

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
 Data File : VN072165.D
 Acq On : 22 Apr 2022 18:35
 Operator : JC\MD
 Sample : N2522-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NWB-1756

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: VN072165.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.681	276	288	308	rBV	273036	758652	3.65%	0.442%
2	4.281	377	390	428	rBV	6974415	20800593	100.00%	12.130%
3	4.569	430	439	453	rVB	227493	729818	3.51%	0.426%
4	7.334	899	909	925	rBV	228687	628897	3.02%	0.367%
5	8.022	1015	1026	1031	rBV	327285	807138	3.88%	0.471%
6	8.081	1031	1036	1048	rVB	603737	1357601	6.53%	0.792%
7	8.434	1086	1096	1109	rBV2	345753	906011	4.36%	0.528%
8	8.963	1177	1186	1206	rVB	913333	1874774	9.01%	1.093%
9	9.404	1253	1261	1267	rBV	149660	314388	1.51%	0.183%
10	10.328	1410	1418	1426	rBV	429767	778695	3.74%	0.454%
11	10.439	1426	1437	1442	rVV	1562364	2878979	13.84%	1.679%
12	10.504	1442	1448	1461	rVB	4936927	9150651	43.99%	5.336%
13	11.080	1537	1546	1557	rBV2	109969	230584	1.11%	0.134%
14	11.739	1650	1658	1668	rBV	1485603	2653830	12.76%	1.548%
15	11.839	1668	1675	1685	rVV	2521496	4237132	20.37%	2.471%
16	11.945	1685	1693	1713	rVB	7666644	14200160	68.27%	8.281%
17	12.274	1738	1749	1757	rBV	6014668	9973389	47.95%	5.816%
18	12.574	1794	1800	1807	rVB	395304	626468	3.01%	0.365%
19	12.727	1820	1826	1836	rVB	1219758	2014794	9.69%	1.175%
20	12.916	1845	1858	1863	rBV	1057493	1709492	8.22%	0.997%
21	12.980	1863	1869	1872	rVV	4060218	6856618	32.96%	3.999%
22	13.010	1872	1874	1878	rVV	1721331	2461415	11.83%	1.435%
23	13.057	1878	1882	1888	rVB	2013807	3181410	15.29%	1.855%
24	13.216	1897	1909	1916	rBV	2655382	4282805	20.59%	2.498%
25	13.363	1925	1934	1949	rBV	8111027	12777109	61.43%	7.451%
26	13.498	1949	1957	1961	rVB2	224884	412375	1.98%	0.240%
27	13.557	1961	1967	1972	rBV	408959	622822	2.99%	0.363%
28	13.610	1972	1976	1980	rBV	172206	245535	1.18%	0.143%
29	13.698	1987	1991	1998	rVB	3528167	6020142	28.94%	3.511%
30	13.868	2008	2020	2025	rBV2	4340958	8973776	43.14%	5.233%
31	13.998	2036	2042	2047	rBV3	301298	502508	2.42%	0.293%
32	14.063	2047	2053	2058	rBV	647770	1023099	4.92%	0.597%
33	14.127	2058	2064	2066	rVV	925214	1447011	6.96%	0.844%
34	14.157	2066	2069	2073	rVV	854427	1237423	5.95%	0.722%
35	14.210	2073	2078	2084	rVB	1301448	1907750	9.17%	1.113%
36	14.274	2084	2089	2092	rBV	530508	817905	3.93%	0.477%

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37	14.321	2092	2097	2104	rVB	1652165	2676334	12.87%	1.561%
38	14.457	2114	2120	2125	rBV2	520284	833050	4.00%	0.486%
39	14.533	2125	2133	2136	rBV	720906	1165828	5.60%	0.680%
40	14.574	2136	2140	2144	rVB	1074483	1525146	7.33%	0.889%
41	14.616	2144	2147	2151	rVB	220993	263852	1.27%	0.154%
42	14.757	2162	2171	2174	rBV2	292211	543037	2.61%	0.317%
43	14.804	2174	2179	2184	rBV	2009620	3239806	15.58%	1.889%
44	14.927	2191	2200	2206	rBV2	4115416	8279385	39.80%	4.828%
45	14.980	2206	2209	2218	rVB2	199085	415910	2.00%	0.243%
46	15.080	2218	2226	2233	rBV	3767787	6463568	31.07%	3.769%
47	15.192	2239	2245	2248	rBV2	391839	734008	3.53%	0.428%
48	15.227	2248	2251	2256	rVB	333807	523297	2.52%	0.305%
49	15.310	2256	2265	2274	rVB3	690751	1759706	8.46%	1.026%
50	15.498	2291	2297	2302	rVV	608916	1030372	4.95%	0.601%
51	15.557	2302	2307	2312	rVV	517009	826042	3.97%	0.482%
52	15.633	2312	2320	2340	rVB3	498996	2136189	10.27%	1.246%
53	15.798	2340	2348	2356	rBV	532484	1069791	5.14%	0.624%
54	16.010	2368	2384	2395	rBV	1397957	3747011	18.01%	2.185%
55	16.186	2406	2414	2429	rVB3	346012	1010401	4.86%	0.589%
56	16.339	2430	2440	2457	rBV	897542	2001680	9.62%	1.167%
57	16.539	2463	2474	2480	rBV2	270426	700841	3.37%	0.409%
58	16.609	2480	2486	2501	rVB	455300	1131474	5.44%	0.660%

