

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
 Data File : VN072167.D
 Acq On : 22 Apr 2022 19:24
 Operator : JC\MD
 Sample : N2522-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NWB-1766

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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_N\methods\82N033022W.M
 Title : SW846 8260

Signal : TIC: VN072167.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.240	37	43	65	rVB3	50839	187574	1.58%	0.346%
2	3.693	277	290	317	rBV	4426678	11909415	100.00%	21.948%
3	4.287	381	391	400	rBV	76956	228993	1.92%	0.422%
4	4.569	428	439	455	rBV2	215125	696282	5.85%	1.283%
5	7.645	954	962	983	rBV	91488	252264	2.12%	0.465%
6	8.022	1015	1026	1031	rBV	388896	946905	7.95%	1.745%
7	8.081	1031	1036	1051	rVB	701798	1590755	13.36%	2.932%
8	8.439	1085	1097	1110	rBV2	432646	1217708	10.22%	2.244%
9	8.651	1111	1133	1136	rBV4	33463	184943	1.55%	0.341%
10	8.963	1177	1186	1207	rVB	1058136	2184062	18.34%	4.025%
11	10.327	1410	1418	1426	rBV2	118595	229769	1.93%	0.423%
12	10.439	1426	1437	1442	rVV	1759580	3203515	26.90%	5.904%
13	10.504	1442	1448	1462	rVB	1153031	2137949	17.95%	3.940%
14	10.686	1471	1479	1496	rBV	455622	962150	8.08%	1.773%
15	11.739	1648	1658	1669	rBV	1696757	3046431	25.58%	5.614%
16	11.945	1683	1693	1706	rVB	362850	724436	6.08%	1.335%
17	12.274	1738	1749	1759	rBV	500367	848594	7.13%	1.564%
18	12.727	1819	1826	1836	rBV	1386862	2306279	19.37%	4.250%
19	12.980	1862	1869	1872	rBV	224053	398111	3.34%	0.734%
20	13.010	1872	1874	1878	rVV	152367	217837	1.83%	0.401%
21	13.051	1878	1881	1888	rVB	199440	322037	2.70%	0.593%
22	13.216	1902	1909	1916	rBV	259364	423060	3.55%	0.780%
23	13.363	1928	1934	1941	rBV	395311	624530	5.24%	1.151%
24	13.668	1980	1986	1997	rBV2	1560656	3412073	28.65%	6.288%
25	13.768	1997	2003	2009	rVV	147309	252871	2.12%	0.466%
26	13.833	2009	2014	2015	rVV	104316	139598	1.17%	0.257%
27	13.868	2015	2020	2023	rVV2	444341	799539	6.71%	1.473%
28	13.904	2023	2026	2036	rVB2	306723	562448	4.72%	1.037%
29	14.063	2047	2053	2059	rBV	170717	270647	2.27%	0.499%
30	14.127	2059	2064	2066	rBV	136207	223328	1.88%	0.412%
31	14.157	2066	2069	2073	rVB	156573	212080	1.78%	0.391%
32	14.210	2073	2078	2084	rVB	324413	487556	4.09%	0.899%
33	14.274	2084	2089	2092	rBV	82999	135948	1.14%	0.251%
34	14.321	2092	2097	2103	rVB2	345642	551200	4.63%	1.016%
35	14.451	2114	2119	2125	rVB2	157704	246506	2.07%	0.454%
36	14.533	2125	2133	2136	rBV	217023	369530	3.10%	0.681%

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Max Peaks: 100

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Peak Location: TOP

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

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37	14.574	2136	2140	2144	rVV	359508	539191	4.53%	0.994%
38	14.757	2162	2171	2174	rBV3	92324	182391	1.53%	0.336%
39	14.804	2174	2179	2184	rBV2	337398	563042	4.73%	1.038%
40	14.845	2184	2186	2191	rVB	130656	172135	1.45%	0.317%
41	14.921	2191	2199	2206	rBV3	961198	2206165	18.52%	4.066%
42	14.980	2206	2209	2216	rVB3	73526	146070	1.23%	0.269%
43	15.080	2219	2226	2233	rBV	698645	1221530	10.26%	2.251%
44	15.192	2233	2245	2248	rBV3	242231	586933	4.93%	1.082%
45	15.304	2256	2264	2274	rVB3	305352	760777	6.39%	1.402%
46	15.498	2293	2297	2302	rVV	103116	188518	1.58%	0.347%
47	15.557	2302	2307	2312	rVV	232234	428691	3.60%	0.790%
48	15.633	2312	2320	2340	rVB3	221490	952260	8.00%	1.755%
49	15.798	2340	2348	2356	rBV	208091	432217	3.63%	0.797%
50	16.009	2379	2384	2394	rVB	516039	1125338	9.45%	2.074%
51	16.186	2406	2414	2428	rVB3	156769	484339	4.07%	0.893%
52	16.339	2430	2440	2458	rBV	420570	946431	7.95%	1.744%
53	16.545	2463	2475	2481	rBV3	144833	394613	3.31%	0.727%
54	16.615	2481	2487	2501	rVB3	155505	424348	3.56%	0.782%

