

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
 Data File : VX020737.D
 Acq On : 25 Jan 2021 20:06
 Operator : JC/SP
 Sample : M1192-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 41623

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
 Title : SW846 8260

Signal : TIC: VX020737.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.471	59	63	74	rVB	295845	341978	1.35%	0.306%
2	2.136	158	172	175	rBV	10148275	25262851	100.00%	22.622%
3	5.465	706	718	727	rBV2	206689	595434	2.36%	0.533%
4	5.623	736	744	772	rVB	339698	1030986	4.08%	0.923%
5	6.032	799	811	817	rBV	219064	569920	2.26%	0.510%
6	6.111	817	824	834	rVB	413971	1045857	4.14%	0.937%
7	6.830	932	942	951	rBV	536838	1260953	4.99%	1.129%
8	7.269	1000	1014	1031	rBV	503917	1102328	4.36%	0.987%
9	7.440	1032	1042	1049	rVB	105312	257938	1.02%	0.231%
10	8.696	1242	1248	1253	rVB	1110694	1690722	6.69%	1.514%
11	8.763	1253	1259	1266	rBV	7324801	11353283	44.94%	10.166%
12	10.092	1472	1477	1483	rBV	1094627	1485612	5.88%	1.330%
13	10.232	1495	1500	1510	rBV	1087170	1386317	5.49%	1.241%
14	10.342	1510	1518	1524	rBV	8819527	11773314	46.60%	10.543%
15	10.683	1568	1574	1585	rVB	4368884	5695134	22.54%	5.100%
16	11.116	1640	1645	1651	rBV	881515	1167154	4.62%	1.045%
17	11.342	1673	1682	1688	rBV	208626	269407	1.07%	0.241%
18	11.415	1689	1694	1702	rVV	3327384	5663604	22.42%	5.072%
19	11.488	1702	1706	1717	rVB	1641344	2014108	7.97%	1.804%
20	11.652	1727	1733	1742	rBV	1509021	1865608	7.38%	1.671%
21	11.793	1750	1756	1761	rBV	5807253	7102984	28.12%	6.360%
22	12.006	1783	1791	1795	rBV	241832	321100	1.27%	0.288%
23	12.061	1795	1800	1805	rVB	1191803	1522718	6.03%	1.364%
24	12.128	1805	1811	1816	rBV	1842956	2301679	9.11%	2.061%
25	12.280	1832	1836	1839	rBV	1291495	1644191	6.51%	1.472%
26	12.354	1844	1848	1857	rVB2	977324	1414805	5.60%	1.267%
27	12.500	1868	1872	1878	rVB	378750	463738	1.84%	0.415%
28	12.567	1878	1883	1885	rBV	600498	740220	2.93%	0.663%
29	12.591	1885	1887	1892	rVB	572878	621076	2.46%	0.556%
30	12.652	1892	1897	1900	rBV	876856	1034981	4.10%	0.927%
31	12.738	1906	1911	1919	rBV3	482444	906235	3.59%	0.811%
32	12.884	1930	1935	1940	rVB2	312251	428237	1.70%	0.383%
33	12.957	1940	1947	1951	rBV	519171	706892	2.80%	0.633%
34	13.000	1951	1954	1960	rVB	910563	1059337	4.19%	0.949%
35	13.207	1980	1988	1992	rBV2	659033	936903	3.71%	0.839%
36	13.292	1998	2002	2004	rBV2	173587	256760	1.02%	0.230%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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37	13.329	2004	2008	2013	rVB2	1625452	2149383	8.51%	1.925%
38	13.463	2025	2030	2038	rVB	1073611	1409990	5.58%	1.263%
39	13.585	2046	2050	2052	rBV	202285	264823	1.05%	0.237%
40	13.622	2052	2056	2061	rVV2	286303	472975	1.87%	0.424%
41	13.701	2067	2069	2075	rVB2	297851	397166	1.57%	0.356%
42	13.817	2081	2088	2094	rVB	917896	1205386	4.77%	1.079%
43	13.890	2094	2100	2106	rVB	460124	655214	2.59%	0.587%
44	13.963	2106	2112	2117	rBV2	489753	700581	2.77%	0.627%
45	14.115	2133	2137	2141	rBV	355000	424023	1.68%	0.380%
46	14.268	2156	2162	2166	rBV	1286539	1895223	7.50%	1.697%
47	14.414	2182	2186	2190	rVB	336087	418216	1.66%	0.374%
48	14.518	2198	2203	2209	rBV2	885617	1194213	4.73%	1.069%
49	14.676	2226	2229	2232	rVB	649989	763717	3.02%	0.684%
50	14.713	2232	2235	2240	rVB	248410	319321	1.26%	0.286%
51	14.817	2249	2252	2256	rVV	302477	359487	1.42%	0.322%
52	14.945	2268	2273	2279	rVB2	151176	258680	1.02%	0.232%
53	15.225	2315	2319	2327	rVB2	321162	677666	2.68%	0.607%
54	15.530	2360	2369	2373	rVV3	232729	420573	1.66%	0.377%
55	15.664	2385	2391	2394	rVV	227898	393227	1.56%	0.352%

