

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
 Data File : VN076791.D
 Acq On : 23 Feb 2023 14:10
 Operator : JC\MD
 Sample : 01664-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-10-0222

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\82N021323W.M
 Title : SW846 8260

Signal : TIC: VN076791.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.837	143	153	163	rVB	24028	75165	2.43%	0.211%
2	3.254	216	224	238	rVB	145530	359044	11.61%	1.006%
3	4.437	411	425	435	rBV6	16223	58942	1.91%	0.165%
4	5.289	556	570	574	rBV3	30538	99358	3.21%	0.278%
5	5.395	576	588	602	rVV4	142600	612599	19.81%	1.716%
6	5.525	602	610	624	rVB3	44758	155195	5.02%	0.435%
7	5.807	642	658	679	rBV	869054	3091889	100.00%	8.660%
8	6.807	817	828	836	rBV4	15299	47594	1.54%	0.133%
9	7.272	892	907	911	rBV2	167963	494962	16.01%	1.386%
10	7.336	911	918	942	rVB	491196	1568933	50.74%	4.394%
11	8.013	1025	1033	1043	rVB3	27827	70048	2.27%	0.196%
12	8.172	1050	1060	1064	rBV	222977	529529	17.13%	1.483%
13	8.230	1064	1070	1083	rVB2	437315	1508542	48.79%	4.225%
14	8.448	1099	1107	1114	rBV3	22833	52086	1.68%	0.146%
15	8.583	1122	1130	1142	rBV2	258862	608467	19.68%	1.704%
16	8.730	1147	1155	1159	rVV6	82432	221934	7.18%	0.622%
17	8.789	1160	1165	1172	rVV4	109709	274853	8.89%	0.770%
18	8.854	1172	1176	1185	rVB3	47405	104332	3.37%	0.292%
19	9.107	1210	1219	1232	rBV	663328	1448801	46.86%	4.058%
20	9.201	1233	1235	1243	rVB3	27722	41641	1.35%	0.117%
21	9.342	1251	1259	1261	rBV4	34443	67368	2.18%	0.189%
22	9.389	1261	1267	1275	rVB	100646	211650	6.85%	0.593%
23	9.572	1285	1298	1299	rBV3	118075	227002	7.34%	0.636%
24	9.607	1299	1304	1319	rVB	220930	499375	16.15%	1.399%
25	9.777	1327	1333	1340	rVB3	46367	91376	2.96%	0.256%
26	9.936	1351	1360	1363	rBV6	24469	64990	2.10%	0.182%
27	9.966	1363	1365	1370	rVV4	26790	47282	1.53%	0.132%
28	10.019	1370	1374	1380	rVB3	33156	61754	2.00%	0.173%
29	10.107	1382	1389	1393	rBV	98040	201131	6.51%	0.563%
30	10.148	1393	1396	1409	rVB4	43113	95957	3.10%	0.269%
31	10.454	1441	1448	1460	rVB2	136558	297187	9.61%	0.832%
32	10.566	1460	1467	1478	rBV	980930	1885865	60.99%	5.282%
33	10.760	1492	1500	1508	rVV2	123982	265863	8.60%	0.745%
34	10.830	1508	1512	1520	rVB2	81073	153540	4.97%	0.430%
35	10.983	1531	1538	1547	rBV4	72680	189314	6.12%	0.530%
36	11.060	1547	1551	1561	rVB2	47102	95601	3.09%	0.268%

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37	11.171	1561	1570	1578	rBV2	37921	86491	2.80%	0.242%
38	11.242	1578	1582	1588	rBV5	19977	40001	1.29%	0.112%
39	11.577	1633	1639	1647	rVB5	17093	32811	1.06%	0.092%
40	11.866	1680	1688	1695	rBV	867779	1547832	50.06%	4.335%
41	11.966	1695	1705	1711	rVB	189108	350429	11.33%	0.981%
42	12.077	1718	1724	1735	rVB	351246	593706	19.20%	1.663%
43	12.401	1768	1779	1788	rBV	133791	246313	7.97%	0.690%
44	12.695	1825	1829	1835	rVB2	29765	46904	1.52%	0.131%
45	12.848	1847	1855	1864	rBV	645909	1133354	36.66%	3.174%
46	13.036	1883	1887	1894	rBV2	26764	44496	1.44%	0.125%
47	13.130	1894	1903	1915	rVB	124475	245308	7.93%	0.687%
48	13.336	1930	1938	1948	rBV	1811175	2998494	96.98%	8.398%
49	13.483	1957	1963	1969	rBV	638926	1022211	33.06%	2.863%
50	13.536	1969	1972	1978	rVB2	53403	83096	2.69%	0.233%
51	13.671	1990	1995	2000	rBV3	25502	36464	1.18%	0.102%
52	13.730	2000	2005	2008	rBV3	28126	46048	1.49%	0.129%
53	13.795	2008	2016	2018	rBV	870376	1473078	47.64%	4.126%
54	13.889	2028	2032	2037	rVB4	28754	47624	1.54%	0.133%
55	13.954	2037	2043	2045	rBV2	297056	461950	14.94%	1.294%
56	13.995	2045	2050	2055	rVV	895409	1654990	53.53%	4.635%
57	14.036	2055	2057	2064	rVB	170673	217753	7.04%	0.610%
58	14.118	2064	2071	2076	rBV	94681	161685	5.23%	0.453%
59	14.183	2076	2082	2088	rBV2	183482	327761	10.60%	0.918%
60	14.248	2088	2093	2095	rVV2	79430	136527	4.42%	0.382%
61	14.277	2095	2098	2102	rVV	140911	206320	6.67%	0.578%
62	14.330	2102	2107	2113	rVB	686667	1088104	35.19%	3.048%
63	14.395	2113	2118	2121	rBV2	90096	142556	4.61%	0.399%
64	14.448	2121	2127	2136	rVV2	401614	704282	22.78%	1.973%
65	14.571	2144	2148	2155	rVB2	105855	182465	5.90%	0.511%
66	14.654	2155	2162	2166	rBV	428994	737405	23.85%	2.065%
67	14.695	2166	2169	2174	rVV	349080	540288	17.47%	1.513%
68	14.736	2174	2176	2180	rVB2	48641	59587	1.93%	0.167%
69	14.877	2192	2200	2204	rBV4	47411	94132	3.04%	0.264%
70	14.930	2204	2209	2213	rVV3	99823	189993	6.14%	0.532%
71	14.971	2213	2216	2220	rVV	43769	65677	2.12%	0.184%
72	15.054	2220	2230	2236	rVV2	497902	1157948	37.45%	3.243%
73	15.118	2236	2241	2247	rVB6	67523	150036	4.85%	0.420%
74	15.207	2252	2256	2261	rVB3	44025	75835	2.45%	0.212%
75	15.324	2270	2276	2289	rBV3	153785	442147	14.30%	1.238%
76	15.436	2289	2295	2303	rVB3	145537	362330	11.72%	1.015%

