

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
 Data File : VN073144.D
 Acq On : 26 Jun 2022 06:58
 Operator : JC\MD
 Sample : N3426-04 40X
 Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 B2202-30

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\82N061722W.M

Title : SW846 8260

Signal : TIC: VN073144.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.405	82	88	94	rBV2	17890	31163	1.26%	0.097%
2	3.081	194	203	215	rBV	147443	331045	13.43%	1.027%
3	3.457	258	267	279	rBV	69669	180982	7.34%	0.561%
4	3.657	293	301	310	rVB3	16406	38125	1.55%	0.118%
5	3.928	337	347	360	rVB	20157	52865	2.14%	0.164%
6	4.187	382	391	402	rBV3	7885	26190	1.06%	0.081%
7	5.004	517	530	532	rBV2	46940	136394	5.53%	0.423%
8	5.063	532	540	557	rVB2	125976	464121	18.83%	1.440%
9	5.534	608	620	634	rBV3	89574	310768	12.61%	0.964%
10	5.857	663	675	689	rBV4	14559	59015	2.39%	0.183%
11	6.040	693	706	722	rBV3	56471	174298	7.07%	0.541%
12	6.198	724	733	743	rBV4	12136	36979	1.50%	0.115%
13	6.322	743	754	758	rBV2	18085	53354	2.16%	0.166%
14	6.381	760	764	775	rVB3	22326	57785	2.34%	0.179%
15	6.528	775	789	798	rBV3	13991	43212	1.75%	0.134%
16	6.657	801	811	819	rVB8	8980	29788	1.21%	0.092%
17	6.816	827	838	851	rBV5	24994	81680	3.31%	0.253%
18	6.963	851	863	872	rBV3	56099	173143	7.02%	0.537%
19	7.069	872	881	891	rVV2	81069	239414	9.71%	0.743%
20	7.234	901	909	924	rVB2	31769	86292	3.50%	0.268%
21	7.769	994	1000	1009	rVB2	29327	71050	2.88%	0.220%
22	7.951	1020	1031	1035	rBV2	242653	585003	23.73%	1.815%
23	8.016	1035	1042	1050	rVV2	534990	1391534	56.45%	4.317%
24	8.104	1050	1057	1069	rVB2	144006	369969	15.01%	1.148%
25	8.228	1069	1078	1087	rBV	121241	282895	11.48%	0.878%
26	8.369	1093	1102	1115	rBV	251253	584249	23.70%	1.812%
27	8.545	1115	1132	1143	rVV	479282	1336678	54.22%	4.147%
28	8.645	1143	1149	1158	rVV4	28732	75159	3.05%	0.233%
29	8.757	1158	1168	1183	rVB3	58124	150706	6.11%	0.468%
30	8.898	1183	1192	1214	rBV	685811	1539692	62.46%	4.777%
31	9.057	1214	1219	1226	rVV4	15200	29941	1.21%	0.093%
32	9.175	1226	1239	1250	rVB9	15268	64536	2.62%	0.200%
33	9.392	1260	1276	1281	rBV3	123900	334743	13.58%	1.038%
34	9.445	1281	1285	1293	rVV2	79113	147610	5.99%	0.458%
35	9.522	1293	1298	1303	rVV2	15105	36367	1.48%	0.113%
36	9.575	1303	1307	1313	rVB2	25837	48583	1.97%	0.151%

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37	9.669	1313	1323	1330	rBV4	18408	45284	1.84%	0.140%
38	9.816	1341	1348	1358	rVB	206182	418142	16.96%	1.297%
39	9.951	1358	1371	1377	rBV3	234074	523127	21.22%	1.623%
40	10.016	1377	1382	1386	rBV	54431	89016	3.61%	0.276%
41	10.151	1395	1405	1418	rBV2	80028	202929	8.23%	0.630%
42	10.304	1426	1431	1436	rVB2	65515	112640	4.57%	0.349%
43	10.380	1436	1444	1450	rBV	890313	1680675	68.18%	5.214%
44	10.439	1450	1454	1460	rVB2	52682	99402	4.03%	0.308%
45	10.563	1467	1475	1481	rVB3	57404	134083	5.44%	0.416%
46	10.822	1509	1519	1528	rVB8	31615	91923	3.73%	0.285%
47	10.898	1528	1532	1537	rVB5	16097	25956	1.05%	0.081%
48	10.986	1539	1547	1553	rBV7	11157	26904	1.09%	0.083%
49	11.092	1560	1565	1573	rVB3	17862	38922	1.58%	0.121%
50	11.463	1619	1628	1640	rVB3	47719	120215	4.88%	0.373%
51	11.569	1640	1646	1651	rBV	36881	66200	2.69%	0.205%
52	11.680	1656	1665	1674	rBV	955821	1700473	68.98%	5.275%
53	11.780	1674	1682	1691	rVB	378199	637163	25.85%	1.977%
54	11.886	1692	1700	1712	rBV	1257968	2465072	100.00%	7.647%
55	12.216	1745	1756	1765	rBV	375739	657439	26.67%	2.040%
56	12.516	1801	1807	1813	rBV	43556	77559	3.15%	0.241%
57	12.669	1823	1833	1843	rVB	792253	1414748	57.39%	4.389%
58	12.857	1860	1865	1870	rVV	170802	268438	10.89%	0.833%
59	12.921	1870	1876	1879	rVV	691026	1150747	46.68%	3.570%
60	12.951	1879	1881	1885	rVV	288981	431611	17.51%	1.339%
61	12.998	1885	1889	1897	rVB	408093	651195	26.42%	2.020%
62	13.157	1909	1916	1924	rBV	240693	412656	16.74%	1.280%
63	13.251	1929	1932	1935	rVV2	21026	31440	1.28%	0.098%
64	13.304	1935	1941	1948	rVV	1192034	1894864	76.87%	5.878%
65	13.427	1955	1962	1967	rBV3	23129	64107	2.60%	0.199%
66	13.498	1970	1974	1978	rVV2	34862	57519	2.33%	0.178%
67	13.545	1978	1982	1986	rVV7	20308	39428	1.60%	0.122%
68	13.610	1986	1993	2014	rVB2	1083572	2301401	93.36%	7.140%
69	13.774	2016	2021	2022	rBV	79631	97311	3.95%	0.302%
70	13.810	2022	2027	2030	rVV2	275220	513089	20.81%	1.592%
71	13.845	2030	2033	2044	rVV2	244490	468946	19.02%	1.455%
72	13.933	2044	2048	2053	rVB2	33289	50676	2.06%	0.157%
73	14.004	2054	2060	2065	rVV	56734	90359	3.67%	0.280%
74	14.068	2065	2071	2073	rVV	123841	197665	8.02%	0.613%
75	14.092	2073	2075	2080	rVV	118493	179683	7.29%	0.557%
76	14.151	2080	2085	2092	rVV	204526	322926	13.10%	1.002%

