

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
 Data File : VN077480.D
 Acq On : 25 Apr 2023 19:34
 Operator : JC\MD
 Sample : 02460-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-5

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\82N042423W.M
 Title : SW846 8260

Signal : TIC: VN077480.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.525	92	100	110	rBV	2787439	4690413	6.53%	0.839%
2	3.254	213	224	251	rBV	33297193	71857175	100.00%	12.849%
3	3.654	280	292	312	rBV	6425738	16403096	22.83%	2.933%
4	4.154	367	377	392	rVB	681310	1922223	2.68%	0.344%
5	4.431	410	424	448	rVB	1885519	6637934	9.24%	1.187%
6	5.160	528	548	556	rBV8	174192	737204	1.03%	0.132%
7	5.290	556	570	572	rVV	3034144	8415005	11.71%	1.505%
8	5.354	574	581	619	rVB4	8221781	36719621	51.10%	6.566%
9	5.825	644	661	684	rBV	6350745	21956633	30.56%	3.926%
10	6.137	701	714	732	rVB3	333643	1087967	1.51%	0.195%
11	6.325	732	746	763	rBV	3469361	10623103	14.78%	1.899%
12	6.666	783	804	816	rBV	1097701	3499244	4.87%	0.626%
13	6.807	816	828	837	rBV2	705048	2233780	3.11%	0.399%
14	7.089	866	876	890	rBV2	925240	2767803	3.85%	0.495%
15	7.231	890	900	908	rVV	918119	2694888	3.75%	0.482%
16	7.336	908	918	939	rVV	8257701	24690869	34.36%	4.415%
17	8.013	1025	1033	1041	rVB	1066556	2650576	3.69%	0.474%
18	8.172	1052	1060	1063	rBV3	546534	1277432	1.78%	0.228%
19	8.260	1063	1075	1082	rVV2	4475943	17304942	24.08%	3.094%
20	8.336	1082	1088	1099	rVB	3314008	8315143	11.57%	1.487%
21	8.454	1099	1108	1114	rBV	2736015	6233930	8.68%	1.115%
22	8.519	1114	1119	1125	rVB	506355	1029264	1.43%	0.184%
23	8.583	1125	1130	1139	rBV2	409508	1014683	1.41%	0.181%
24	8.725	1145	1154	1159	rBV3	1949188	5169755	7.19%	0.924%
25	8.795	1161	1166	1172	rVV	1921258	4337975	6.04%	0.776%
26	8.860	1172	1177	1186	rVB	1660212	3630073	5.05%	0.649%
27	8.972	1186	1196	1208	rVB3	1298705	3191509	4.44%	0.571%
28	9.107	1208	1219	1226	rBV3	2891624	7546144	10.50%	1.349%
29	9.266	1240	1246	1252	rVB	449689	833367	1.16%	0.149%
30	9.336	1252	1258	1262	rBV	384174	772531	1.08%	0.138%
31	9.389	1262	1267	1286	rVB	612700	1392032	1.94%	0.249%
32	9.607	1286	1304	1319	rBV2	6005504	16744706	23.30%	2.994%
33	9.783	1327	1334	1340	rVB	1117938	2170290	3.02%	0.388%
34	9.942	1354	1361	1363	rBV	464801	907302	1.26%	0.162%
35	10.013	1369	1373	1382	rVB2	796084	1666273	2.32%	0.298%
36	10.107	1382	1389	1393	rBV	1693594	3371610	4.69%	0.603%

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Integrator: RTE

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Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

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Stop Thrs : 0

Peak Location: TOP

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37	10.154	1393	1397	1402	rVV2	1104050	2086720	2.90%	0.373%
38	10.342	1419	1429	1441	rVB3	794658	2005430	2.79%	0.359%
39	10.454	1441	1448	1453	rBV	1043074	2012564	2.80%	0.360%
40	10.572	1461	1468	1474	rBV	2659855	5351173	7.45%	0.957%
41	10.754	1492	1499	1506	rBV3	713003	1671835	2.33%	0.299%
42	10.913	1520	1526	1532	rBV3	385126	815999	1.14%	0.146%
43	10.983	1532	1538	1549	rVV2	1248650	3320601	4.62%	0.594%
44	11.866	1680	1688	1695	rBV	2055679	3829026	5.33%	0.685%
45	11.966	1695	1705	1715	rVB	9034352	15937989	22.18%	2.850%
46	12.077	1715	1724	1735	rBV	3023950	5654192	7.87%	1.011%
47	12.695	1819	1829	1837	rBV	5551783	9284641	12.92%	1.660%
48	12.848	1848	1855	1865	rVB	1423457	2669931	3.72%	0.477%
49	13.036	1879	1887	1895	rBV	15114327	25333091	35.25%	4.530%
50	13.130	1895	1903	1908	rVV	3939139	6792465	9.45%	1.215%
51	13.177	1908	1911	1918	rVB	1327187	1973004	2.75%	0.353%
52	13.336	1931	1938	1948	rVB	2395816	4024100	5.60%	0.720%
53	13.483	1956	1963	1970	rVB	6722544	10853819	15.10%	1.941%
54	13.613	1977	1985	1991	rBV3	1430650	3559838	4.95%	0.637%
55	13.795	2010	2016	2017	rBV	1946908	2706271	3.77%	0.484%
56	13.818	2018	2020	2028	rVB	2712882	4061058	5.65%	0.726%
57	13.954	2036	2043	2046	rBV	4398824	7107192	9.89%	1.271%
58	13.995	2046	2050	2055	rVV	6361801	10369827	14.43%	1.854%
59	14.118	2066	2071	2076	rVB	938533	1455796	2.03%	0.260%
60	14.183	2076	2082	2088	rBV	1684720	2712208	3.77%	0.485%
61	14.248	2088	2093	2096	rBV	1214867	2039240	2.84%	0.365%
62	14.330	2102	2107	2114	rVV	10775941	17042058	23.72%	3.047%
63	14.401	2114	2119	2122	rVV	1928562	3240488	4.51%	0.579%
64	14.448	2122	2127	2134	rVB2	7072247	12338328	17.17%	2.206%
65	14.577	2143	2149	2154	rBV2	627684	1016281	1.41%	0.182%
66	14.654	2154	2162	2167	rBV	6533890	11041191	15.37%	1.974%
67	14.695	2167	2169	2173	rVV	1271503	1660179	2.31%	0.297%
68	14.736	2173	2176	2180	rVV	667620	913301	1.27%	0.163%
69	14.930	2204	2209	2214	rBV	5344650	8760722	12.19%	1.566%
70	15.048	2220	2229	2235	rBV2	8332025	18233120	25.37%	3.260%
71	15.095	2235	2237	2249	rVV3	709391	1521988	2.12%	0.272%
72	15.201	2249	2255	2261	rVV3	1132024	1991091	2.77%	0.356%
73	15.301	2261	2272	2276	rVV2	2627092	6659293	9.27%	1.191%
74	15.342	2276	2279	2283	rVV	1219621	1953440	2.72%	0.349%
75	15.412	2283	2291	2300	rVB4	2376904	5577269	7.76%	0.997%
76	15.589	2314	2321	2327	rVB	5468914	9402291	13.08%	1.681%

