

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
 Data File : VN078581.D
 Acq On : 19 Jul 2023 16:43
 Operator : JC\MD
 Sample : 03668-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 N48719

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\82N071023W.M
 Title : SW846 8260

Signal : TIC: VN078581.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.295	54	58	65	rBV2	35227	57879	1.46%	0.102%
2	5.283	555	566	568	rBV4	24226	59768	1.50%	0.106%
3	5.347	568	577	592	rVB4	90608	350263	8.81%	0.620%
4	5.824	643	658	672	rVB2	86348	290176	7.30%	0.513%
5	6.324	729	743	754	rBV	176946	530073	13.34%	0.938%
6	6.665	794	801	813	rVB4	22075	64867	1.63%	0.115%
7	7.336	904	915	926	rVV2	437430	1226730	30.86%	2.170%
8	8.018	1023	1031	1039	rVB4	30087	73913	1.86%	0.131%
9	8.171	1047	1057	1062	rBV	362701	889327	22.37%	1.573%
10	8.236	1062	1068	1081	rVV3	770587	3038276	76.44%	5.375%
11	8.330	1081	1084	1094	rVB5	93682	224854	5.66%	0.398%
12	8.453	1097	1105	1111	rBV2	166380	378397	9.52%	0.669%
13	8.518	1111	1116	1121	rVV2	84714	205314	5.17%	0.363%
14	8.588	1121	1128	1141	rVV	330784	777398	19.56%	1.375%
15	8.730	1143	1152	1158	rVV6	245516	640563	16.12%	1.133%
16	8.794	1158	1163	1168	rVV2	224928	521100	13.11%	0.922%
17	8.859	1168	1174	1184	rVB	364657	846228	21.29%	1.497%
18	8.971	1185	1193	1205	rVB2	100031	223598	5.63%	0.396%
19	9.106	1207	1216	1233	rBV	913926	2061553	51.87%	3.647%
20	9.388	1255	1264	1274	rVB7	21762	53179	1.34%	0.094%
21	9.606	1285	1301	1318	rBV	1729835	3974717	100.00%	7.032%
22	9.782	1323	1331	1339	rBV2	203890	403327	10.15%	0.714%
23	9.871	1339	1346	1352	rVB2	67900	132037	3.32%	0.234%
24	9.947	1352	1359	1360	rBV3	57585	98880	2.49%	0.175%
25	10.012	1366	1370	1381	rVB5	105619	222943	5.61%	0.394%
26	10.153	1390	1394	1399	rVV6	26939	56308	1.42%	0.100%
27	10.206	1399	1403	1408	rVV4	37249	69513	1.75%	0.123%
28	10.347	1418	1427	1437	rBV4	78232	190105	4.78%	0.336%
29	10.459	1439	1446	1451	rBV2	126867	251486	6.33%	0.445%
30	10.571	1457	1465	1472	rBV2	1534069	2993783	75.32%	5.296%
31	10.747	1489	1495	1503	rVB5	137009	341840	8.60%	0.605%
32	10.918	1518	1524	1530	rBV2	201168	377197	9.49%	0.667%
33	11.006	1530	1539	1547	rVB5	126001	316914	7.97%	0.561%
34	11.094	1548	1554	1559	rVB9	23339	52775	1.33%	0.093%
35	11.259	1575	1582	1586	rVV4	58157	138994	3.50%	0.246%
36	11.306	1586	1590	1594	rVV4	52913	106592	2.68%	0.189%

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37	11.347	1594	1597	1601	rVV3	62293	104655	2.63%	0.185%
38	11.424	1601	1610	1615	rVV	250482	549878	13.83%	0.973%
39	11.476	1615	1619	1625	rVV3	122813	220480	5.55%	0.390%
40	11.582	1632	1637	1642	rVB5	34786	60644	1.53%	0.107%
41	11.682	1648	1654	1664	rBV7	51601	118153	2.97%	0.209%
42	11.794	1667	1673	1679	rBV3	34193	69902	1.76%	0.124%
43	11.871	1679	1686	1696	rVB	1207878	2170874	54.62%	3.840%
44	11.971	1696	1703	1713	rBV	384380	703098	17.69%	1.244%
45	12.076	1713	1721	1725	rVV	118808	249121	6.27%	0.441%
46	12.118	1726	1728	1740	rVB7	76705	196326	4.94%	0.347%
47	12.282	1749	1756	1765	rVB5	29450	66431	1.67%	0.118%
48	12.400	1765	1776	1782	rBV3	132496	317121	7.98%	0.561%
49	12.582	1802	1807	1812	rVV4	79032	149038	3.75%	0.264%
50	12.700	1819	1827	1834	rVB	369398	620475	15.61%	1.098%
51	12.853	1846	1853	1863	rVB	991494	1710015	43.02%	3.025%
52	12.935	1863	1867	1874	rVB5	32133	55841	1.40%	0.099%
53	13.041	1876	1885	1890	rBV	387301	646393	16.26%	1.144%
54	13.100	1890	1895	1898	rVV	213029	354038	8.91%	0.626%
55	13.135	1898	1901	1905	rVV	230270	378726	9.53%	0.670%
56	13.176	1905	1908	1915	rVB	272023	412398	10.38%	0.730%
57	13.341	1928	1936	1943	rBV	465751	768457	19.33%	1.359%
58	13.435	1944	1952	1956	rBV6	36620	97900	2.46%	0.173%
59	13.488	1956	1961	1966	rVB	158723	266852	6.71%	0.472%
60	13.618	1975	1983	1988	rBV2	131652	279019	7.02%	0.494%
61	13.676	1988	1993	1998	rVB2	112183	169419	4.26%	0.300%
62	13.735	1998	2003	2007	rBV	103932	162531	4.09%	0.288%
63	13.800	2007	2014	2024	rVB2	1173490	2779496	69.93%	4.917%
64	13.888	2024	2029	2034	rVB2	74955	130611	3.29%	0.231%
65	13.953	2034	2040	2043	rBV	309673	517395	13.02%	0.915%
66	14.000	2043	2048	2063	rVB3	778419	1991515	50.10%	3.523%
67	14.123	2063	2069	2074	rVB	170160	277352	6.98%	0.491%
68	14.188	2074	2080	2084	rBV	125480	204135	5.14%	0.361%
69	14.247	2085	2090	2093	rBV	134926	216245	5.44%	0.383%
70	14.276	2093	2095	2099	rVB2	85076	99794	2.51%	0.177%
71	14.335	2099	2105	2111	rVB	783487	1197840	30.14%	2.119%
72	14.400	2111	2116	2119	rBV	228835	398110	10.02%	0.704%
73	14.447	2119	2124	2132	rVB2	1039826	1831740	46.08%	3.240%
74	14.517	2132	2136	2141	rVB5	44891	76067	1.91%	0.135%
75	14.582	2141	2147	2152	rBV3	169555	286681	7.21%	0.507%
76	14.659	2152	2160	2163	rBV	493501	802420	20.19%	1.420%

