

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071219\
 Data File : VX010862.D
 Acq On : 13 Jul 2019 01:45
 Operator : JC/SP
 Sample : K3791-09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0595

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	2.891	280	285	298	rVB	135520	295812	2.69%	0.365%
2	3.172	321	331	345	rVB	141849	333552	3.04%	0.412%
3	3.525	380	389	401	rBV	78536	182965	1.67%	0.226%
4	4.403	523	533	548	rVB	229495	674611	6.14%	0.833%
5	5.501	700	713	719	rBV	139687	399606	3.64%	0.494%
6	5.586	719	727	734	rVV3	224067	726717	6.61%	0.898%
7	5.659	734	739	747	rVB	195554	503387	4.58%	0.622%
8	5.933	773	784	796	rBV5	110332	367091	3.34%	0.454%
9	6.061	796	805	818	rVB	140684	360218	3.28%	0.445%
10	6.263	825	838	846	rBV3	70002	227929	2.07%	0.282%
11	6.360	846	854	862	rVV2	85559	249289	2.27%	0.308%
12	6.458	862	870	883	rVB	108007	300996	2.74%	0.372%
13	6.860	926	936	947	rBV	355717	845294	7.69%	1.044%
14	7.464	1021	1035	1046	rBV	881197	1986974	18.08%	2.455%
15	7.732	1073	1079	1088	rVB	80786	164585	1.50%	0.203%
16	8.031	1121	1128	1143	rVB2	59508	178097	1.62%	0.220%
17	8.177	1143	1152	1160	rBV2	53441	119021	1.08%	0.147%
18	8.665	1225	1232	1235	rBV2	180214	313234	2.85%	0.387%
19	8.713	1235	1240	1253	rVB	786258	1300661	11.84%	1.607%
20	8.835	1254	1260	1265	rBV	97496	169919	1.55%	0.210%
21	9.036	1288	1293	1300	rVB	200091	329798	3.00%	0.407%
22	9.165	1305	1314	1320	rBV	91871	155768	1.42%	0.192%
23	9.573	1374	1381	1385	rBV5	81109	147413	1.34%	0.182%
24	9.622	1385	1389	1393	rVB	187161	248459	2.26%	0.307%
25	9.677	1393	1398	1402	rBV	144885	205783	1.87%	0.254%
26	10.109	1464	1469	1475	rVB	766093	1017508	9.26%	1.257%
27	10.183	1475	1481	1486	rVB	81328	122216	1.11%	0.151%
28	10.250	1486	1492	1498	rBV	387232	499087	4.54%	0.617%
29	10.359	1501	1510	1516	rBV	901198	1311283	11.93%	1.620%
30	10.640	1545	1556	1564	rBV2	59090	156013	1.42%	0.193%
31	10.835	1584	1588	1593	rVB	88314	121967	1.11%	0.151%
32	11.018	1613	1618	1623	rBV	714045	912179	8.30%	1.127%
33	11.134	1632	1637	1642	rBV	702097	878607	8.00%	1.085%
34	11.359	1668	1674	1681	rBV	1182224	1424439	12.96%	1.760%

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35	11.432	1681	1686	1688	rBV	4039897	5167905	47.03%	6.385%
36	11.506	1694	1698	1706	rVB	3126833	3785473	34.45%	4.677%
37	11.664	1719	1724	1732	rVB	3337366	4143822	37.71%	5.119%
38	11.804	1742	1747	1753	rVB	8687955	10988452	100.00%	13.576%
39	11.920	1758	1766	1767	rBV	123085	174970	1.59%	0.216%
40	11.945	1767	1770	1775	rVB	325044	388115	3.53%	0.479%
41	12.024	1776	1783	1786	rBV	715272	851311	7.75%	1.052%
42	12.073	1786	1791	1797	rVB2	1045786	1608478	14.64%	1.987%
43	12.140	1797	1802	1808	rBV	3940514	4754460	43.27%	5.874%
44	12.231	1812	1817	1822	rVB	168742	213163	1.94%	0.263%
45	12.298	1822	1828	1830	rBV2	997891	1329783	12.10%	1.643%
46	12.322	1830	1832	1835	rVV	1009962	1008362	9.18%	1.246%
47	12.371	1835	1840	1849	rVB2	1427717	2108319	19.19%	2.605%
48	12.457	1849	1854	1859	rBV	410200	498197	4.53%	0.615%
49	12.518	1859	1864	1870	rVB	1068571	1284433	11.69%	1.587%
50	12.585	1870	1875	1877	rBV	1072864	1405298	12.79%	1.736%
51	12.609	1877	1879	1883	rVB	1271333	1259341	11.46%	1.556%
52	12.664	1884	1888	1892	rVV	1204100	1436879	13.08%	1.775%
53	12.700	1892	1894	1899	rVB	307852	320699	2.92%	0.396%
54	12.761	1899	1904	1911	rBV4	812257	1686095	15.34%	2.083%
55	12.847	1914	1918	1922	rBV2	187435	243258	2.21%	0.301%
56	12.902	1922	1927	1932	rVB2	947993	1256609	11.44%	1.552%
57	12.975	1932	1939	1942	rBV	1107325	1423941	12.96%	1.759%
58	13.017	1942	1946	1952	rVB	1765395	2121687	19.31%	2.621%
59	13.078	1952	1956	1961	rBV3	238617	422527	3.85%	0.522%
60	13.194	1972	1975	1977	rBV2	186118	228331	2.08%	0.282%
61	13.225	1977	1980	1983	rVB	272150	311479	2.83%	0.385%
62	13.261	1983	1986	1989	rVB2	102210	137848	1.25%	0.170%
63	13.304	1989	1993	1995	rBV3	210996	258375	2.35%	0.319%
64	13.341	1995	1999	2005	rVB2	2253050	3073729	27.97%	3.797%
65	13.426	2010	2013	2017	rVV	126720	172276	1.57%	0.213%
66	13.475	2017	2021	2029	rVB	805345	1068028	9.72%	1.319%
67	13.560	2030	2035	2038	rBV2	136597	210363	1.91%	0.260%
68	13.597	2038	2041	2044	rVV	143409	209562	1.91%	0.259%
69	13.633	2044	2047	2053	rVV3	419214	840327	7.65%	1.038%
70	13.694	2054	2057	2059	rVV2	191658	298021	2.71%	0.368%
71	13.719	2059	2061	2067	rVB2	309266	461826	4.20%	0.571%