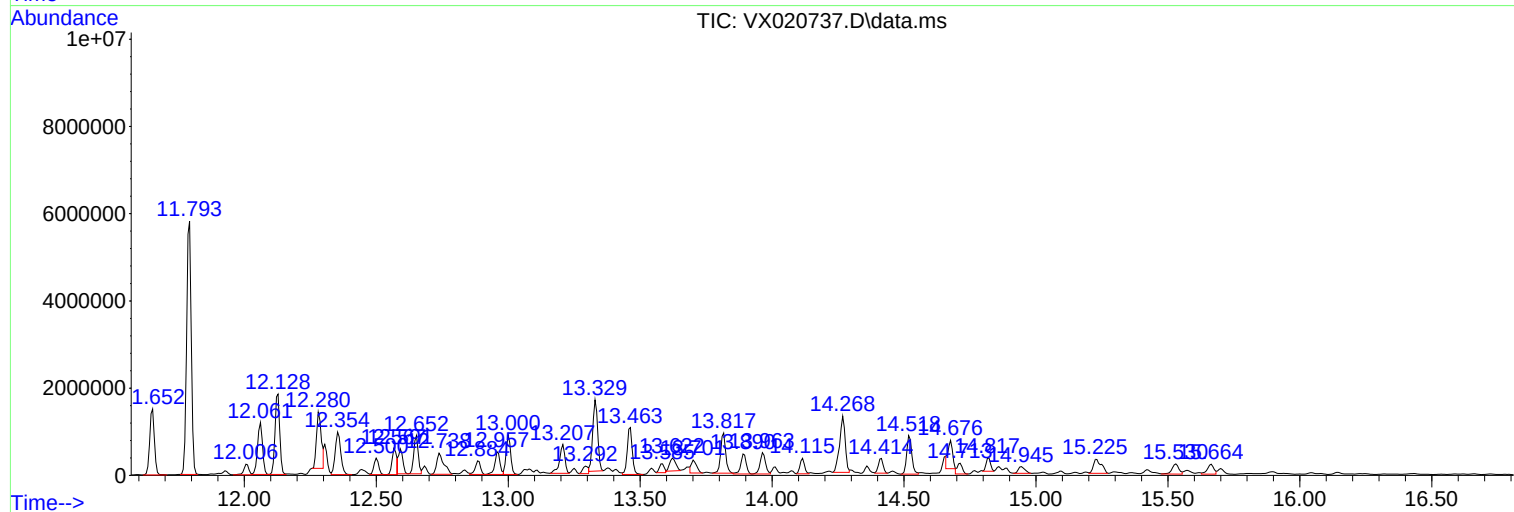
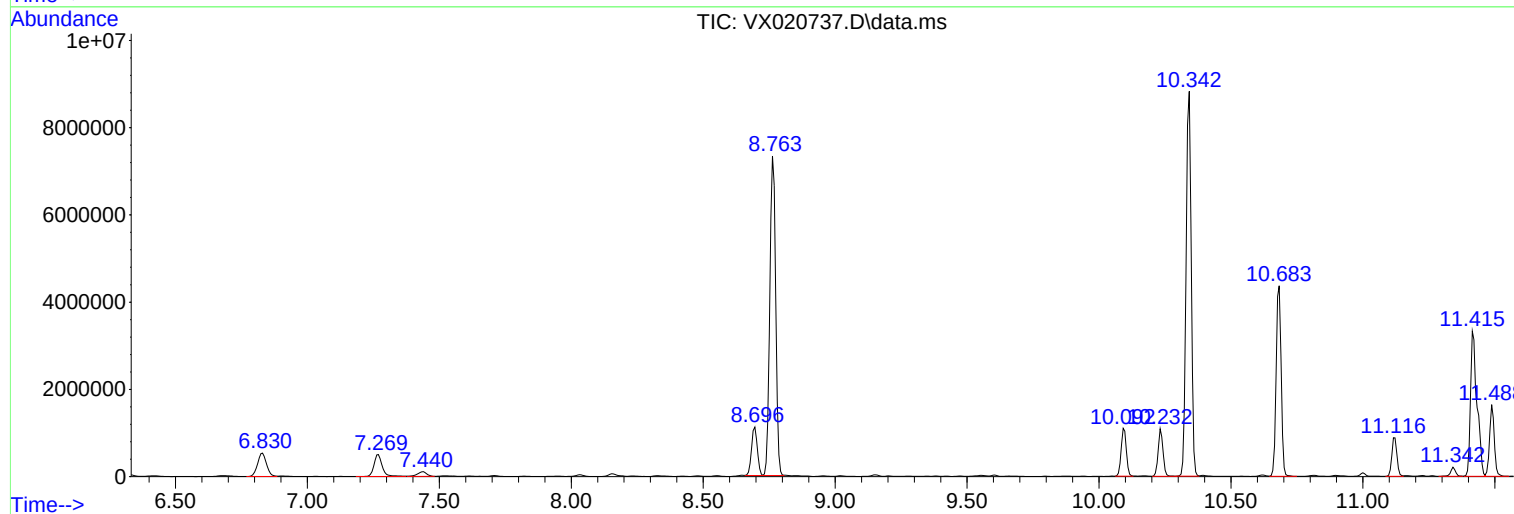
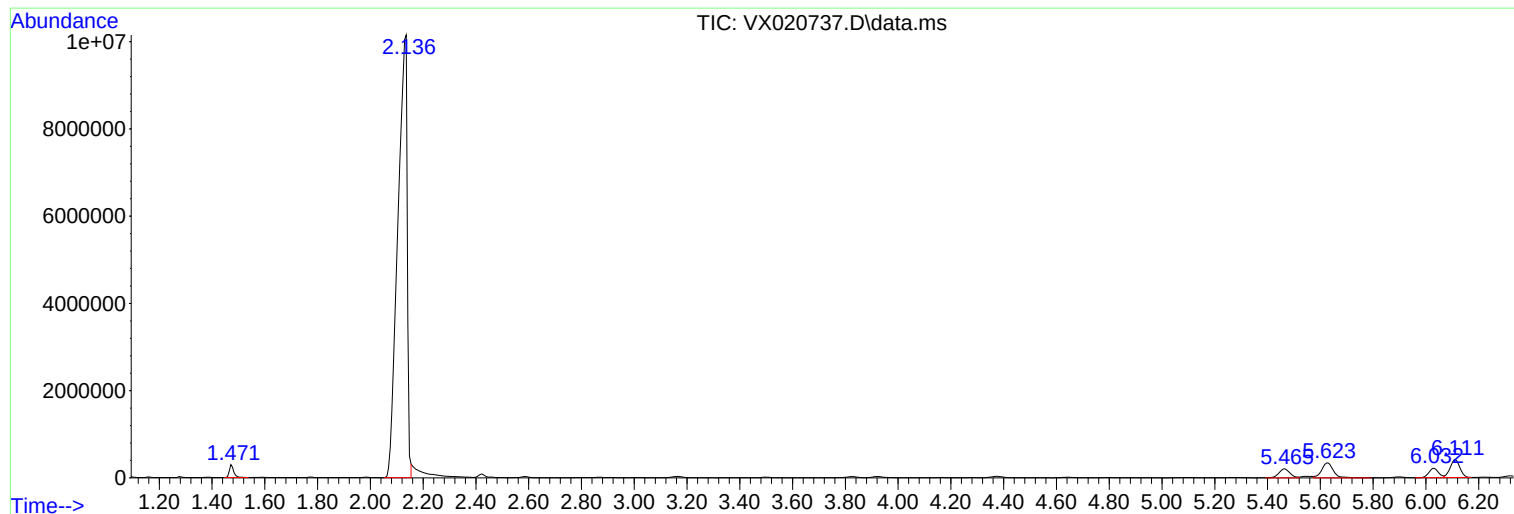
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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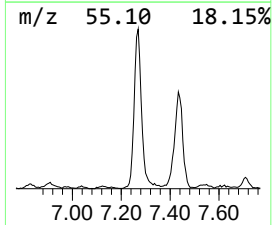
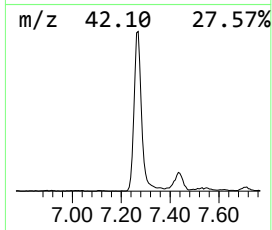
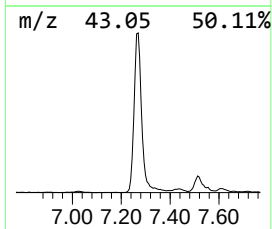
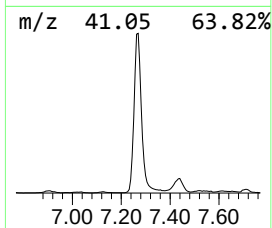
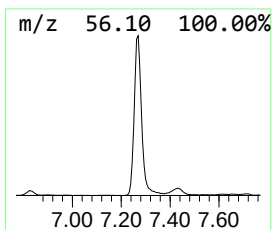
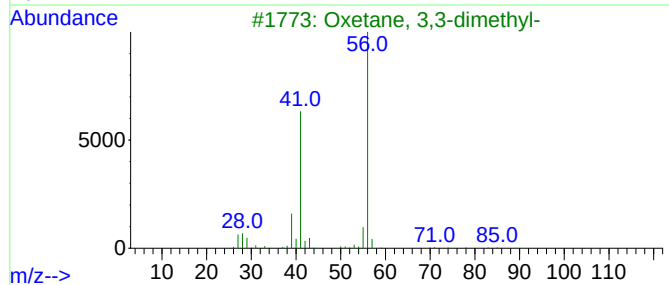
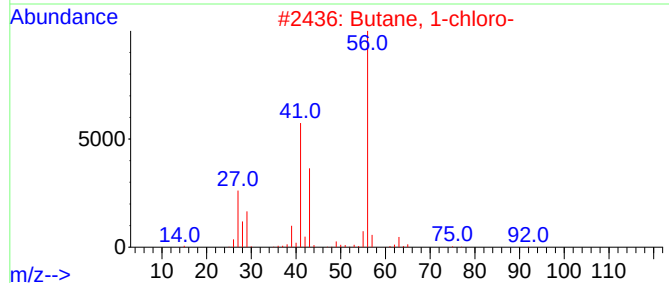
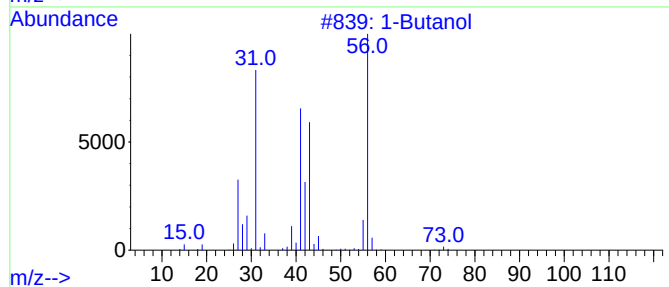
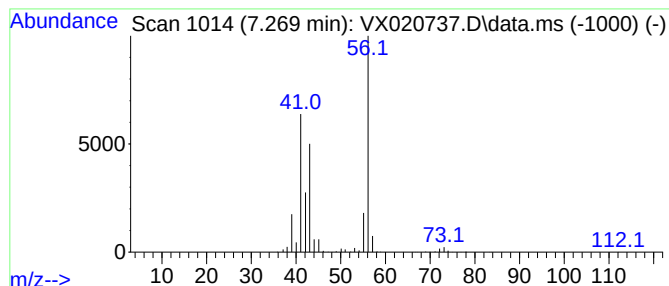
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

