 2019 18:09
Operator : JC/SP
Sample : K3664-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0579

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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, 3-methyl-	3.17	55.3	ug/l	709804	1	5.67	642229	50.0
Cyclopentane, met...	4.41	118.9	ug/l	1527780	1	5.67	642229	50.0
Benzene, 1,2,3-tr...	12.14	180.2	ug/l	3945920	4	12.08	1094610	50.0
Benzene, 1-propenyl-	12.30	41.9	ug/l	917275	4	12.08	1094610	50.0
Benzene, 1-ethyl-...	12.66	41.9	ug/l	918097	4	12.08	1094610	50.0
Benzene, (2-methy...	12.76	39.6	ug/l	868072	4	12.08	1094610	50.0
Benzene, 1,2,3,4-...	13.02	45.8	ug/l	1003430	4	12.08	1094610	50.0
o-Cymene	13.34	86.8	ug/l	1900370	4	12.08	1094610	50.0
Naphthalene, 1,2,...	13.47	36.0	ug/l	788063	4	12.08	1094610	50.0
Naphthalene, 1-me...	14.69	40.6	ug/l	887975	4	12.08	1094610	50.0