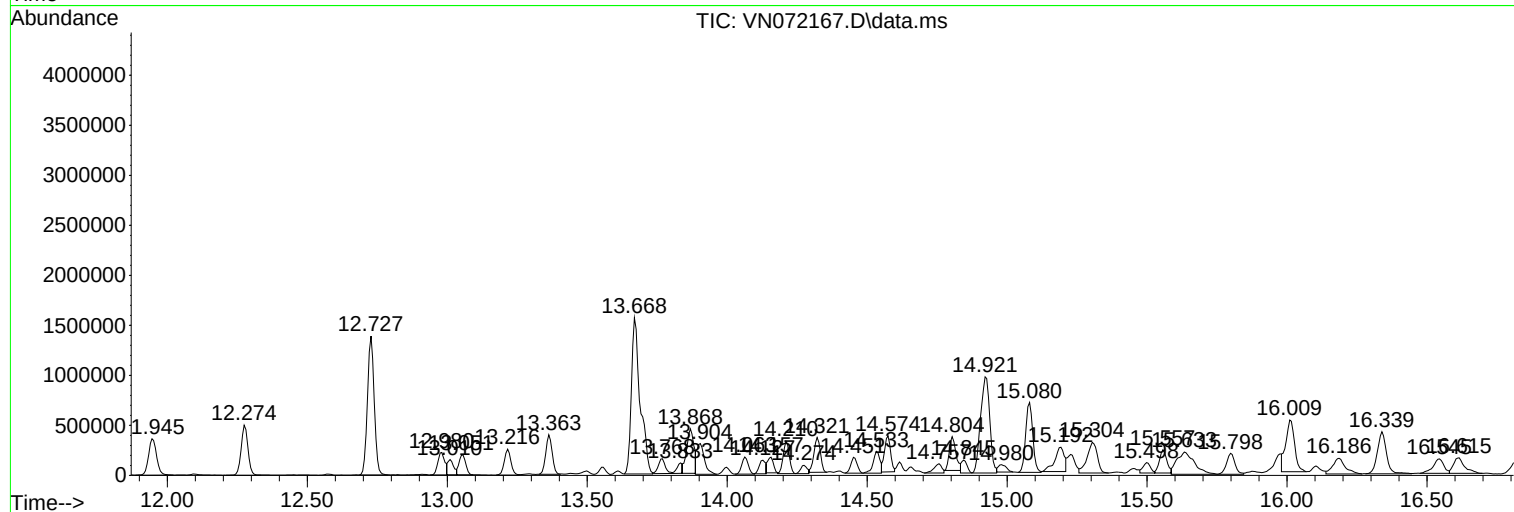
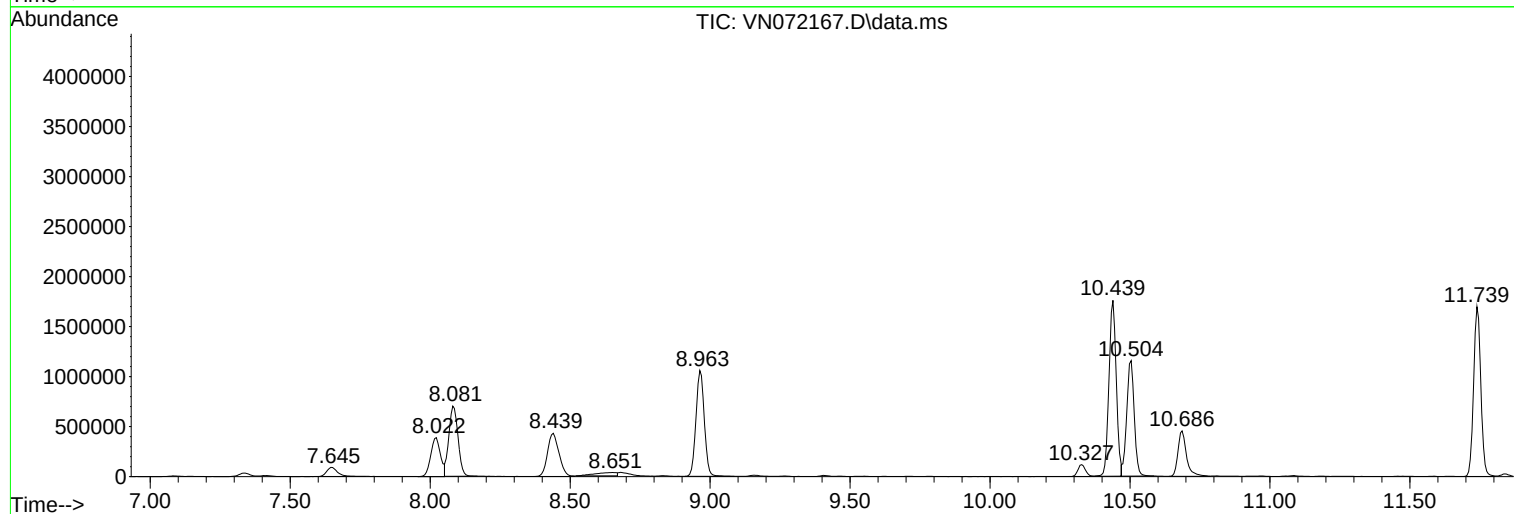
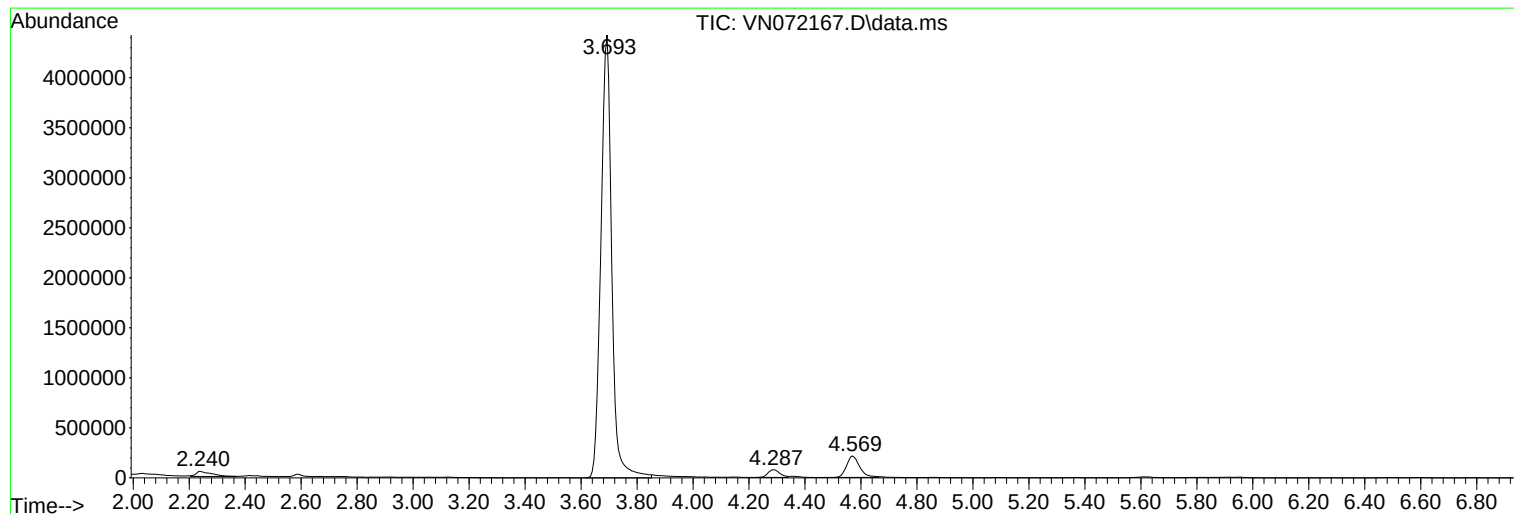
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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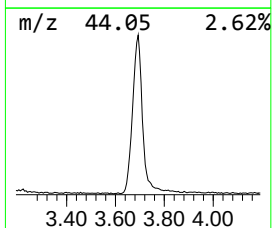
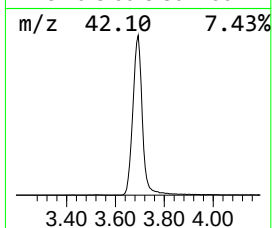
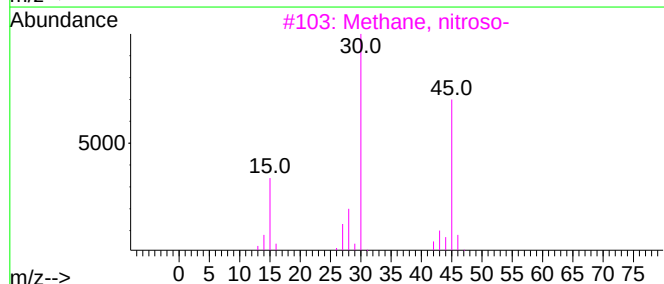
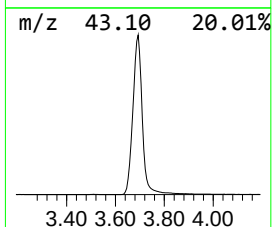
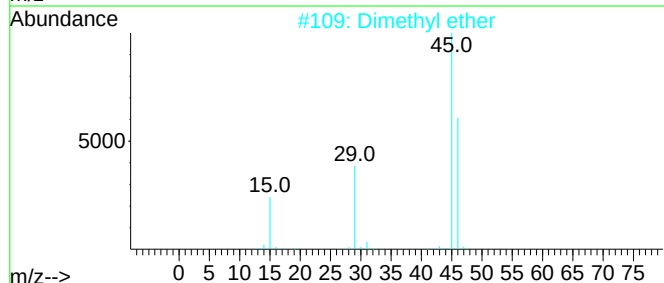
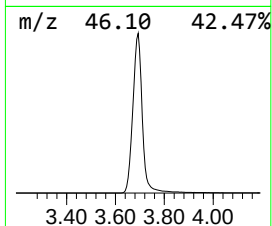
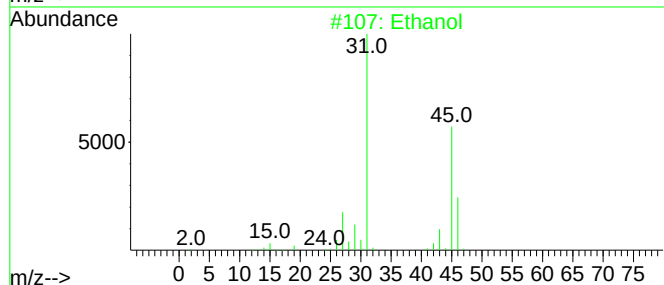
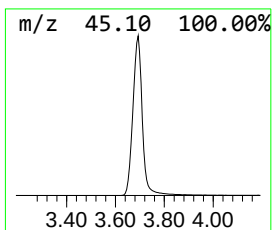
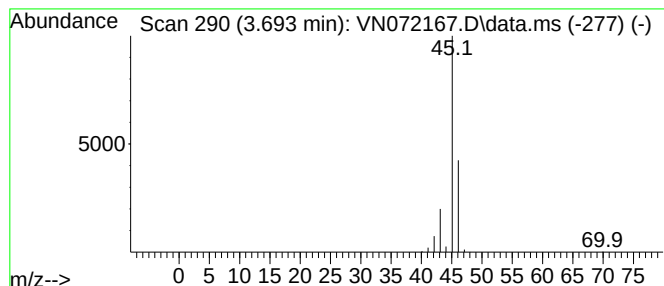
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.693	374.33 ug/l	11909400	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol			46	C2H6O	000064-17-5	83
2	Dimethyl ether			46	C2H6O	000115-10-6	9
3	Methane, nitroso-			45	CH3NO	000865-40-7	4
4	Formic acid			46	CH2O2	000064-18-6	4
5	Oxalic acid			90	C2H2O4	000144-62-7	4