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Integrator: RTE
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Stop Thrs : 0 Peak Location: TOP

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Title : SW846 8260

77	15.677	2332	2336	2340	rVV	633435	999643	1.39%	0.179%
78	15.724	2340	2344	2356	rVB	680886	1370212	1.91%	0.245%
79	15.836	2356	2363	2370	rBV	1065397	1790855	2.49%	0.320%
80	15.989	2380	2389	2401	rBV3	540654	1658674	2.31%	0.297%
81	16.154	2410	2417	2430	rVB3	402060	942646	1.31%	0.169%
82	16.495	2468	2475	2485	rVB	2483622	4650874	6.47%	0.832%
83	16.654	2495	2502	2517	rVB	1187064	2374905	3.31%	0.425%

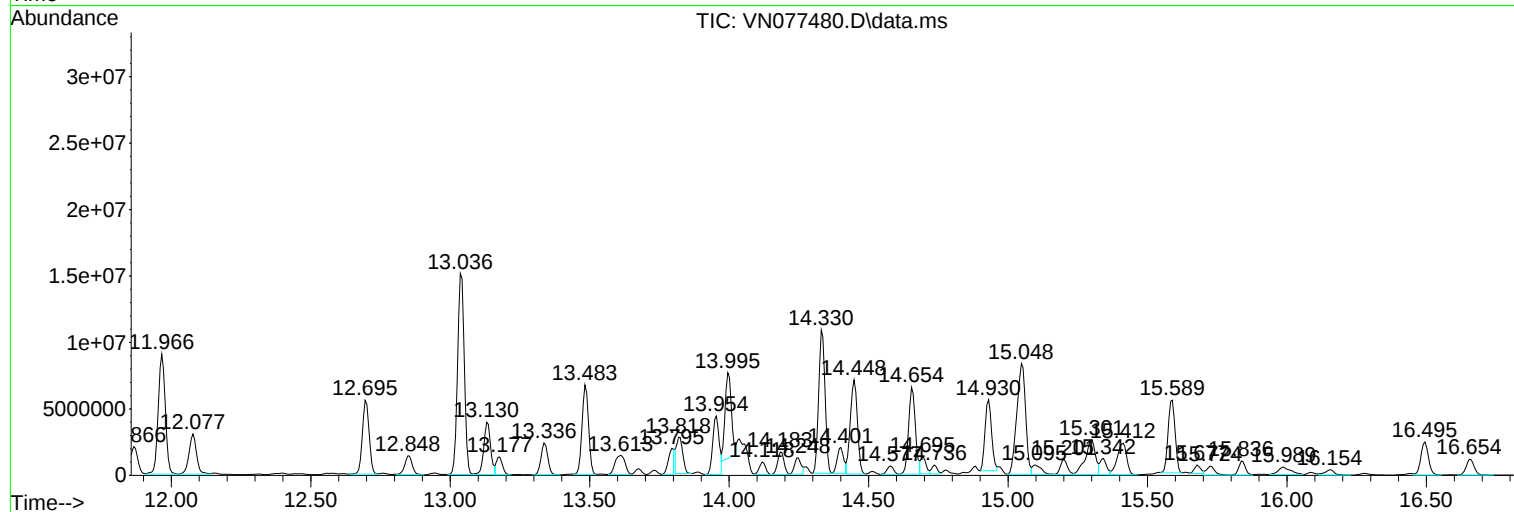
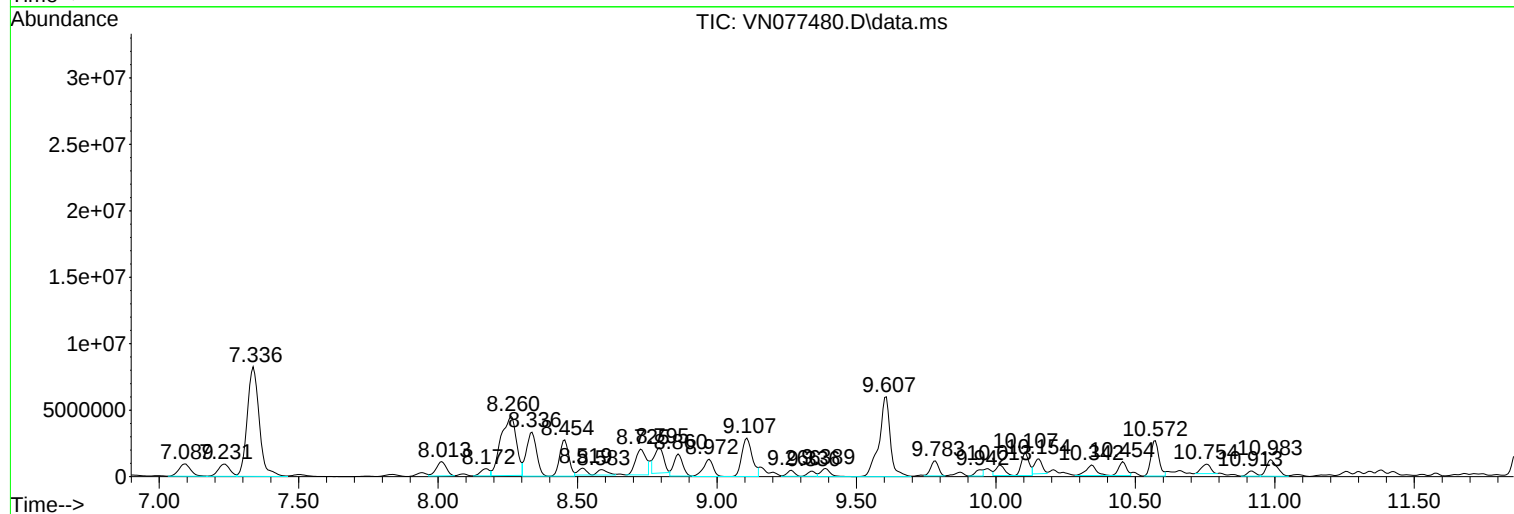
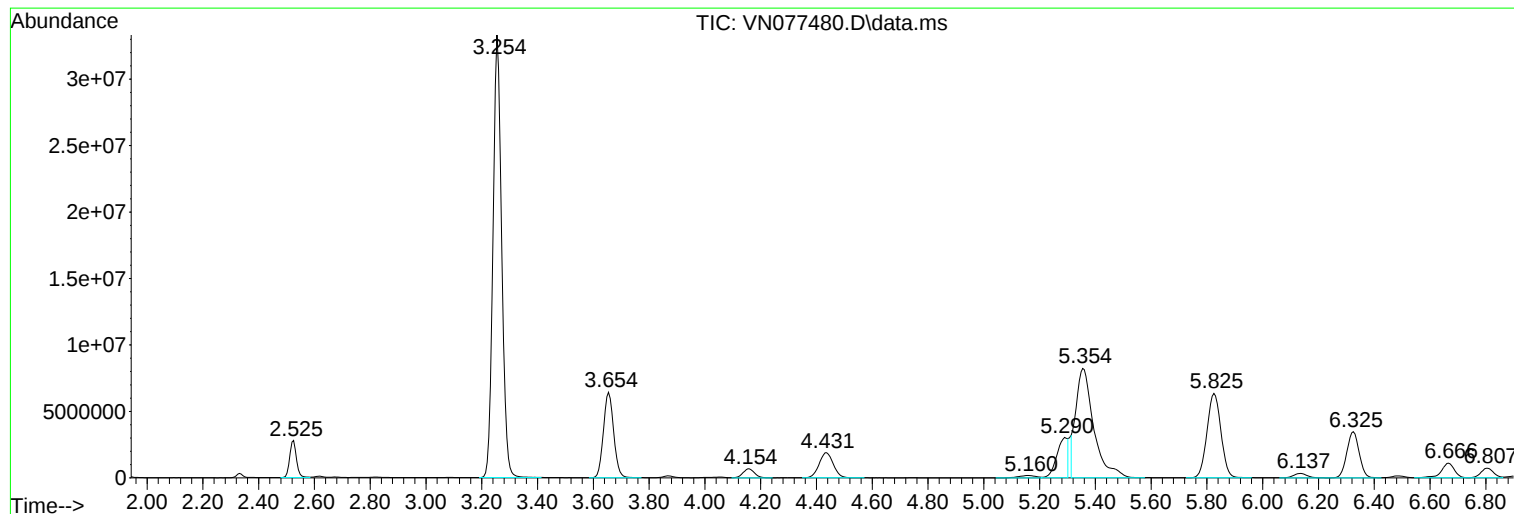
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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



