

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
 Data File : VN074786.D
 Acq On : 06 Oct 2022 16:16
 Operator : JC\MD
 Sample : N4910-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3

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\82N100522W.M
 Title : SW846 8260

Signal : TIC: VN074786.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.318	38	45	52	rBV	256882	720271	4.09%	0.332%
2	2.530	77	81	87	rBV2	152494	332579	1.89%	0.153%
3	3.253	196	204	222	rVB	1462719	3840160	21.83%	1.770%
4	3.659	262	273	286	rBV2	637410	1579769	8.98%	0.728%
5	5.365	539	563	579	rBV3	938527	4521551	25.70%	2.084%
6	5.830	628	642	657	rBV	642820	2176032	12.37%	1.003%
7	6.330	713	727	744	rBV	1015550	3088394	17.56%	1.423%
8	6.671	779	785	797	rVB2	87184	267129	1.52%	0.123%
9	7.236	872	881	889	rBV3	99080	280649	1.60%	0.129%
10	7.341	889	899	910	rVV	2092447	5846660	33.24%	2.694%
11	7.453	910	918	925	rVV	122479	329877	1.88%	0.152%
12	8.018	1007	1014	1024	rVB2	147182	370297	2.11%	0.171%
13	8.183	1032	1042	1045	rBV	423687	1022957	5.82%	0.471%
14	8.265	1045	1056	1064	rVV2	2533749	8851185	50.32%	4.079%
15	8.335	1064	1068	1079	rVB2	496574	1202480	6.84%	0.554%
16	8.453	1079	1088	1095	rBV	1079689	2478833	14.09%	1.142%
17	8.518	1095	1099	1105	rVV2	394208	827275	4.70%	0.381%
18	8.612	1105	1115	1126	rVV	6159527	14973749	85.12%	6.900%
19	8.730	1126	1135	1141	rVV4	1172417	2939243	16.71%	1.354%
20	8.800	1141	1147	1152	rVV	1033752	2280128	12.96%	1.051%
21	8.865	1152	1158	1168	rVB	1655443	3694957	21.01%	1.703%
22	8.977	1168	1177	1189	rVB	1470654	3073980	17.48%	1.417%
23	9.106	1191	1199	1206	rBV3	933308	2108524	11.99%	0.972%
24	9.159	1206	1208	1220	rVB3	167064	298657	1.70%	0.138%
25	9.606	1268	1284	1302	rBV	7345373	17590333	100.00%	8.106%
26	9.782	1302	1314	1321	rBV	969239	1867999	10.62%	0.861%
27	9.871	1321	1329	1336	rVB2	532525	1070287	6.08%	0.493%
28	10.012	1345	1353	1362	rVB4	787924	1731212	9.84%	0.798%
29	10.106	1363	1369	1373	rBV2	282678	538663	3.06%	0.248%
30	10.212	1381	1387	1401	rVB2	795305	1916602	10.90%	0.883%
31	10.347	1401	1410	1420	rVB3	746285	1749710	9.95%	0.806%
32	10.459	1421	1429	1433	rBV	426459	851951	4.84%	0.393%
33	10.571	1440	1448	1453	rBV	3923383	7615115	43.29%	3.509%
34	10.612	1453	1455	1465	rVV2	977240	2048366	11.64%	0.944%
35	10.694	1465	1469	1472	rVV2	286942	503594	2.86%	0.232%
36	10.747	1472	1478	1486	rVV5	1965080	4337798	24.66%	1.999%

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37	10.865	1492	1498	1501	rVV4	144019	296767	1.69%	0.137%
38	10.918	1501	1507	1513	rVB	1301064	2398403	13.63%	1.105%
39	11.006	1513	1522	1529	rBV4	807722	2276332	12.94%	1.049%
40	11.094	1532	1537	1542	rVB5	186102	381436	2.17%	0.176%
41	11.176	1542	1551	1557	rBV4	275017	589189	3.35%	0.272%
42	11.265	1557	1566	1569	rBV3	230444	661545	3.76%	0.305%
43	11.300	1569	1572	1576	rVV2	208809	323000	1.84%	0.149%
44	11.347	1576	1580	1583	rVV2	390772	653398	3.71%	0.301%
45	11.423	1583	1593	1597	rVV	1691613	3645785	20.73%	1.680%
46	11.476	1597	1602	1607	rVV	1222810	2145949	12.20%	0.989%
47	11.529	1607	1611	1615	rVV4	178161	269384	1.53%	0.124%
48	11.576	1615	1619	1624	rVB3	431874	691781	3.93%	0.319%
49	11.647	1624	1631	1633	rBV3	304483	577796	3.28%	0.266%
50	11.676	1633	1636	1644	rVB3	423968	742948	4.22%	0.342%
51	11.747	1644	1648	1652	rBV	326636	460239	2.62%	0.212%
52	11.871	1662	1669	1674	rBV	1312377	2373308	13.49%	1.094%
53	11.965	1680	1685	1690	rBV	1608584	2718788	15.46%	1.253%
54	12.070	1695	1703	1709	rBV	4560145	9249298	52.58%	4.262%
55	12.282	1733	1739	1748	rVB6	242777	600616	3.41%	0.277%
56	12.376	1748	1755	1763	rBV4	568332	1540187	8.76%	0.710%
57	12.482	1769	1773	1778	rVB2	376921	636665	3.62%	0.293%
58	12.582	1778	1790	1794	rBV3	975351	2345589	13.33%	1.081%
59	12.694	1805	1809	1815	rBV	454098	710349	4.04%	0.327%
60	12.853	1830	1836	1845	rVB2	1432449	2683561	15.26%	1.237%
61	12.935	1845	1850	1857	rBV3	321067	539739	3.07%	0.249%
62	13.035	1860	1867	1872	rBV2	895800	1703535	9.68%	0.785%
63	13.100	1872	1878	1882	rBV	2335594	4238193	24.09%	1.953%
64	13.135	1882	1884	1886	rVV	1079414	1311568	7.46%	0.604%
65	13.176	1886	1891	1897	rVB	2823771	5079623	28.88%	2.341%
66	13.335	1910	1918	1925	rBV	1062068	2171378	12.34%	1.001%
67	13.400	1925	1929	1932	rBV2	282067	421105	2.39%	0.194%
68	13.482	1938	1943	1950	rVB	4936496	8138442	46.27%	3.750%
69	13.617	1956	1966	1971	rVB4	402685	884927	5.03%	0.408%
70	13.676	1971	1976	1981	rBV2	1103636	1899234	10.80%	0.875%
71	13.723	1981	1984	1991	rVV3	413868	814211	4.63%	0.375%
72	13.794	1991	1996	1998	rVV	1789702	2856546	16.24%	1.316%
73	13.823	1998	2001	2006	rVB	1832713	2686629	15.27%	1.238%
74	13.953	2017	2023	2025	rBV	580134	980144	5.57%	0.452%
75	13.988	2025	2029	2032	rVV2	1725861	3127343	17.78%	1.441%
76	14.023	2032	2035	2045	rVB3	1535320	3113390	17.70%	1.435%

