

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
 Data File : VN077478.D
 Acq On : 25 Apr 2023 18:45
 Operator : JC\MD
 Sample : 02460-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-4

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: VN077478.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.519	88	99	109	rBV	1161698	1929929	4.75%	0.473%
2	3.254	213	224	243	rBV	18258576	40641714	100.00%	9.960%
3	3.654	281	292	307	rBV	2171871	5506888	13.55%	1.350%
4	4.160	367	378	392	rVB2	394022	1102905	2.71%	0.270%
5	4.437	409	425	446	rBV2	1173079	4091802	10.07%	1.003%
6	5.160	531	548	556	rBV6	104278	438075	1.08%	0.107%
7	5.290	556	570	573	rVV2	2487650	7775079	19.13%	1.905%
8	5.354	574	581	618	rVB4	5519667	23626897	58.13%	5.790%
9	5.825	645	661	680	rBV	4436714	15306795	37.66%	3.751%
10	6.137	699	714	728	rVB2	242461	803154	1.98%	0.197%
11	6.325	733	746	759	rBV	1298271	3925878	9.66%	0.962%
12	6.666	785	804	815	rBV	721186	2240201	5.51%	0.549%
13	6.807	815	828	838	rBV2	596026	1865823	4.59%	0.457%
14	7.090	865	876	889	rBV4	771318	2303758	5.67%	0.565%
15	7.237	889	901	909	rVV	1977066	5849644	14.39%	1.434%
16	7.337	909	918	939	rVV	5043031	15456443	38.03%	3.788%
17	7.837	994	1003	1012	rVB4	141862	432371	1.06%	0.106%
18	7.942	1012	1021	1025	rBV3	284128	692569	1.70%	0.170%
19	8.013	1025	1033	1041	rVB2	787809	1953230	4.81%	0.479%
20	8.172	1051	1060	1063	rBV2	507129	1197916	2.95%	0.294%
21	8.231	1063	1070	1073	rBV2	2576203	5751385	14.15%	1.409%
22	8.337	1081	1088	1099	rVB	5580544	14209129	34.96%	3.482%
23	8.454	1099	1108	1115	rBV	2599631	5944458	14.63%	1.457%
24	8.513	1115	1118	1125	rVB	352085	686891	1.69%	0.168%
25	8.584	1125	1130	1139	rBV2	383463	947866	2.33%	0.232%
26	8.731	1145	1155	1157	rBV4	1837391	4368882	10.75%	1.071%
27	8.772	1157	1162	1172	rVV3	2685486	8814939	21.69%	2.160%
28	8.860	1172	1177	1186	rVB	1331771	3014482	7.42%	0.739%
29	8.972	1186	1196	1205	rVB2	861368	2315828	5.70%	0.568%
30	9.107	1209	1219	1225	rBV3	2901745	7477802	18.40%	1.833%
31	9.266	1240	1246	1252	rVB	501922	926640	2.28%	0.227%
32	9.336	1252	1258	1262	rBV	405755	823007	2.03%	0.202%
33	9.389	1262	1267	1285	rVB	749354	1689077	4.16%	0.414%
34	9.601	1287	1303	1319	rBV3	4414580	13758559	33.85%	3.372%
35	9.731	1319	1325	1328	rBV	229474	439522	1.08%	0.108%
36	9.783	1328	1334	1340	rVV	945969	1858257	4.57%	0.455%

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37	9.872	1340	1349	1355	rVB2	391485	893896	2.20%	0.219%
38	10.019	1368	1374	1382	rVB	2067703	4355311	10.72%	1.067%
39	10.107	1382	1389	1392	rBV2	1492089	2945614	7.25%	0.722%
40	10.148	1392	1396	1402	rVV	3105765	6170300	15.18%	1.512%
41	10.207	1402	1406	1409	rVV3	401664	716997	1.76%	0.176%
42	10.242	1409	1412	1420	rVB4	349879	824140	2.03%	0.202%
43	10.342	1420	1429	1441	rBV3	851677	2129431	5.24%	0.522%
44	10.454	1441	1448	1452	rBV	939882	1821036	4.48%	0.446%
45	10.495	1452	1455	1461	rVB	559542	912606	2.25%	0.224%
46	10.572	1461	1468	1474	rBV2	2707315	5387429	13.26%	1.320%
47	10.660	1479	1483	1487	rVV4	294599	628042	1.55%	0.154%
48	10.760	1491	1500	1505	rVV4	908067	2522179	6.21%	0.618%
49	10.919	1521	1527	1532	rBV4	467090	1006367	2.48%	0.247%
50	10.983	1532	1538	1549	rVB3	1673869	4299853	10.58%	1.054%
51	11.078	1549	1554	1561	rVB4	201622	496854	1.22%	0.122%
52	11.254	1577	1584	1588	rBV3	420105	896776	2.21%	0.220%
53	11.301	1588	1592	1595	rVV2	321466	497657	1.22%	0.122%
54	11.342	1595	1599	1602	rVV2	314923	533787	1.31%	0.131%
55	11.378	1602	1605	1610	rVV2	396771	737125	1.81%	0.181%
56	11.425	1610	1613	1619	rVB3	321144	489355	1.20%	0.120%
57	11.866	1681	1688	1694	rBV	1898651	3499318	8.61%	0.858%
58	11.966	1699	1705	1715	rVB	3240209	5520931	13.58%	1.353%
59	12.077	1715	1724	1734	rBV	1706325	3626637	8.92%	0.889%
60	12.695	1820	1829	1836	rBV	3442307	5968476	14.69%	1.463%
61	12.848	1845	1855	1865	rVB	1356882	2583861	6.36%	0.633%
62	13.036	1878	1887	1894	rVV	8588466	13903142	34.21%	3.407%
63	13.101	1894	1898	1899	rVV3	393070	551394	1.36%	0.135%
64	13.130	1899	1903	1907	rVV	1845292	3069859	7.55%	0.752%
65	13.177	1907	1911	1918	rVB	902295	1540343	3.79%	0.377%
66	13.336	1930	1938	1947	rBV	2229442	3788292	9.32%	0.928%
67	13.483	1956	1963	1970	rVB	2759018	4342729	10.69%	1.064%
68	13.613	1977	1985	1991	rBV3	1208474	3043838	7.49%	0.746%
69	13.671	1991	1995	2001	rVB	348877	538437	1.32%	0.132%
70	13.795	2009	2016	2019	rBV	1681166	3168708	7.80%	0.777%
71	13.954	2037	2043	2046	rBV	3575216	5557378	13.67%	1.362%
72	13.995	2046	2050	2054	rVV	2563439	3617918	8.90%	0.887%
73	14.036	2054	2057	2059	rBV2	641500	800823	1.97%	0.196%
74	14.118	2066	2071	2077	rVB	824100	1258112	3.10%	0.308%
75	14.183	2077	2082	2088	rBV	1596770	2572067	6.33%	0.630%
76	14.248	2088	2093	2101	rVV	728628	1351074	3.32%	0.331%

