

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
 Data File : VN073568.D
 Acq On : 21 Jul 2022 20:27
 Operator : JC\MD
 Sample : N3806-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FRAC-TANK-N48964

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\82N062822W.M

Title : SW846 8260

Signal : TIC: VN073568.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.552	106	113	119	rBV2	12430	24804	1.79%	0.144%
2	3.640	289	298	309	rBV	32365	87877	6.35%	0.509%
3	4.234	389	399	418	rBV	18009	63002	4.56%	0.365%
4	4.510	437	446	463	rBV2	18711	58672	4.24%	0.340%
5	6.504	777	785	802	rVB4	17785	58561	4.23%	0.339%
6	7.075	873	882	891	rVB3	15870	42275	3.06%	0.245%
7	7.269	904	915	926	rBV3	15730	46358	3.35%	0.269%
8	7.951	1021	1031	1036	rBV	179488	429084	31.03%	2.486%
9	8.016	1036	1042	1058	rVB2	328724	770149	55.69%	4.462%
10	8.369	1091	1102	1113	rBV	189835	439611	31.79%	2.547%
11	8.781	1163	1172	1181	rBV3	6700	19845	1.43%	0.115%
12	8.904	1184	1193	1207	rBV	455904	969156	70.08%	5.615%
13	9.351	1264	1269	1275	rBV2	7908	14790	1.07%	0.086%
14	9.498	1286	1294	1302	rVB3	6839	16851	1.22%	0.098%
15	10.275	1418	1426	1435	rBV2	19778	37050	2.68%	0.215%
16	10.381	1435	1444	1451	rVV	658322	1197597	86.60%	6.939%
17	10.445	1451	1455	1466	rVB	51070	95420	6.90%	0.553%
18	10.916	1526	1535	1540	rVB5	6531	13893	1.00%	0.080%
19	11.686	1658	1666	1676	rBV	656488	1163858	84.16%	6.743%
20	11.786	1677	1683	1692	rVV	32595	57737	4.17%	0.335%
21	11.886	1694	1700	1712	rVV	106168	201923	14.60%	1.170%
22	12.227	1751	1758	1766	rBV3	107117	217950	15.76%	1.263%
23	12.522	1798	1808	1813	rBV4	10554	17858	1.29%	0.103%
24	12.674	1827	1834	1844	rBV	532460	894344	64.67%	5.182%
25	12.863	1860	1866	1871	rBV	23821	36972	2.67%	0.214%
26	12.922	1871	1876	1880	rVV	117033	199672	14.44%	1.157%
27	12.957	1880	1882	1885	rVV	47071	59712	4.32%	0.346%
28	12.998	1885	1889	1897	rVB	78806	132429	9.58%	0.767%
29	13.163	1909	1917	1924	rBV3	78442	156997	11.35%	0.910%
30	13.310	1934	1942	1952	rBV	309420	528194	38.19%	3.060%
31	13.439	1957	1964	1970	rVB3	11236	23490	1.70%	0.136%
32	13.504	1970	1975	1979	rBV3	19217	30498	2.21%	0.177%
33	13.551	1979	1983	1987	rBV4	10712	14740	1.07%	0.085%
34	13.616	1987	1994	2005	rBV2	676829	1382933	100.00%	8.012%
35	13.710	2006	2010	2016	rVB2	40258	61440	4.44%	0.356%
36	13.774	2016	2021	2022	rBV3	27036	30569	2.21%	0.177%

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37	13.816	2022	2028	2031	rVV2	151608	279906	20.24%	1.622%
38	13.851	2031	2034	2044	rVB3	95920	182999	13.23%	1.060%
39	13.939	2044	2049	2055	rBV3	26244	43507	3.15%	0.252%
40	14.010	2055	2061	2066	rBV	36866	57564	4.16%	0.334%
41	14.069	2066	2071	2074	rBV	66068	115314	8.34%	0.668%
42	14.098	2074	2076	2081	rVB	65051	88616	6.41%	0.513%
43	14.157	2081	2086	2091	rVB	100113	148245	10.72%	0.859%
44	14.221	2091	2097	2100	rBV2	32710	54810	3.96%	0.318%
45	14.269	2100	2105	2111	rVB2	109538	181220	13.10%	1.050%
46	14.398	2122	2127	2132	rBV	45417	67799	4.90%	0.393%
47	14.480	2134	2141	2144	rBV	88643	148944	10.77%	0.863%
48	14.516	2144	2147	2152	rVB	126952	182647	13.21%	1.058%
49	14.563	2152	2155	2158	rVB2	29226	36166	2.62%	0.210%
50	14.604	2158	2162	2169	rVB3	23882	45024	3.26%	0.261%
51	14.704	2169	2179	2182	rBV4	30307	56496	4.09%	0.327%
52	14.751	2182	2187	2191	rBV	133235	222236	16.07%	1.288%
53	14.792	2191	2194	2198	rVB3	54605	77476	5.60%	0.449%
54	14.863	2198	2206	2213	rBV3	297295	676694	48.93%	3.921%
55	14.916	2213	2215	2224	rVB4	30562	61637	4.46%	0.357%
56	15.021	2226	2233	2240	rBV	270640	458233	33.13%	2.655%
57	15.127	2241	2251	2255	rBV5	98785	242659	17.55%	1.406%
58	15.245	2264	2271	2280	rVB4	114697	272232	19.69%	1.577%
59	15.327	2282	2285	2290	rVB6	9877	17588	1.27%	0.102%
60	15.433	2291	2303	2309	rBV	233139	494419	35.75%	2.865%
61	15.492	2309	2313	2318	rVV	90085	156783	11.34%	0.908%
62	15.568	2318	2326	2334	rVV4	78804	285731	20.66%	1.655%
63	15.633	2334	2337	2346	rVB7	42116	73657	5.33%	0.427%
64	15.733	2346	2354	2360	rBV	79814	159528	11.54%	0.924%
65	15.810	2362	2367	2372	rVB9	11557	21570	1.56%	0.125%
66	15.945	2373	2390	2400	rBV	234711	616879	44.61%	3.574%
67	16.033	2400	2405	2410	rBV3	29673	55488	4.01%	0.321%
68	16.115	2411	2419	2433	rVB5	61450	194723	14.08%	1.128%
69	16.262	2435	2444	2452	rBV3	169668	386029	27.91%	2.237%
70	16.462	2469	2478	2483	rBV5	92888	214746	15.53%	1.244%
71	16.533	2483	2490	2498	rVV	328891	683313	49.41%	3.959%
72	16.739	2516	2525	2532	rBV	246836	602875	43.59%	3.493%