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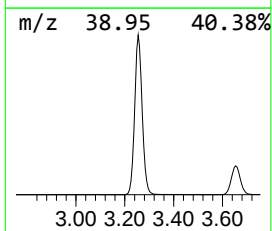
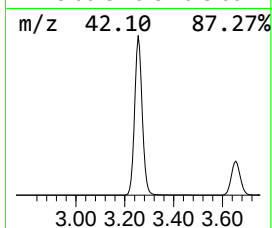
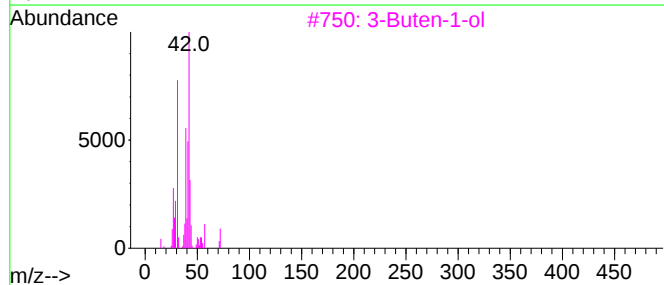
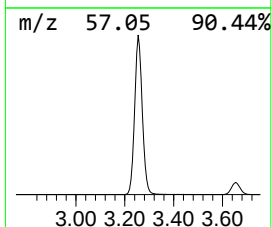
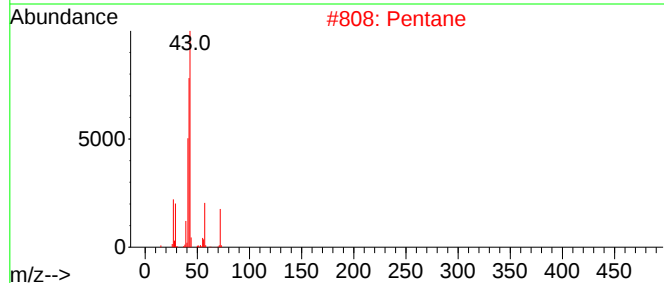
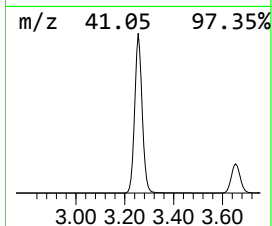
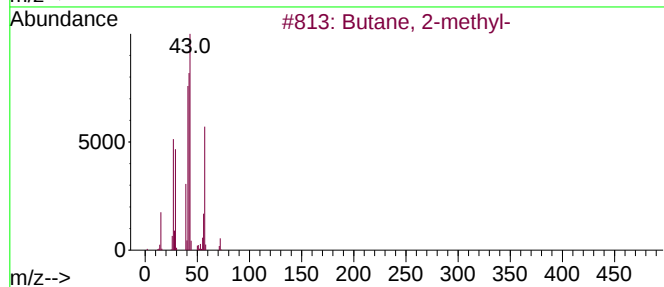
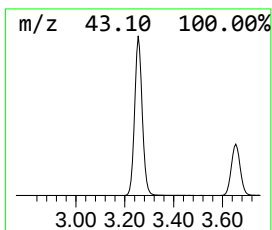
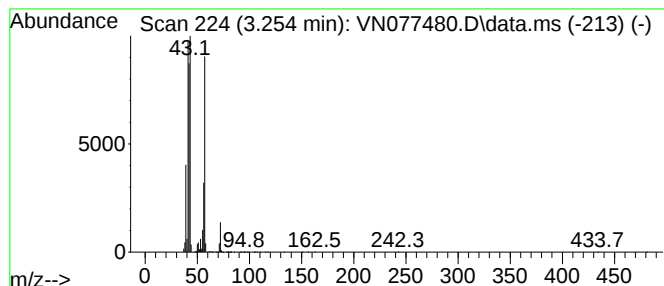
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.254	207.62 ug/l	71857200	Pentafluorobenzene	8.231

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-			72	C5H12	000078-78-4	64
2	Pentane			72	C5H12	000109-66-0	37
3	3-Buten-1-ol			72	C4H8O	000627-27-0	33
4	2-Butene-1,4-diol, (Z)-			88	C4H8O2	006117-80-2	28
5	1-Propene, 2-methyl-			56	C4H8	000115-11-7	27



