

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
 Data File : VN071016.D
 Acq On : 18 Feb 2022 21:27
 Operator : JC\MD
 Sample : N1611-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1070

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\82N021522W.M
 Title : SW846 8260

Signal : TIC: VN071016.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.587	96	102	111	rBV	137119	244486	1.57%	0.124%
2	3.681	276	288	313	rBV	1176740	3201565	20.53%	1.622%
3	4.287	381	391	403	rBV	53492	163066	1.05%	0.083%
4	8.016	1013	1025	1031	rBV	543631	1352446	8.67%	0.685%
5	8.081	1031	1036	1057	rVB	1115151	2503386	16.05%	1.268%
6	8.433	1084	1096	1108	rBV	644633	1562283	10.02%	0.792%
7	8.963	1177	1186	1207	rBV	1552690	3221862	20.66%	1.632%
8	10.439	1423	1437	1443	rBV	2393027	4408229	28.26%	2.234%
9	10.498	1443	1447	1461	rVB	491773	925726	5.94%	0.469%
10	11.739	1650	1658	1669	rBV	2117667	3760317	24.11%	1.905%
11	11.839	1669	1675	1684	rBV	509534	875053	5.61%	0.443%
12	11.945	1684	1693	1704	rBV	2324638	4488064	28.77%	2.274%
13	12.274	1737	1749	1758	rBV	2143643	3866462	24.79%	1.959%
14	12.363	1759	1764	1769	rVB	131721	200451	1.29%	0.102%
15	12.480	1781	1784	1794	rVV5	120726	308614	1.98%	0.156%
16	12.574	1794	1800	1805	rVV	298654	527064	3.38%	0.267%
17	12.651	1805	1813	1816	rVV4	178147	441377	2.83%	0.224%
18	12.680	1816	1818	1820	rVV	186883	199081	1.28%	0.101%
19	12.727	1820	1826	1837	rVB	1812574	3229609	20.71%	1.636%
20	12.916	1850	1858	1863	rBV	974249	1645552	10.55%	0.834%
21	12.974	1863	1868	1872	rVV	4032317	6892721	44.19%	3.492%
22	13.010	1872	1874	1877	rVV	1831070	2579390	16.54%	1.307%
23	13.051	1877	1881	1887	rVV2	3129870	5659534	36.29%	2.868%
24	13.104	1887	1890	1895	rVB2	172915	257161	1.65%	0.130%
25	13.215	1900	1909	1916	rBV	2751251	4751585	30.46%	2.408%
26	13.280	1916	1920	1926	rVB2	418873	715928	4.59%	0.363%
27	13.363	1927	1934	1946	rBV	9775501	15597143	100.00%	7.903%
28	13.492	1947	1956	1961	rVB2	381176	769921	4.94%	0.390%
29	13.557	1961	1967	1971	rBV2	940784	1606563	10.30%	0.814%
30	13.604	1971	1975	1979	rBV2	426938	649405	4.16%	0.329%
31	13.668	1981	1986	1988	rBV	1994248	3085126	19.78%	1.563%
32	13.698	1988	1991	1996	rVB	4095880	6236813	39.99%	3.160%
33	13.739	1996	1998	2005	rVB2	283272	409644	2.63%	0.208%
34	13.868	2009	2020	2024	rBV2	5113613	10360307	66.42%	5.249%
35	13.992	2036	2041	2046	rBV2	2489732	4008868	25.70%	2.031%
36	14.062	2046	2053	2058	rVV	1042978	1735973	11.13%	0.880%

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Integrator: RTE

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

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37	14.127	2059	2064	2066	rVV	1370015	2136440	13.70%	1.082%
38	14.157	2066	2069	2073	rVB	1468671	2075058	13.30%	1.051%
39	14.210	2073	2078	2083	rVB	2037673	3012492	19.31%	1.526%
40	14.274	2083	2089	2092	rBV2	563292	830869	5.33%	0.421%
41	14.321	2092	2097	2103	rVB	2736352	4297362	27.55%	2.177%
42	14.386	2103	2108	2114	rVB6	236747	518793	3.33%	0.263%
43	14.457	2114	2120	2124	rBV3	1066283	2130672	13.66%	1.080%
44	14.504	2124	2128	2131	rBV3	661078	1005245	6.45%	0.509%
45	14.539	2131	2134	2137	rVV2	885563	1229258	7.88%	0.623%
46	14.574	2137	2140	2144	rVB	1773208	2315822	14.85%	1.173%
47	14.615	2144	2147	2151	rVB	804282	961633	6.17%	0.487%
48	14.657	2151	2154	2161	rVB3	545788	908862	5.83%	0.461%
49	14.721	2161	2165	2167	rBV3	256467	383334	2.46%	0.194%
50	14.757	2167	2171	2174	rVB2	488164	604747	3.88%	0.306%
51	14.804	2174	2179	2182	rBV	3014299	4720085	30.26%	2.392%
52	14.839	2182	2185	2191	rVB2	2562091	4102504	26.30%	2.079%
53	14.927	2191	2200	2206	rBV2	6433902	14066723	90.19%	7.127%
54	15.080	2219	2226	2233	rVB	5592936	9685762	62.10%	4.908%
55	15.151	2234	2238	2239	rBV3	286355	372761	2.39%	0.189%
56	15.186	2239	2244	2248	rVV3	1034340	1873945	12.01%	0.949%
57	15.227	2249	2251	2256	rVB	560890	770967	4.94%	0.391%
58	15.304	2257	2264	2273	rVB3	1293815	3331894	21.36%	1.688%
59	15.386	2274	2278	2285	rVB5	456190	997582	6.40%	0.505%
60	15.462	2285	2291	2294	rBV2	764467	1518715	9.74%	0.770%
61	15.498	2294	2297	2302	rVB	762387	1175081	7.53%	0.595%
62	15.557	2302	2307	2312	rBV	1007221	1599738	10.26%	0.811%
63	15.609	2312	2316	2318	rBV2	670091	1047948	6.72%	0.531%
64	15.698	2326	2331	2340	rVB	1934389	3795673	24.34%	1.923%
65	15.798	2340	2348	2356	rBV2	1272380	3024707	19.39%	1.533%
66	15.880	2358	2362	2368	rVB4	330973	588133	3.77%	0.298%
67	16.015	2371	2385	2394	rVB	2790105	7304281	46.83%	3.701%
68	16.104	2397	2400	2405	rVV4	143123	217843	1.40%	0.110%
69	16.186	2407	2414	2421	rVB3	607499	1555085	9.97%	0.788%
70	16.339	2430	2440	2446	rBV	1841658	4400967	28.22%	2.230%
71	16.456	2456	2460	2465	rBV	255989	443829	2.85%	0.225%
72	16.539	2468	2474	2481	rVV4	780863	1898067	12.17%	0.962%
73	16.609	2481	2486	2491	rVV	723087	1582493	10.15%	0.802%
74	16.674	2492	2497	2510	rVB	1180160	2437510	15.63%	1.235%