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72	13.834	2071	2080	2087	rVB2	623171	1091429	9.93%	1.348%
73	13.908	2087	2092	2097	rVB2	130307	191584	1.74%	0.237%
74	13.987	2097	2105	2108	rBV2	243576	390257	3.55%	0.482%
75	14.127	2124	2128	2131	rBV2	91561	127343	1.16%	0.157%
76	14.279	2146	2153	2159	rVB3	227503	491579	4.47%	0.607%
77	14.377	2165	2169	2173	rVV	91694	111207	1.01%	0.137%
78	14.426	2173	2177	2181	rVB2	99752	134005	1.22%	0.166%
79	14.536	2190	2195	2204	rBV2	260667	364271	3.32%	0.450%
80	14.694	2213	2221	2225	rBV	1379271	1726927	15.72%	2.134%
81	14.834	2239	2244	2253	rBV	897906	1164814	10.60%	1.439%
82	15.548	2354	2361	2366	rBV	139494	224554	2.04%	0.277%
83	15.682	2374	2383	2387	rBV	154877	273123	2.49%	0.337%
84	15.724	2387	2390	2397	rVB	102231	150991	1.37%	0.187%
85	15.913	2411	2421	2431	rVB2	41722	117646	1.07%	0.145%

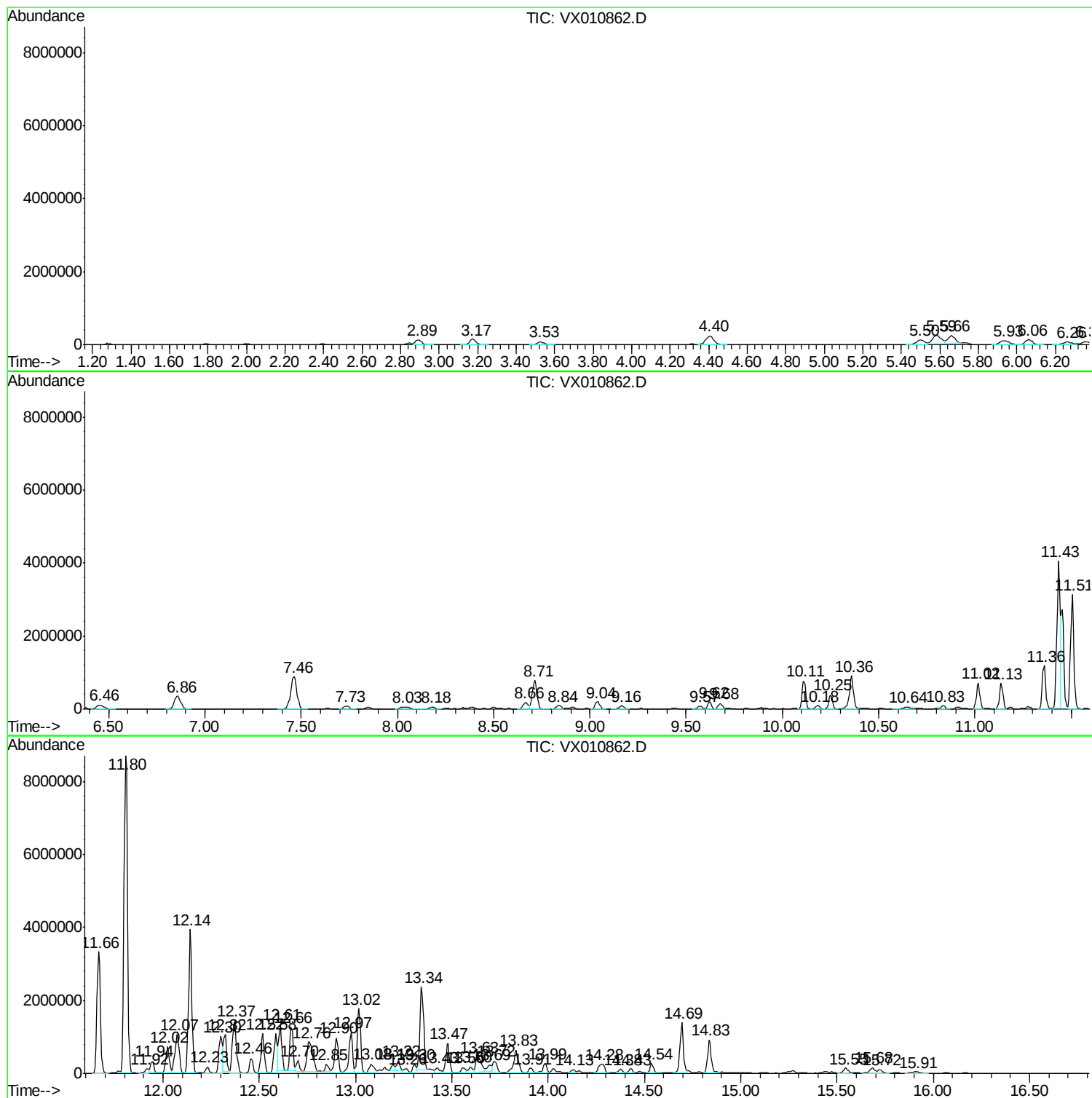
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



