

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
 Data File : VN074250.D
 Acq On : 01 Sep 2022 18:22
 Operator : JC\MD
 Sample : N4329-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 LB-3R

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\82N083022W.M
 Title : SW846 8260

Signal : TIC: VN074250.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.398	73	87	92	rBV	84394	187034	3.74%	0.193%
2	3.087	195	204	221	rVB	792750	1788207	35.80%	1.848%
3	3.451	255	266	280	rBV2	227039	536689	10.74%	0.555%
4	4.187	378	391	409	rBV2	110418	407482	8.16%	0.421%
5	4.998	514	529	532	rBV2	538323	1712038	34.28%	1.769%
6	5.063	533	540	573	rVB2	1164594	4627306	92.64%	4.781%
7	5.534	603	620	635	rBV	974306	3389432	67.86%	3.502%
8	6.381	751	764	776	rBV2	78274	228961	4.58%	0.237%
9	6.522	776	788	797	rBV4	57085	182489	3.65%	0.189%
10	6.810	828	837	848	rBV3	72381	232072	4.65%	0.240%
11	6.969	853	864	871	rBV2	231645	692607	13.87%	0.716%
12	7.069	871	881	903	rVB	1492801	4459162	89.28%	4.607%
13	7.704	981	989	993	rBV5	34669	95394	1.91%	0.099%
14	7.775	995	1001	1010	rVV3	88379	245170	4.91%	0.253%
15	7.951	1021	1031	1034	rVV	255098	584397	11.70%	0.604%
16	8.022	1034	1043	1050	rVV3	1304518	3989138	79.87%	4.122%
17	8.104	1050	1057	1069	rVV2	929264	2506034	50.17%	2.589%
18	8.228	1069	1078	1084	rVV2	542848	1299534	26.02%	1.343%
19	8.286	1084	1088	1095	rVV3	146550	353945	7.09%	0.366%
20	8.380	1095	1104	1116	rVV4	407540	1244732	24.92%	1.286%
21	8.504	1116	1125	1128	rVV3	616969	1485006	29.73%	1.534%
22	8.557	1129	1134	1143	rVV4	772155	2498130	50.01%	2.581%
23	8.639	1143	1148	1158	rVV2	492452	1162825	23.28%	1.201%
24	8.727	1158	1163	1178	rVB7	50348	178922	3.58%	0.185%
25	8.904	1183	1193	1204	rBV2	759319	1809802	36.23%	1.870%
26	9.057	1214	1219	1226	rVV	51753	106090	2.12%	0.110%
27	9.139	1226	1233	1237	rVV2	69025	150417	3.01%	0.155%
28	9.175	1237	1239	1250	rVB4	38664	87461	1.75%	0.090%
29	9.392	1259	1276	1285	rBV2	1548574	4350834	87.11%	4.495%
30	9.533	1293	1300	1302	rBV3	50217	94643	1.89%	0.098%
31	9.574	1302	1307	1315	rVB	286020	558939	11.19%	0.578%
32	9.669	1315	1323	1329	rBV4	73418	157726	3.16%	0.163%
33	9.739	1329	1335	1337	rBV2	96491	181345	3.63%	0.187%
34	9.816	1343	1348	1358	rVB	403088	830059	16.62%	0.858%
35	9.910	1358	1364	1367	rBV	433540	817739	16.37%	0.845%
36	9.951	1367	1371	1378	rVB	677716	1317252	26.37%	1.361%

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37	10.145	1395	1404	1405	rBV3	142505	262846	5.26%	0.272%
38	10.257	1417	1423	1427	rBV	130218	238748	4.78%	0.247%
39	10.304	1427	1431	1437	rVB3	82524	150895	3.02%	0.156%
40	10.380	1437	1444	1450	rBV	1145794	2211734	44.28%	2.285%
41	10.439	1451	1454	1459	rVV	176664	338455	6.78%	0.350%
42	10.492	1459	1463	1468	rVV4	116887	237401	4.75%	0.245%
43	10.574	1468	1477	1487	rVB5	193682	549305	11.00%	0.568%
44	10.721	1497	1502	1508	rBV2	95712	191921	3.84%	0.198%
45	10.798	1508	1515	1525	rVV4	360835	873334	17.48%	0.902%
46	10.892	1525	1531	1538	rVB9	33367	92375	1.85%	0.095%
47	11.033	1545	1555	1557	rBV4	55183	115470	2.31%	0.119%
48	11.192	1578	1582	1586	rVV3	141390	250129	5.01%	0.258%
49	11.227	1586	1588	1596	rVB4	64162	121576	2.43%	0.126%
50	11.492	1625	1633	1637	rBV7	41182	91277	1.83%	0.094%
51	11.557	1639	1644	1650	rBV3	72482	151917	3.04%	0.157%
52	11.686	1658	1666	1674	rBV	1134686	2056127	41.17%	2.124%
53	11.786	1674	1683	1693	rVV2	457433	964895	19.32%	0.997%
54	11.892	1693	1701	1712	rVV	269839	670058	13.42%	0.692%
55	11.986	1712	1717	1730	rVB3	67445	145783	2.92%	0.151%
56	12.215	1746	1756	1765	rBV	100876	251262	5.03%	0.260%
57	12.515	1801	1807	1815	rVB	1105570	1830281	36.64%	1.891%
58	12.668	1825	1833	1842	rBV2	834736	1424256	28.51%	1.472%
59	12.780	1843	1852	1857	rVV7	66474	158339	3.17%	0.164%
60	12.857	1857	1865	1873	rVV	3410429	4994786	100.00%	5.161%
61	12.957	1873	1882	1886	rVV2	112052	241249	4.83%	0.249%
62	12.998	1886	1889	1897	rVB3	51479	92935	1.86%	0.096%
63	13.157	1909	1916	1924	rBV	647334	1083650	21.70%	1.120%
64	13.304	1937	1941	1947	rVB	74322	120135	2.41%	0.124%
65	13.439	1954	1964	1969	rBV2	278287	684515	13.70%	0.707%
66	13.498	1969	1974	1979	rVB2	70784	102068	2.04%	0.105%
67	13.610	1987	1993	2005	rVB2	756608	1476383	29.56%	1.525%
68	13.774	2015	2021	2024	rBV	763718	1187618	23.78%	1.227%
69	13.815	2024	2028	2032	rVB	1662183	2389958	47.85%	2.469%
70	13.862	2032	2036	2037	rBV	202471	242971	4.86%	0.251%
71	13.939	2045	2049	2055	rVB	179356	279918	5.60%	0.289%
72	14.009	2055	2061	2066	rBV	97127	152761	3.06%	0.158%
73	14.068	2066	2071	2079	rVB2	92784	140287	2.81%	0.145%
74	14.157	2079	2086	2091	rBV	2032832	3156149	63.19%	3.261%
75	14.215	2091	2096	2100	rVV	396015	670412	13.42%	0.693%
76	14.262	2100	2104	2112	rVV2	1459395	2414766	48.35%	2.495%