Sum of corrected areas: 197363655

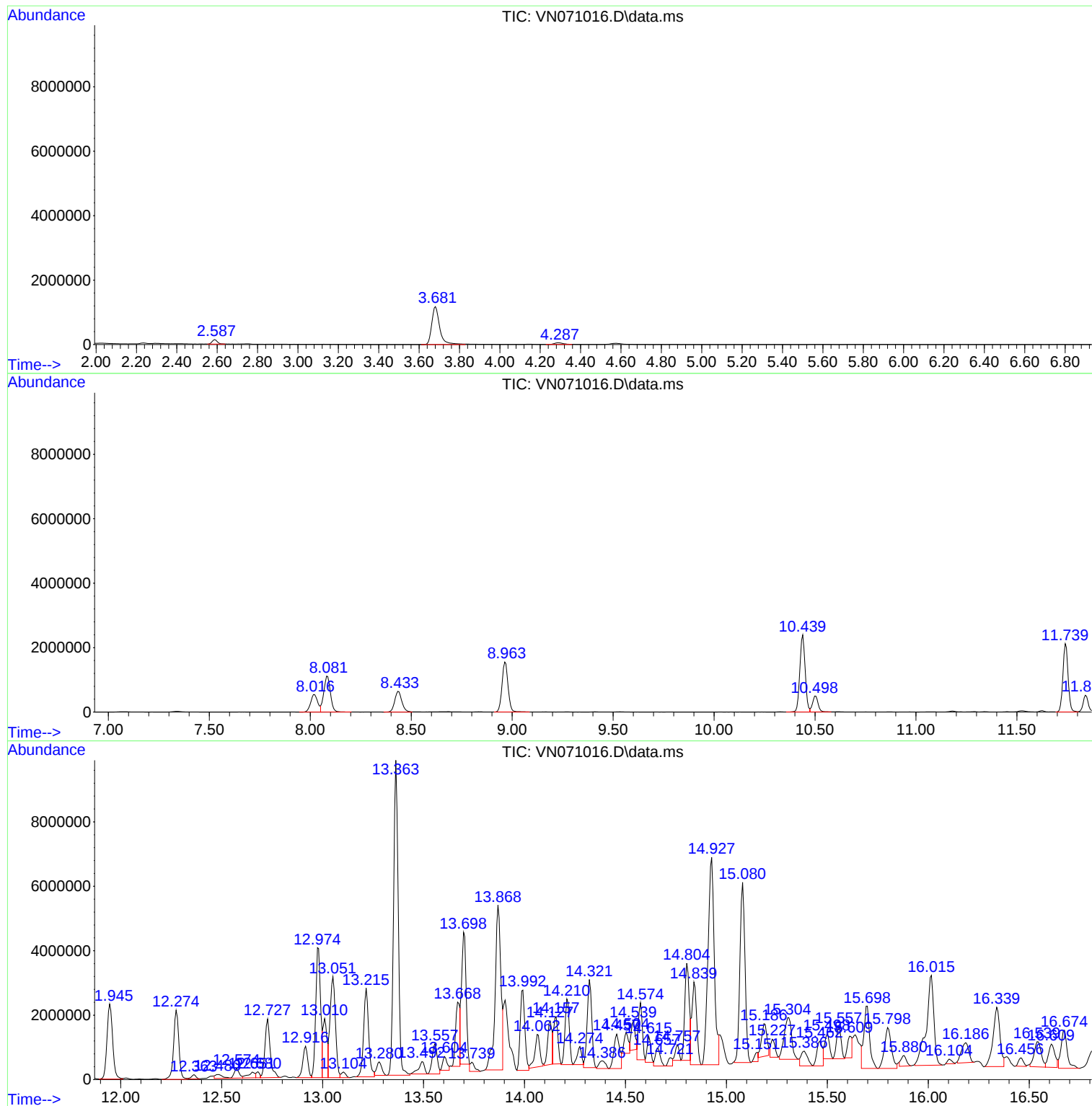
Instrument :
MSVOA_N
ClientSampleId :
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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
Quant Title : SW846 8260

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



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Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
1070