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77	14.112	2045	2050	2057	rBV3	781187	1576039	8.96%	0.726%
78	14.188	2057	2063	2068	rVV	710017	1171616	6.66%	0.540%
79	14.247	2068	2073	2075	rVV	871313	1372236	7.80%	0.632%
80	14.276	2075	2078	2083	rVB	1035472	1531075	8.70%	0.706%
81	14.335	2083	2088	2093	rBV	1230999	1837612	10.45%	0.847%
82	14.400	2095	2099	2101	rBV	259171	350121	1.99%	0.161%
83	14.447	2101	2107	2112	rVB3	1797251	3067197	17.44%	1.413%
84	14.576	2124	2129	2134	rBV2	565462	976963	5.55%	0.450%
85	14.629	2134	2138	2139	rVV2	313884	356904	2.03%	0.164%
86	14.659	2140	2143	2146	rVV2	597704	934429	5.31%	0.431%
87	14.694	2146	2149	2153	rVB	918544	1214046	6.90%	0.559%
88	14.735	2153	2156	2160	rVB2	414457	480688	2.73%	0.222%
89	14.776	2160	2163	2166	rBV2	254267	375030	2.13%	0.173%
90	14.882	2173	2181	2185	rBV3	383677	711510	4.04%	0.328%
91	14.935	2185	2190	2192	rVV4	525134	998511	5.68%	0.460%
92	14.964	2192	2195	2201	rVB2	678255	1088699	6.19%	0.502%
93	15.041	2201	2208	2225	rBV2	1768347	5367642	30.51%	2.474%
94	15.217	2232	2238	2244	rBV3	212316	427211	2.43%	0.197%
95	15.323	2244	2256	2260	rBV3	578132	1521928	8.65%	0.701%
96	15.447	2269	2277	2283	rVB2	723910	1804358	10.26%	0.831%
97	15.594	2297	2302	2305	rBV4	164706	313197	1.78%	0.144%
98	15.641	2305	2310	2316	rVB2	516019	1007343	5.73%	0.464%
99	15.753	2323	2329	2334	rBV3	176325	318891	1.81%	0.147%
100	15.947	2354	2362	2368	rBV5	138105	334027	1.90%	0.154%

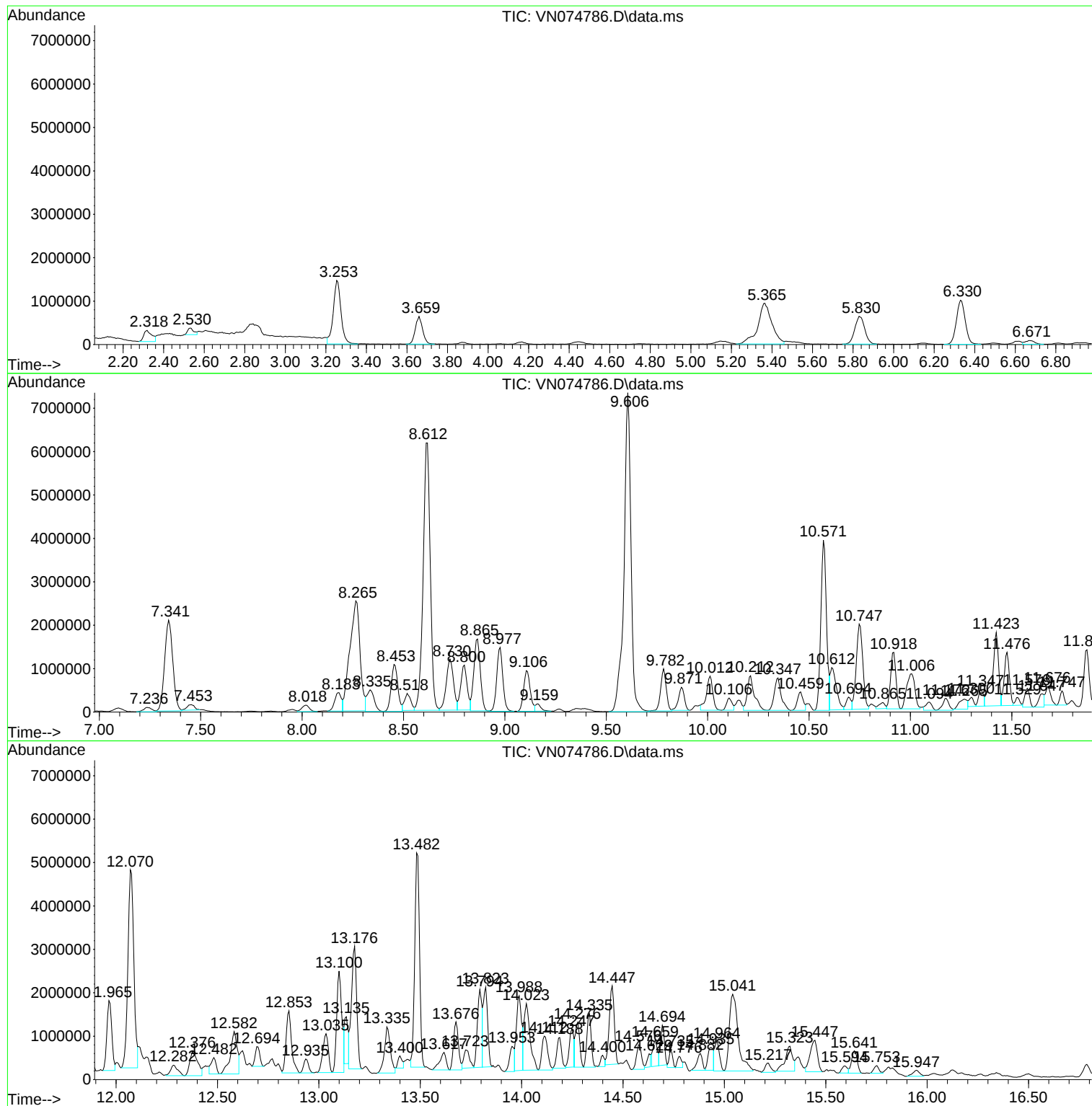
Sum of corrected areas: 217002559

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100522W.M
Quant Title : SW846 8260

