

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
 Data File : VN071194.D
 Acq On : 08 Mar 2022 16:53
 Operator : JC\MD
 Sample : N1815-01 10X
 Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 P-1

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\82N030722W.M
 Title : SW846 8260

Signal : TIC: VN071194.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.075	512	525	526	rBV	40415	94818	1.65%	0.108%
2	5.140	526	536	555	rVB3	158815	629053	10.93%	0.714%
3	5.610	601	616	630	rBV2	135490	482738	8.39%	0.548%
4	6.116	688	702	717	rBV	194641	595572	10.35%	0.676%
5	6.887	821	833	844	rBV4	27532	91139	1.58%	0.104%
6	7.040	847	859	868	rBV3	100982	320395	5.57%	0.364%
7	7.145	868	877	898	rVB	188787	587946	10.22%	0.668%
8	7.840	988	995	1003	rVB2	47998	121023	2.10%	0.137%
9	8.016	1014	1025	1029	rBV	525703	1285322	22.34%	1.460%
10	8.081	1029	1036	1044	rVV2	1452102	3889150	67.59%	4.417%
11	8.169	1044	1051	1062	rVB	291903	755684	13.13%	0.858%
12	8.292	1062	1072	1080	rBV2	678265	1596583	27.75%	1.813%
13	8.434	1088	1096	1109	rVB	613086	1352307	23.50%	1.536%
14	8.610	1109	1126	1136	rBV	990332	3141472	54.59%	3.568%
15	8.704	1136	1142	1153	rVB	212584	495090	8.60%	0.562%
16	8.822	1153	1162	1174	rVB	514161	1088267	18.91%	1.236%
17	8.963	1176	1186	1200	rBV	1648156	3537695	61.48%	4.018%
18	9.457	1254	1270	1275	rBV	1144733	2779877	48.31%	3.157%
19	9.510	1275	1279	1287	rVB	398540	740081	12.86%	0.841%
20	9.586	1287	1292	1296	rBV	44338	89958	1.56%	0.102%
21	9.633	1296	1300	1306	rVB	134569	244870	4.26%	0.278%
22	9.728	1306	1316	1323	rVB3	169769	415285	7.22%	0.472%
23	9.875	1334	1341	1351	rVB2	675826	1404739	24.41%	1.595%
24	9.969	1351	1357	1359	rBV2	121613	205091	3.56%	0.233%
25	10.016	1359	1365	1370	rVB	653789	1190017	20.68%	1.352%
26	10.075	1370	1375	1379	rBV	705103	1280771	22.26%	1.455%
27	10.210	1389	1398	1410	rBV	790059	1903939	33.09%	2.162%
28	10.322	1410	1417	1419	rBV3	93611	167615	2.91%	0.190%
29	10.369	1419	1425	1430	rVV2	325959	615717	10.70%	0.699%
30	10.439	1430	1437	1449	rVV	2826973	5754292	100.00%	6.535%
31	10.557	1452	1457	1460	rVV4	100353	184181	3.20%	0.209%
32	10.622	1460	1468	1475	rVV4	603375	1344351	23.36%	1.527%
33	10.728	1482	1486	1490	rVV4	65322	122479	2.13%	0.139%
34	10.775	1490	1494	1502	rVB	282495	530306	9.22%	0.602%
35	10.875	1502	1511	1521	rBV5	279244	865343	15.04%	0.983%
36	10.963	1521	1526	1531	rVB	181719	316410	5.50%	0.359%

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37	11.051	1531	1541	1546	rBV	331944	664833	11.55%	0.755%
38	11.151	1546	1558	1565	rBV	292212	813380	14.14%	0.924%
39	11.222	1565	1570	1572	rVV3	147911	263224	4.57%	0.299%
40	11.257	1573	1576	1577	rVV2	155271	199608	3.47%	0.227%
41	11.286	1577	1581	1586	rVV	387587	735313	12.78%	0.835%
42	11.339	1586	1590	1596	rVV	514747	974611	16.94%	1.107%
43	11.392	1596	1599	1604	rVV3	146571	263679	4.58%	0.299%
44	11.451	1604	1609	1613	rVV3	352425	617588	10.73%	0.701%
45	11.522	1613	1621	1633	rVB4	817911	2151593	37.39%	2.444%
46	11.622	1633	1638	1650	rBV	733768	1331272	23.14%	1.512%
47	11.739	1650	1658	1667	rBV	2264574	4077103	70.85%	4.631%
48	11.839	1668	1675	1684	rVV	2593005	4657290	80.94%	5.290%
49	11.922	1684	1689	1694	rVV4	193782	421613	7.33%	0.479%
50	11.986	1694	1700	1704	rVV2	356080	710692	12.35%	0.807%
51	12.027	1704	1707	1712	rVB2	212299	329244	5.72%	0.374%
52	12.086	1712	1717	1720	rBV6	55969	100389	1.74%	0.114%
53	12.251	1738	1745	1751	rBV3	422319	929531	16.15%	1.056%
54	12.363	1758	1764	1769	rVB	471126	778145	13.52%	0.884%
55	12.451	1769	1779	1781	rBV5	375302	833671	14.49%	0.947%
56	12.574	1795	1800	1804	rBV	225984	451795	7.85%	0.513%
57	12.627	1804	1809	1810	rVV4	121778	200751	3.49%	0.228%
58	12.651	1810	1813	1816	rVV	170293	230656	4.01%	0.262%
59	12.727	1821	1826	1830	rBV	1562908	2452617	42.62%	2.786%
60	12.810	1836	1840	1845	rVB4	110200	183363	3.19%	0.208%
61	12.916	1846	1858	1863	rVV	804802	1682718	29.24%	1.911%
62	12.974	1863	1868	1872	rVV7	65161	124257	2.16%	0.141%
63	13.051	1872	1881	1887	rVV5	412960	1110619	19.30%	1.261%
64	13.104	1887	1890	1896	rVB4	161542	247174	4.30%	0.281%
65	13.192	1901	1905	1909	rBV2	171600	316690	5.50%	0.360%
66	13.239	1911	1913	1916	rVV3	78128	91994	1.60%	0.104%
67	13.280	1916	1920	1928	rVB2	470808	727854	12.65%	0.827%
68	13.363	1930	1934	1938	rVB	205196	305445	5.31%	0.347%
69	13.410	1938	1942	1946	rVB	133434	184671	3.21%	0.210%
70	13.463	1947	1951	1953	rBV2	130212	208205	3.62%	0.236%
71	13.557	1964	1967	1971	rVB3	218265	323308	5.62%	0.367%
72	13.604	1971	1975	1977	rBV	253949	308500	5.36%	0.350%
73	13.633	1977	1980	1981	rBV	128568	110397	1.92%	0.125%
74	13.669	1981	1986	1990	rBV	1738891	2944819	51.18%	3.345%
75	13.739	1996	1998	2005	rVB2	328107	411619	7.15%	0.468%
76	13.833	2008	2014	2017	rVV	427424	726542	12.63%	0.825%