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77	14.330	2101	2107	2114	rVV	9108327	14306103	35.20%	3.506%
78	14.401	2114	2119	2122	rVV	1200959	2144414	5.28%	0.526%
79	14.448	2122	2127	2134	rVV2	5984350	10247070	25.21%	2.511%
80	14.577	2144	2149	2154	rBV2	344092	547067	1.35%	0.134%
81	14.654	2154	2162	2167	rBV	5643310	9395895	23.12%	2.303%
82	14.695	2167	2169	2173	rVV	896010	1173610	2.89%	0.288%
83	14.736	2173	2176	2180	rVV	777513	1090157	2.68%	0.267%
84	14.777	2180	2183	2190	rVB2	323740	510371	1.26%	0.125%
85	14.883	2196	2201	2204	rBV2	486791	712397	1.75%	0.175%
86	14.930	2204	2209	2214	rBV	3975259	6566592	16.16%	1.609%
87	15.048	2220	2229	2235	rBV3	5676646	12820930	31.55%	3.142%
88	15.095	2235	2237	2249	rVV3	720104	1428901	3.52%	0.350%
89	15.201	2249	2255	2261	rVV2	732866	1351211	3.32%	0.331%
90	15.301	2261	2272	2276	rVV2	2463442	6265732	15.42%	1.535%
91	15.342	2276	2279	2283	rVV	1065333	1747020	4.30%	0.428%
92	15.413	2283	2291	2301	rVB3	2182159	5107740	12.57%	1.252%
93	15.583	2315	2320	2326	rVB	2533845	4379056	10.77%	1.073%
94	15.677	2332	2336	2340	rBV	498606	791284	1.95%	0.194%
95	15.724	2340	2344	2357	rVB	580603	1174832	2.89%	0.288%
96	15.842	2357	2364	2370	rBV	916442	1583111	3.90%	0.388%
97	15.989	2380	2389	2401	rVV4	440499	1434446	3.53%	0.352%
98	16.154	2410	2417	2429	rVB2	332287	866117	2.13%	0.212%
99	16.495	2469	2475	2486	rVB	1180036	2212026	5.44%	0.542%
100	16.660	2495	2503	2511	rVB	724726	1472177	3.62%	0.361%

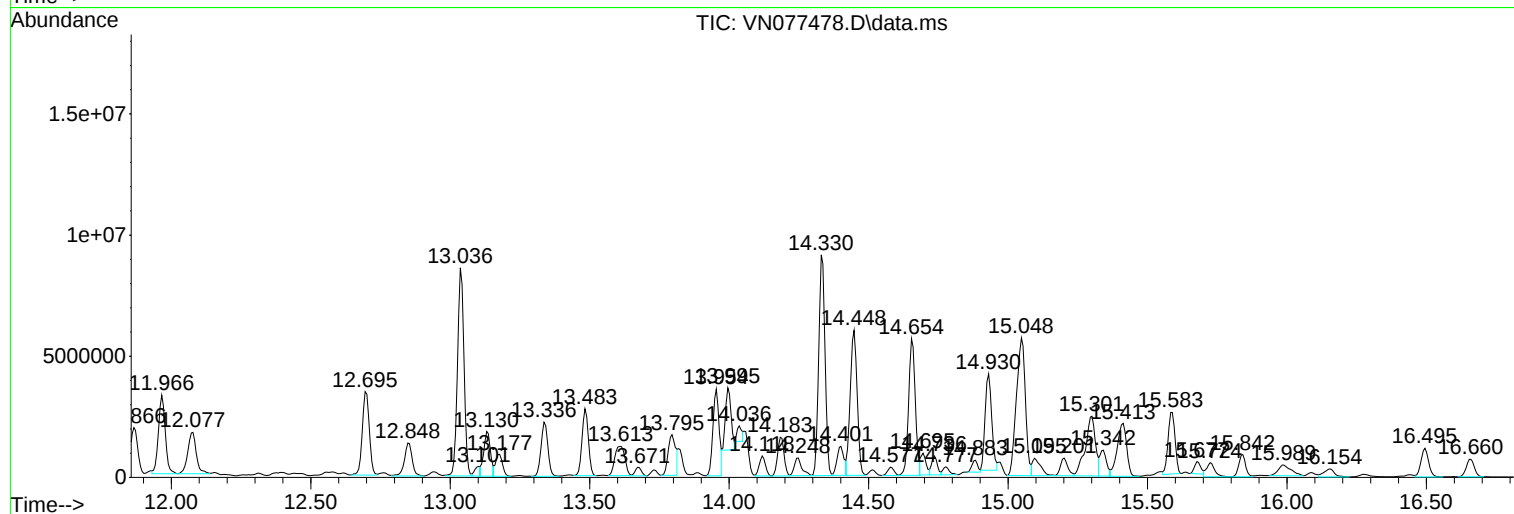
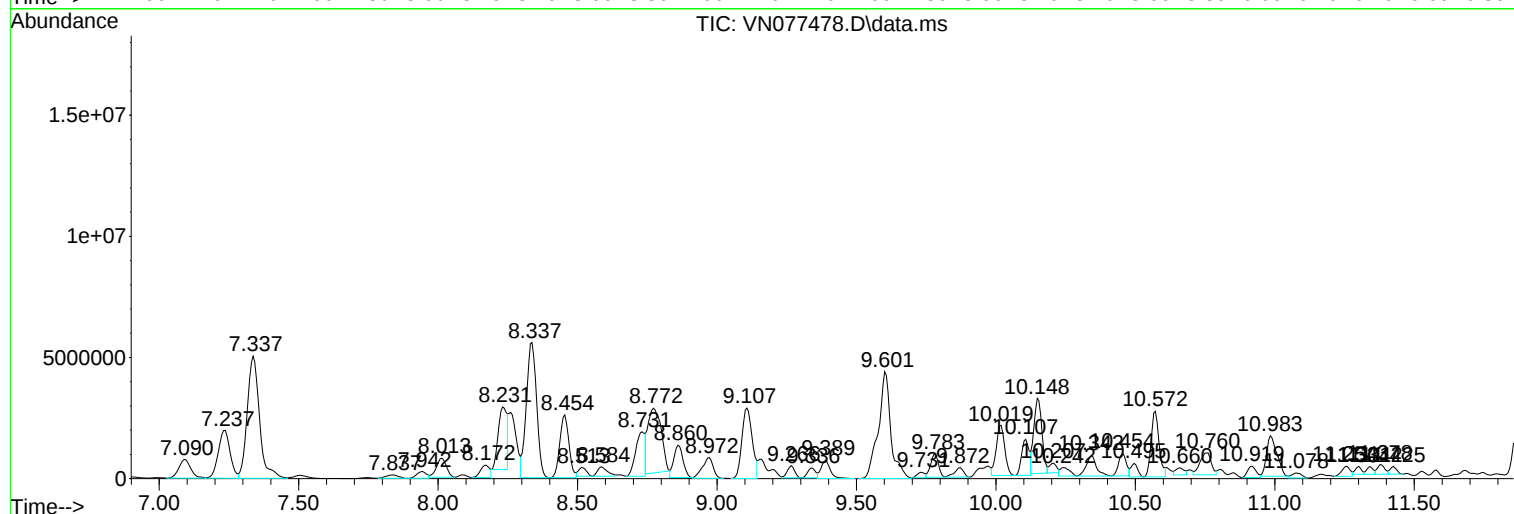
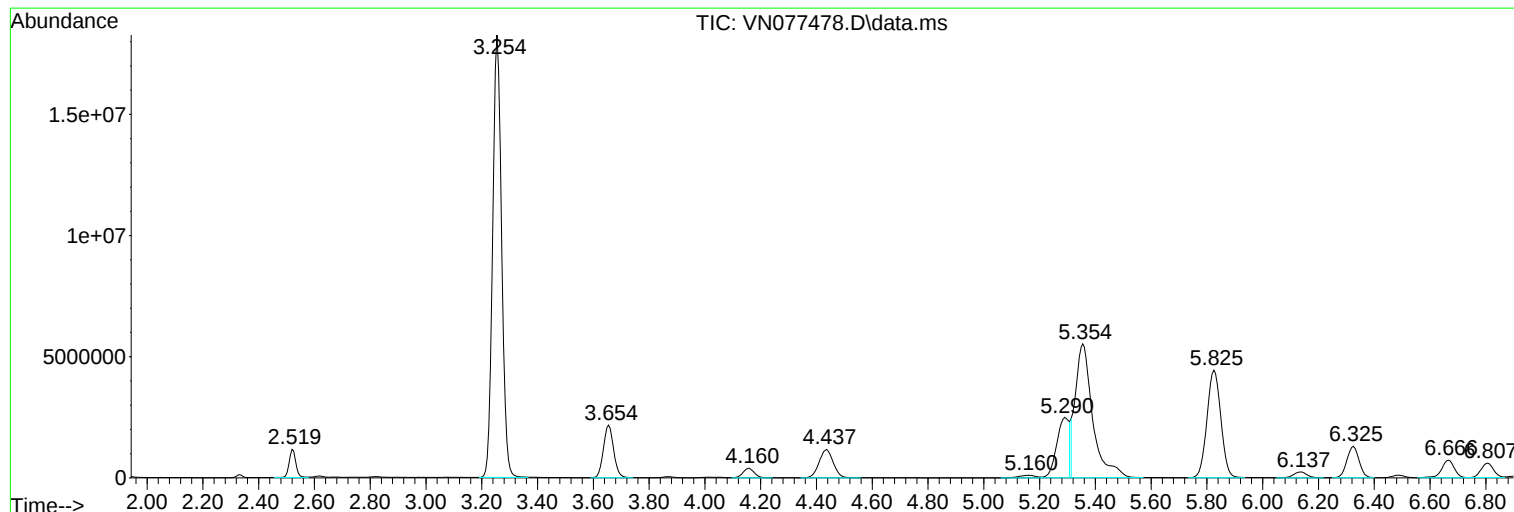
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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