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



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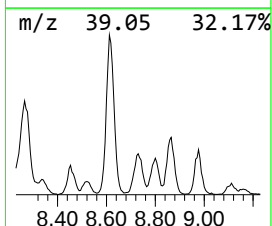
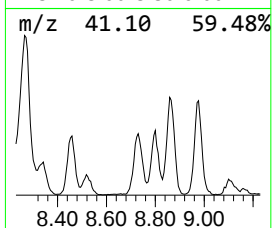
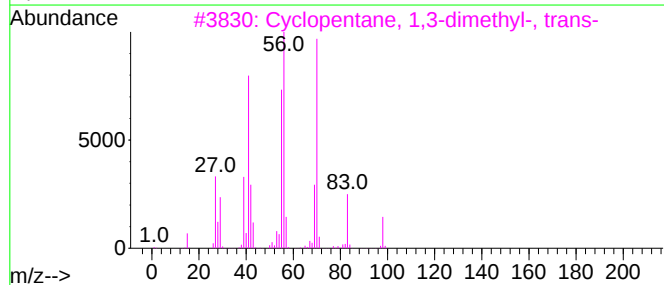
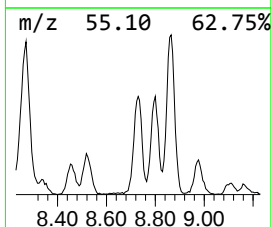
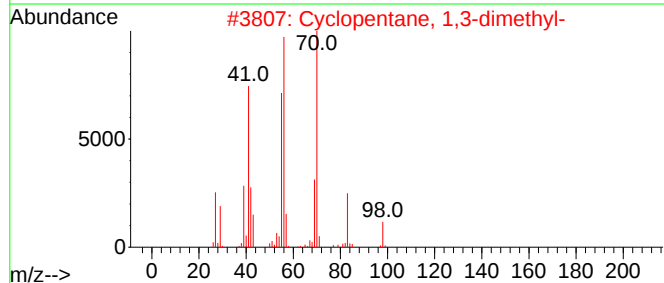
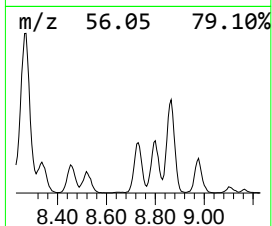
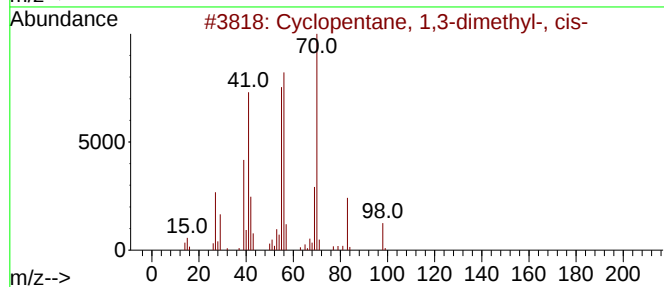
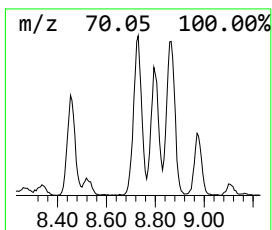
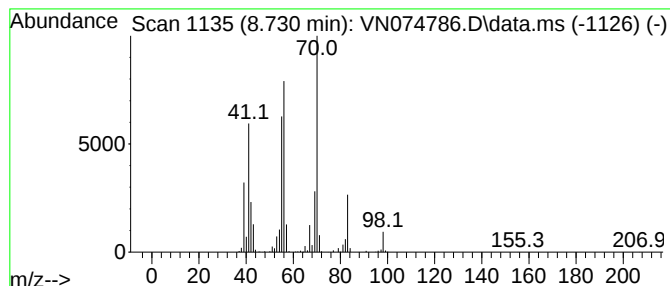
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100522W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, 1,3-dimethyl-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.730	69.70 ug/l	2939240	1,4-Difluorobenzene	9.106

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	95
2			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	94
3			Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	90
4			1-Pentene, 3,4-dimethyl-	98	C7H14	007385-78-6	72
5			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	70



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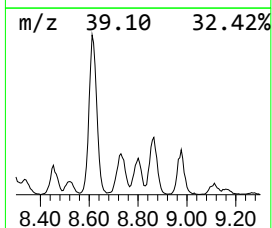
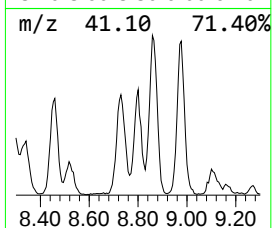
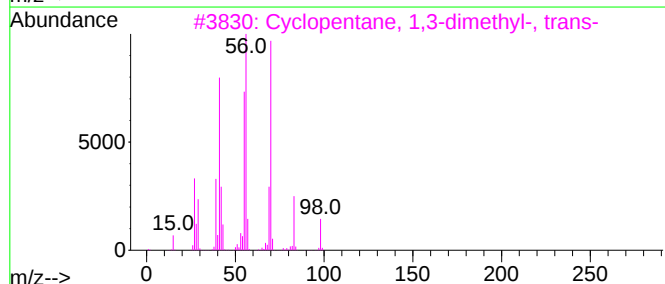
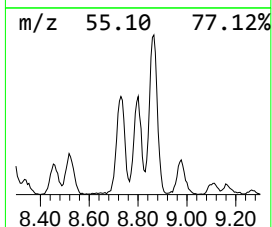
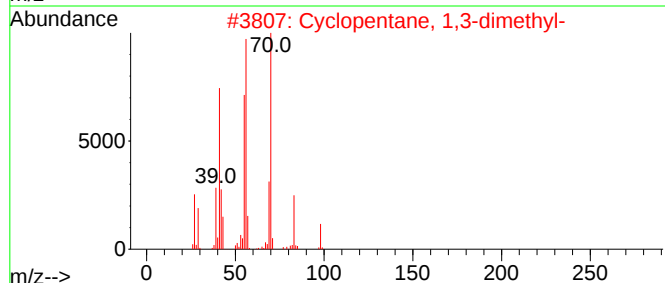
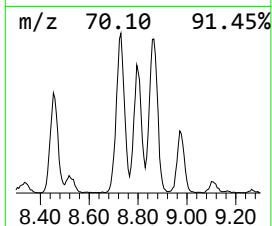
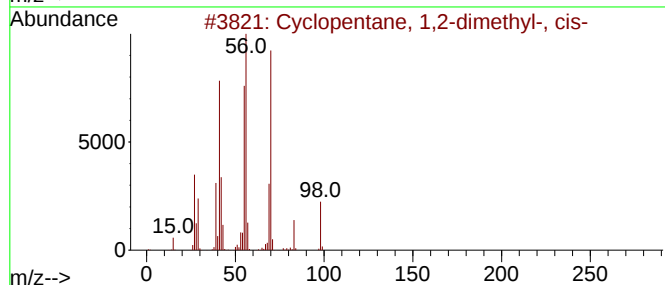
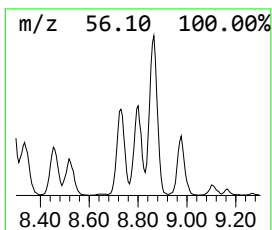
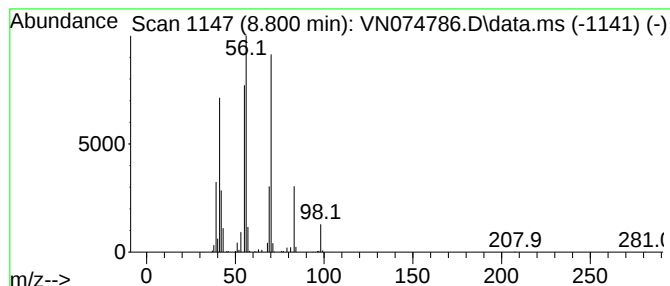
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100522W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, 1,2-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.800	54.07 ug/l	2280130	1,4-Difluorobenzene	9.106

