

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
 Data File : VN073114.D
 Acq On : 25 Jun 2022 16:43
 Operator : JC\MD
 Sample : N3454-07
 Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-14A-1A-1-GR

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: VN073114.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.310	233	242	255	rVB	129502	312286	1.54%	0.094%
2	4.981	518	526	553	rVB3	223054	967050	4.76%	0.291%
3	5.469	594	609	624	rBV	217866	754657	3.72%	0.227%
4	5.981	685	696	714	rVB	187437	576111	2.84%	0.173%
5	7.034	864	875	893	rVB2	390115	1203679	5.93%	0.362%
6	8.004	1016	1040	1048	rBV5	1268373	4254389	20.94%	1.280%
7	8.080	1048	1053	1066	rVV	229643	602408	2.97%	0.181%
8	8.210	1066	1075	1081	rVV	561373	1313350	6.47%	0.395%
9	8.269	1081	1085	1093	rVV2	182510	421201	2.07%	0.127%
10	8.363	1093	1101	1111	rVV2	260023	642290	3.16%	0.193%
11	8.486	1112	1122	1128	rVV2	451321	1128123	5.55%	0.339%
12	8.557	1128	1134	1139	rVV2	382723	871216	4.29%	0.262%
13	8.622	1139	1145	1157	rVV	641577	1475218	7.26%	0.444%
14	8.745	1159	1166	1176	rVB	243815	496449	2.44%	0.149%
15	8.898	1181	1192	1205	rBV	645218	1480606	7.29%	0.445%
16	9.386	1260	1275	1281	rBV	3076963	7606264	37.45%	2.288%
17	9.433	1281	1283	1294	rVB	235255	423098	2.08%	0.127%
18	9.569	1299	1306	1312	rVV	261592	511324	2.52%	0.154%
19	9.657	1312	1321	1329	rVB	462192	1028285	5.06%	0.309%
20	9.804	1339	1346	1354	rVB2	588662	1183432	5.83%	0.356%
21	9.951	1366	1371	1375	rBV	199024	336060	1.65%	0.101%
22	10.010	1375	1381	1385	rBV	672697	1307537	6.44%	0.393%
23	10.145	1394	1404	1416	rBV3	716605	1993895	9.82%	0.600%
24	10.257	1417	1423	1427	rBV2	148208	320782	1.58%	0.096%
25	10.369	1434	1442	1447	rBV2	2454988	5049590	24.86%	1.519%
26	10.551	1466	1473	1481	rVB4	808084	2163863	10.65%	0.651%
27	10.674	1483	1494	1496	rBV4	253807	554700	2.73%	0.167%
28	10.716	1496	1501	1508	rVB	1151538	2088936	10.28%	0.628%
29	10.810	1509	1517	1522	rBV3	652718	1467775	7.23%	0.442%
30	10.898	1528	1532	1537	rVB	345463	547771	2.70%	0.165%
31	10.986	1537	1547	1554	rBV	974035	1777183	8.75%	0.535%
32	11.086	1554	1564	1572	rBV	582847	1620641	7.98%	0.488%
33	11.157	1572	1576	1579	rVV2	404089	667728	3.29%	0.201%
34	11.227	1579	1588	1592	rVV	1641449	3237438	15.94%	0.974%
35	11.280	1592	1597	1603	rVV	2456346	4396038	21.64%	1.322%
36	11.333	1603	1606	1610	rVB2	302767	395235	1.95%	0.119%