Peak Number 1 1-Butanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.269	43.71 ug/l	1102330	1,4-Difluorobenzene	6.830

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butanol	74	C4H10O	000071-36-3	90
2			Butane, 1-chloro-	92	C4H9Cl	000109-69-3	45
3			Oxetane, 3,3-dimethyl-	86	C5H10O	006921-35-3	33
4			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	33
5			3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	33



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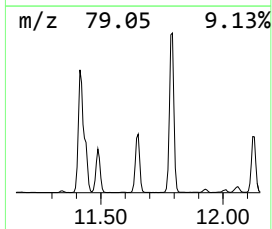
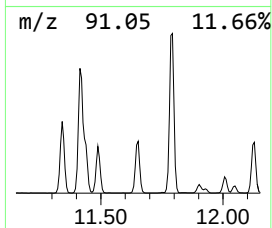
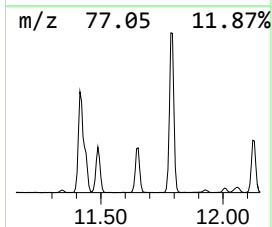
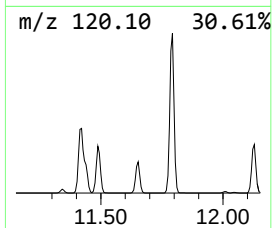
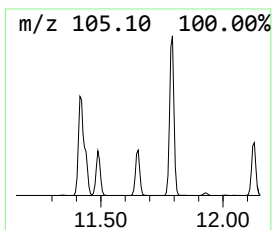
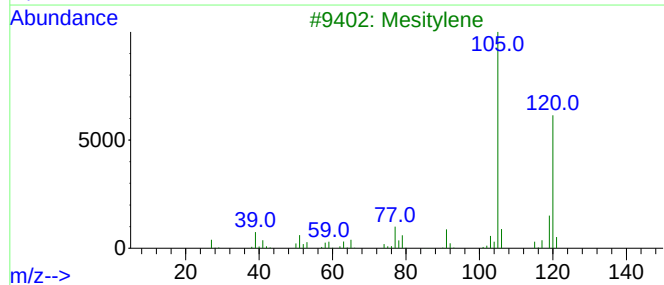
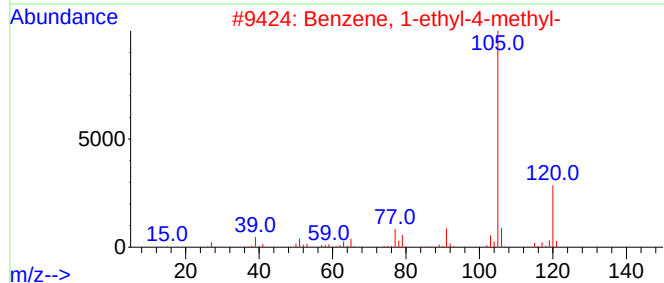
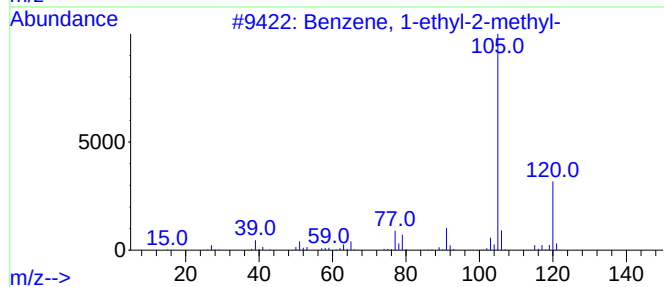
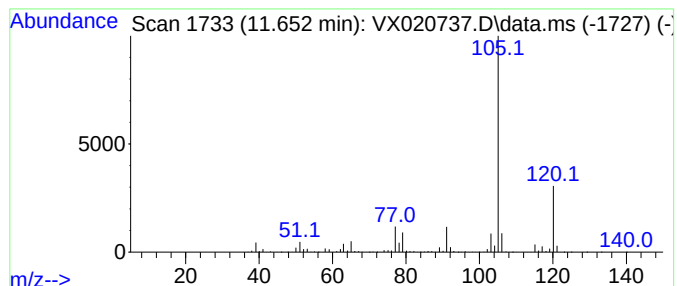
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.652	61.26 ug/l	1865610	1,4-Dichlorobenzene-d4	12.061