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77	14.398	2122	2127	2132	rBV3	65671	103617	2.07%	0.107%
78	14.480	2132	2141	2150	rBV	1259363	2137683	42.80%	2.209%
79	14.557	2150	2154	2158	rVV	115171	158546	3.17%	0.164%
80	14.698	2174	2178	2181	rBV3	96661	143717	2.88%	0.148%
81	14.745	2181	2186	2192	rVV	1048364	1773461	35.51%	1.832%
82	14.792	2192	2194	2198	rVB2	107121	133772	2.68%	0.138%
83	14.868	2198	2207	2213	rBV3	1955199	4052534	81.14%	4.187%
84	14.921	2213	2216	2223	rVB3	171843	302410	6.05%	0.312%
85	15.021	2227	2233	2238	rBV2	189397	326880	6.54%	0.338%
86	15.127	2240	2251	2255	rVV2	614117	1499541	30.02%	1.549%
87	15.168	2255	2258	2263	rVV	295512	516174	10.33%	0.533%
88	15.245	2263	2271	2282	rVB2	575687	1329854	26.62%	1.374%
89	15.433	2297	2303	2309	rVV	624439	1139998	22.82%	1.178%
90	15.492	2309	2313	2317	rVV2	93103	168832	3.38%	0.174%
91	15.545	2317	2322	2326	rVV3	180119	395849	7.93%	0.409%
92	15.598	2328	2331	2343	rVB2	170807	319846	6.40%	0.330%
93	15.727	2346	2353	2361	rBV	281345	555538	11.12%	0.574%
94	15.903	2373	2383	2387	rBV3	188174	487482	9.76%	0.504%
95	16.033	2399	2405	2410	rBV2	80669	168118	3.37%	0.174%
96	16.109	2410	2418	2435	rVB3	171982	469858	9.41%	0.485%
97	16.262	2435	2444	2454	rBV4	129839	319117	6.39%	0.330%
98	16.462	2466	2478	2482	rBV5	76226	199437	3.99%	0.206%
99	16.533	2482	2490	2503	rVB	398603	946761	18.95%	0.978%
100	16.733	2515	2524	2533	rBV	764921	1775432	35.55%	1.834%

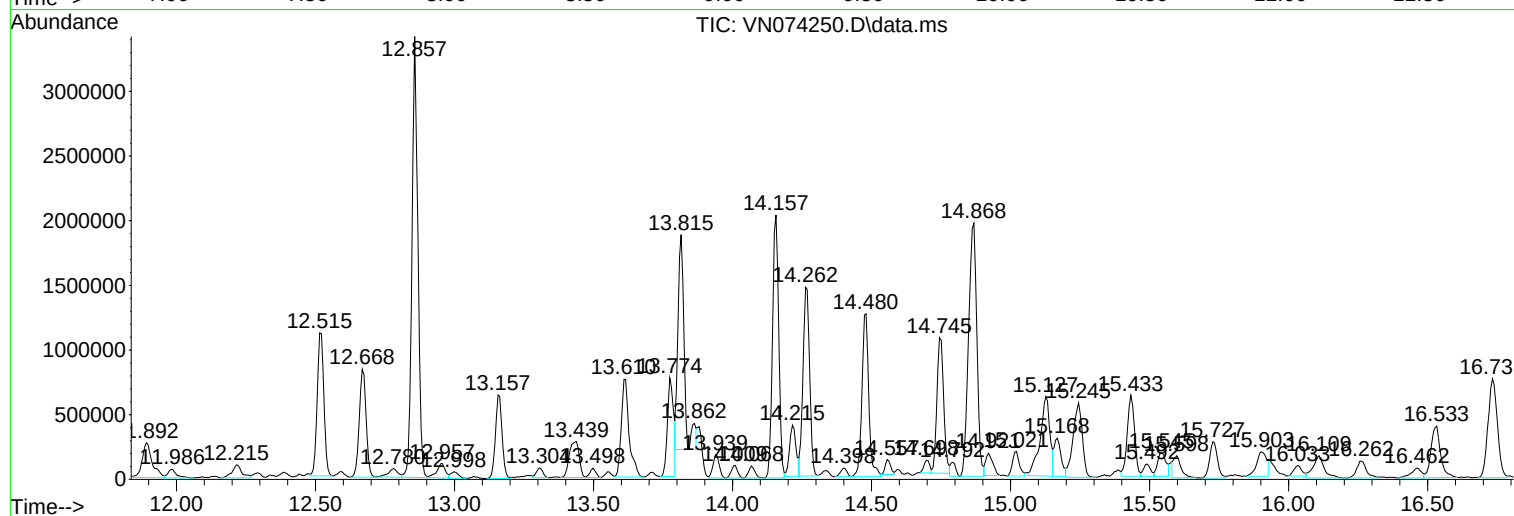
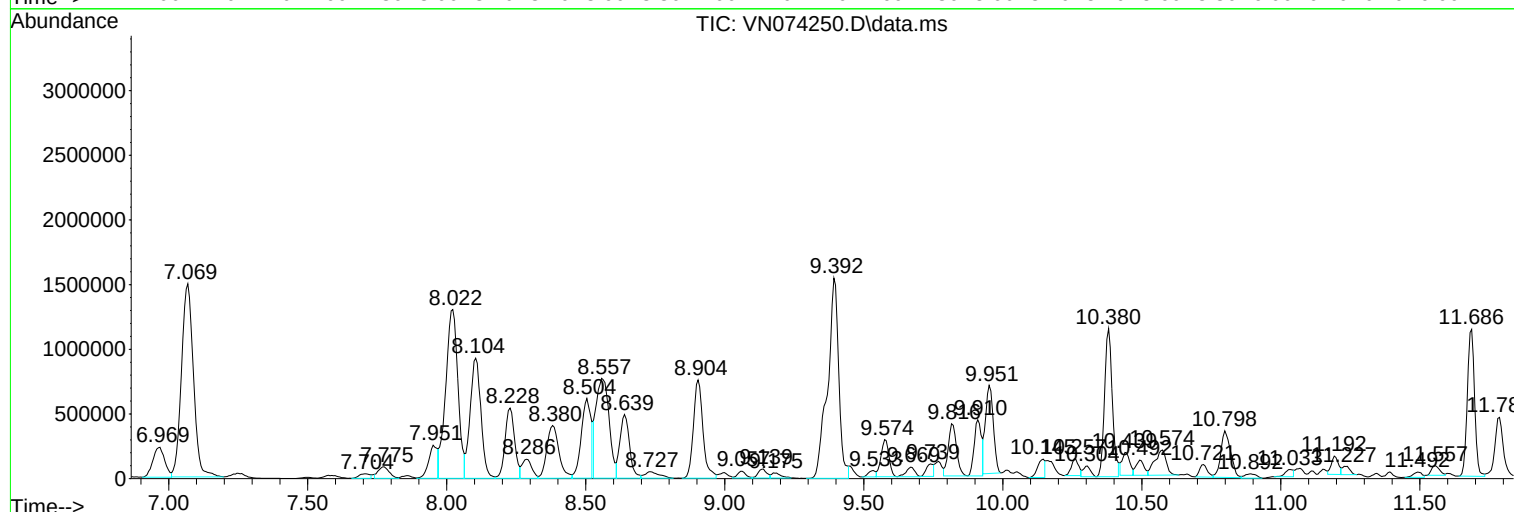
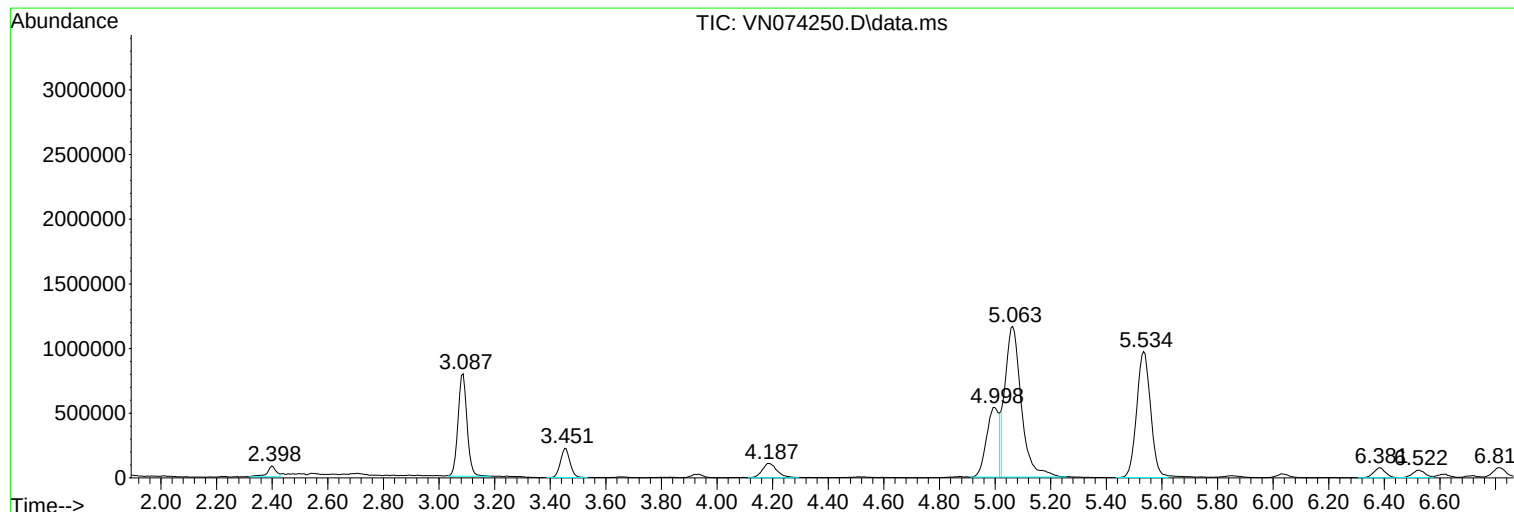
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
Quant Title : SW846 8260