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Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
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37	11.392	1610	1616	1620	rBV2	1027352	1804694	8.88%	0.543%
38	11.480	1620	1631	1640	rVB5	1113076	3911006	19.25%	1.177%
39	11.563	1640	1645	1650	rBV	1309815	2185671	10.76%	0.657%
40	11.686	1659	1666	1670	rBV	881305	1632812	8.04%	0.491%
41	11.774	1676	1681	1684	rBV2	405984	696629	3.43%	0.210%
42	11.816	1684	1688	1691	rVV	421902	509589	2.51%	0.153%
43	11.880	1691	1699	1703	rBV5	998049	3087978	15.20%	0.929%
44	11.927	1703	1707	1711	rVV	1082374	1664510	8.19%	0.501%
45	11.968	1711	1714	1719	rVB2	1319222	2075956	10.22%	0.624%
46	12.027	1719	1724	1729	rBV2	393388	808307	3.98%	0.243%
47	12.116	1731	1739	1744	rVB4	543499	1295907	6.38%	0.390%
48	12.192	1744	1752	1755	rBV4	2459397	5679891	27.96%	1.709%
49	12.304	1763	1771	1776	rVB2	2121968	3694474	18.19%	1.111%
50	12.357	1776	1780	1781	rBV3	637512	887862	4.37%	0.267%
51	12.392	1781	1786	1788	rVV2	2120831	3577363	17.61%	1.076%
52	12.421	1788	1791	1802	rVB4	2454414	5765966	28.39%	1.735%
53	12.516	1802	1807	1811	rBV3	1081144	2045038	10.07%	0.615%
54	12.557	1812	1814	1818	rVV5	344456	493277	2.43%	0.148%
55	12.668	1829	1833	1837	rBV3	1327487	2220123	10.93%	0.668%
56	12.704	1837	1839	1843	rVV2	544824	823423	4.05%	0.248%
57	12.751	1843	1847	1852	rVV4	803453	1690176	8.32%	0.508%
58	12.833	1856	1861	1869	rVV2	3216474	7419386	36.53%	2.232%
59	12.921	1869	1876	1879	rVV4	2260721	4488352	22.10%	1.350%
60	12.998	1879	1889	1894	rVV5	5056208	13565246	66.78%	4.081%
61	13.045	1894	1897	1902	rVB3	1367558	2233188	10.99%	0.672%
62	13.157	1908	1916	1923	rBV3	2391821	6720810	33.09%	2.022%
63	13.221	1923	1927	1935	rVB3	2672844	4168718	20.52%	1.254%
64	13.304	1935	1941	1947	rBV	8380464	13888601	68.37%	4.178%
65	13.439	1955	1964	1968	rVB4	1395233	2906881	14.31%	0.874%
66	13.498	1968	1974	1979	rBV4	3387329	6507195	32.03%	1.957%
67	13.545	1979	1982	1990	rVV3	1652376	2923500	14.39%	0.879%
68	13.639	1994	1998	2003	rVB	4436185	7151083	35.20%	2.151%
69	13.774	2014	2021	2023	rBV2	1488672	3227489	15.89%	0.971%
70	13.804	2023	2026	2030	rVV2	3063003	4827401	23.76%	1.452%
71	13.851	2030	2034	2043	rVB3	2633383	5864678	28.87%	1.764%
72	13.933	2043	2048	2053	rBV2	3625054	6158278	30.32%	1.853%
73	14.010	2057	2061	2065	rVV	1612301	3088388	15.20%	0.929%
74	14.068	2067	2071	2073	rVV	2424033	3810867	18.76%	1.146%
75	14.098	2073	2076	2081	rVV	2747948	4631334	22.80%	1.393%
76	14.151	2081	2085	2090	rVB	3495217	5282215	26.00%	1.589%

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77	14.262	2098	2104	2109	rVB3	4462681	7705793	37.94%	2.318%
78	14.333	2109	2116	2121	rBV6	1145442	3008588	14.81%	0.905%
79	14.398	2121	2127	2130	rBV3	1679348	3267903	16.09%	0.983%
80	14.445	2130	2135	2138	rBV2	2367455	3779382	18.61%	1.137%
81	14.515	2144	2147	2151	rVB	2752295	3578029	17.61%	1.076%
82	14.557	2151	2154	2158	rBV	2139051	2568607	12.65%	0.773%
83	14.604	2158	2162	2169	rVB4	2026124	3555707	17.50%	1.070%
84	14.662	2169	2172	2173	rBV3	483243	527467	2.60%	0.159%
85	14.698	2175	2178	2182	rVB	1868370	2676169	13.17%	0.805%
86	14.762	2182	2189	2190	rBV3	1254688	2555874	12.58%	0.769%
87	14.786	2191	2193	2198	rVB	2041108	2741832	13.50%	0.825%
88	14.851	2198	2204	2211	rBV3	6203218	13840972	68.14%	4.164%
89	14.915	2211	2215	2220	rVB2	1826638	3333634	16.41%	1.003%
90	15.127	2247	2251	2255	rBV3	1961128	3105532	15.29%	0.934%
91	15.239	2264	2270	2276	rVB4	2576952	5613049	27.63%	1.689%
92	15.398	2291	2297	2299	rBV4	1291832	2659276	13.09%	0.800%
93	15.427	2299	2302	2308	rVB	2178182	3523158	17.34%	1.060%
94	15.539	2316	2321	2325	rBV	1664633	2825689	13.91%	0.850%
95	15.721	2346	2352	2360	rBV7	1213151	3817609	18.79%	1.148%
96	15.809	2362	2367	2371	rBV4	991093	1881960	9.26%	0.566%
97	16.104	2410	2417	2423	rBV5	923784	2914199	14.35%	0.877%
98	16.251	2434	2442	2449	rBV6	2037360	5598385	27.56%	1.684%
99	16.527	2482	2489	2508	rVB2	5499861	16443082	80.95%	4.946%
100	16.733	2516	2524	2533	rVB	8523281	20313122	100.00%	6.111%