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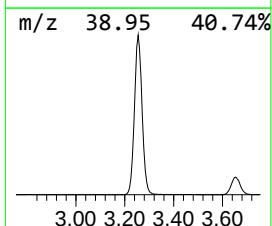
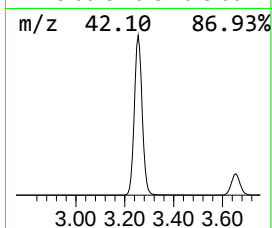
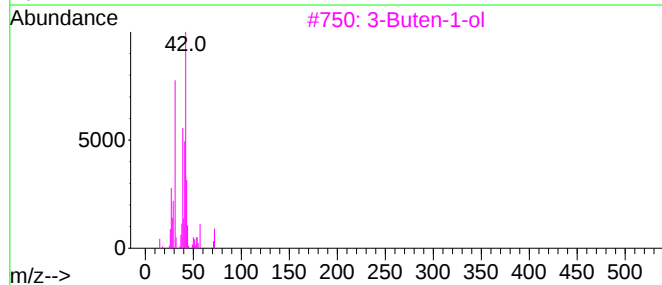
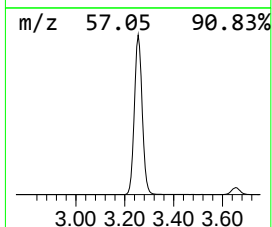
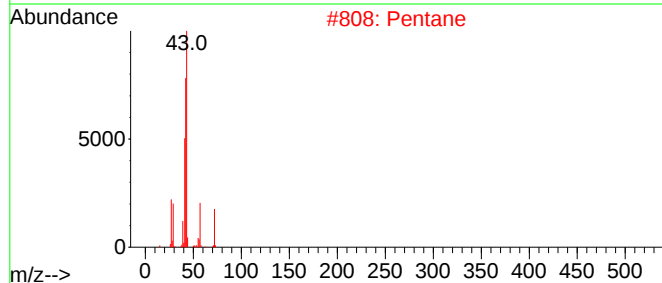
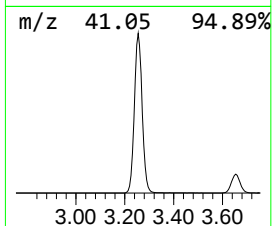
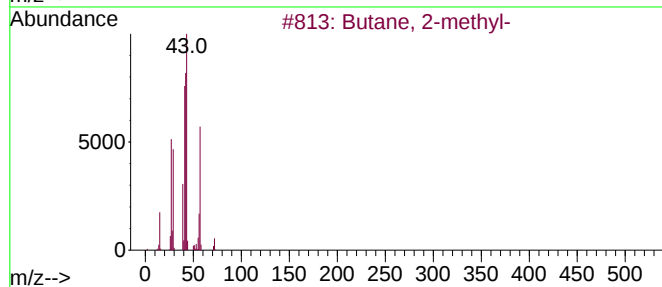
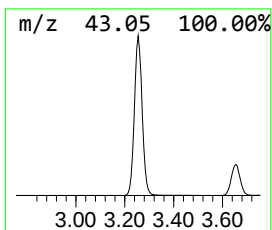
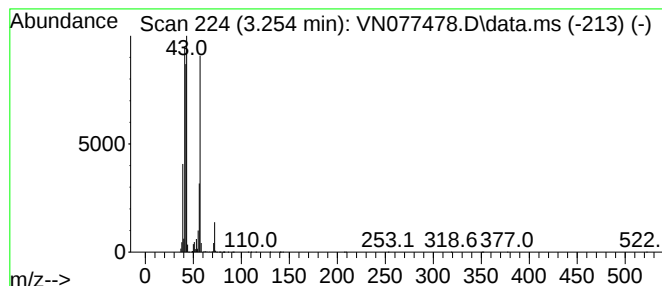
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.254	353.32 ug/l	40641700	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-Butene			56	C4H8	000106-98-9	27