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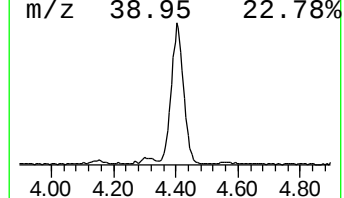
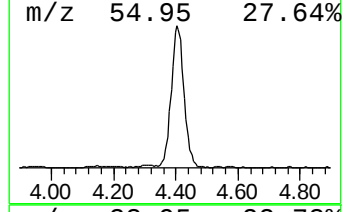
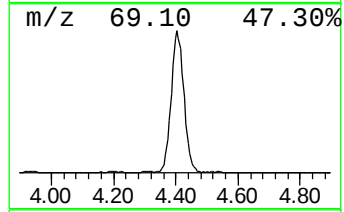
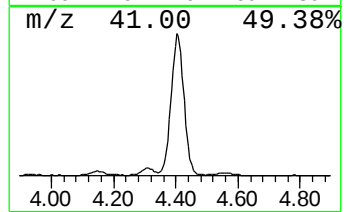
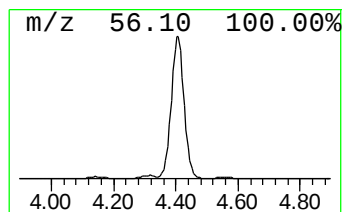
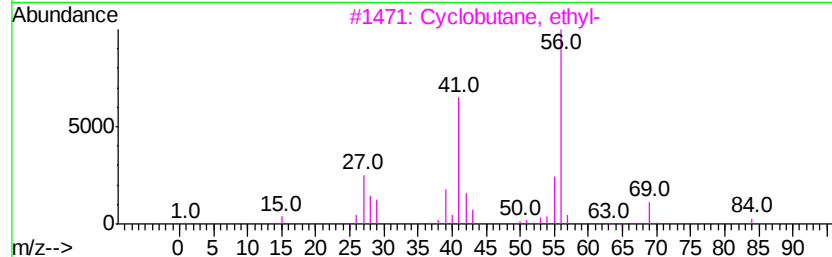
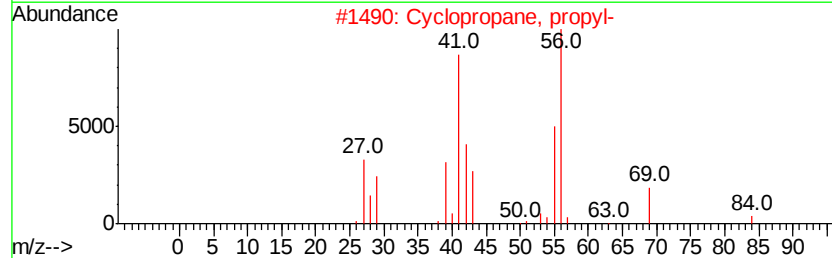
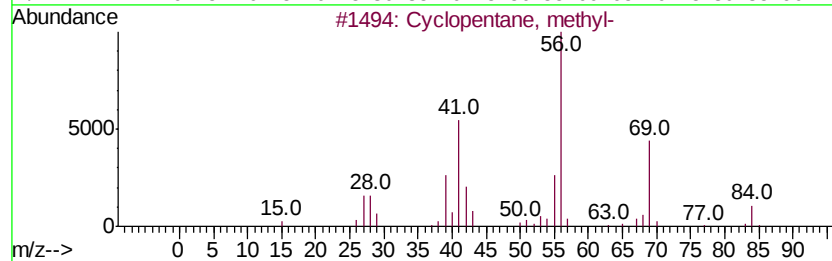
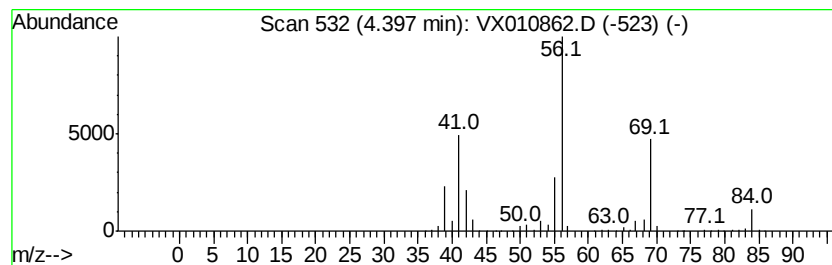
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	67.01 ug/l	674611	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopropane, propyl-	84	C6H12	002415-72-7	78
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	46



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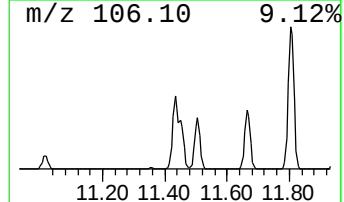
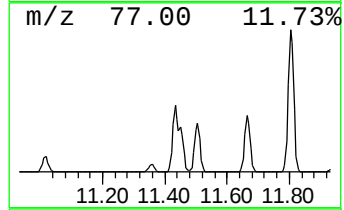
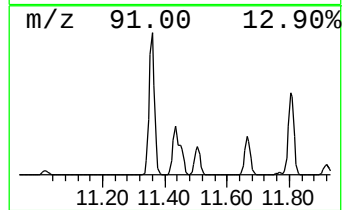
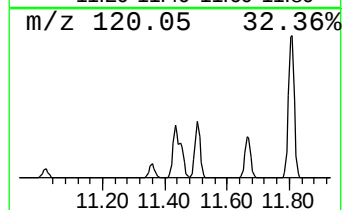
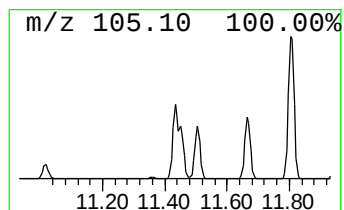
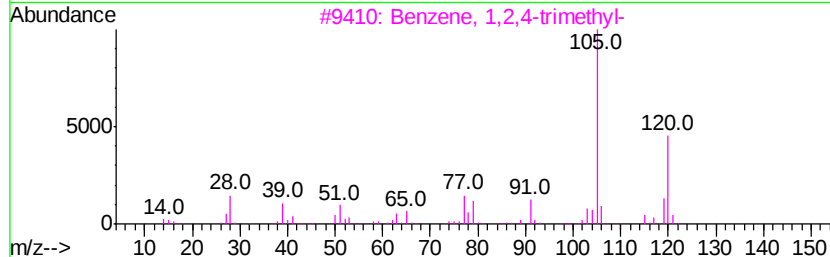
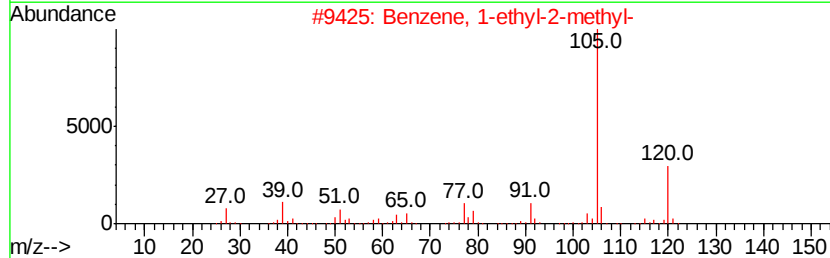
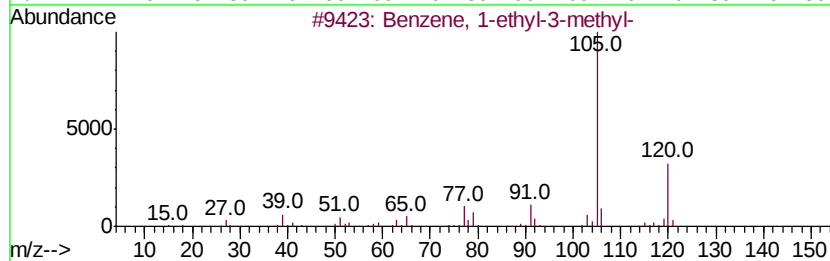
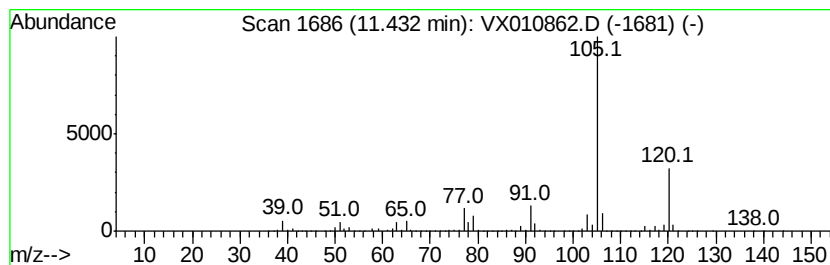
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	160.65 ug/l	5167910	1,4-Dichlorobenzene-d4	12.08