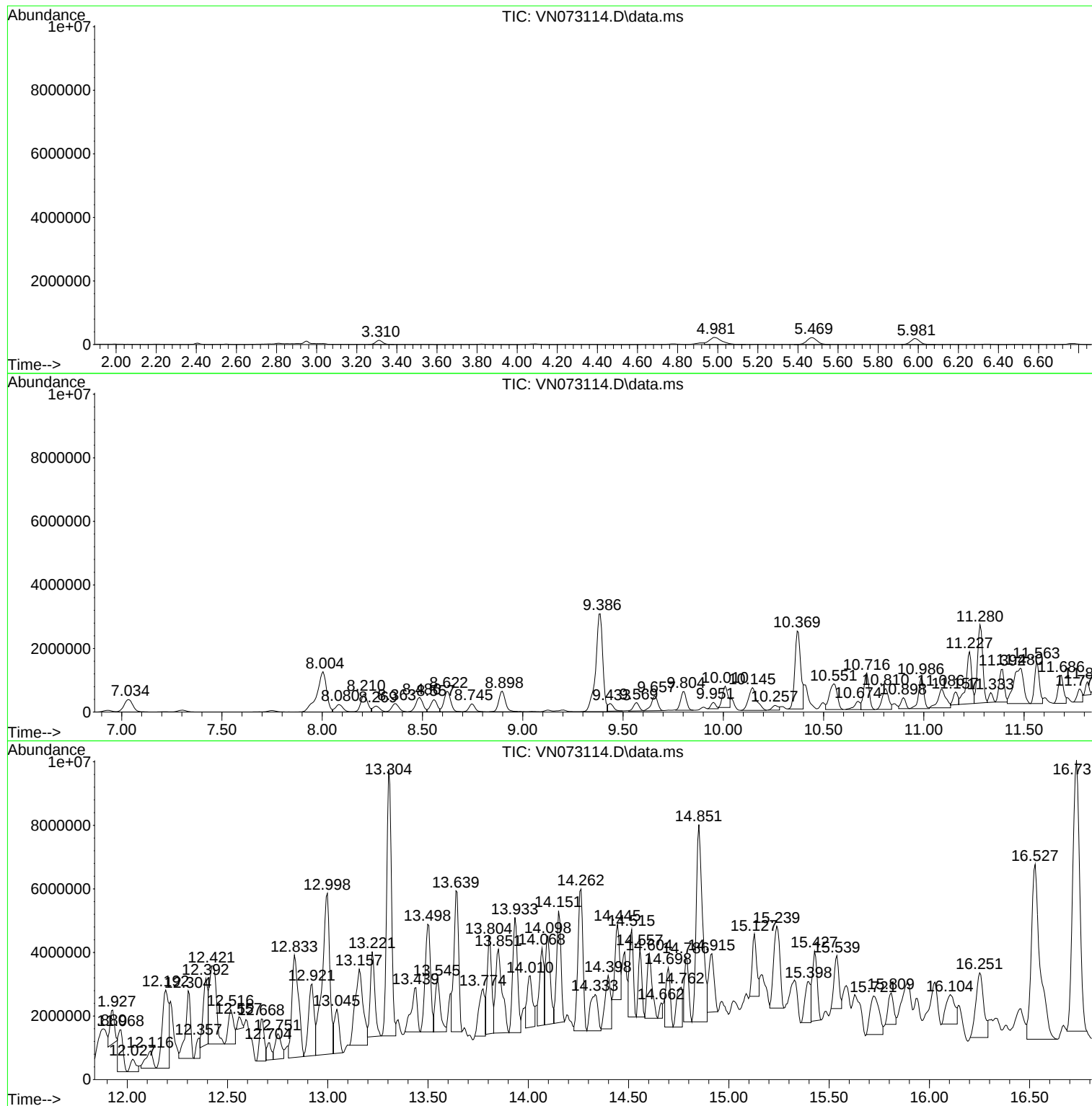
Sum of corrected areas: 332425888

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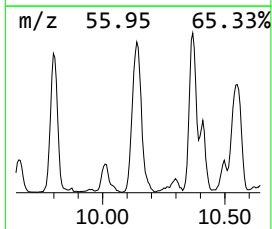
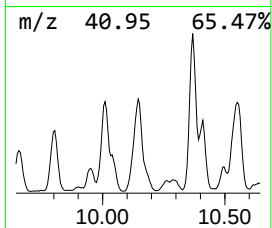
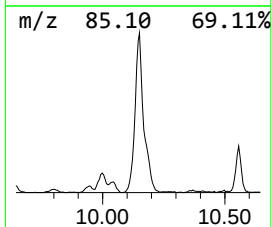
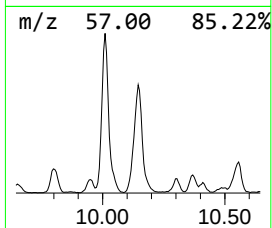
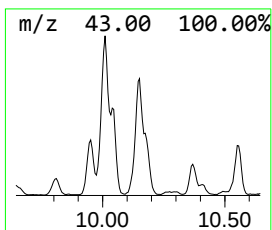
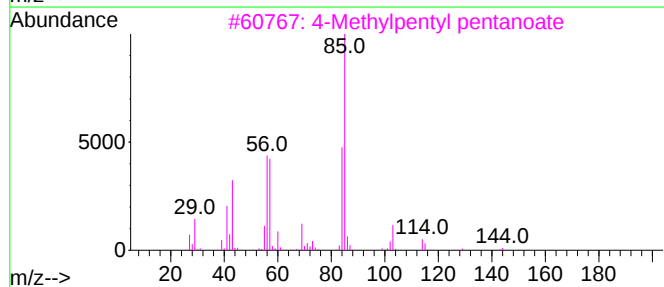
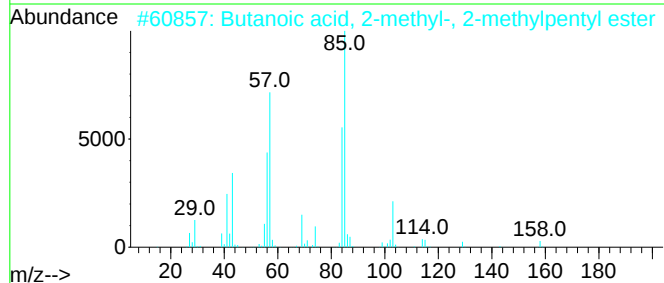
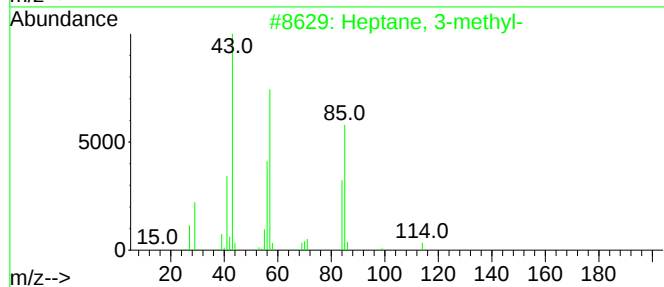
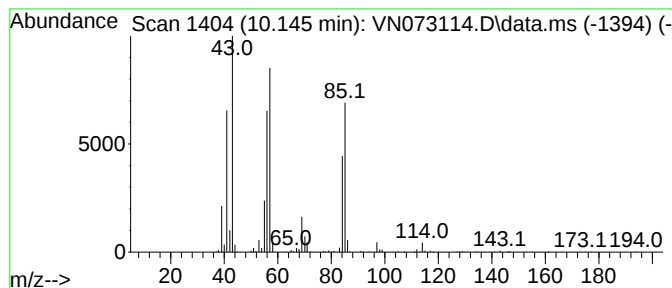
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.145	67.33 ug/l	1993900	1,4-Difluorobenzene	8.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	83
2			Butanoic acid, 2-methyl-, 2-meth...	186	C11H22O2	083783-88-4	64
3			4-Methylpentyl pentanoate	186	C11H22O2	035852-47-2	64
4			Heptane, 4-ethyl-	128	C9H20	002216-32-2	59
5			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	59