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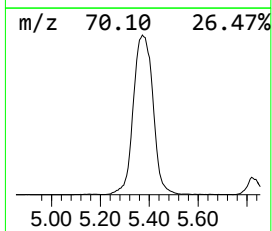
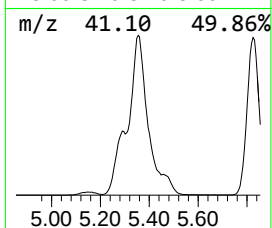
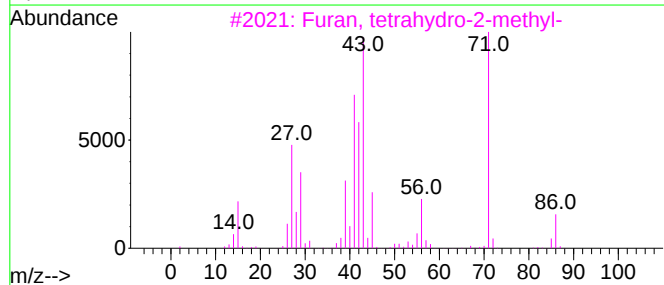
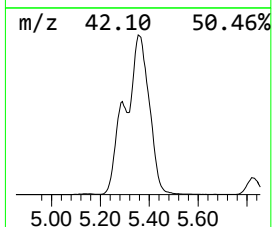
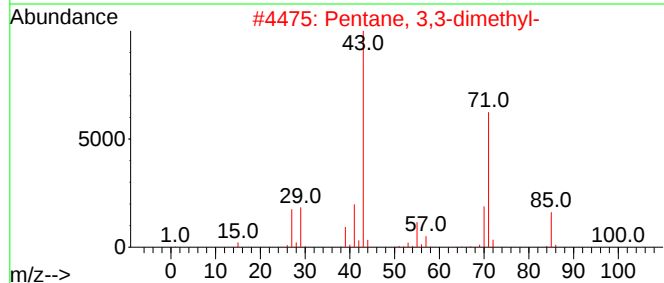
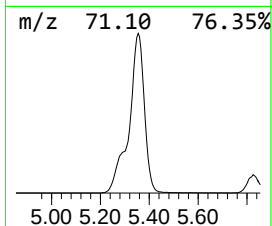
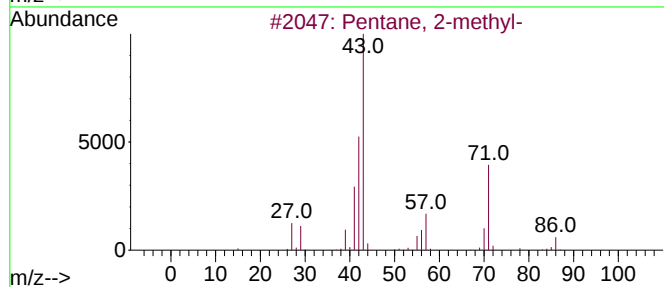
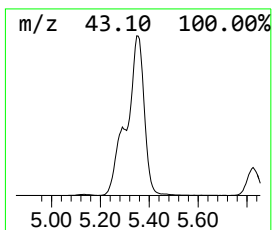
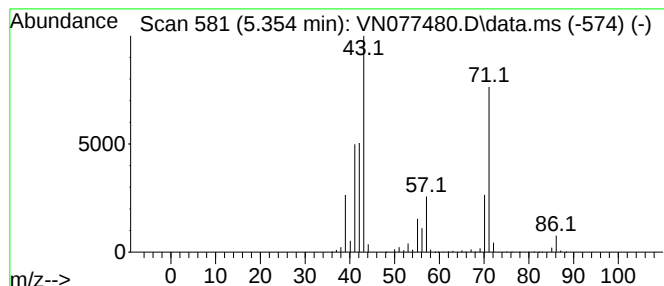
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.354	106.10 ug/l	36719600	Pentafluorobenzene	8.231

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	78
2			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	45
3			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	38
4			Pentane, 3-bromo-	150	C5H11Br	001809-10-5	37
5			3-Penten-2-ol	86	C5H10O	001569-50-2	37



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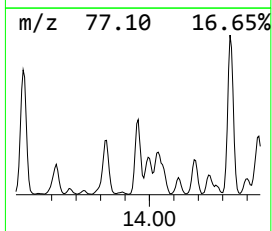
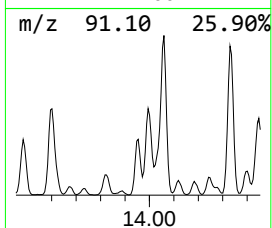
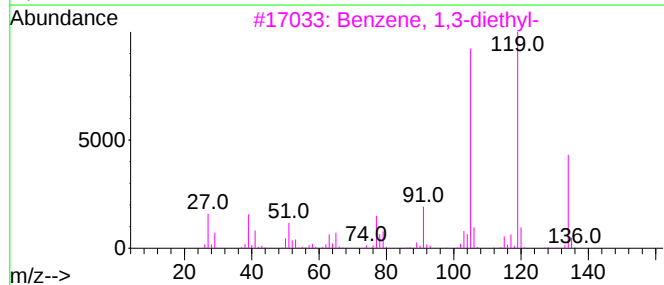
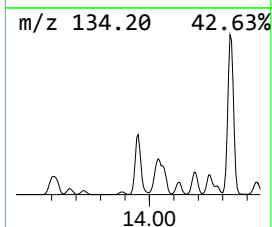
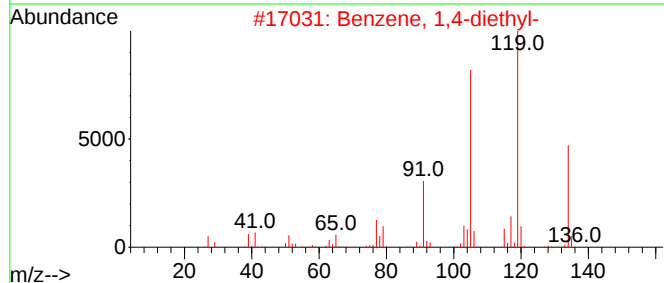
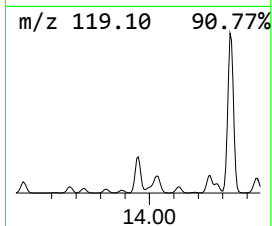
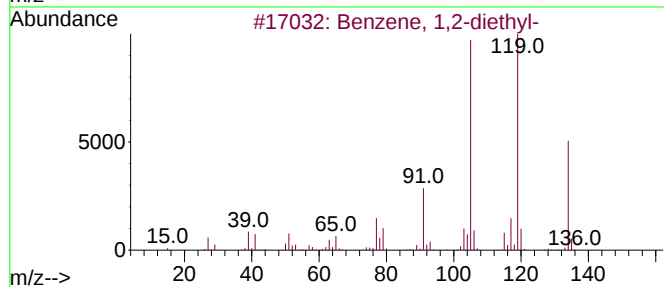
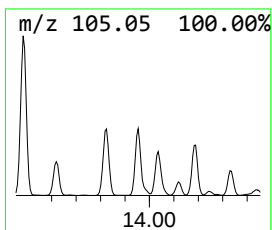
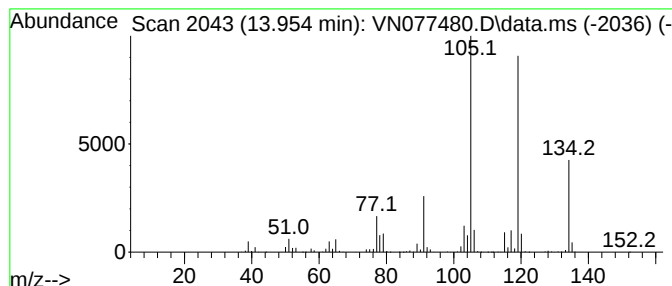
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.954	131.31 ug/l	7107190	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			o-Cymene	134	C10H14	000527-84-4	81