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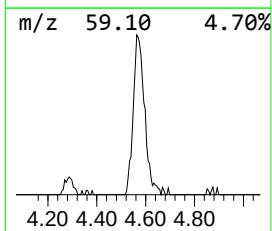
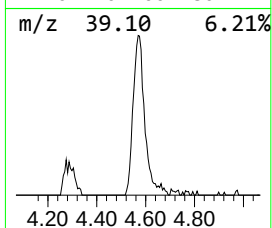
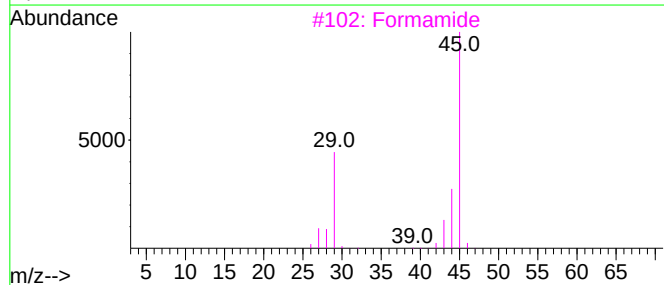
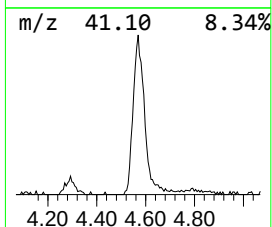
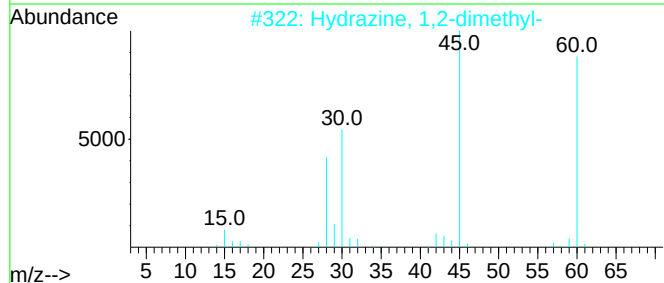
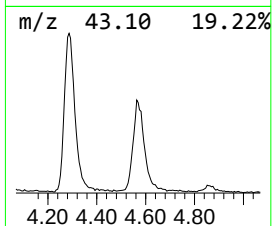
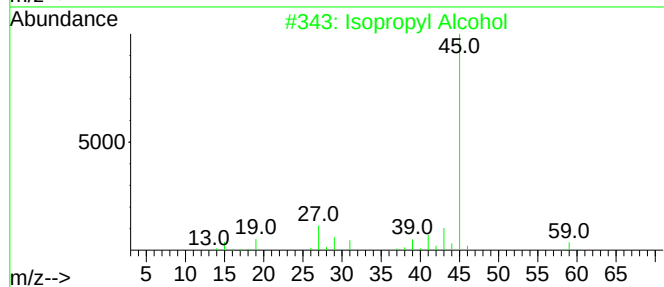
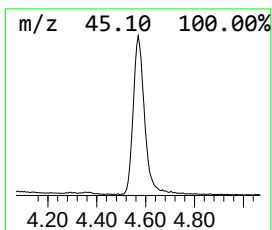
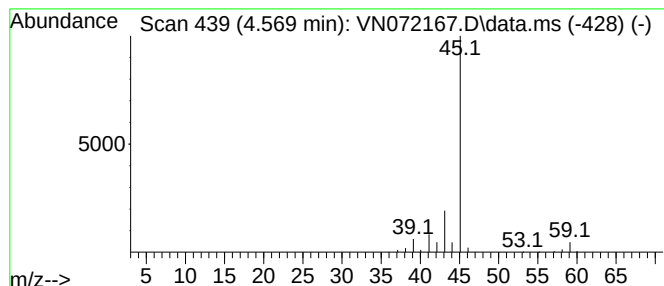
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Isopropyl Alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.569	21.89 ug/l	696282	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	90
2			Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3			Formamide	45	CH3NO	000075-12-7	5
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	4
5			Hydrazine, ethyl-	60	C2H8N2	000624-80-6	4