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77	15.642	2325	2330	2338	rVB2	68457	131575	4.26%	0.369%
78	15.748	2338	2348	2354	rVV2	71481	150686	4.87%	0.422%
79	15.806	2354	2358	2371	rVV8	30696	78720	2.55%	0.220%
80	15.948	2375	2382	2388	rVB3	26869	54372	1.76%	0.152%
81	16.130	2404	2413	2421	rVV2	34610	100959	3.27%	0.283%
82	16.201	2421	2425	2430	rVV6	17547	34360	1.11%	0.096%
83	16.265	2430	2436	2440	rVB5	20421	39518	1.28%	0.111%

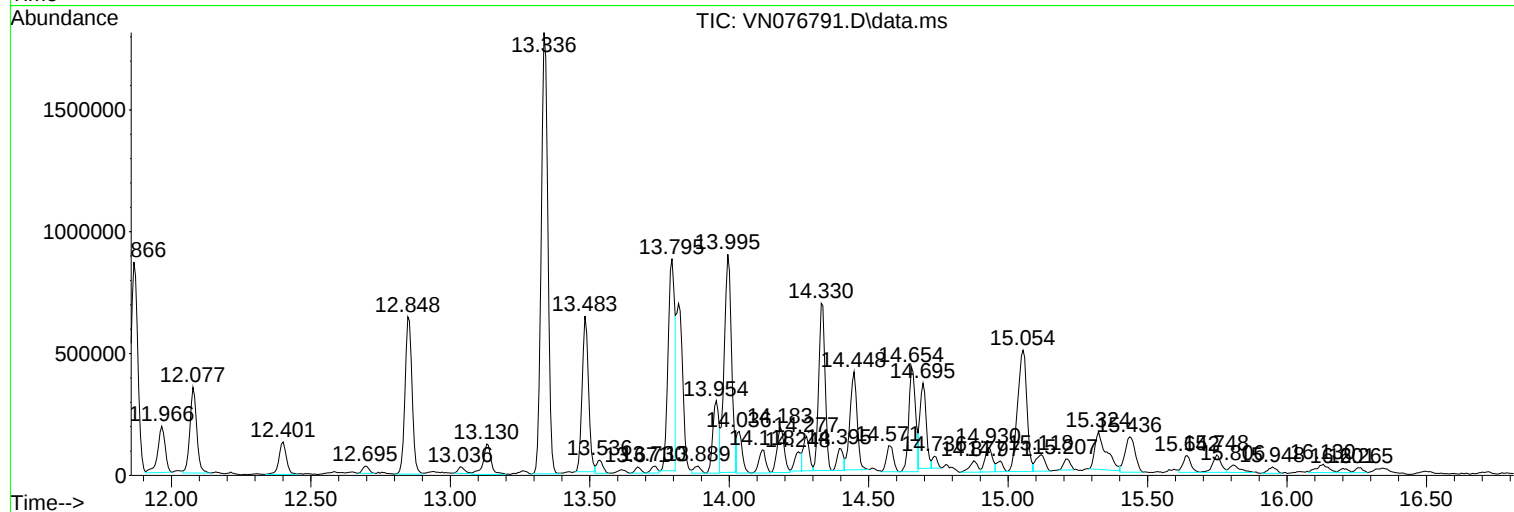
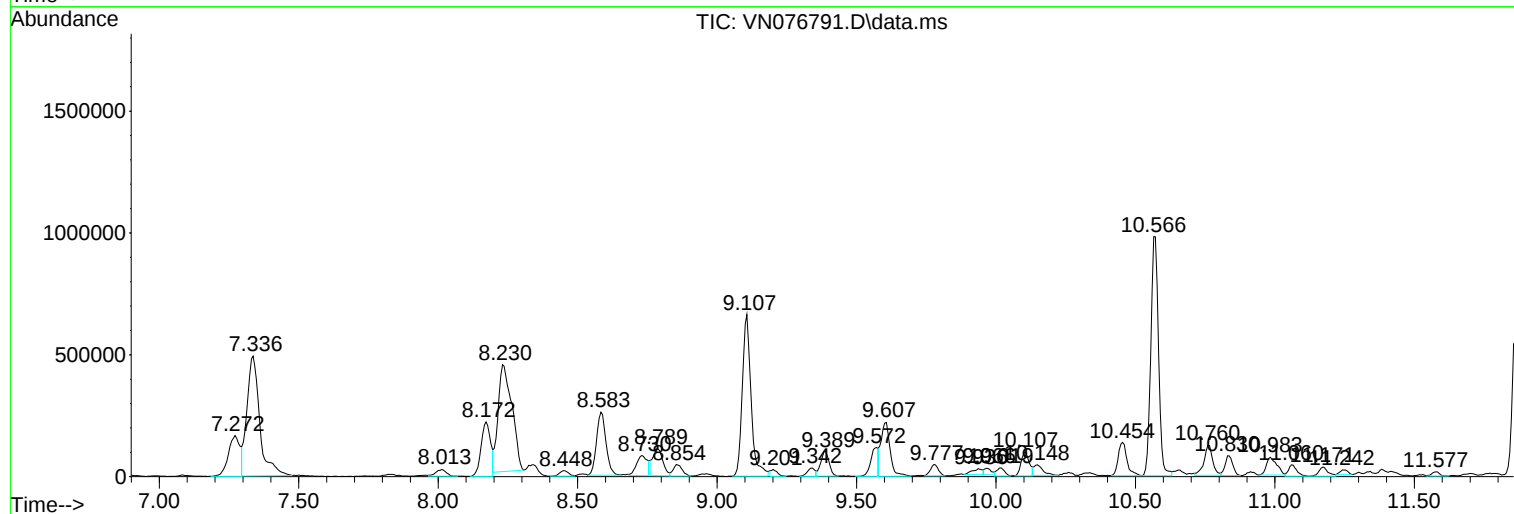
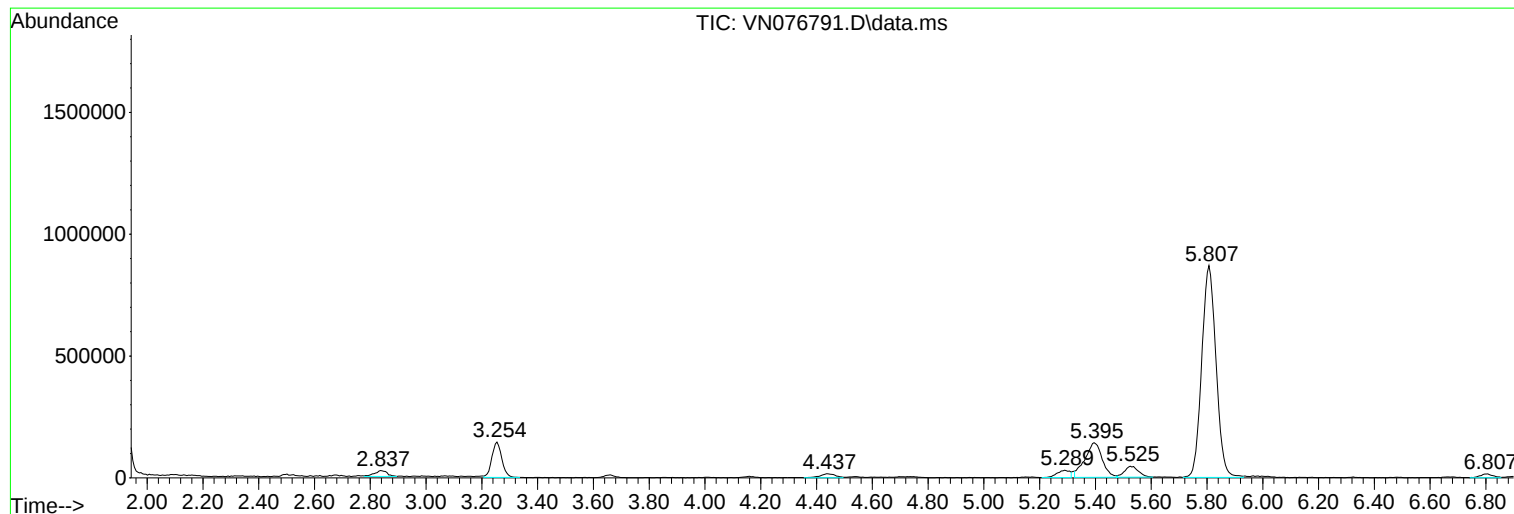
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.M
Quant Title : SW846 8260