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



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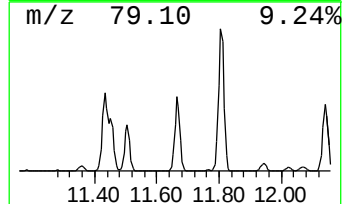
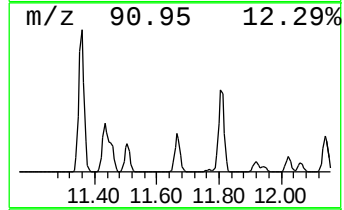
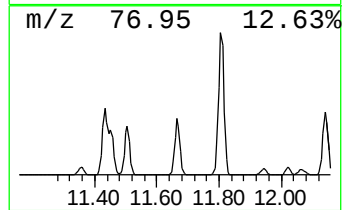
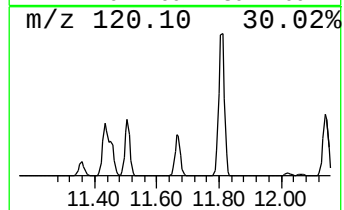
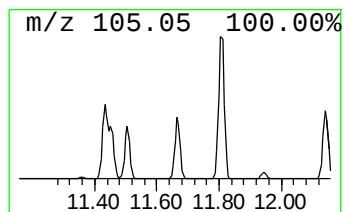
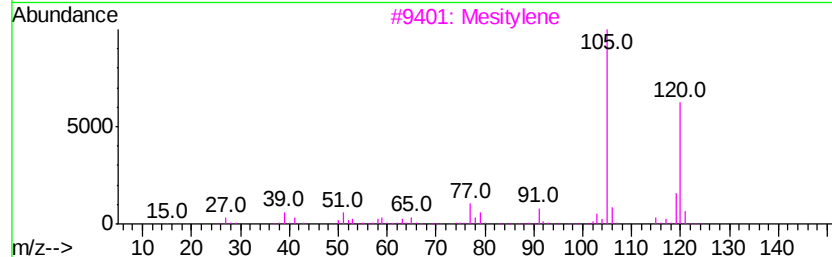
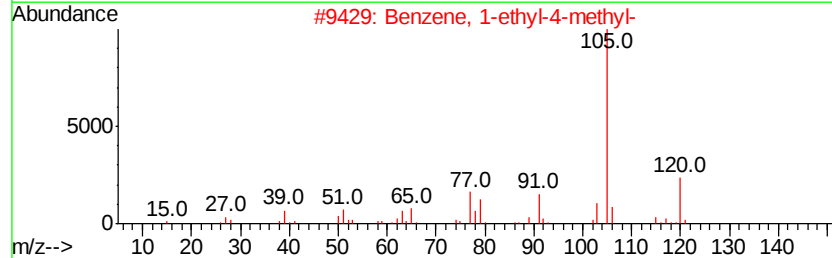
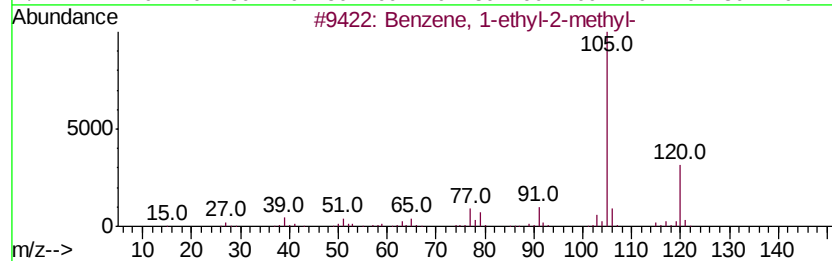
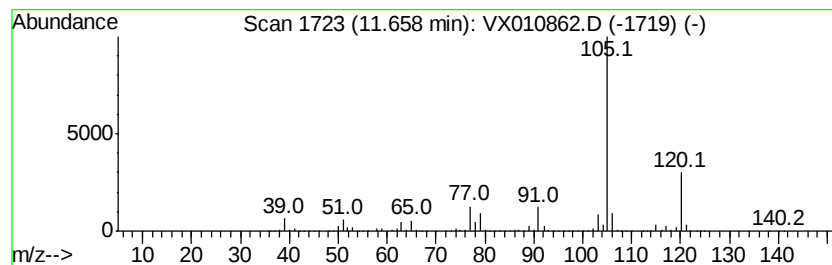
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	128.81 ug/l	4143820	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 94
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, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91