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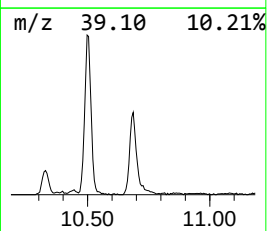
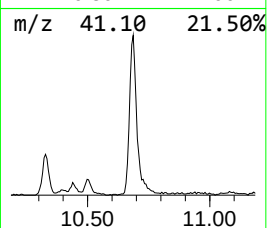
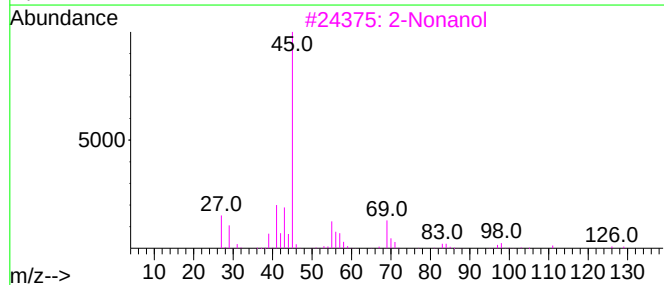
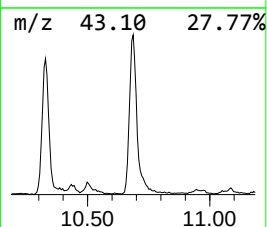
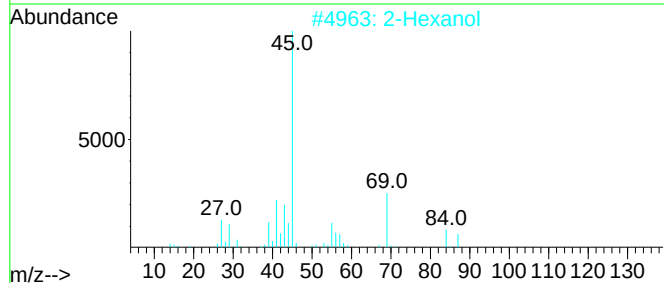
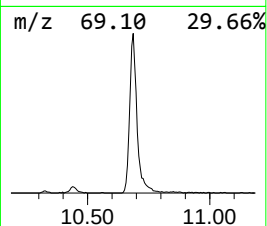
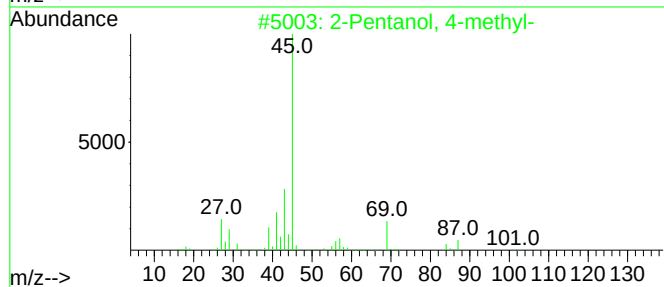
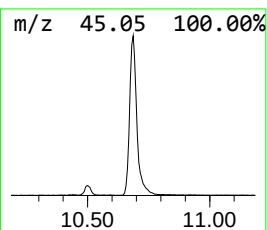
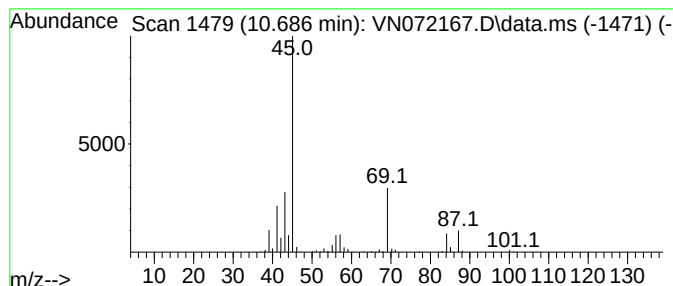
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanol, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.686	15.79 ug/l	962150	Chlorobenzene-d5	11.739

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-			102	C6H14O	000108-11-2	83
2	2-Hexanol			102	C6H14O	000626-93-7	78
3	2-Nonanol			144	C9H20O	000628-99-9	64
4	2-Octanol			130	C8H18O	000123-96-6	64
5	2-Octanol, (S)-			130	C8H18O	006169-06-8	56



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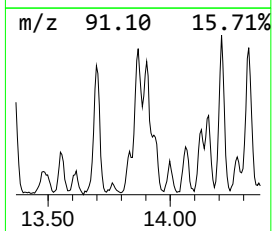
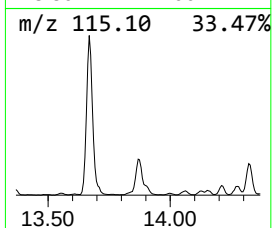
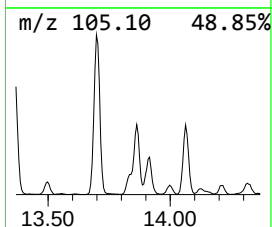
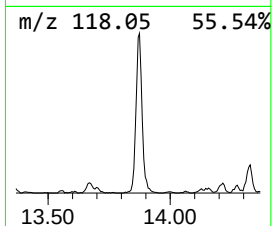
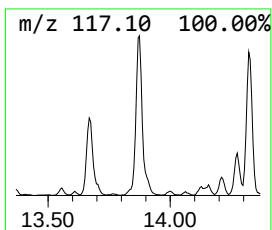
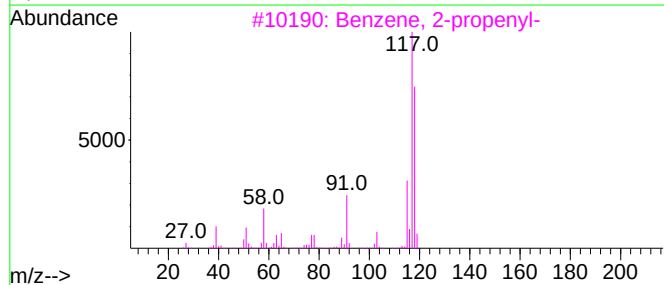
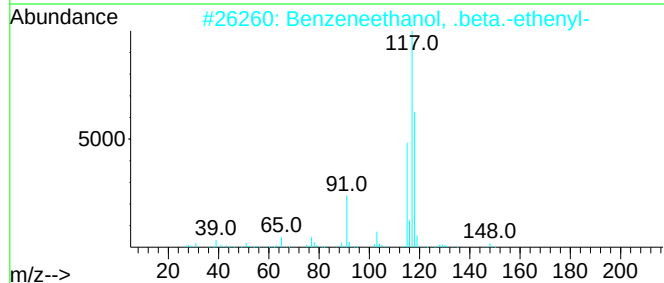
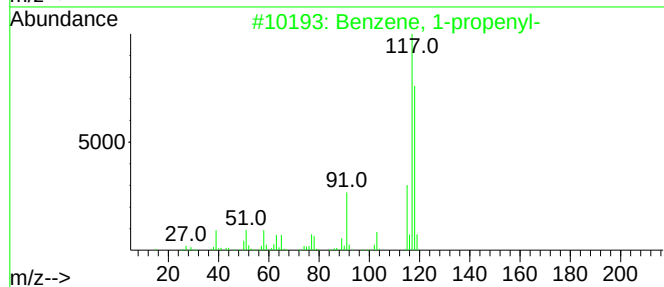
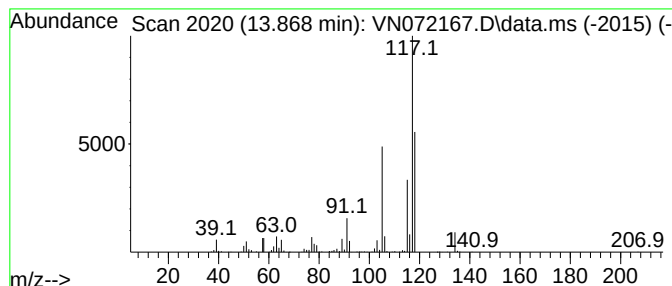
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-propenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.868	11.72 ug/l	799539	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	76
2			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	72
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	68
5			1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	64