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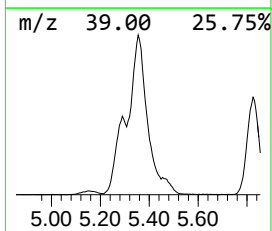
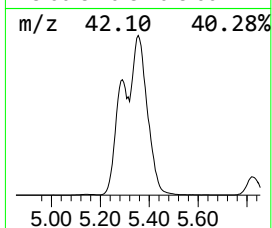
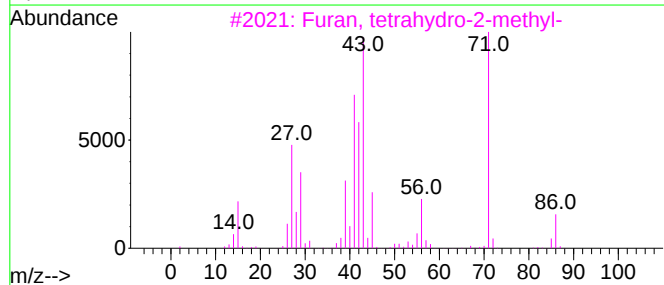
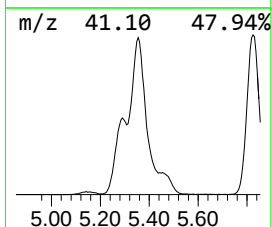
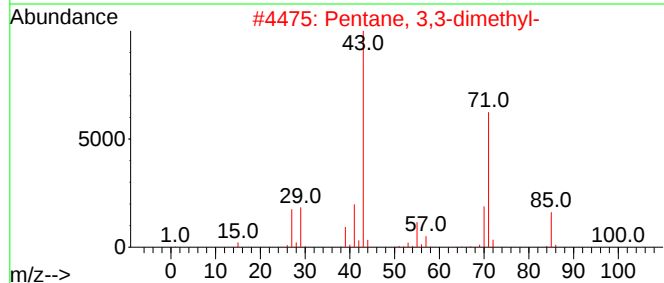
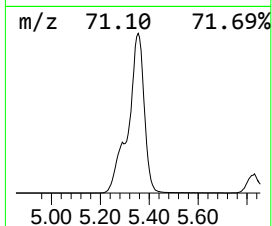
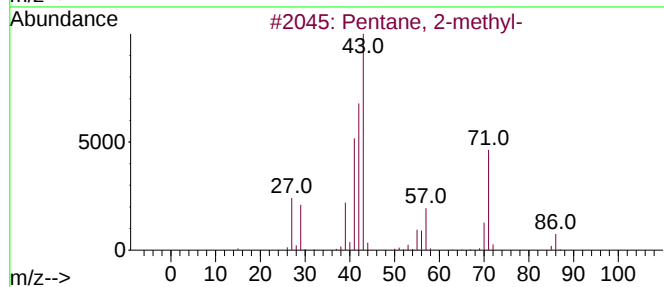
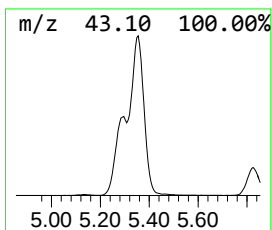
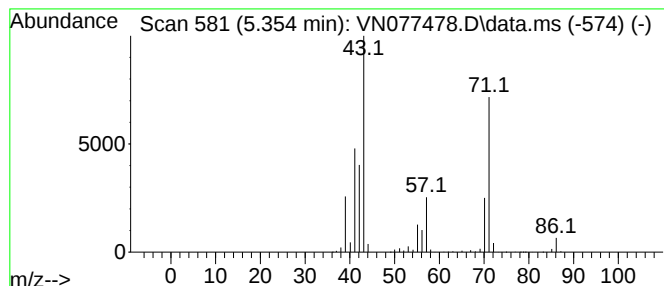
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.354	205.40 ug/l	23626900	Pentafluorobenzene	8.231

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	86
2			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	50
3			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	45
4			Butane, 2-methyl-	72	C5H12	000078-78-4	38
5			3-Penten-2-ol	86	C5H10O	001569-50-2	35



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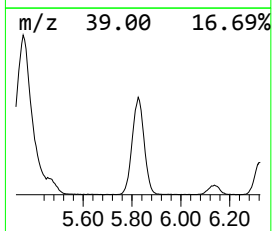
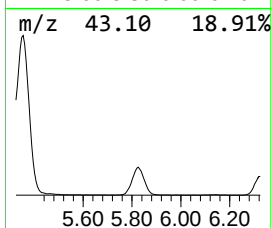
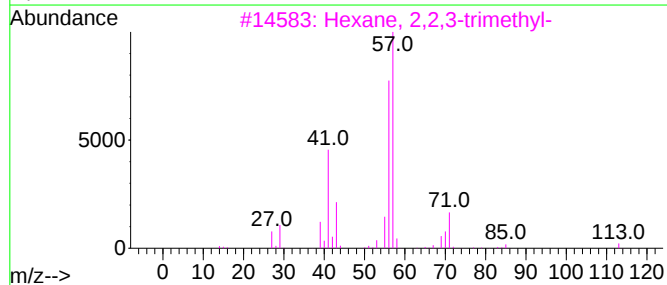
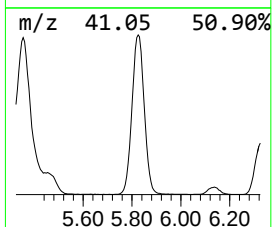
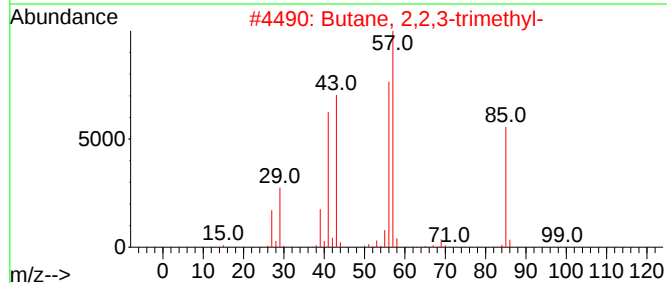
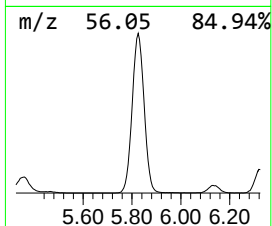
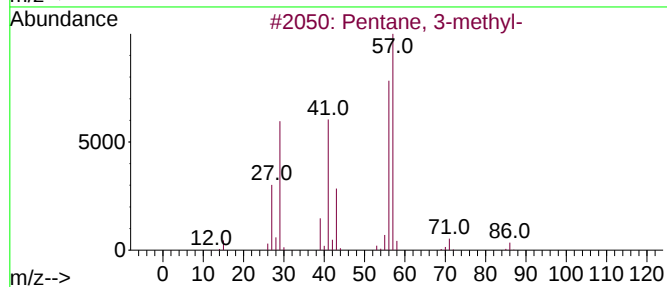
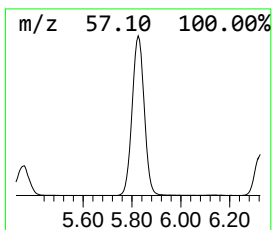
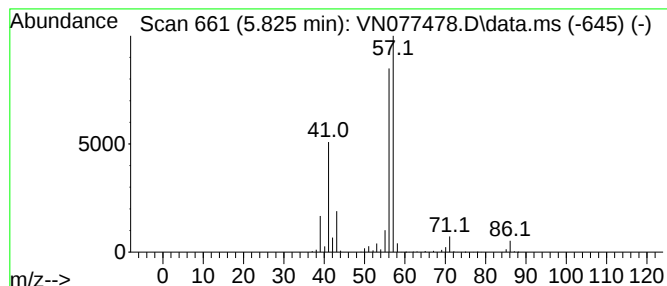
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.825	133.07 ug/l	15306800	Pentafluorobenzene	8.231