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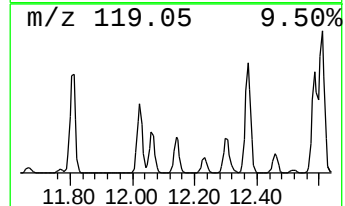
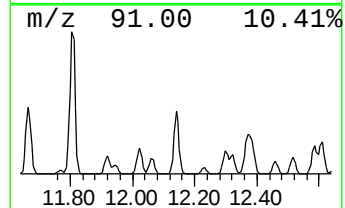
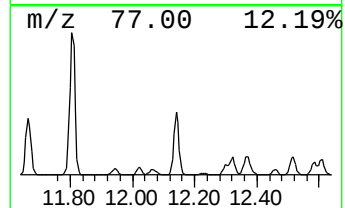
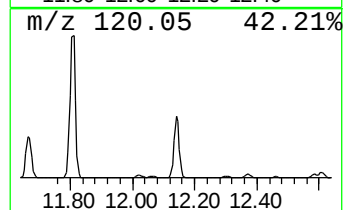
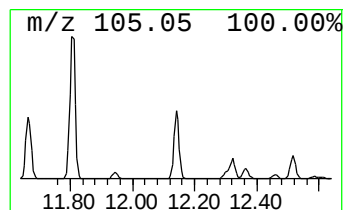
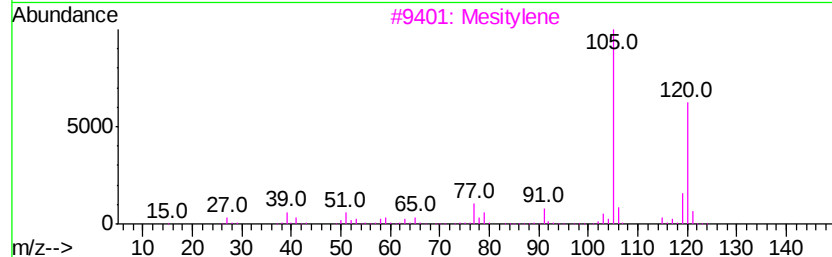
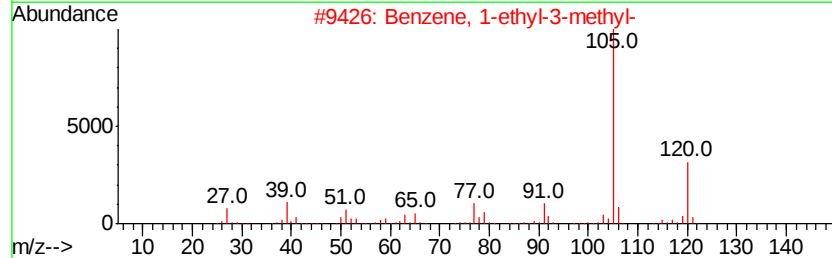
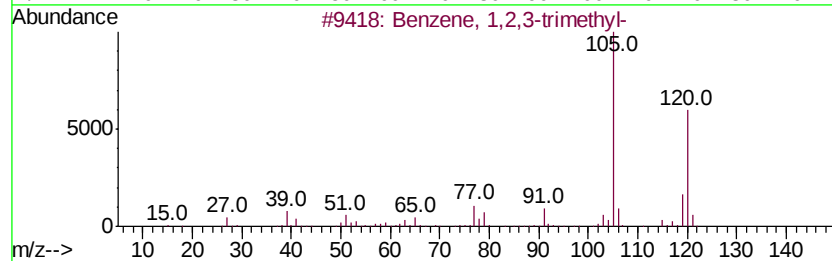
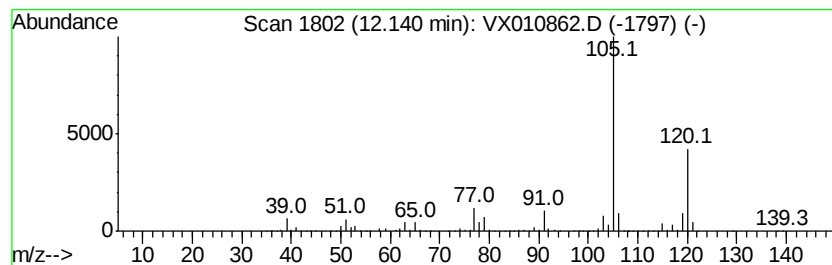
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	147.79 ug/l	4754460	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010862.D
Acq On : 13 Jul 2019 01:45
Operator : JC/SP
Sample : K3791-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

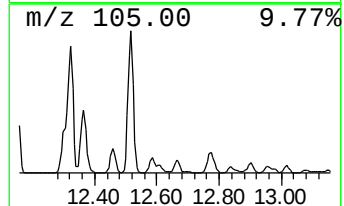
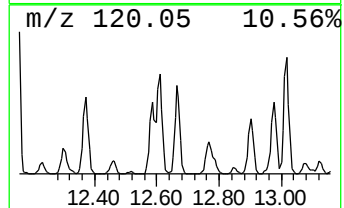
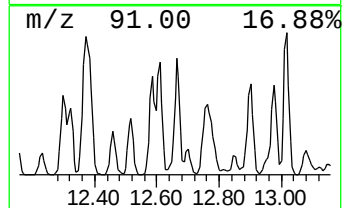
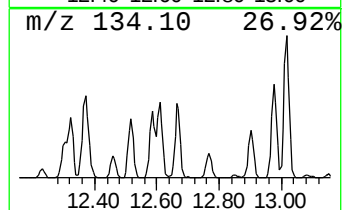
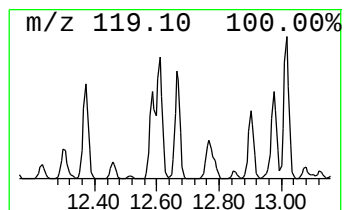
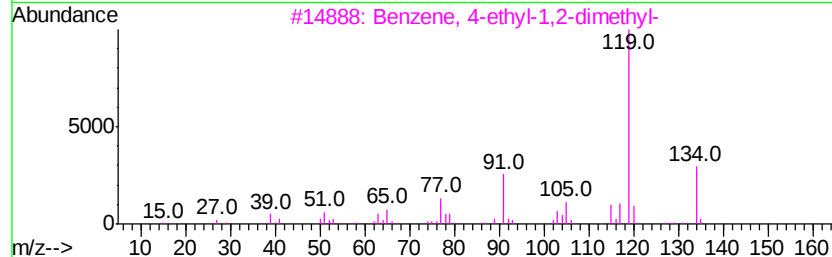
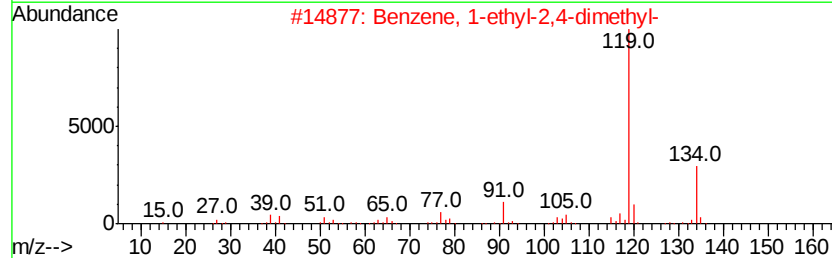
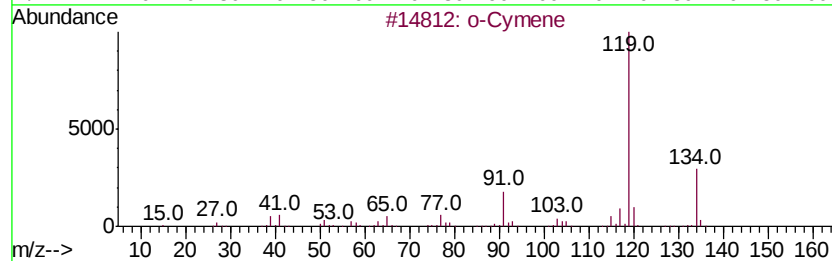
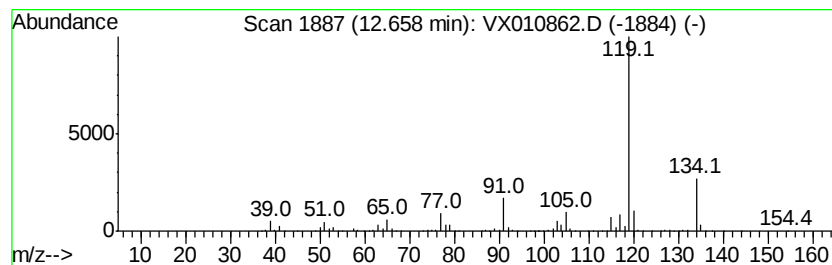
Instrument :
MSVOA_X
ClientSampleId :
FL-0595