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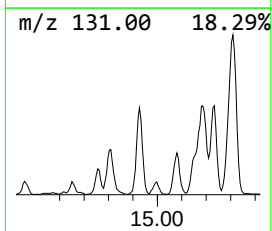
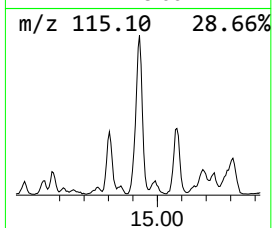
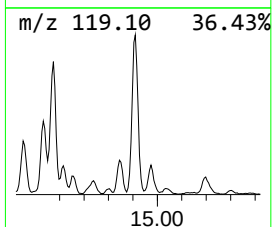
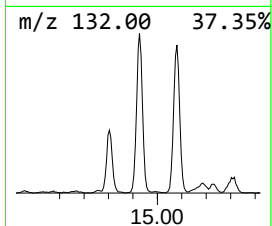
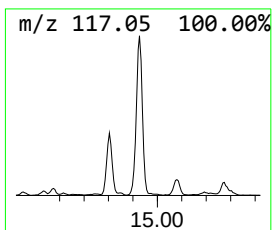
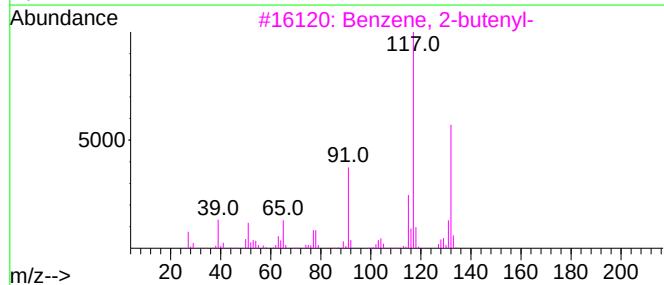
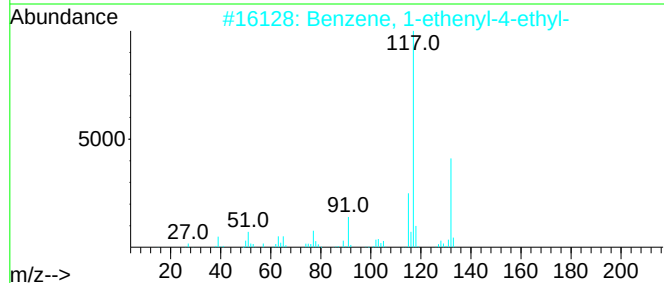
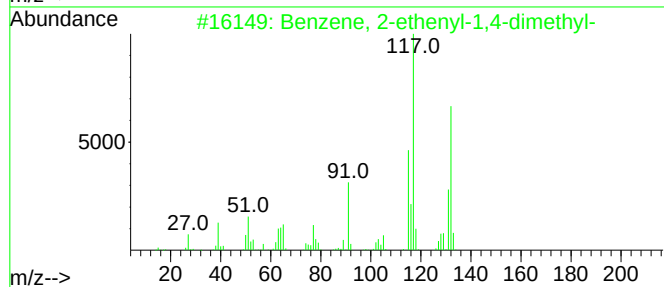
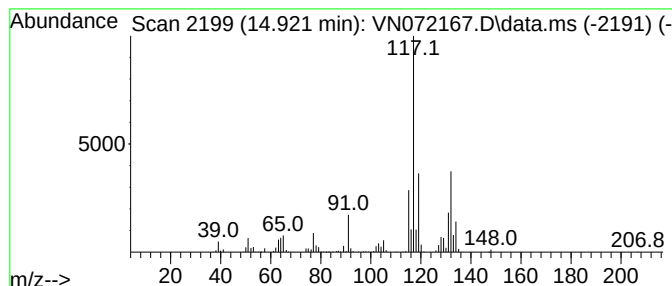
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.921	32.33 ug/l	2206170	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072167.D
Acq On : 22 Apr 2022 19:24
Operator : JC\MD
Sample : N2522-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1766