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77	14.694	2163	2166	2171	rVB	274518	400958	10.09%	0.709%
78	14.741	2171	2174	2177	rVB2	84940	103652	2.61%	0.183%
79	14.776	2177	2180	2183	rBV	49186	68318	1.72%	0.121%
80	14.806	2183	2185	2190	rVB2	59102	77976	1.96%	0.138%
81	14.882	2190	2198	2202	rBV3	84377	170598	4.29%	0.302%
82	14.929	2202	2206	2212	rBV2	430259	744812	18.74%	1.318%
83	15.053	2218	2227	2234	rBV3	1536240	3877158	97.55%	6.859%
84	15.112	2234	2237	2246	rVB4	71434	154202	3.88%	0.273%
85	15.212	2248	2254	2262	rBV2	531189	1001974	25.21%	1.773%
86	15.323	2262	2273	2277	rBV3	310946	750279	18.88%	1.327%
87	15.364	2277	2280	2286	rVB3	139510	227413	5.72%	0.402%
88	15.441	2286	2293	2303	rVB3	402071	1007118	25.34%	1.782%
89	15.588	2313	2318	2322	rBV3	59816	105346	2.65%	0.186%
90	15.647	2322	2328	2333	rVB	314202	586556	14.76%	1.038%
91	15.706	2333	2338	2342	rBV2	80419	137535	3.46%	0.243%
92	15.753	2342	2346	2348	rBV2	96410	138672	3.49%	0.245%
93	15.947	2370	2379	2387	rBV	195509	419025	10.54%	0.741%
94	16.029	2389	2393	2399	rVV6	40165	93960	2.36%	0.166%
95	16.129	2401	2410	2413	rVV4	195358	481510	12.11%	0.852%
96	16.170	2413	2417	2427	rVV3	249934	571893	14.39%	1.012%
97	16.264	2427	2433	2439	rVV4	86581	222406	5.60%	0.393%
98	16.347	2439	2447	2461	rVB2	164953	502903	12.65%	0.890%
99	16.511	2463	2475	2484	rBV3	211785	547723	13.78%	0.969%
100	16.717	2498	2510	2514	rBV6	66236	187204	4.71%	0.331%