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



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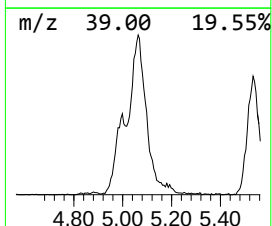
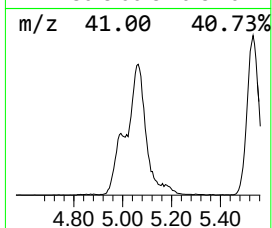
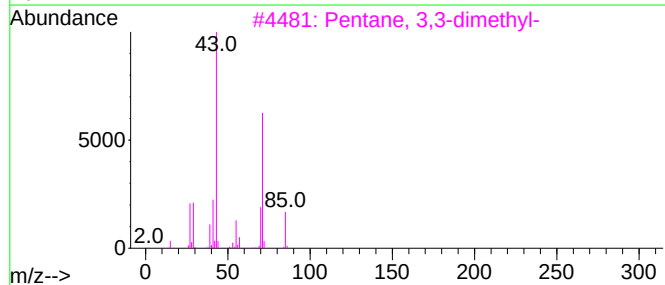
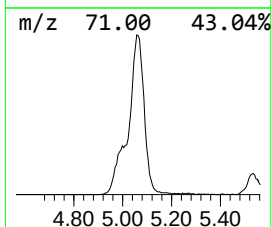
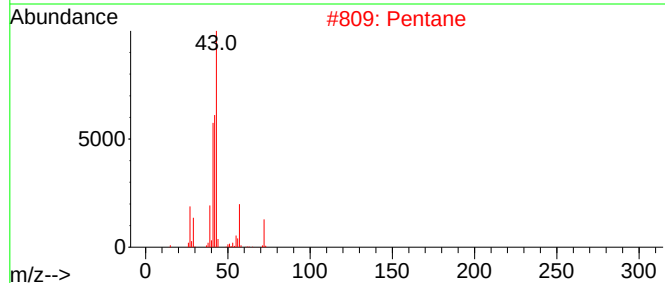
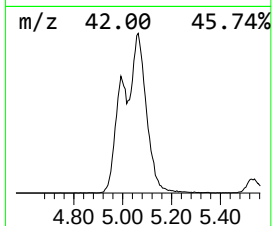
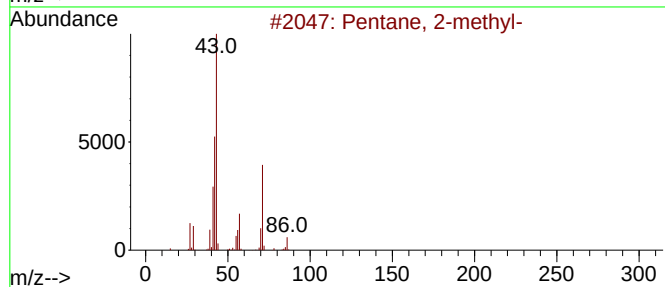
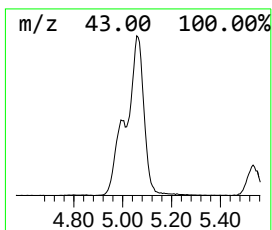
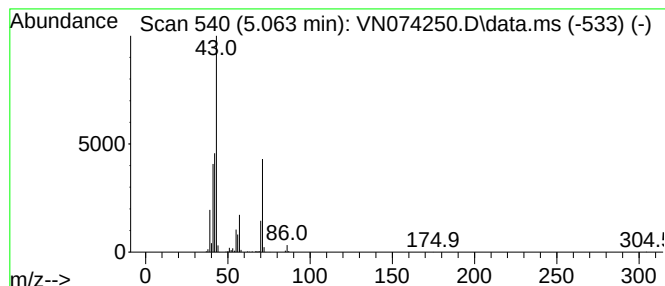
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.063	58.00 ug/l	4627310	Pentafluorobenzene	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2			Pentane	72	C5H12	000109-66-0	47
3			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	28
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	28
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28