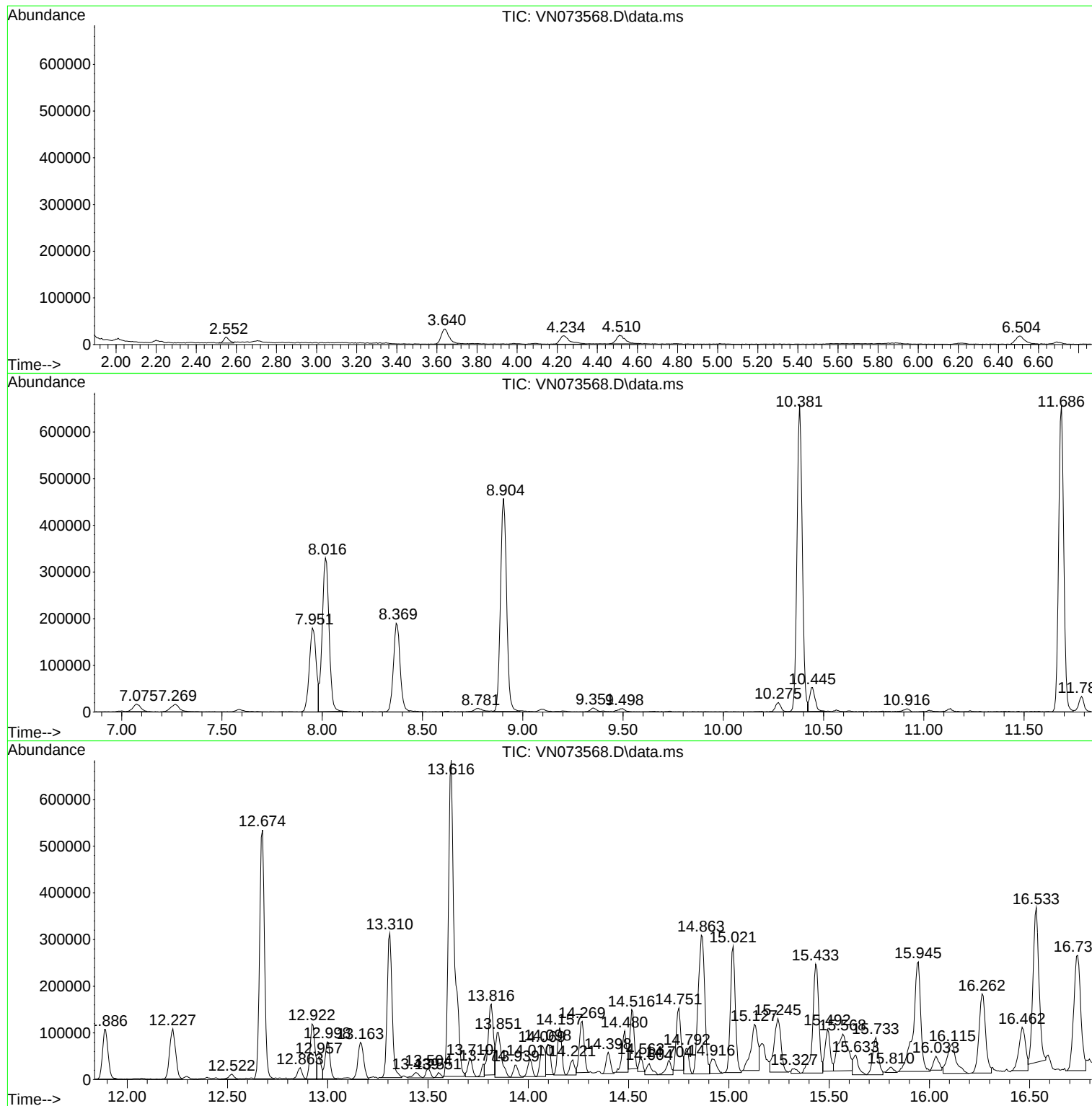
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ClientSampleId :
FRAC-TANK-N48964

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

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



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ClientSampleId :
FRAC-TANK-N48964

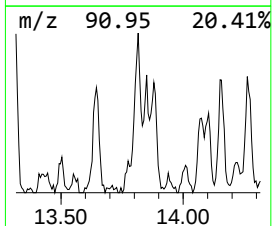
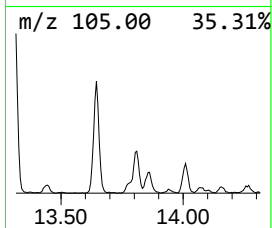
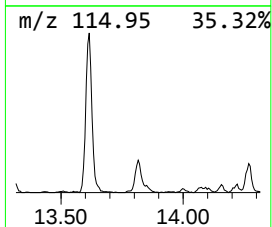
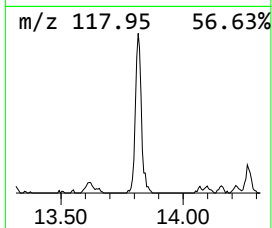
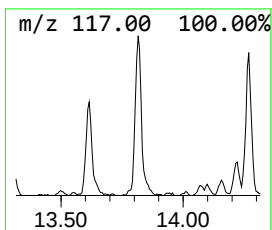
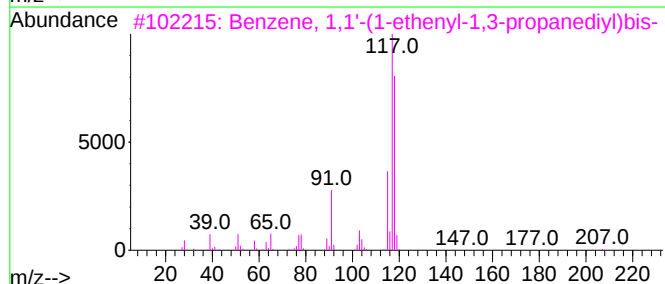
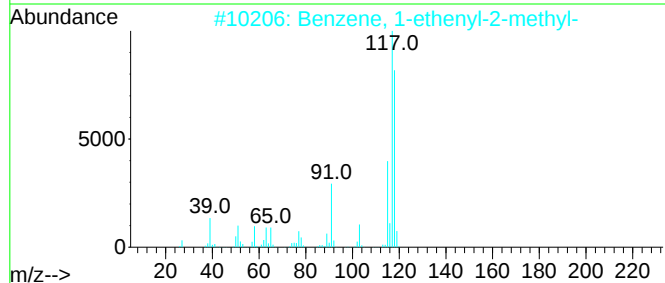
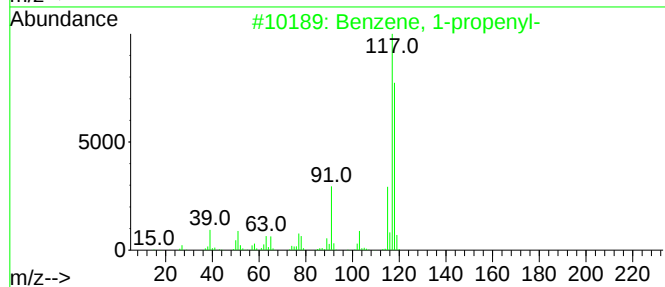
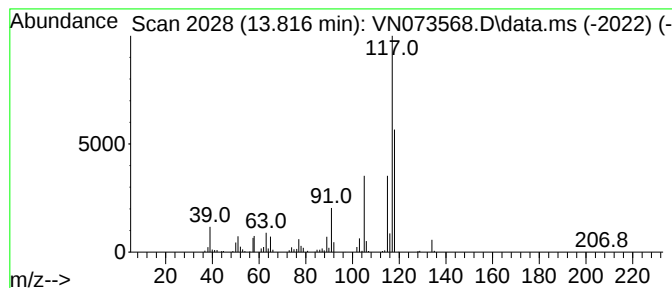
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	10.12 ug/l	279906	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	92
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
3			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	72
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
5			Indane	118	C9H10	000496-11-7	70