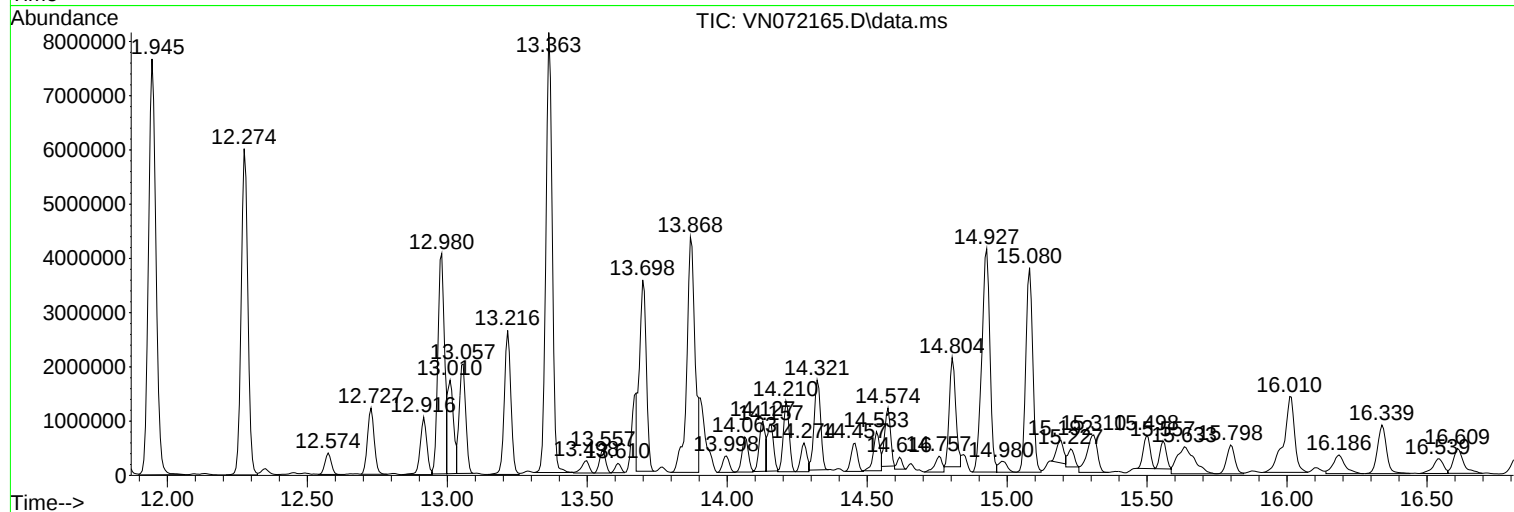
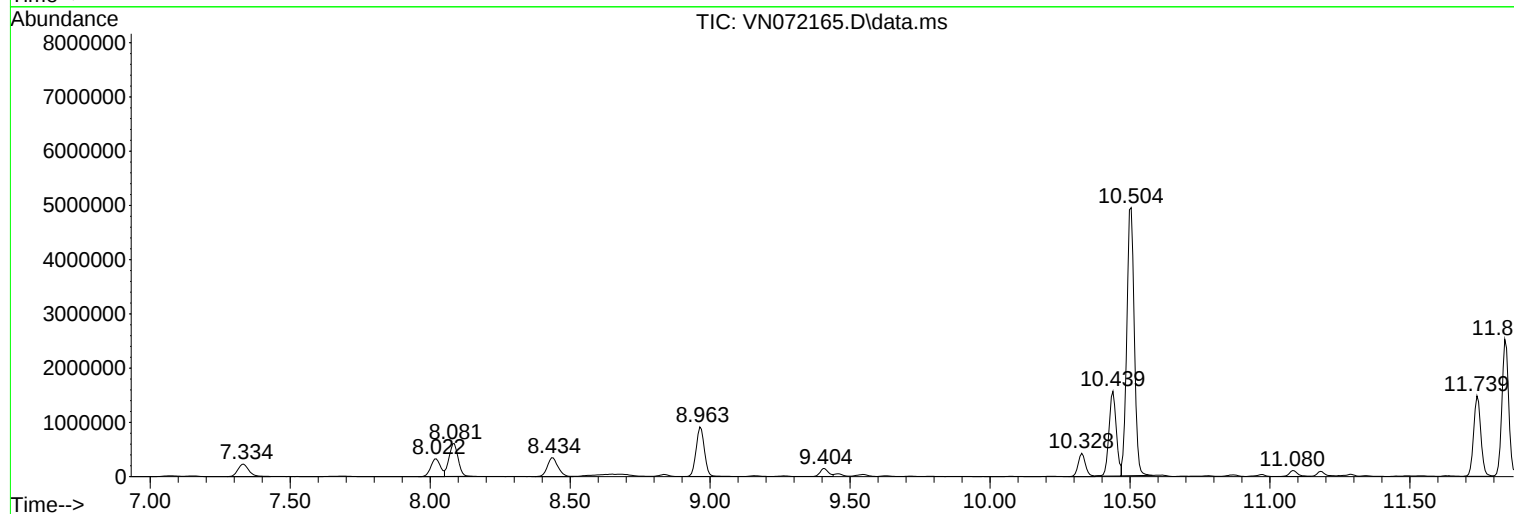
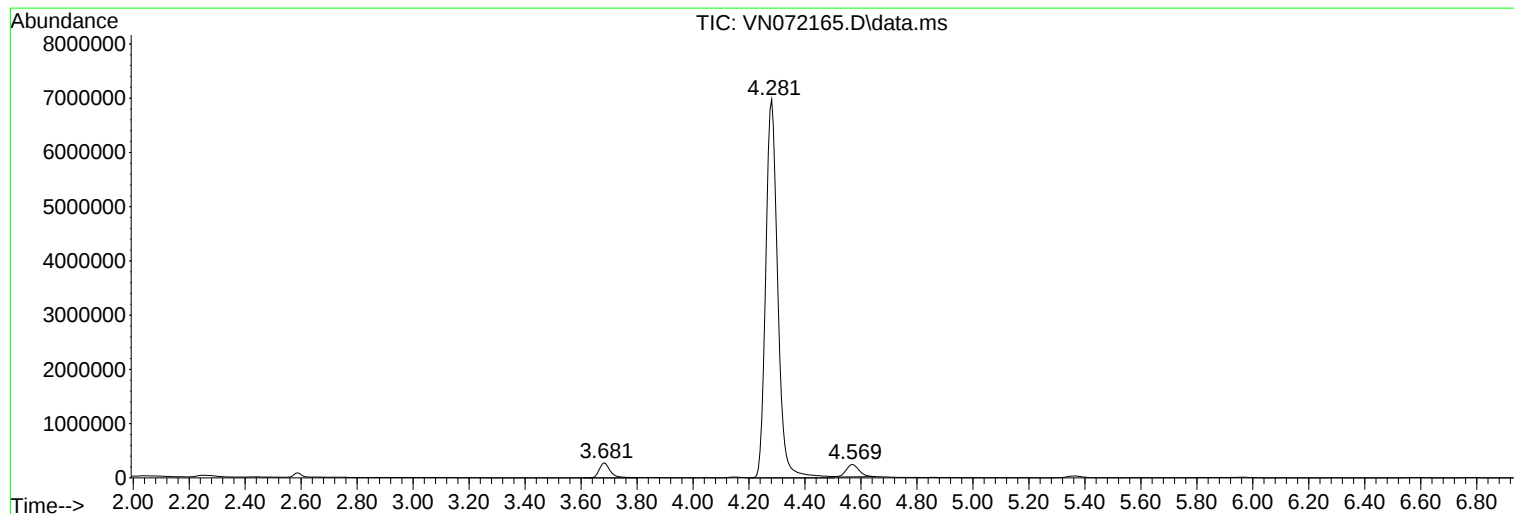
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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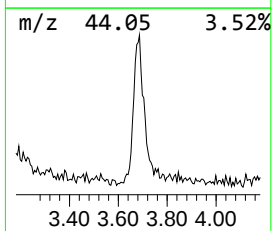
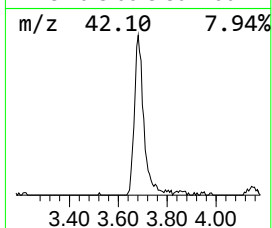
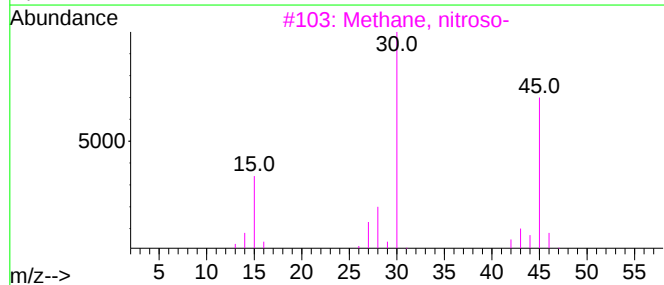
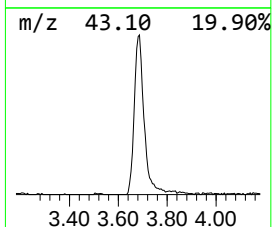
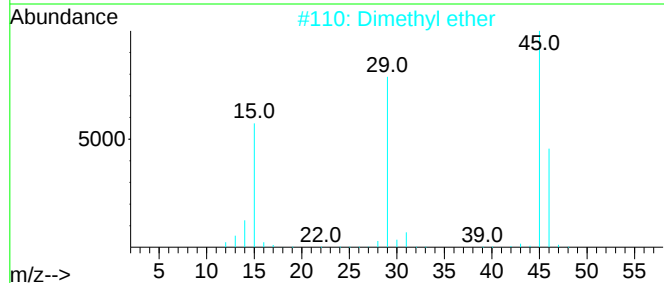
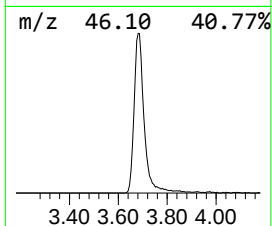
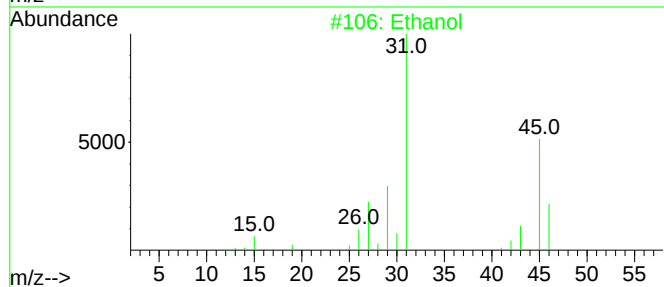
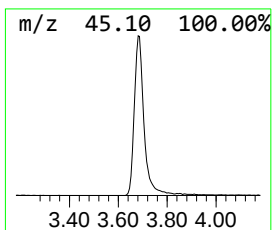
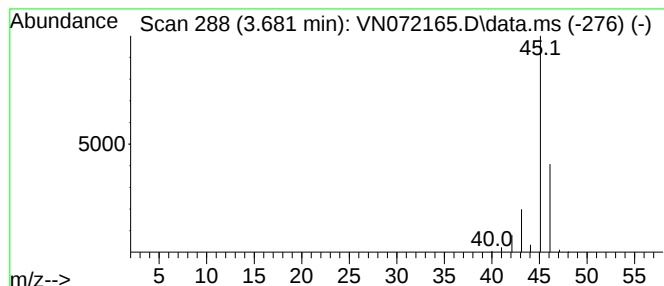
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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