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



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Instrument :
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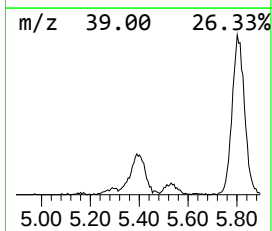
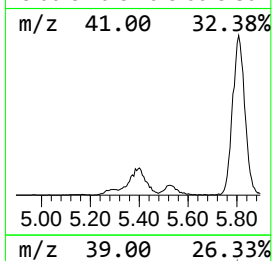
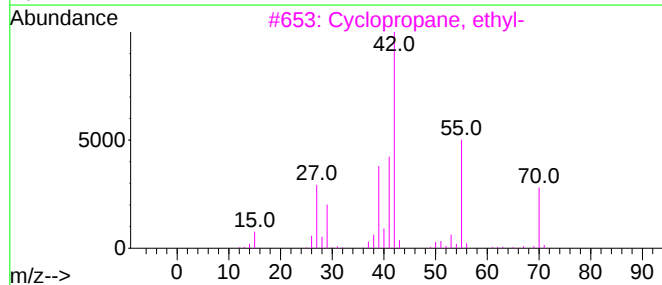
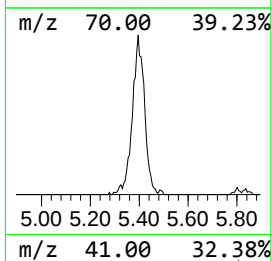
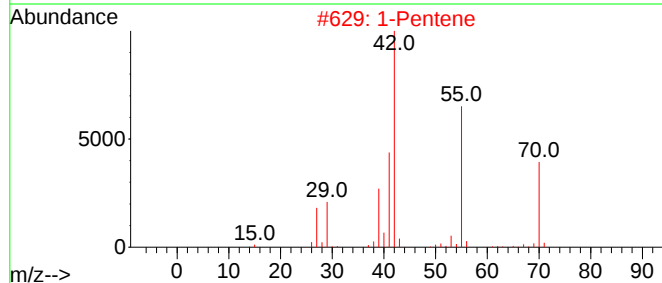
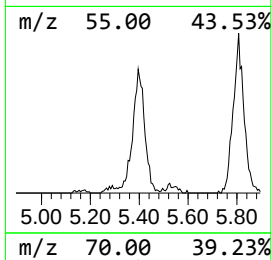
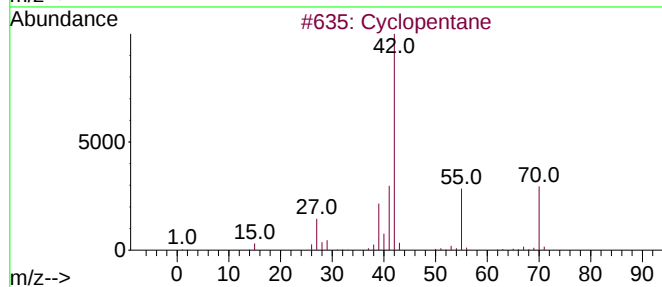
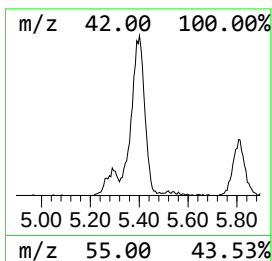
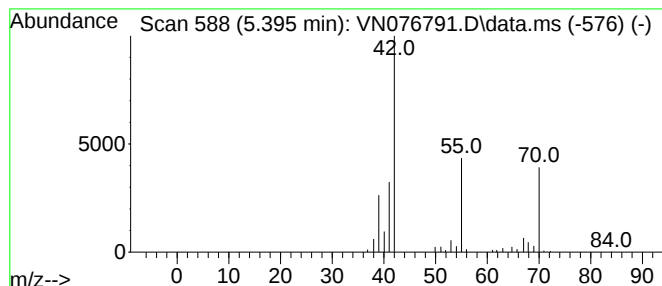
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.395	20.30 ug/l	612599	Pentafluorobenzene	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane	70	C5H10	000287-92-3	72
2			1-Pentene	70	C5H10	000109-67-1	64
3			Cyclopropane, ethyl-	70	C5H10	001191-96-4	45
4			Isobutyronitrile	69	C4H7N	000078-82-0	9
5			Cyclobutane, methyl-	70	C5H10	000598-61-8	9