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77	14.216	2092	2096	2099	rVV3	20624	32101	1.30%	0.100%
78	14.263	2099	2104	2109	rVB2	86784	140166	5.69%	0.435%
79	14.398	2121	2127	2132	rBV	45476	77592	3.15%	0.241%
80	14.474	2132	2140	2143	rVV	117586	211026	8.56%	0.655%
81	14.515	2143	2147	2151	rVV	159974	254898	10.34%	0.791%
82	14.557	2151	2154	2158	rVV2	65415	100024	4.06%	0.310%
83	14.598	2158	2161	2164	rVV	32373	46264	1.88%	0.144%
84	14.627	2164	2166	2172	rVV5	15710	28467	1.15%	0.088%
85	14.698	2174	2178	2181	rVV	31353	44505	1.81%	0.138%
86	14.745	2181	2186	2190	rVV	99054	176855	7.17%	0.549%
87	14.786	2190	2193	2198	rVV3	56348	88958	3.61%	0.276%
88	14.863	2198	2206	2211	rVV4	139974	346169	14.04%	1.074%
89	14.915	2211	2215	2224	rVV3	36877	84146	3.41%	0.261%
90	15.015	2227	2232	2240	rVV8	16820	38902	1.58%	0.121%
91	15.127	2240	2251	2263	rVV4	60061	201354	8.17%	0.625%
92	15.233	2263	2269	2279	rVB4	44896	115777	4.70%	0.359%
93	15.433	2292	2303	2310	rVB	136761	284275	11.53%	0.882%
94	15.539	2316	2321	2326	rBV9	18270	38937	1.58%	0.121%
95	15.627	2333	2336	2345	rVB7	22421	40720	1.65%	0.126%
96	15.721	2347	2352	2359	rBV2	26556	58454	2.37%	0.181%
97	15.898	2379	2382	2391	rVB5	15779	34645	1.41%	0.107%
98	16.104	2410	2417	2423	rBV9	15713	36467	1.48%	0.113%
99	16.527	2482	2489	2495	rBV	72251	157129	6.37%	0.487%
100	16.727	2516	2523	2533	rBV	36885	89545	3.63%	0.278%

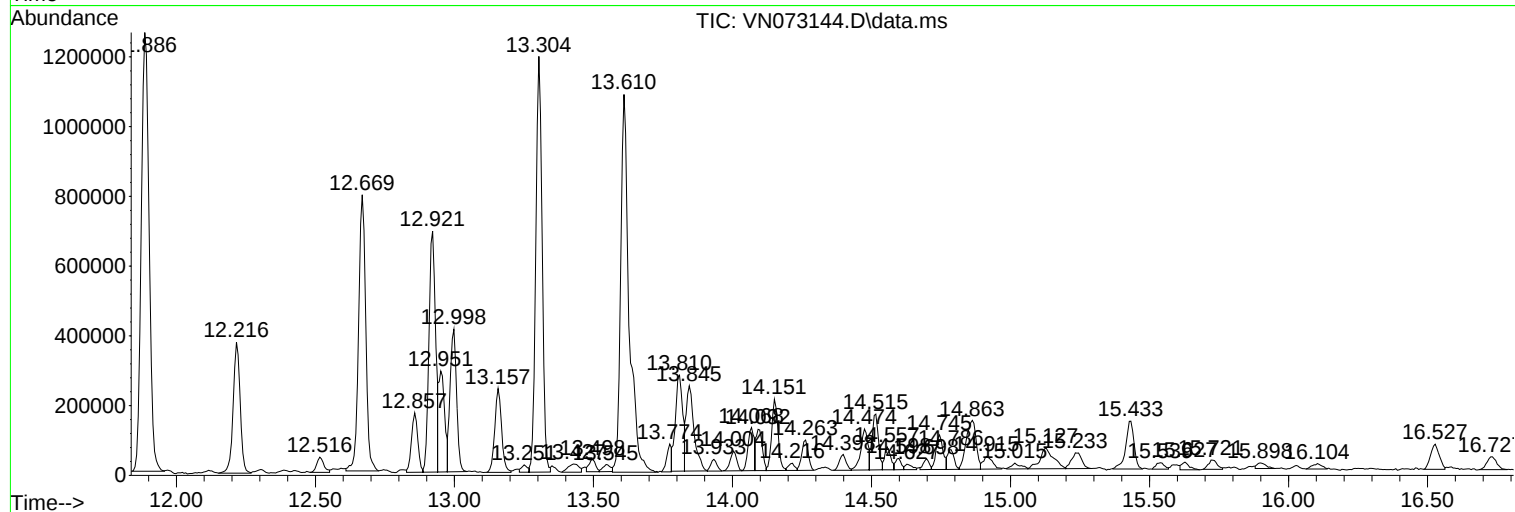
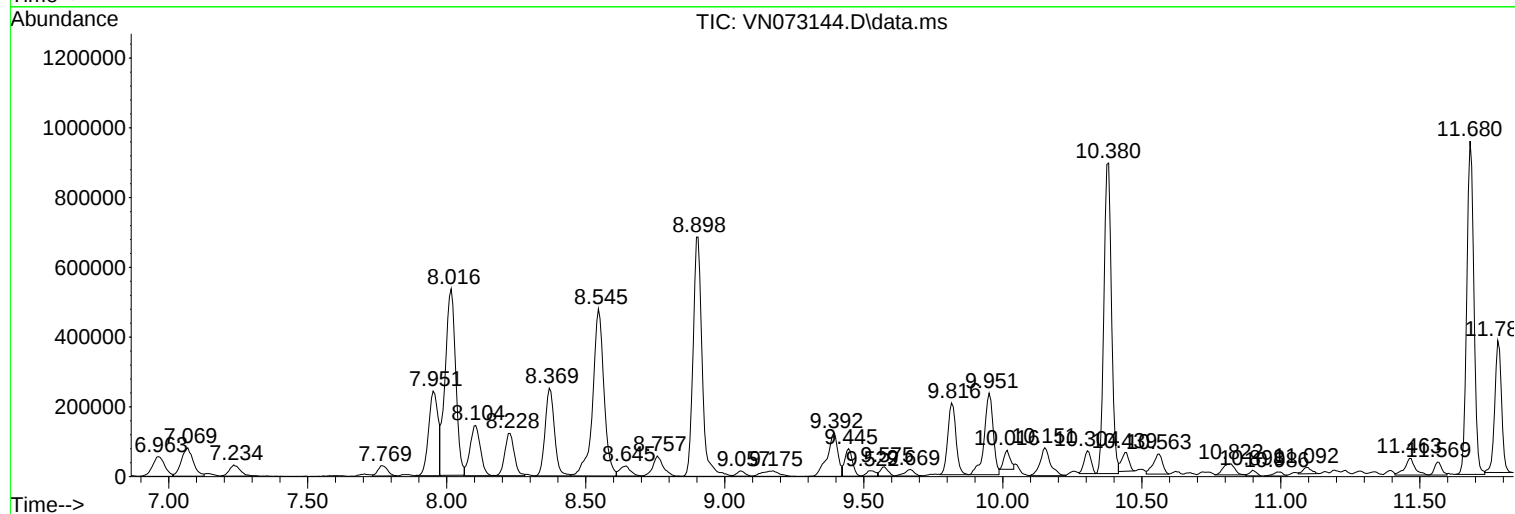
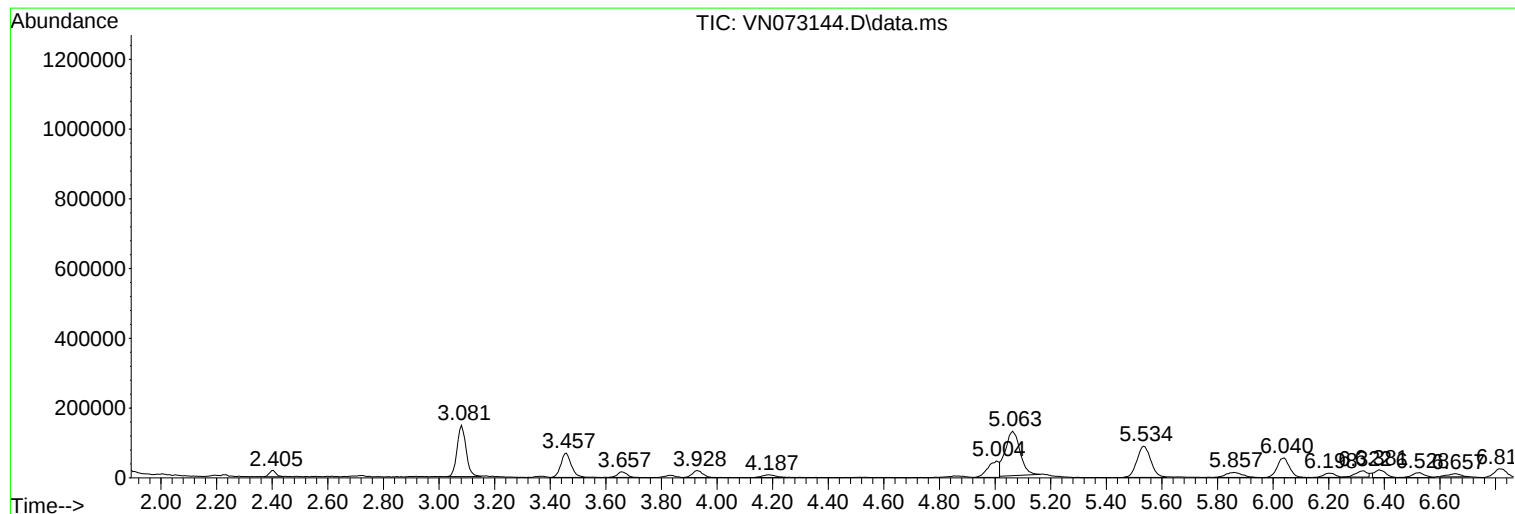
Sum of corrected areas: 32234637