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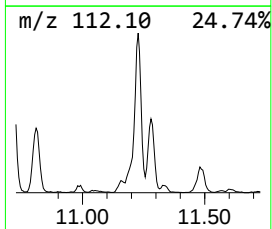
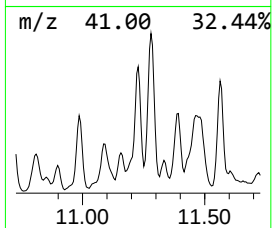
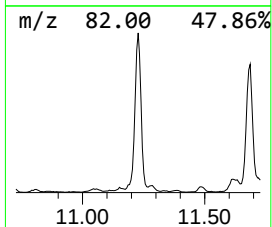
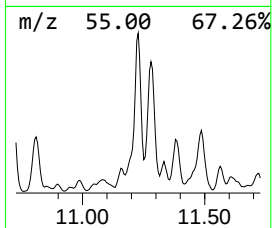
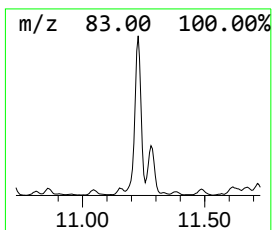
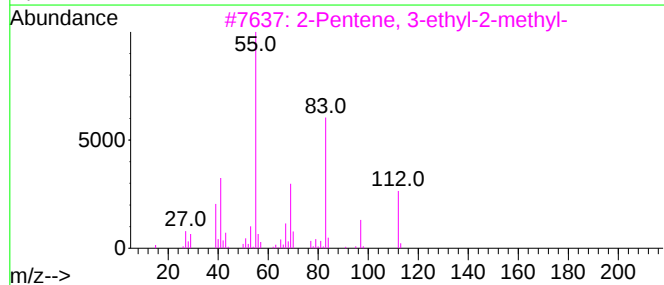
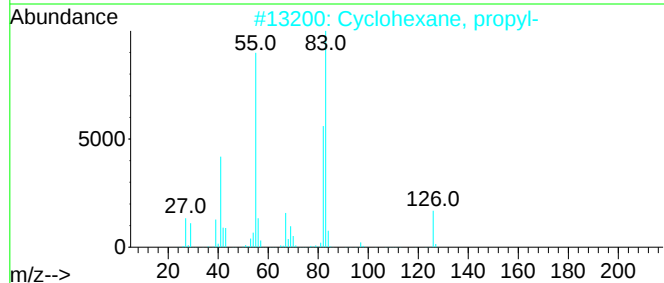
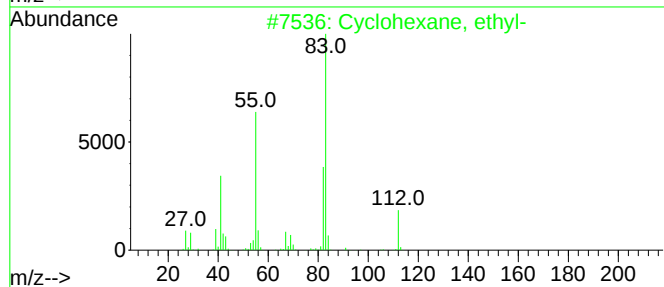
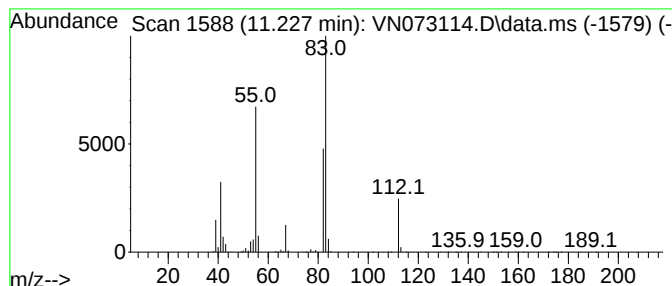
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.227	99.14 ug/l	3237440	Chlorobenzene-d5	11.686

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	91
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
3			2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	59
4			2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	58
5			3-Ethyl-3-hexene	112	C8H16	016789-51-8	53



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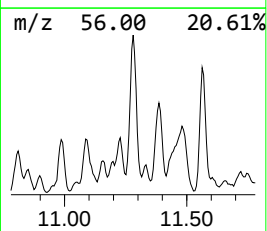
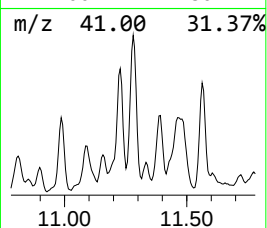
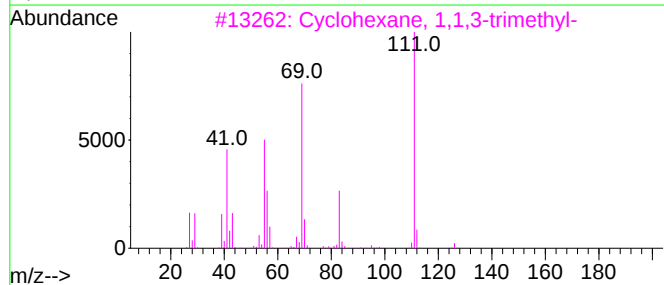
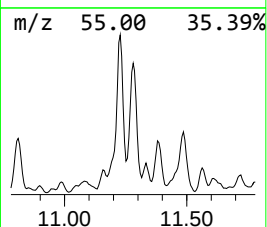
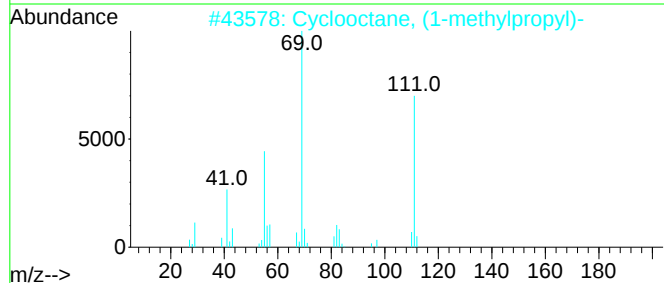
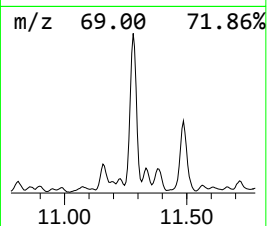
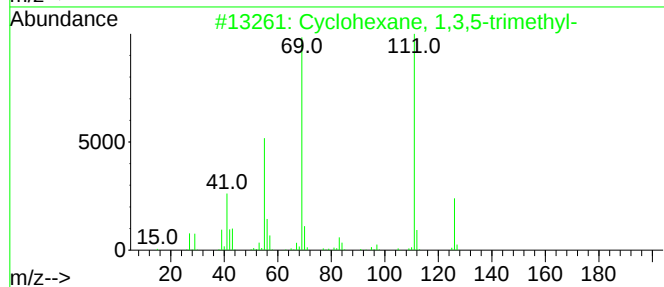
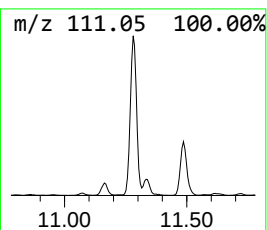
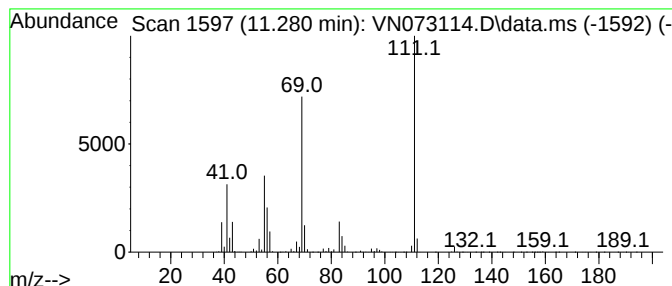
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1,3,5-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.280	134.62 ug/l	4396040	Chlorobenzene-d5	11.686