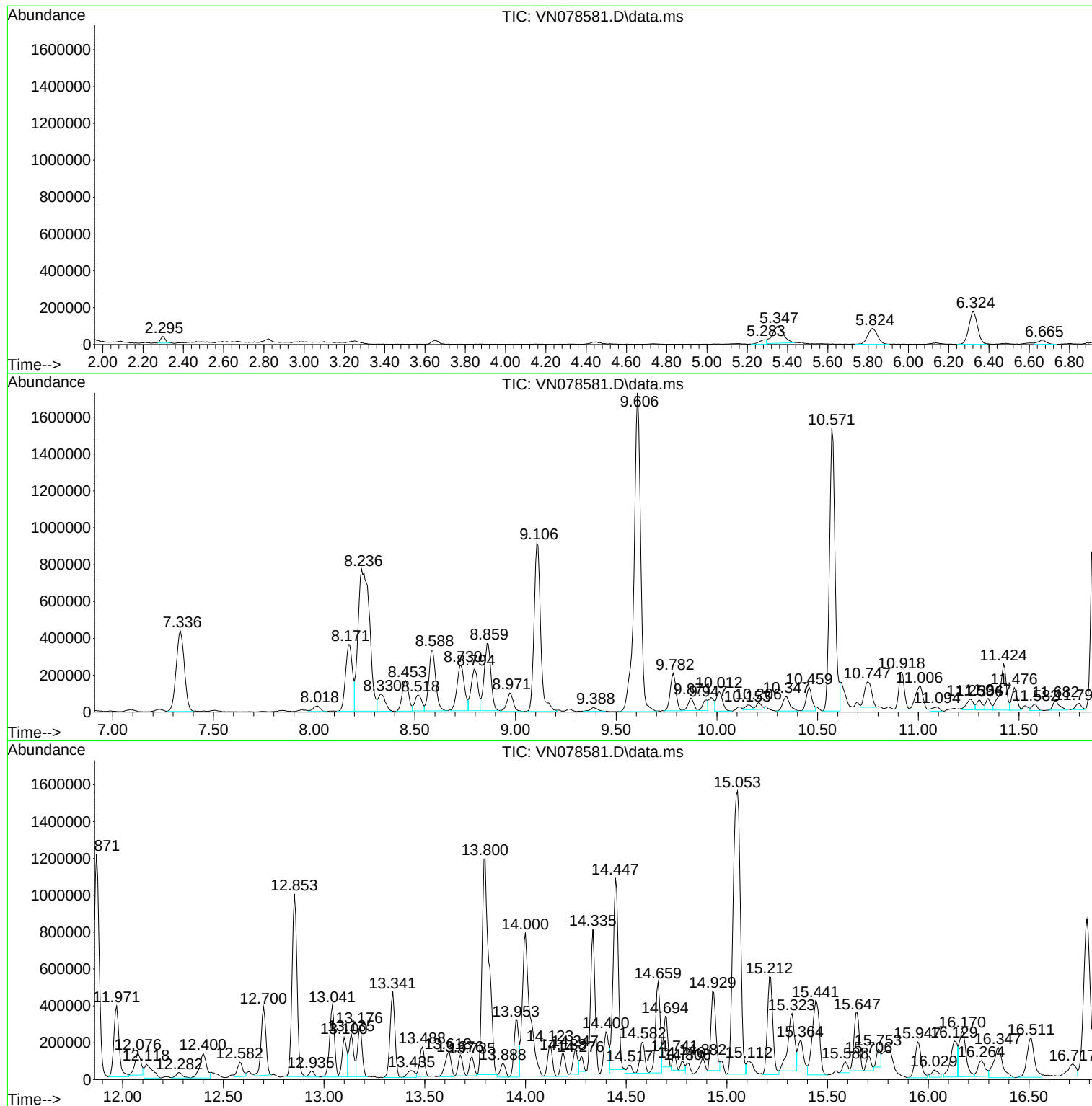
Sum of corrected areas: 56527244

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
Quant Title : SW846 8260

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



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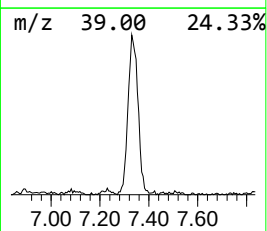
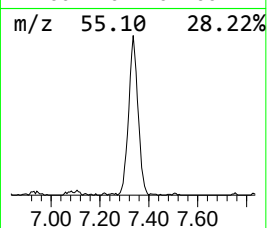
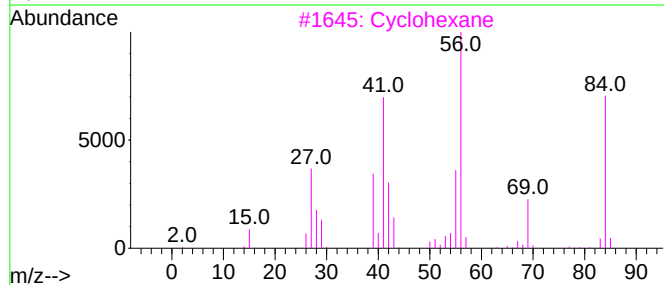
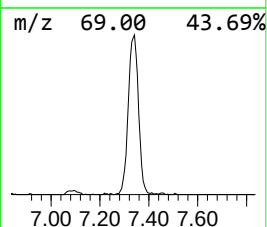
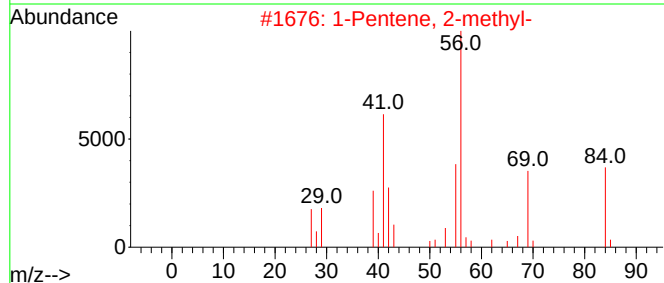
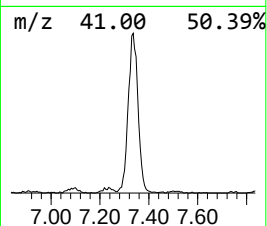
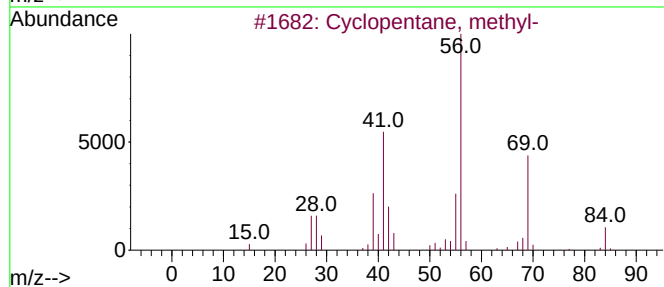
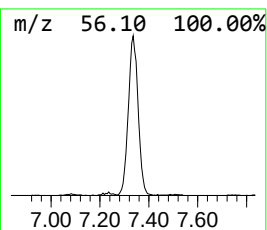
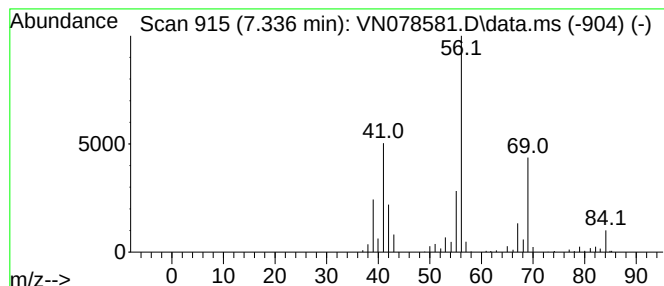
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.336	20.19 ug/l	1226730	Pentafluorobenzene	8.236

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	83
3		Cyclohexane	84	C6H12	000110-82-7	80
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64



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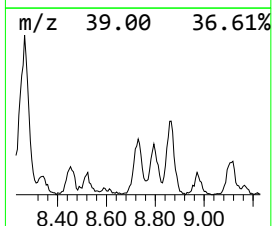
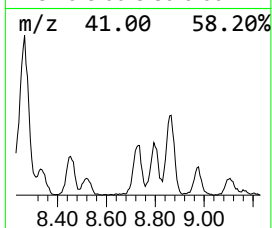
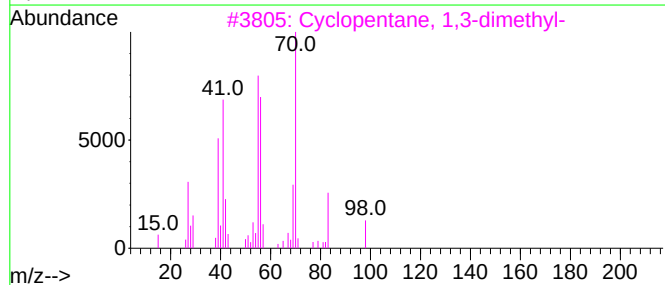
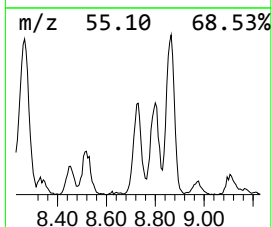
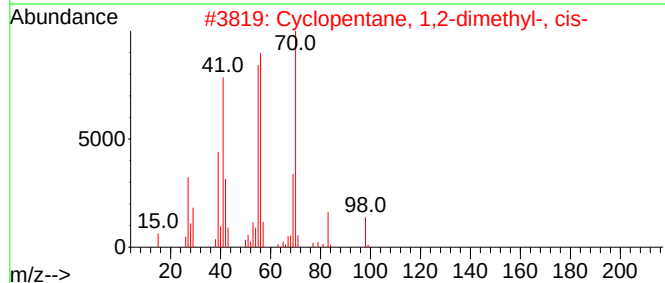
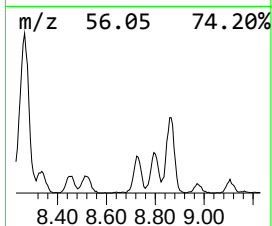
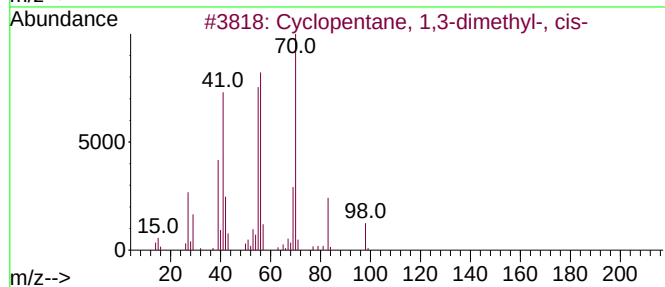
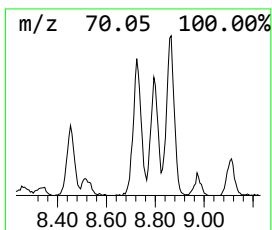
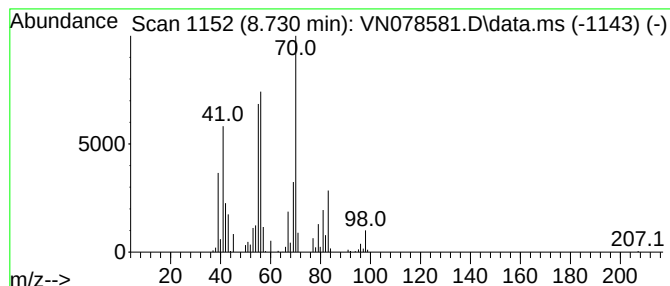
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, 1,3-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.730	15.54 ug/l	640563	1,4-Difluorobenzene	9.112