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	91
2			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	91
3			Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	91
4			Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90
5			Cycloheptane	98	C7H14	000291-64-5	86



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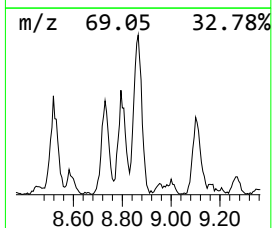
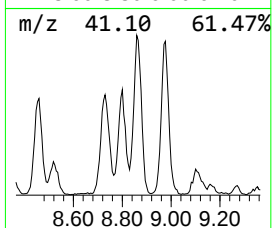
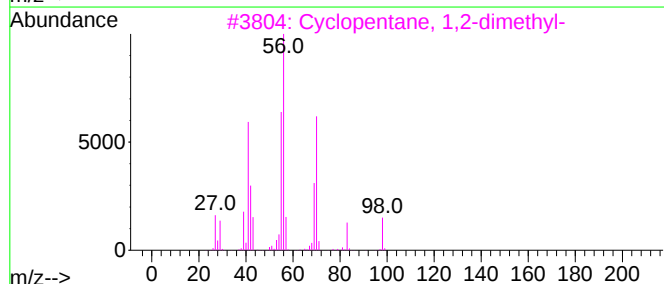
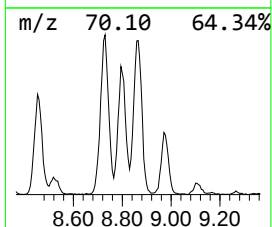
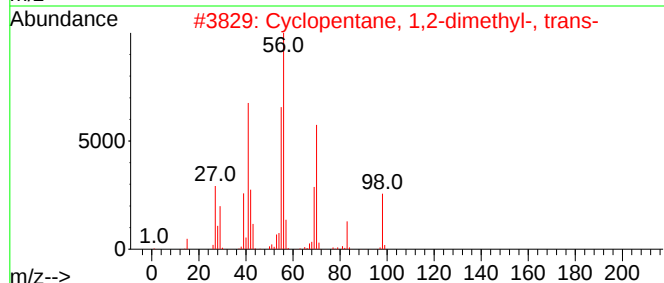
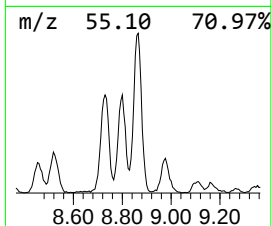
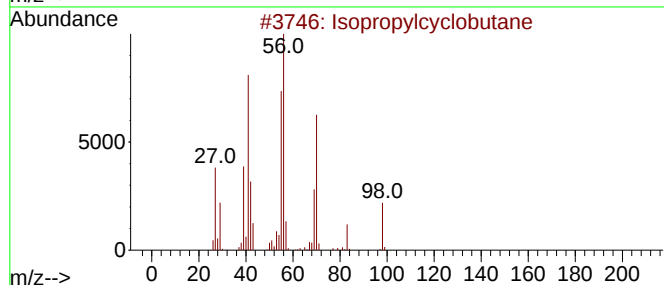
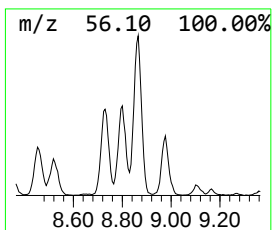
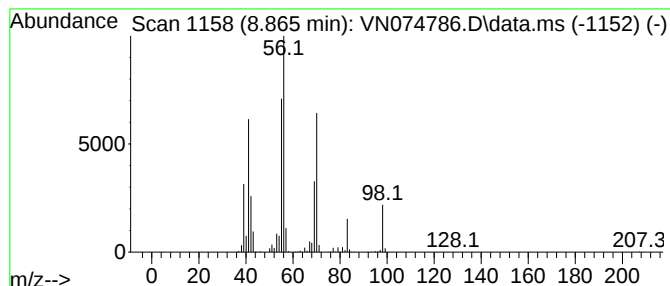
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Quant Title : SW846 8260

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

Peak Number 3 Isopropylcyclobutane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.865	87.62 ug/l	3694960	1,4-Difluorobenzene	9.106

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropylcyclobutane	98	C7H14	000872-56-0	95
2			Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	95
3			Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	90
4			1-Heptene	98	C7H14	000592-76-7	90
5			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

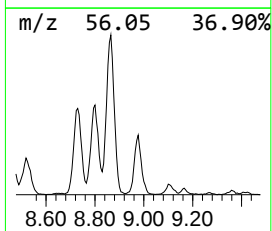
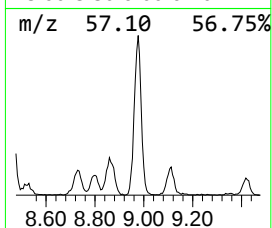
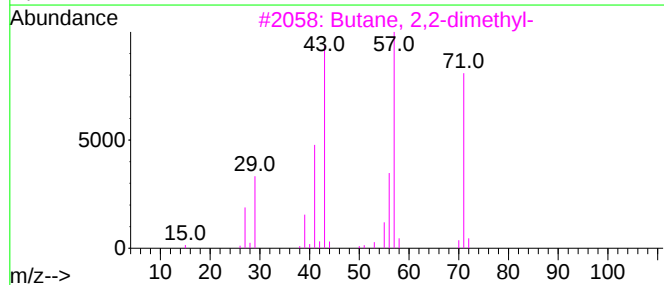
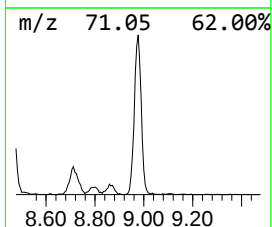
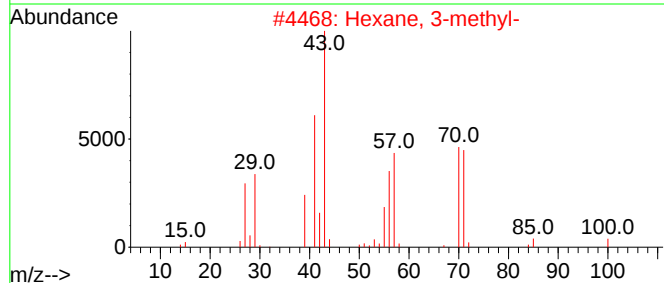
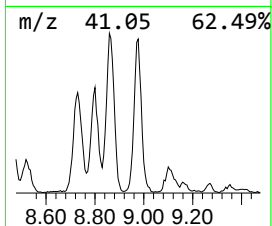
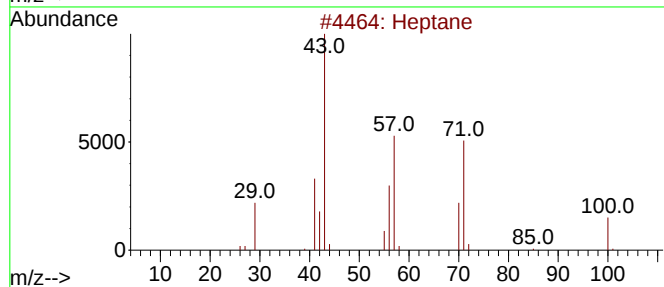
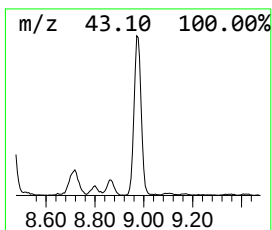
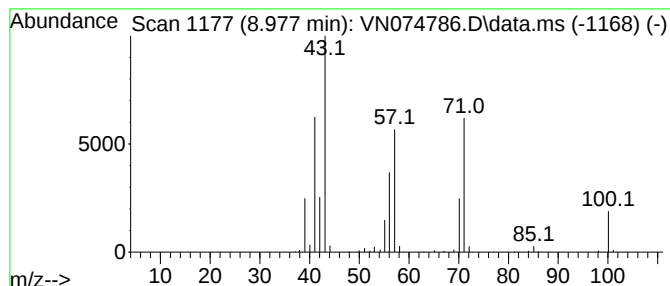
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Quant Title : SW846 8260