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	72
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4			Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	43
5			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

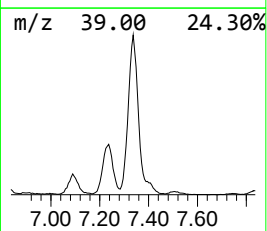
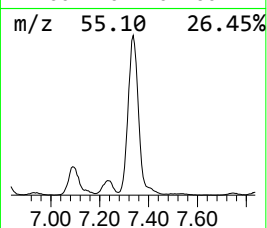
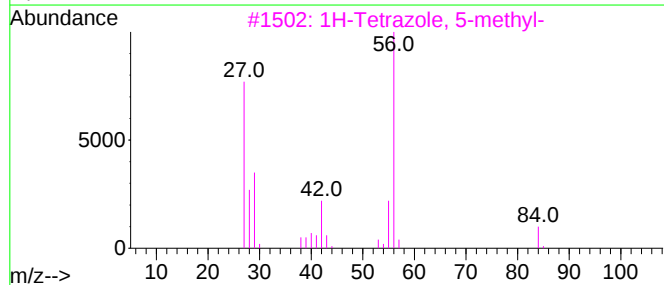
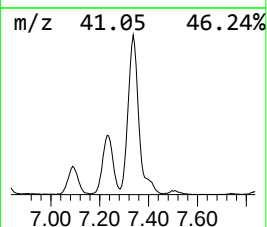
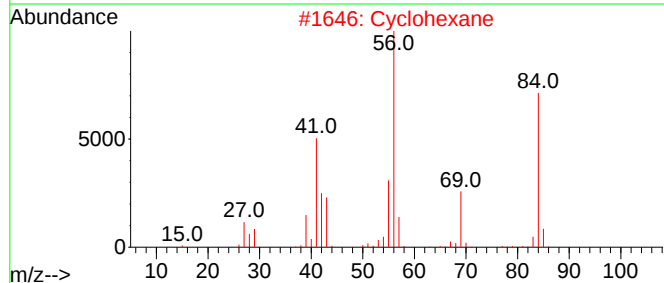
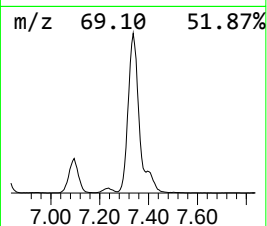
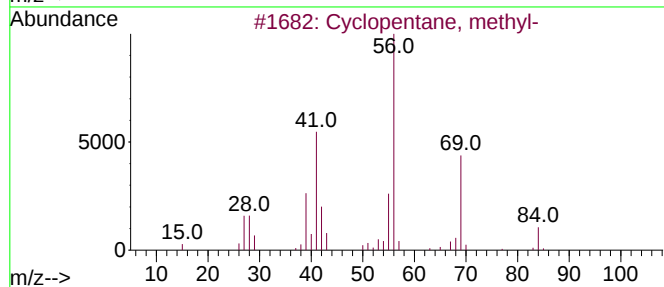
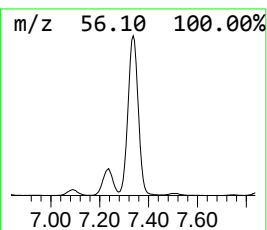
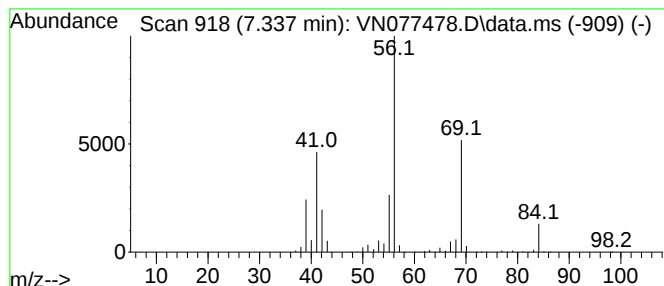
Instrument :
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.337	134.37 ug/l	15456400	Pentafluorobenzene	8.231	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclohexane	84	C6H12	000110-82-7	78
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-4

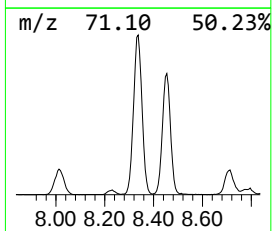
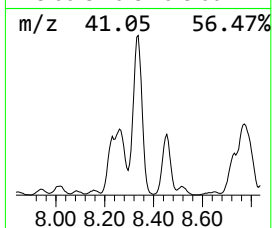
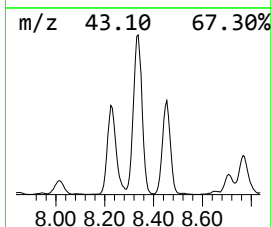
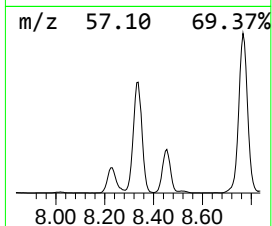
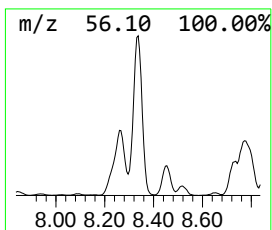
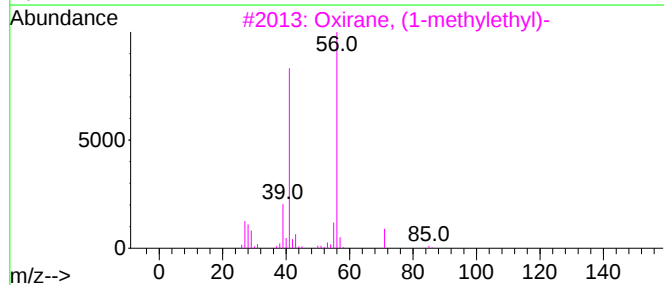
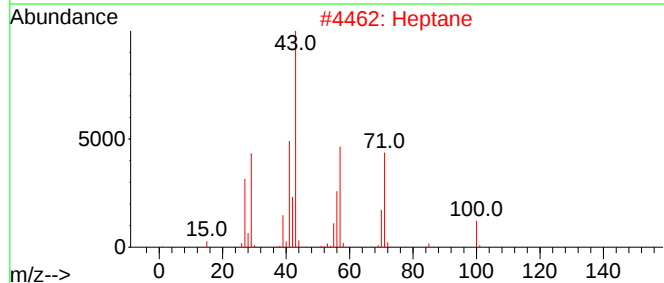
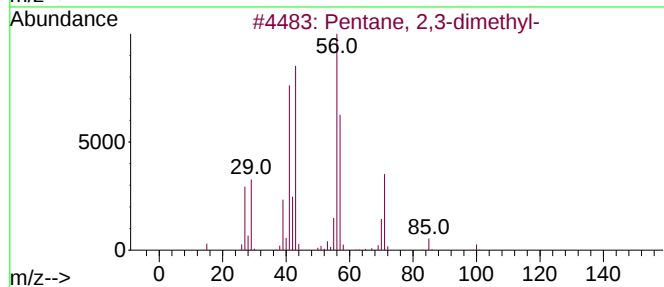
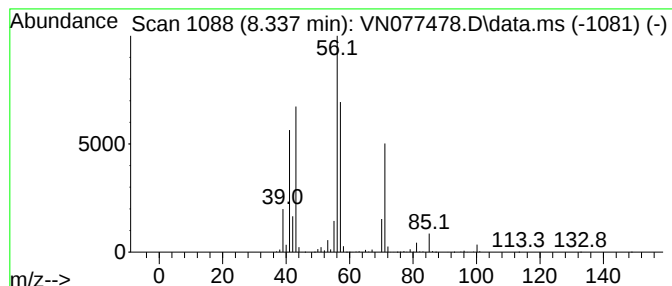
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.337	123.53 ug/l	14209100	Pentafluorobenzene	8.231