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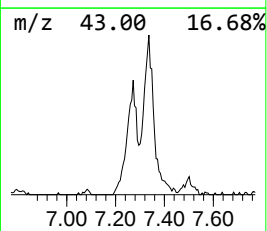
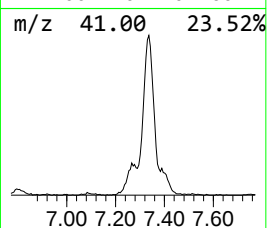
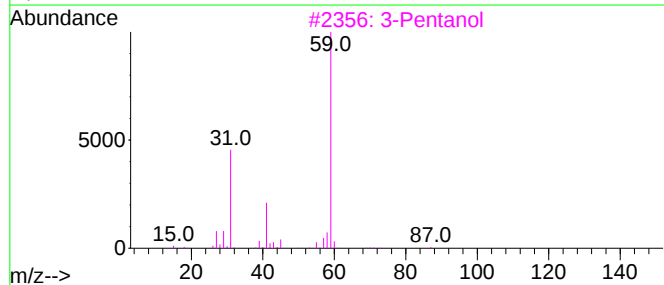
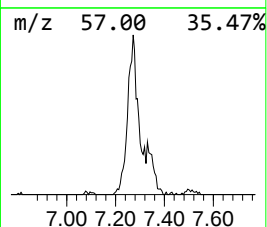
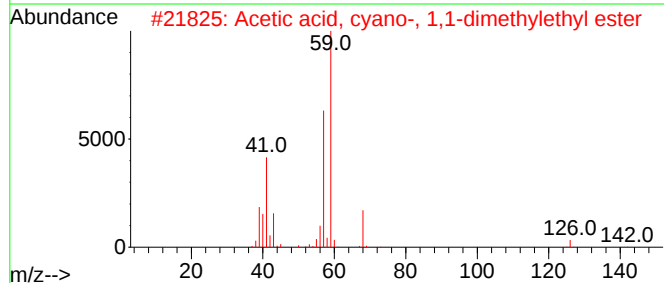
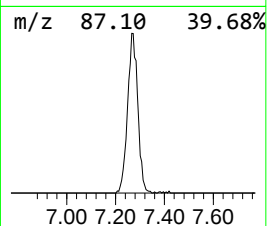
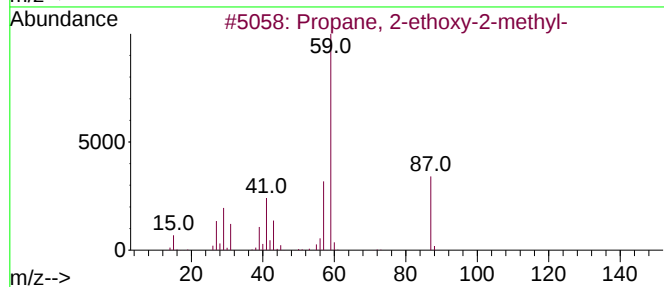
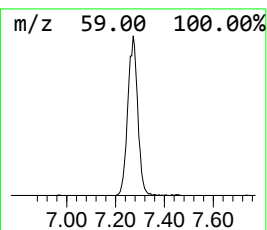
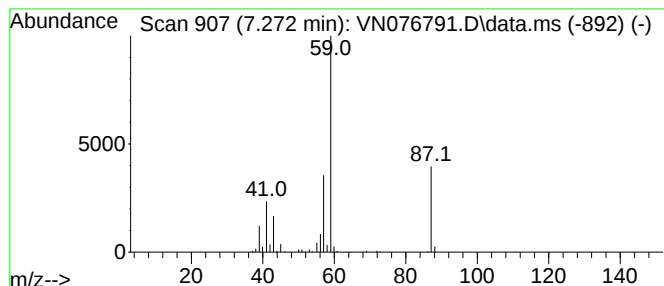
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.M
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propane, 2-ethoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.272	16.41 ug/l	494962	Pentafluorobenzene	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	16
3			3-Pentanol	88	C5H12O	000584-02-1	9
4			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9
5			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	9



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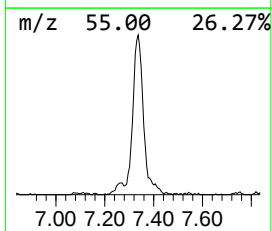
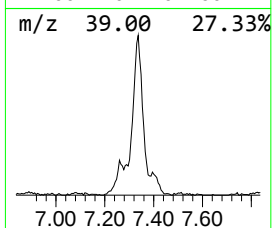
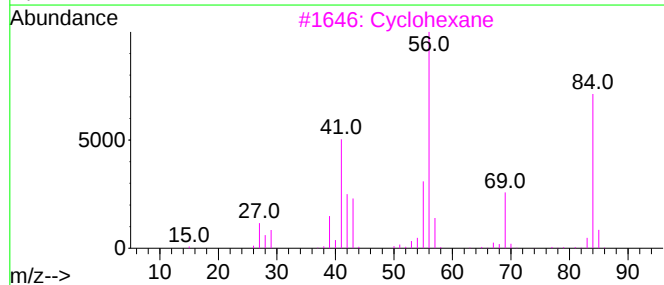
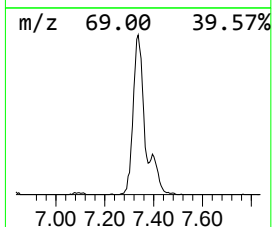
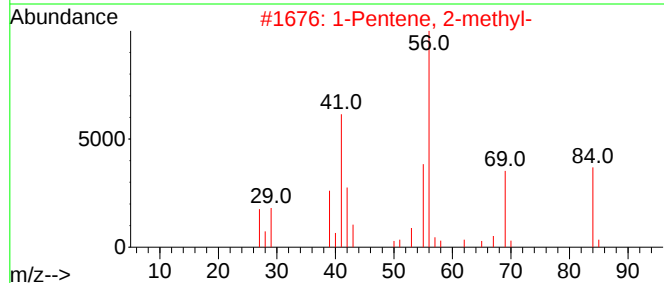
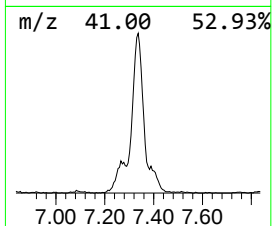
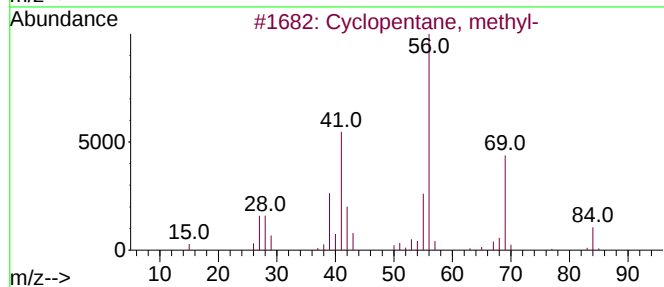
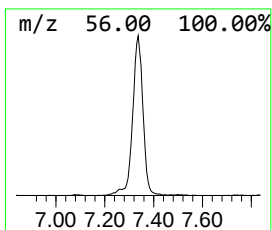
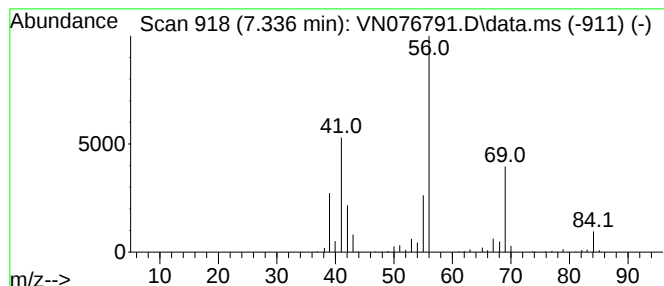
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.336	52.00 ug/l	1568930	Pentafluorobenzene	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80
3			Cyclohexane	84	C6H12	000110-82-7	78
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