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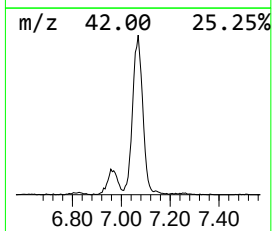
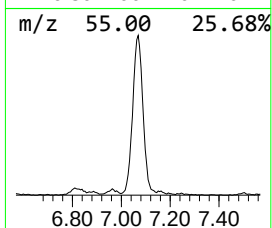
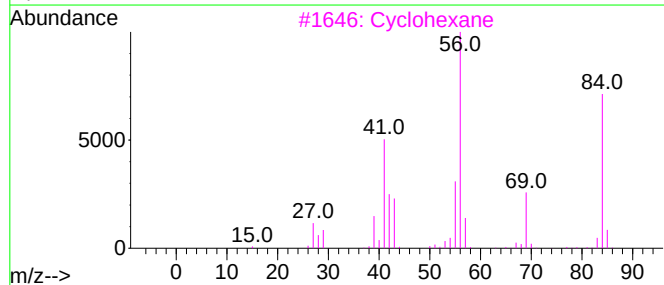
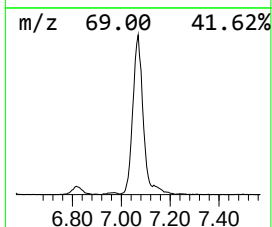
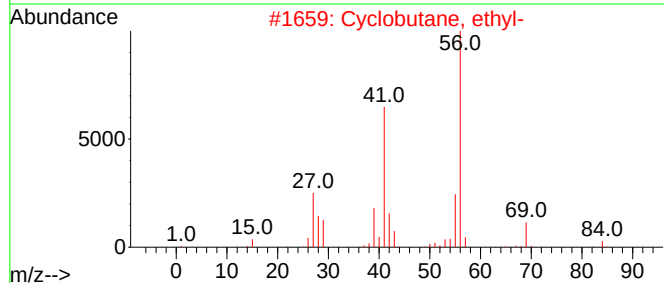
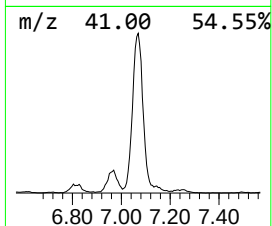
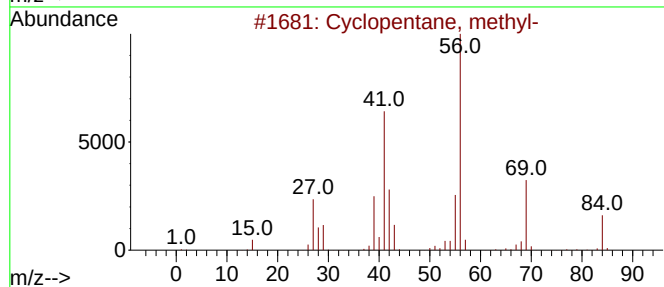
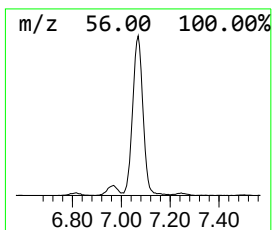
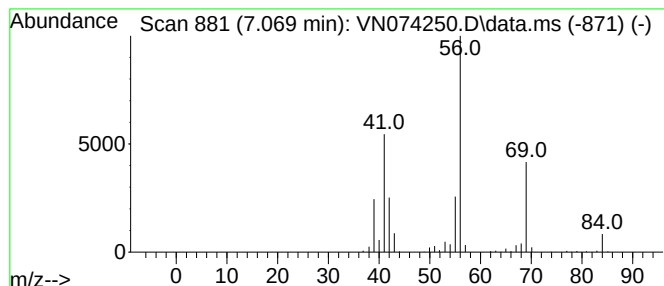
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Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.069	55.89 ug/l	4459160	Pentafluorobenzene	8.022

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3		Cyclohexane	84	C6H12	000110-82-7	64
4		Cyclopropane, propyl-	84	C6H12	002415-72-7	56
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50



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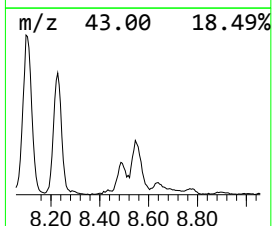
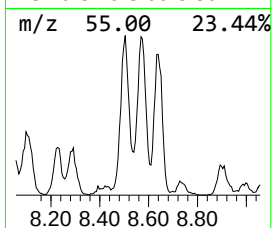
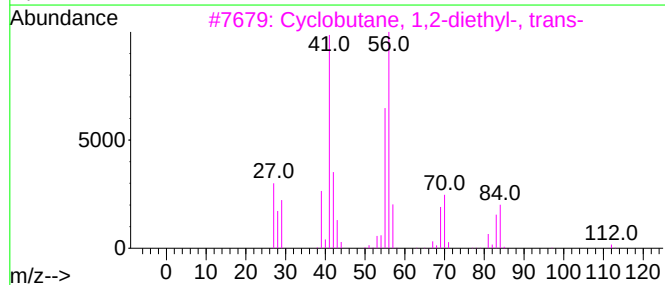
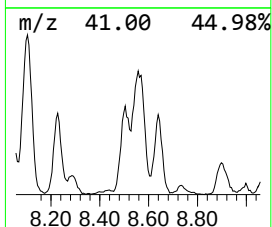
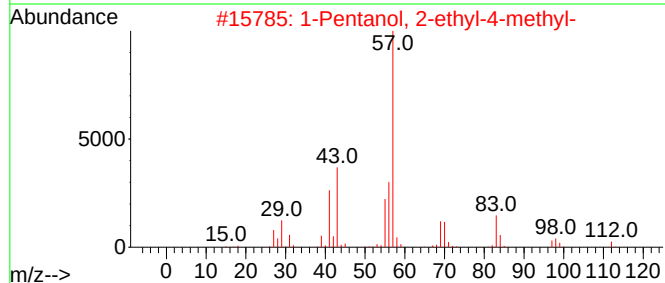
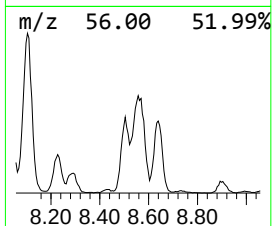
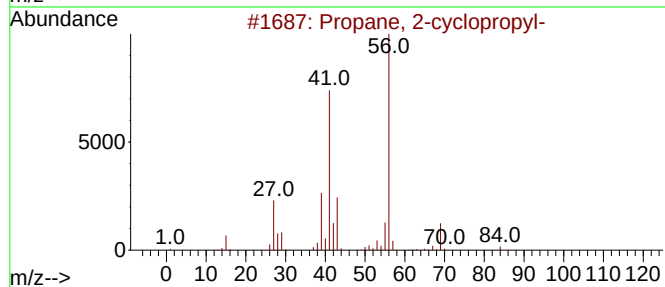
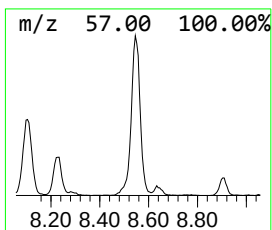
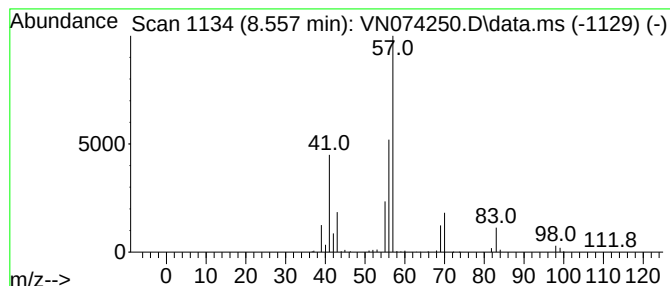
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Quant Title : SW846 8260