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

Peak Number 4 Heptane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.977	72.89 ug/l	3073980	1,4-Difluorobenzene	9.106

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	59
3			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	38
4			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	35
5			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

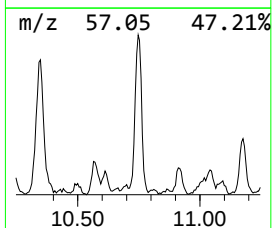
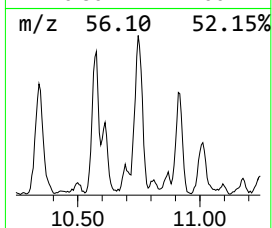
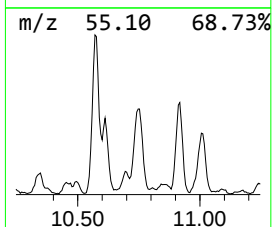
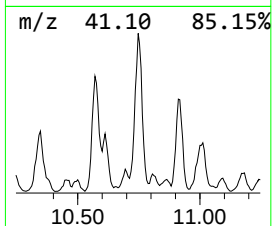
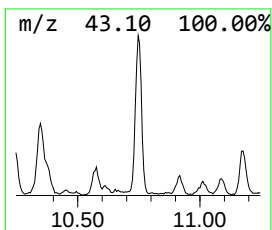
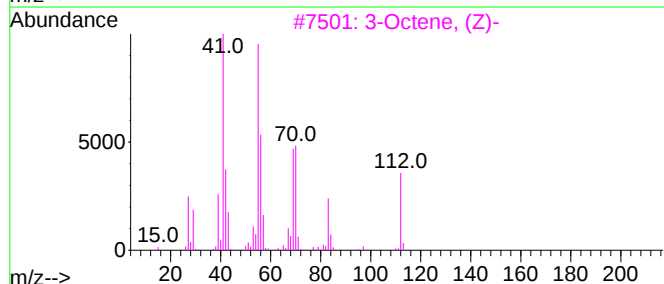
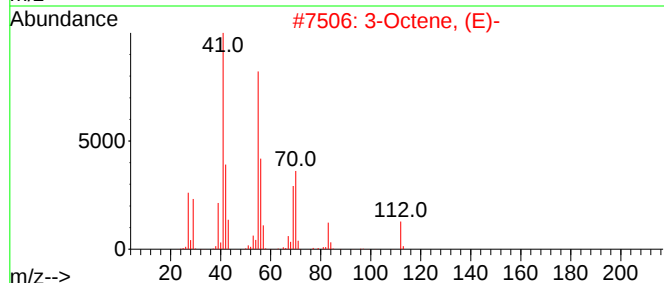
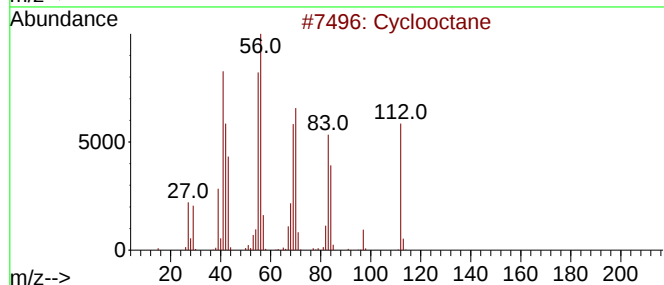
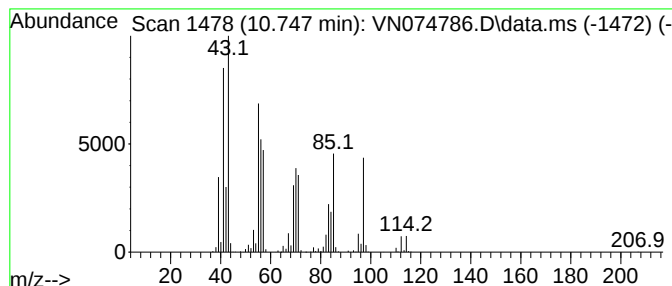
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Quant Title : SW846 8260

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

Peak Number 5 unknown10.747 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.747	91.39 ug/l	4337800	Chlorobenzene-d5	11.871

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclooctane	112	C8H16	000292-64-8	46
2			3-Octene, (E)-	112	C8H16	014919-01-8	45
3			3-Octene, (Z)-	112	C8H16	014850-22-7	45
4			2-Heptenal, (Z)-	112	C7H12O	057266-86-1	43
5			Cycloheptane, methyl-	112	C8H16	004126-78-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