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			Heptane	100	C7H16	000142-82-5	43
3			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	43
4			Hexane, 3-methyl-	100	C7H16	000589-34-4	38
5			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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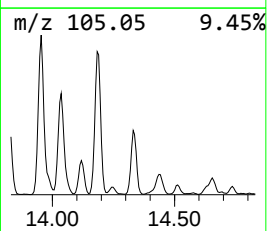
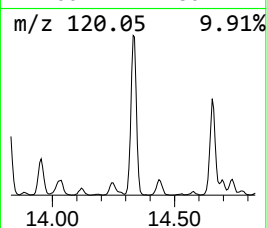
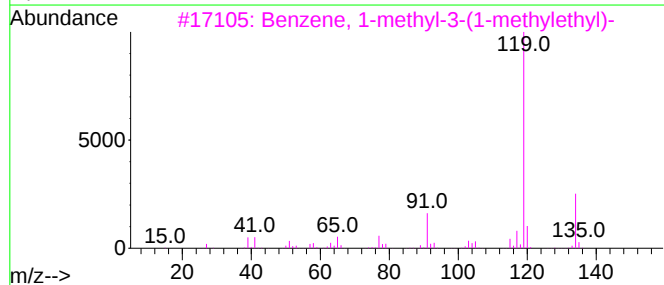
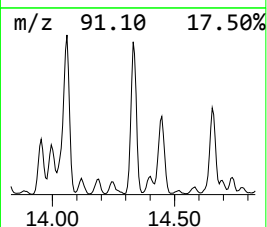
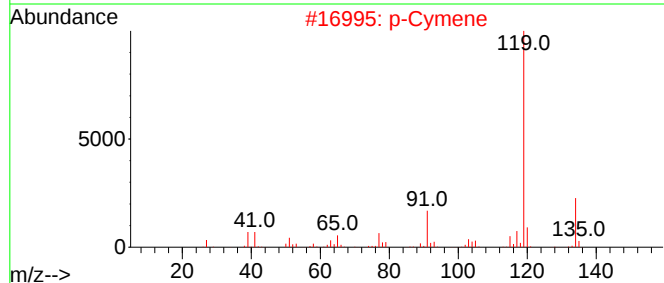
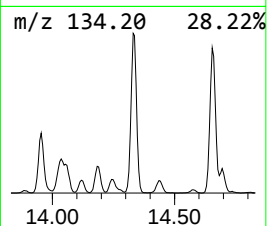
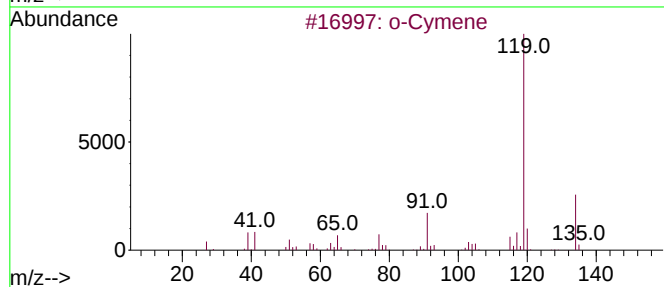
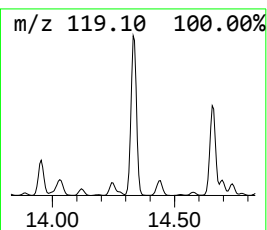
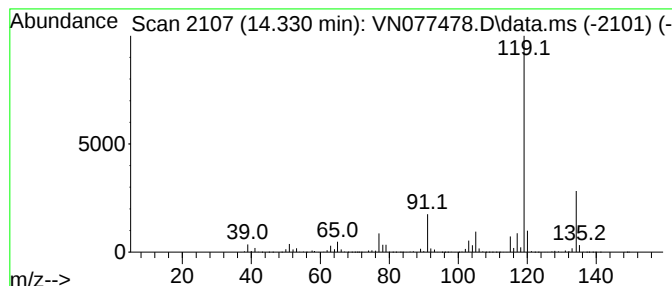
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Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.330	225.74 ug/l	14306100	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			p-Cymene	134	C10H14	000099-87-6	95
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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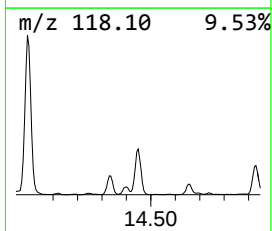
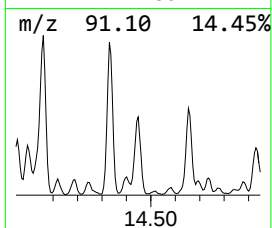
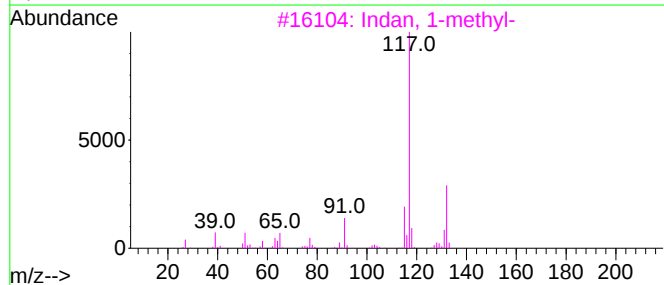
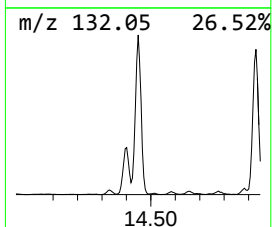
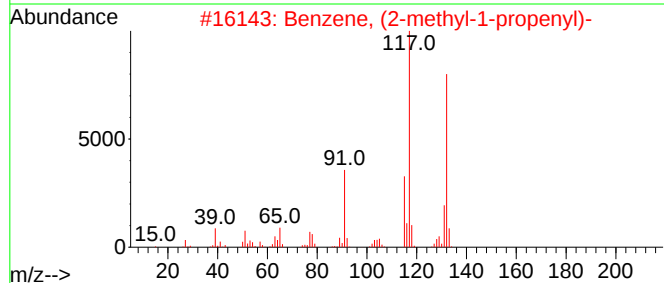
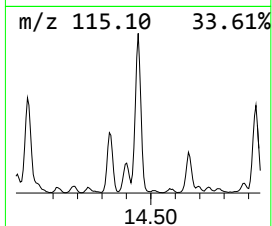
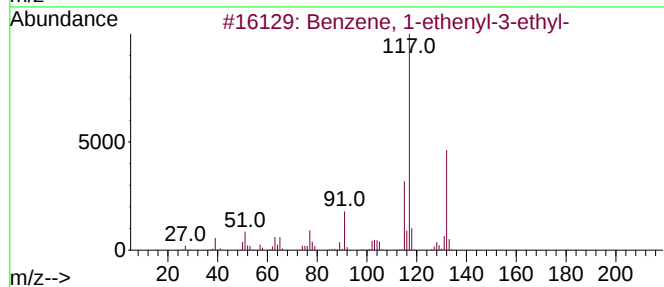
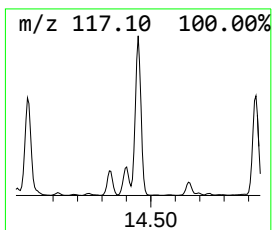
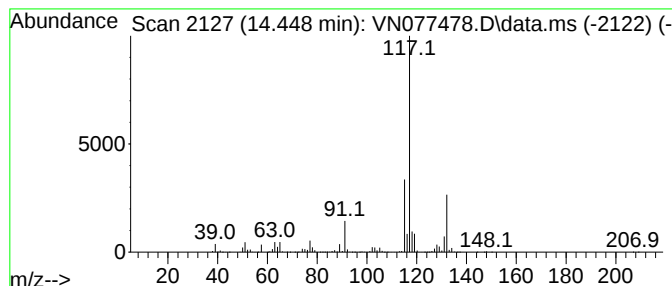