Peak Number 1 Ethanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.681	27.94 ug/l	758652	Pentafluorobenzene	8.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol	46	C2H6O	000064-17-5	90
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			L-Lactic acid	90	C3H6O3	000079-33-4	4



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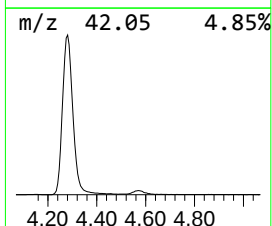
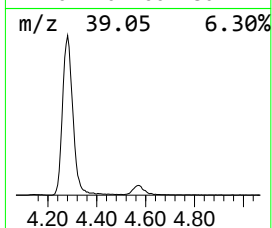
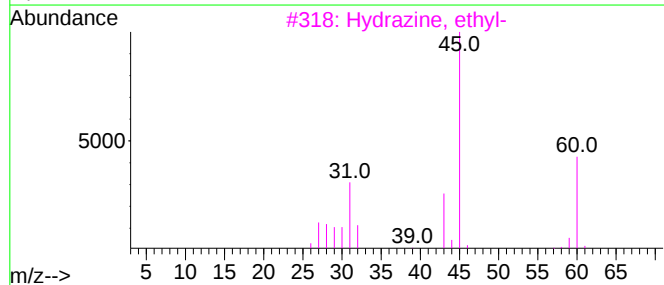
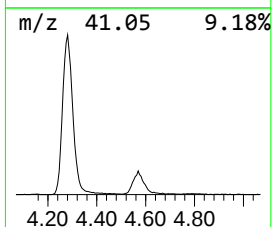
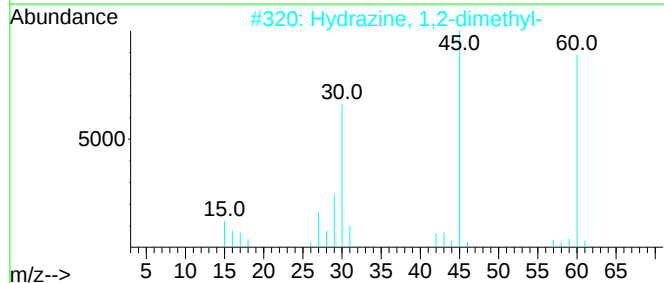
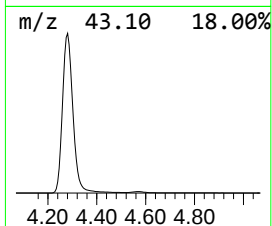
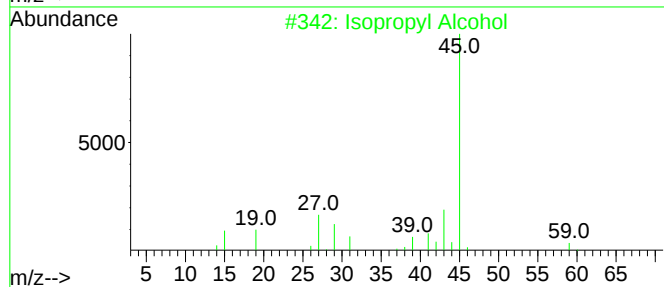
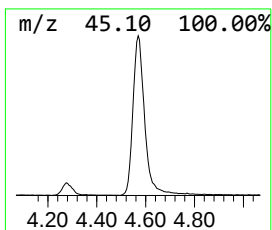
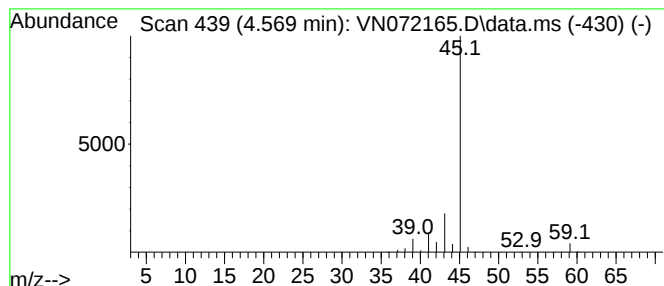
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Isopropyl Alcohol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.569	26.88 ug/l	729818	Pentafluorobenzene	8.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	72
2			Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3			Hydrazine, ethyl-	60	C2H8N2	000624-80-6	9
4			4-Penten-2-ol	86	C5H10O	000625-31-0	9
5			Formamide	45	CH3NO	000075-12-7	7