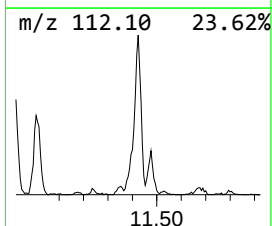
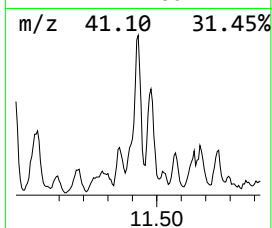
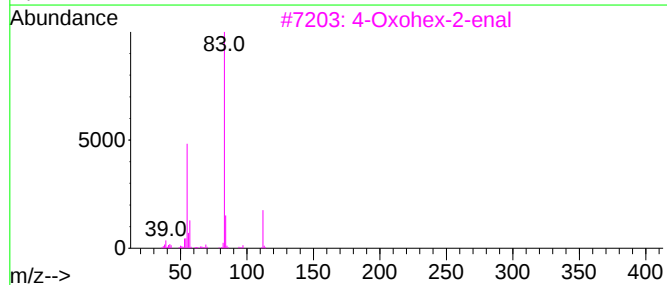
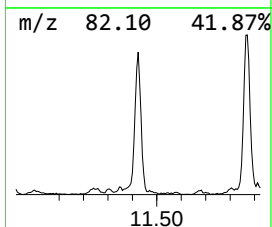
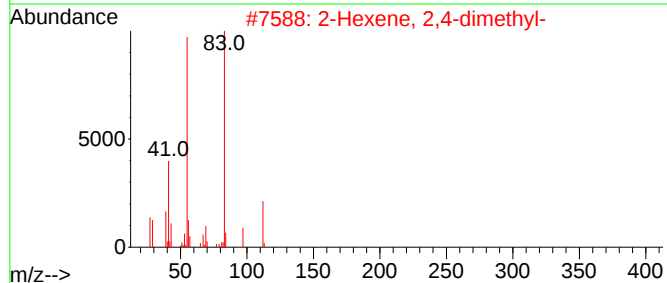
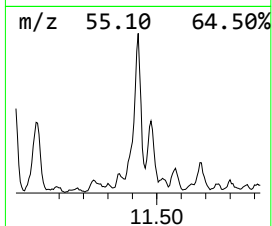
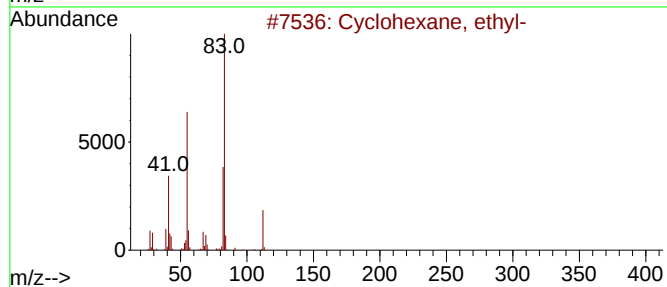
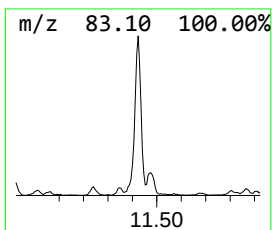
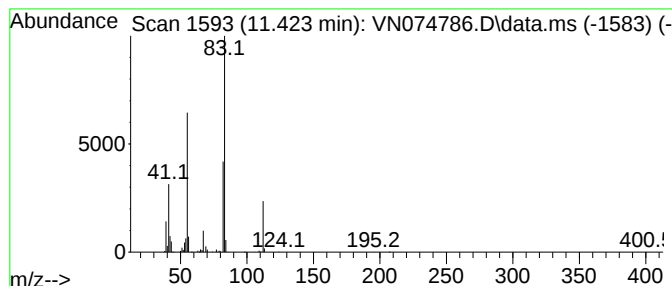
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Quant Title : SW846 8260

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

Peak Number 6 Cyclohexane, ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.424	76.81 ug/l	3645790	Chlorobenzene-d5	11.871

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	90
2			2-Hexene, 2,4-dimethyl-	112	C8H16	014255-23-3	53
3			4-Oxohex-2-enal	112	C6H8O2	020697-55-6	53
4			trans-4,4-Dimethyl-2-hexene	112	C8H16	019550-83-5	50
5			Cyclohexane, bromo-	162	C6H11Br	000108-85-0	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

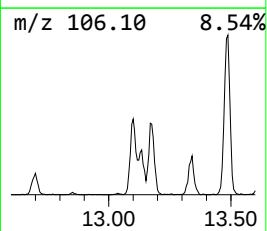
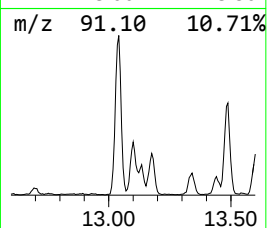
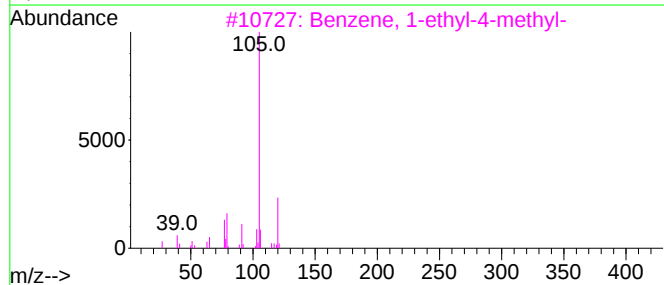
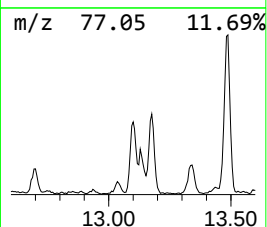
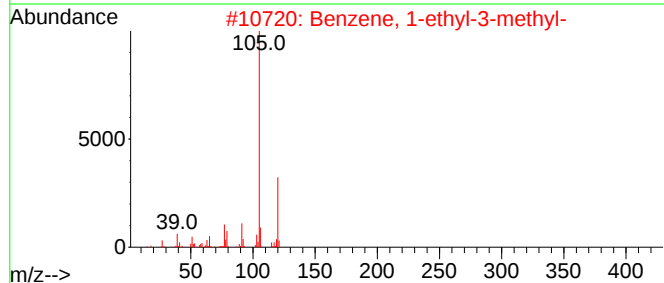
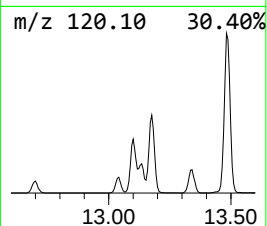
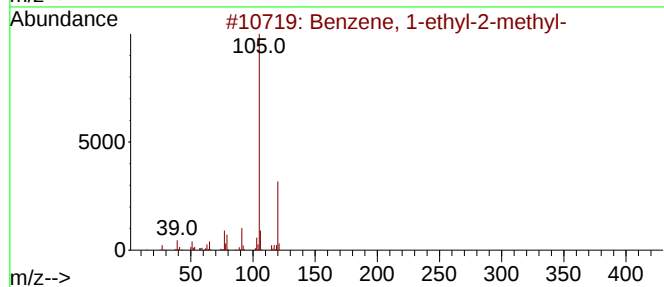
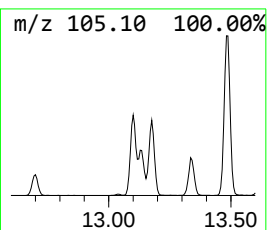
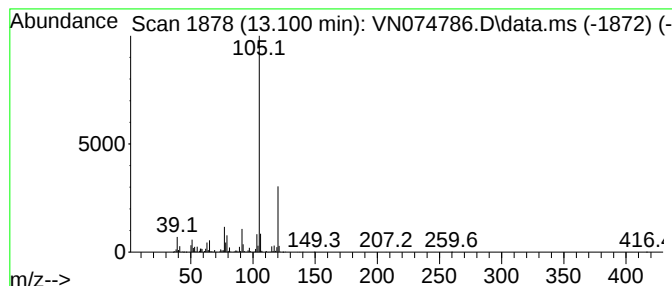
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	74.18 ug/l	4238190	1,4-Dichlorobenzene-d4	13.794