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

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



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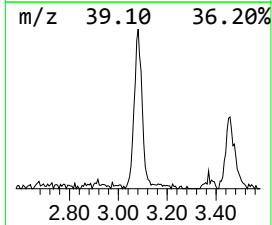
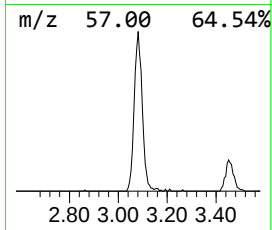
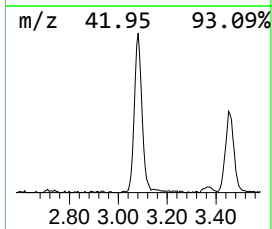
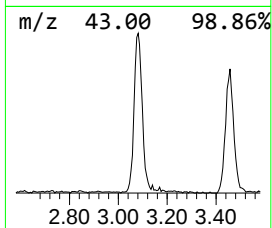
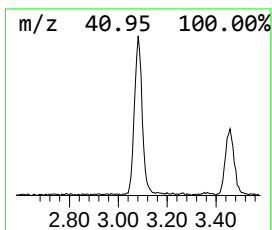
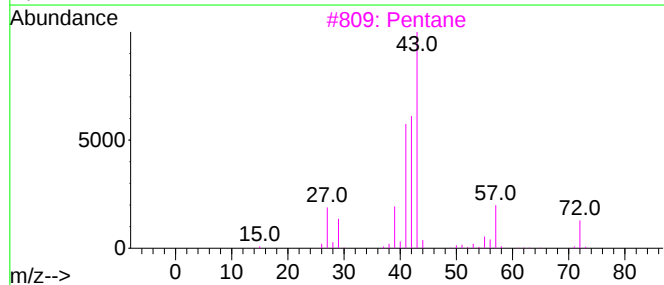
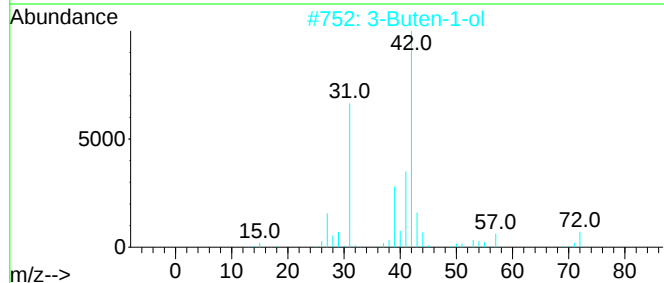
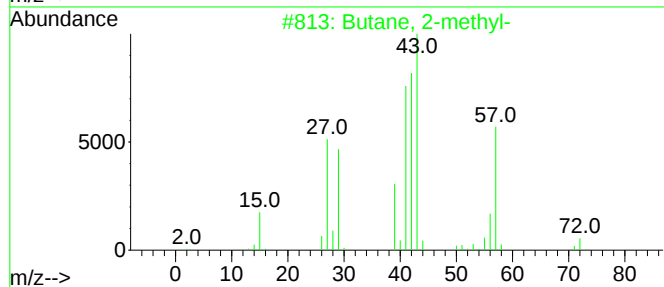
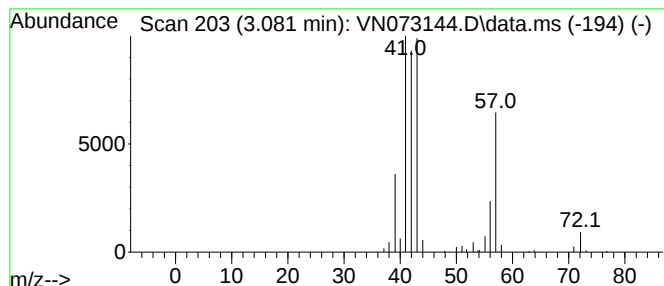
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.081	11.90 ug/l	331045	Pentafluorobenzene	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			3-Buten-1-ol	72	C4H8O	000627-27-0	47
3			Pentane	72	C5H12	000109-66-0	40
4			2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	28
5			Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	10