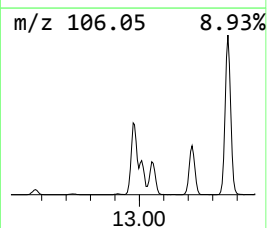
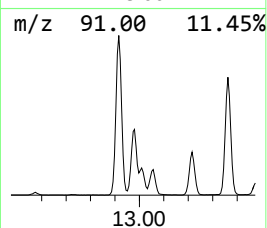
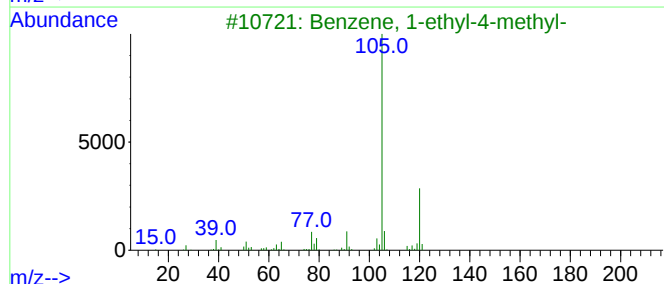
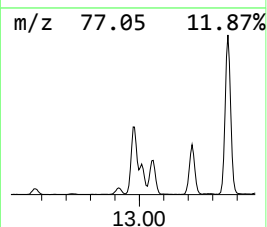
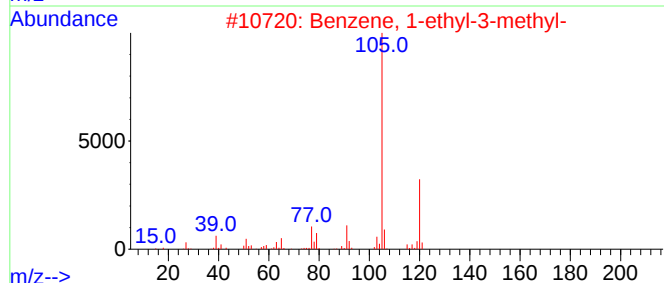
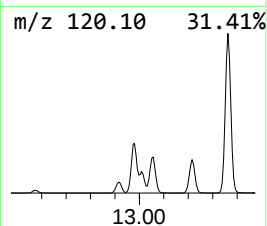
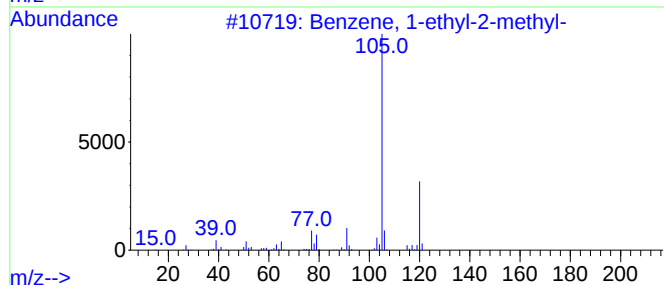
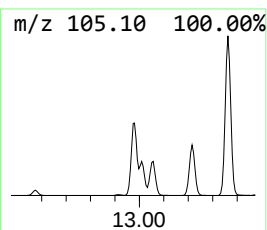
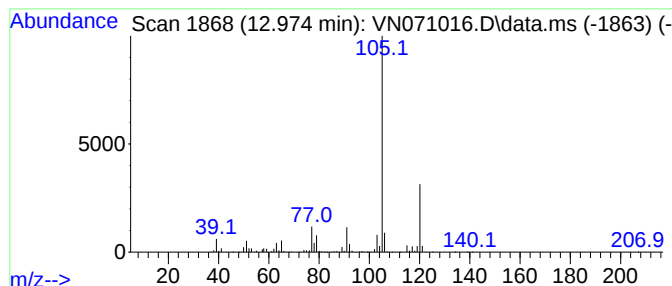
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.974	111.71 ug/l	6892720	1,4-Dichlorobenzene-d4	13.668

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			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



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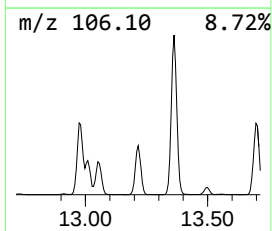
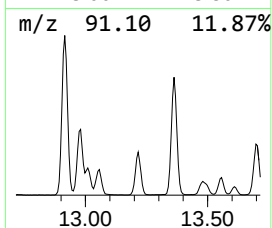
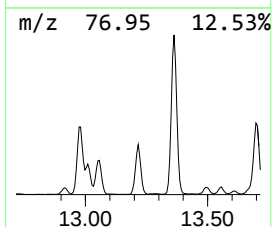
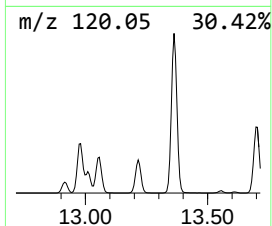
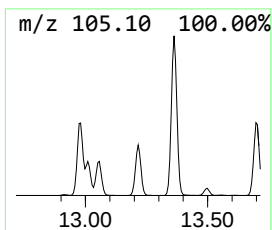
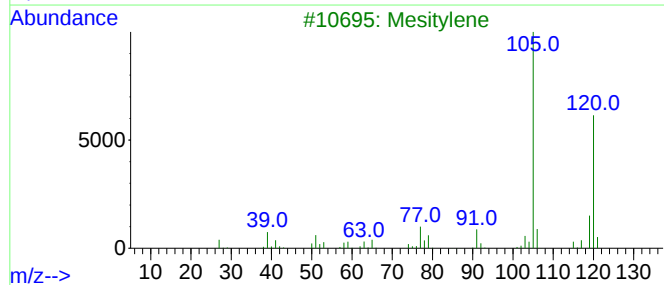
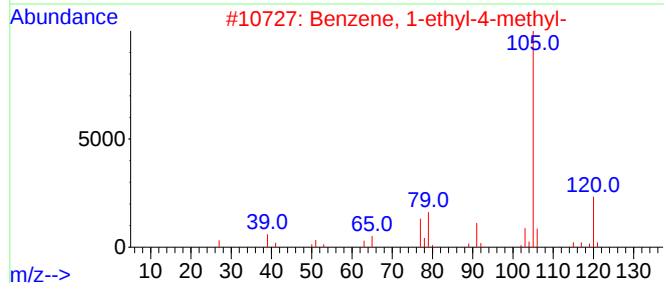
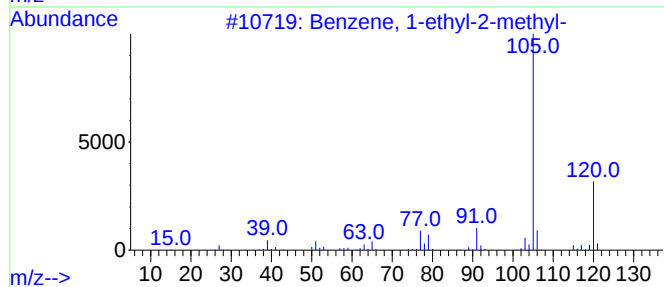
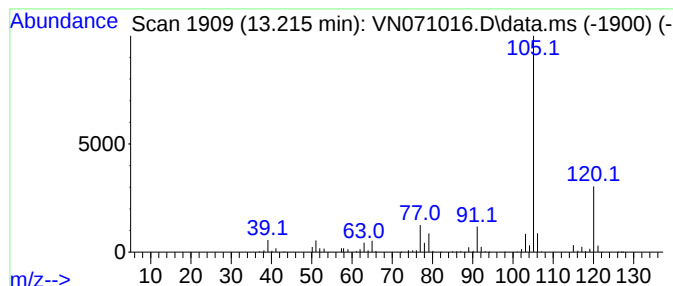
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	77.01 ug/l	4751590	1,4-Dichlorobenzene-d4	13.668