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	94
2			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	93
3			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	87
4			Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	81
5			1-Heptanethiol	132	C7H16S	001639-09-4	59



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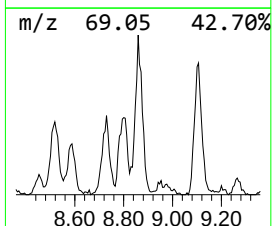
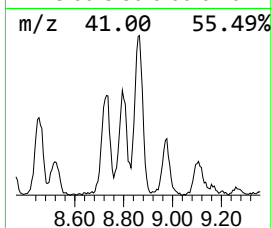
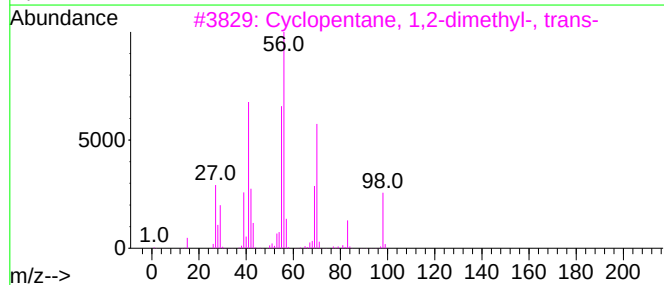
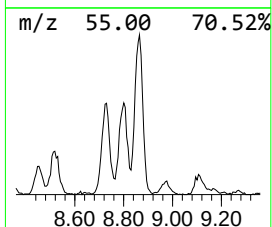
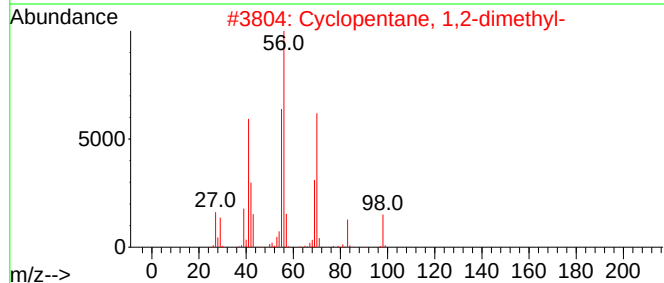
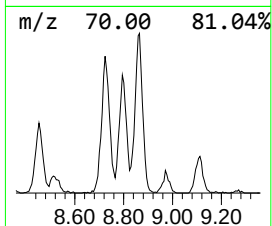
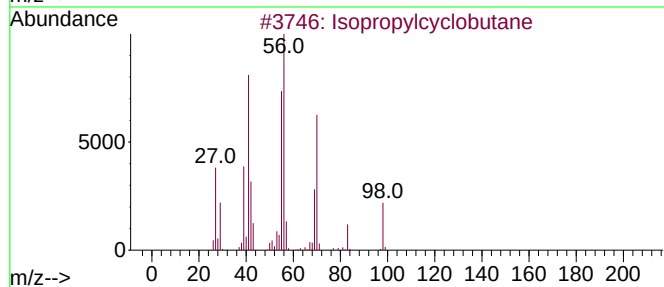
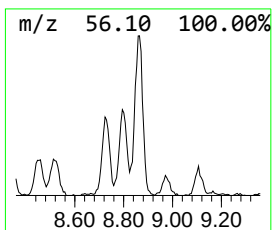
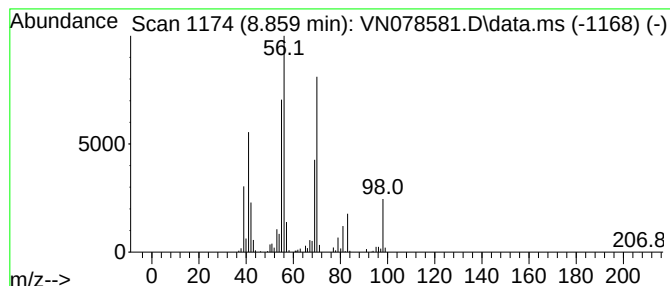
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
Quant Title : SW846 8260

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

Peak Number 3 Isopropylcyclobutane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.859	20.52 ug/l	846228	1,4-Difluorobenzene	9.112

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropylcyclobutane	98	C7H14	000872-56-0	93
2			Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
3			Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
4			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	80
5			1-Heptene	98	C7H14	000592-76-7	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48719