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	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-3-methyl-	120	C9H12	000620-14-4	91



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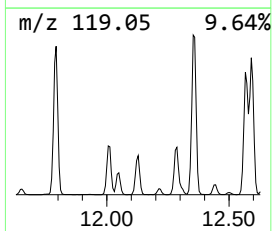
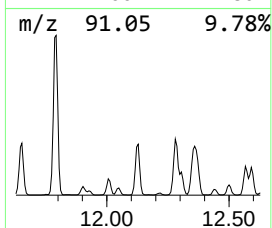
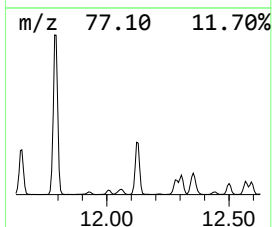
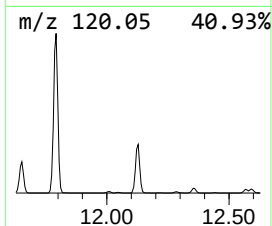
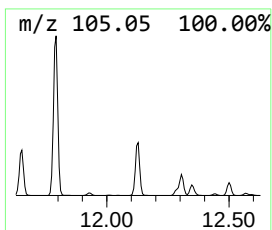
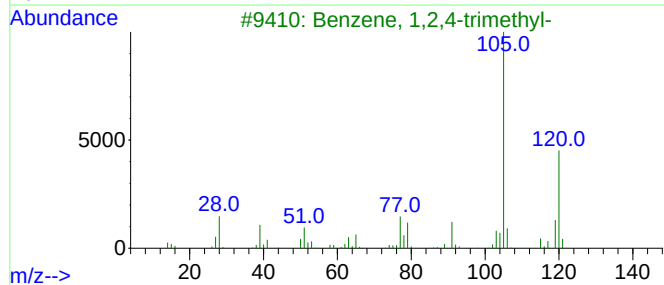
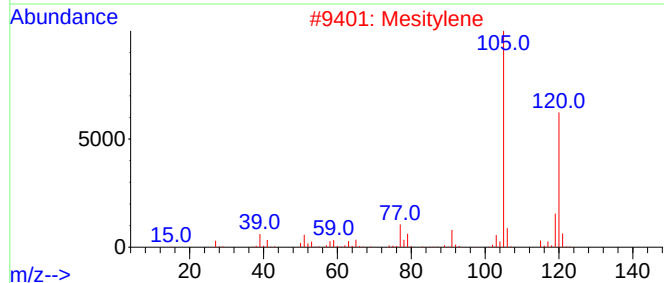
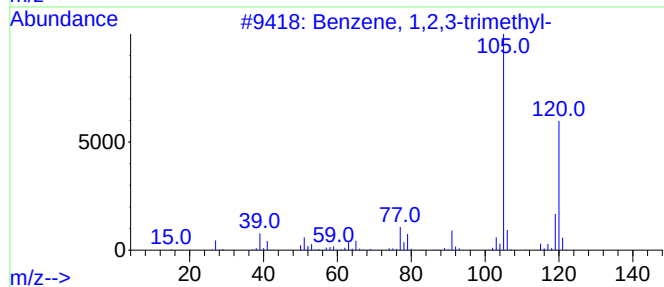
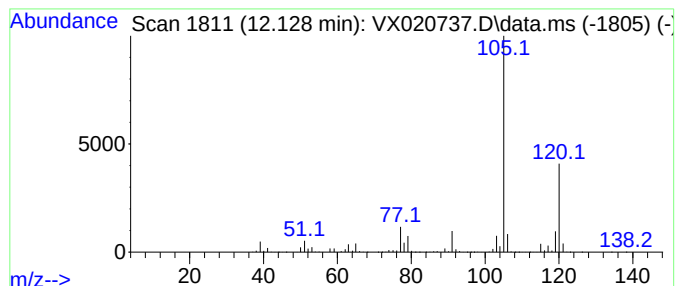
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.128	75.58 ug/l	2301680	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	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	94
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



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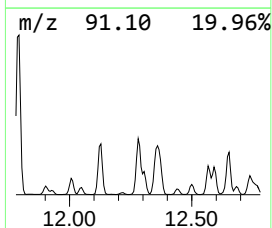
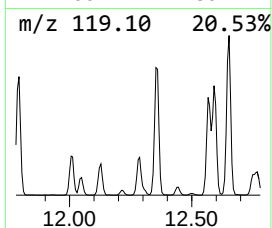
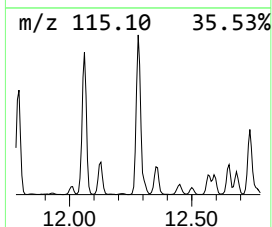
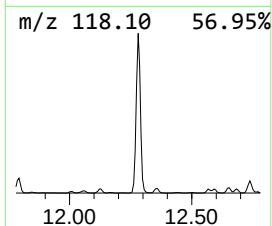
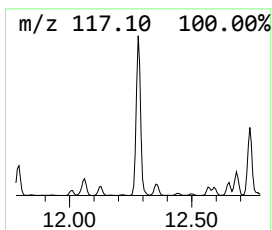
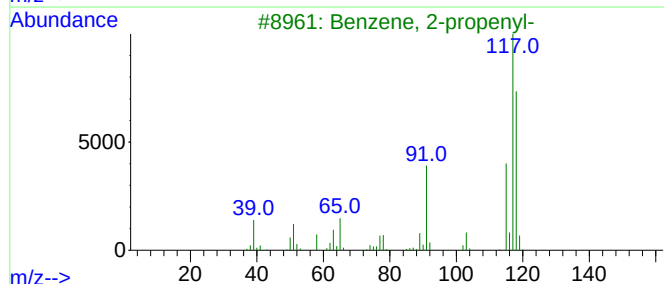
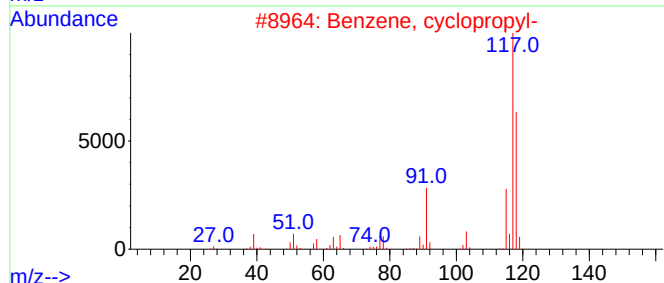
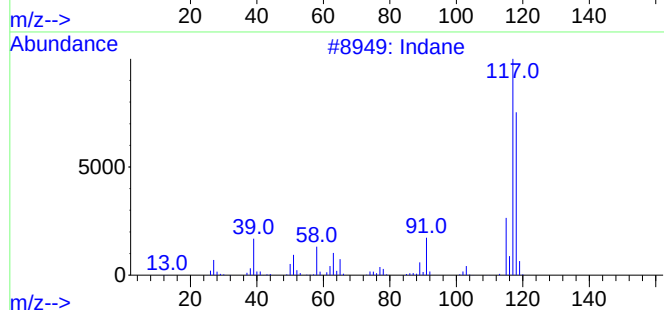
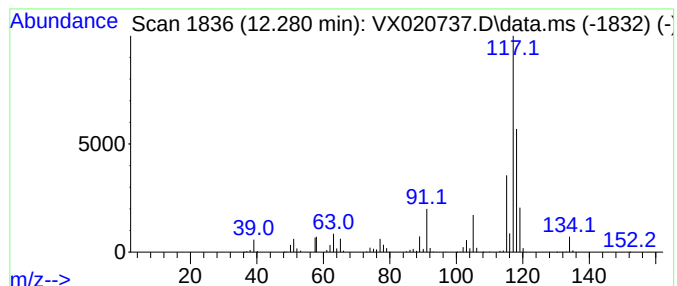
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.280	53.99 ug/l	1644190	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	70
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4	Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	64
5	Deltacyclene	118	C9H10	007785-10-6	52