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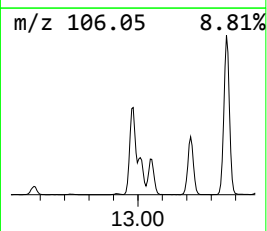
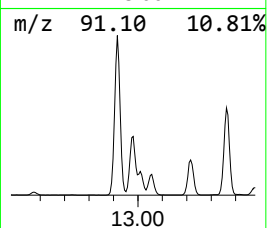
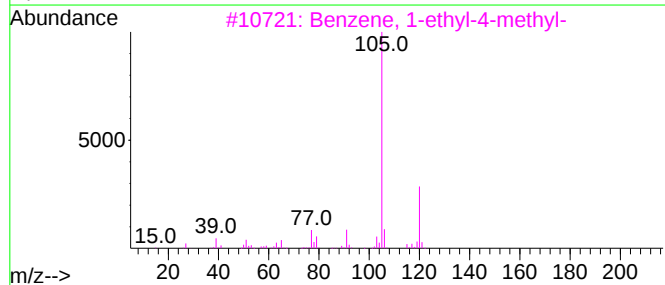
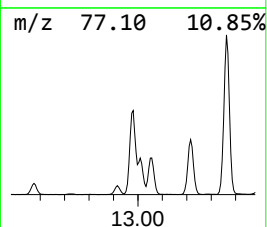
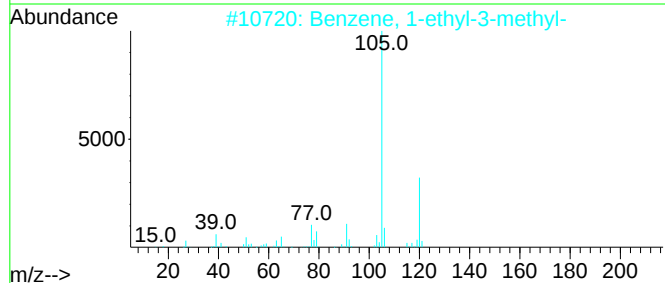
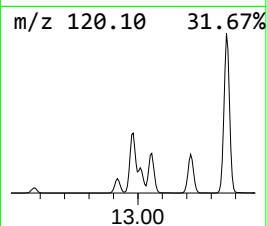
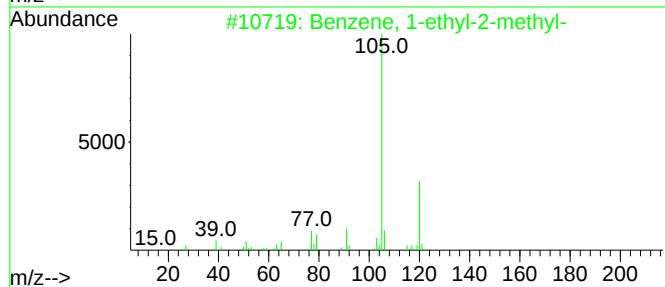
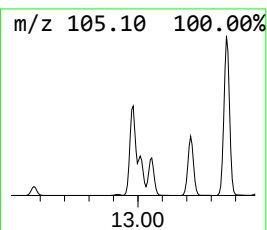
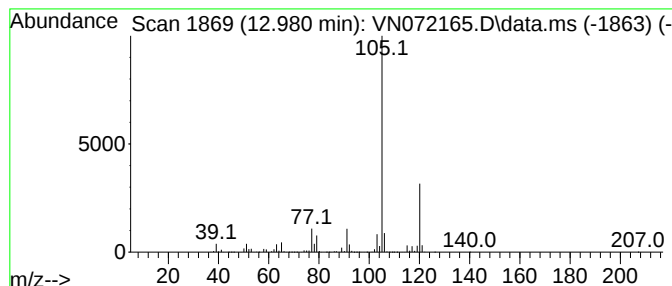
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	56.95 ug/l	6856620	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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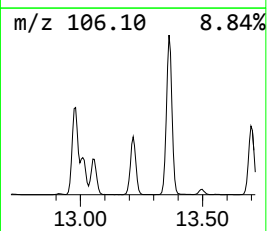
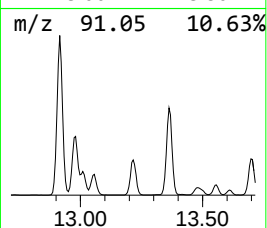
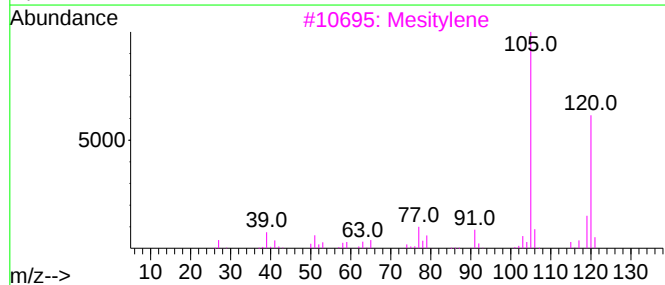
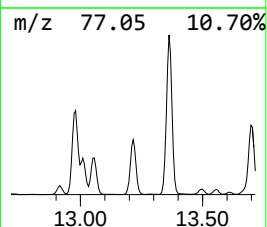
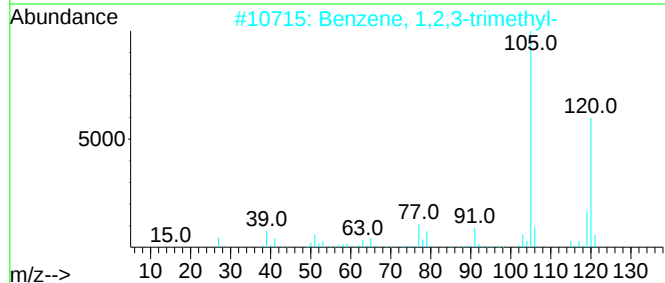
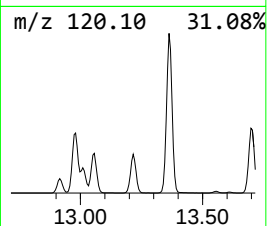
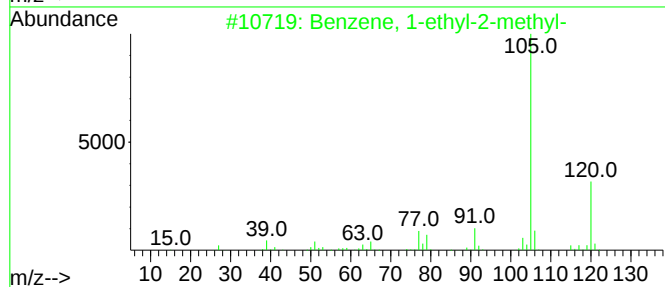
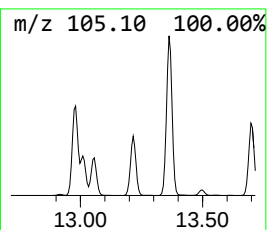
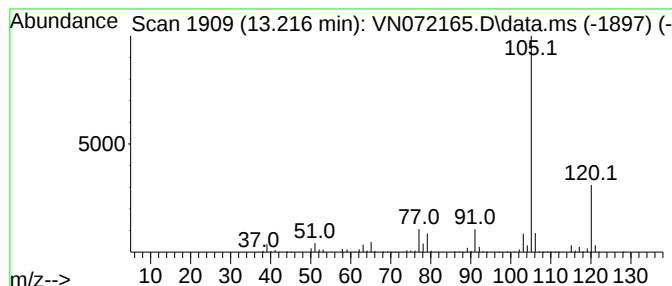
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	35.57 ug/l	4282810	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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NWB-1756