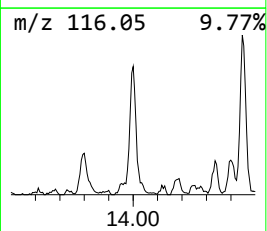
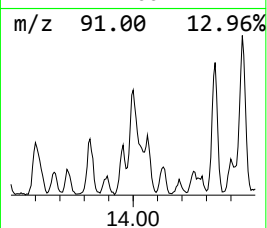
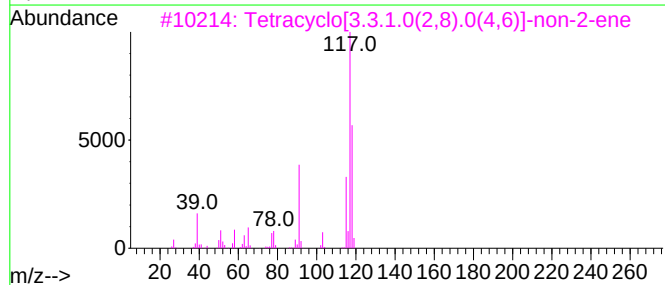
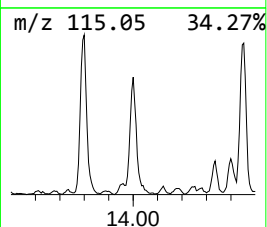
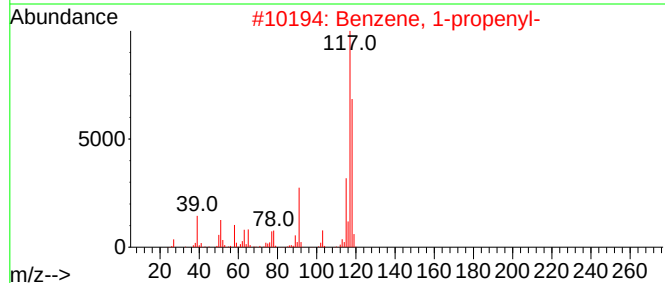
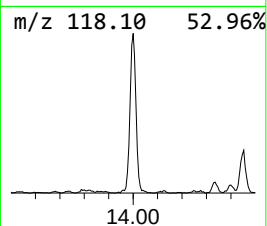
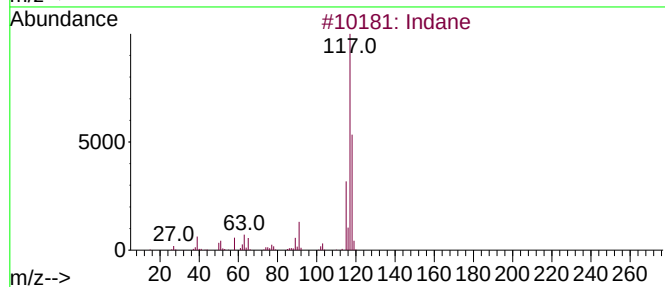
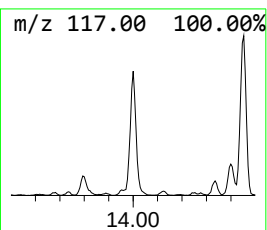
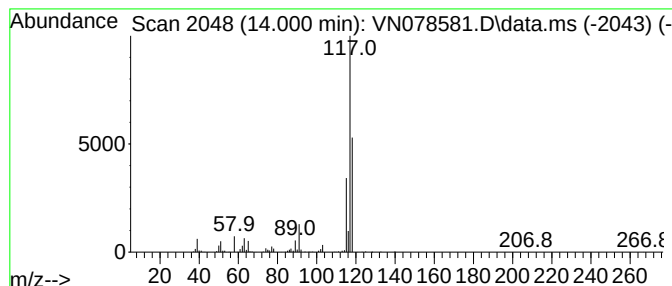
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.000	35.83 ug/l	1991520	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	80
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
4			Deltacyclene	118	C9H10	007785-10-6	72
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
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Operator : JC\MD
Sample : 03668-05
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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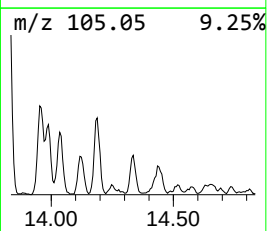
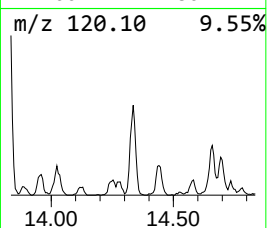
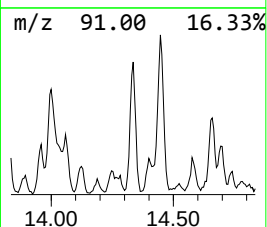
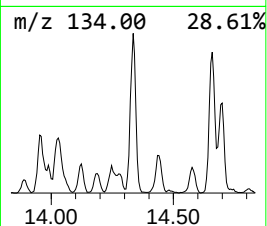
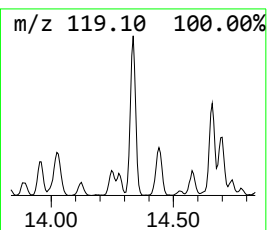
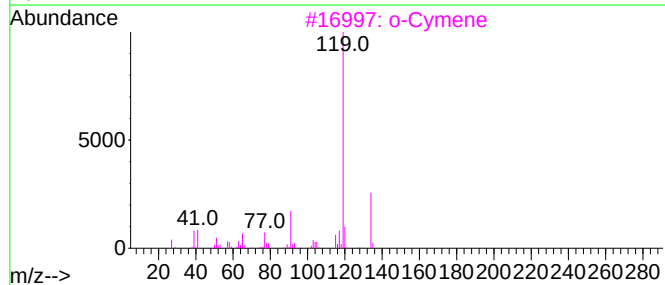
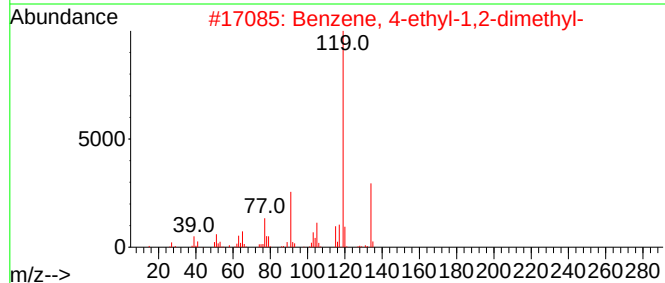
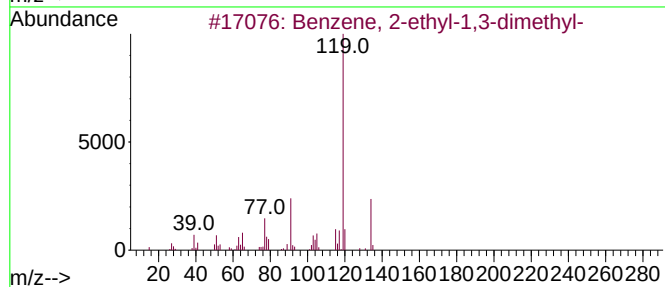
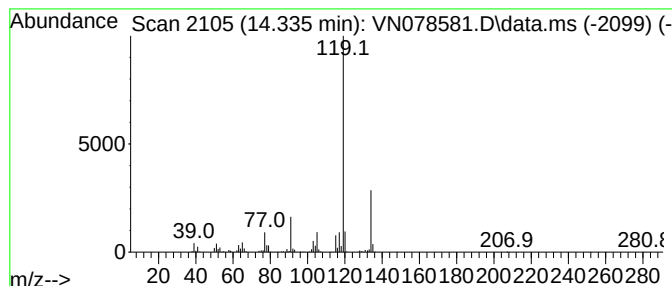