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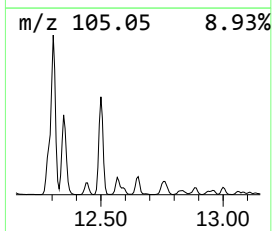
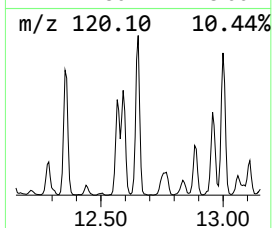
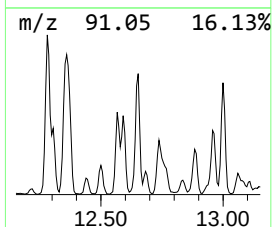
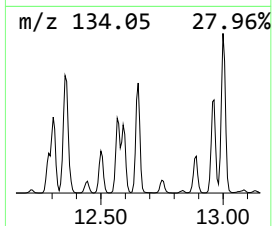
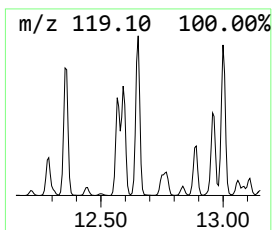
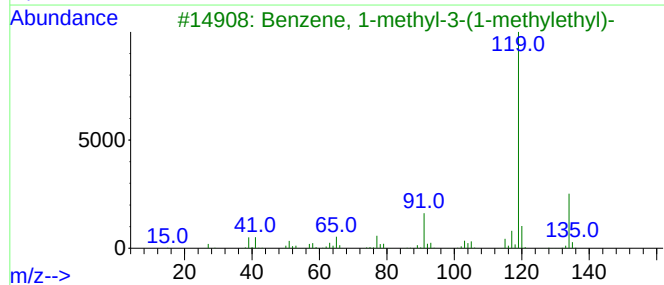
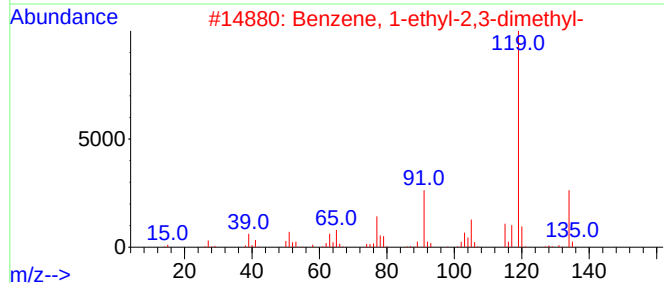
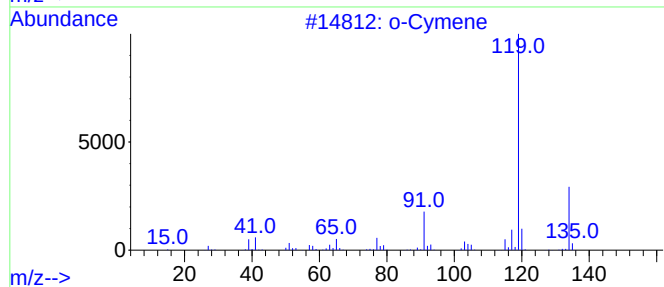
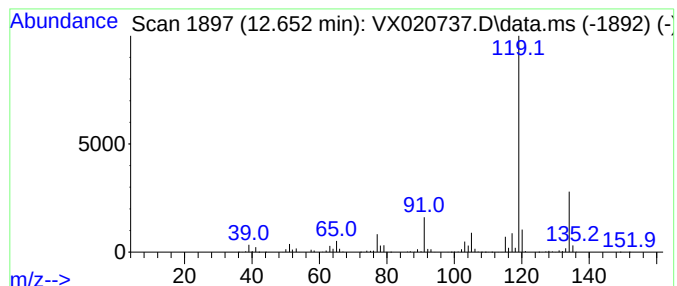
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Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.652	33.98 ug/l	1034980	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5			p-Cymene	134	C10H14	000099-87-6	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
Data File : VX020737.D
Acq On : 25 Jan 2021 20:06
Operator : JC/SP
Sample : M1192-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
41623

