

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071219\
 Data File : VX010840.D
 Acq On : 12 Jul 2019 16:49
 Operator : JC/SP
 Sample : K3786-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0589

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	2.001	134	139	146	rVB	42771	62356	1.17%	0.147%
2	2.891	280	285	300	rVB	58459	145616	2.72%	0.342%
3	3.172	322	331	343	rBV	72702	166071	3.10%	0.390%
4	3.525	380	389	398	rBV	64983	148744	2.78%	0.350%
5	4.403	523	533	548	rVB	262011	771823	14.43%	1.815%
6	5.501	699	713	718	rBV	121334	339506	6.35%	0.798%
7	5.580	718	726	734	rVV	293187	976834	18.26%	2.297%
8	5.659	734	739	747	rVV	208812	615123	11.50%	1.446%
9	5.732	747	751	764	rVB3	62482	158727	2.97%	0.373%
10	5.933	774	784	796	rBV4	73392	249363	4.66%	0.586%
11	6.061	796	805	819	rVB	117898	305656	5.71%	0.719%
12	6.263	827	838	847	rBV3	67384	232344	4.34%	0.546%
13	6.360	847	854	861	rVV3	71343	195647	3.66%	0.460%
14	6.458	861	870	885	rVB	103055	289989	5.42%	0.682%
15	6.860	925	936	946	rBV	309530	731743	13.68%	1.720%
16	7.464	1021	1035	1047	rBV	1115131	2553624	47.73%	6.004%
17	7.732	1072	1079	1089	rVB	117060	233254	4.36%	0.548%
18	7.848	1091	1098	1106	rVB	96316	194085	3.63%	0.456%
19	8.031	1120	1128	1140	rVB3	110869	258854	4.84%	0.609%
20	8.390	1182	1187	1193	rVB2	72455	129316	2.42%	0.304%
21	8.579	1214	1218	1225	rVB3	30292	59295	1.11%	0.139%
22	8.665	1225	1232	1235	rBV2	215564	382785	7.15%	0.900%
23	8.713	1235	1240	1253	rVV	695284	1169172	21.85%	2.749%
24	8.841	1253	1261	1265	rVV	201694	350453	6.55%	0.824%
25	8.908	1269	1272	1279	rVB	58928	91240	1.71%	0.215%
26	9.037	1288	1293	1305	rVB	433906	700092	13.09%	1.646%
27	9.165	1305	1314	1319	rBV	166879	266604	4.98%	0.627%
28	9.573	1374	1381	1385	rBV3	126934	212844	3.98%	0.500%
29	9.622	1385	1389	1393	rVB	206109	274127	5.12%	0.645%
30	9.677	1393	1398	1403	rBV	432655	625620	11.69%	1.471%
31	9.890	1426	1433	1447	rVB2	57526	107652	2.01%	0.253%
32	10.110	1464	1469	1475	rBV	649441	871633	16.29%	2.049%
33	10.183	1475	1481	1486	rVV	113355	169396	3.17%	0.398%
34	10.250	1486	1492	1498	rBV	1409666	1829409	34.19%	4.301%

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35	10.359	1501	1510	1515	rBV	1886702	2423461	45.30%	5.698%
36	10.512	1530	1535	1543	rVB	44196	62011	1.16%	0.146%
37	10.646	1546	1557	1564	rBV4	70139	194510	3.64%	0.457%
38	10.835	1585	1588	1596	rVB	53535	69172	1.29%	0.163%
39	10.914	1596	1601	1607	rBV3	37434	63455	1.19%	0.149%
40	11.018	1613	1618	1630	rVB	772279	991013	18.52%	2.330%
41	11.134	1632	1637	1642	rBV	567292	730028	13.64%	1.716%
42	11.274	1657	1660	1665	rVB2	41790	58733	1.10%	0.138%
43	11.359	1668	1674	1681	rVV	908479	1117840	20.89%	2.628%
44	11.451	1684	1689	1694	rVV	1288268	1558458	29.13%	3.664%
45	11.506	1694	1698	1706	rVB	78974	103140	1.93%	0.242%
46	11.664	1719	1724	1733	rBV	1462196	1790790	33.47%	4.210%
47	11.804	1742	1747	1753	rVB	4524029	5350246	100.00%	12.579%
48	11.920	1758	1766	1767	rBV	67476	95870	1.79%	0.225%
49	11.945	1767	1770	1778	rVB	151422	190158	3.55%	0.447%
50	12.073	1785	1791	1797	rBV2	753569	1102062	20.60%	2.591%
51	12.140	1797	1802	1808	rVV	1831273	2190492	40.94%	5.150%
52	12.231	1812	1817	1821	rBV	46426	56571	1.06%	0.133%
53	12.298	1821	1828	1833	rBV	596368	726751	13.58%	1.709%
54	12.365	1833	1839	1849	rVB2	220827	366014	6.84%	0.861%
55	12.457	1849	1854	1859	rBV	95457	121151	2.26%	0.285%
56	12.518	1859	1864	1869	rBV	221362	269285	5.03%	0.633%
57	12.585	1870	1875	1877	rBV	247428	325718	6.09%	0.766%
58	12.609	1877	1879	1883	rVV	295461	304689	5.69%	0.716%
59	12.664	1883	1888	1891	rVV	328133	403817	7.55%	0.949%
60	12.701	1891	1894	1898	rVV	125339	156285	2.92%	0.367%
61	12.755	1898	1903	1910	rVB2	281854	462520	8.64%	1.087%
62	12.902	1921	1927	1932	rVB2	179665	251396	4.70%	0.591%
63	12.975	1932	1939	1942	rBV	213453	293487	5.49%	0.690%
64	13.018	1942	1946	1951	rVB	385771	465492	8.70%	1.094%
65	13.078	1951	1956	1962	rVB3	72073	100312	1.87%	0.236%
66	13.152	1963	1968	1971	rBV2	37771	60201	1.13%	0.142%
67	13.194	1971	1975	1977	rBV2	66424	77951	1.46%	0.183%
68	13.225	1977	1980	1983	rVB	110246	111011	2.07%	0.261%
69	13.304	1989	1993	1995	rBV2	74820	91491	1.71%	0.215%
70	13.341	1995	1999	2005	rVB2	645196	910612	17.02%	2.141%
71	13.475	2017	2021	2026	rVB	188683	249829	4.67%	0.587%