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.448	161.69 ug/l	10247100	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	90
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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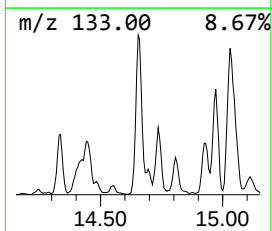
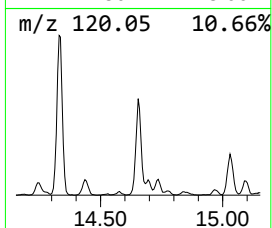
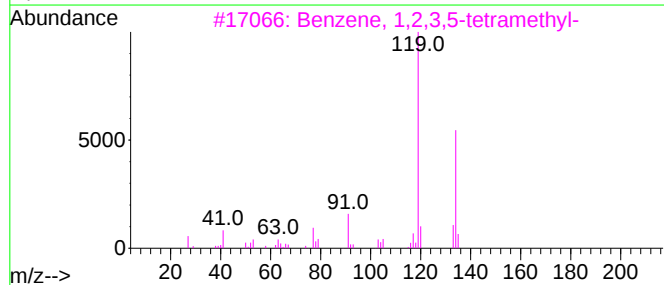
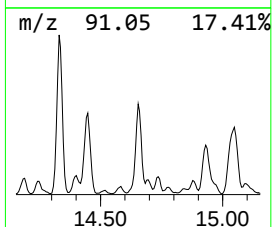
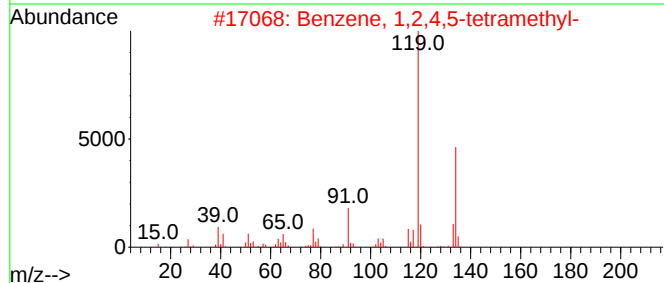
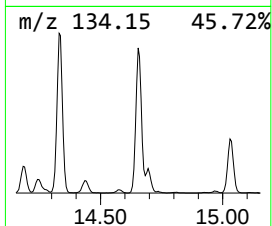
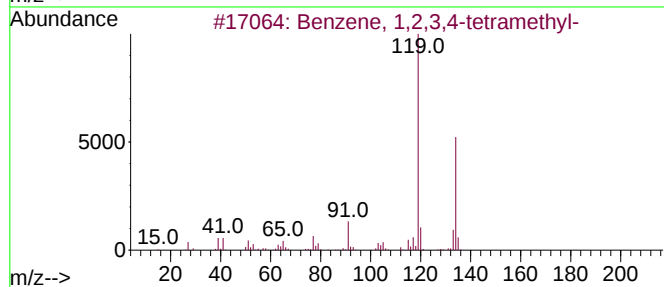
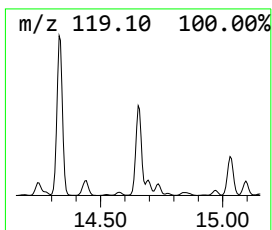
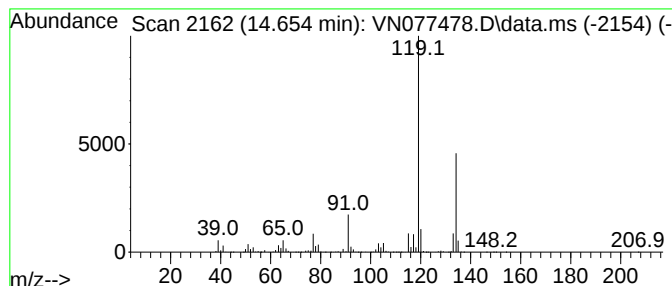
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.654	148.26 ug/l	9395900	1,4-Dichlorobenzene-d4	13.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	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 : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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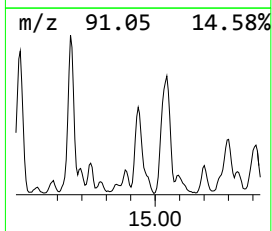
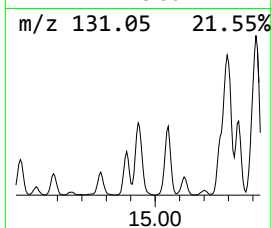
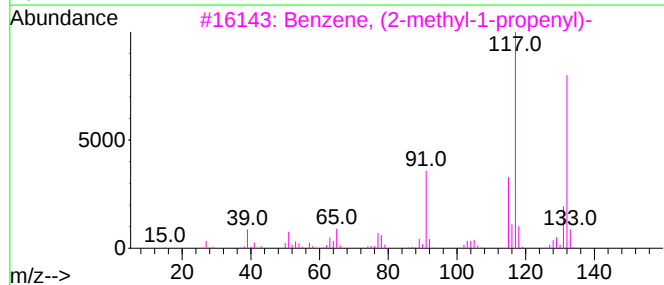
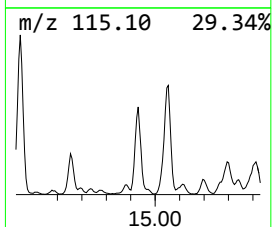
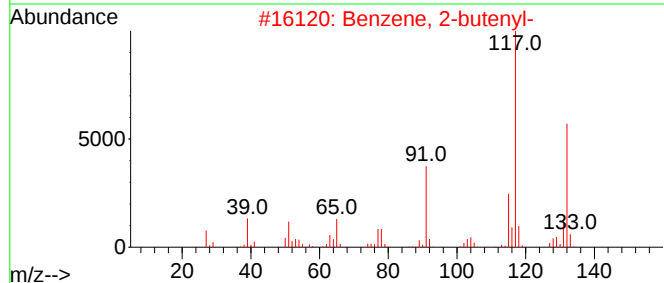
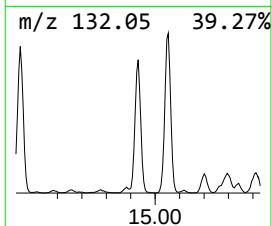
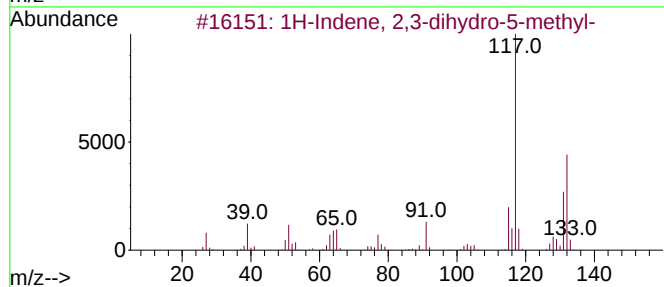
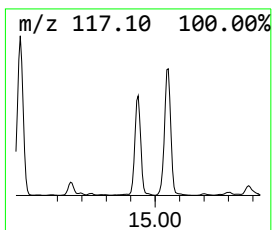
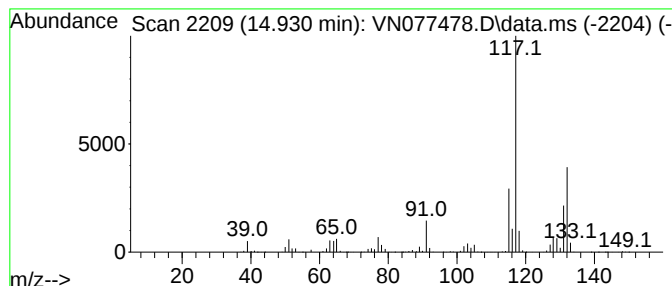
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.930	103.62 ug/l	6566590	1,4-Dichlorobenzene-d4	13.795