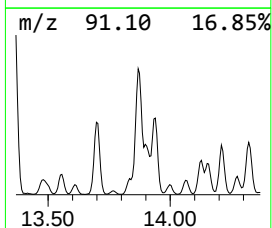
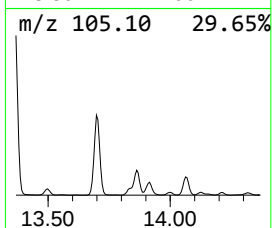
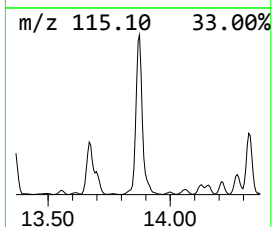
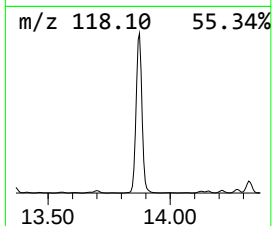
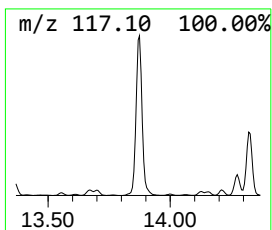
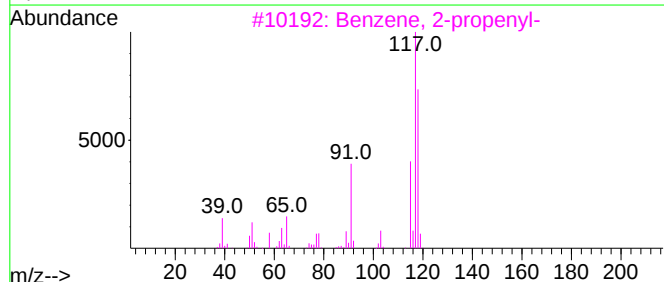
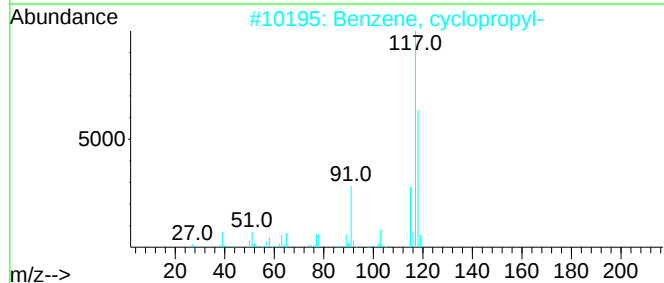
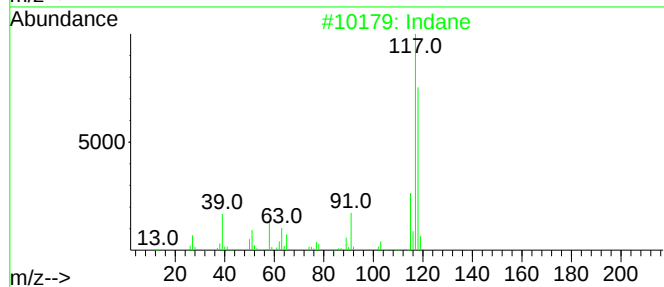
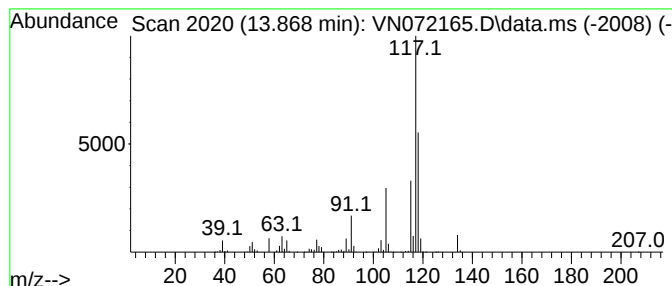
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Quant Title : SW846 8260

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

Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.868	74.53 ug/l	8973780	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	89
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4	Delta-cyclene	118	C9H10	007785-10-6	62
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072165.D
Acq On : 22 Apr 2022 18:35
Operator : JC\MD
Sample : N2522-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1756