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72	13.560	2030	2035	2038	rBV3	43618	65916	1.23%	0.155%
73	13.597	2038	2041	2044	rBV	53875	68979	1.29%	0.162%
74	13.639	2044	2048	2053	rBV3	107701	182977	3.42%	0.430%
75	13.719	2059	2061	2070	rVB2	110180	162081	3.03%	0.381%
76	13.834	2070	2080	2088	rBV	497424	755427	14.12%	1.776%
77	13.987	2097	2105	2108	rBV3	65266	113818	2.13%	0.268%
78	14.127	2124	2128	2132	rBV3	39649	58621	1.10%	0.138%
79	14.267	2146	2151	2160	rVB2	75135	149241	2.79%	0.351%
80	14.426	2173	2177	2181	rVB	43030	56845	1.06%	0.134%
81	14.536	2190	2195	2204	rBV2	52739	86844	1.62%	0.204%
82	14.694	2213	2221	2225	rBV	409920	511983	9.57%	1.204%
83	14.834	2239	2244	2253	rBV	281554	390166	7.29%	0.917%
84	15.548	2353	2361	2365	rBV	49309	82320	1.54%	0.194%
85	15.682	2374	2383	2387	rBV	46915	82607	1.54%	0.194%

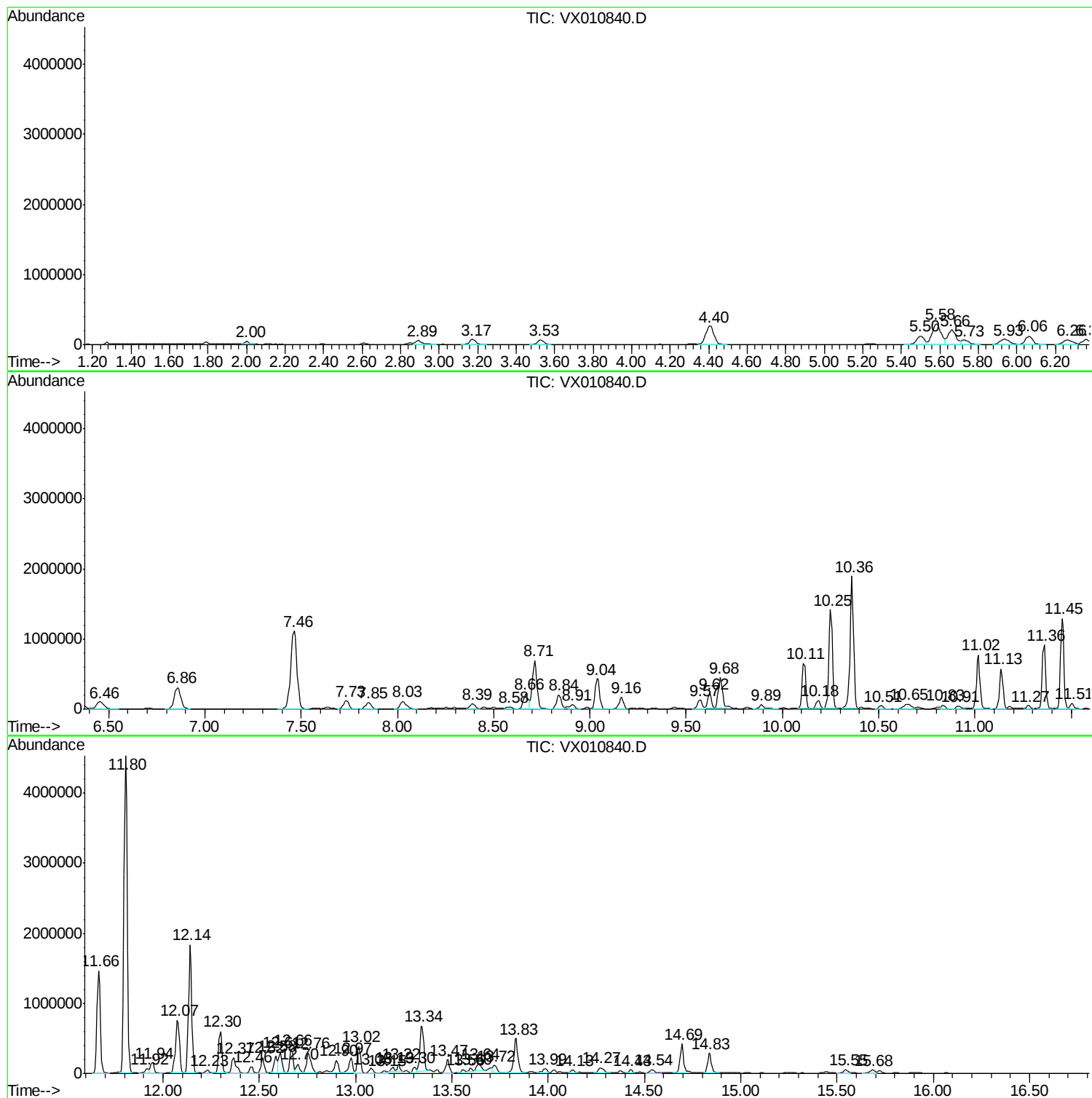
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