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MW-10-0222

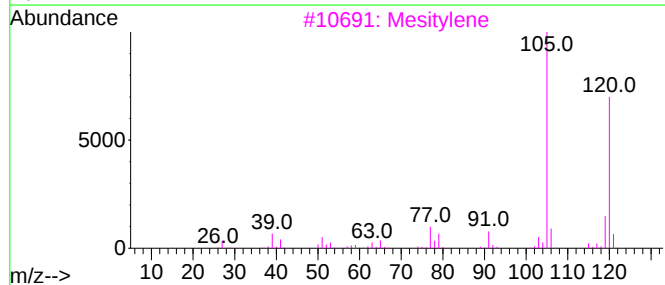
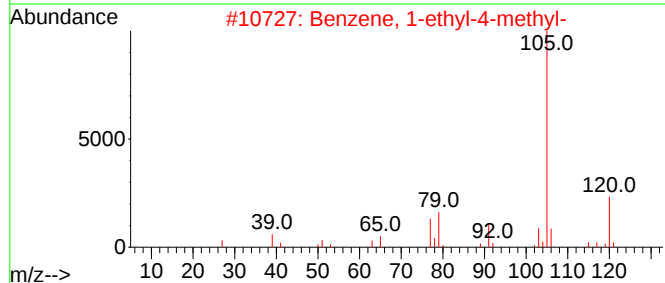
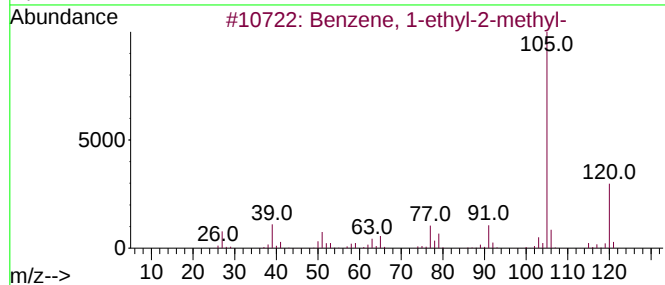
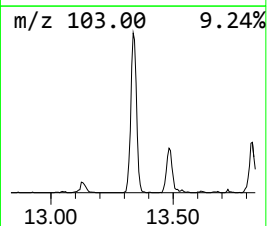
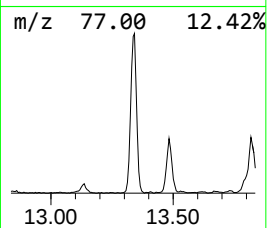
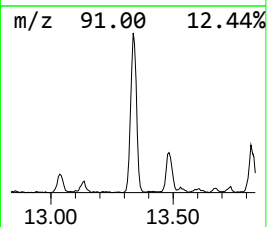
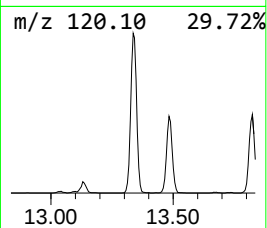
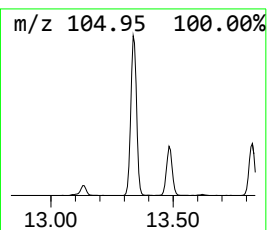
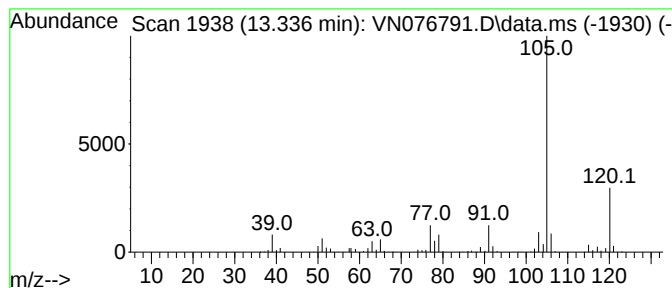
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.336	101.78 ug/l	2998490	1,4-Dichlorobenzene-d4	13.795

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-4-methyl-	120	C9H12	000622-96-8	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

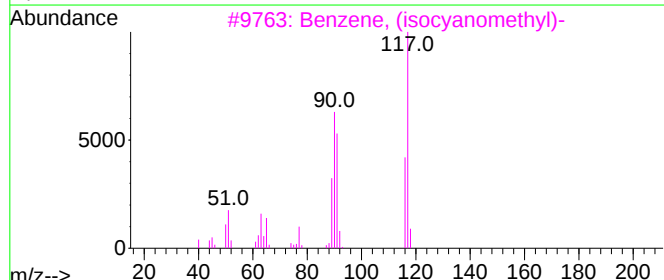
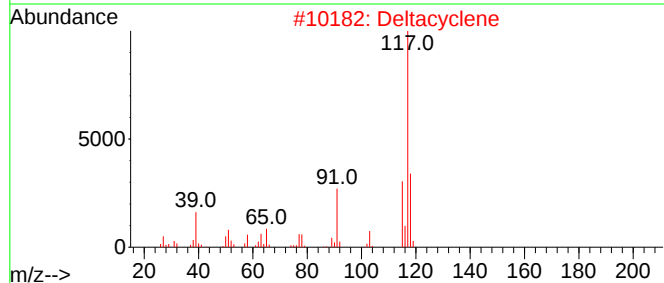
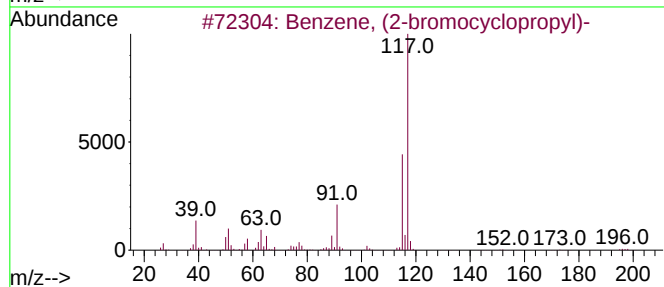
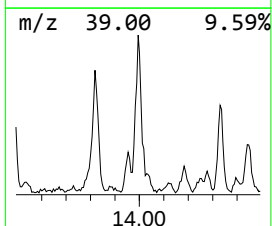
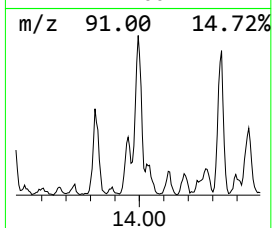
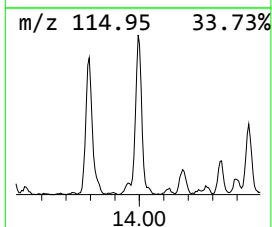
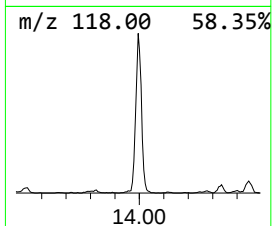
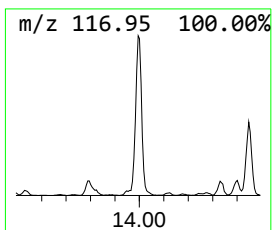
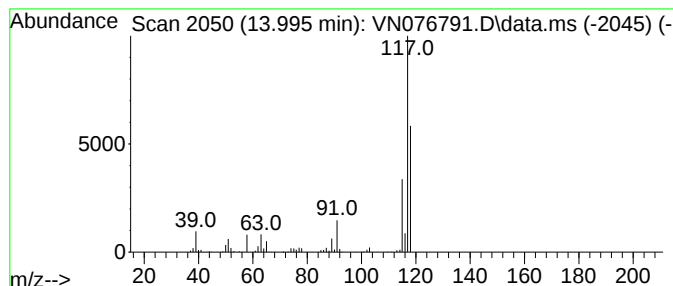
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-bromocyclopropyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.995	56.17 ug/l	1654990	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	59
2			Deltacyclene	118	C9H10	007785-10-6	53
3			Benzene, (isocyanomethyl)-	117	C8H7N	010340-91-7	47
4			m-Aminophenylacetylene	117	C8H7N	054060-30-9	43
5			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