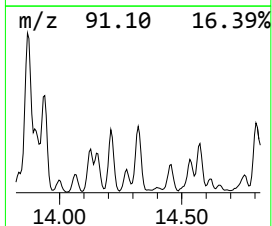
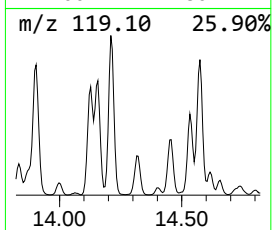
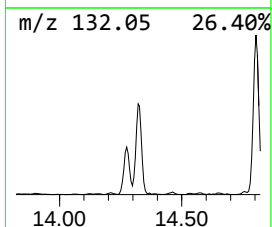
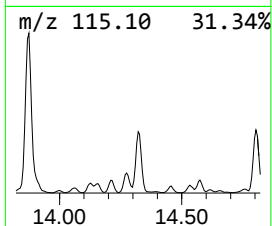
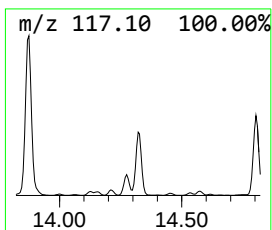
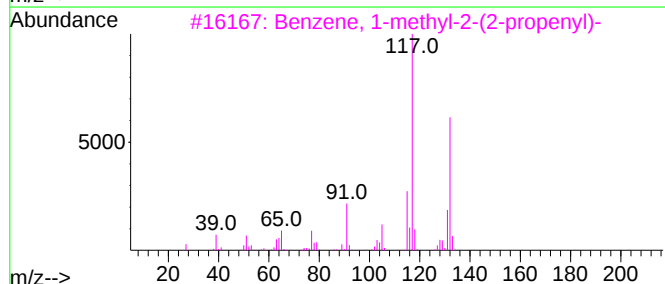
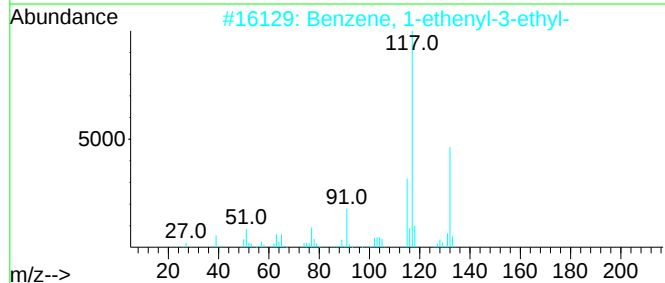
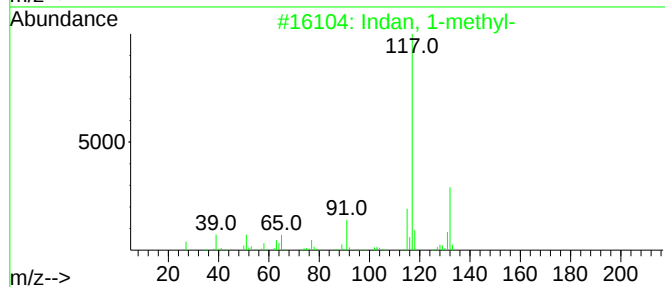
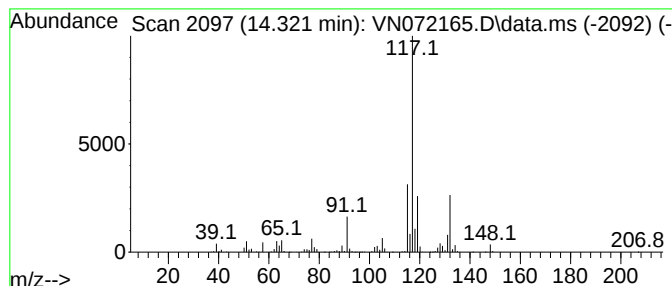
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Quant Title : SW846 8260

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.321	22.23 ug/l	2676330	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	68
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072165.D
Acq On : 22 Apr 2022 18:35
Operator : JC\MD
Sample : N2522-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1756

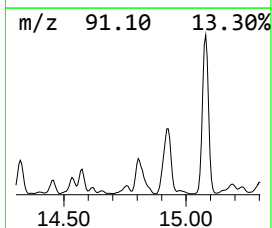
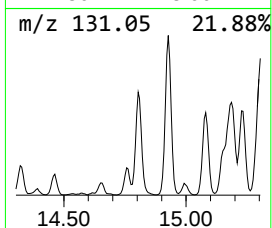
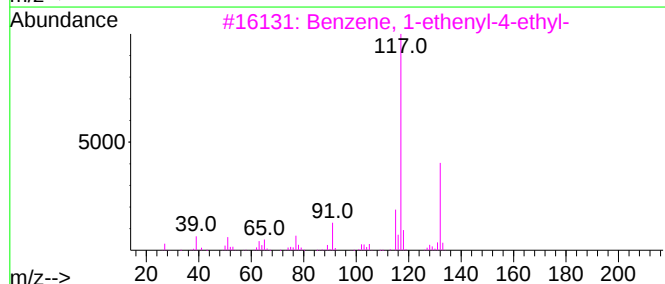
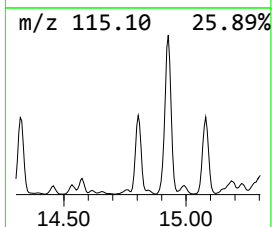
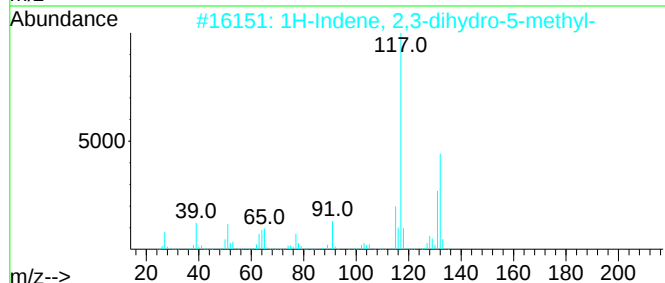
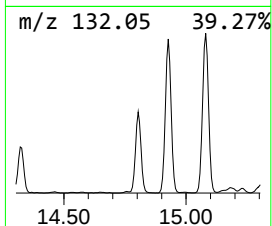
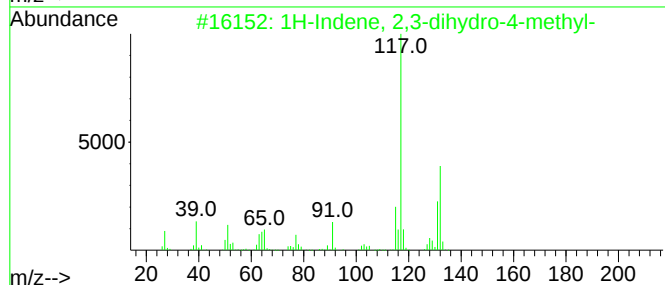
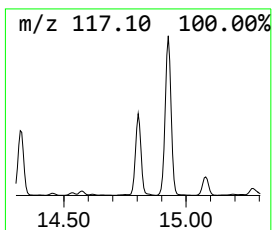
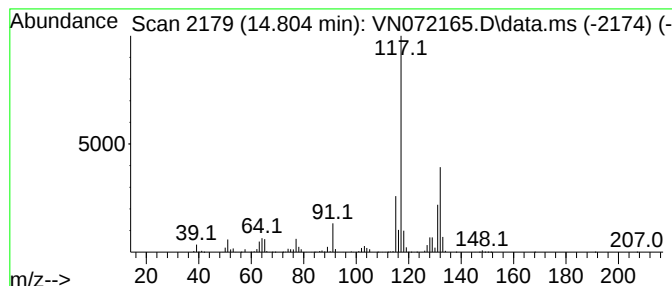
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	26.91 ug/l	3239810	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072165.D
Acq On : 22 Apr 2022 18:35
Operator : JC\MD
Sample : N2522-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1756