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77	13.869	2017	2020	2025	rVV	574711	941947	16.37%	1.070%
78	13.916	2025	2028	2036	rVB3	298978	724690	12.59%	0.823%
79	13.992	2036	2041	2046	rBV3	378743	683156	11.87%	0.776%
80	14.063	2050	2053	2058	rVV	212216	296678	5.16%	0.337%
81	14.210	2073	2078	2083	rBV	1111957	1714161	29.79%	1.947%
82	14.321	2092	2097	2102	rVB2	495760	832761	14.47%	0.946%
83	14.380	2103	2107	2114	rVB5	145193	274995	4.78%	0.312%
84	14.457	2114	2120	2124	rBV3	292184	635008	11.04%	0.721%
85	14.510	2125	2129	2130	rVV3	295117	441218	7.67%	0.501%
86	14.533	2130	2133	2138	rVB	538992	859502	14.94%	0.976%
87	14.616	2143	2147	2151	rVV	359811	500424	8.70%	0.568%
88	14.663	2151	2155	2161	rVB6	137501	234751	4.08%	0.267%
89	14.804	2175	2179	2184	rBV2	315097	577986	10.04%	0.656%
90	14.845	2184	2186	2191	rVB2	154601	211340	3.67%	0.240%
91	14.921	2192	2199	2204	rVV4	710191	1716589	29.83%	1.950%
92	14.968	2204	2207	2215	rVB2	280292	563017	9.78%	0.639%
93	15.080	2223	2226	2231	rBV5	52495	83092	1.44%	0.094%
94	15.192	2240	2245	2257	rVB3	294523	681965	11.85%	0.775%
95	15.310	2258	2265	2272	rVB4	193663	429363	7.46%	0.488%
96	15.463	2285	2291	2293	rBV4	99907	176218	3.06%	0.200%
97	15.504	2293	2298	2309	rVB	311274	623617	10.84%	0.708%
98	15.610	2312	2316	2320	rVB5	54960	93479	1.62%	0.106%
99	15.798	2341	2348	2356	rBV3	97780	231322	4.02%	0.263%
100	16.615	2481	2487	2494	rBV2	44905	103114	1.79%	0.117%