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	64
2			Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	64
3			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	62
4			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	59
5			1-Aminocyclopropanecarboxylic ac...	156	C7H12N2O2	1000375-50-8	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14A-1A-1-GR

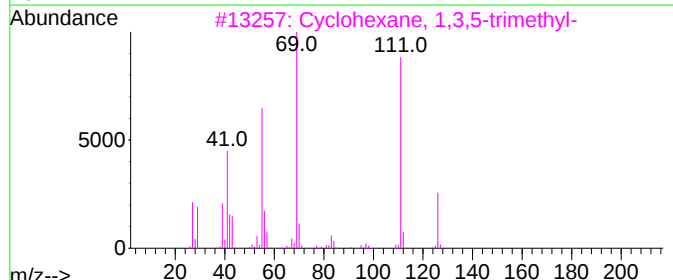
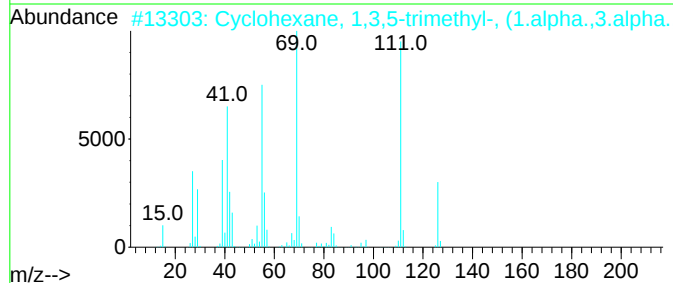
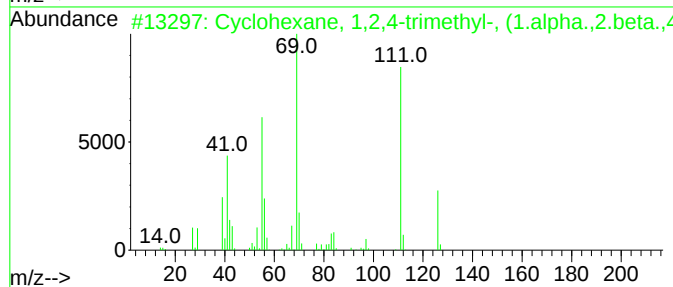
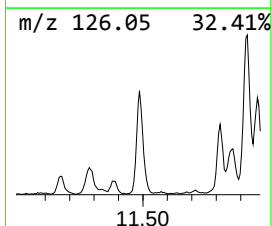
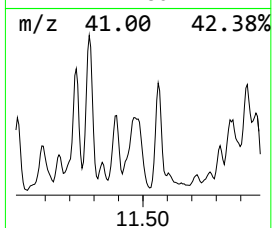
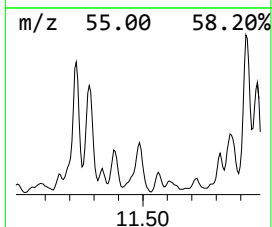
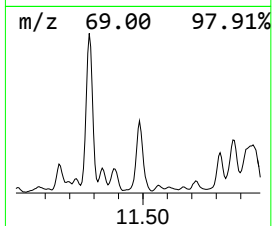
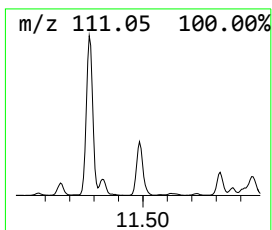
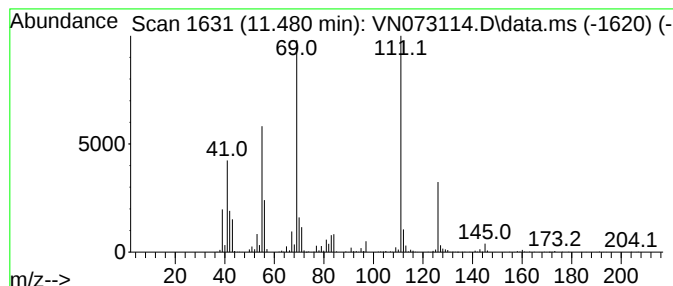
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, 1,2,4-trimethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.480	119.76 ug/l	3911010	Chlorobenzene-d5	11.686