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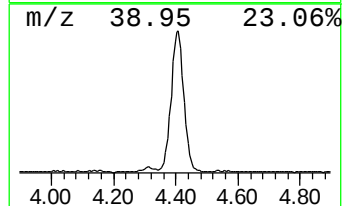
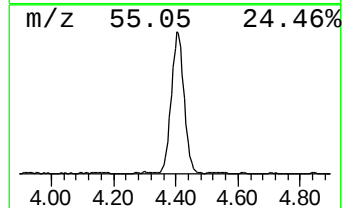
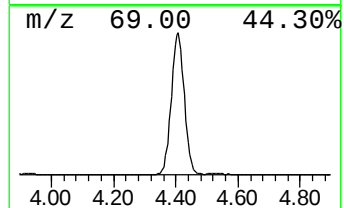
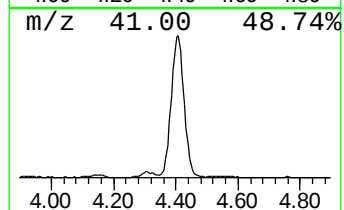
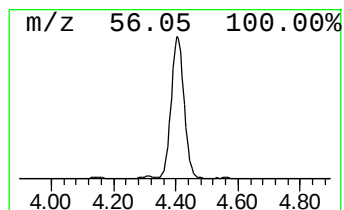
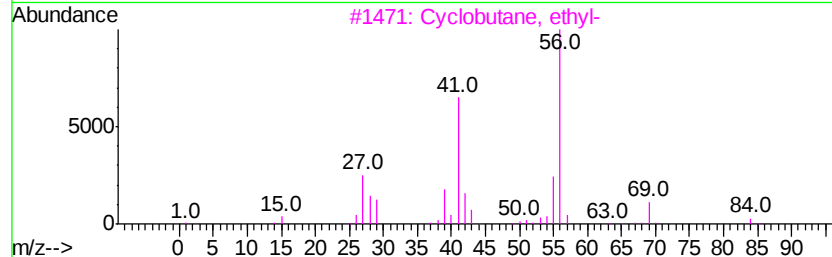
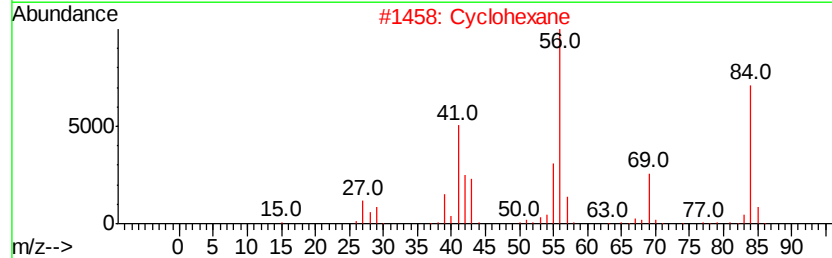
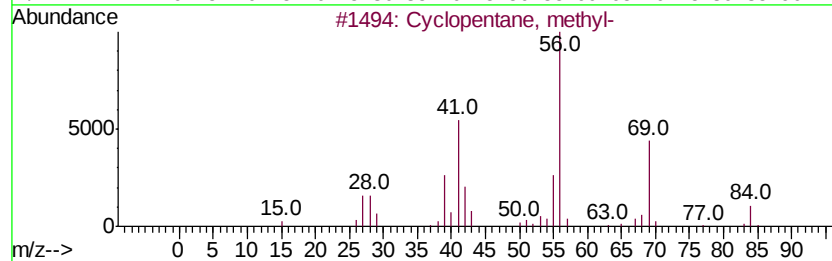
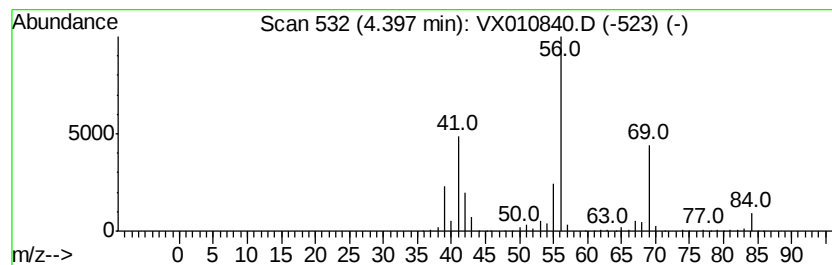
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	62.74 ug/l	771823	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		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	46



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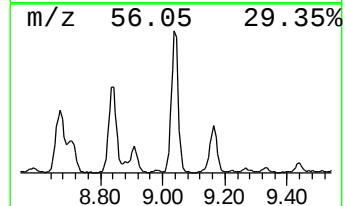
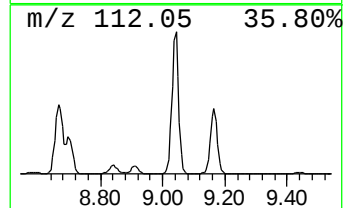
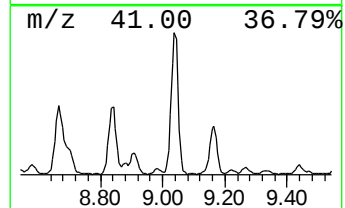
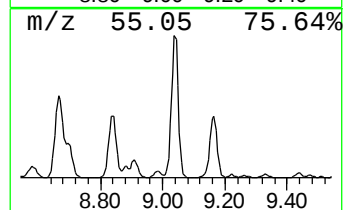
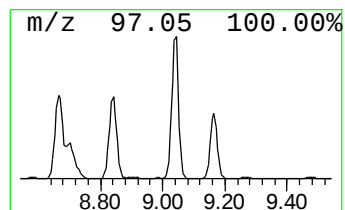
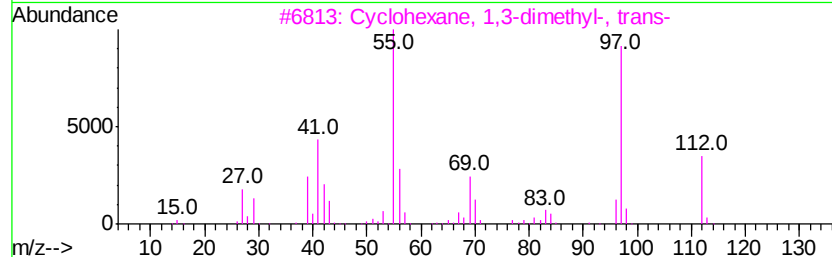
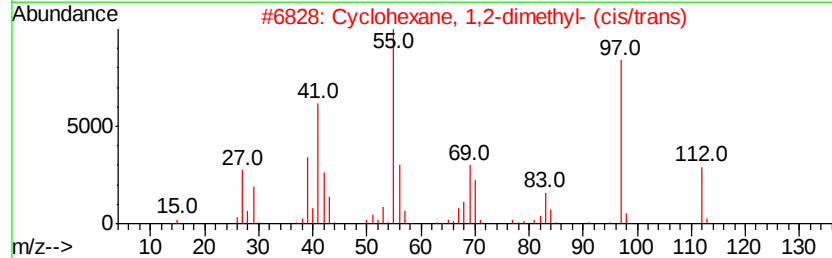
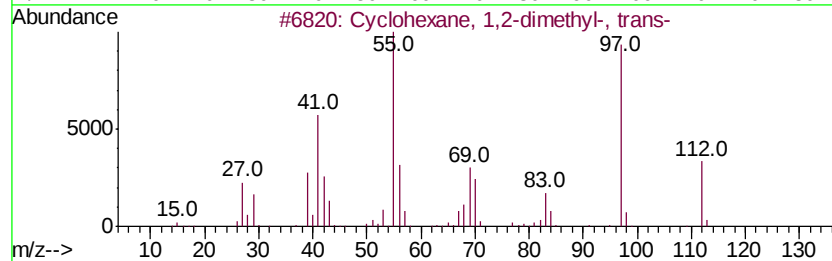
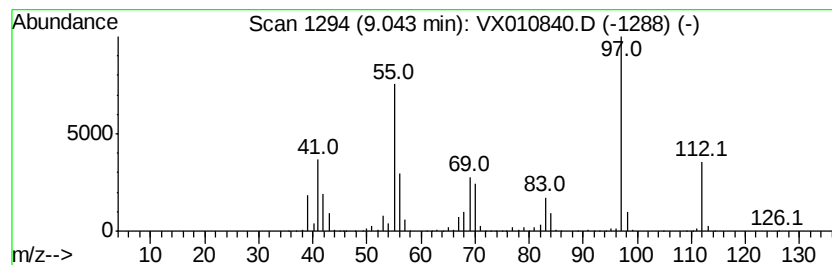
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.04	40.16 ug/l	700092	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	96
2		Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	91
3		Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
4		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	81
5		Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	76