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ClientSampleId :
MW-5

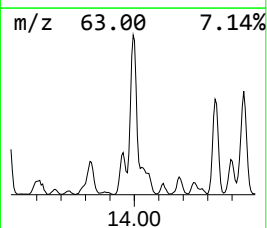
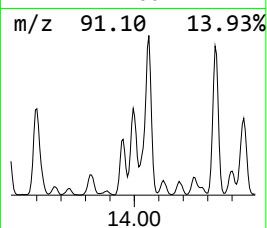
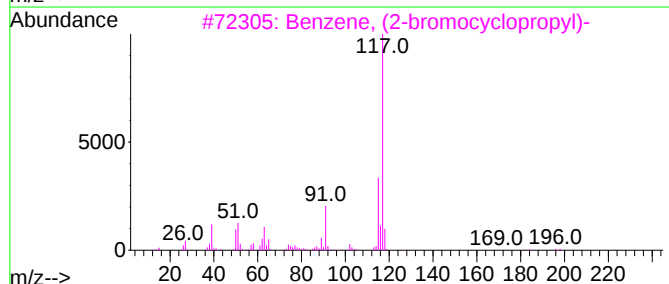
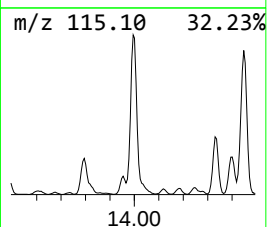
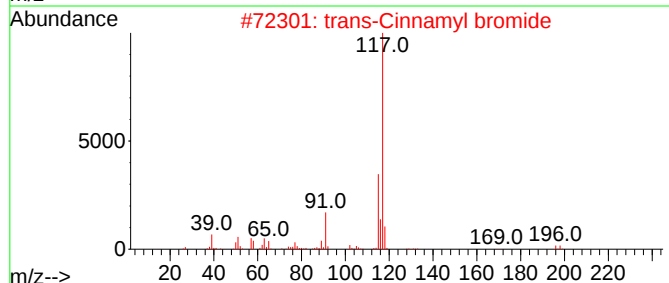
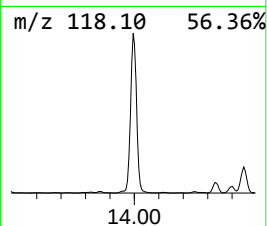
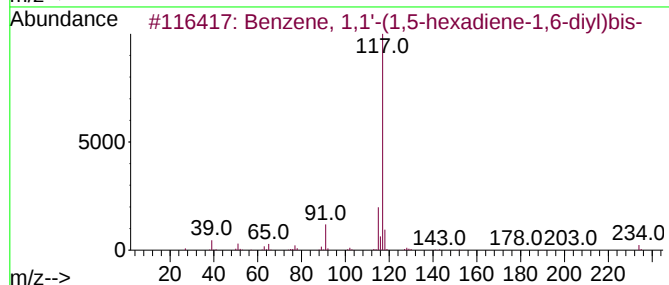
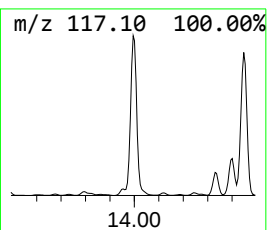
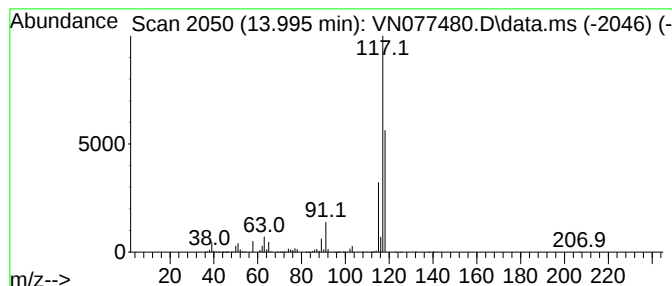
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.995	191.59 ug/l	10369800	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	45
3			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
4			Indole	117	C8H7N	000120-72-9	43
5			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

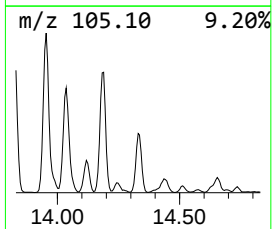
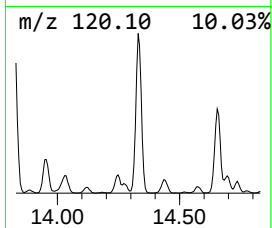
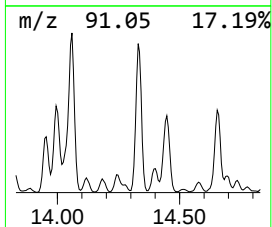
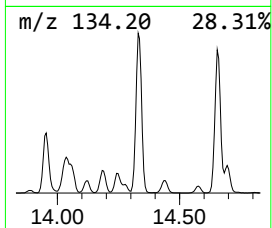
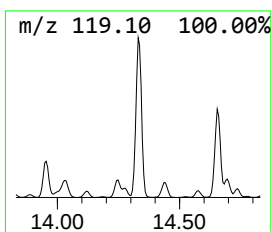
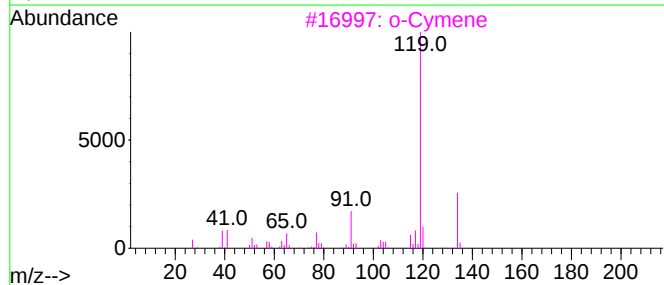
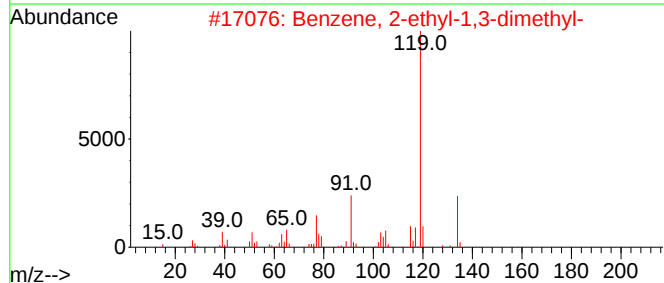
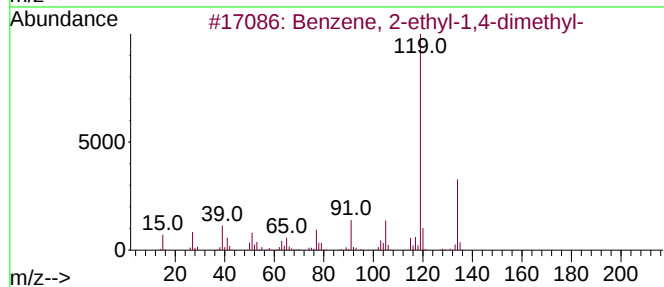
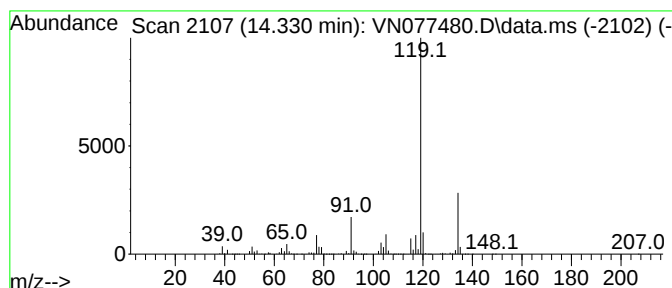
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.330	314.86 ug/l	17042100	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