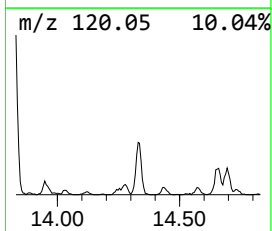
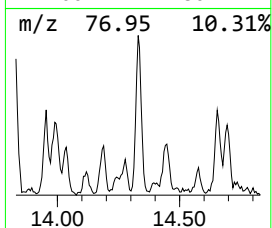
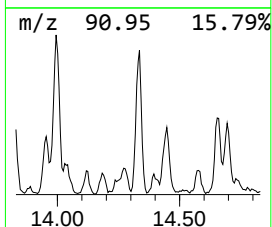
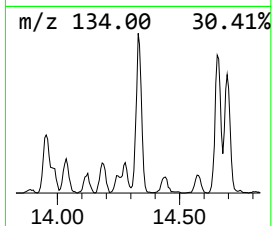
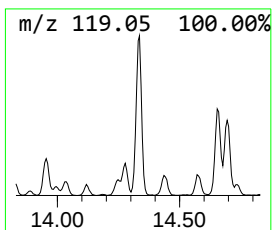
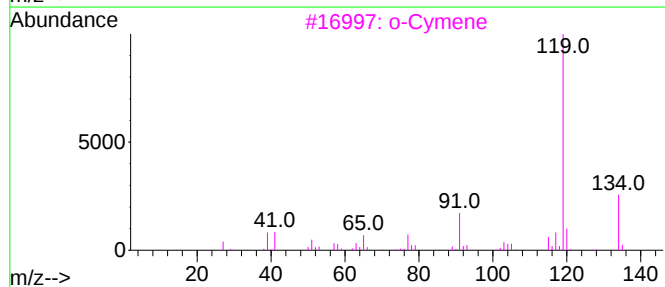
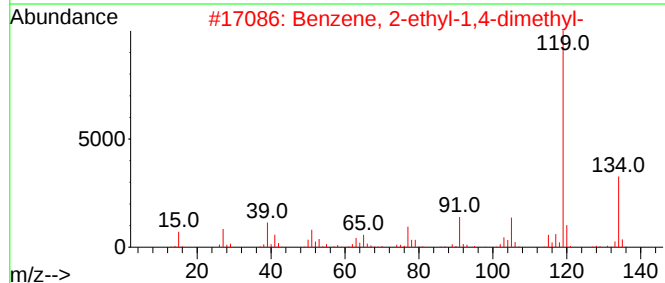
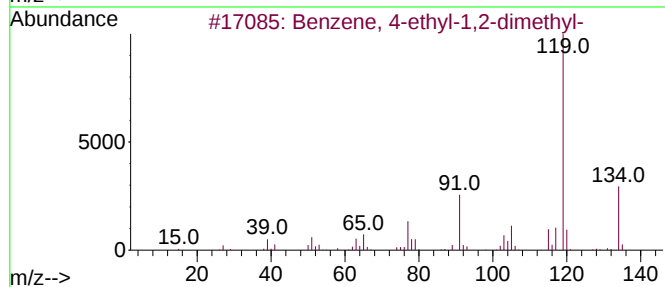
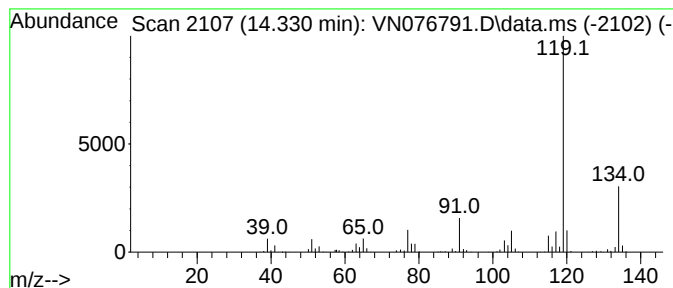
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.330	36.93 ug/l	1088100	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
3			o-Cymene	134	C10H14	000527-84-4	94
4			p-Cymene	134	C10H14	000099-87-6	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

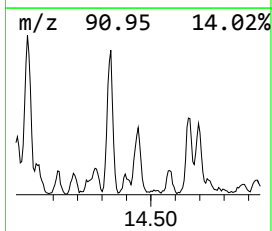
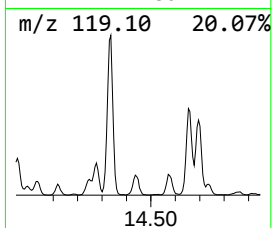
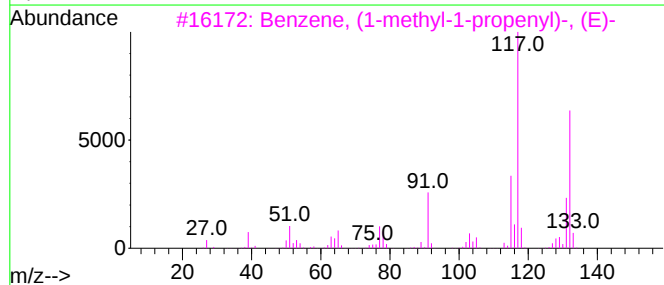
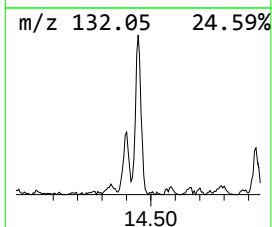
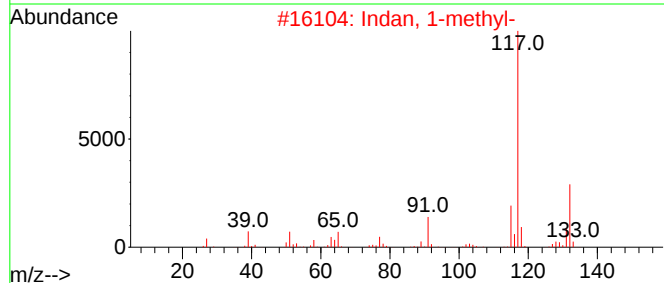
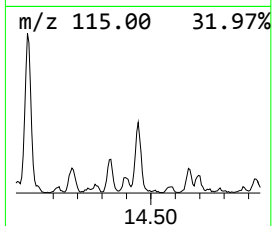
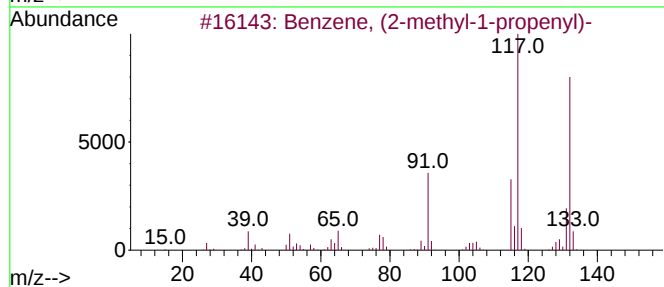
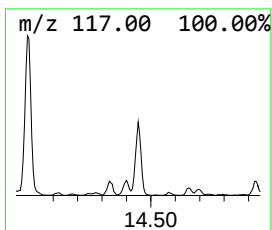
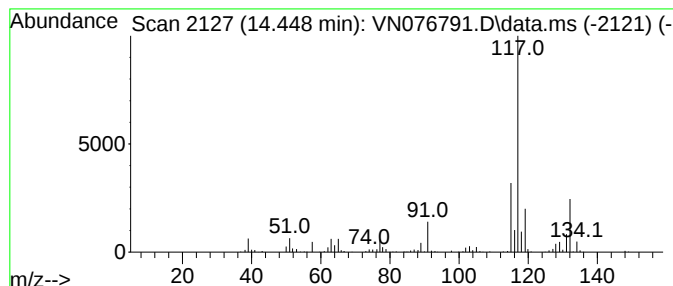
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-propenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.448	23.91 ug/l	704282	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
2			Indan, 1-methyl-	132	C10H12	000767-58-8	76
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	72
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