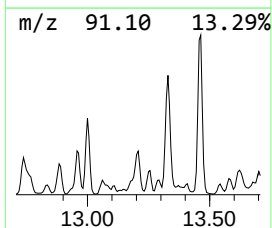
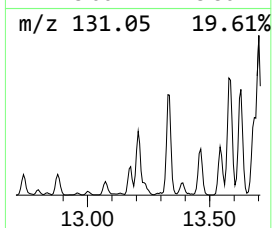
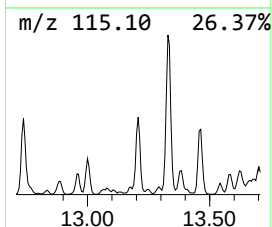
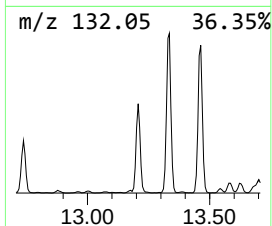
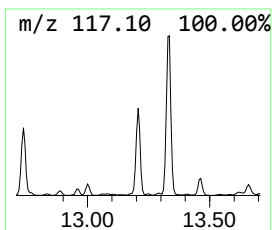
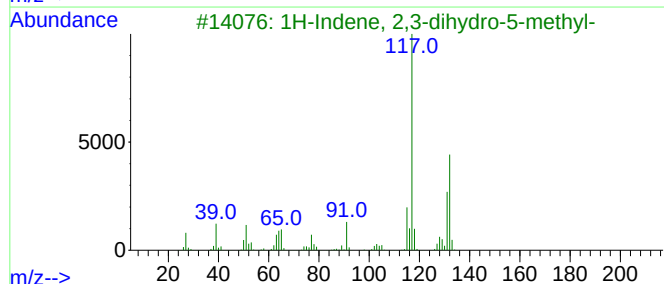
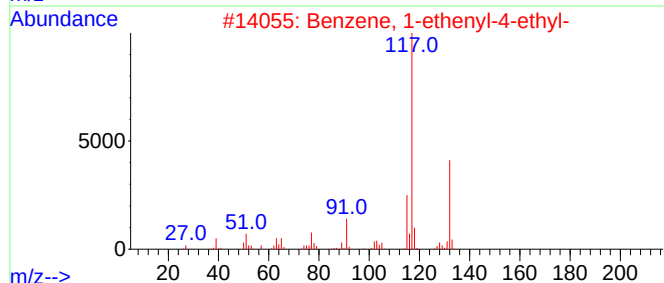
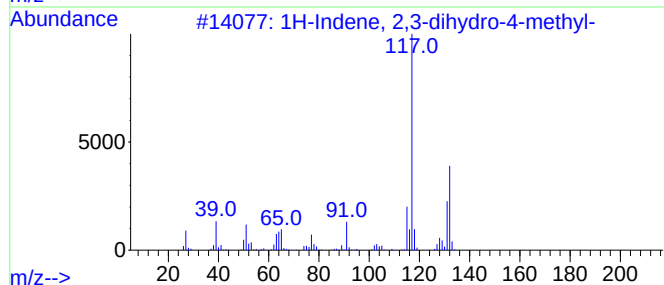
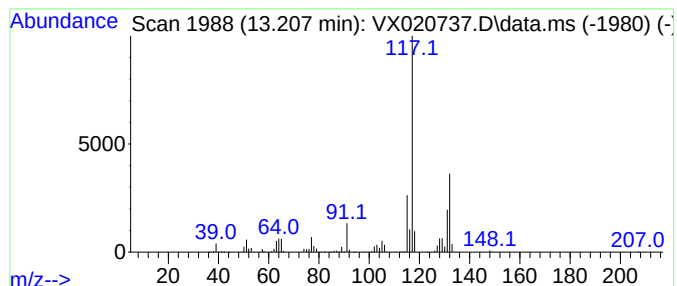
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.207	30.76 ug/l	936903	1,4-Dichlorobenzene-d4			12.061

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
Data File : VX020737.D
Acq On : 25 Jan 2021 20:06
Operator : JC/SP
Sample : M1192-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
41623

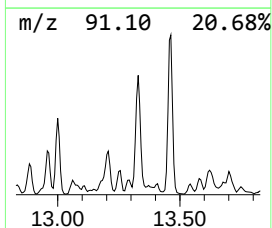
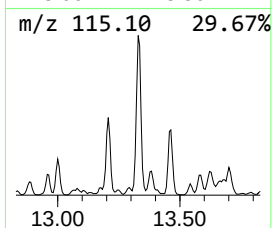
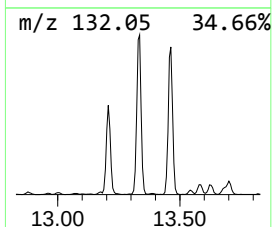
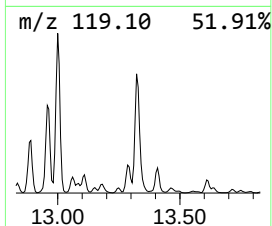
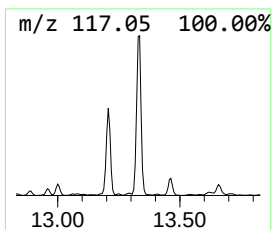
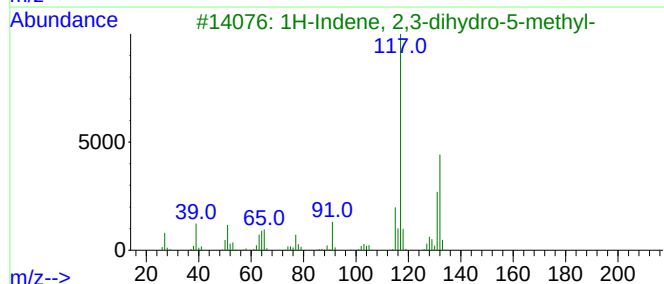
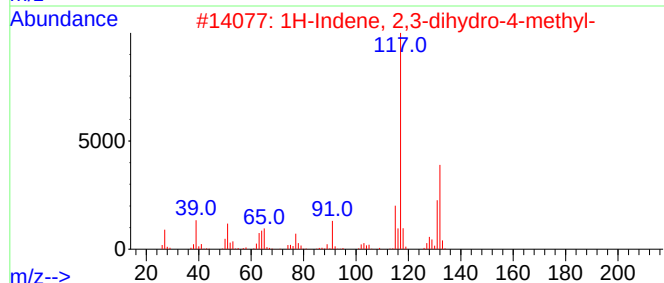
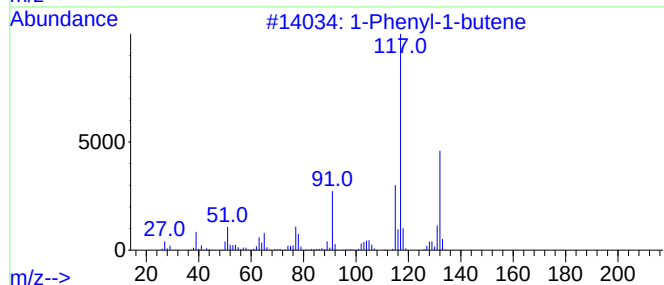
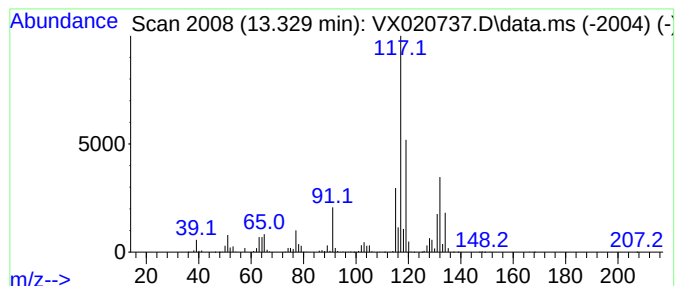
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.329	70.58 ug/l	2149380	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	93
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	92
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	89
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
Data File : VX020737.D
Acq On : 25 Jan 2021 20:06
Operator : JC/SP
Sample : M1192-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
41623