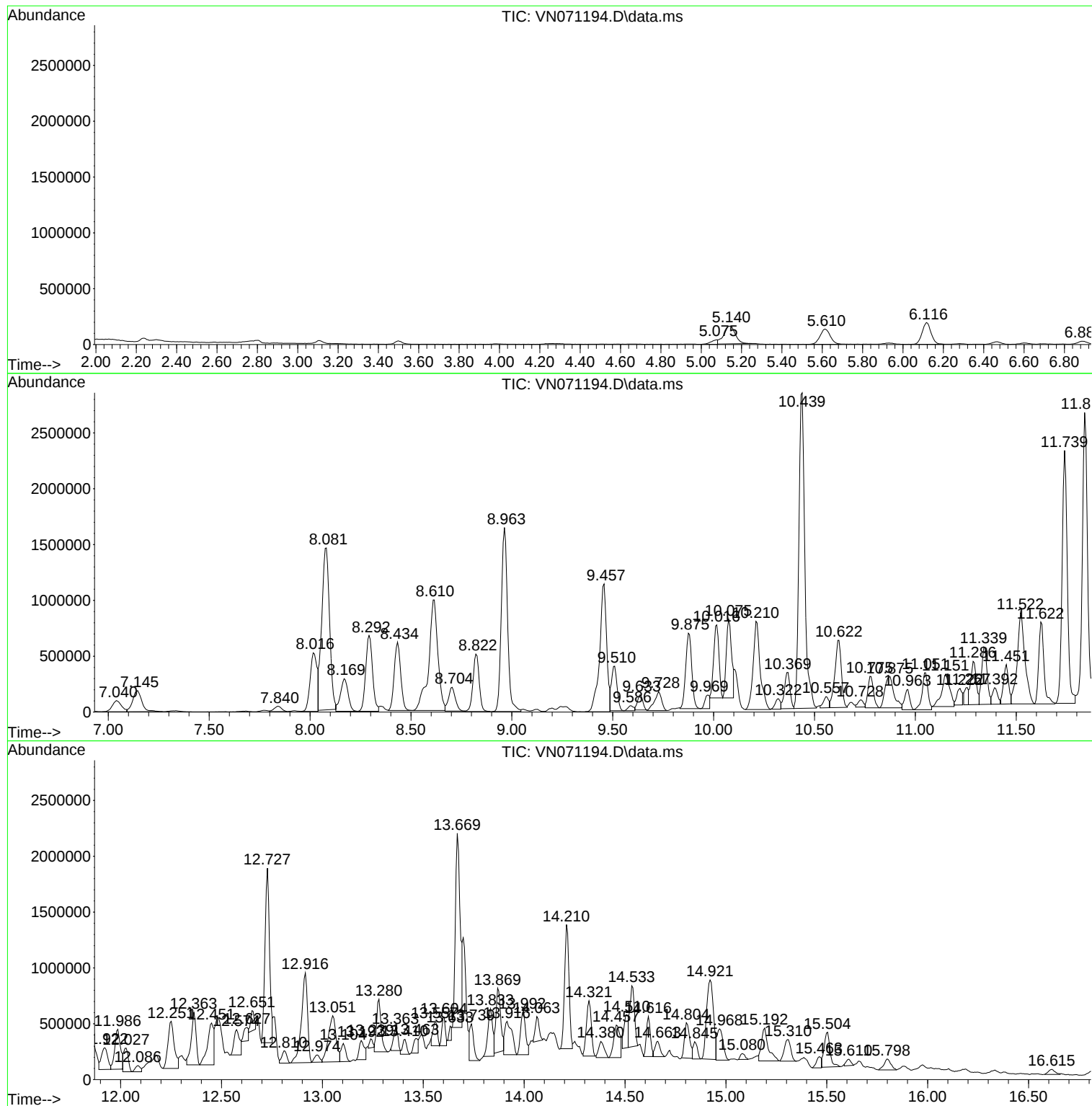
Sum of corrected areas: 88046742

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



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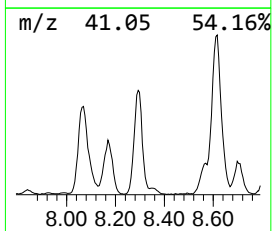
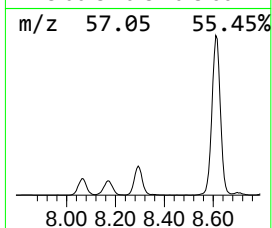
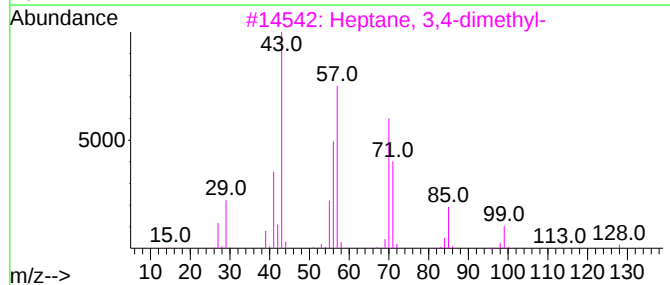
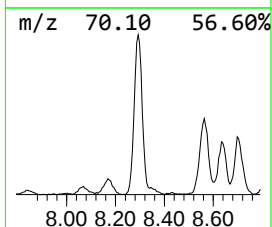
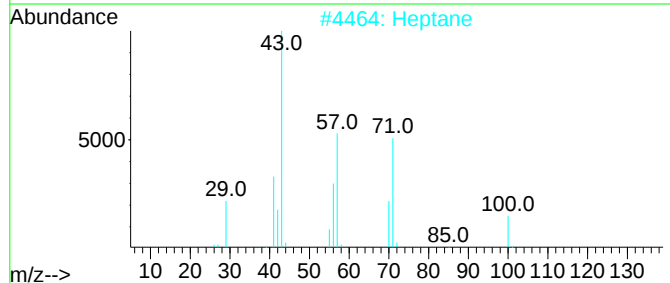
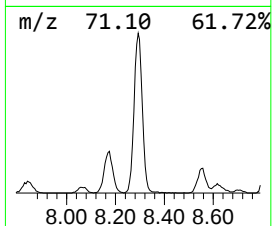
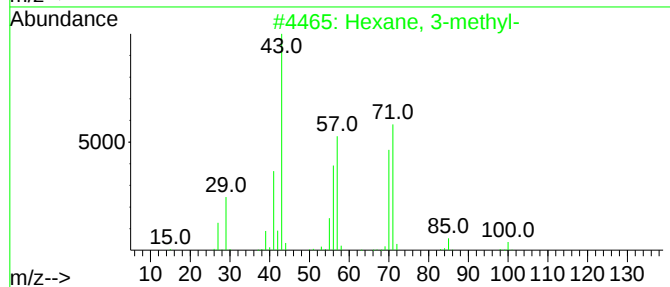
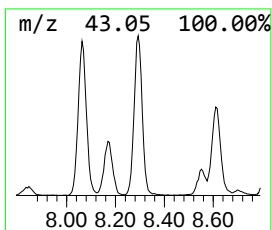
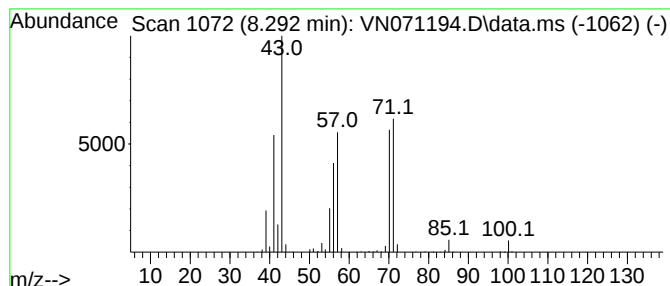
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.292	20.53 ug/l	1596580	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	64
4			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53
5			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	53



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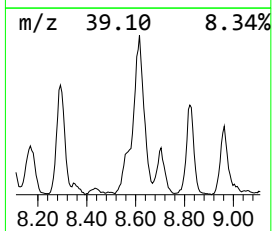
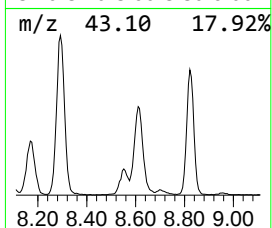
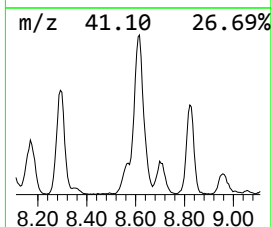
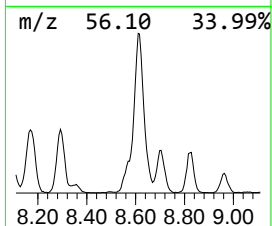
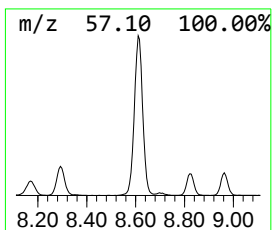
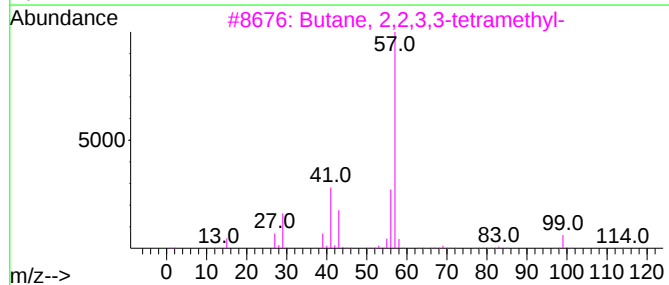
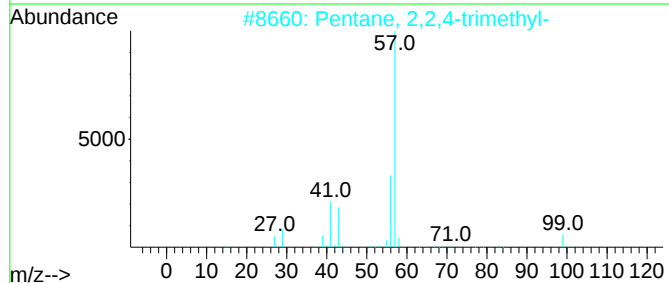
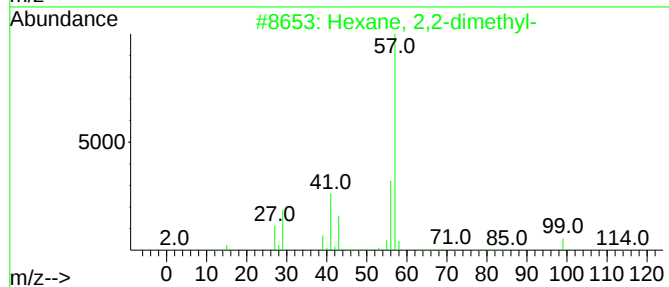
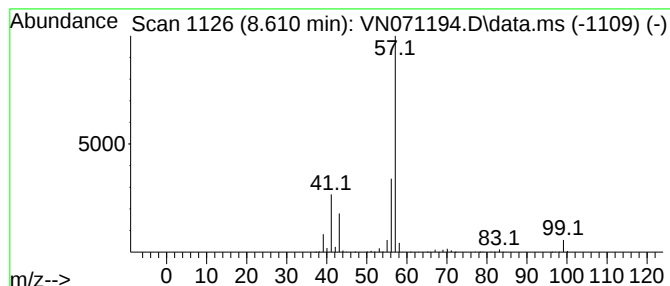
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Hexane, 2,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.610	44.40 ug/l	3141470	1,4-Difluorobenzene	8.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
4			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59