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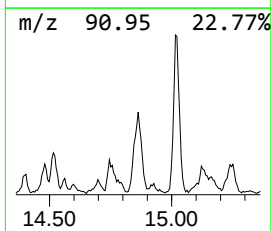
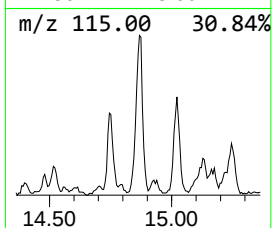
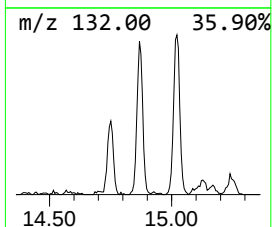
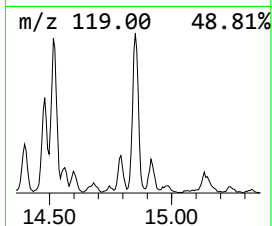
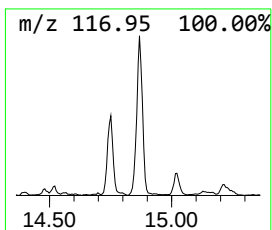
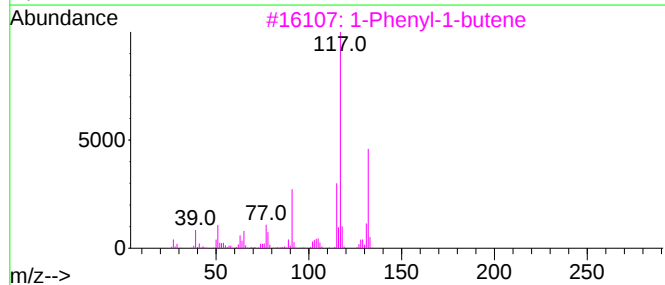
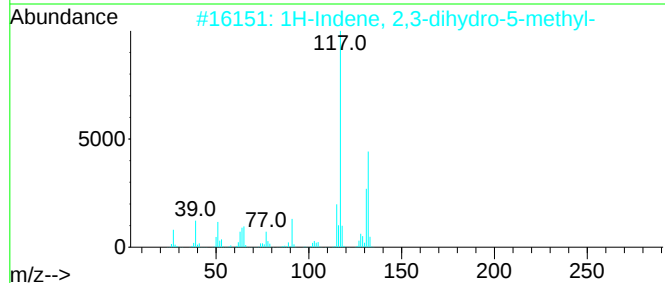
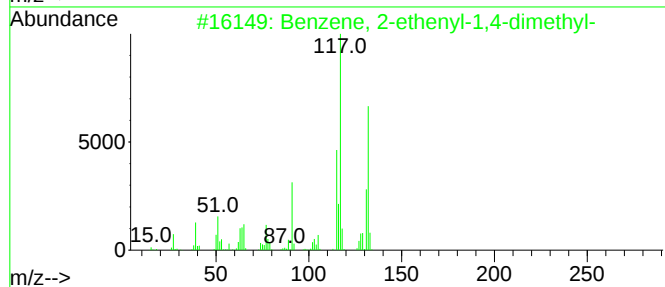
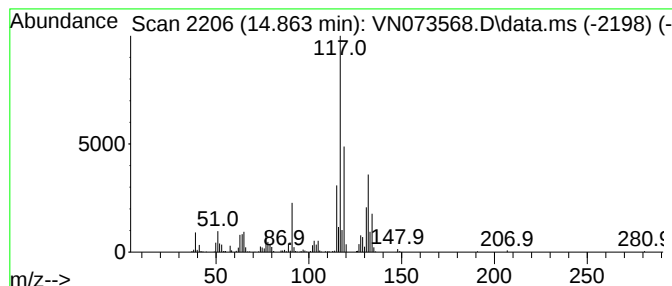
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.863	24.47 ug/l	676694	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4			2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76