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	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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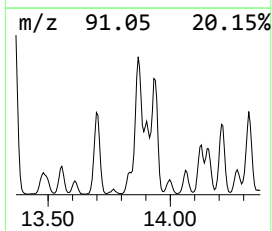
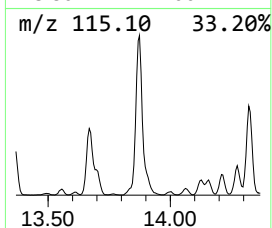
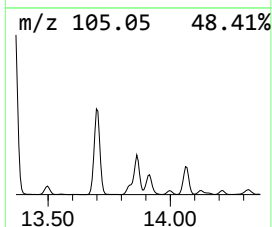
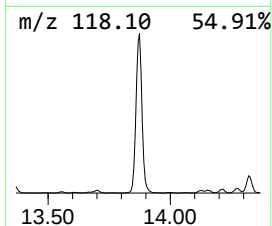
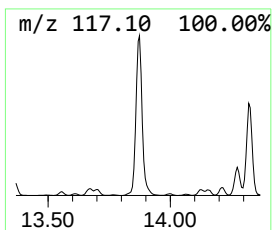
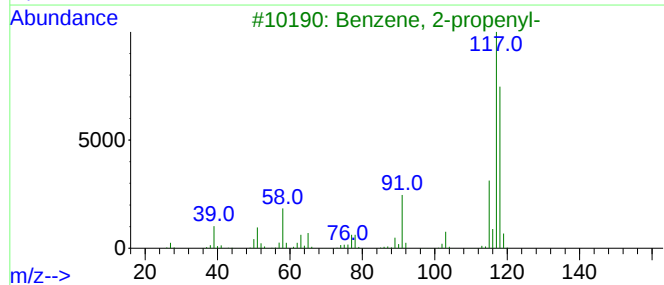
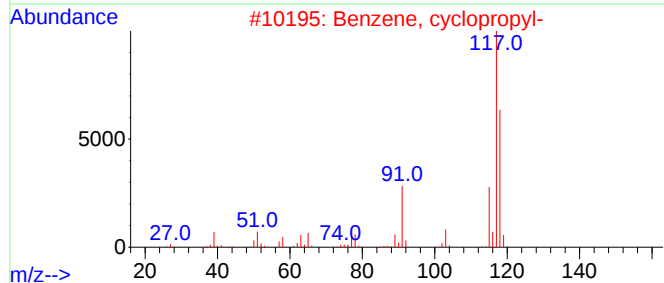
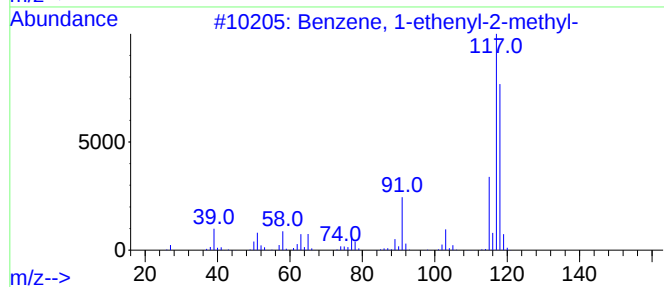
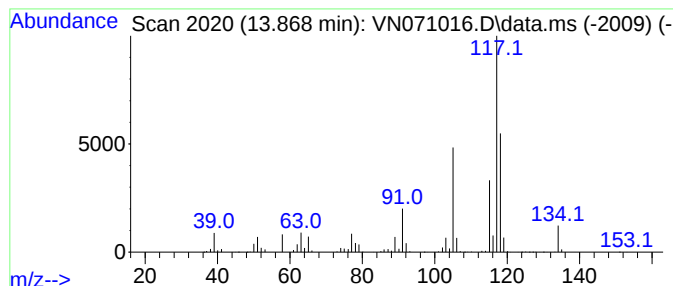
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.868	167.91 ug/l	10360300	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4			Indane	118	C9H10	000496-11-7	76
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64



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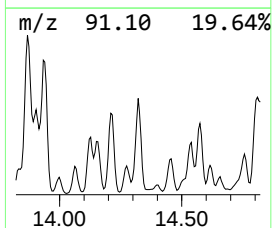
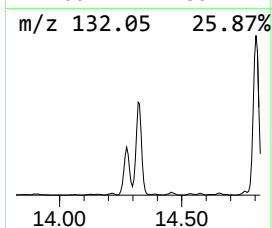
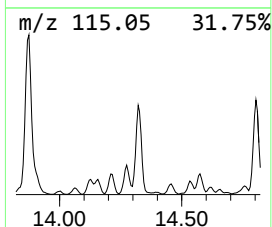
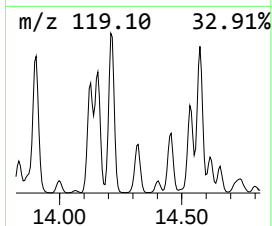
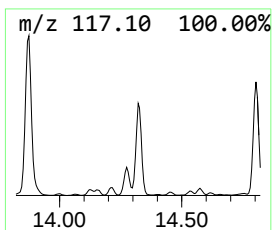
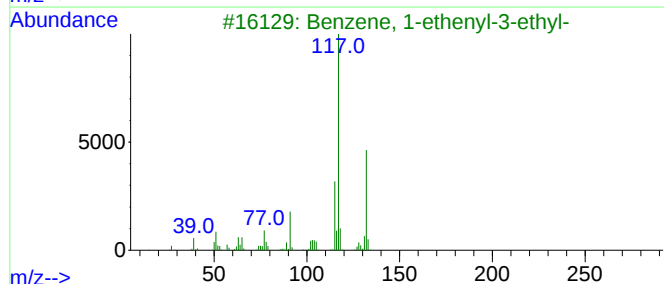
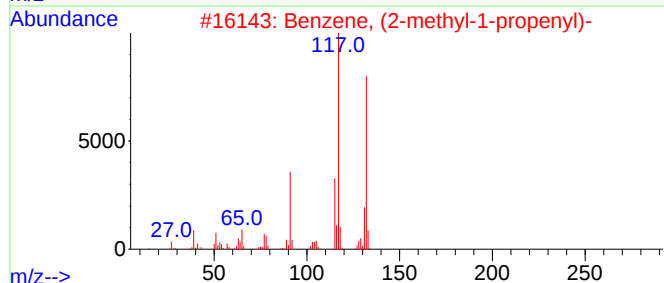
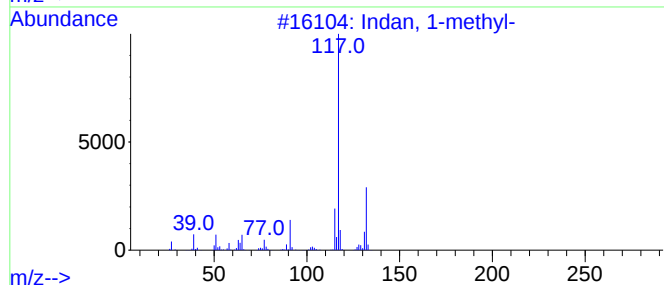
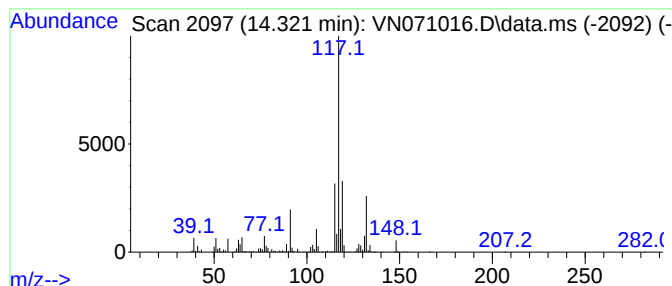
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Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.321	69.65 ug/l	4297360	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	76
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	74
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
1070