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

Peak Number 3 unknown8.557 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.557	69.02 ug/l	2498130	1,4-Difluorobenzene	8.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	43
2			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	38
3			Cyclobutane, 1,2-diethyl-, trans-	112	C8H16	019341-98-1	38
4			5-Methyl-2-hexanol, chlorodifluo...	228	C9H15ClF2O2	1010376-27-1	38
5			Cyclobutane, ethyl-	84	C6H12	004806-61-5	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

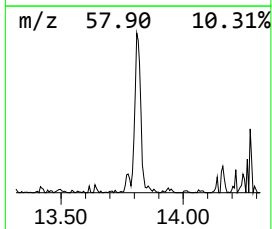
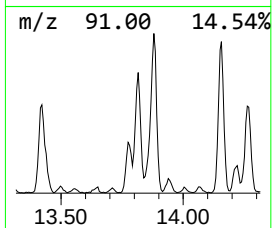
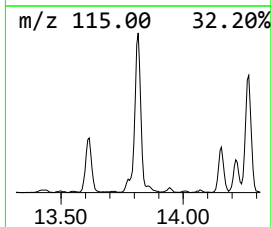
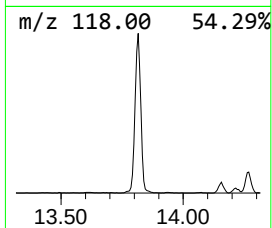
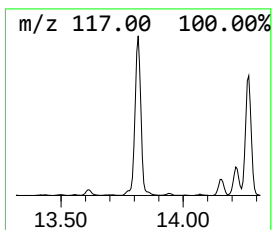
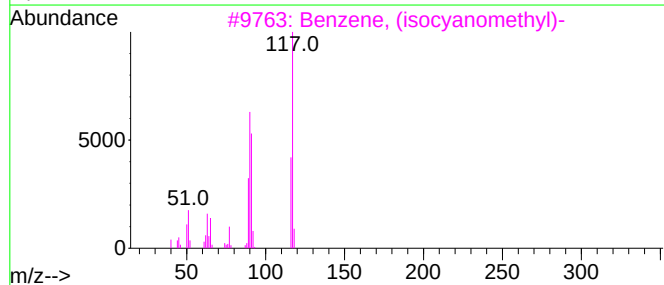
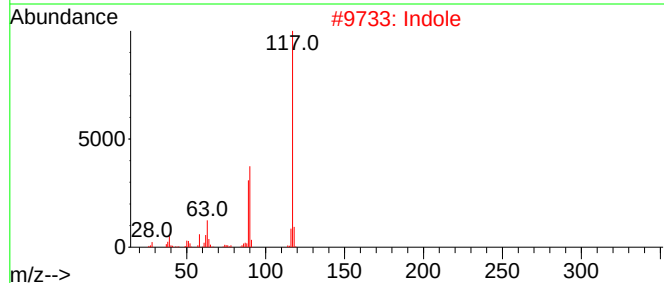
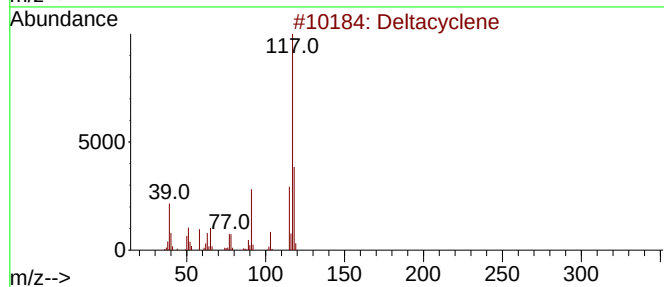
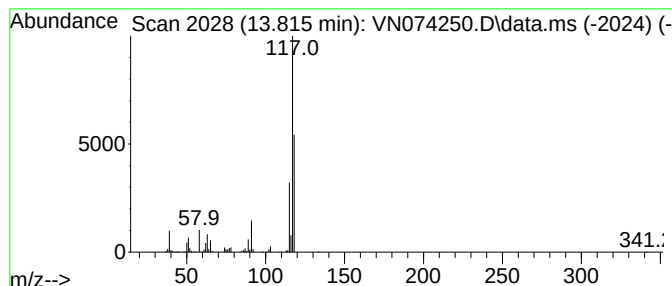
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Quant Title : SW846 8260

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

Peak Number 4 Indole Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.815	80.94 ug/l	2389960	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Deltacyclene	118	C9H10	007785-10-6	59
2			Indole	117	C8H7N	000120-72-9	50
3			Benzene, (isocyanomethyl)-	117	C8H7N	010340-91-7	47
4			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	43
5			4-Ethynylaniline	117	C8H7N	014235-81-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

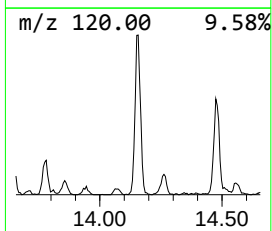
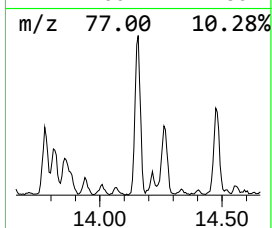
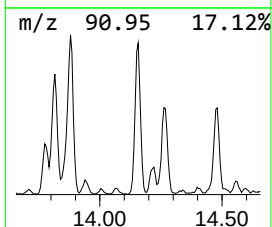
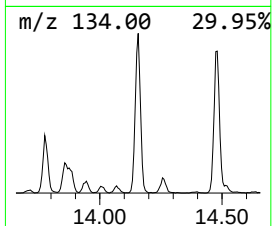
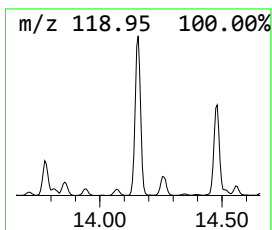
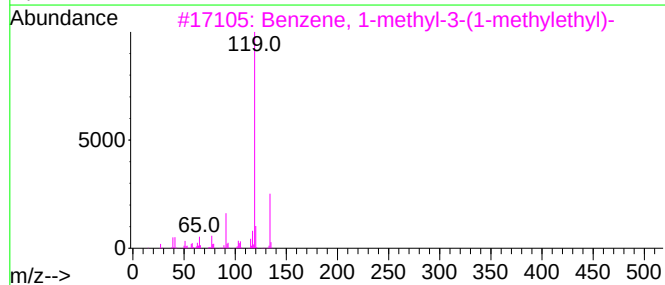
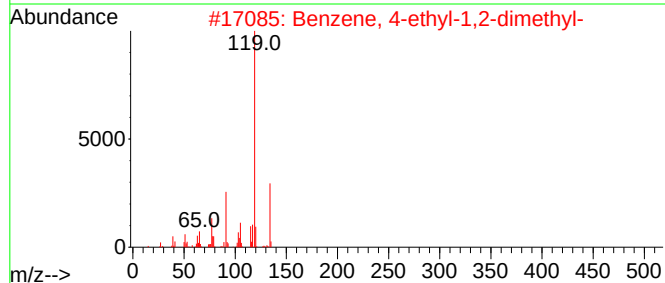
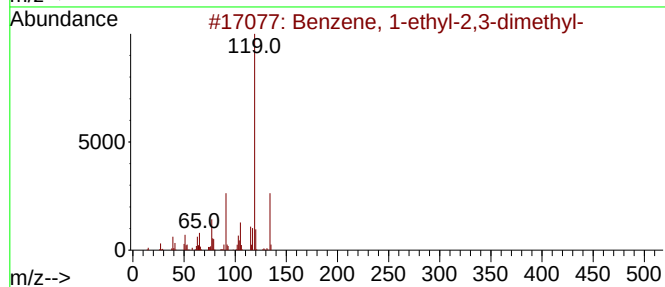
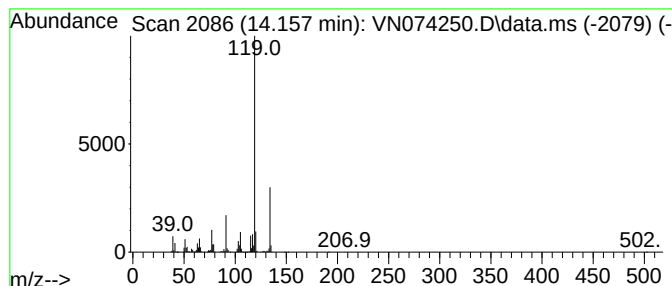
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.156	106.89 ug/l	3156150	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