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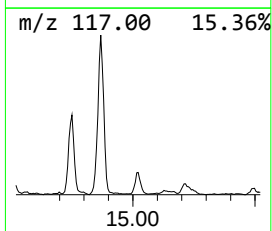
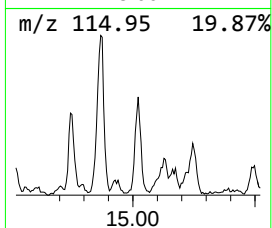
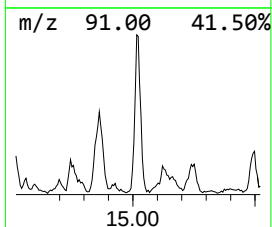
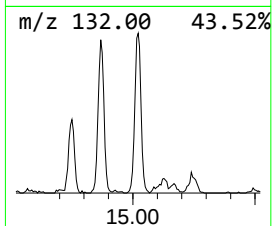
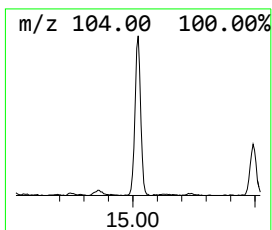
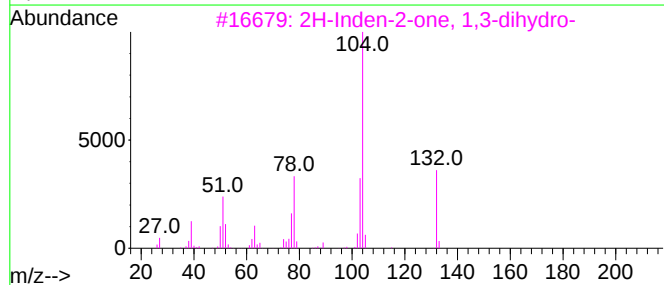
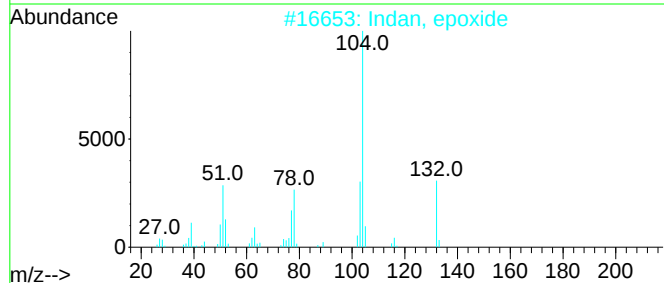
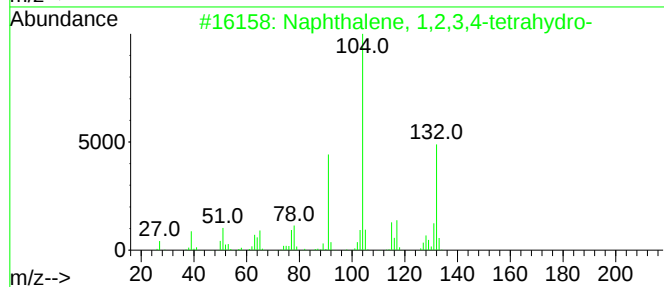
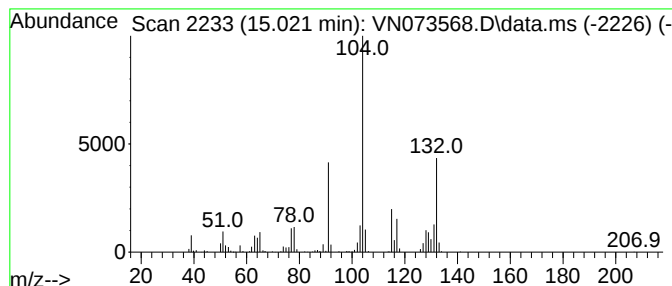
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Peak Number 3 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.021	16.57 ug/l	458233	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Indan, epoxide	132	C9H8O	000768-22-9	60
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	49
4			N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	38
5			3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	35



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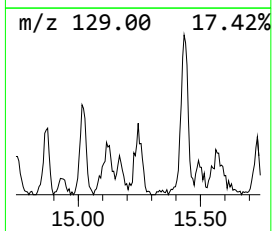
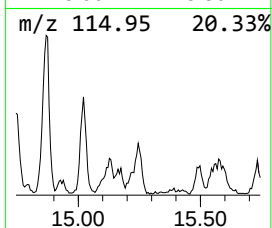
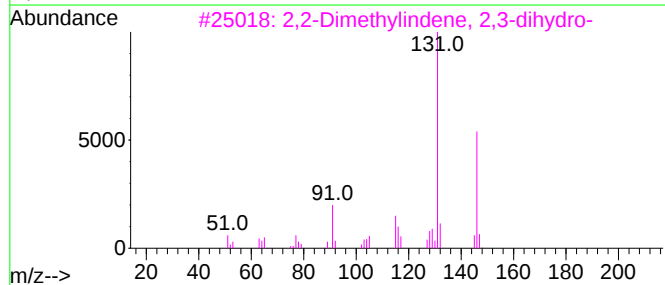
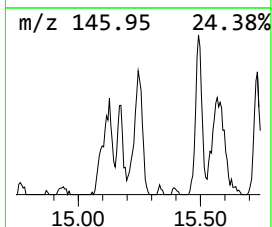
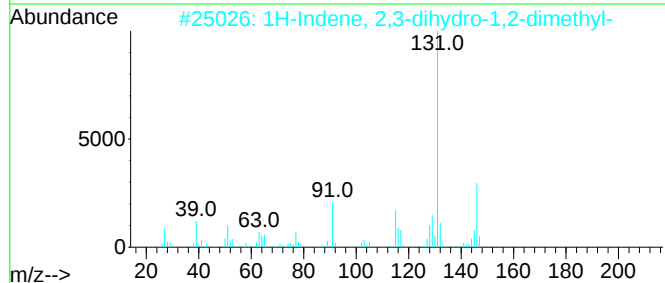
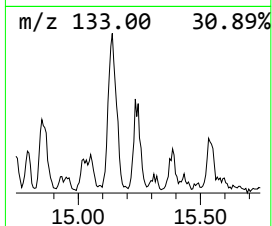
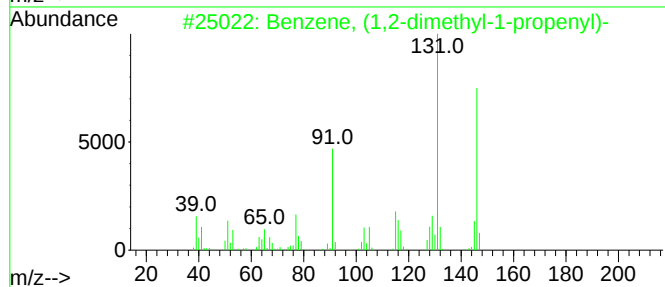
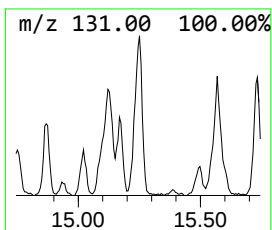
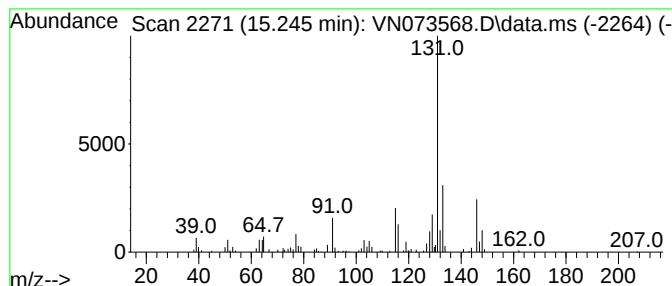
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.245	9.84 ug/l	272232	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	81
5			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FRAC-TANK-N48964