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	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

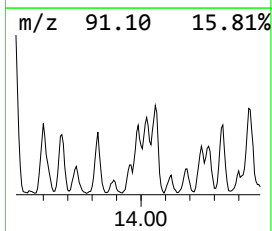
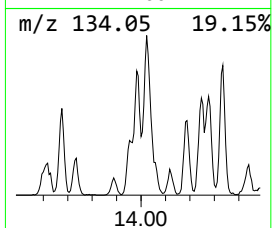
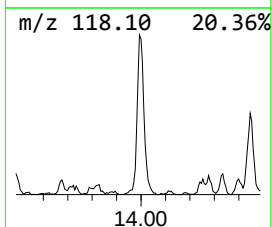
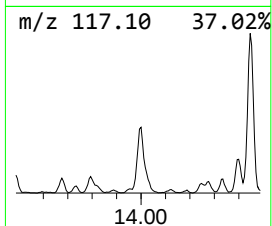
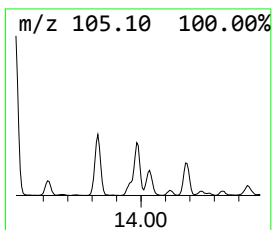
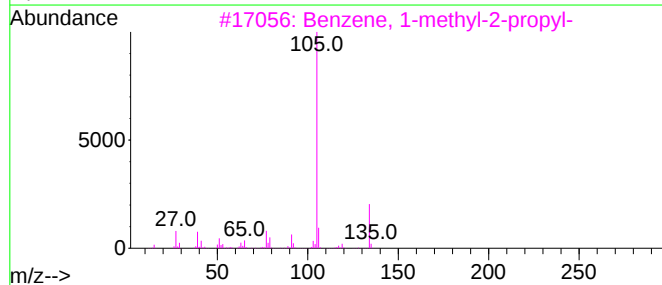
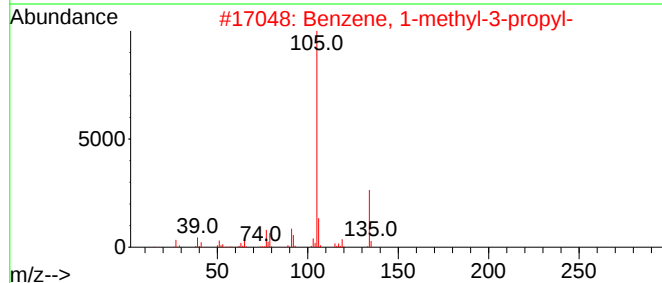
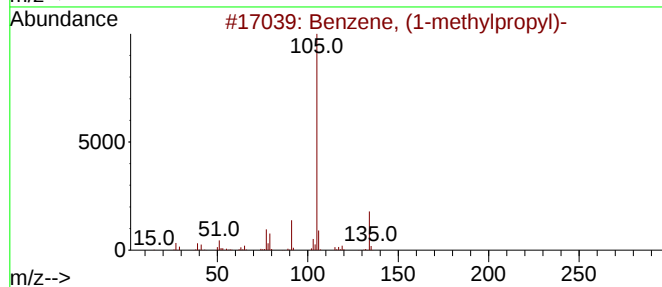
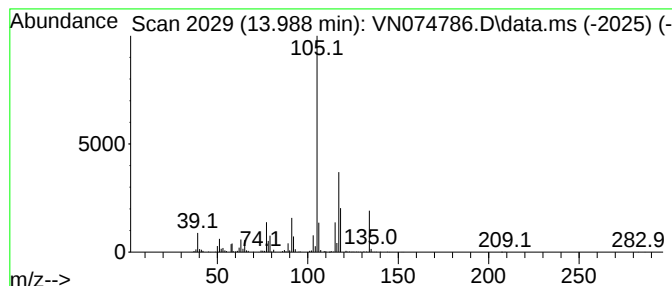
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methylpropyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.988	54.74 ug/l	3127340	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	55
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	43
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	43
5			2-Indanol	134	C9H10O	004254-29-9	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

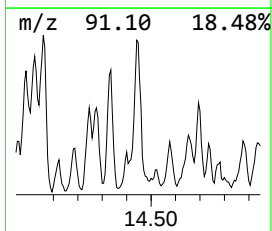
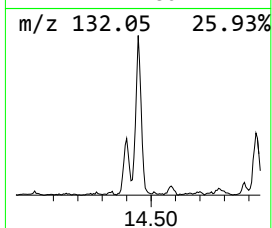
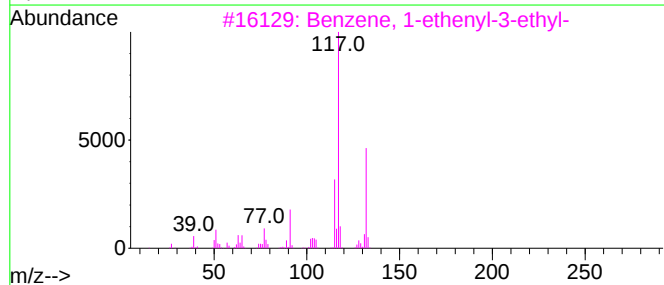
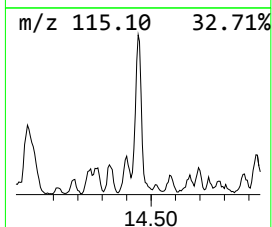
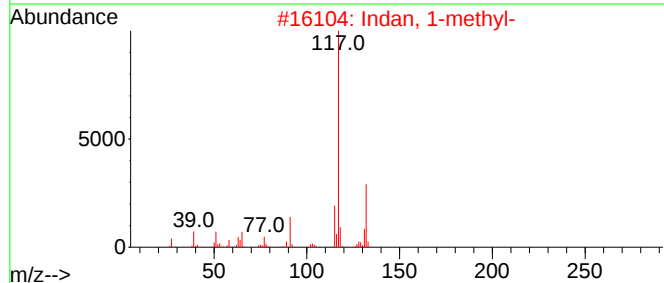
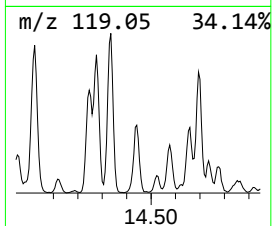
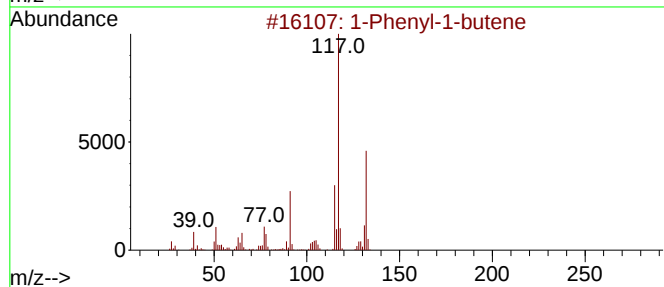
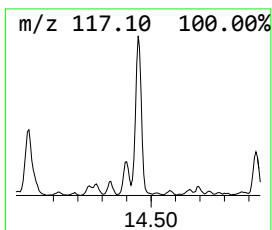
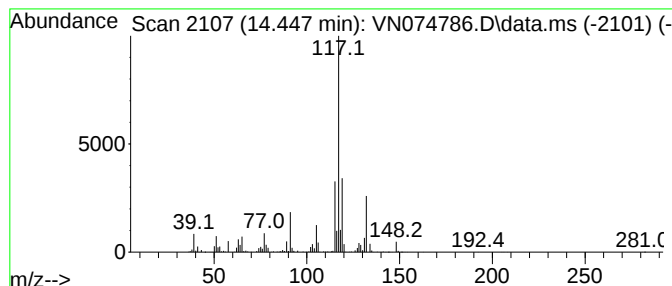
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Quant Title : SW846 8260