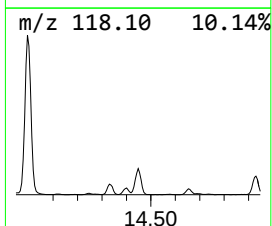
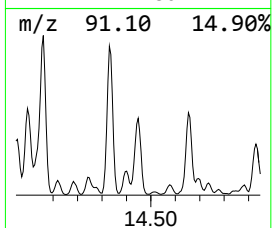
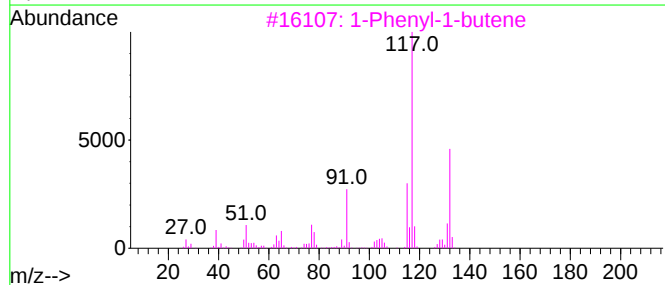
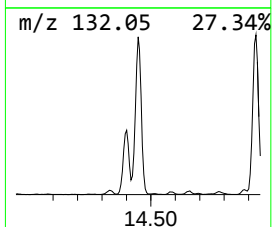
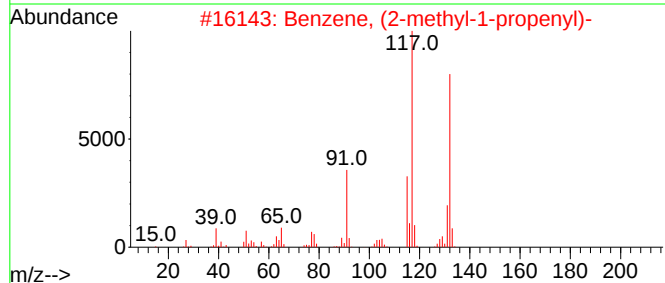
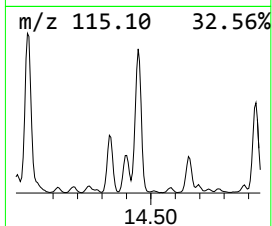
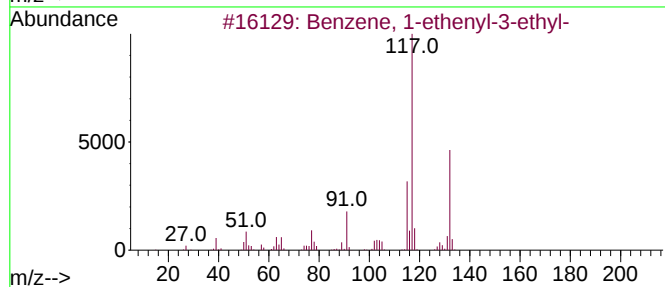
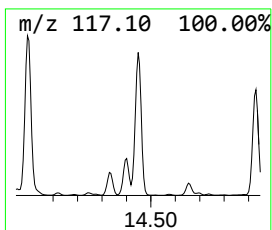
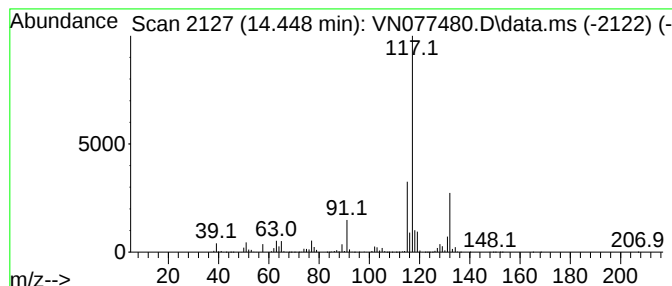
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.448	227.96 ug/l	12338300	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
4			Indan, 1-methyl-	132	C10H12	000767-58-8	83
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

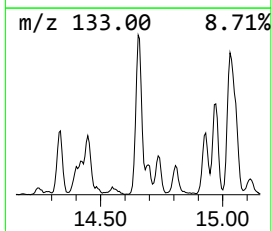
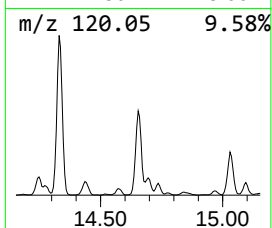
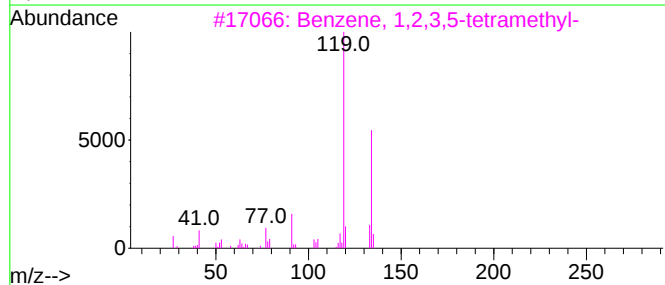
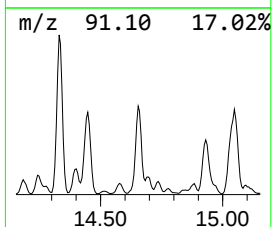
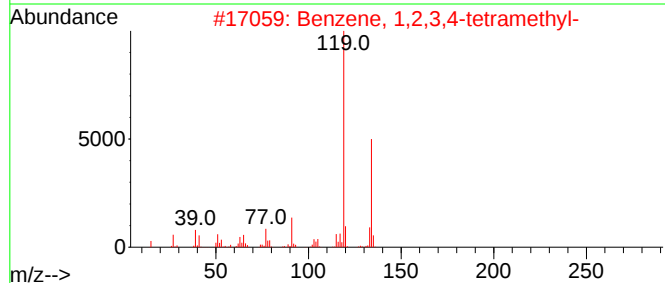
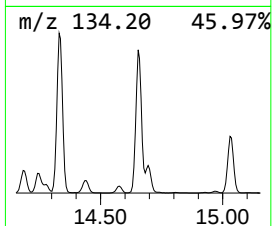
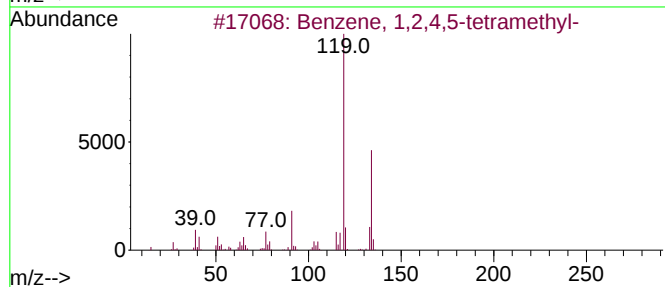
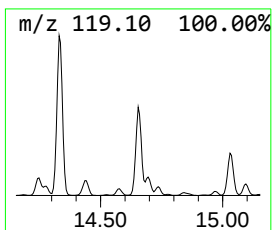
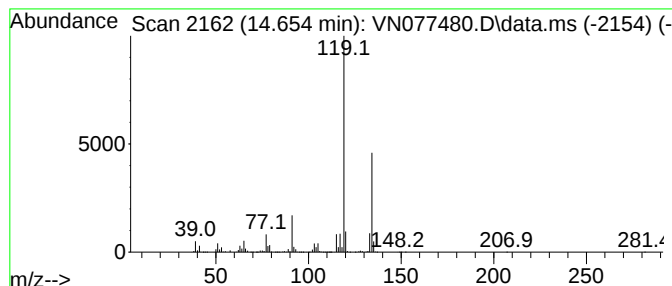
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.654	203.99 ug/l	11041200	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	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\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