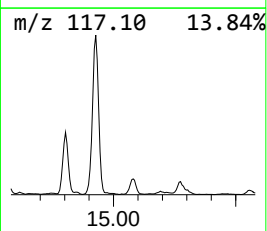
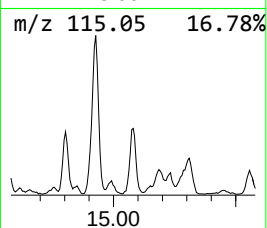
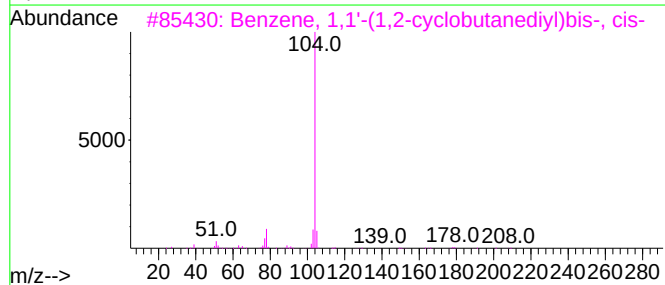
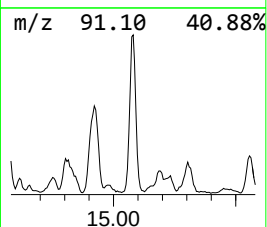
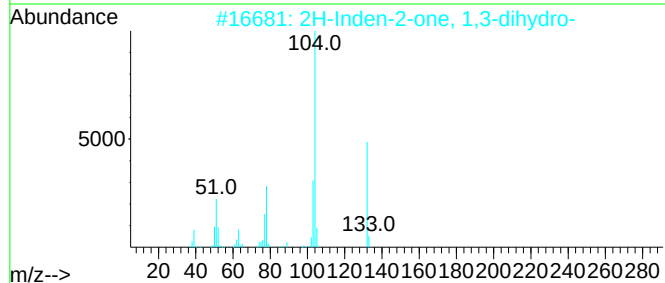
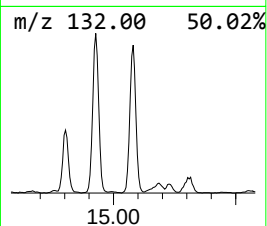
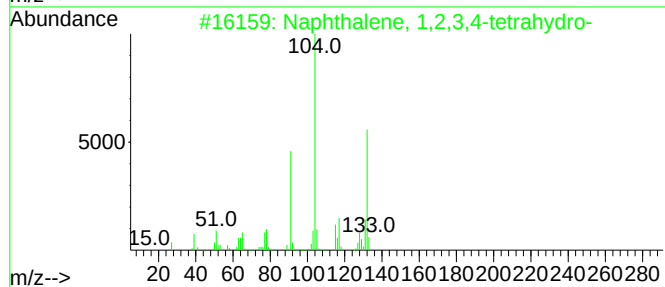
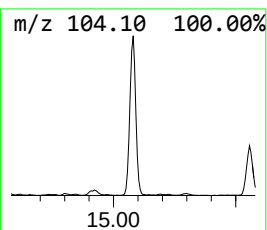
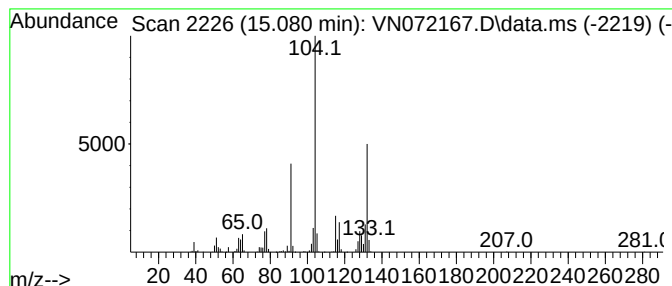
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.080	17.90 ug/l	1221530	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	62
3			Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	43
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	37
5			2-Propenoic acid, 2-phenylethyl ...	176	C11H12O2	003530-36-7	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072167.D
Acq On : 22 Apr 2022 19:24
Operator : JC\MD
Sample : N2522-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1766

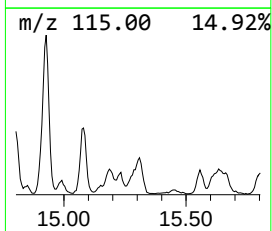
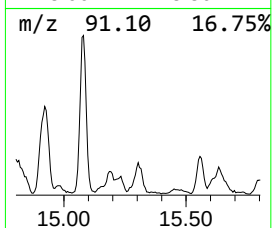
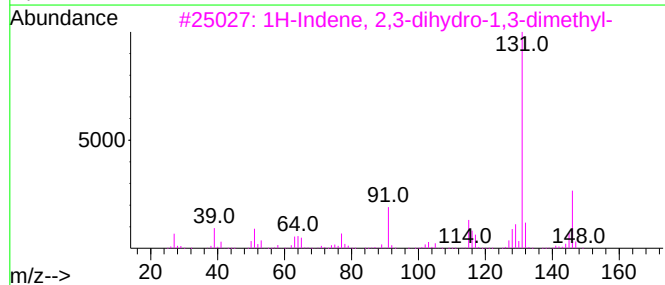
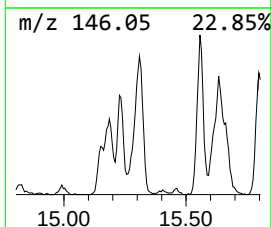
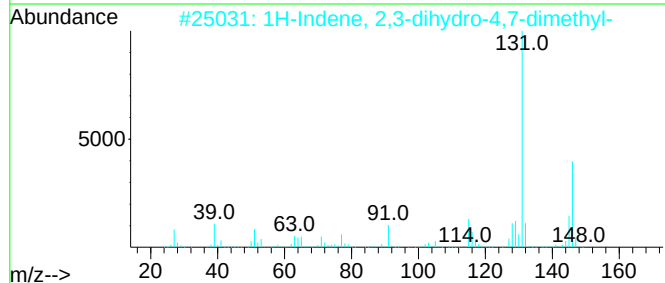
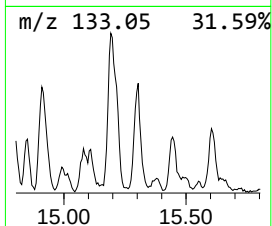
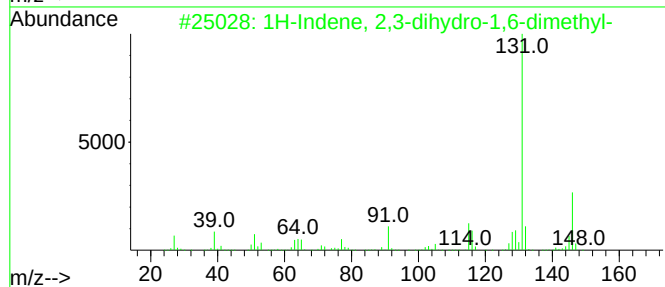
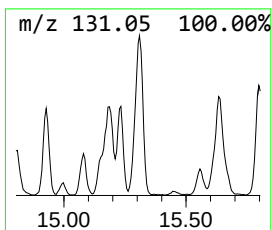
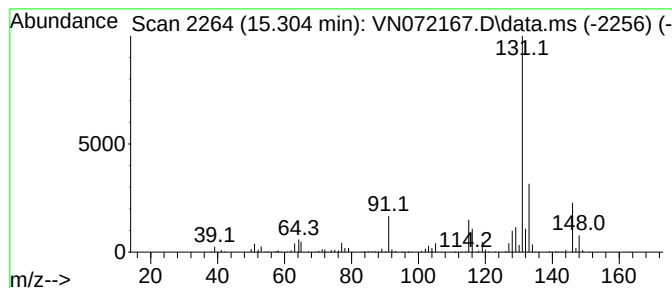
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.304	11.15 ug/l	760777	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
5			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072167.D
Acq On : 22 Apr 2022 19:24
Operator : JC\MD
Sample : N2522-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