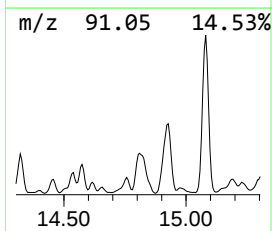
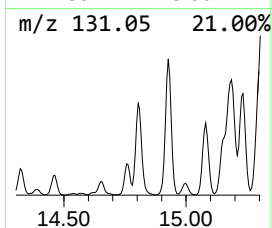
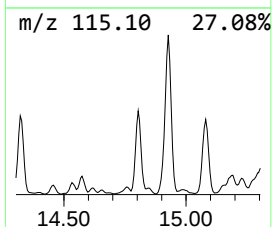
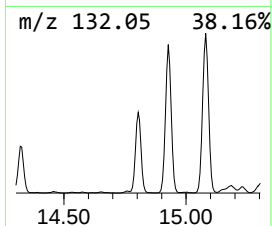
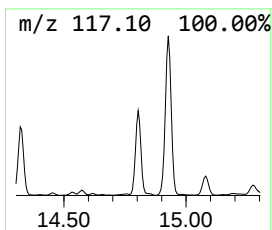
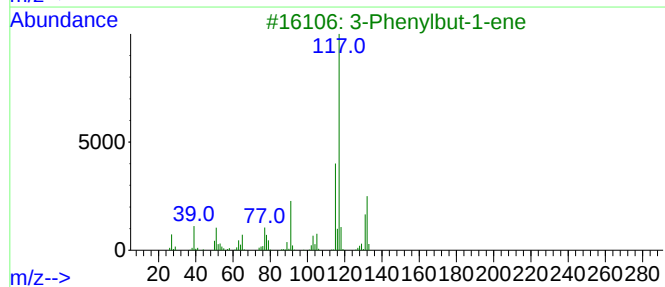
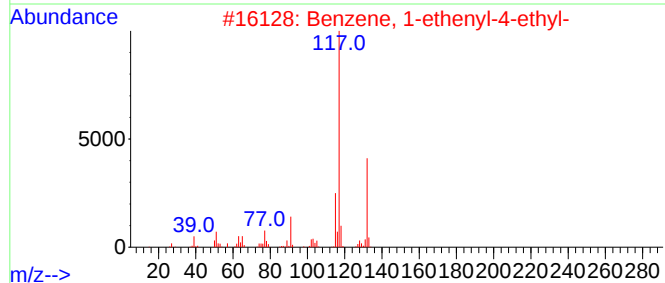
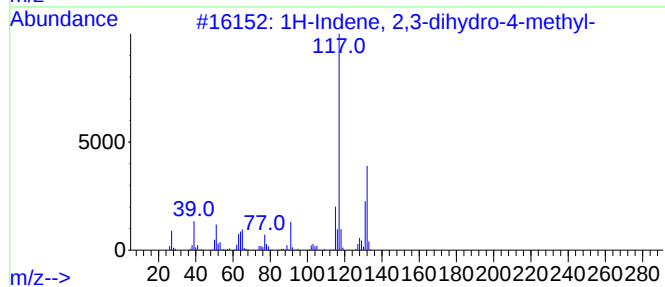
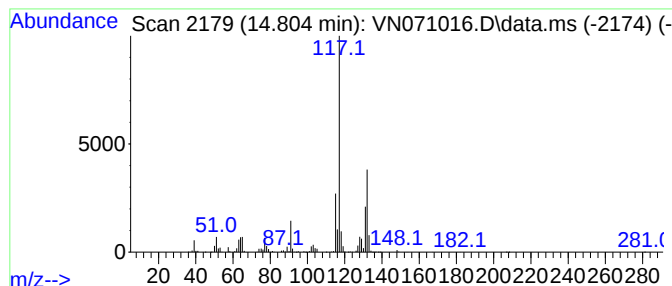
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	76.50 ug/l	4720090	1,4-Dichlorobenzene-d4	13.668

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	93
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
5			Indan, 1-methyl-	132	C10H12	000767-58-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
1070