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	95
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	93
3			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	90
4			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	87
5			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14A-1A-1-GR

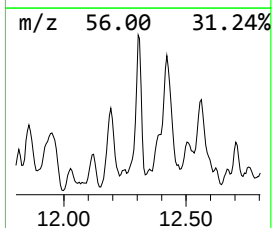
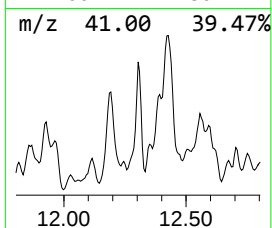
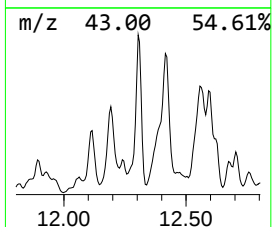
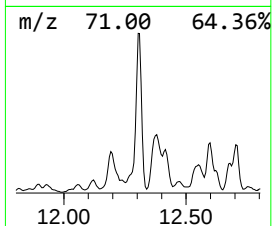
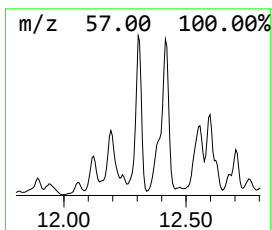
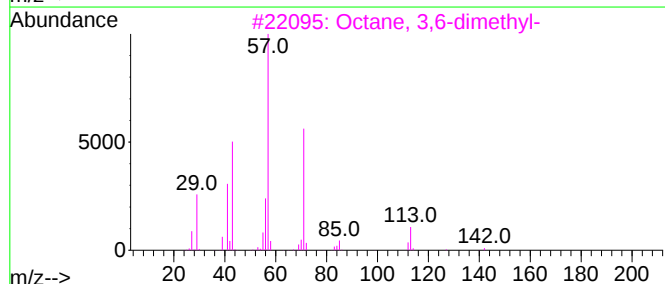
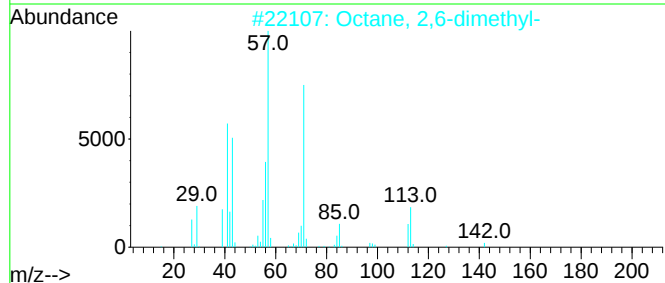
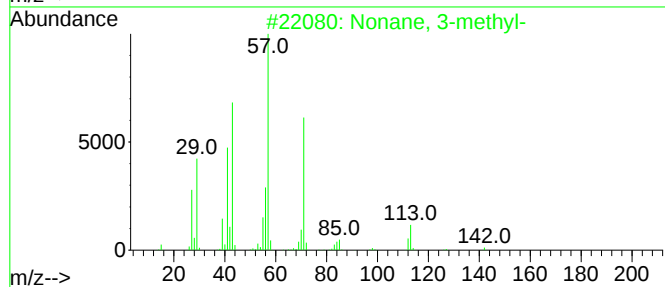
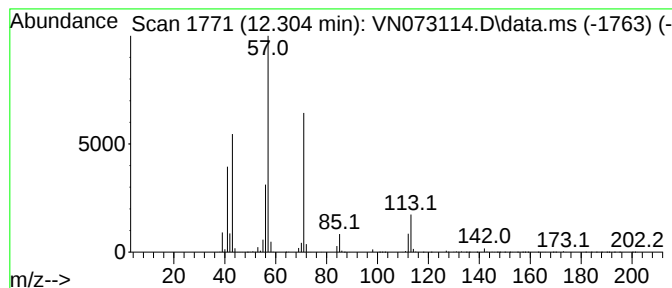
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Quant Title : SW846 8260

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

Peak Number 5 Nonane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.304	113.13 ug/l	3694470	Chlorobenzene-d5	11.686