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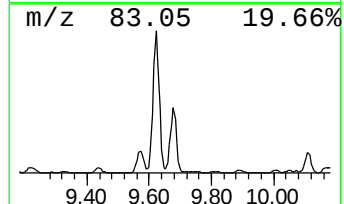
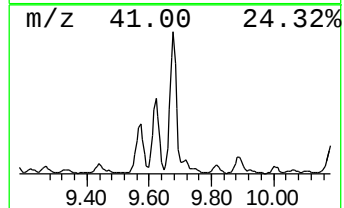
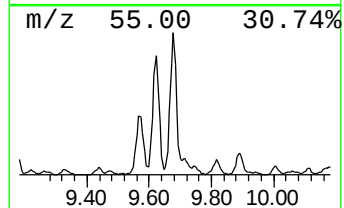
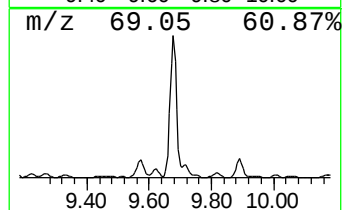
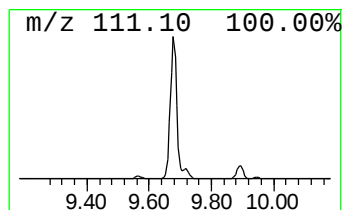
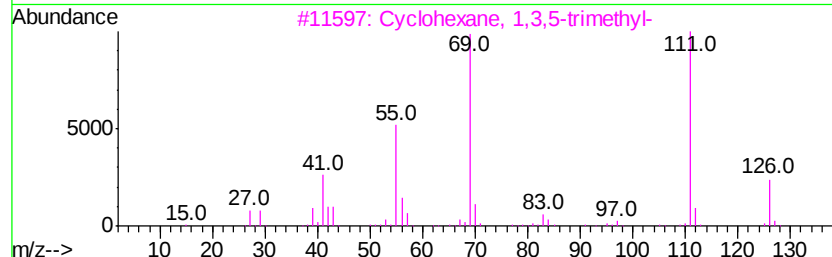
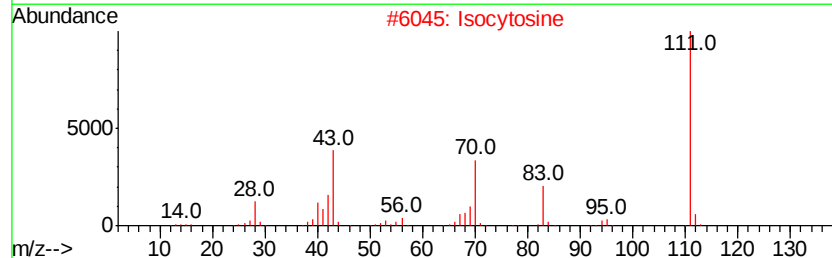
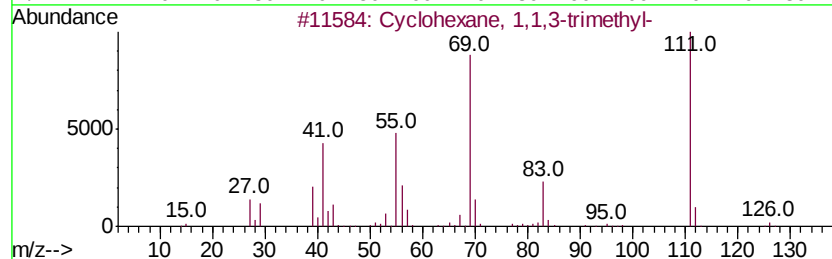
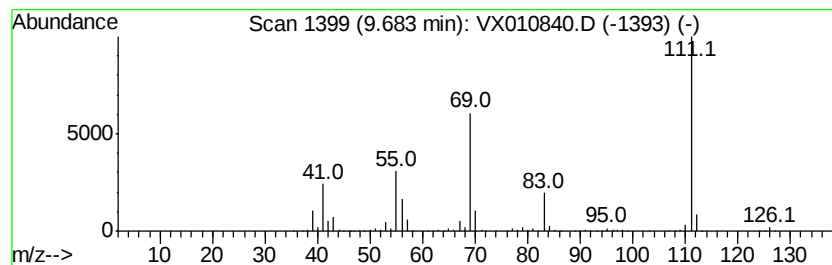
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.68	35.89 ug/l	625620	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	97
2		Isocytosine	111	C4H5N3O	000108-53-2	72
3		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
4		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-27-3	64
5		Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	64