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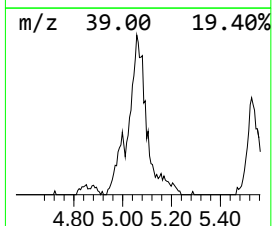
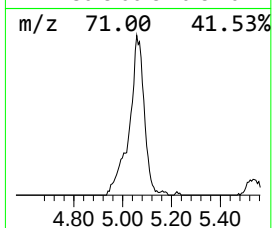
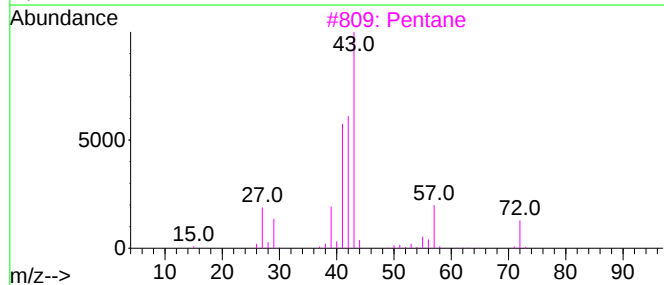
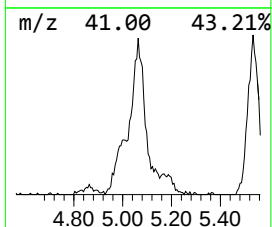
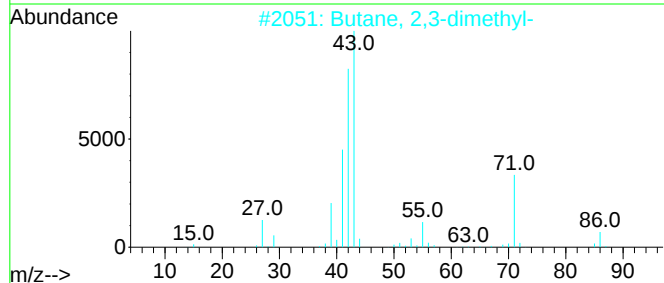
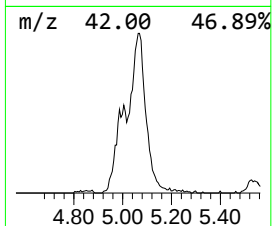
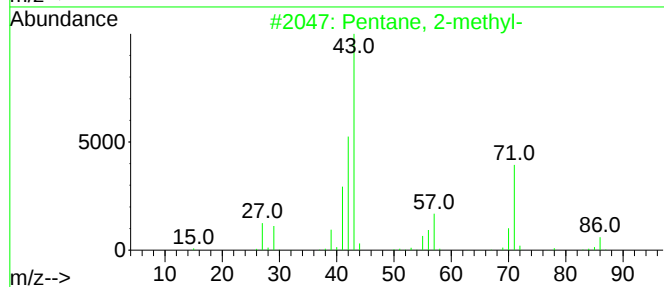
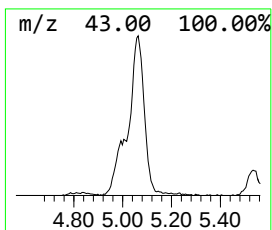
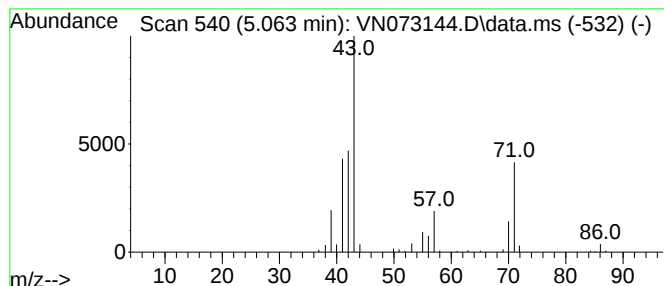
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Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.063	16.68 ug/l	464121	Pentafluorobenzene	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	50
3			Pentane	72	C5H12	000109-66-0	47
4			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38
5			Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	33



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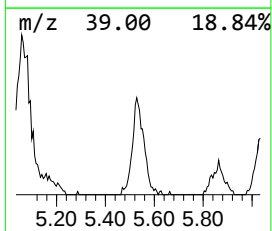
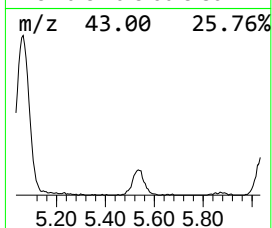
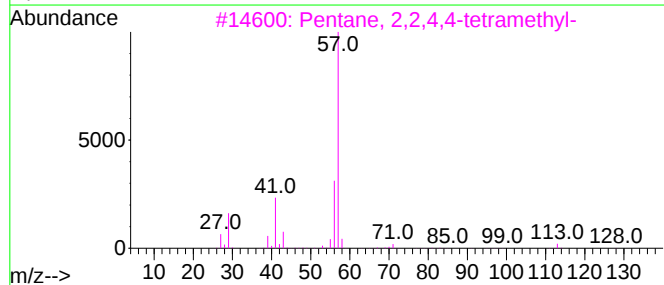
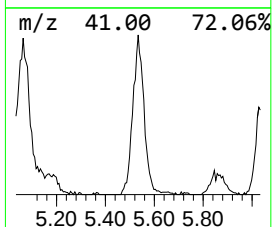
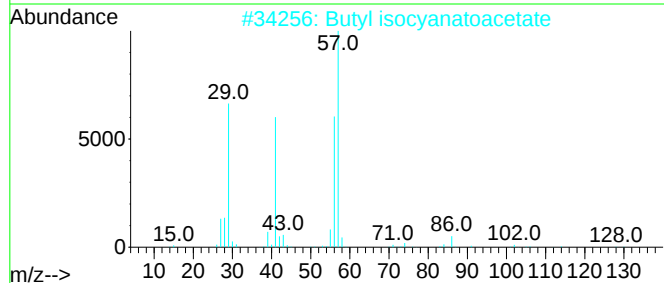
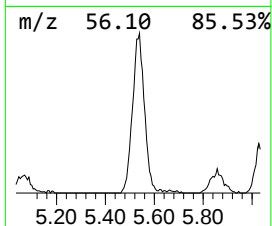
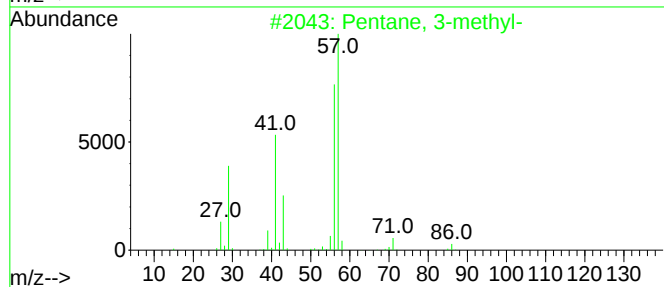
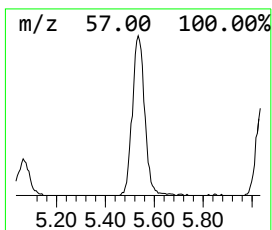
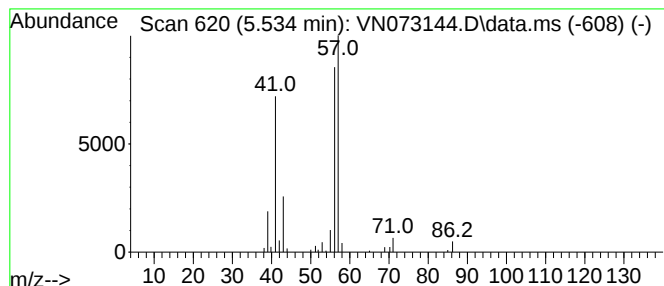
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Quant Title : SW846 8260

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.
5.534	11.17 ug/l	310768	Pentafluorobenzene	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	64
3			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45
4			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
5			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