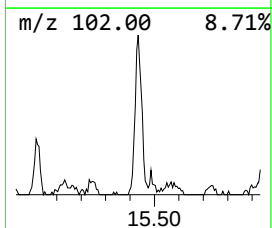
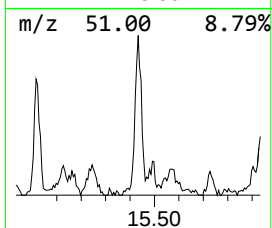
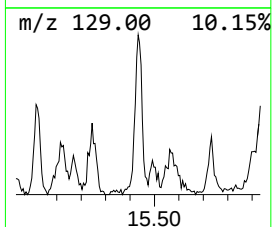
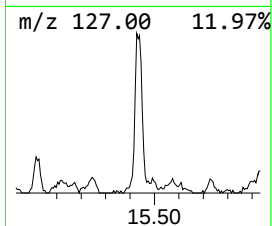
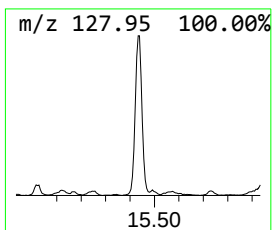
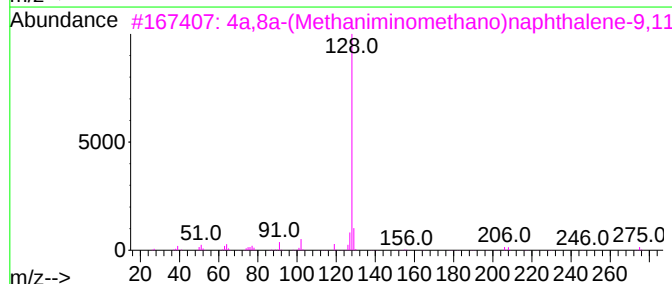
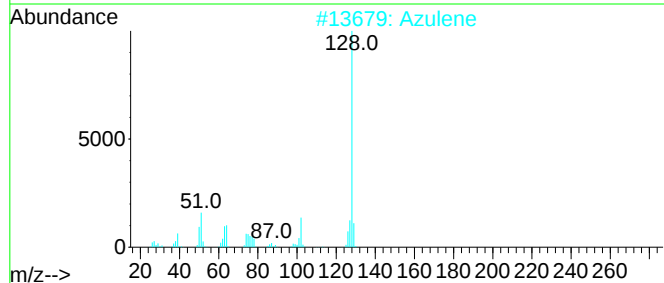
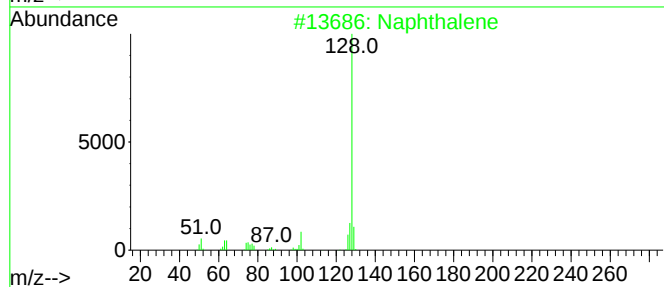
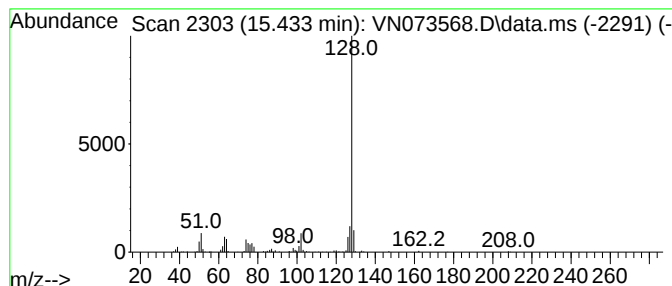
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Quant Title : SW846 8260

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

Peak Number 5 4a,8a-(Methaniminomethano)n... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.433	17.88 ug/l	494419	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	94
3			4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	83
4			1H-Cyclopenta[c]thiophene, hexah...	128	C7H12S	053907-80-5	47
5			[4.2.2]Propella-2,4,7,9-tetraene	128	C10H8	088090-34-0	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FRAC-TANK-N48964

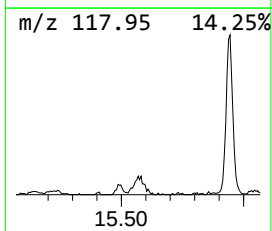
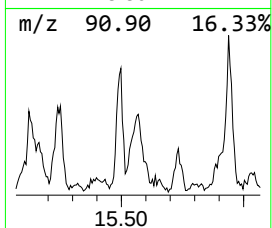
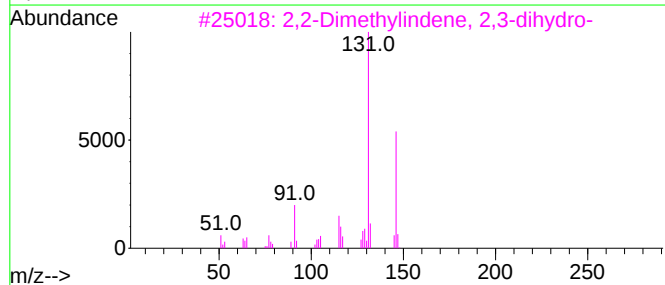
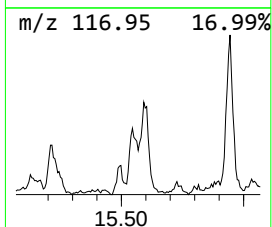
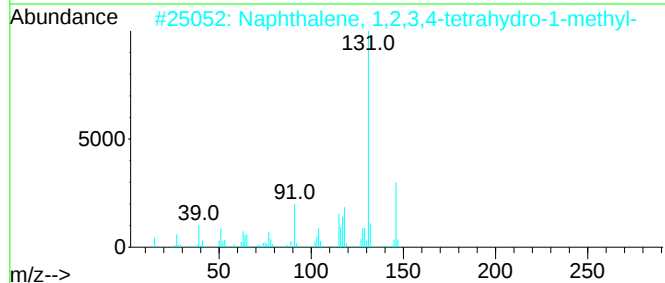
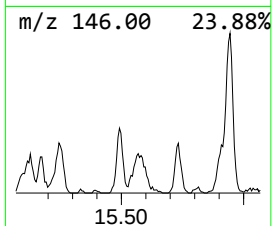
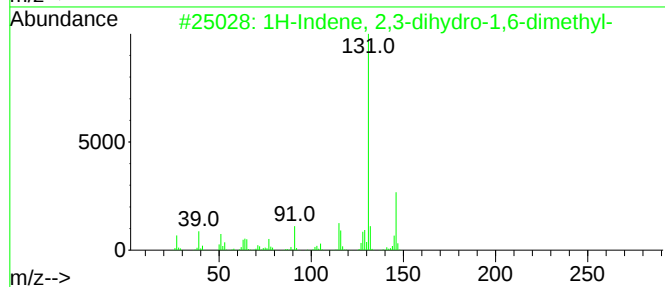
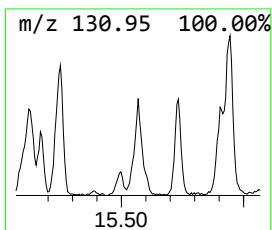
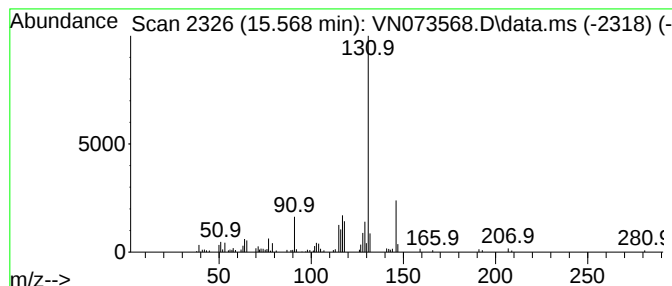
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.568	10.33 ug/l	285731	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	93
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