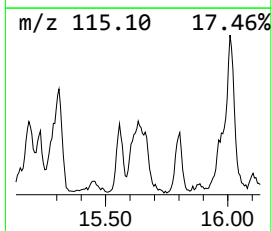
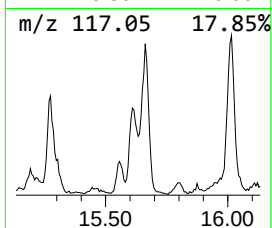
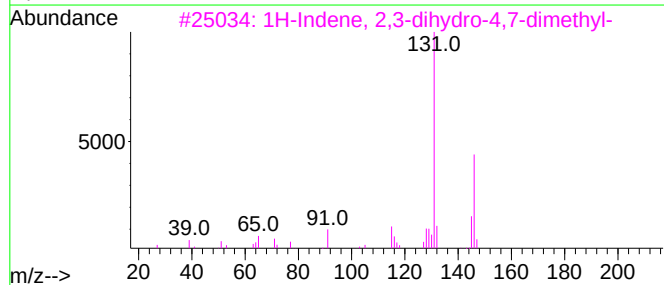
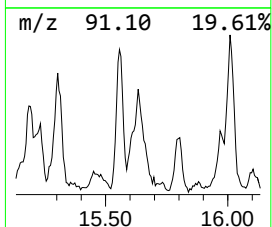
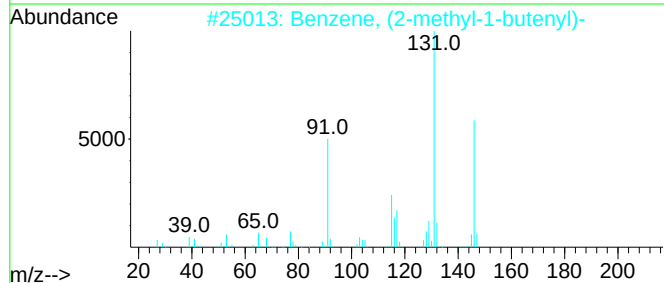
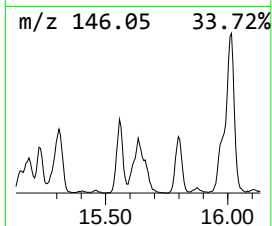
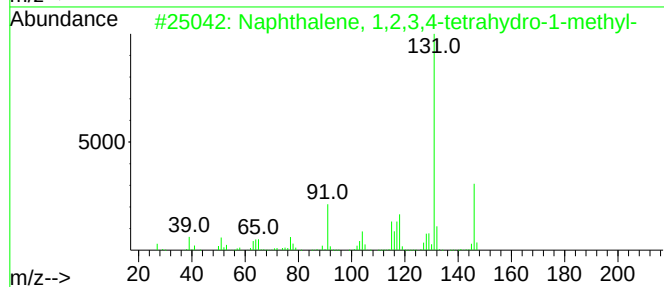
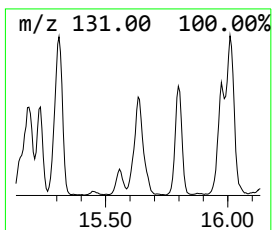
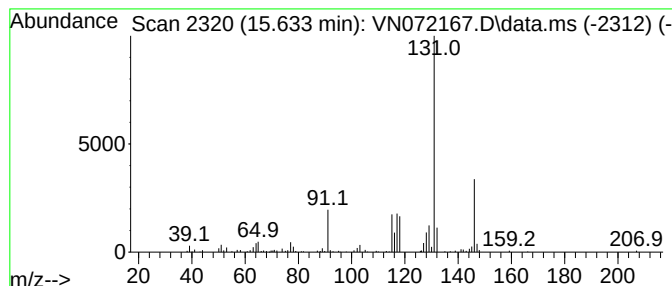
Instrument :
MSVOA_N
ClientSampleId :
NWB-1766

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.633	13.95 ug/l	952260	1,4-Dichlorobenzene-d4		13.668	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001559-81-5	94
2	Benzene, (2-methyl-1-butenyl)-		146	C11H14	056253-64-6	91
3	1H-Indene, 2,3-dihydro-4,7-dimet...		146	C11H14	006682-71-9	90
4	1H-Indene, 2,3-dihydro-1,6-dimet...		146	C11H14	017059-48-2	90
5	2,2-Dimethylindene, 2,3-dihydro-		146	C11H14	020836-11-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072167.D
Acq On : 22 Apr 2022 19:24
Operator : JC\MD
Sample : N2522-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1766

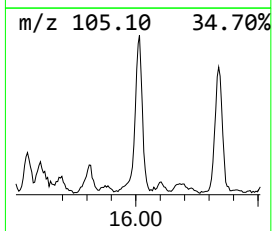
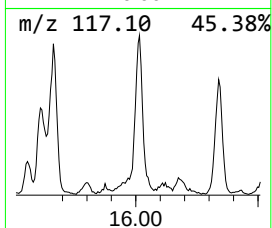
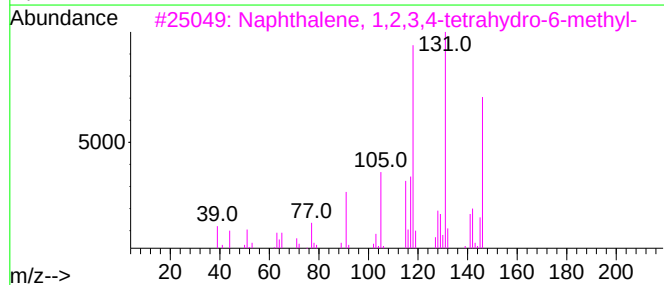
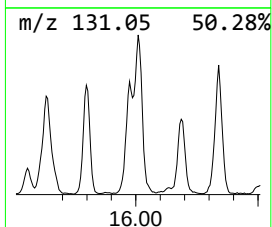
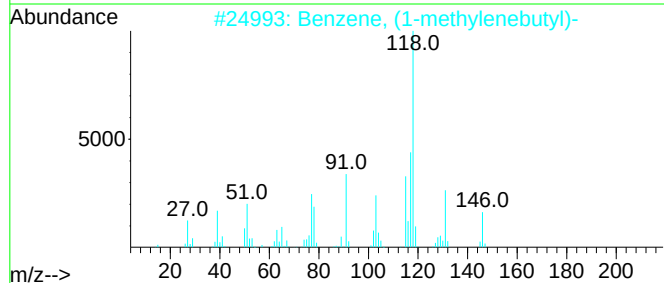
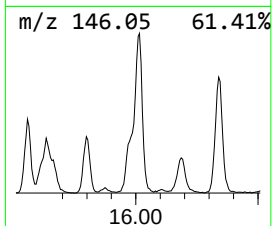
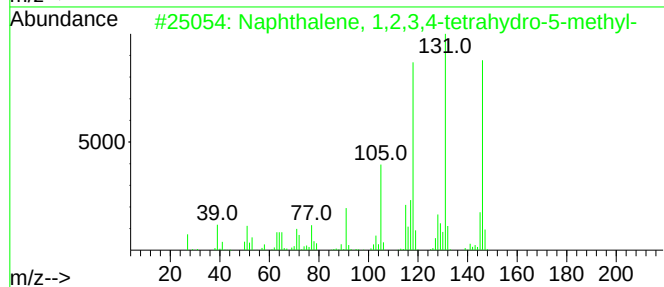
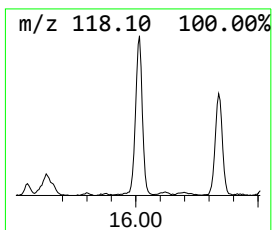
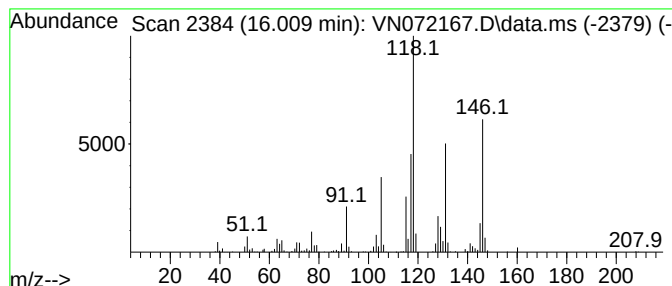
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.009	16.49 ug/l	1125340	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	91
2			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	87
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	76
4			5H-Benzocycloheptene,6,7,8,9-tet...	146	C11H14	001075-16-7	49
5			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072167.D
Acq On : 22 Apr 2022 19:24
Operator : JC\MD
Sample : N2522-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1766