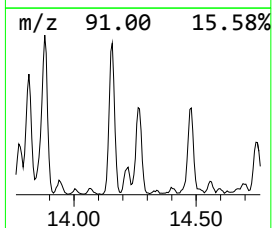
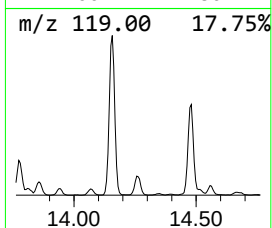
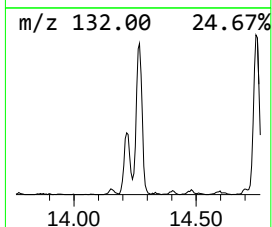
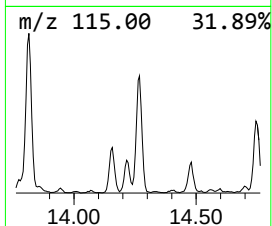
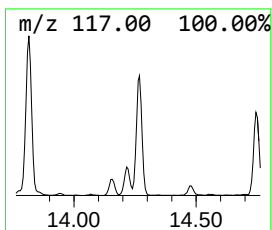
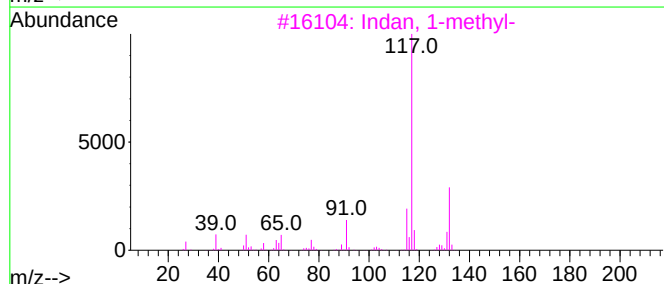
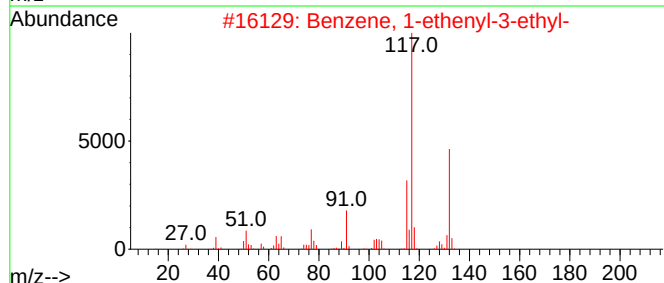
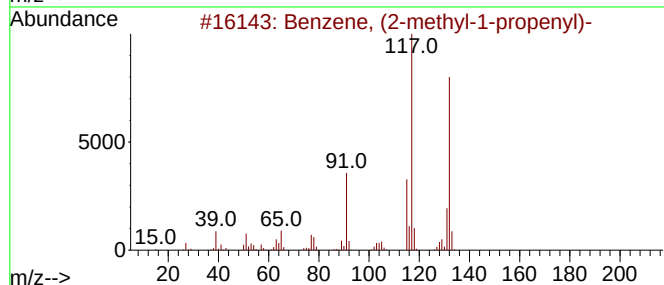
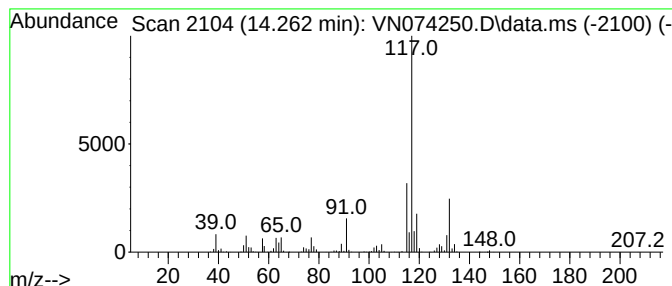
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-propen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.262	81.78 ug/l	2414770	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
3			Indan, 1-methyl-	132	C10H12	000767-58-8	81
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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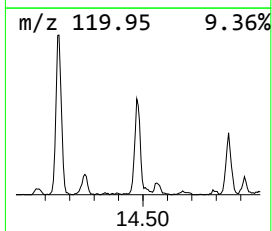
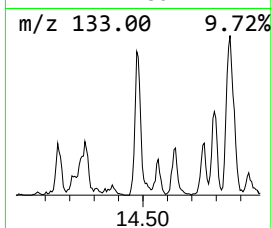
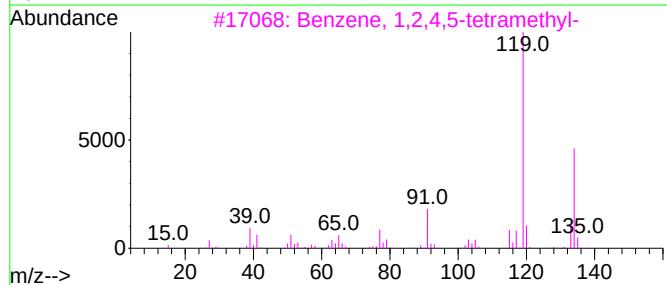
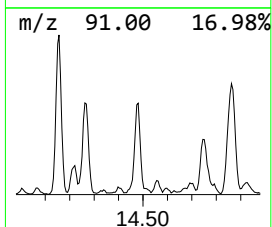
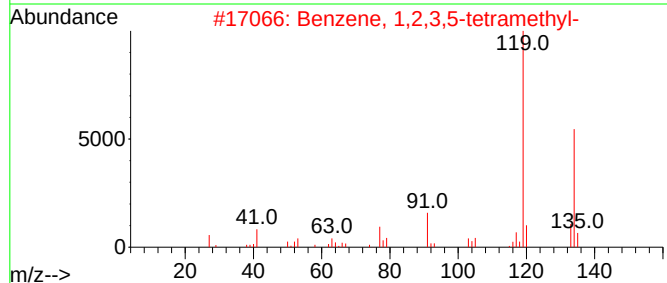
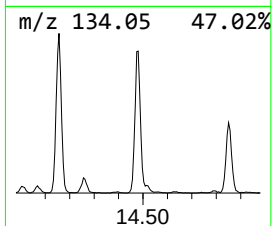
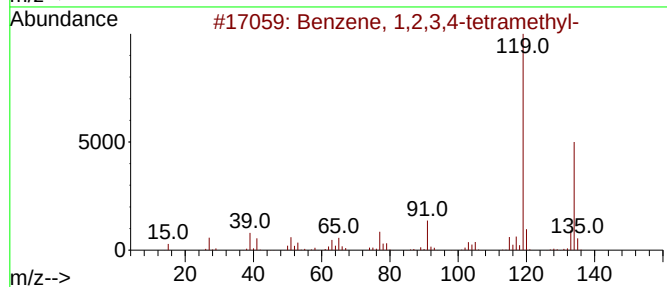
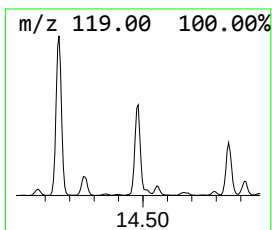
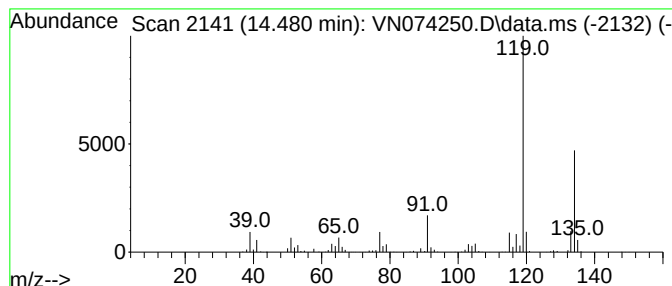
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.480	72.40 ug/l	2137680	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