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.335	21.55 ug/l	1197840	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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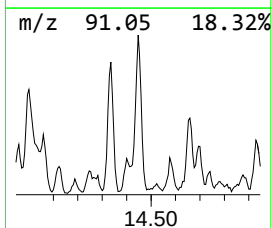
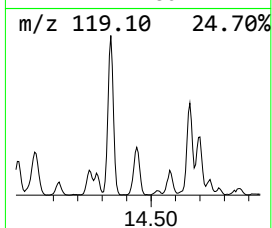
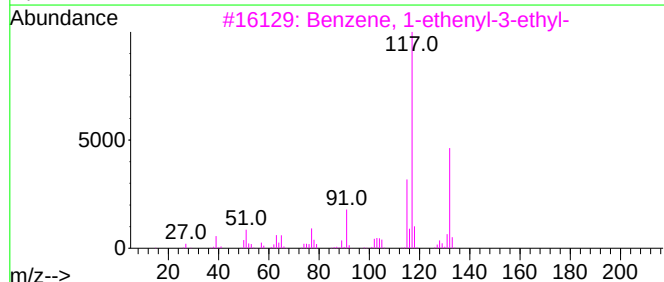
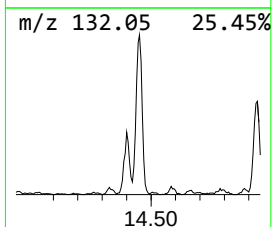
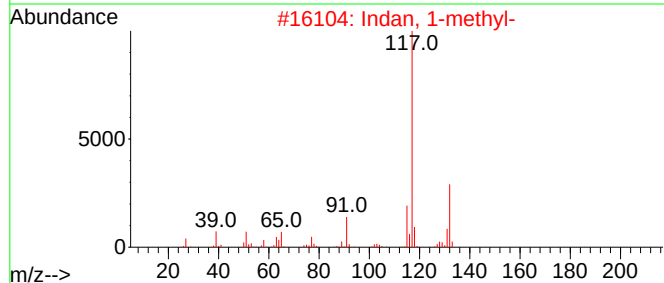
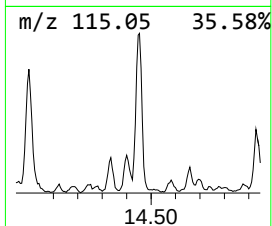
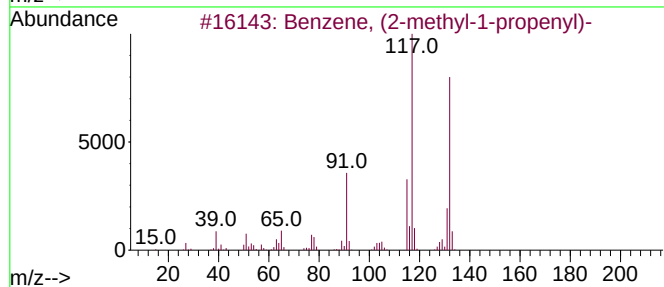
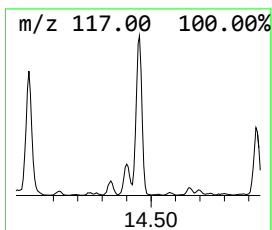
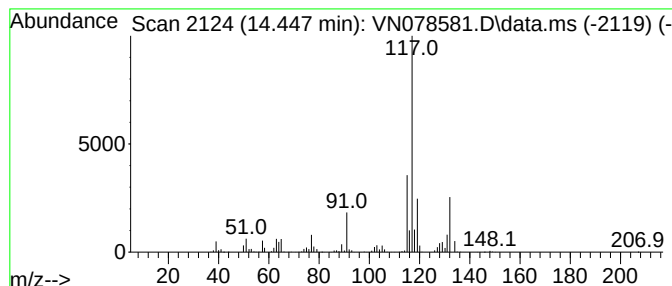
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.447	32.95 ug/l	1831740	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	80
2			Indan, 1-methyl-	132	C10H12	000767-58-8	76
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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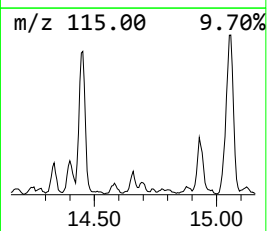
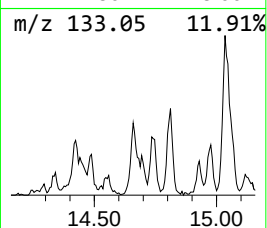
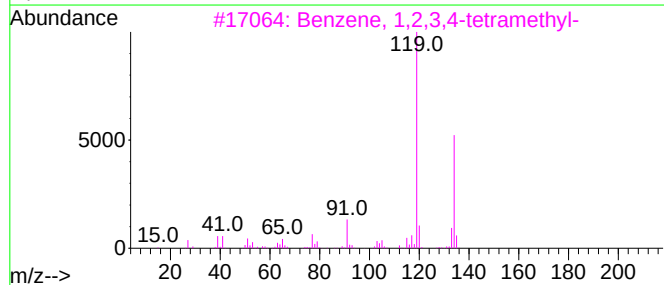
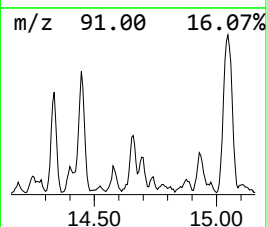
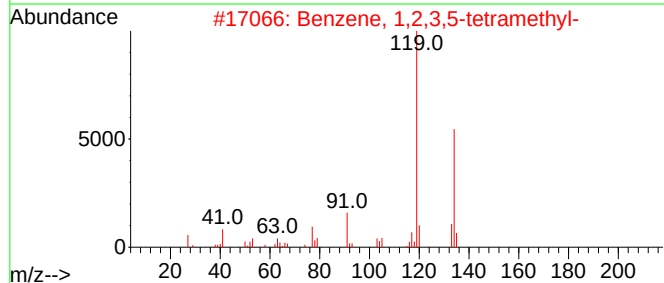
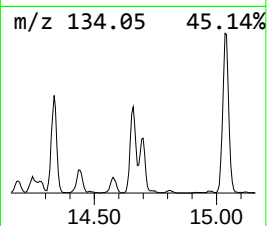
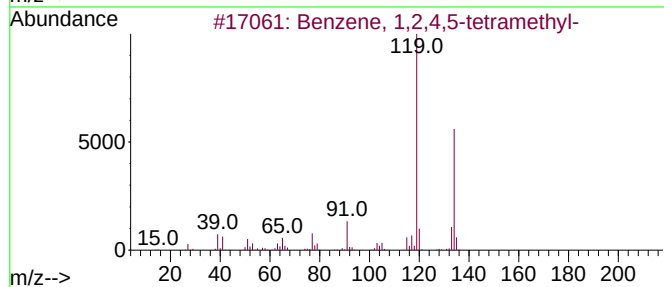
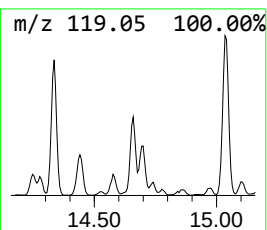
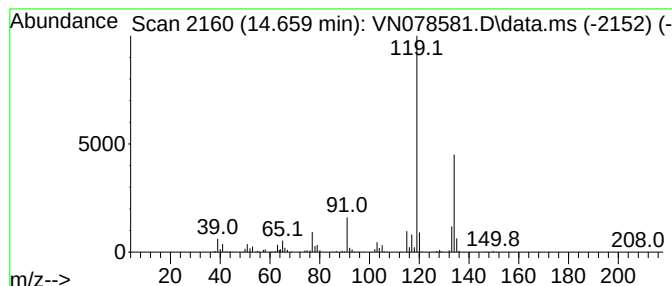
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.659	14.43 ug/l	802420	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48719