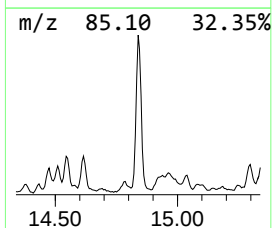
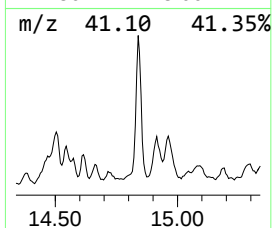
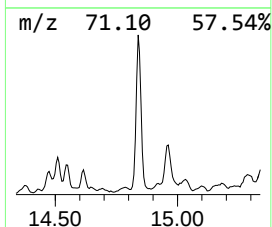
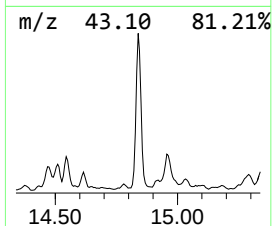
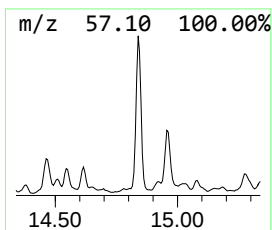
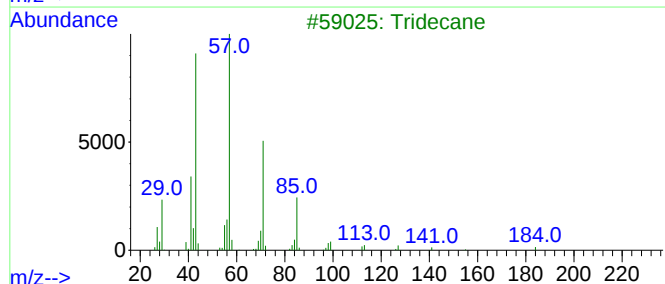
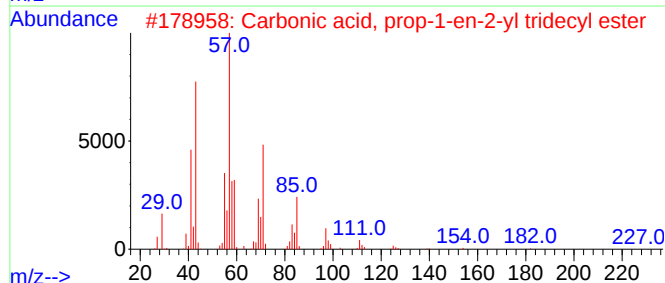
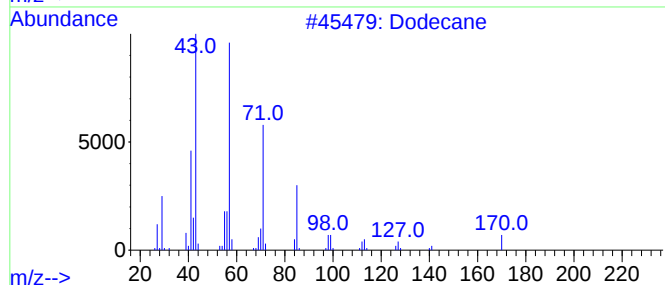
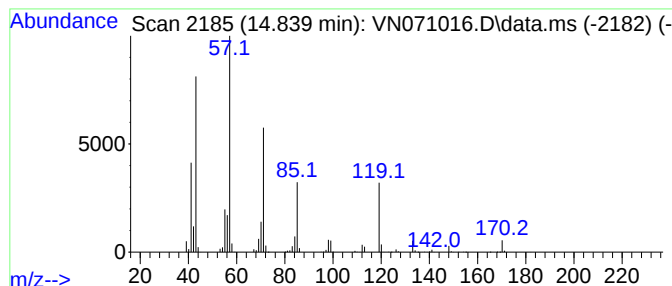
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Quant Title : SW846 8260