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 5 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	44.67 ug/l	1436880	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 95
2		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 95
3		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 95
4		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 94
5		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010862.D
Acq On : 13 Jul 2019 01:45
Operator : JC/SP
Sample : K3791-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

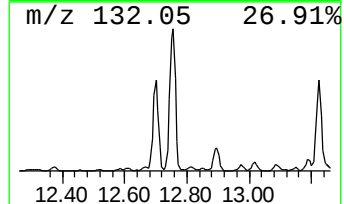
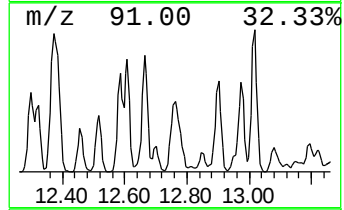
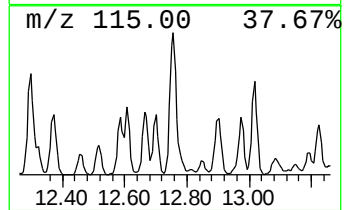
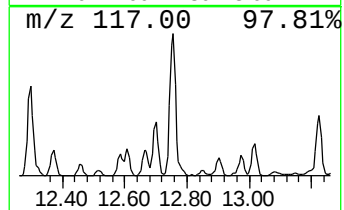
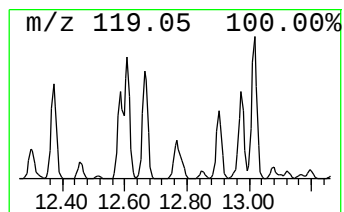
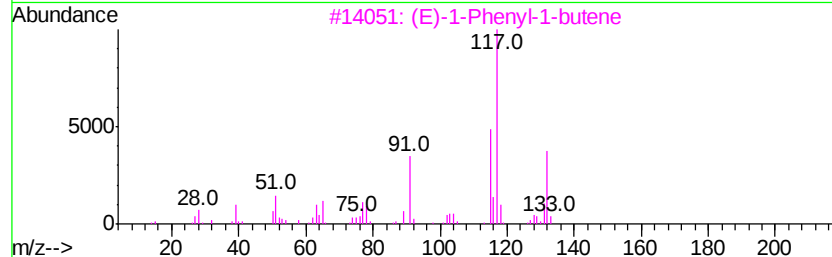
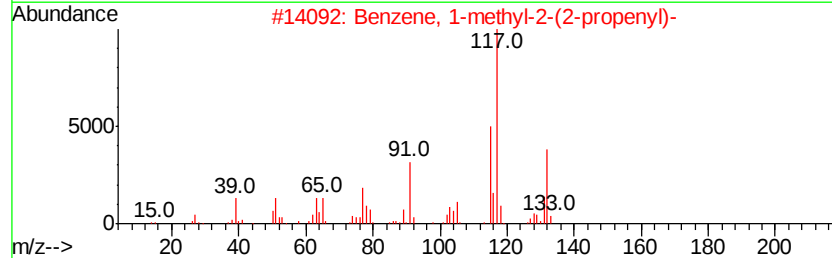
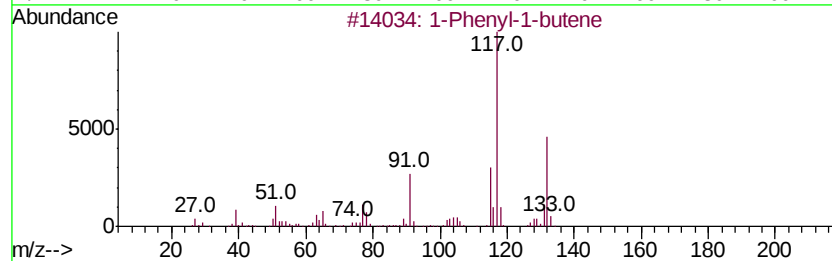
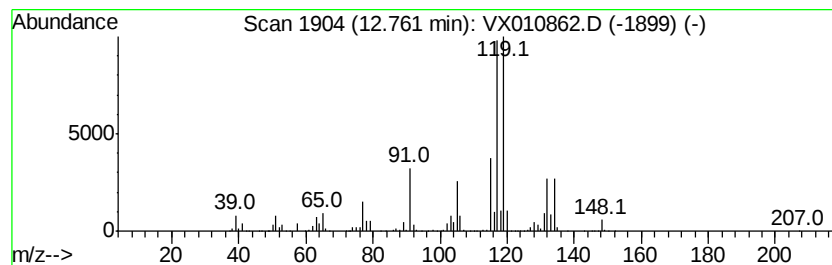
Instrument :
MSVOA_X
ClientSampleId :
FL-0595

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 6 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	52.41 ug/l	1686100	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132 C10H12	000824-90-8	89
2	Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8	80
3	(E)-1-Phenyl-1-butene	132 C10H12	001005-64-7	64
4	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	50
5	Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010862.D
Acq On : 13 Jul 2019 01:45
Operator : JC/SP
Sample : K3791-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0595