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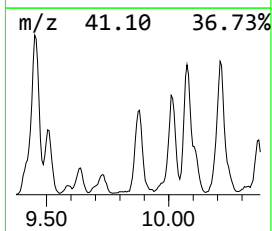
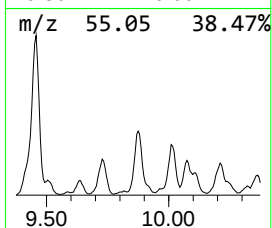
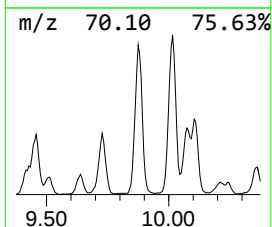
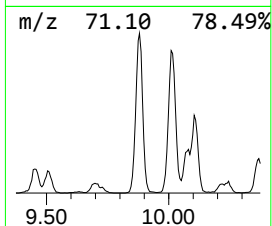
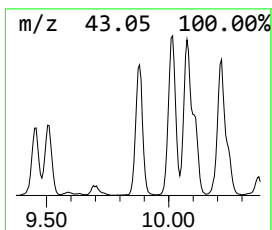
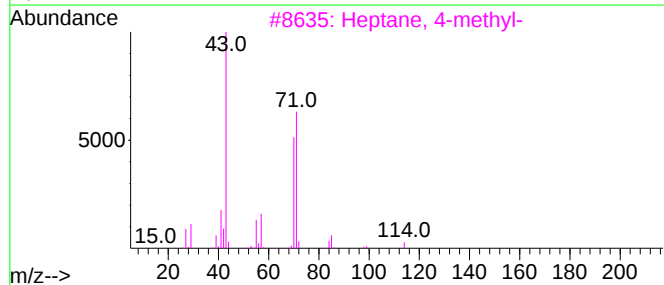
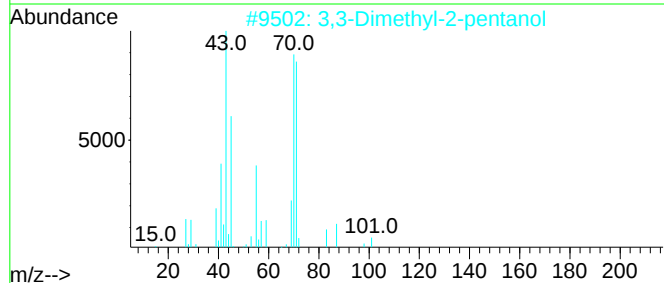
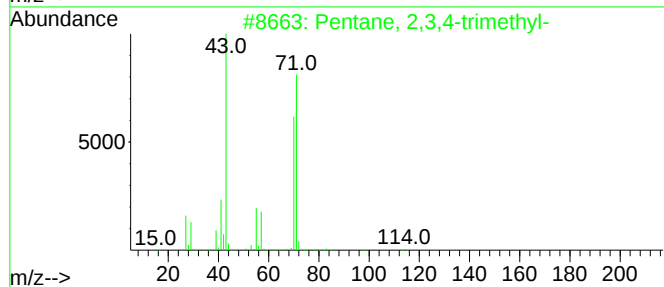
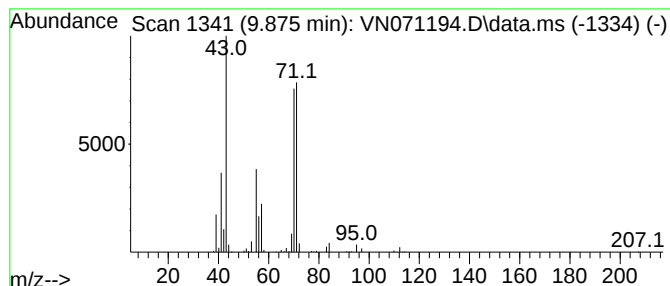
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.875	19.85 ug/l	1404740	1,4-Difluorobenzene	8.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	86
2			3,3-Dimethyl-2-pentanol	116	C7H16O	019781-24-9	78
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	78
4			Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

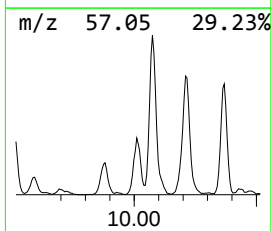
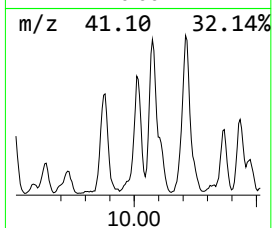
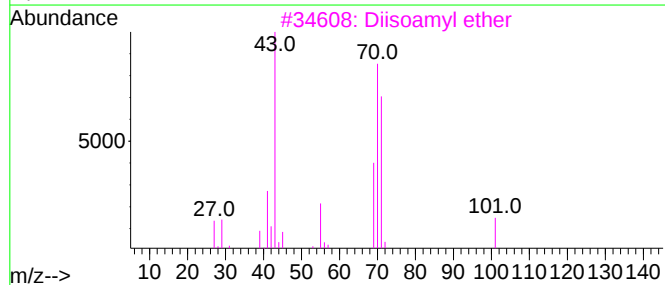
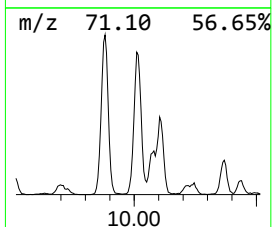
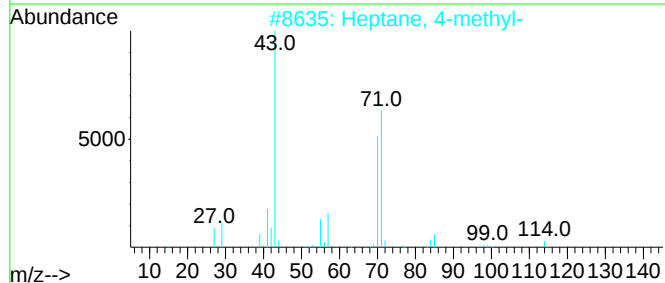
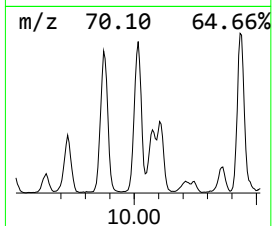
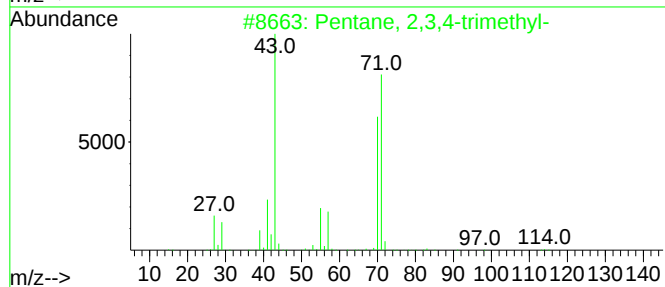
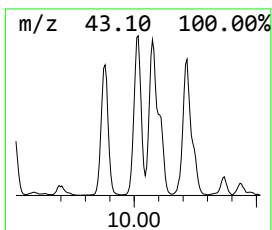
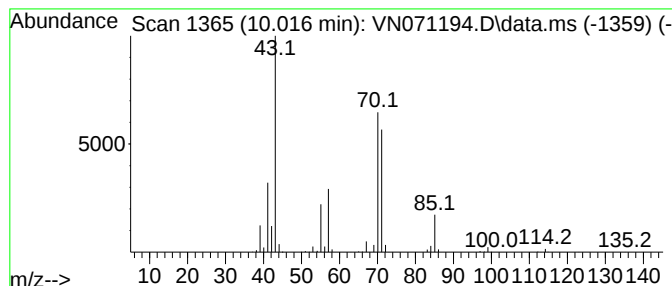
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Quant Title : SW846 8260