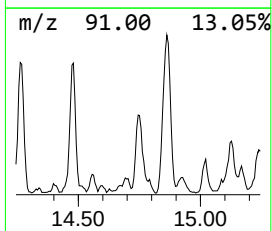
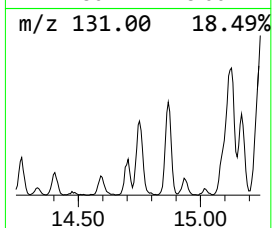
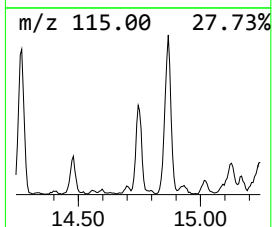
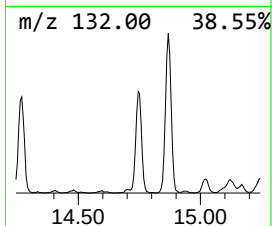
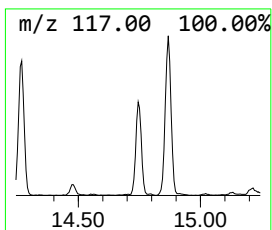
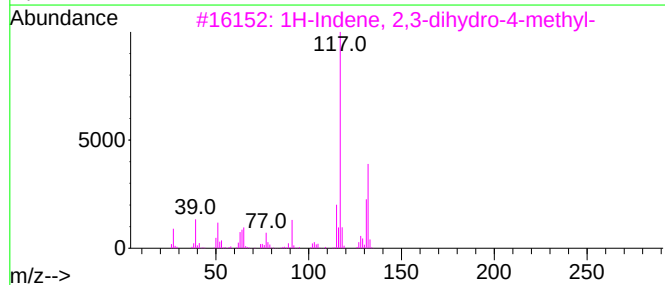
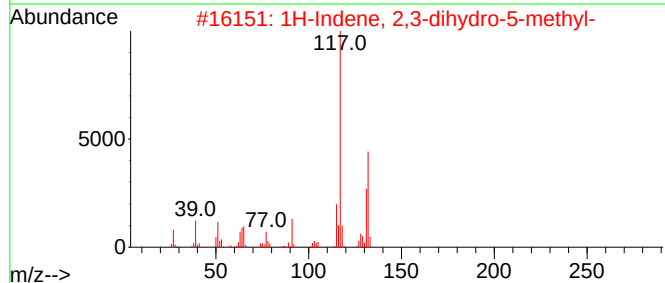
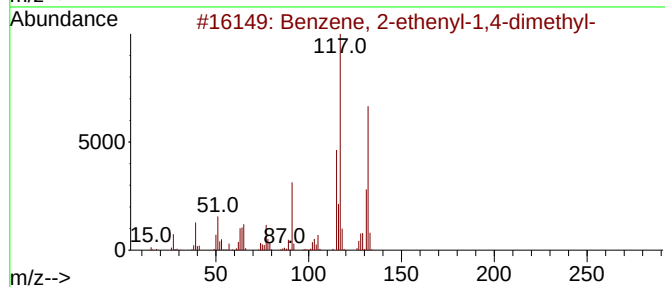
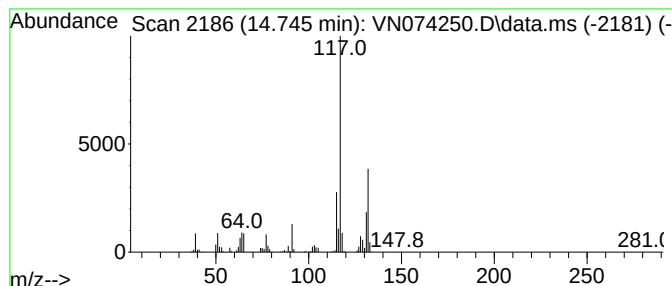
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.745	60.06 ug/l	1773460	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
4			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	91
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
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Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