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.447	53.69 ug/l	3067200	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	81
2			Indan, 1-methyl-	132	C10H12	000767-58-8	76
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	72
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	72
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

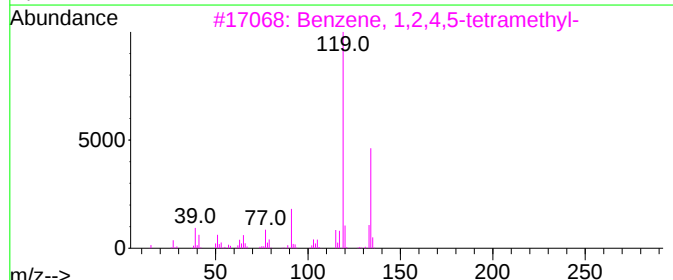
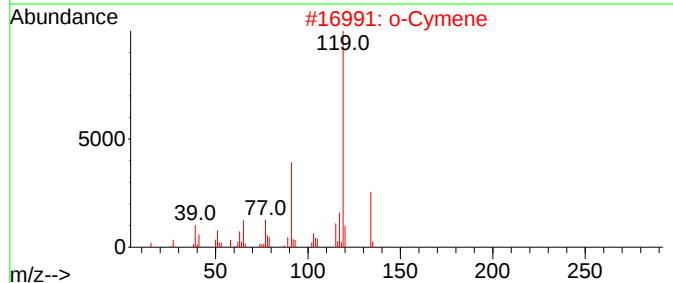
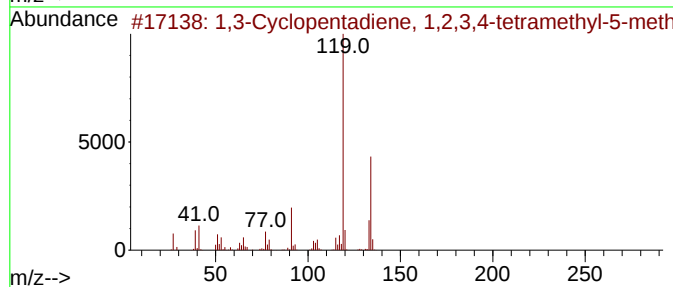
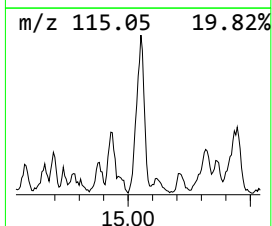
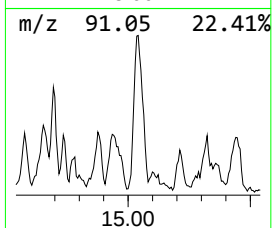
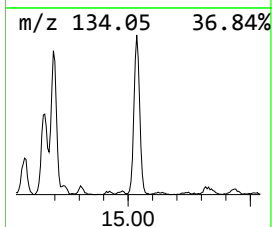
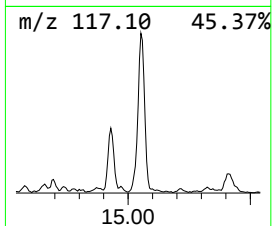
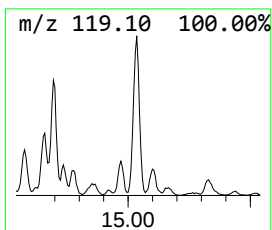
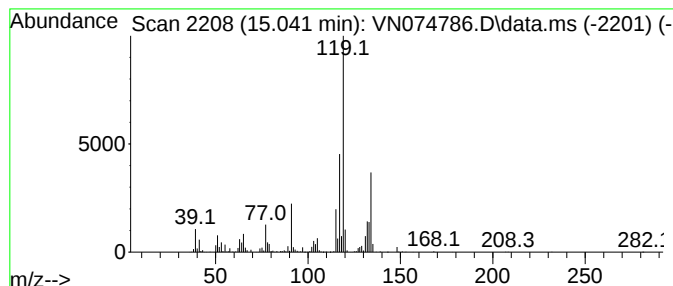
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100522W.M
Quant Title : SW846 8260

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

Peak Number 10 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.041	93.95 ug/l	5367640	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	64
2			o-Cymene	134	C10H14	000527-84-4	64
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4			p-Cymene	134	C10H14	000099-87-6	60
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100622\
Data File : VN074786.D
Acq On : 06 Oct 2022 16:16
Operator : JC\MD
Sample : N4910-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100522W.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
Cyclopentane, 1...	8.730	69.7	ug/l	2939240	2	9.106	2108520	50.0
Cyclopentane, 1...	8.800	54.1	ug/l	2280130	2	9.106	2108520	50.0
Isopropylcyclob...	8.865	87.6	ug/l	3694960	2	9.106	2108520	50.0
Heptane	8.977	72.9	ug/l	3073980	2	9.106	2108520	50.0
unknown10.747	10.747	91.4	ug/l	4337800	3	11.871	2373310	50.0
Cyclohexane, et...	11.424	76.8	ug/l	3645790	3	11.871	2373310	50.0
Benzene, 1-ethy...	13.100	74.2	ug/l	4238190	4	13.794	2856550	50.0
Benzene, (1-met...	13.988	54.7	ug/l	3127340	4	13.794	2856550	50.0
1-Phenyl-1-butene	14.447	53.7	ug/l	3067200	4	13.794	2856550	50.0
1,3-Cyclopentad...	15.041	94.0	ug/l	5367640	4	13.794	2856550	50.0