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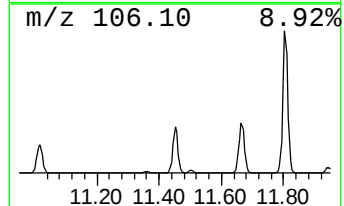
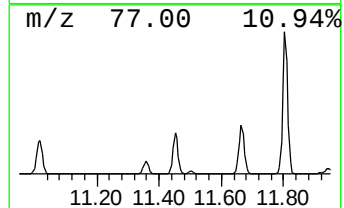
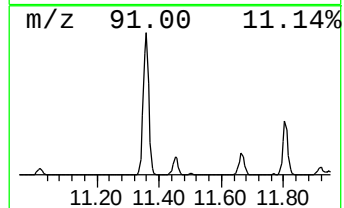
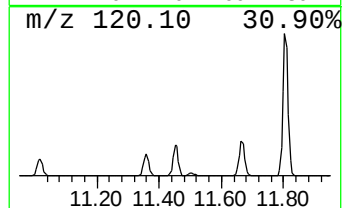
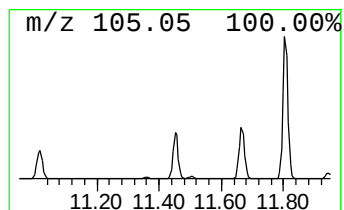
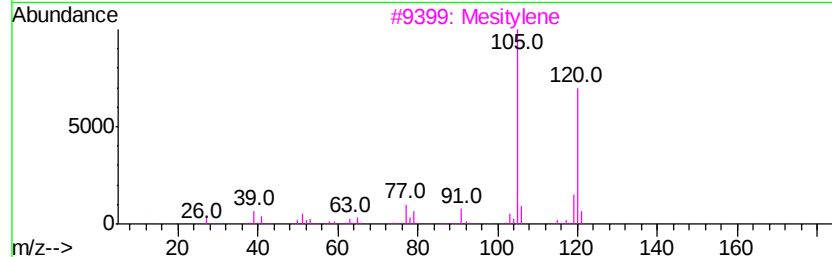
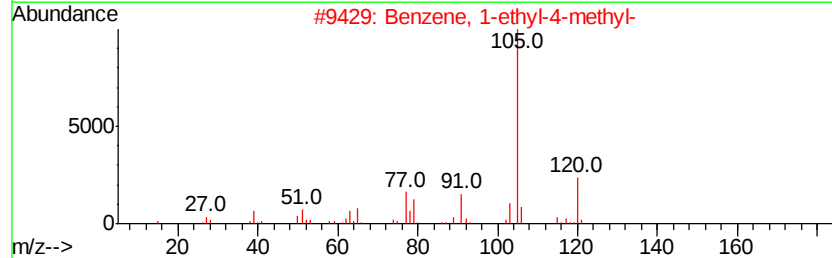
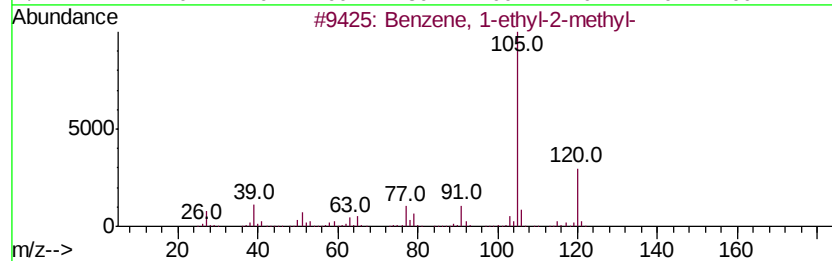
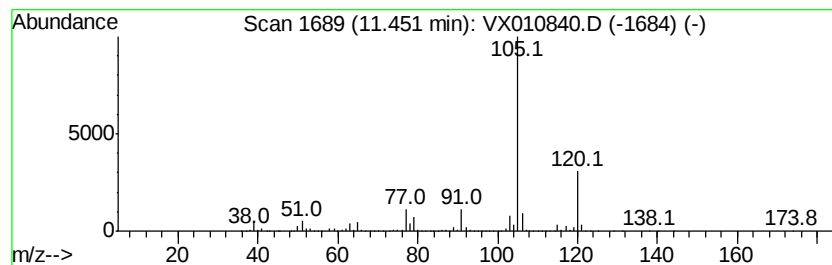
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.45	70.71 ug/l	1558460	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-4-methyl-	120	C9H12	000622-96-8	93
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010840.D
Acq On : 12 Jul 2019 16:49
Operator : JC/SP
Sample : K3786-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0589

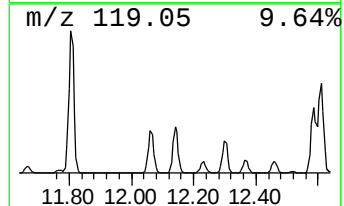
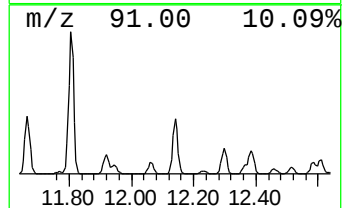
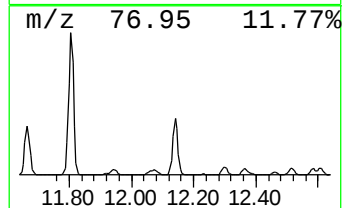
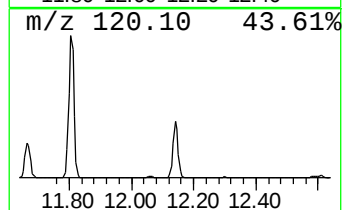
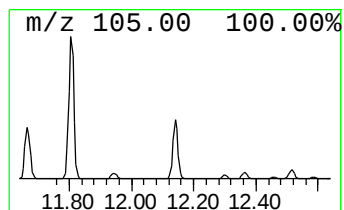
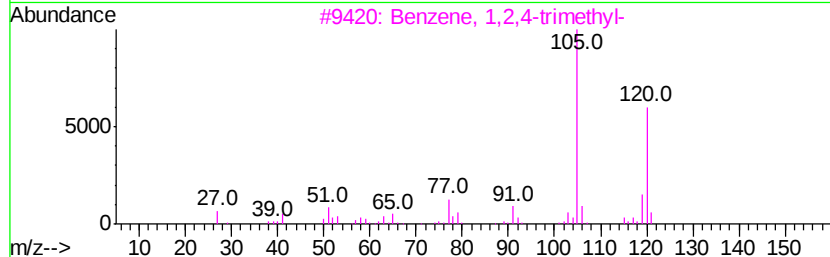
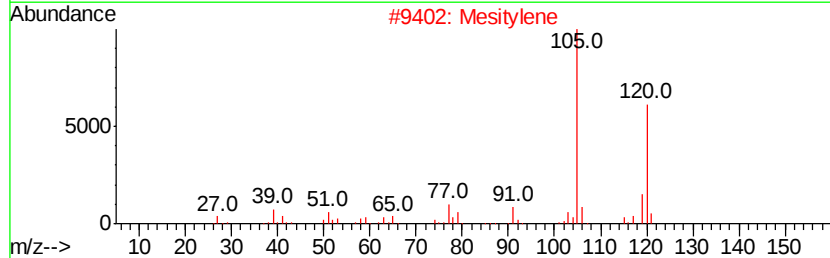
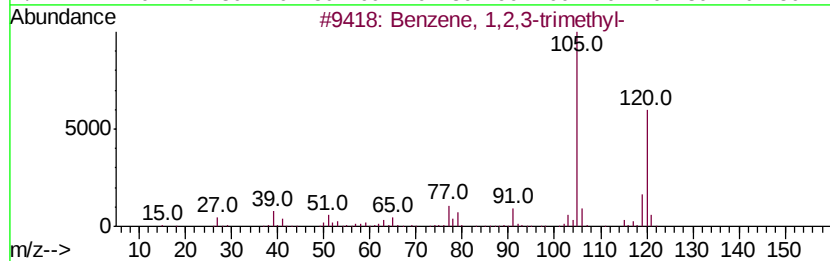
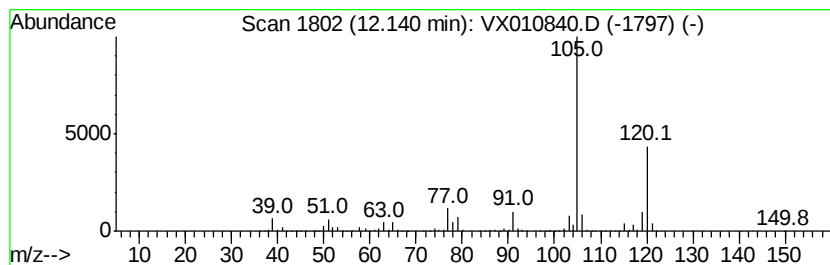
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	99.38 ug/l	2190490	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010840.D
Acq On : 12 Jul 2019 16:49
Operator : JC/SP
Sample : K3786-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