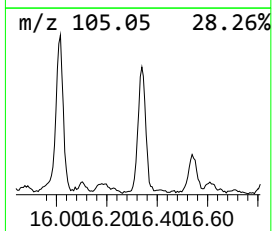
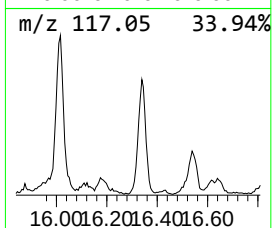
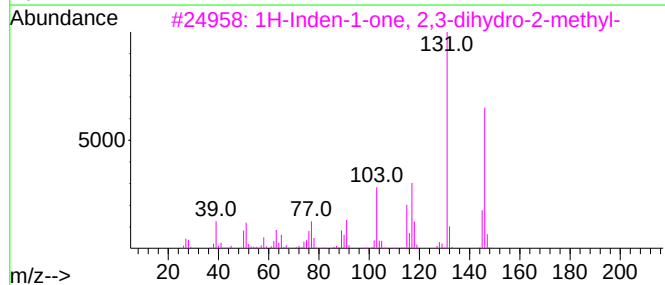
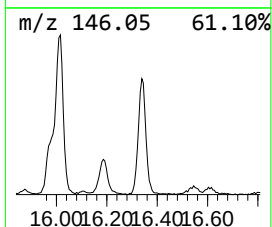
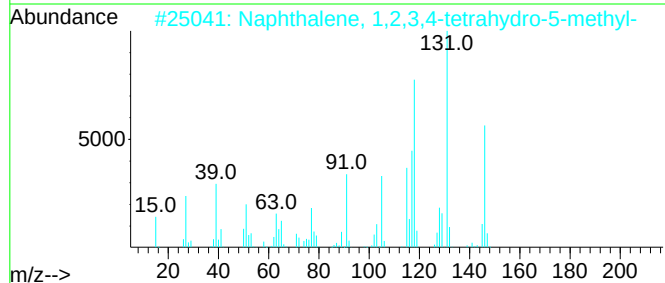
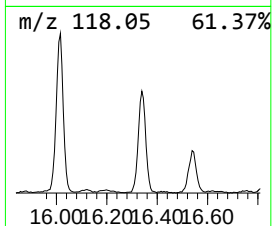
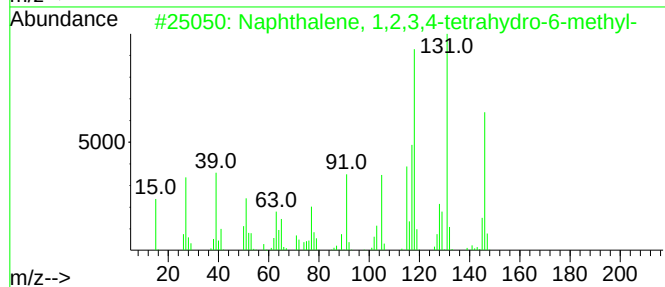
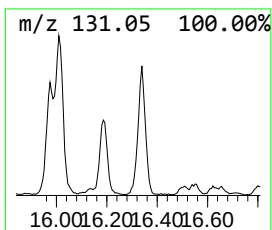
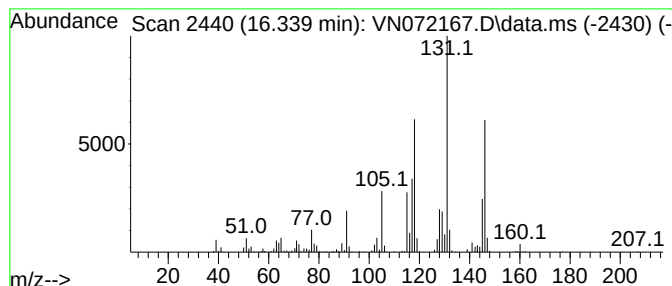
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.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.
16.339	13.87 ug/l	946431	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3			1H-Inden-1-one, 2,3-dihydro-2-methyl-	146	C10H10O	017496-14-9	70
4			1H-1,3,2-Benzodiazaborole, 2-ethyl-	146	C8H11BN2	074663-81-3	68
5			1H-Indene, 2,3-dihydro-4,7-dimethyl-	146	C11H14	006682-71-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072167.D
Acq On : 22 Apr 2022 19:24
Operator : JC\MD
Sample : N2522-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1766

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Isopropyl Alcohol	4.569	21.9	ug/l	696282	1	8.081	1590760	50.0
2-Pentanol, 4-m...	10.686	15.8	ug/l	962150	3	11.739	3046430	50.0
Benzene, 1-prop...	13.868	11.7	ug/l	799539	4	13.668	3412070	50.0
Benzene, 2-ethe...	14.921	32.3	ug/l	2206170	4	13.668	3412070	50.0
Naphthalene, 1,...	15.080	17.9	ug/l	1221530	4	13.668	3412070	50.0
1H-Indene, 2,3-...	15.304	11.2	ug/l	760777	4	13.668	3412070	50.0
Naphthalene, 1,...	15.633	13.9	ug/l	952260	4	13.668	3412070	50.0
Naphthalene, 1,...	16.009	16.5	ug/l	1125340	4	13.668	3412070	50.0
Naphthalene, 1,...	16.339	13.9	ug/l	946431	4	13.668	3412070	50.0