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

Peak Number 4 Heptane, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.016	16.82 ug/l	1190020	1,4-Difluorobenzene	8.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	72
3			Diisoamyl ether	158	C10H22O	000544-01-4	72
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	72
5			3,4-Diethyl hexane	142	C10H22	019398-77-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

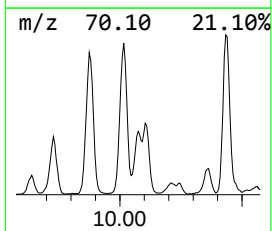
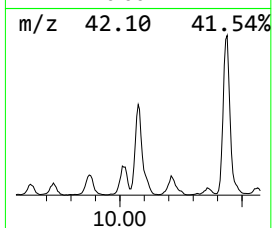
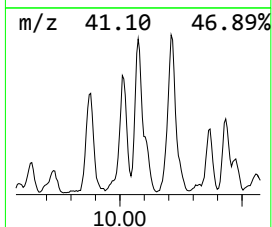
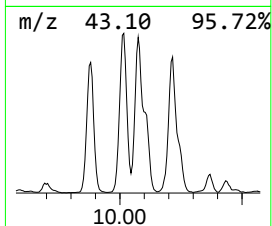
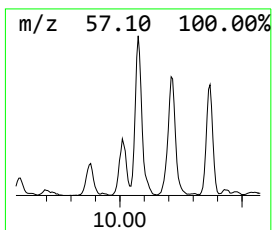
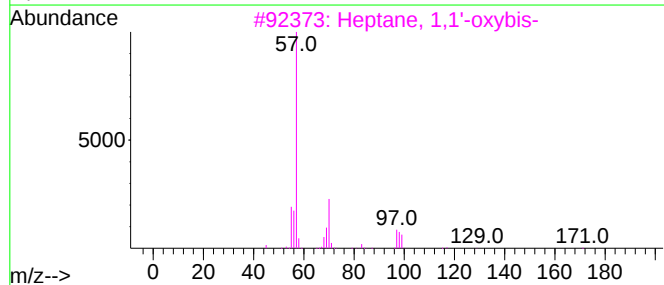
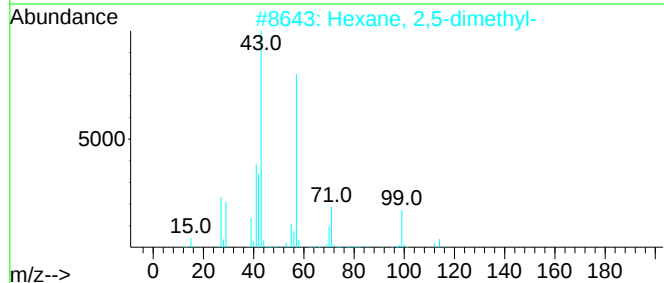
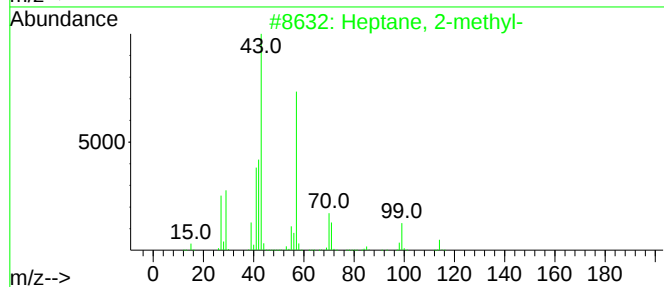
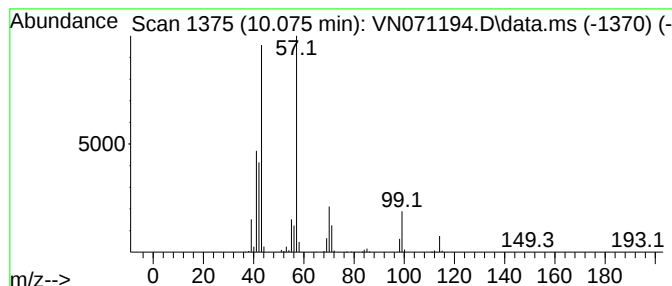
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Quant Title : SW846 8260

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

Peak Number 5 Heptane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.075	18.10 ug/l	1280770	1,4-Difluorobenzene	8.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2-methyl-	114	C8H18	000592-27-8	91
2			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	72
3			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	59
4			Piperazine, 1,4-dimethyl-	114	C6H14N2	000106-58-1	38
5			Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