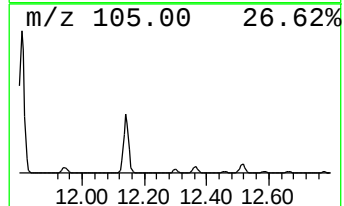
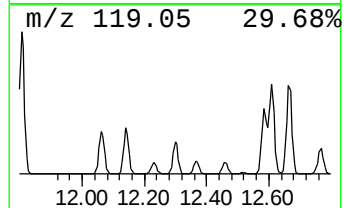
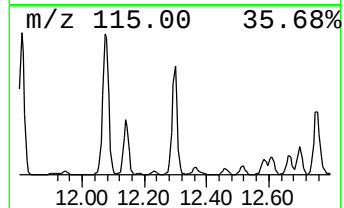
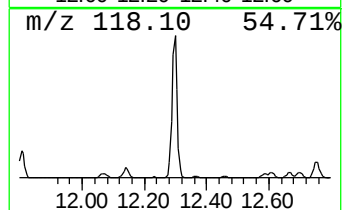
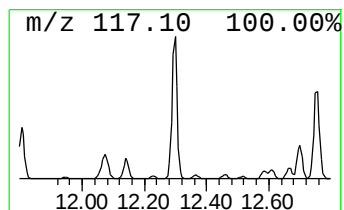
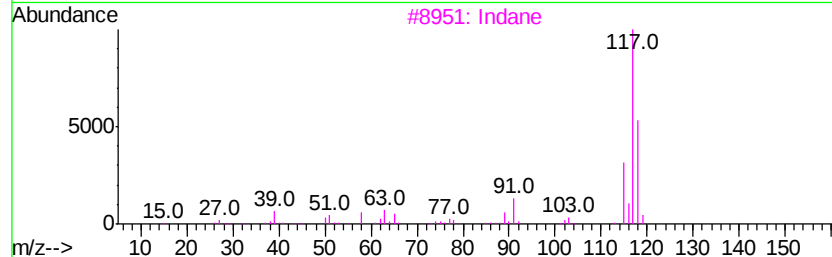
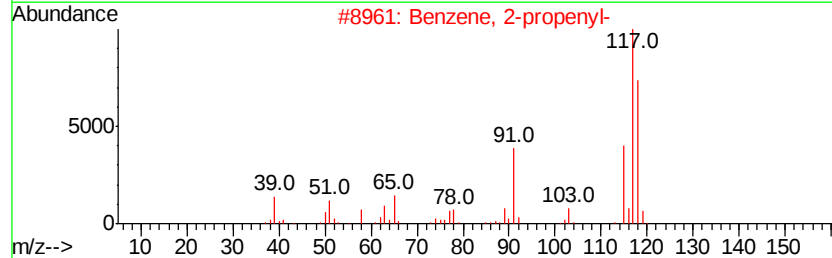
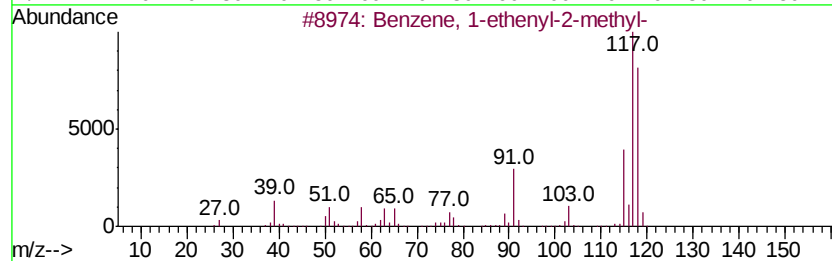
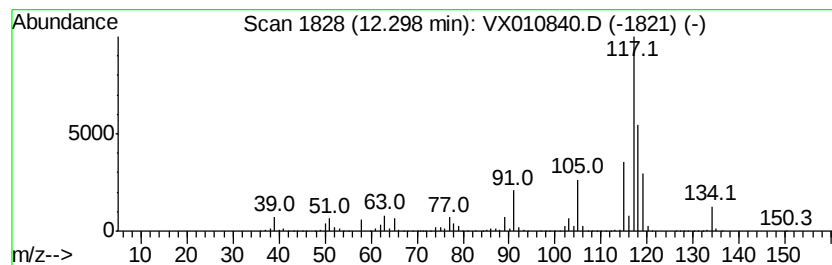
Instrument :
MSVOA_X
ClientSampleId :
FL-0589

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-ethenyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	32.97 ug/l	726751	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 83
2		Benzene, 2-propenyl-	118 C9H10	000300-57-2 64
3		Indane	118 C9H10	000496-11-7 64
4		Benzene, 1-propenyl-	118 C9H10	000637-50-3 64
5		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010840.D
Acq On : 12 Jul 2019 16:49
Operator : JC/SP
Sample : K3786-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0589