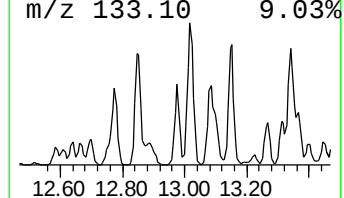
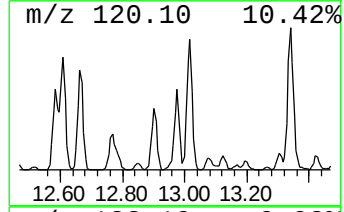
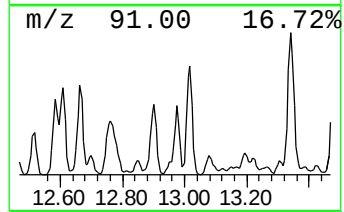
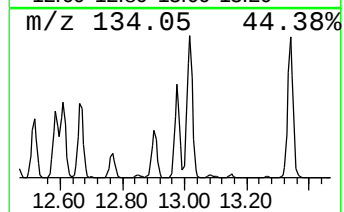
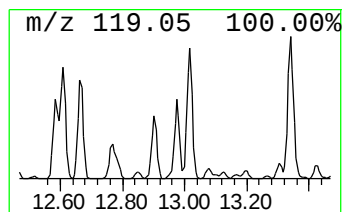
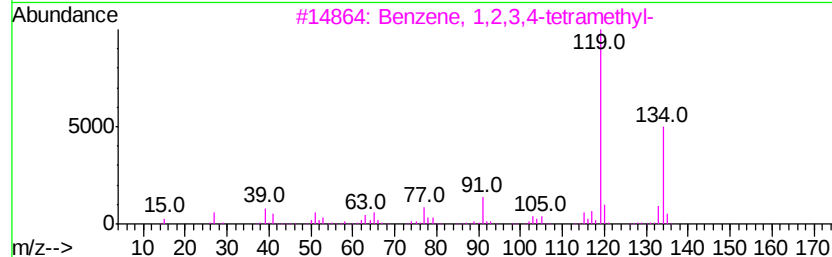
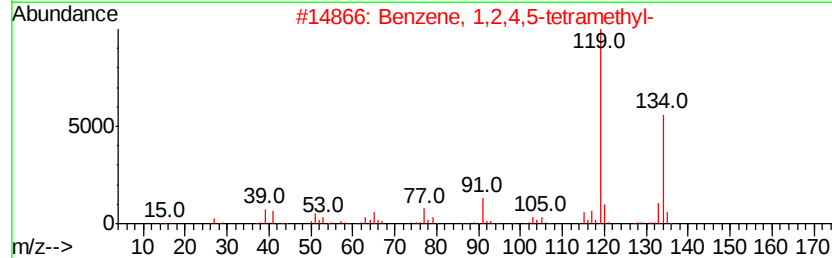
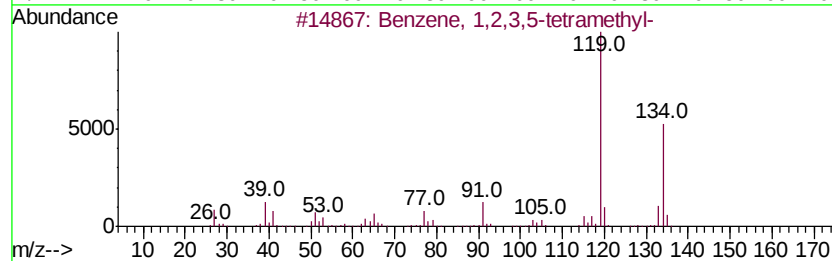
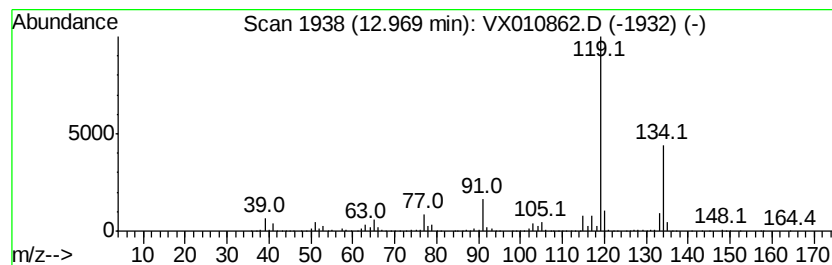
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	44.26 ug/l	1423940	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010862.D
Acq On : 13 Jul 2019 01:45
Operator : JC/SP
Sample : K3791-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0595

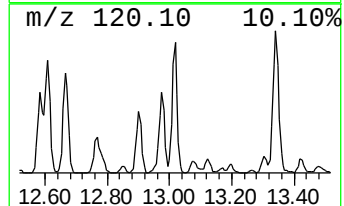
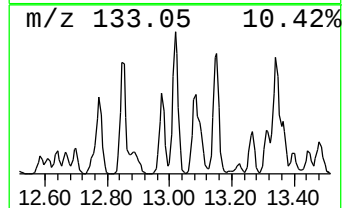
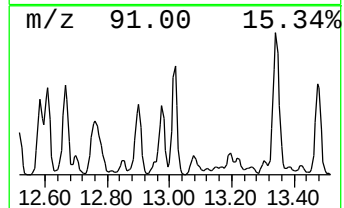
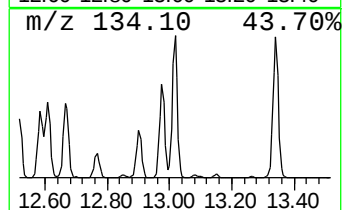
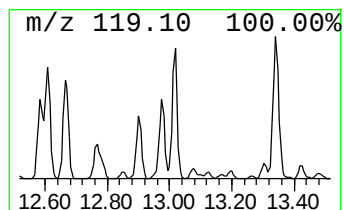
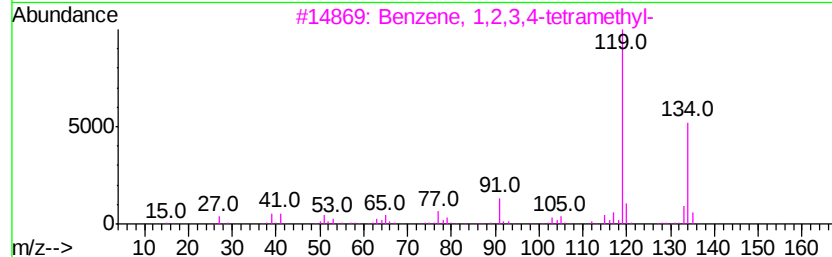
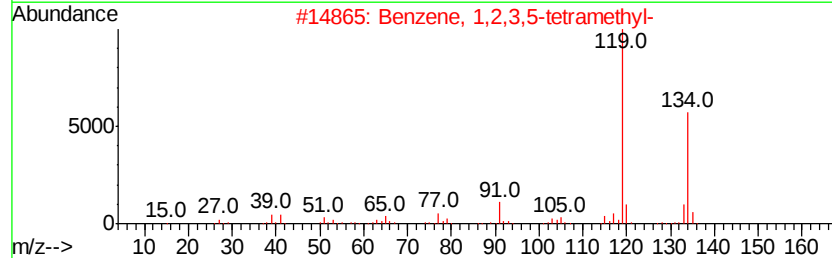
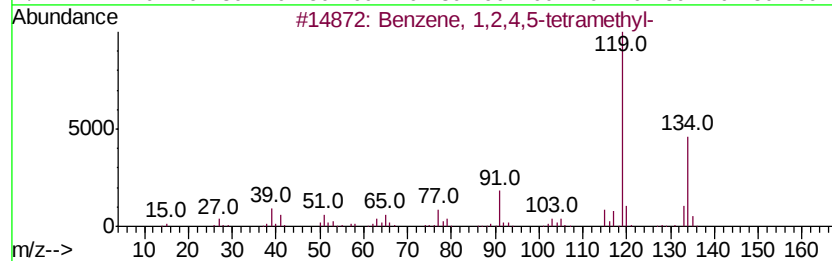
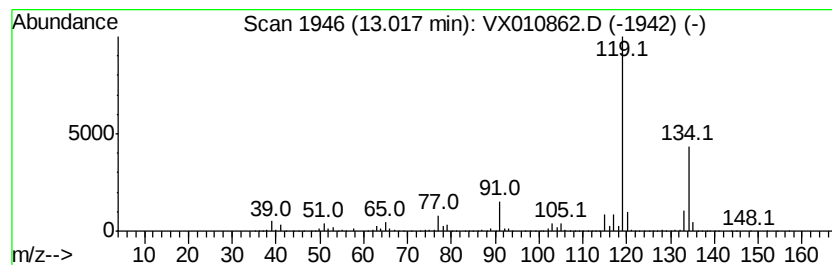
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	65.95 ug/l	2121690	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	97
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-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071219\
Data File : VX010862.D
Acq On : 13 Jul 2019 01:45
Operator : JC/SP
Sample : K3791-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0595