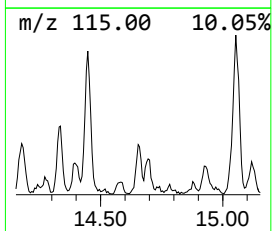
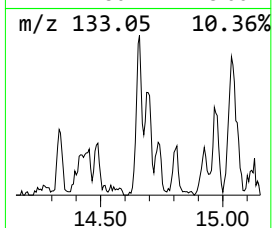
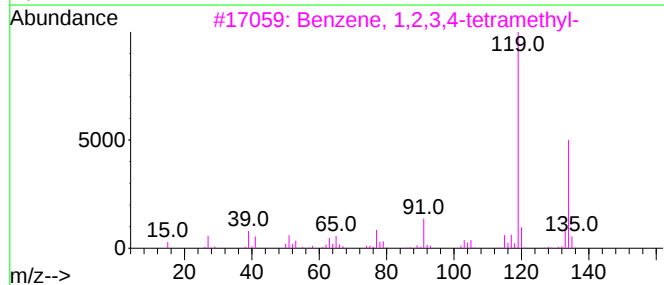
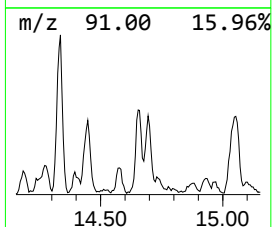
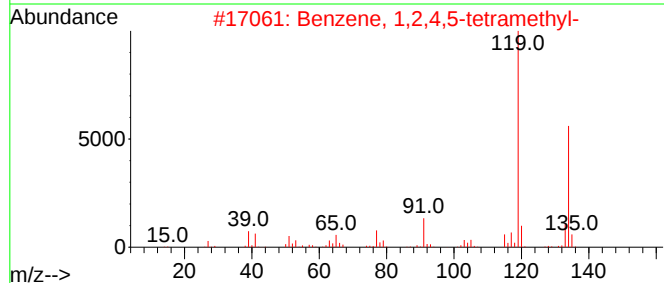
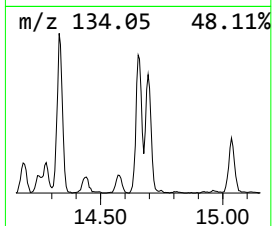
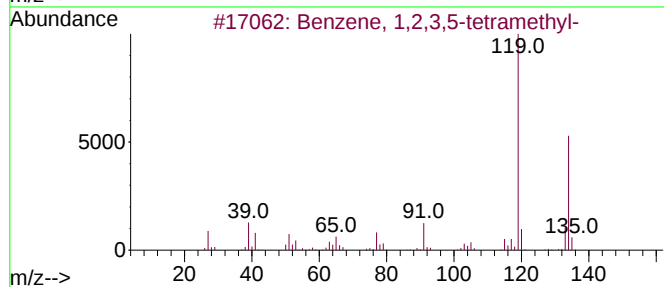
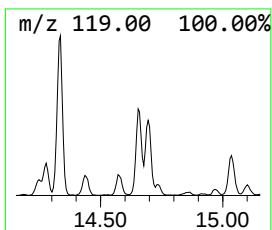
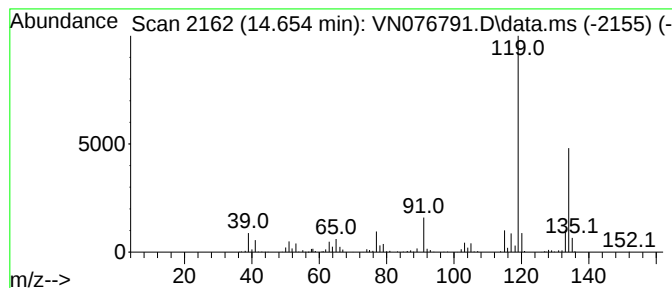
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.654	25.03 ug/l	737405	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