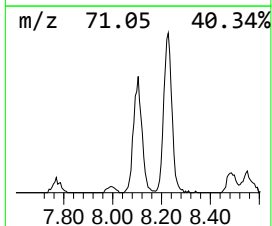
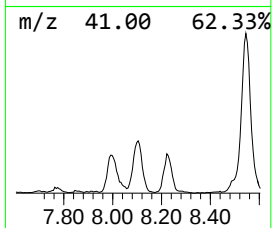
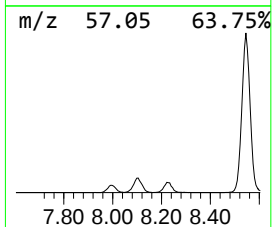
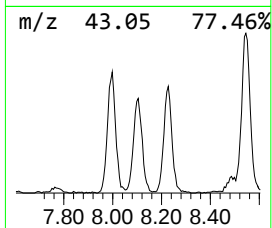
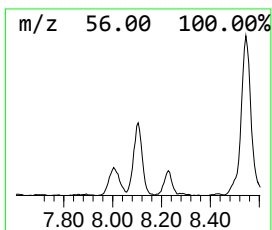
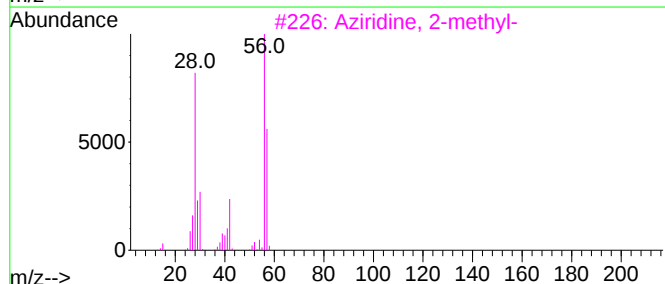
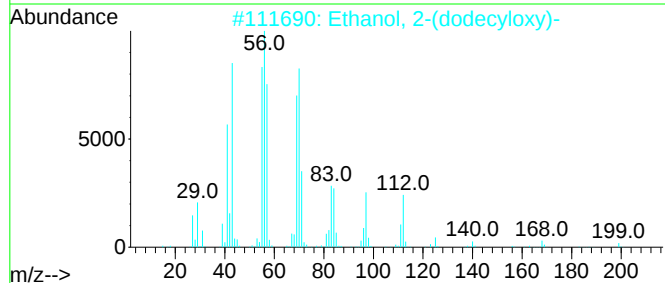
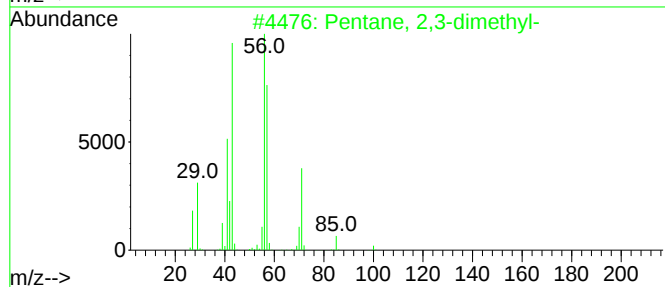
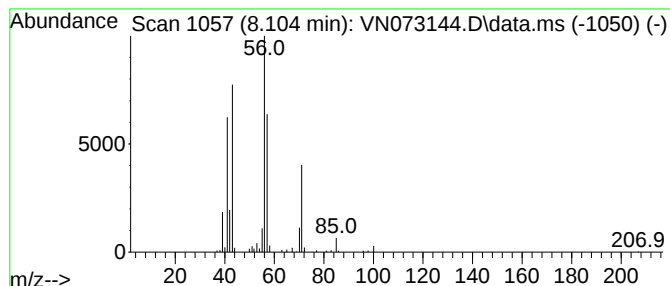
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	13.29 ug/l	369969	Pentafluorobenzene	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	94
2			Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	64
3			Aziridine, 2-methyl-	57	C3H7N	000075-55-8	46
4			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	38
5			Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

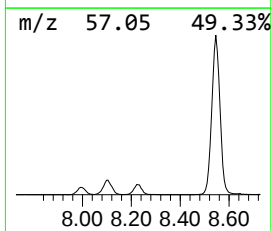
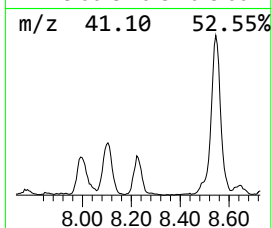
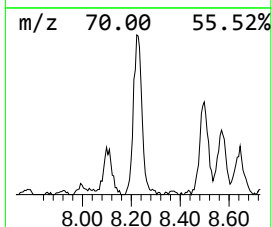
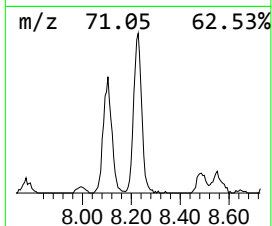
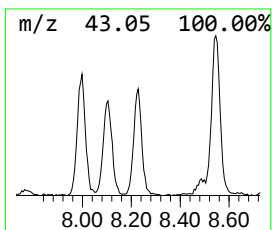
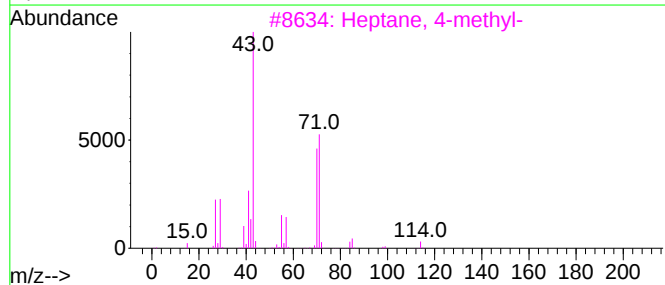
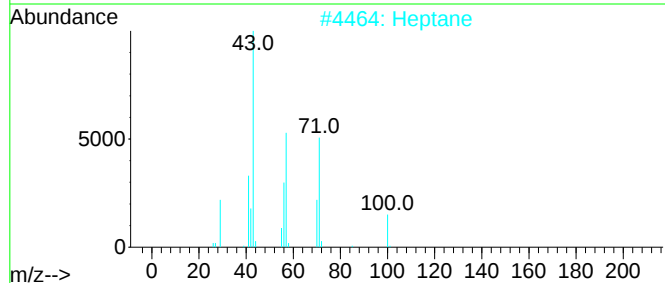
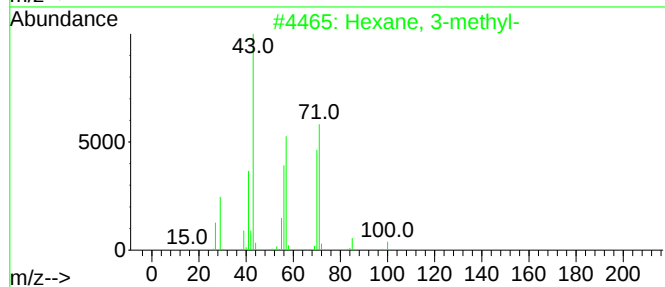
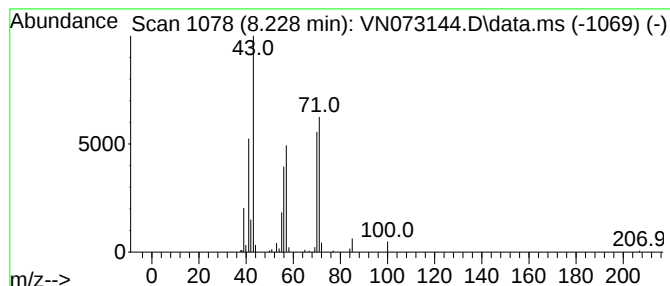
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Quant Title : SW846 8260