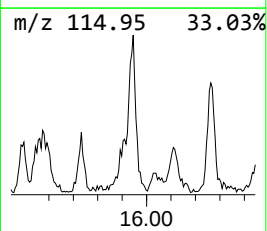
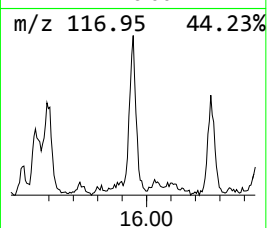
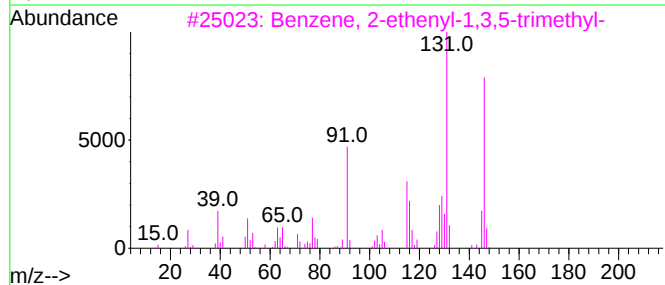
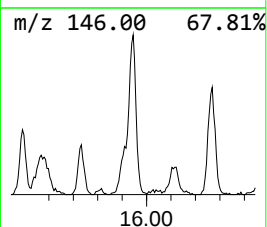
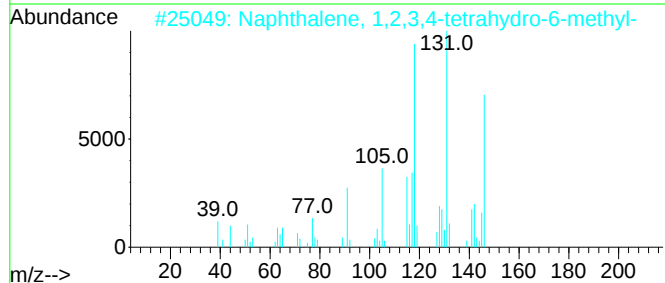
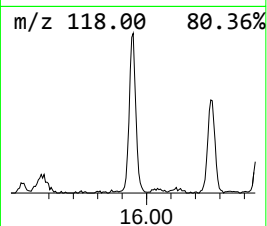
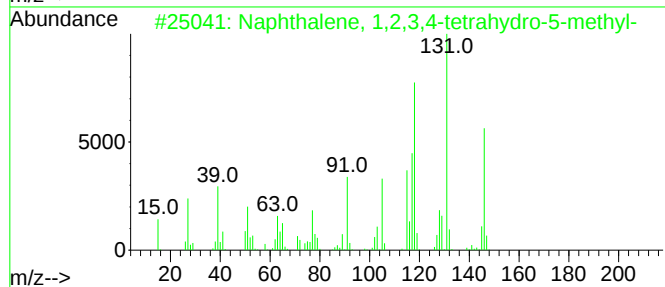
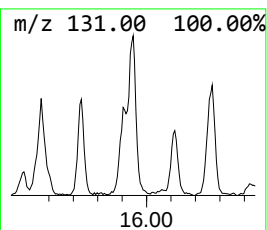
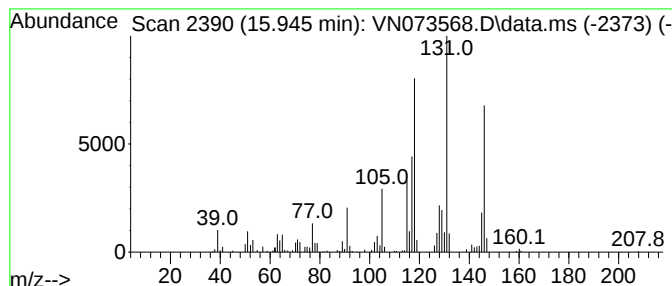
Instrument :
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ClientSampleId :
FRAC-TANK-N48964

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.945	22.30 ug/l	616879	1,4-Dichlorobenzene-d4			13.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	002809-64-5	93
2	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001680-51-9	91
3	Benzene, 2-ethenyl-1,3,5-trimethyl-		146	C11H14	000769-25-5	49
4	Benzene, (1,2-dimethyl-1-propenyl)-		146	C11H14	000769-57-3	49
5	1H-Inden-1-one, 2,3-dihydro-2-me...		146	C10H10O	017496-14-9	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