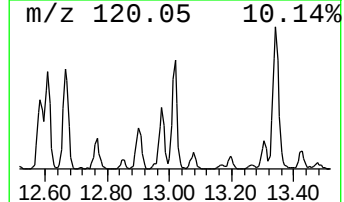
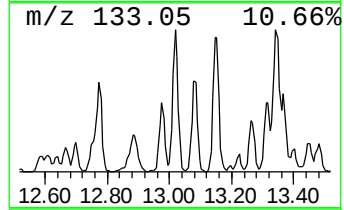
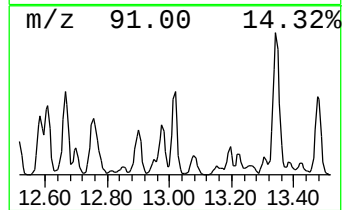
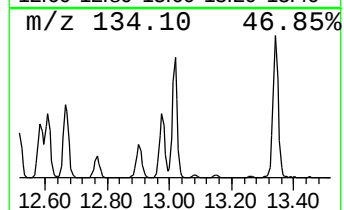
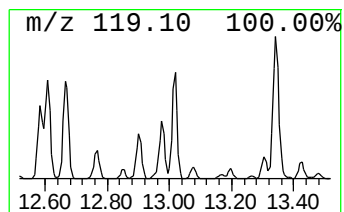
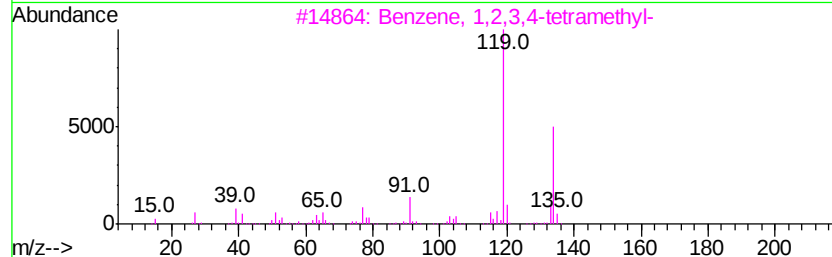
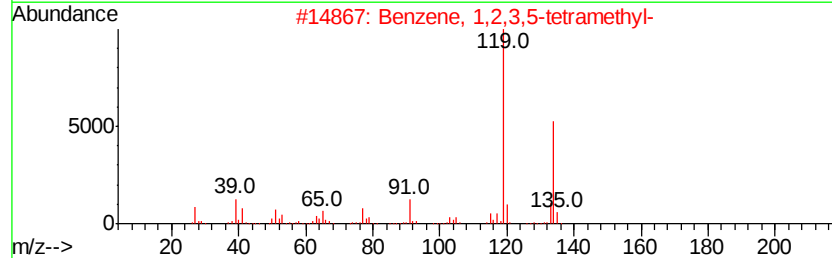
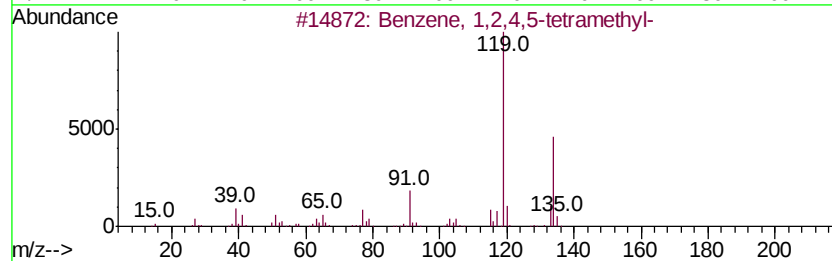
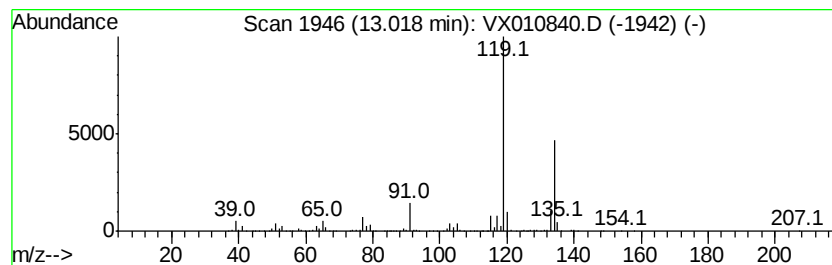
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	21.12 ug/l	465492	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010840.D
Acq On : 12 Jul 2019 16:49
Operator : JC/SP
Sample : K3786-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0589

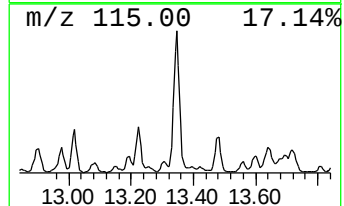
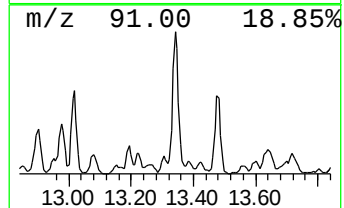
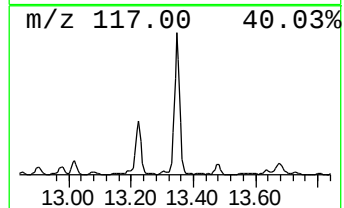
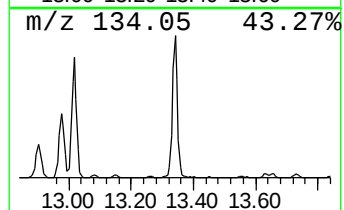
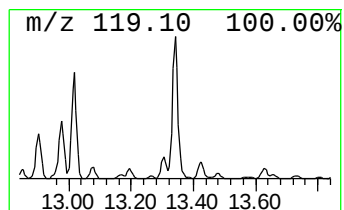
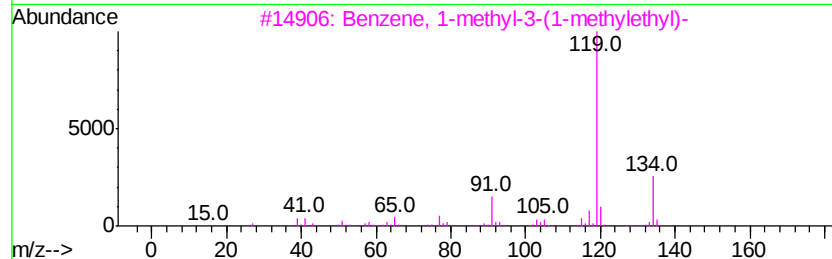
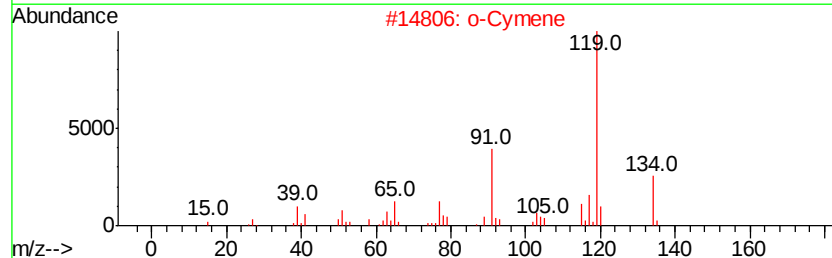
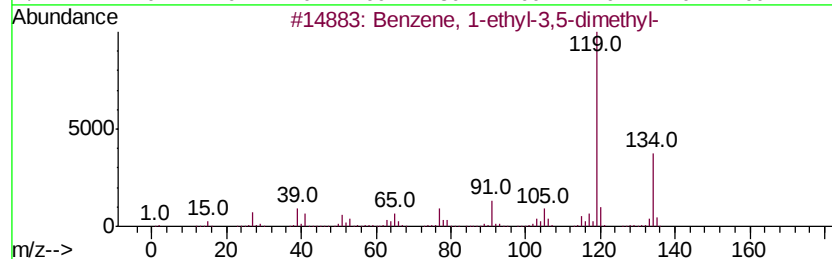
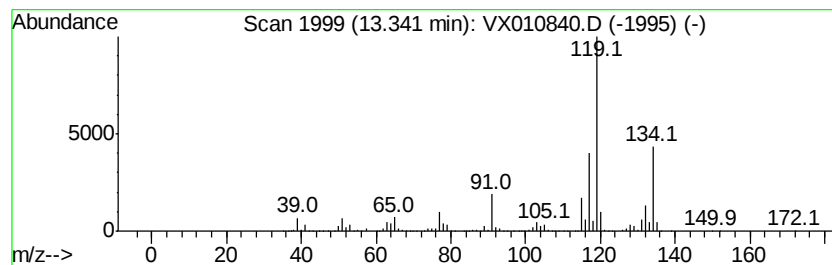
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	41.31 ug/l	910612	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	76
2		o-Cymene	134	C10H14	000527-84-4	70
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	70
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010840.D
Acq On : 12 Jul 2019 16:49
Operator : JC/SP
Sample : K3786-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0589