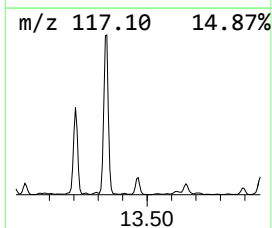
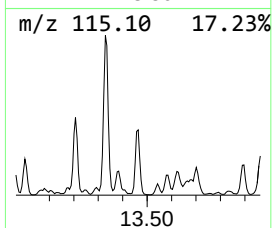
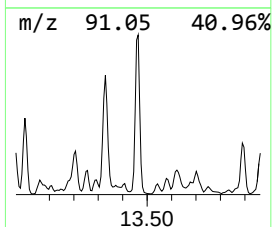
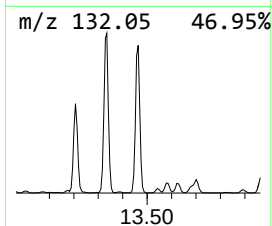
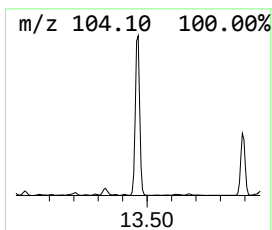
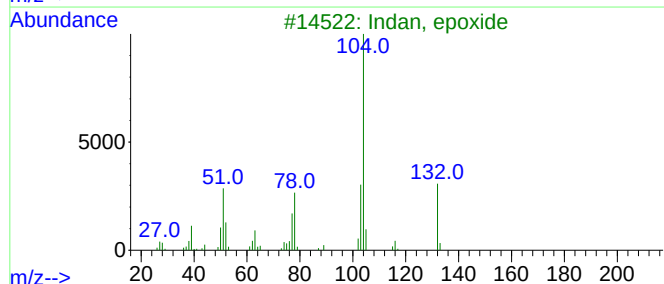
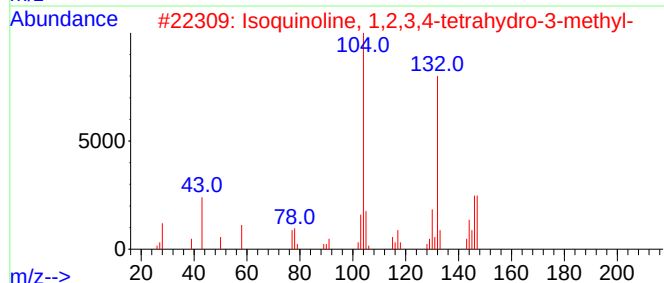
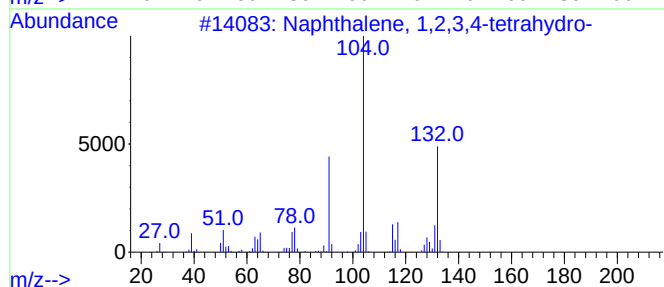
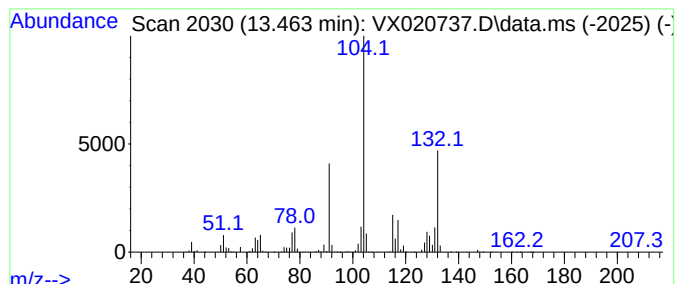
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.463	46.30 ug/l	1409990	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2			Isoquinoline, 1,2,3,4-tetrahydro...	147	C10H13N	029726-60-1	59
3			Indan, epoxide	132	C9H8O	000768-22-9	52
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	40
5			Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
Data File : VX020737.D
Acq On : 25 Jan 2021 20:06
Operator : JC/SP
Sample : M1192-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
41623

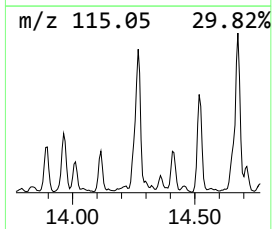
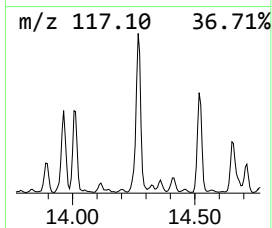
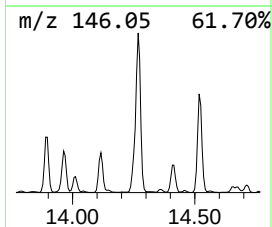
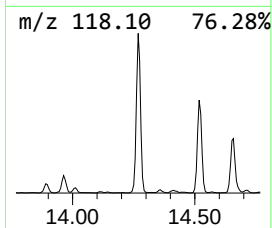
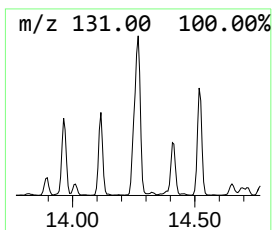
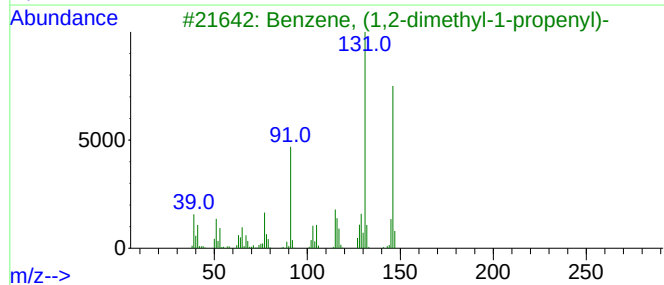
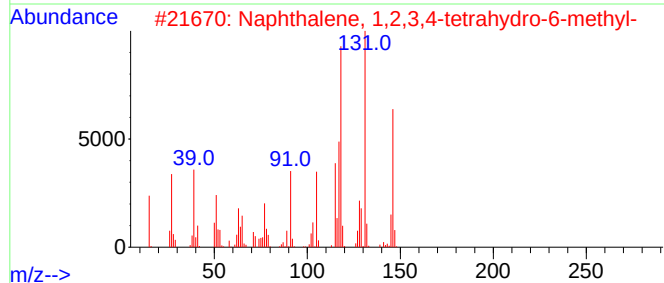
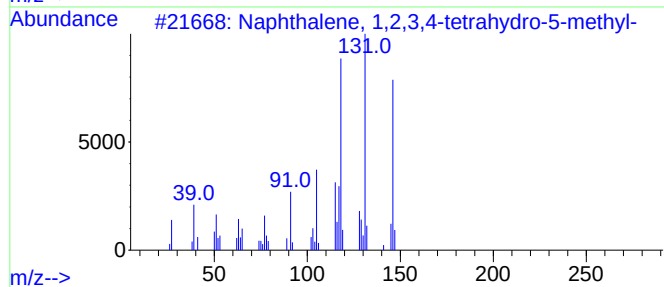
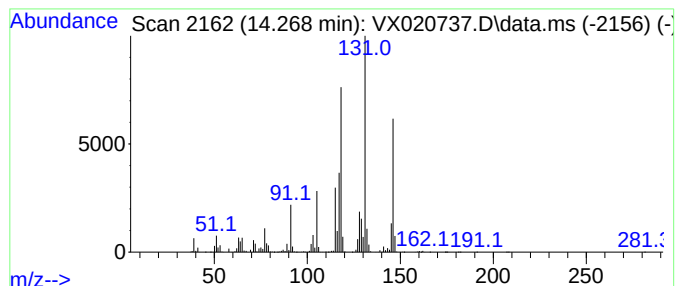
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.268	62.23 ug/l	1895220	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	98
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
Data File : VX020737.D
Acq On : 25 Jan 2021 20:06
Operator : JC/SP
Sample : M1192-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
41623