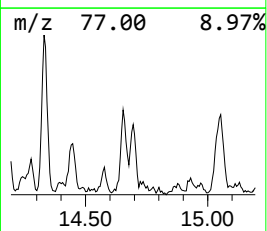
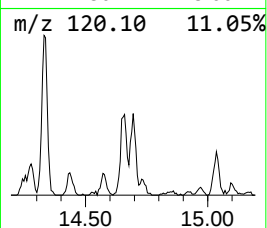
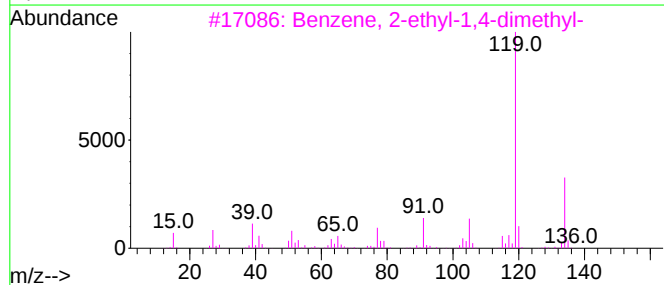
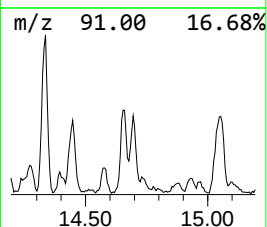
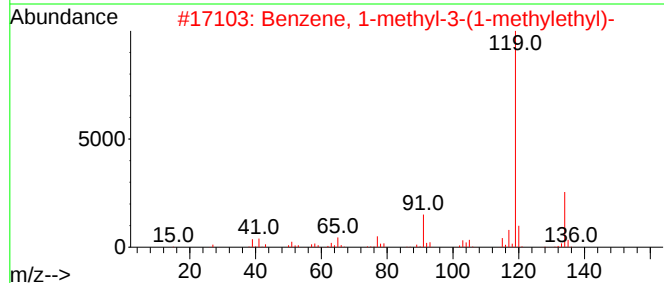
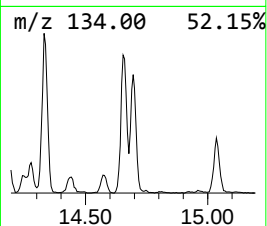
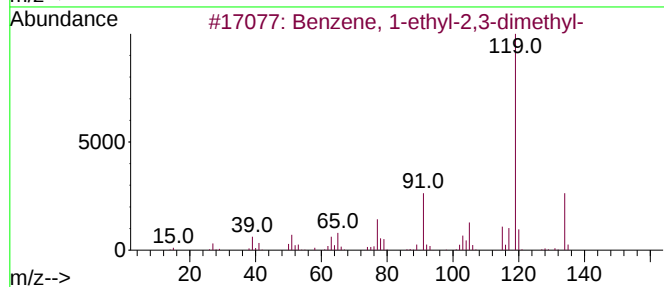
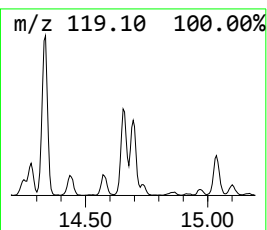
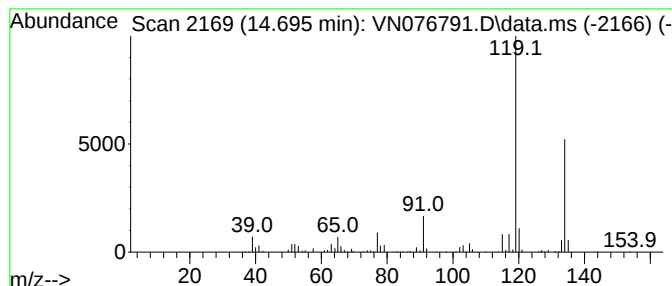
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.695	18.34 ug/l	540288	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	91
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	91
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

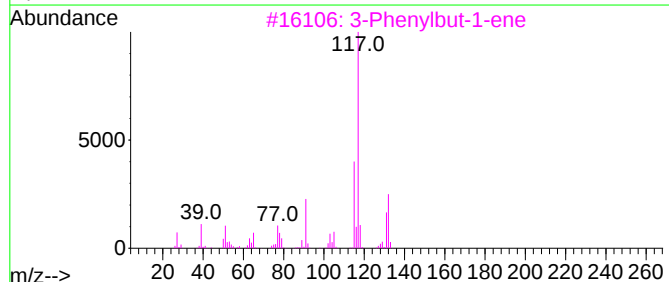
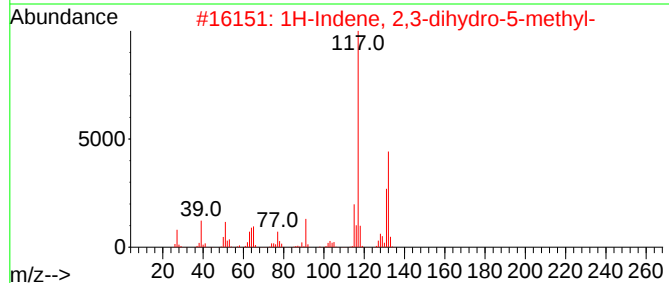
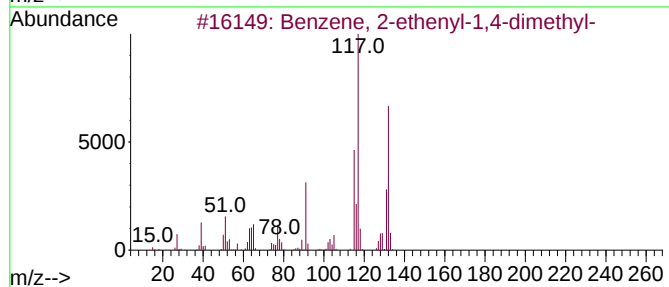
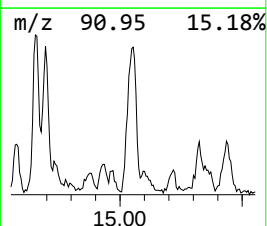
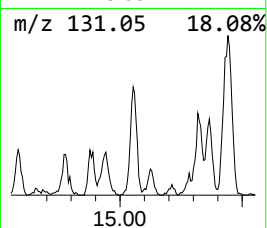
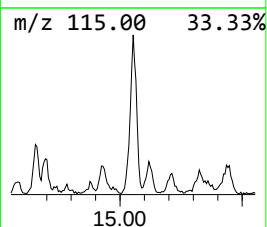
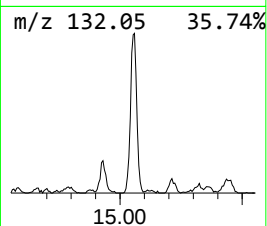
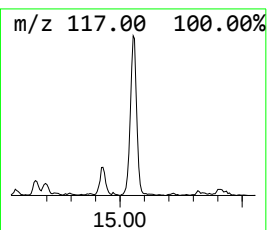
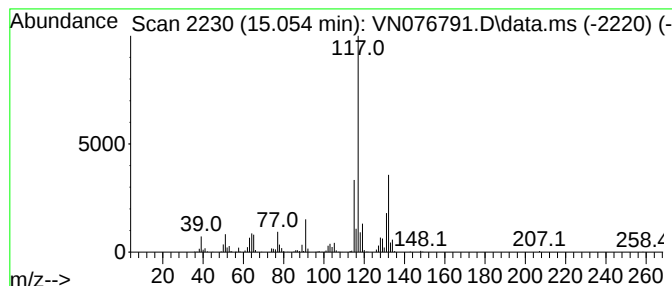
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.054	39.30 ug/l	1157950	1,4-Dichlorobenzene-d4	13.795

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	95
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022323\
Data File : VN076791.D
Acq On : 23 Feb 2023 14:10
Operator : JC\MD
Sample : 01664-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10-0222

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021323W.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	5.395	20.3	ug/l	612599	1	8.230	1508540	50.0
Propane, 2-etho...	7.272	16.4	ug/l	494962	1	8.230	1508540	50.0
Cyclopentane, m...	7.336	52.0	ug/l	1568930	1	8.230	1508540	50.0
Benzene, 1-ethy...	13.336	101.8	ug/l	2998490	4	13.795	1473080	50.0
Benzene, (2-bro...	13.995	56.2	ug/l	1654990	4	13.795	1473080	50.0
Benzene, 4-ethy...	14.330	36.9	ug/l	1088100	4	13.795	1473080	50.0
Benzene, (2-met...	14.448	23.9	ug/l	704282	4	13.795	1473080	50.0
Benzene, 1,2,3,...	14.654	25.0	ug/l	737405	4	13.795	1473080	50.0
Benzene, 1-ethy...	14.695	18.3	ug/l	540288	4	13.795	1473080	50.0
Benzene, 2-ethe...	15.054	39.3	ug/l	1157950	4	13.795	1473080	50.0