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

Peak Number 5 Hexane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.228	10.16 ug/l	282895	Pentafluorobenzene	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	94
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

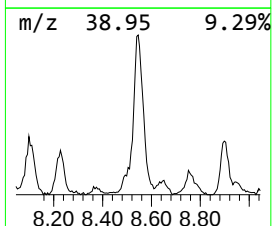
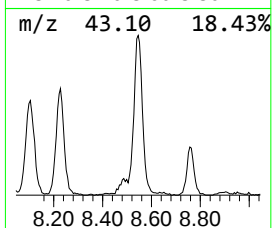
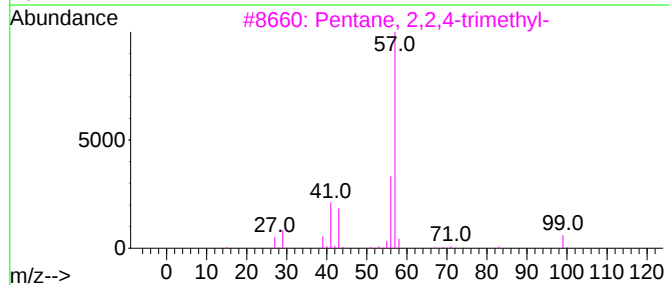
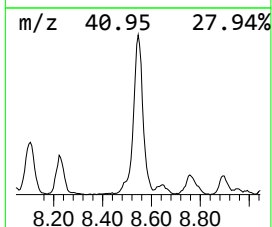
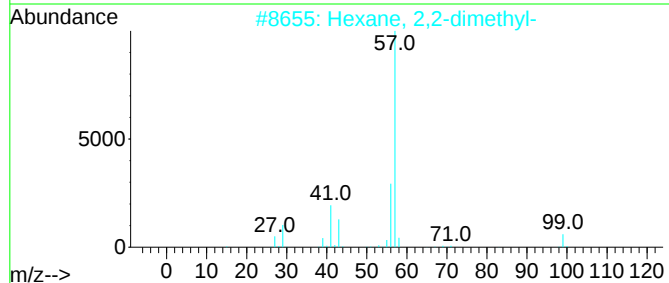
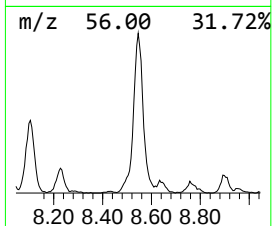
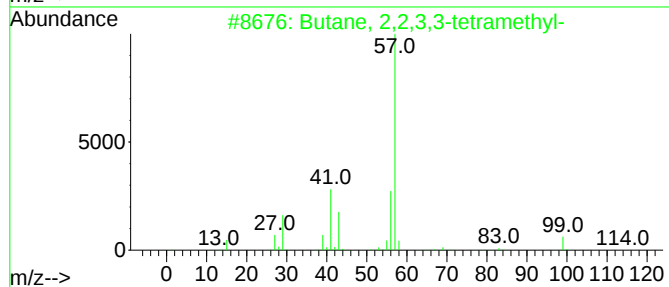
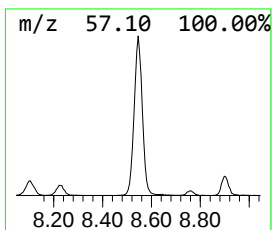
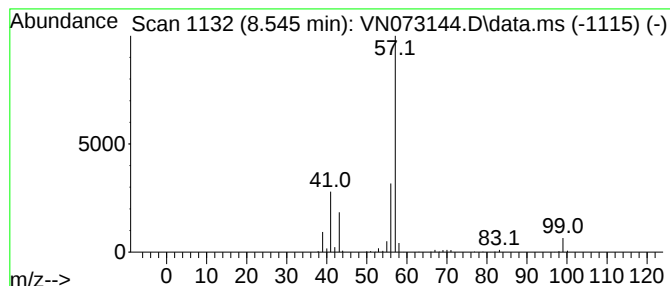
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Quant Title : SW846 8260

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

Peak Number 6 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.545	43.41 ug/l	1336680	1,4-Difluorobenzene	8.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5			Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

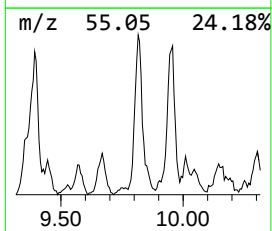
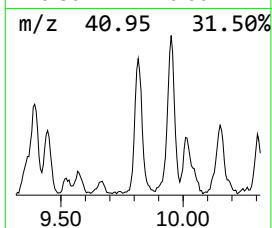
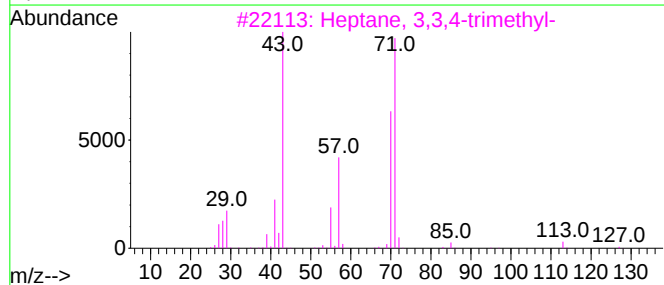
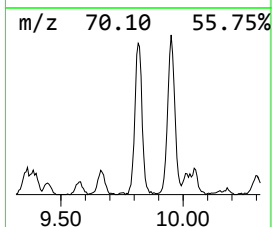
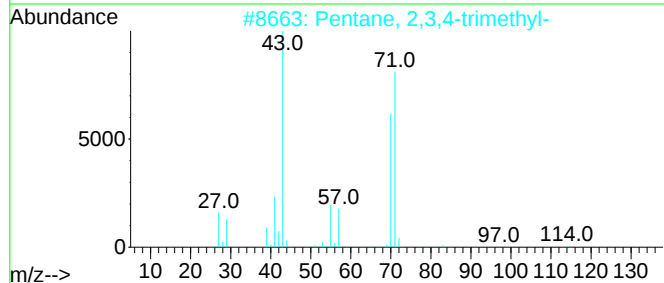
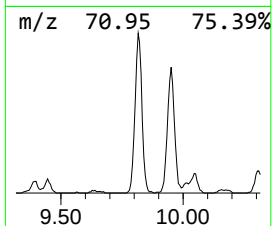
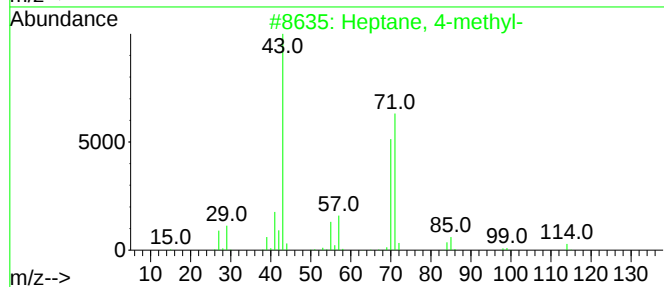
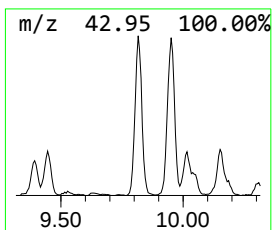
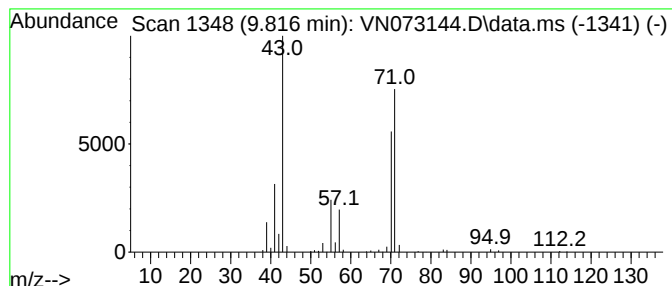
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Quant Title : SW846 8260