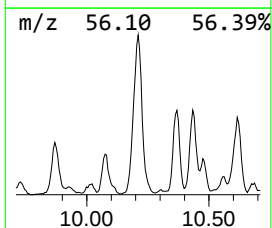
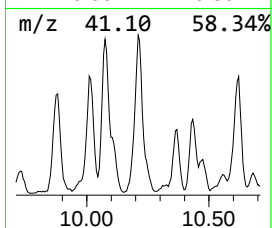
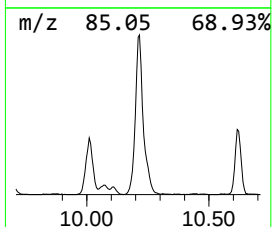
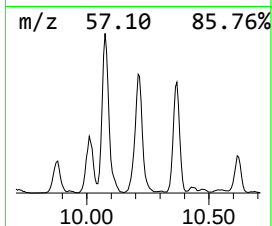
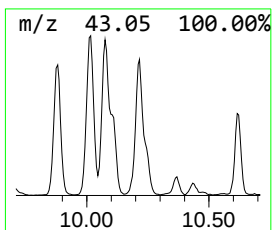
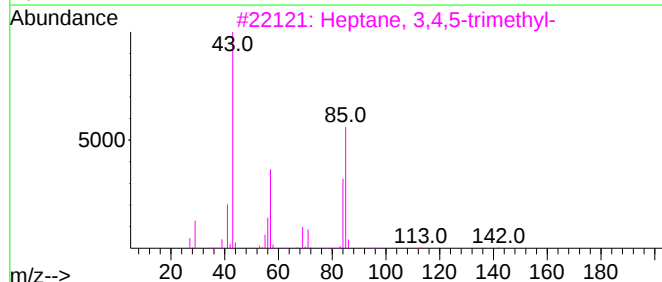
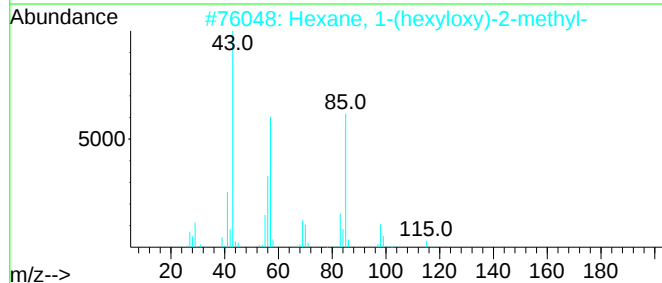
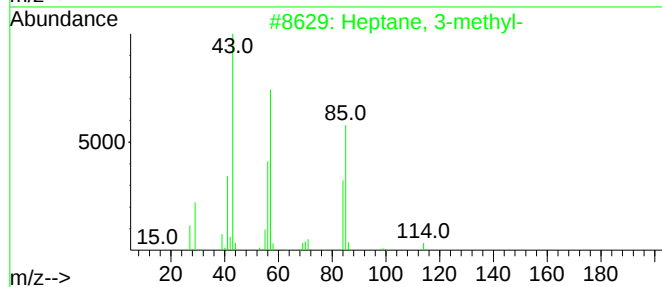
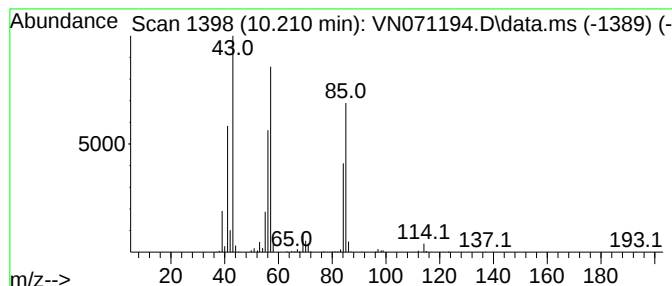
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Quant Title : SW846 8260

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

Peak Number 6 Heptane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.210	26.91 ug/l	1903940	1,4-Difluorobenzene	8.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2			Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	64
3			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	59
4			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	53
5			Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

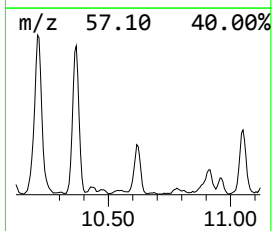
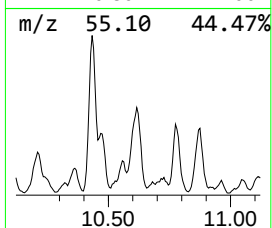
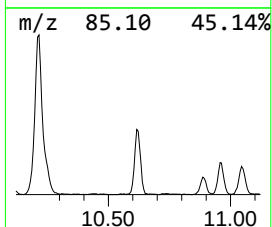
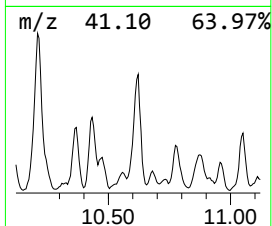
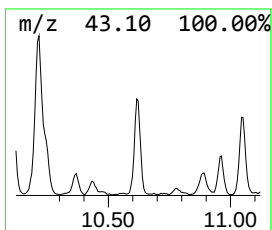
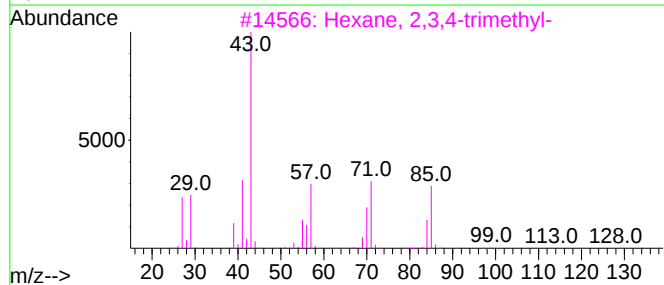
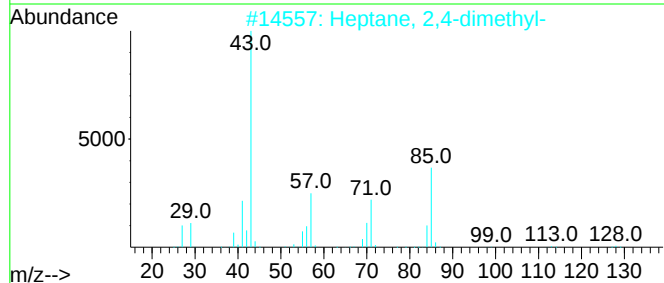
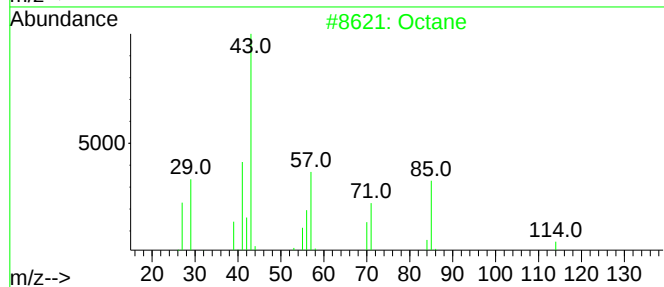
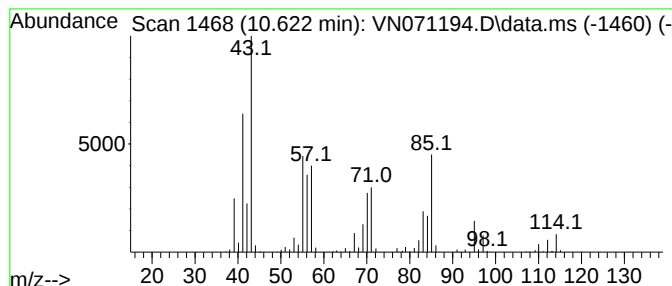
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Quant Title : SW846 8260