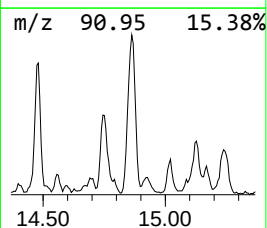
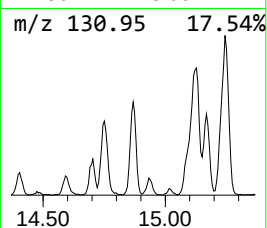
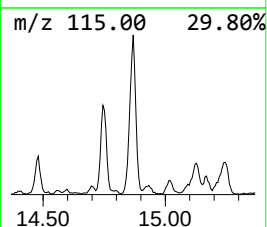
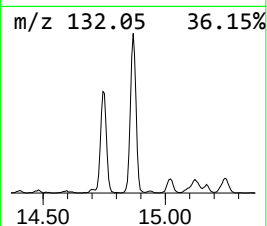
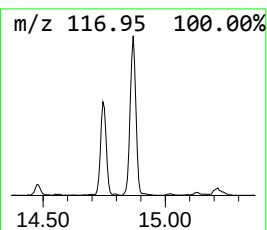
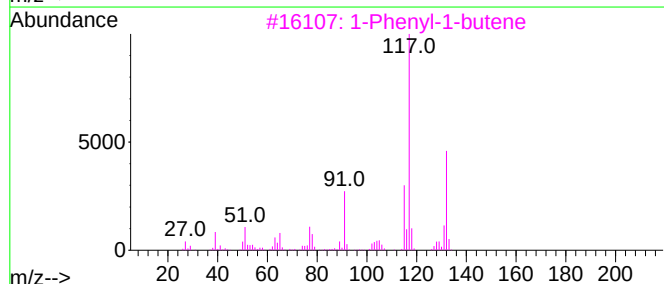
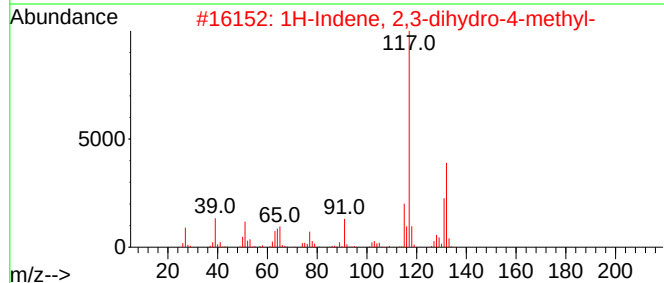
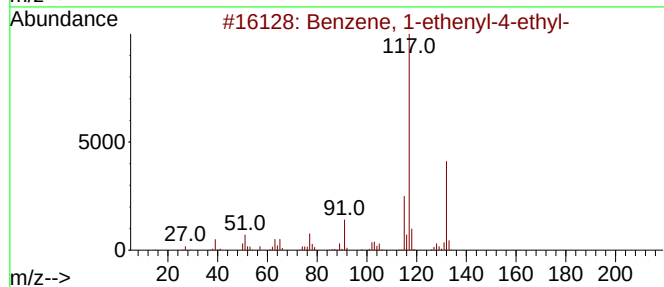
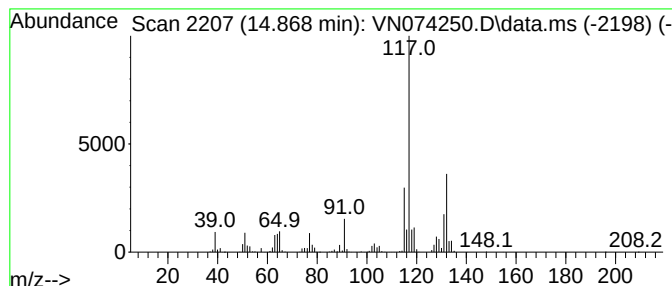
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.868	137.25 ug/l	4052530	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	93
4			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

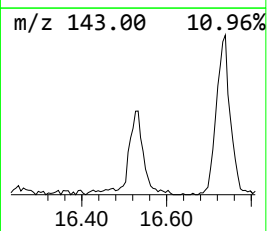
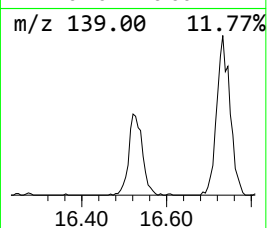
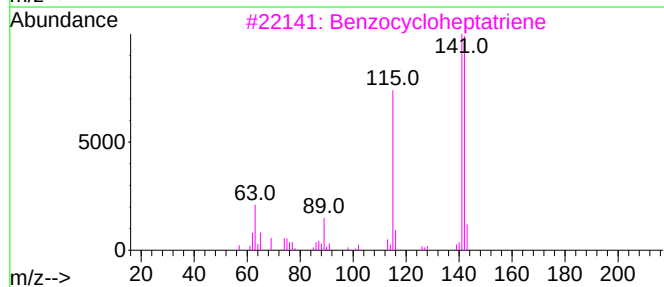
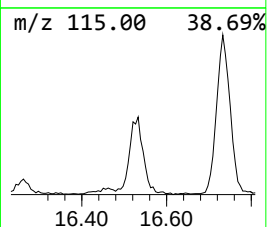
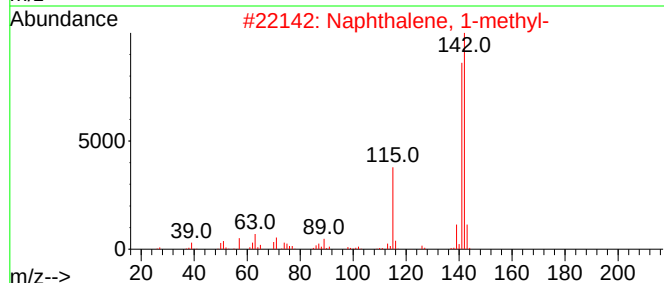
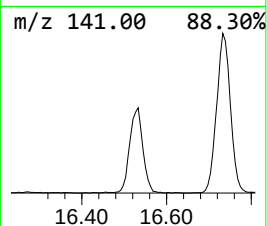
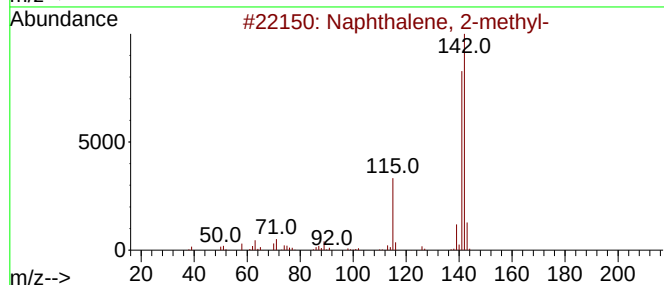
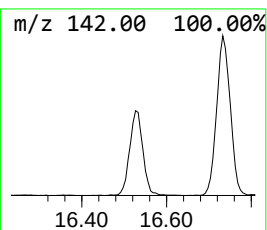
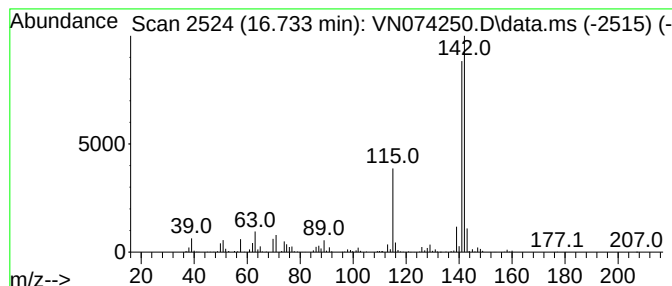
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.733	60.13 ug/l	1775430	1,4-Dichlorobenzene-d4	13.615

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	94
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
Data File : VN074250.D
Acq On : 01 Sep 2022 18:22
Operator : JC\MD
Sample : N4329-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.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
Pentane, 2-methyl-	5.063	58.0	ug/l	4627310	1	8.022	3989140	50.0
Cyclopentane, m...	7.069	55.9	ug/l	4459160	1	8.022	3989140	50.0
unknown8.557	8.557	69.0	ug/l	2498130	2	8.904	1809800	50.0
Indole	13.815	80.9	ug/l	2389960	4	13.615	1476380	50.0
Benzene, 1-ethy...	14.156	106.9	ug/l	3156150	4	13.615	1476380	50.0
Benzene, (2-met...	14.262	81.8	ug/l	2414770	4	13.615	1476380	50.0
Benzene, 1,2,3,...	14.480	72.4	ug/l	2137680	4	13.615	1476380	50.0
Benzene, 2-ethe...	14.745	60.1	ug/l	1773460	4	13.615	1476380	50.0
Benzene, 1-ethe...	14.868	137.3	ug/l	4052530	4	13.615	1476380	50.0
Naphthalene, 2-...	16.733	60.1	ug/l	1775430	4	13.615	1476380	50.0