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

Peak Number 7 Heptane, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.816	13.58 ug/l	418142	1,4-Difluorobenzene	8.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 4-methyl-	114	C8H18	000589-53-7	78
2			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
3			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	78
4			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	64
5			Pentane, 3-ethyl-	100	C7H16	000617-78-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

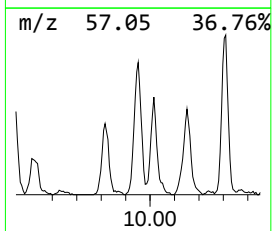
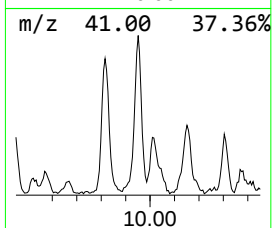
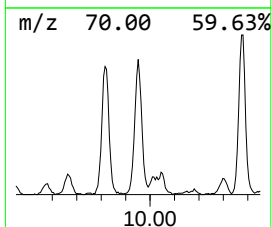
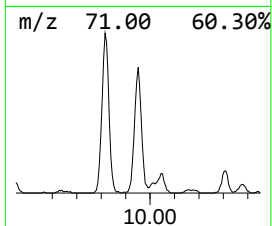
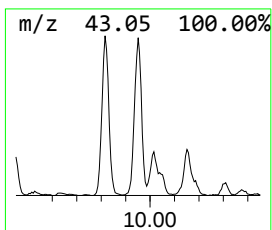
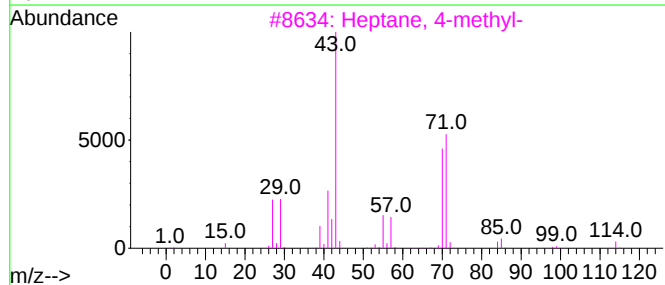
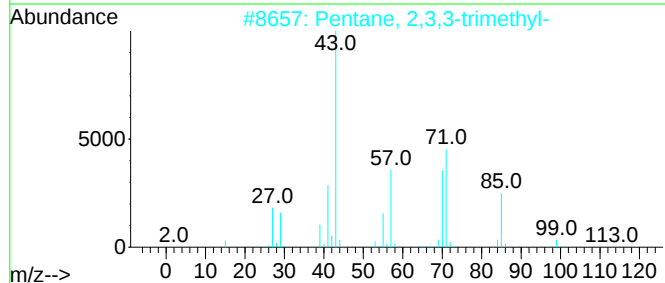
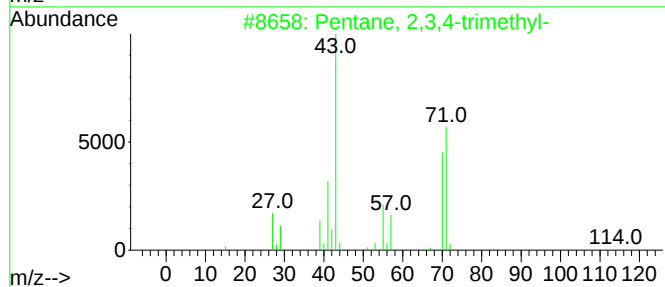
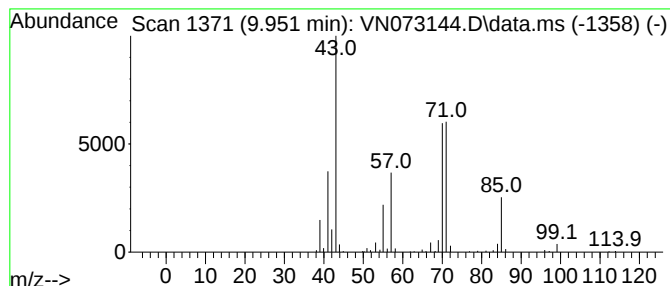
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Quant Title : SW846 8260

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

Peak Number 8 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.951	16.99 ug/l	523127	1,4-Difluorobenzene	8.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	76
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4			Hexane, 3-methyl-	100	C7H16	000589-34-4	59
5			Diisooamyl ether	158	C10H22O	000544-01-4	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