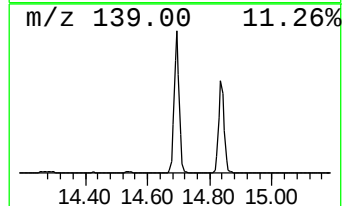
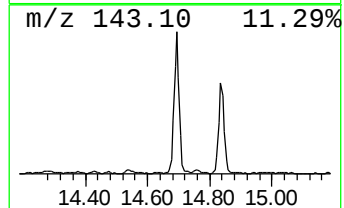
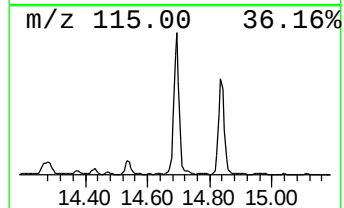
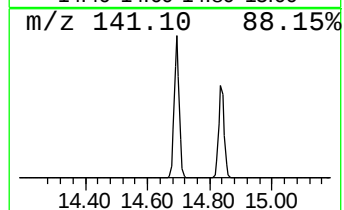
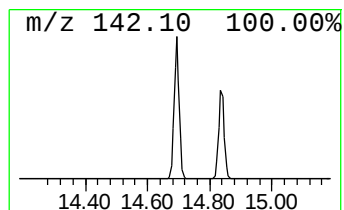
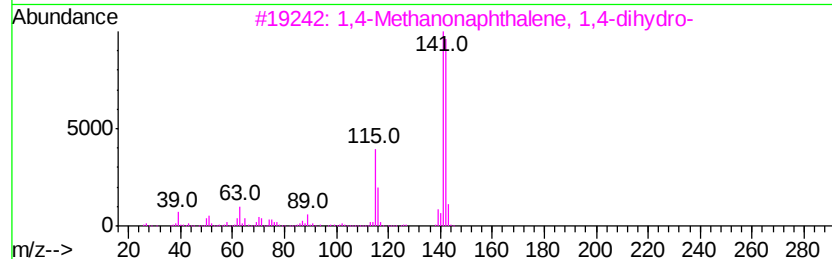
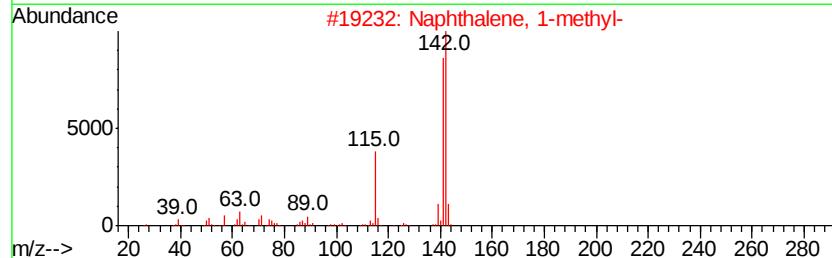
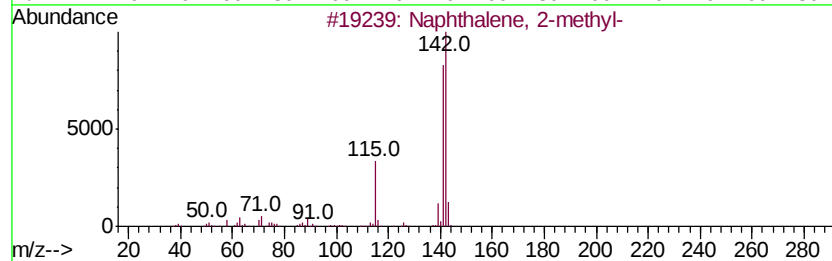
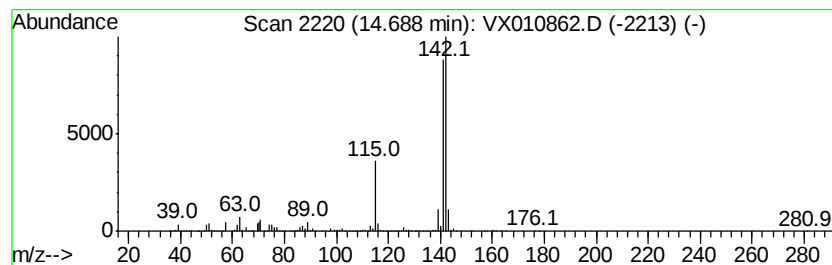
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	53.68 ug/l	1726930	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	94
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\VX071219\
Data File : VX010862.D
Acq On : 13 Jul 2019 01:45
Operator : JC/SP
Sample : K3791-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0595

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	67.0	ug/l	674611	1	5.66	503387	50.0
Benzene, 1-ethyl-...	11.43	160.7	ug/l	5167910	4	12.08	1608480	50.0
Benzene, 1-ethyl-...	11.66	128.8	ug/l	4143820	4	12.08	1608480	50.0
Benzene, 1,2,3-tr...	12.14	147.8	ug/l	4754460	4	12.08	1608480	50.0
o-Cymene	12.66	44.7	ug/l	1436880	4	12.08	1608480	50.0
1-Phenyl-1-butene	12.76	52.4	ug/l	1686100	4	12.08	1608480	50.0
Benzene, 1,2,3,5-...	12.97	44.3	ug/l	1423940	4	12.08	1608480	50.0
Benzene, 1,2,4,5-...	13.02	66.0	ug/l	2121690	4	12.08	1608480	50.0
Naphthalene, 1-me...	14.69	53.7	ug/l	1726930	4	12.08	1608480	50.0