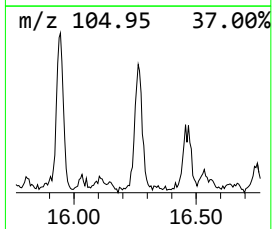
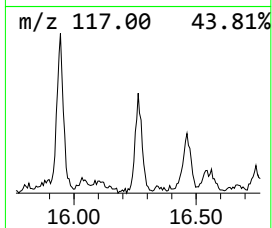
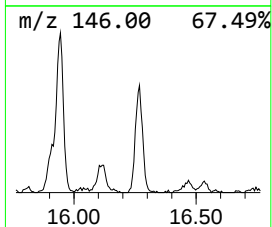
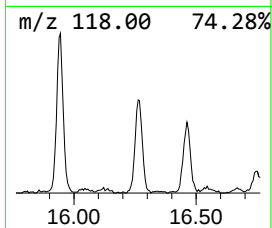
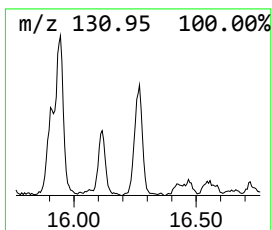
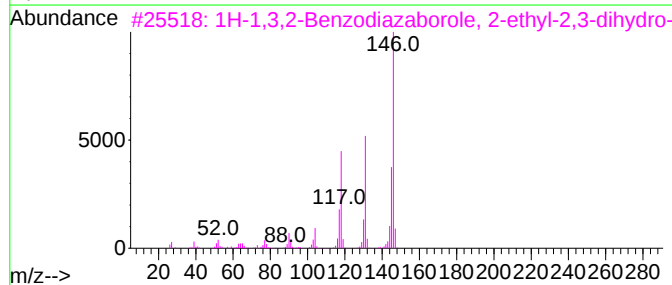
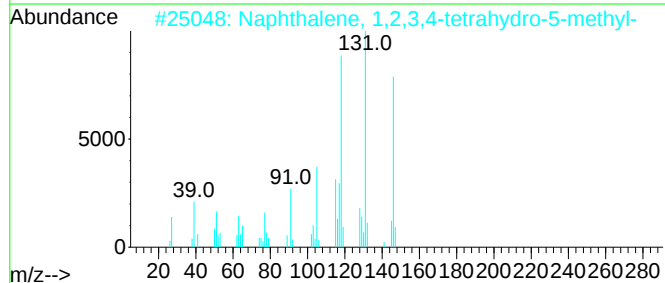
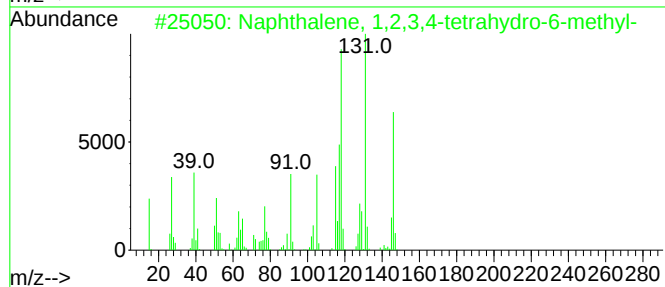
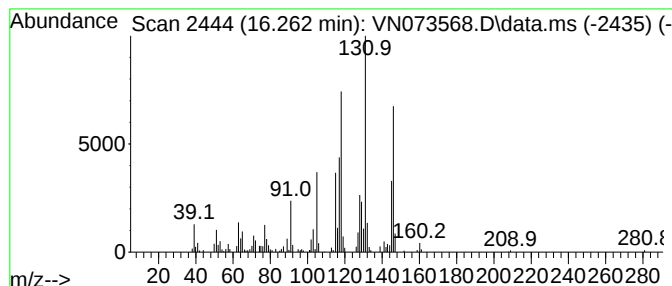
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MSVOA_N
ClientSampleId :
FRAC-TANK-N48964

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.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.
16.262	13.96 ug/l	386029	1,4-Dichlorobenzene-d4			13.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001680-51-9	93
2	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	002809-64-5	93
3	1H-1,3,2-Benzodiazaborole, 2-eth...		146	C8H11BN2	074663-81-3	90
4	2,3,4,5,6,7-Hexahydro-1H-cyclope...		146	C11H14	1010189-31-0	87
5	Benzene, 2-ethenyl-1,3,5-trimethyl-		146	C11H14	000769-25-5	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FRAC-TANK-N48964

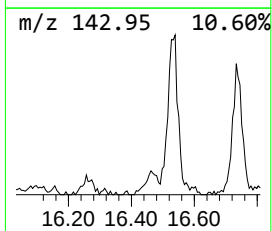
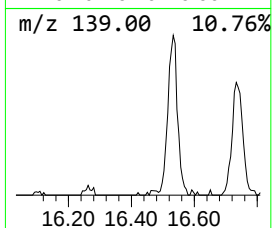
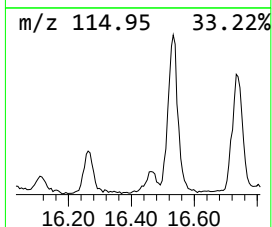
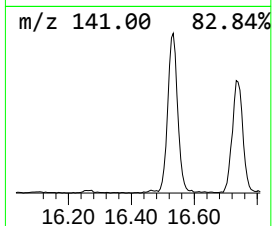
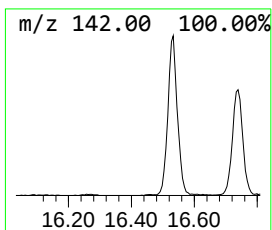
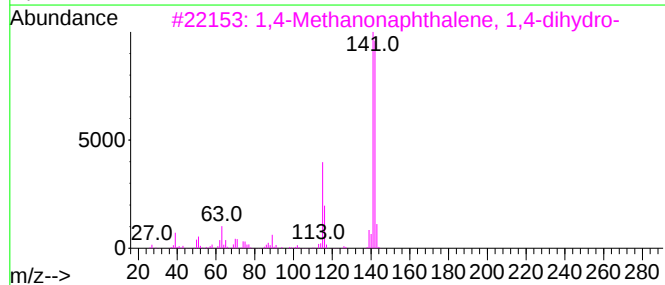
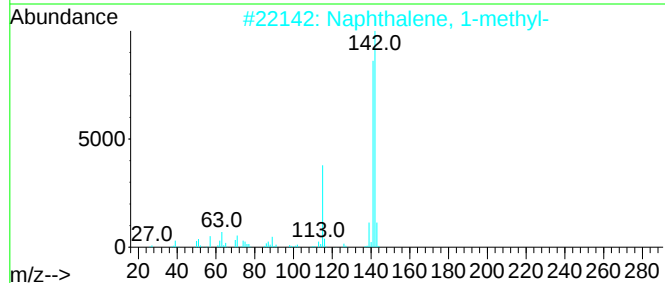
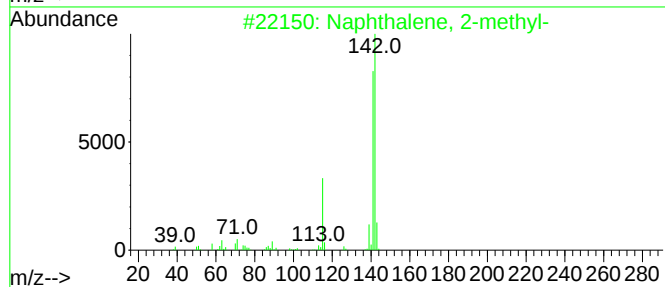
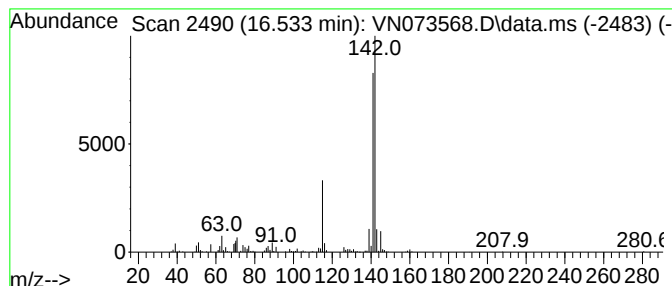
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Quant Title : SW846 8260