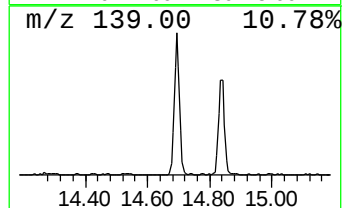
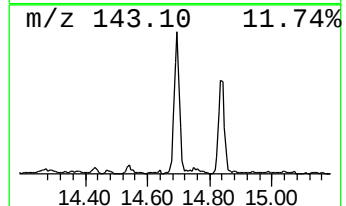
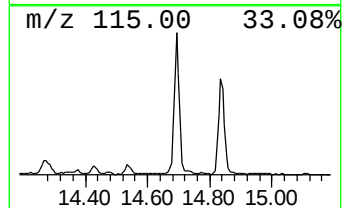
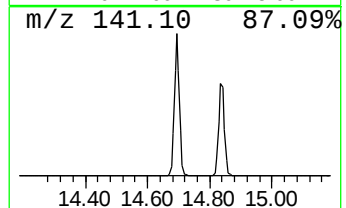
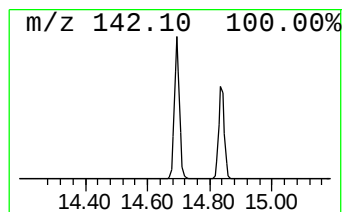
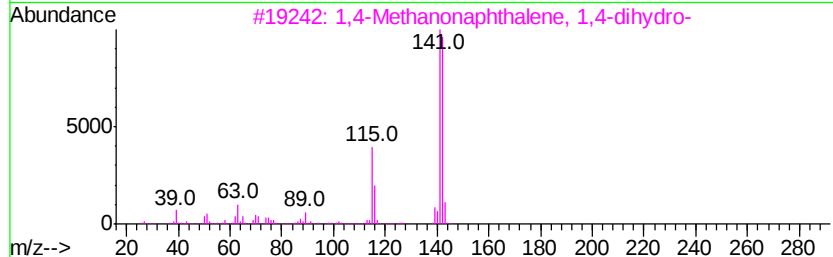
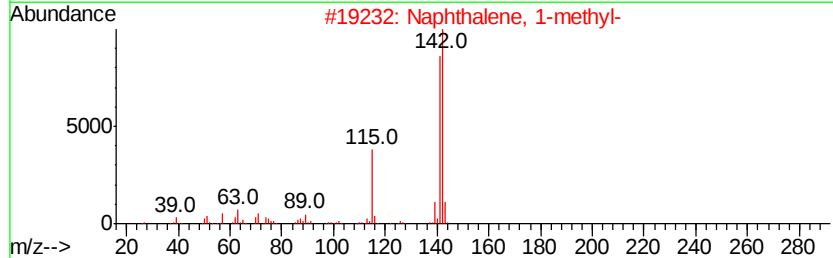
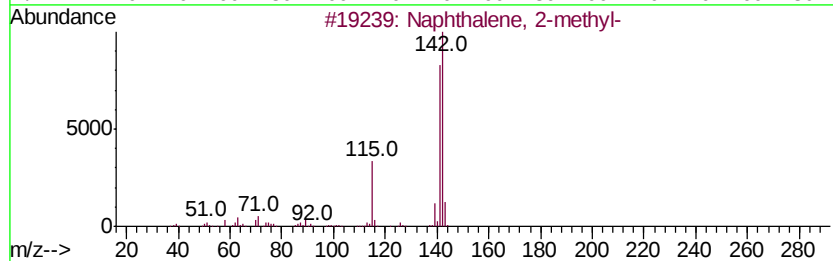
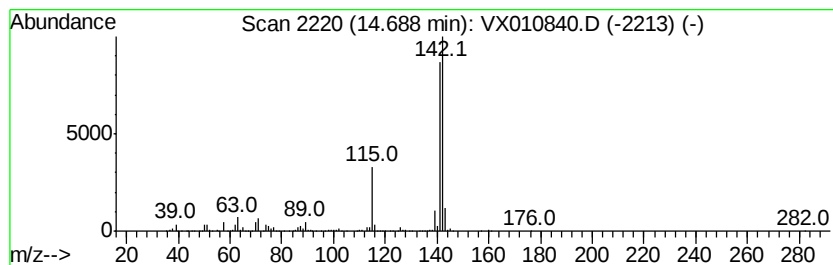
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	23.23 ug/l	511983	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		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	78



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071219\
Data File : VX010840.D
Acq On : 12 Jul 2019 16:49
Operator : JC/SP
Sample : K3786-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0589

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
Cyclopentane, met...	4.40	62.7	ug/l	771823	1	5.67	615123	50.0
Cyclohexane, 1,2-...	9.04	40.2	ug/l	700092	3	10.11	871633	50.0
Cyclohexane, 1,1,...	9.68	35.9	ug/l	625620	3	10.11	871633	50.0
Benzene, 1-ethyl-...	11.45	70.7	ug/l	1558460	4	12.08	1102060	50.0
Benzene, 1,2,3-tr...	12.14	99.4	ug/l	2190490	4	12.08	1102060	50.0
Benzene, 1-etheny...	12.30	33.0	ug/l	726751	4	12.08	1102060	50.0
Benzene, 1,2,4,5-...	13.02	21.1	ug/l	465492	4	12.08	1102060	50.0
Benzene, 1-ethyl-...	13.34	41.3	ug/l	910612	4	12.08	1102060	50.0
Naphthalene, 1-me...	14.69	23.2	ug/l	511983	4	12.08	1102060	50.0