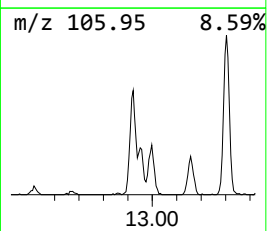
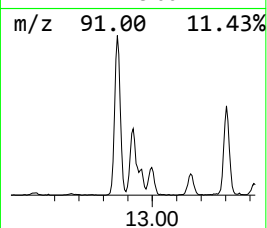
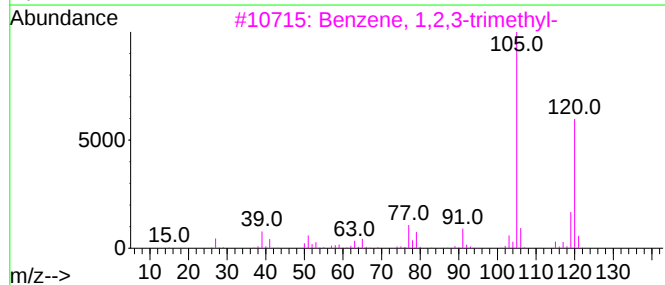
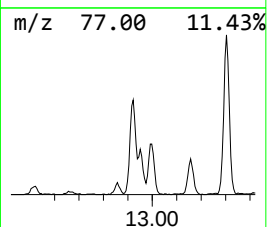
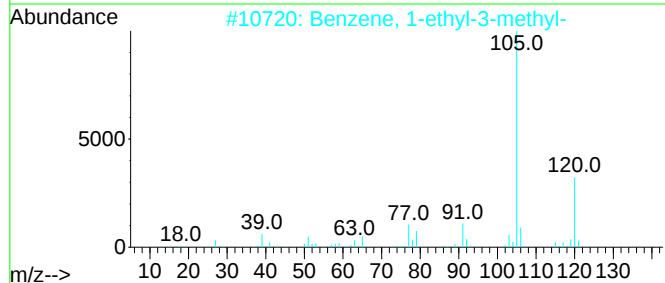
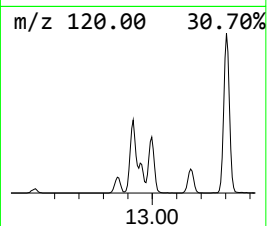
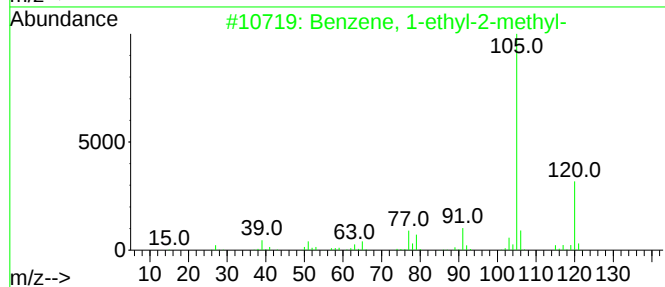
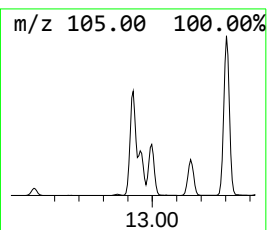
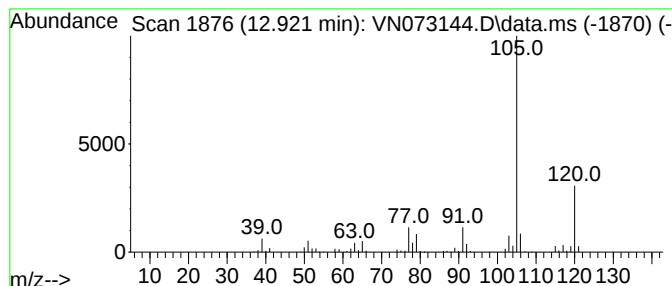
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.922	25.00 ug/l	1150750	1,4-Dichlorobenzene-d4	13.610

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-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

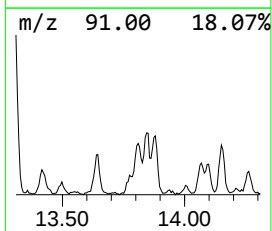
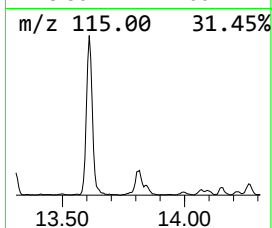
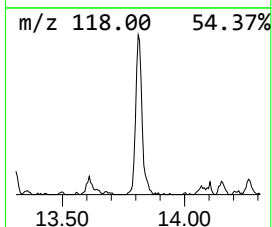
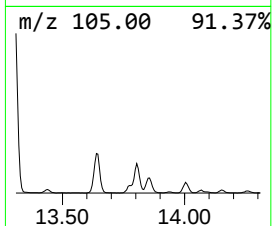
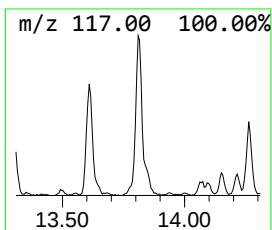
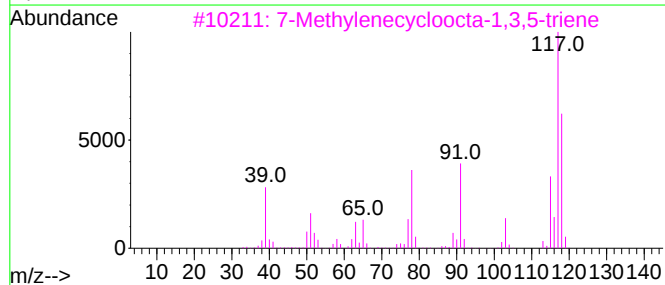
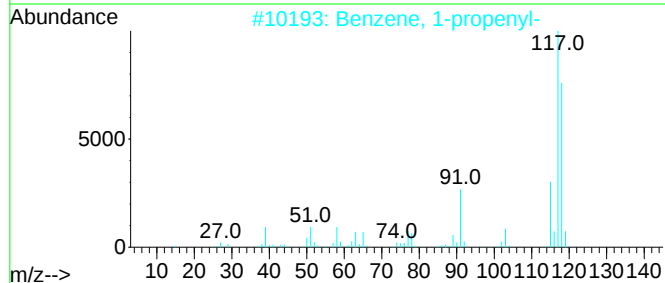
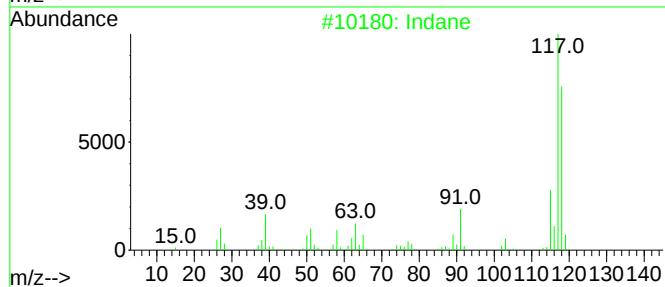
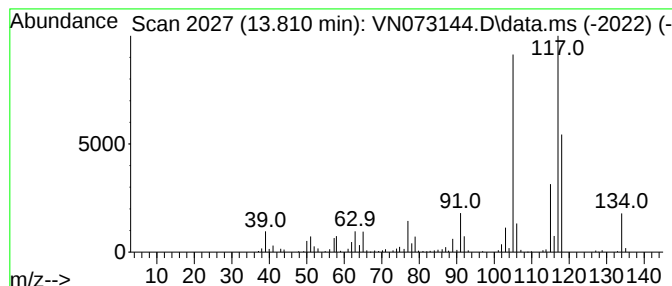
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

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

Peak Number 10 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	11.15 ug/l	513089	1,4-Dichlorobenzene-d4	13.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	64
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	58
3			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	58
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	52
5			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073144.D
Acq On : 26 Jun 2022 06:58
Operator : JC\MD
Sample : N3426-04 40X
Misc : 6.11g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2202-30

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.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-	3.081	11.9	ug/l	331045	1	8.016	1391530	50.0
Pentane, 2-methyl-	5.063	16.7	ug/l	464121	1	8.016	1391530	50.0
Pentane, 3-methyl-	5.534	11.2	ug/l	310768	1	8.016	1391530	50.0
Pentane, 2,3-di...	8.104	13.3	ug/l	369969	1	8.016	1391530	50.0
Hexane, 3-methyl-	8.228	10.2	ug/l	282895	1	8.016	1391530	50.0
Butane, 2,2,3,3...	8.545	43.4	ug/l	1336680	2	8.904	1539690	50.0
Heptane, 4-methyl-	9.816	13.6	ug/l	418142	2	8.904	1539690	50.0
Pentane, 2,3,4-...	9.951	17.0	ug/l	523127	2	8.904	1539690	50.0
Benzene, 1-ethy...	12.922	25.0	ug/l	1150750	4	13.610	2301400	50.0
Indane	13.810	11.2	ug/l	513089	4	13.610	2301400	50.0