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

Peak Number 7 Octane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.622	16.49 ug/l	1344350	Chlorobenzene-d5	11.739

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane			114	C8H18	000111-65-9	83
2	Heptane, 2,4-dimethyl-			128	C9H20	002213-23-2	47
3	Hexane, 2,3,4-trimethyl-			128	C9H20	000921-47-1	43
4	1-Pentanol, 2-methyl-			102	C6H14O	000105-30-6	38
5	Acetic acid, decyl ester			200	C12H24O2	000112-17-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

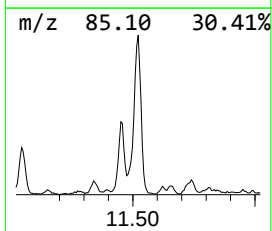
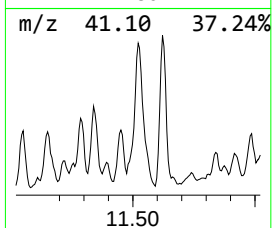
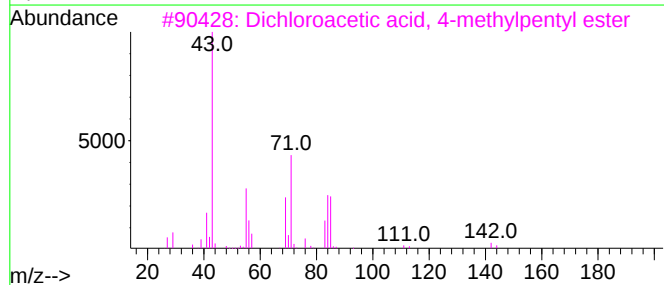
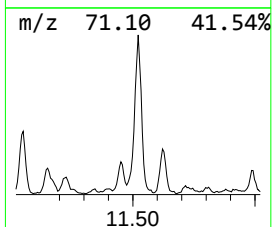
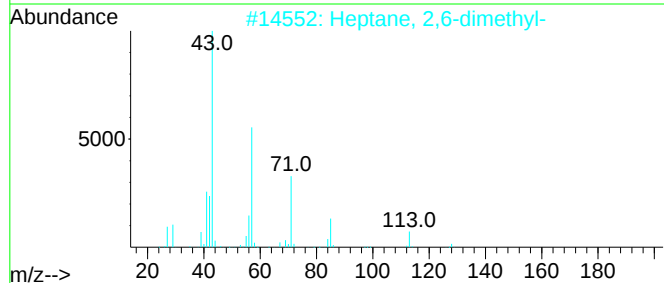
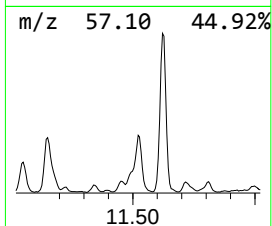
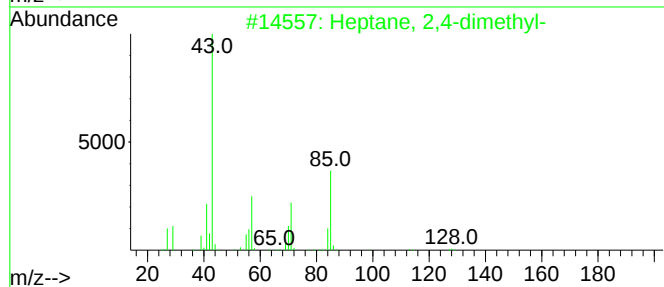
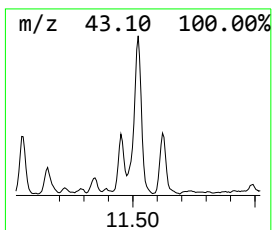
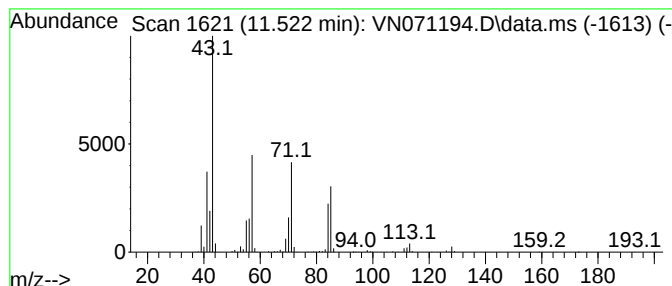
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Quant Title : SW846 8260

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

Peak Number 8 Heptane, 2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.522	26.39 ug/l	2151590	Chlorobenzene-d5	11.739

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
2			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	59
3			Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	59
4			Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	53
5			Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

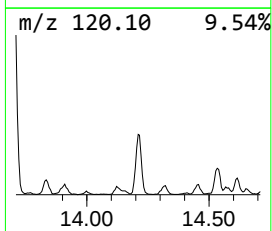
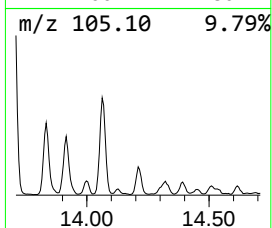
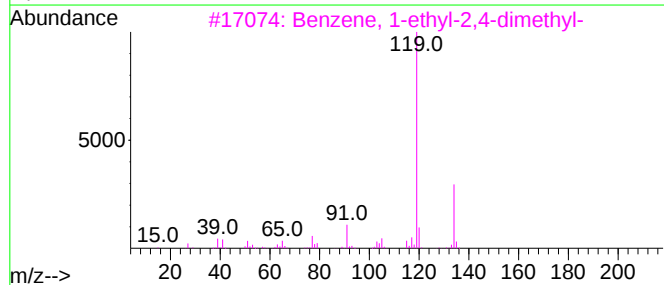
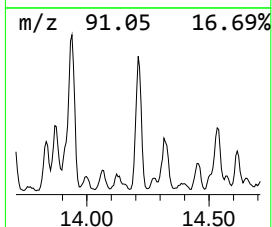
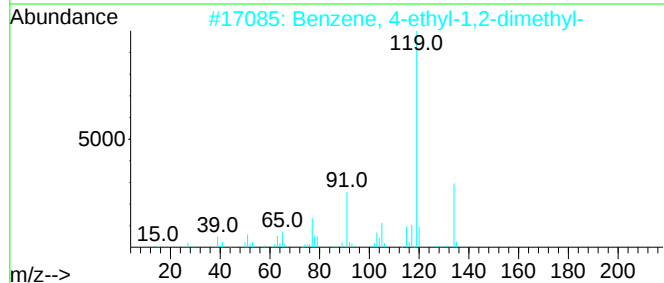
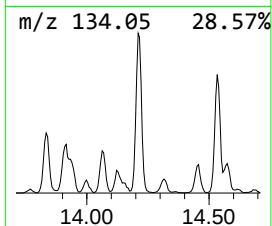
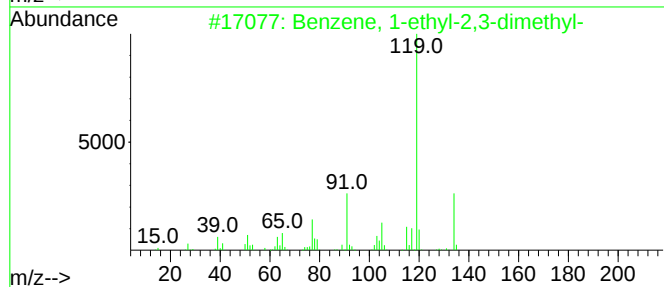
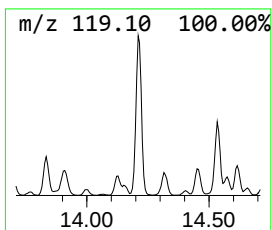
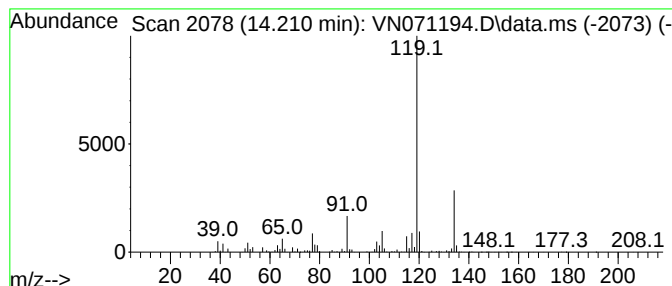
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.210	29.10 ug/l	1714160	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			p-Cymene	134	C10H14	000099-87-6	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P-1