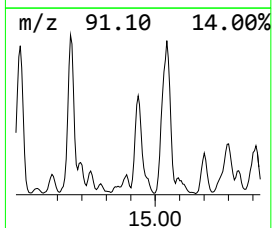
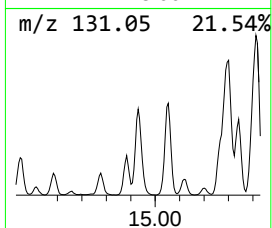
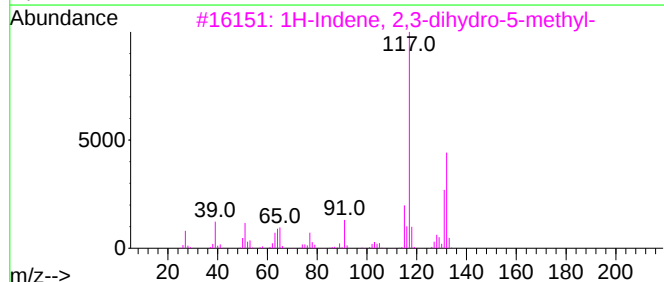
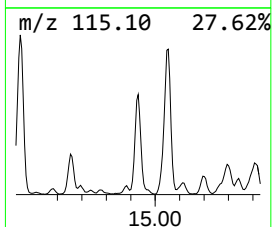
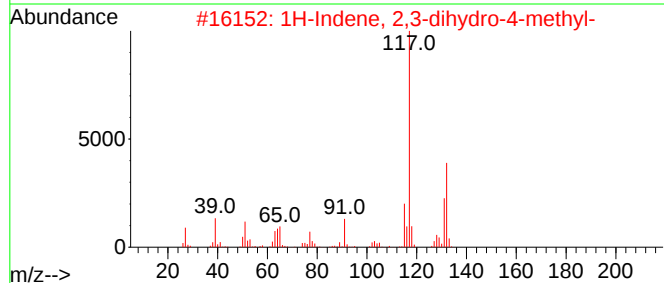
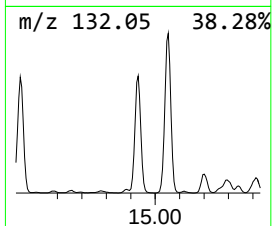
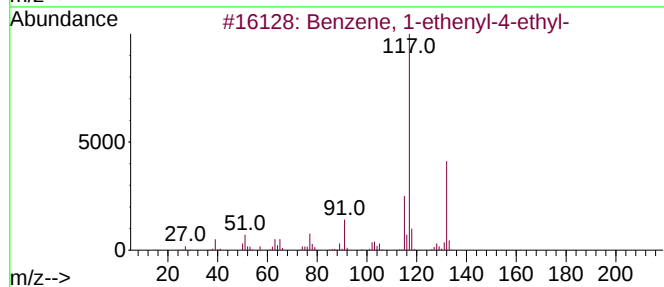
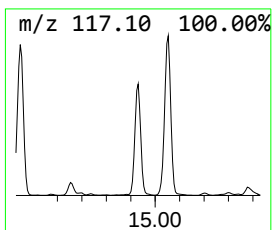
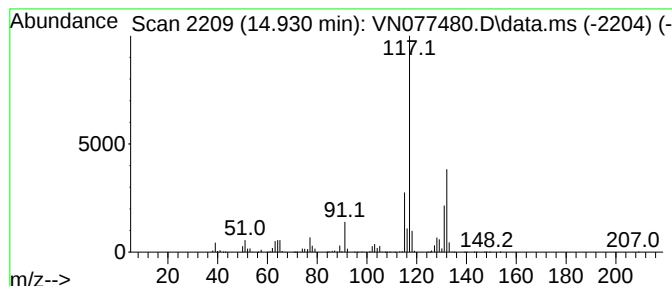
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.930	161.86 ug/l	8760720	1,4-Dichlorobenzene-d4	13.795