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-4

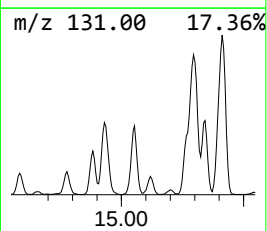
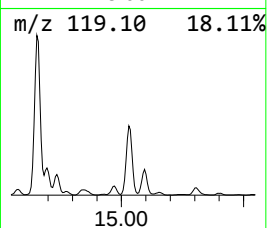
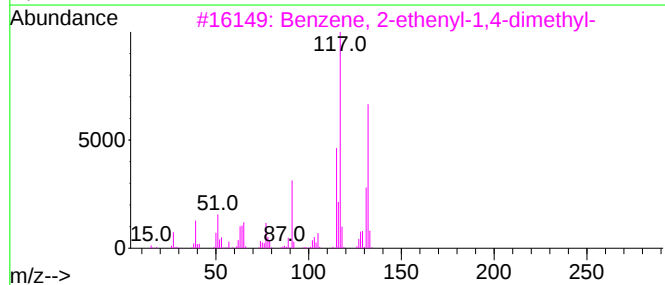
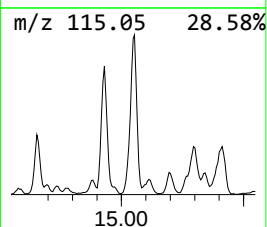
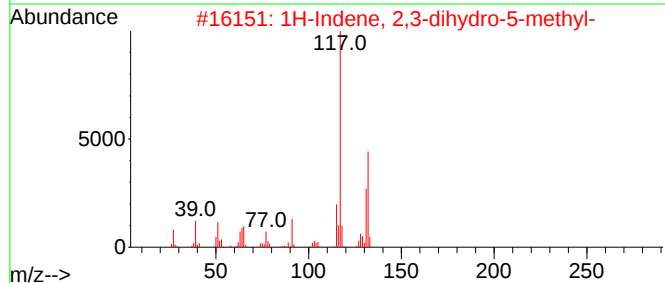
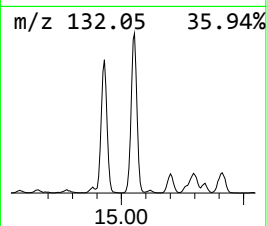
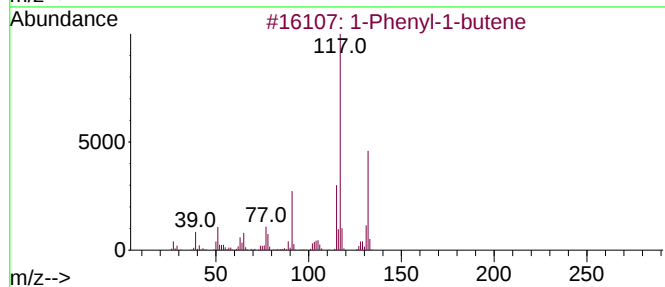
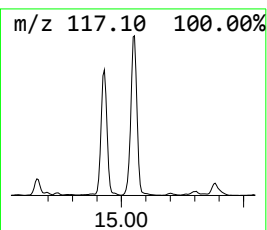
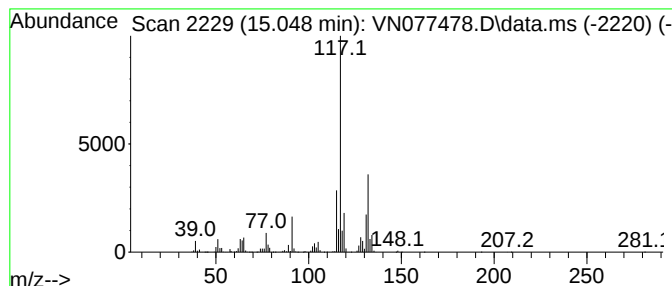
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Quant Title : SW846 8260

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

Peak Number 10 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.048	202.31 ug/l	12820900	1,4-Dichlorobenzene-d4	13.795

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	93
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077478.D
Acq On : 25 Apr 2023 18:45
Operator : JC\MD
Sample : 02460-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-4

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	353.3	ug/l	40641700	1	8.231	5751390	50.0
Pentane, 2-methyl-	5.354	205.4	ug/l	23626900	1	8.231	5751390	50.0
Pentane, 3-methyl-	5.825	133.1	ug/l	15306800	1	8.231	5751390	50.0
Cyclopentane, m...	7.337	134.4	ug/l	15456400	1	8.231	5751390	50.0
Pentane, 2,3-di...	8.337	123.5	ug/l	14209100	1	8.231	5751390	50.0
o-Cymene	14.330	225.7	ug/l	14306100	4	13.795	3168710	50.0
Benzene, 1-ethe...	14.448	161.7	ug/l	10247100	4	13.795	3168710	50.0
Benzene, 1,2,3,...	14.654	148.3	ug/l	9395900	4	13.795	3168710	50.0
1H-Indene, 2,3-...	14.930	103.6	ug/l	6566590	4	13.795	3168710	50.0
1-Phenyl-1-butene	15.048	202.3	ug/l	12820900	4	13.795	3168710	50.0