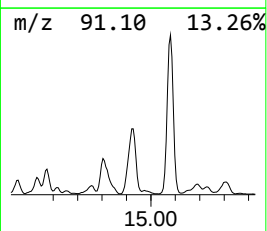
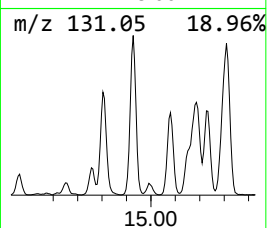
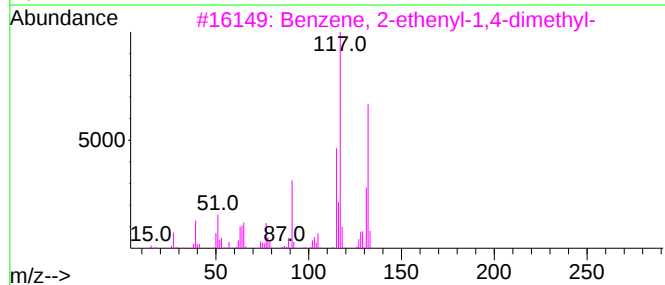
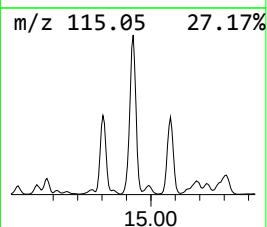
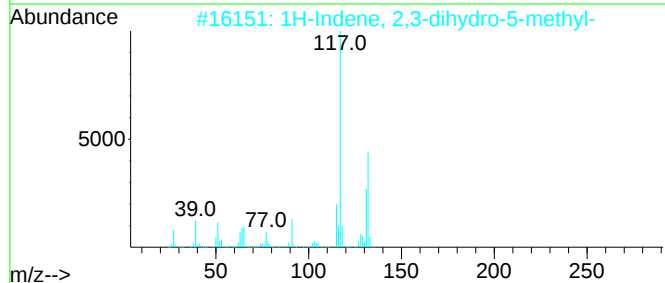
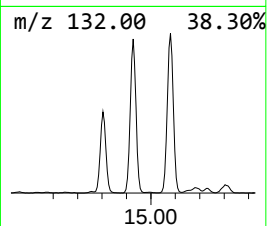
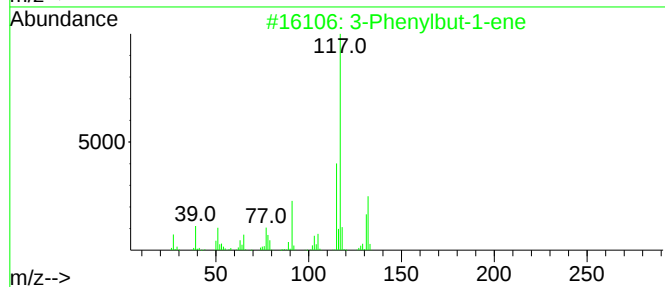
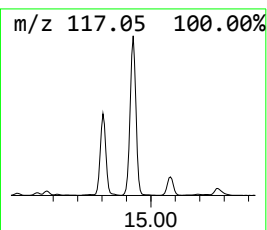
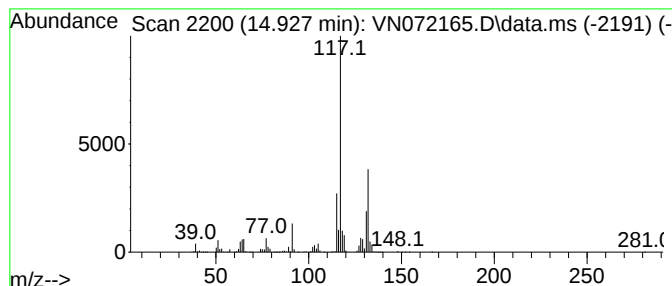
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Quant Title : SW846 8260

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

Peak Number 8 3-Phenylbut-1-ene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	68.76 ug/l	8279390	1,4-Dichlorobenzene-d4	13.669

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Phenylbut-1-ene	132	C10H12	000934-10-1	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072165.D
Acq On : 22 Apr 2022 18:35
Operator : JC\MD
Sample : N2522-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1756

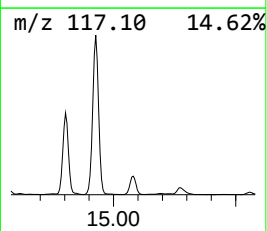
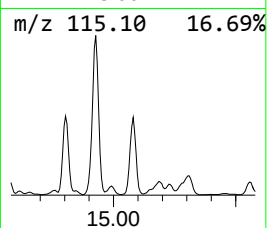
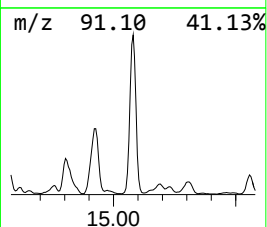
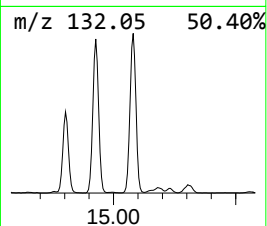
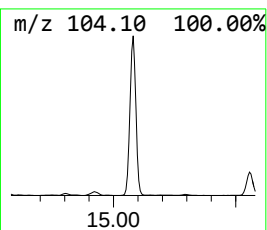
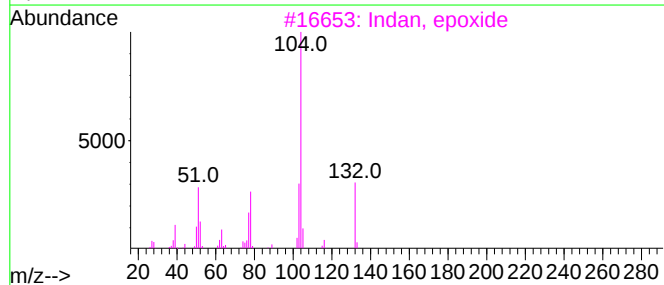
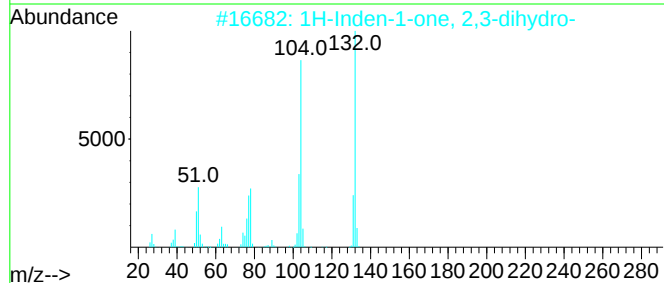
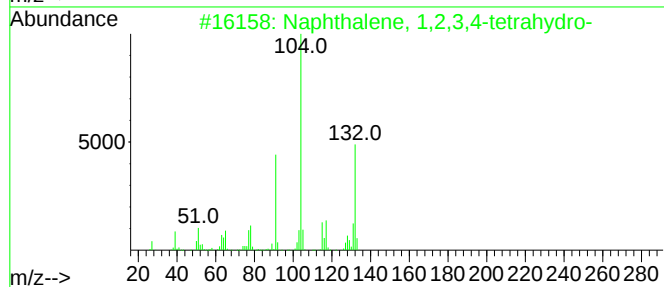
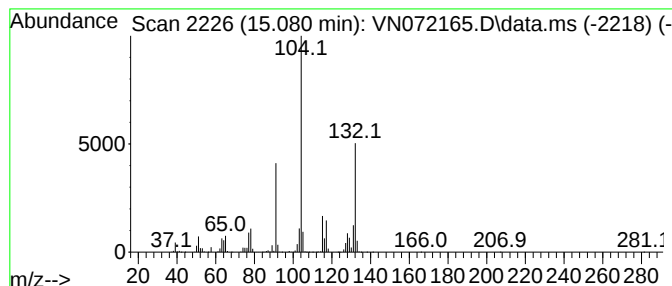
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.080	53.68 ug/l	6463570	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	59
3			Indan, epoxide	132	C9H8O	000768-22-9	58
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
5			3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072165.D
Acq On : 22 Apr 2022 18:35
Operator : JC\MD
Sample : N2522-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1756