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

Peak Number 6 Dodecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.839	66.49 ug/l	4102500	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	89
2			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	58
3			Tridecane	184	C13H28	000629-50-5	53
4			Carbonic acid, dodecyl vinyl ester	256	C15H28O3	1000382-54-8	53
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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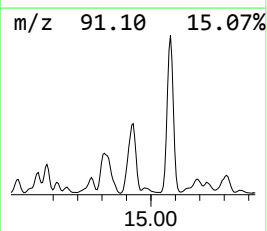
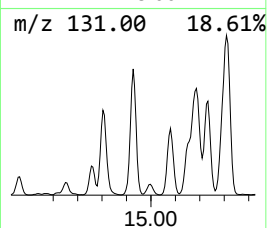
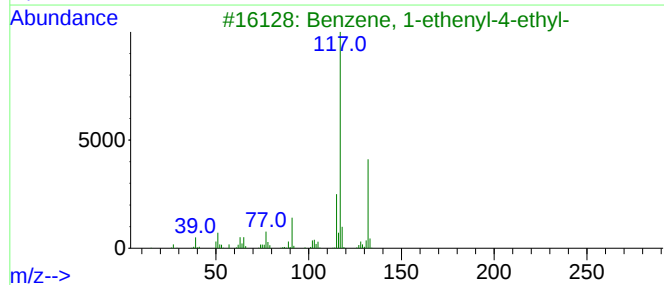
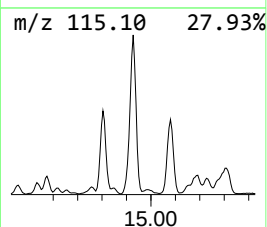
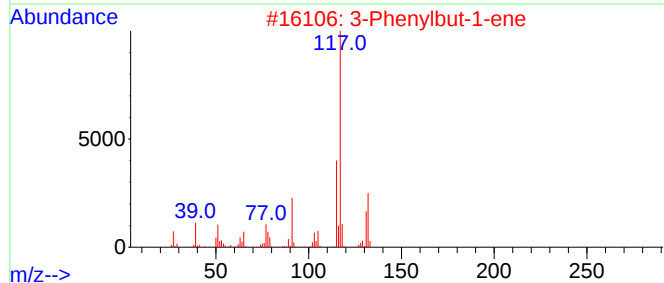
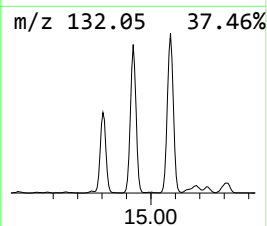
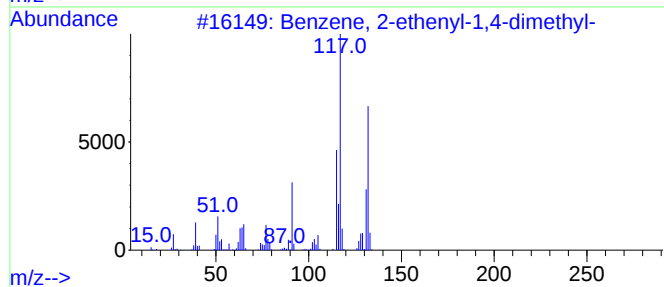
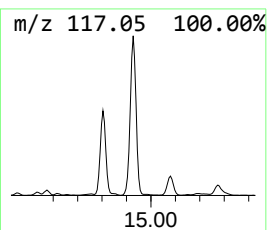
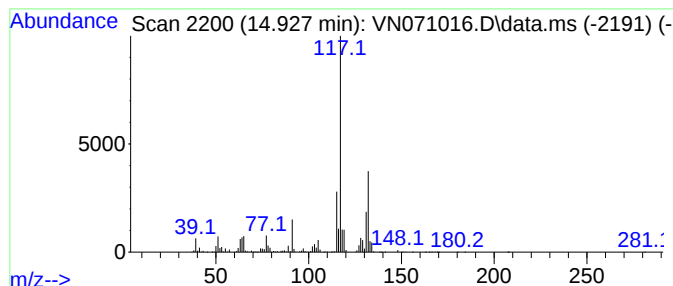
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	227.98 ug/l	14066700	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	96
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	93
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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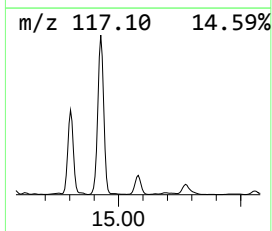
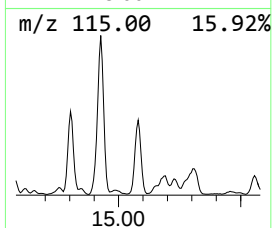
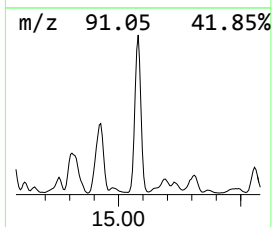
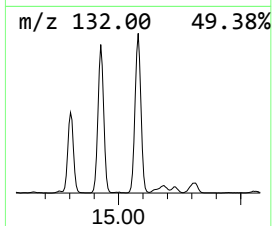
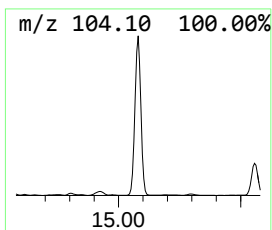
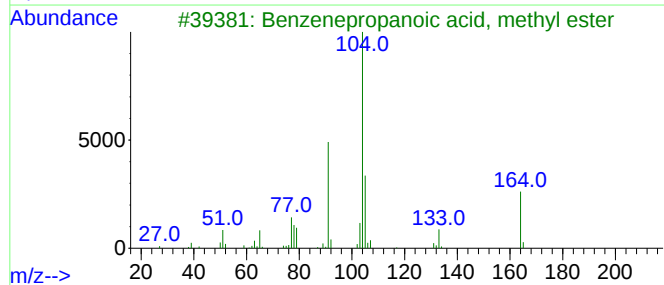
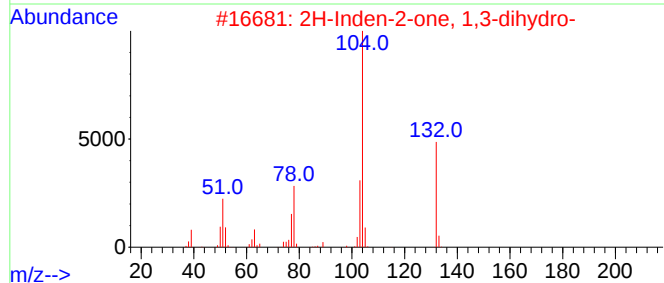
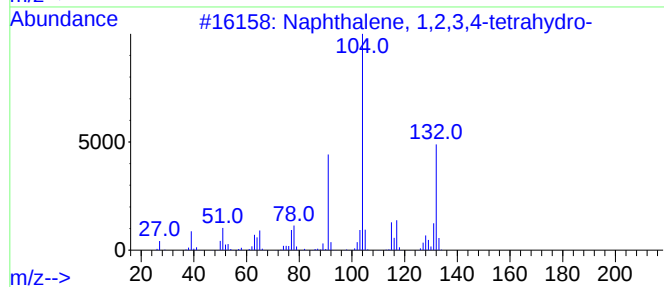
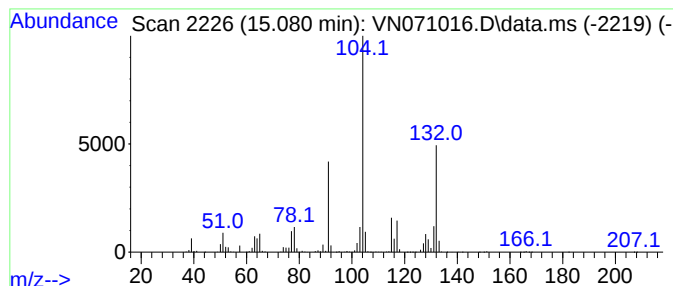
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.080	156.97 ug/l	9685760	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	97
2			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
3			Benzenepropanoic acid, methyl ester	164	C10H12O2	000103-25-3	43
4			5,5-Dimethyl-4-phenyldihydrofura...	190	C12H14O2	013133-96-5	43
5			Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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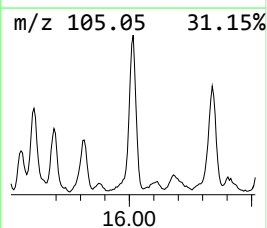
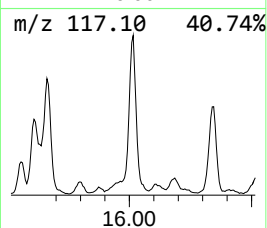
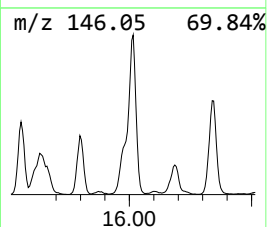