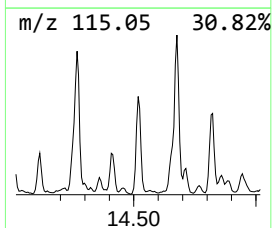
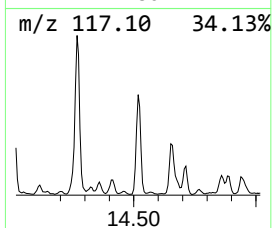
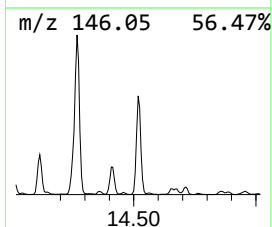
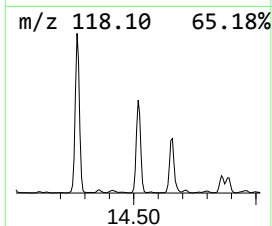
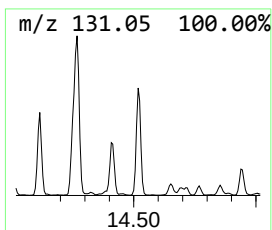
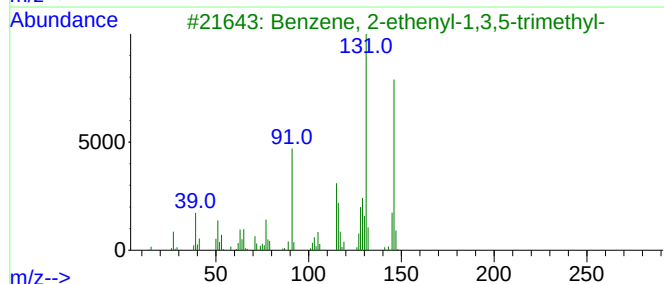
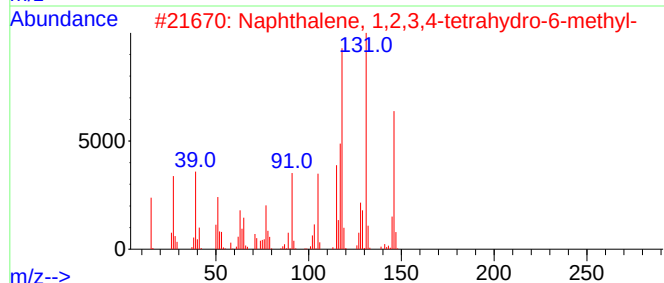
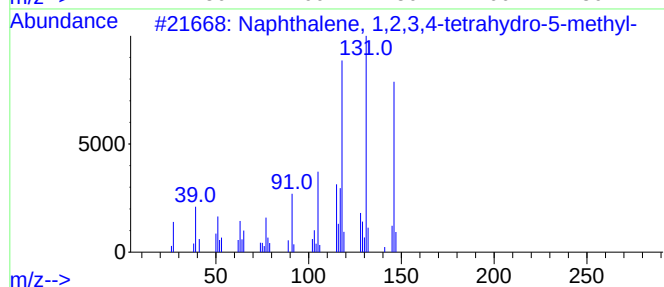
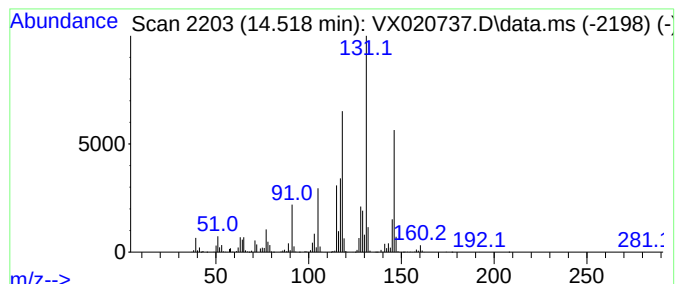
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.518	39.21 ug/l	1194210	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
4			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	76
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012521\
Data File : VX020737.D
Acq On : 25 Jan 2021 20:06
Operator : JC/SP
Sample : M1192-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
41623

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.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
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Benzene, 1-ethy...	11.652	61.3	ug/l	1865610	4	12.061	1522720	50.0
Benzene, 1,2,3-...	12.128	75.6	ug/l	2301680	4	12.061	1522720	50.0
Indane	12.280	54.0	ug/l	1644190	4	12.061	1522720	50.0
o-Cymene	12.652	34.0	ug/l	1034980	4	12.061	1522720	50.0
1H-Indene, 2,3-...	13.207	30.8	ug/l	936903	4	12.061	1522720	50.0
1-Phenyl-1-butene	13.329	70.6	ug/l	2149380	4	12.061	1522720	50.0
Naphthalene, 1,...	13.463	46.3	ug/l	1409990	4	12.061	1522720	50.0
Naphthalene, 1,...	14.268	62.2	ug/l	1895220	4	12.061	1522720	50.0
Naphthalene, 1,...	14.518	39.2	ug/l	1194210	4	12.061	1522720	50.0