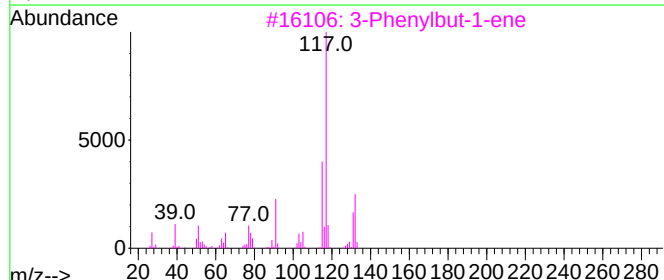
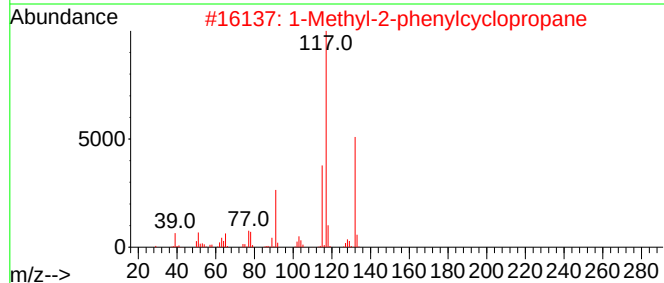
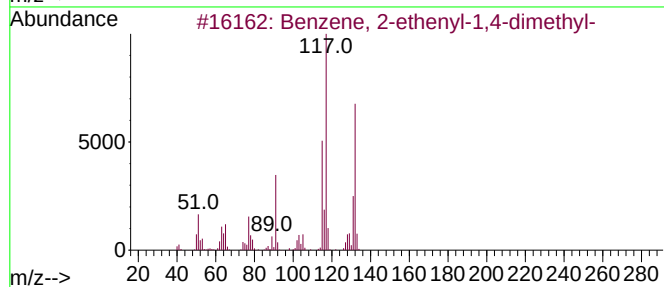
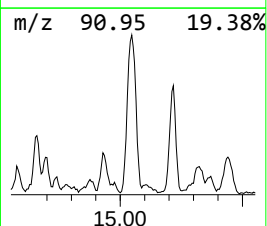
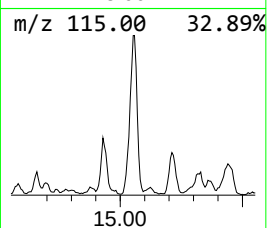
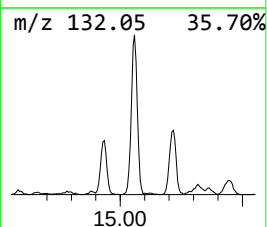
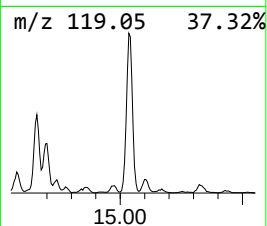
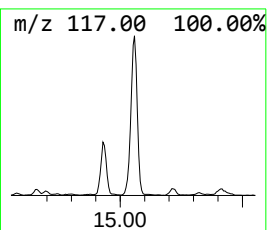
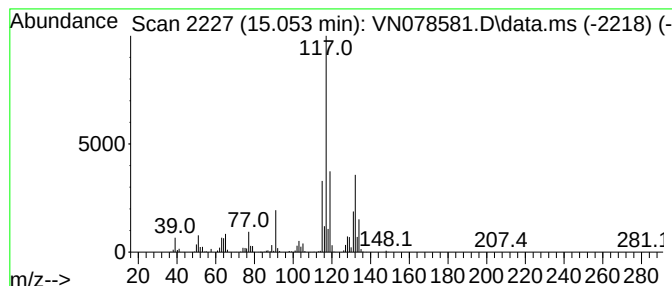
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	69.75 ug/l	3877160	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	92
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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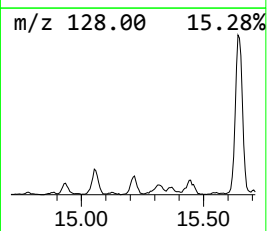
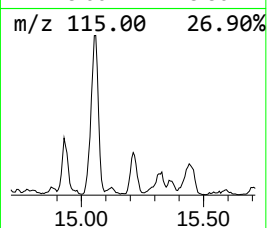
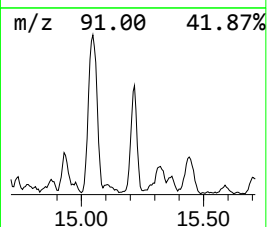
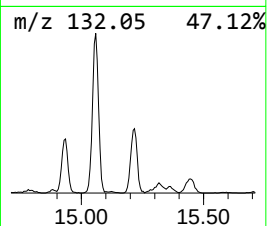
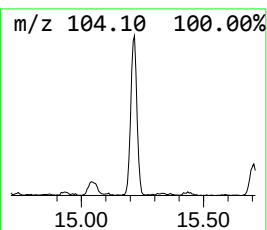
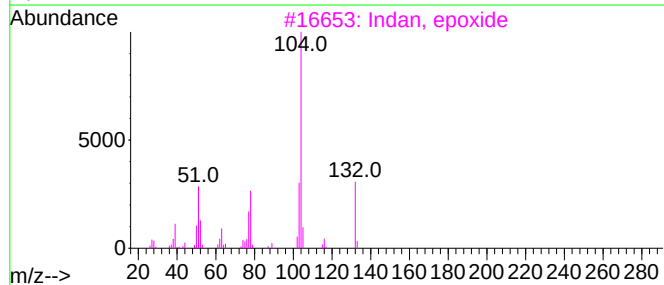
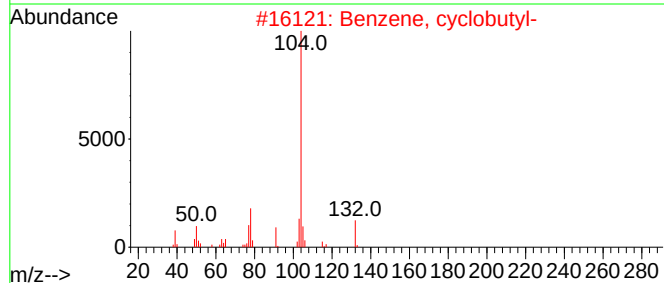
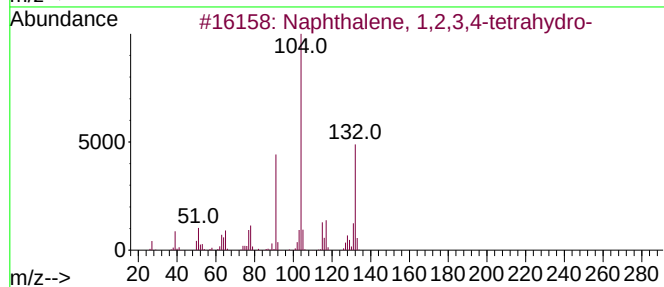
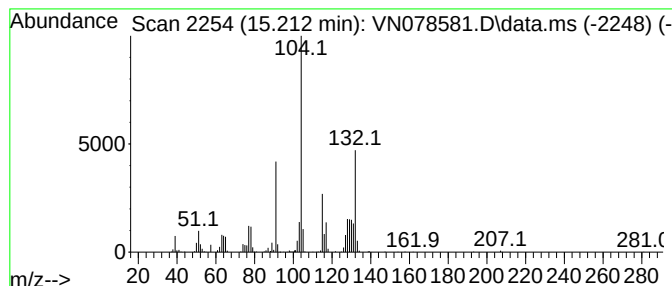
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.212	18.02 ug/l	1001970	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Benzene, cyclobutyl-	132	C10H12	004392-30-7	50
3			Indan, epoxide	132	C9H8O	000768-22-9	49
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	49
5			5,5-Dimethyl-4-phenyldihydrofura...	190	C12H14O2	013133-96-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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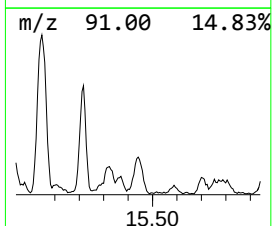