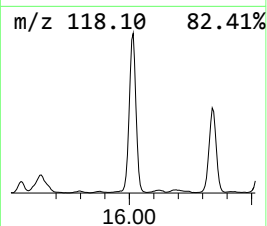
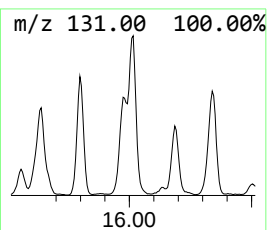
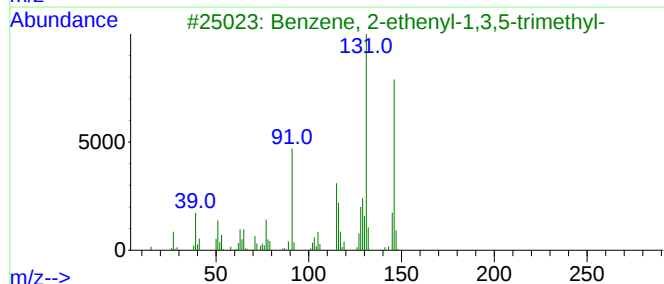
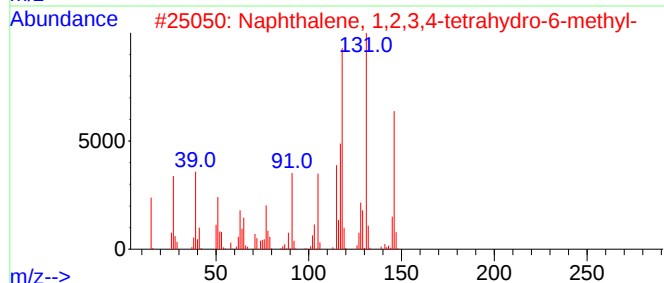
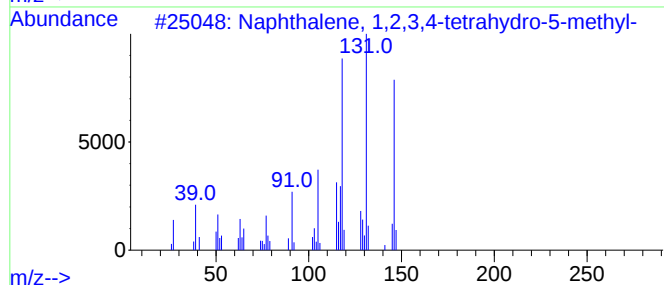
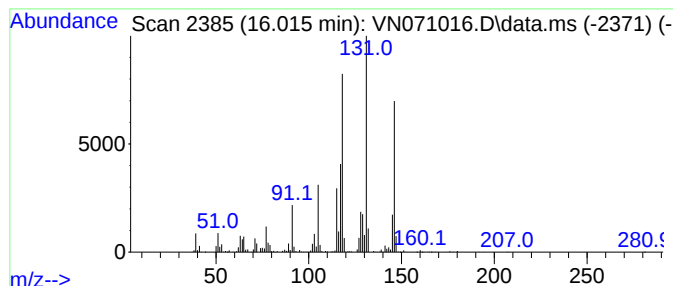
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.015	118.38 ug/l	7304280	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	97
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	97
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	62
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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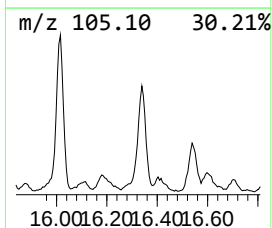
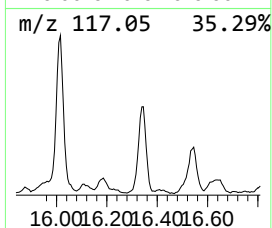
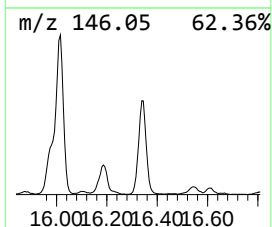
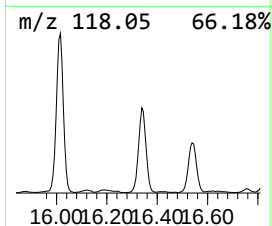
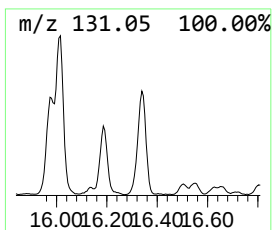
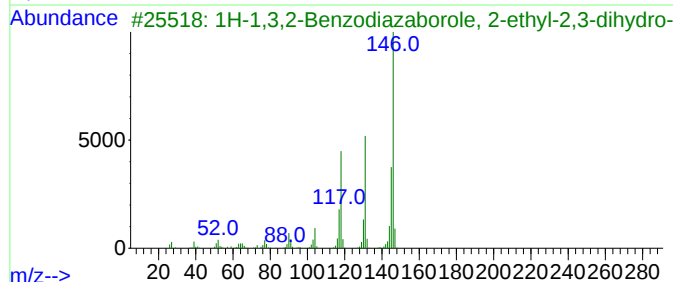
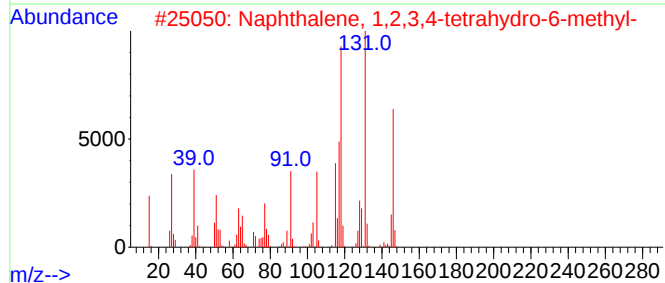
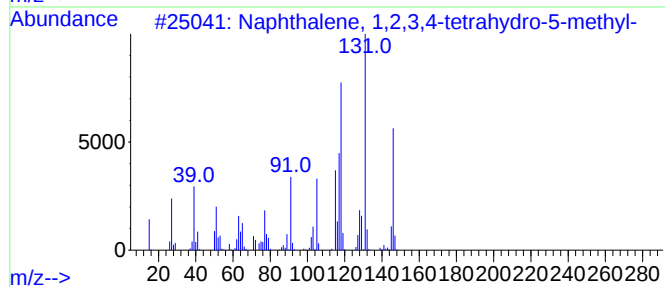
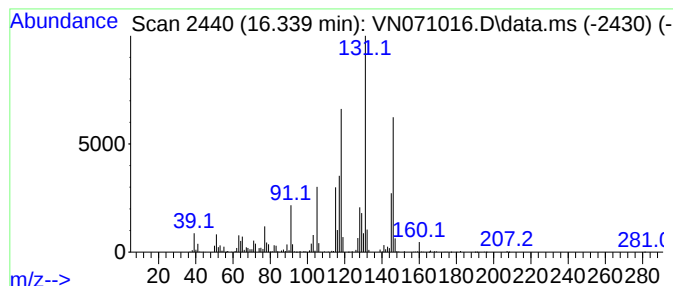
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.339	71.33 ug/l	4400970	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	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	90
4			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	64
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
Data File : VN071016.D
Acq On : 18 Feb 2022 21:27
Operator : JC\MD
Sample : N1611-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.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
Benzene, 1-ethy...	12.974	111.7	ug/l	6892720	4	13.668	3085130	50.0
Benzene, 1-ethy...	13.216	77.0	ug/l	4751590	4	13.668	3085130	50.0
Benzene, 1-ethe...	13.868	167.9	ug/l	10360300	4	13.668	3085130	50.0
Indan, 1-methyl-	14.321	69.7	ug/l	4297360	4	13.668	3085130	50.0
1H-Indene, 2,3-...	14.804	76.5	ug/l	4720090	4	13.668	3085130	50.0
Dodecane	14.839	66.5	ug/l	4102500	4	13.668	3085130	50.0
Benzene, 2-ethe...	14.927	228.0	ug/l	14066700	4	13.668	3085130	50.0
Naphthalene, 1,...	15.080	157.0	ug/l	9685760	4	13.668	3085130	50.0
Naphthalene, 1,...	16.015	118.4	ug/l	7304280	4	13.668	3085130	50.0
Naphthalene, 1,...	16.339	71.3	ug/l	4400970	4	13.668	3085130	50.0