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	94
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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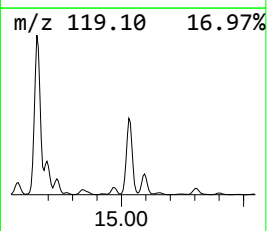
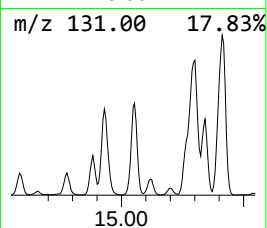
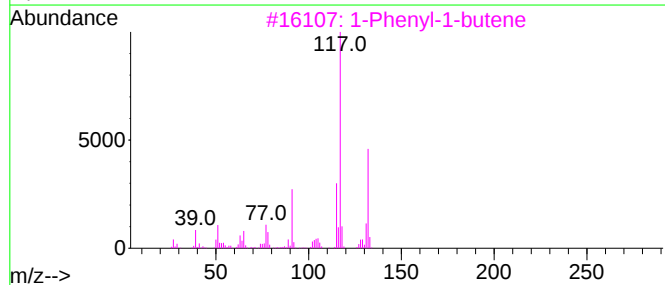
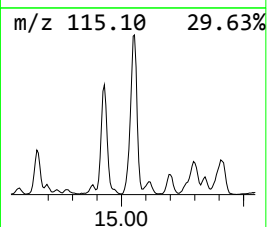
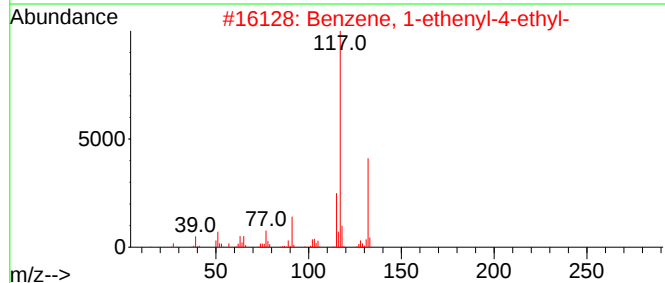
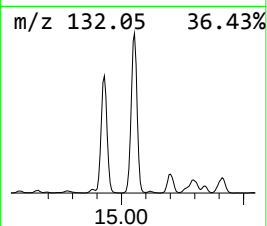
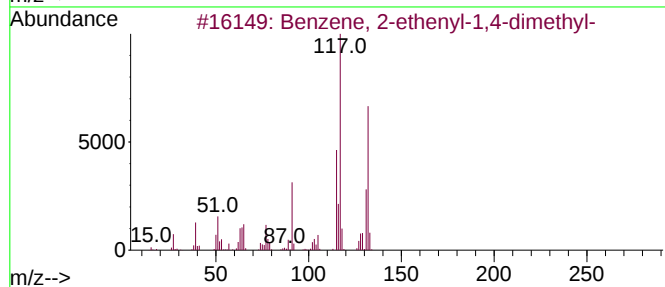
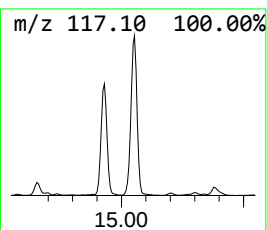
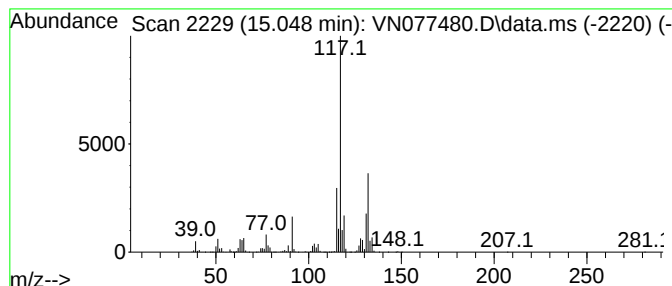
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.048	336.87 ug/l	18233100	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		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	93
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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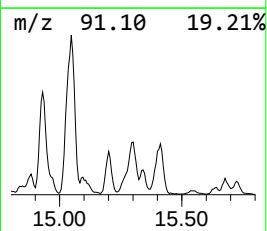
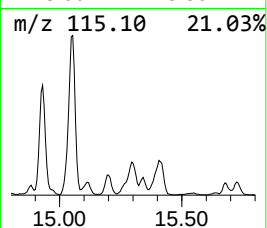
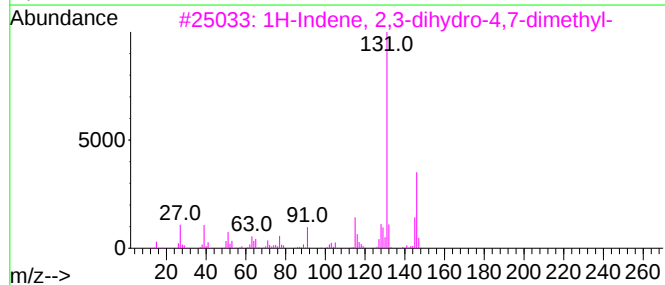
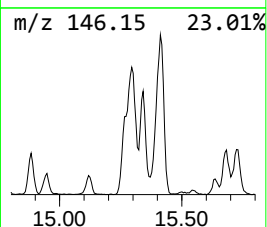
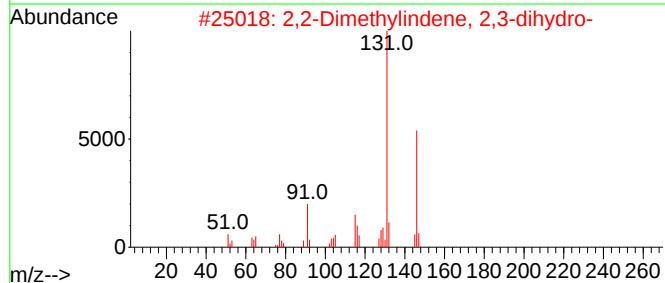
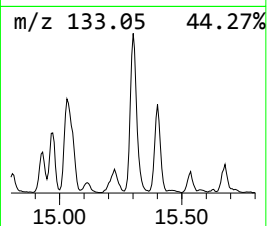
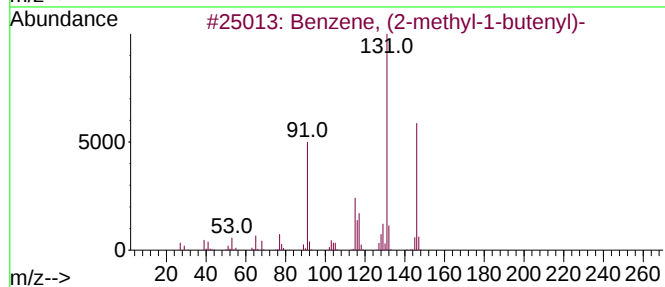
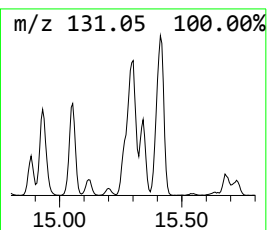
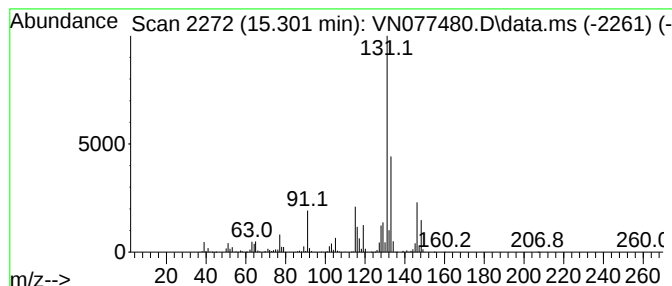
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (2-methyl-1-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.301	123.03 ug/l	6659290	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
2			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	86
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
5			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077480.D
Acq On : 25 Apr 2023 19:34
Operator : JC\MD
Sample : 02460-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.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
Butane, 2-methyl-	3.254	207.6	ug/l	71857200	1	8.231	17304900	50.0
Pentane, 2-methyl-	5.354	106.1	ug/l	36719600	1	8.231	17304900	50.0
Benzene, 1,2-di...	13.954	131.3	ug/l	7107190	4	13.795	2706270	50.0
Benzene, 1,1'-(...	13.995	191.6	ug/l	10369800	4	13.795	2706270	50.0
Benzene, 2-ethy...	14.330	314.9	ug/l	17042100	4	13.795	2706270	50.0
Benzene, 1-ethe...	14.448	228.0	ug/l	12338300	4	13.795	2706270	50.0
Benzene, 1,2,4,...	14.654	204.0	ug/l	11041200	4	13.795	2706270	50.0
Benzene, 1-ethe...	14.930	161.9	ug/l	8760720	4	13.795	2706270	50.0
Benzene, 2-ethe...	15.048	336.9	ug/l	18233100	4	13.795	2706270	50.0
Benzene, (2-met...	15.301	123.0	ug/l	6659290	4	13.795	2706270	50.0