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

Peak Number 9 1,4-Methanonaphthalene, 1,4... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.533	24.71 ug/l	683313	1,4-Dichlorobenzene-d4	13.616

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		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-dihy...	142	C11H10	004453-90-1	93
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FRAC-TANK-N48964

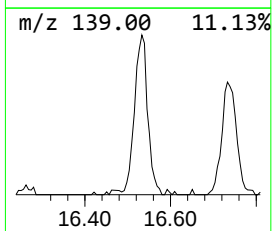
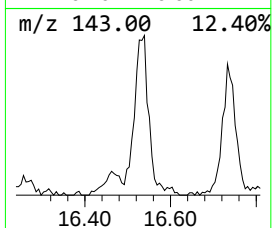
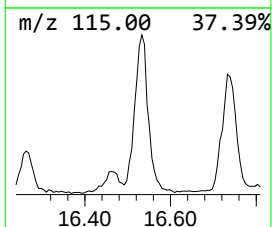
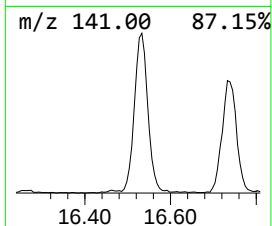
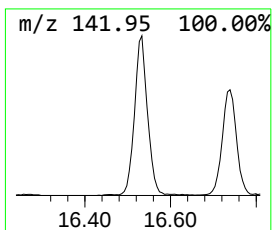
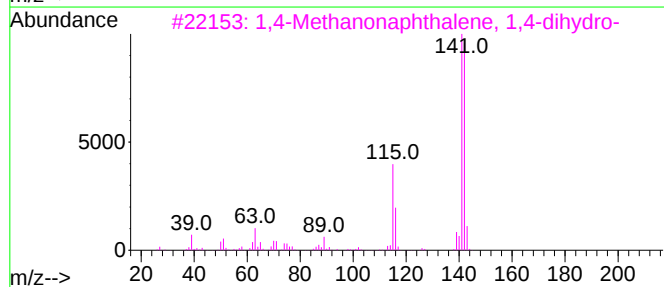
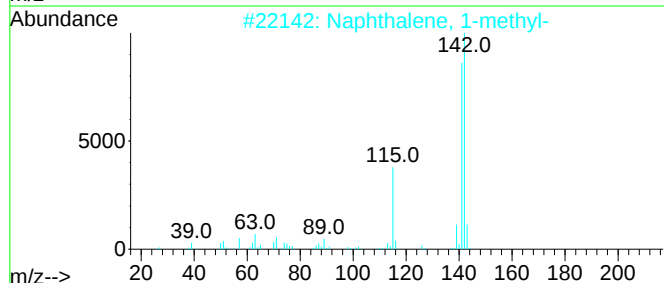
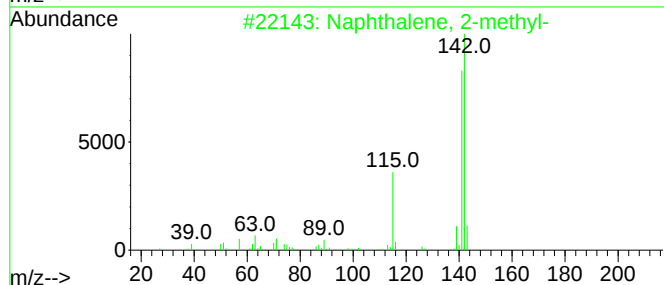
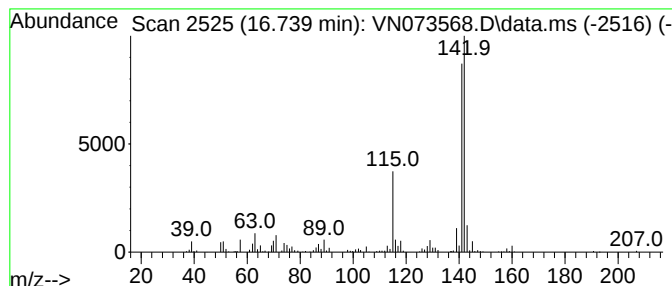
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Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.739	21.80 ug/l	602875	1,4-Dichlorobenzene-d4	13.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073568.D
Acq On : 21 Jul 2022 20:27
Operator : JC\MD
Sample : N3806-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FRAC-TANK-N48964

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.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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Benzene, 2-ethe...	14.863	24.5	ug/l	676694	4	13.616	1382930	50.0
Naphthalene, 1,...	15.021	16.6	ug/l	458233	4	13.616	1382930	50.0
Benzene, (1,2-d...	15.245	9.8	ug/l	272232	4	13.616	1382930	50.0
4a,8a-(Methanim...	15.433	17.9	ug/l	494419	4	13.616	1382930	50.0
1H-Indene, 2,3-...	15.568	10.3	ug/l	285731	4	13.616	1382930	50.0
Naphthalene, 1,...	15.945	22.3	ug/l	616879	4	13.616	1382930	50.0
Naphthalene, 1,...	16.262	14.0	ug/l	386029	4	13.616	1382930	50.0
1,4-Methanonaph...	16.533	24.7	ug/l	683313	4	13.616	1382930	50.0
Benzocyclohepta...	16.739	21.8	ug/l	602875	4	13.616	1382930	50.0