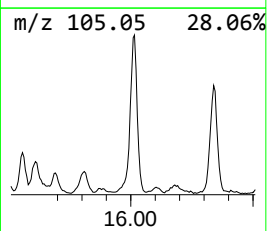
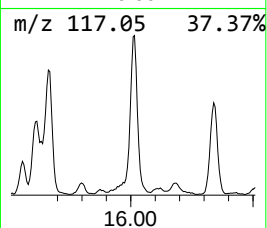
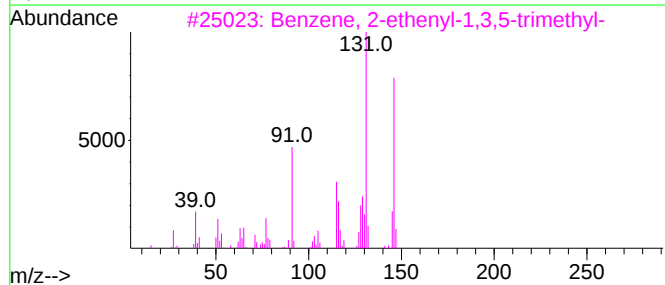
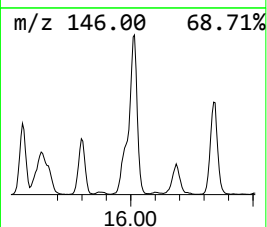
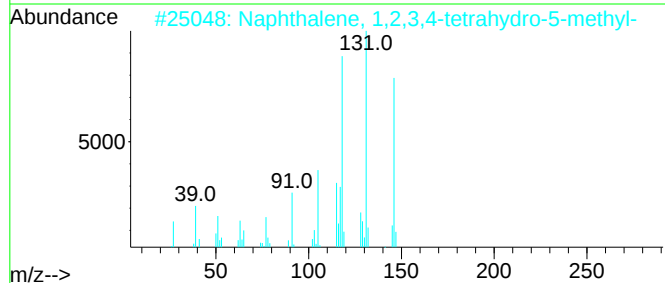
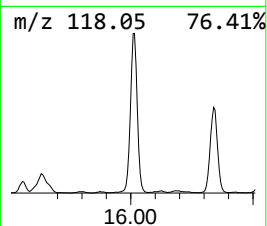
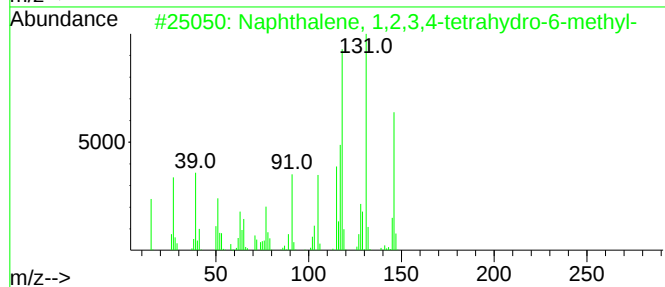
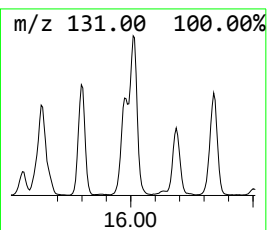
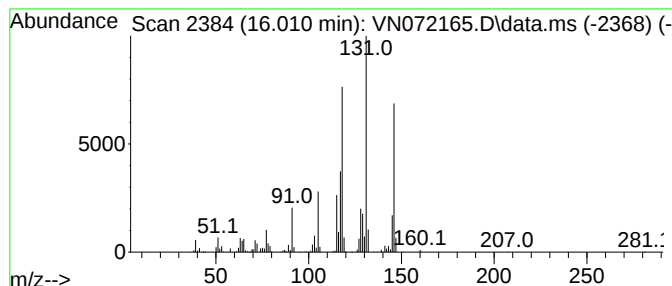
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.009	31.12 ug/l	3747010	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	83
4			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	64
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072165.D
Acq On : 22 Apr 2022 18:35
Operator : JC\MD
Sample : N2522-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NWB-1756

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
Ethanol	3.681	27.9	ug/l	758652	1	8.086	1357600	50.0
Isopropyl Alcohol	4.569	26.9	ug/l	729818	1	8.086	1357600	50.0
Benzene, 1-ethy...	12.980	57.0	ug/l	6856620	4	13.669	6020140	50.0
Benzene, 1,2,3-...	13.216	35.6	ug/l	4282810	4	13.669	6020140	50.0
Indane	13.868	74.5	ug/l	8973780	4	13.669	6020140	50.0
Indan, 1-methyl-	14.321	22.2	ug/l	2676330	4	13.669	6020140	50.0
1H-Indene, 2,3-...	14.804	26.9	ug/l	3239810	4	13.669	6020140	50.0
3-Phenylbut-1-ene	14.927	68.8	ug/l	8279390	4	13.669	6020140	50.0
Naphthalene, 1,...	15.080	53.7	ug/l	6463570	4	13.669	6020140	50.0
Naphthalene, 1,...	16.009	31.1	ug/l	3747010	4	13.669	6020140	50.0