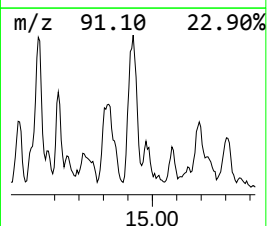
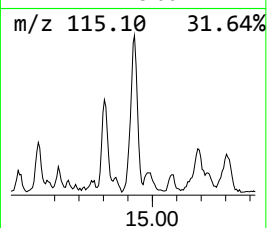
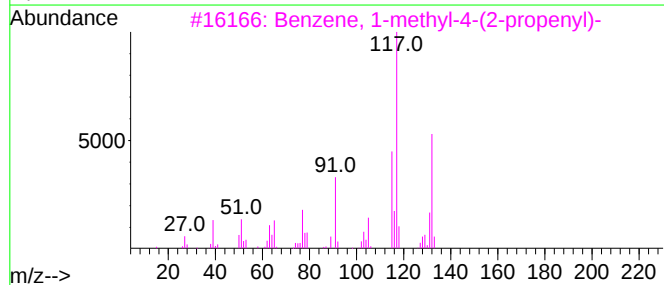
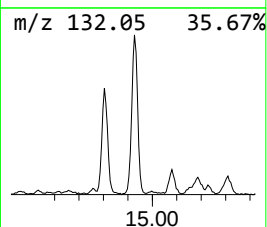
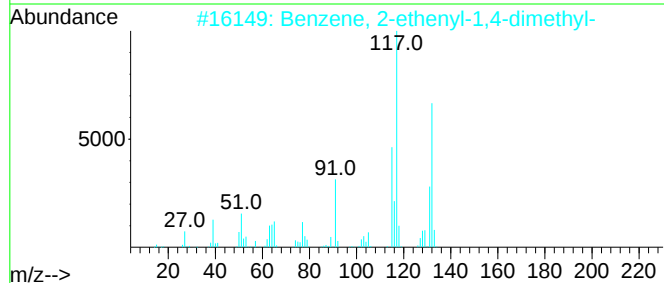
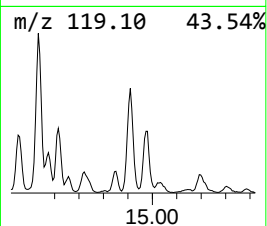
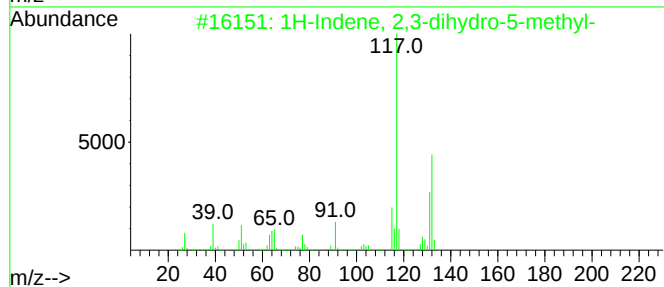
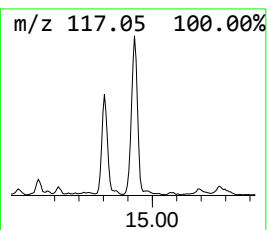
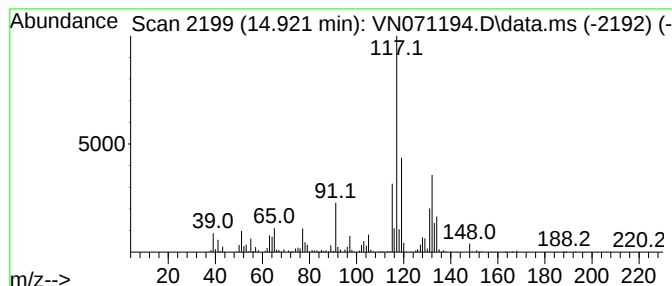
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.921	29.15 ug/l	1716590	1,4-Dichlorobenzene-d4	13.669

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	86
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	86
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030822\
Data File : VN071194.D
Acq On : 08 Mar 2022 16:53
Operator : JC\MD
Sample : N1815-01 10X
Misc : 4.47g/10.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.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
Hexane, 3-methyl-	8.292	20.5	ug/l	1596580	1	8.081	3889150	50.0
Hexane, 2,2-dim...	8.610	44.4	ug/l	3141470	2	8.963	3537700	50.0
Pentane, 2,3,4-...	9.875	19.9	ug/l	1404740	2	8.963	3537700	50.0
Heptane, 4-methyl-	10.016	16.8	ug/l	1190020	2	8.963	3537700	50.0
Heptane, 2-methyl-	10.075	18.1	ug/l	1280770	2	8.963	3537700	50.0
Heptane, 3-methyl-	10.210	26.9	ug/l	1903940	2	8.963	3537700	50.0
Octane	10.622	16.5	ug/l	1344350	3	11.739	4077100	50.0
Heptane, 2,4-di...	11.522	26.4	ug/l	2151590	3	11.739	4077100	50.0
Benzene, 1-ethy...	14.210	29.1	ug/l	1714160	4	13.669	2944820	50.0
1H-Indene, 2,3-...	14.921	29.1	ug/l	1716590	4	13.669	2944820	50.0