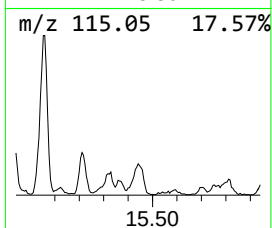
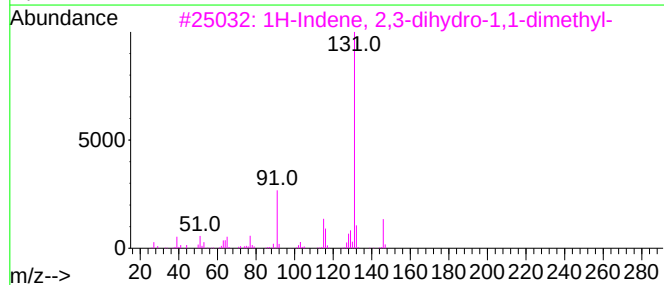
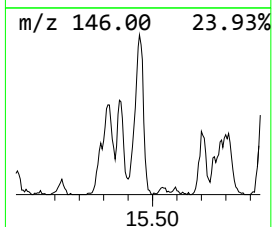
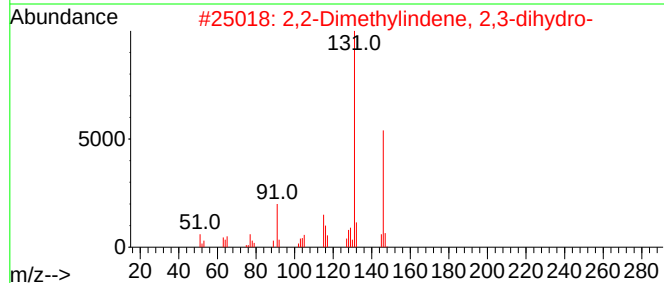
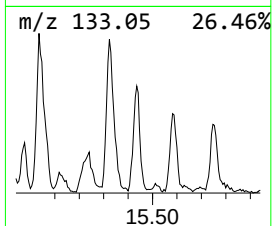
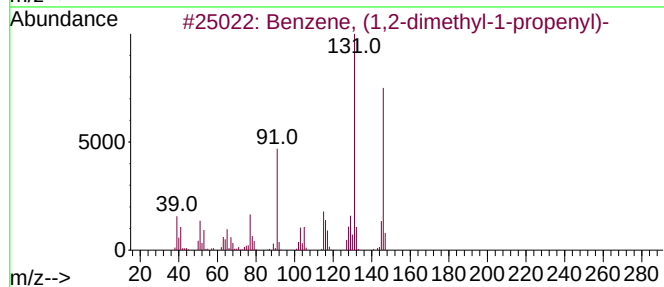
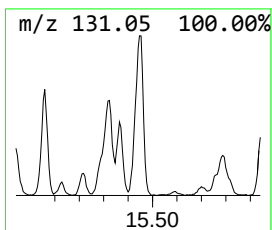
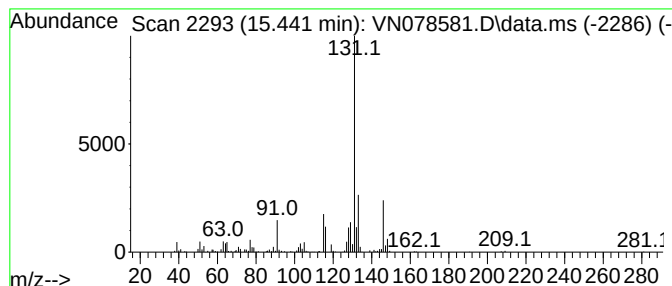
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.441	18.12 ug/l	1007120	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
2			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
3			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071923\
Data File : VN078581.D
Acq On : 19 Jul 2023 16:43
Operator : JC\MD
Sample : 03668-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48719

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.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, m...	7.336	20.2	ug/l	1226730	1	8.236	3038280	50.0
Cyclopentane, 1...	8.730	15.5	ug/l	640563	2	9.112	2061550	50.0
Isopropylcyclob...	8.859	20.5	ug/l	846228	2	9.112	2061550	50.0
Indane	14.000	35.8	ug/l	1991520	4	13.794	2779500	50.0
Benzene, 2-ethy...	14.335	21.6	ug/l	1197840	4	13.794	2779500	50.0
Benzene, (2-met...	14.447	33.0	ug/l	1831740	4	13.794	2779500	50.0
Benzene, 1,2,4,...	14.659	14.4	ug/l	802420	4	13.794	2779500	50.0
Benzene, 2-ethe...	15.053	69.8	ug/l	3877160	4	13.794	2779500	50.0
Naphthalene, 1,...	15.212	18.0	ug/l	1001970	4	13.794	2779500	50.0
Benzene, (1,2-d...	15.441	18.1	ug/l	1007120	4	13.794	2779500	50.0