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	91
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	91
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	87
4			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	72
5			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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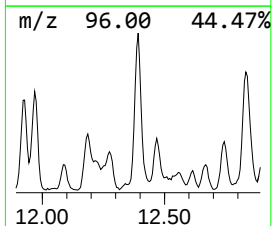
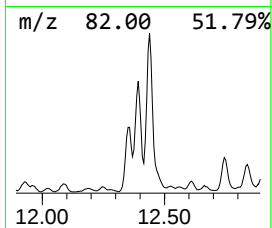
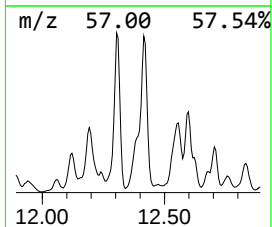
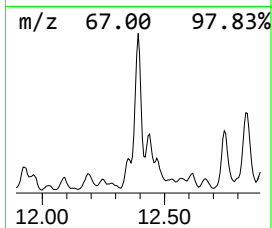
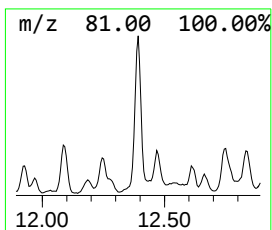
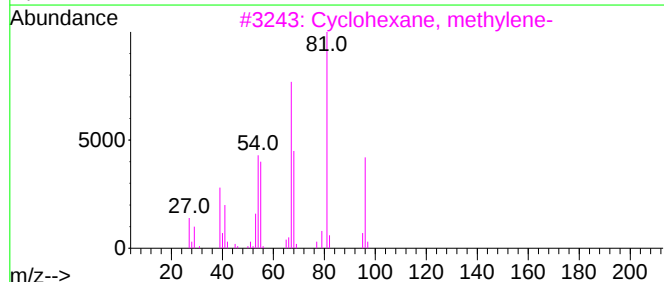
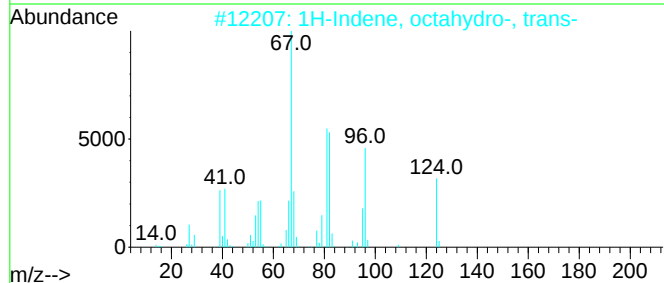
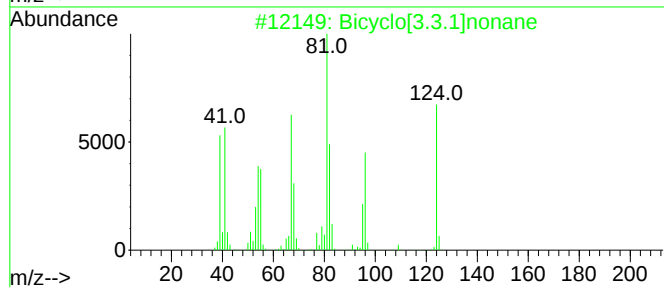
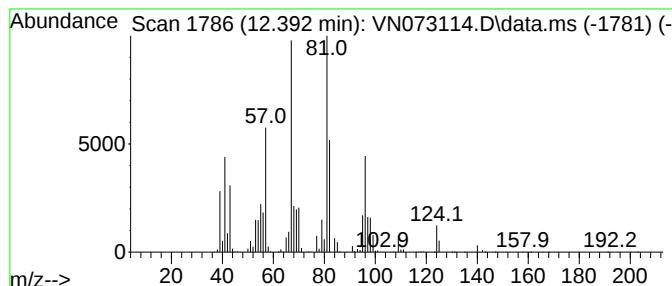
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Quant Title : SW846 8260

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

Peak Number 6 Bicyclo[3.3.1]nonane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.392	109.55 ug/l	3577360	Chlorobenzene-d5	11.686

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.3.1]nonane	124	C9H16	000280-65-9	58
2			1H-Indene, octahydro-, trans-	124	C9H16	003296-50-2	58
3			Cyclohexane, methylene-	96	C7H12	001192-37-6	52
4			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	50
5			1-Methylcyclooctene	124	C9H16	000933-11-9	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14A-1A-1-GR

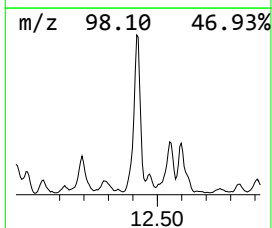
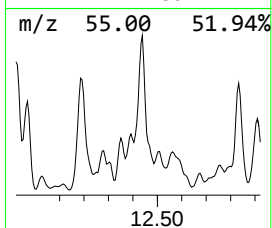
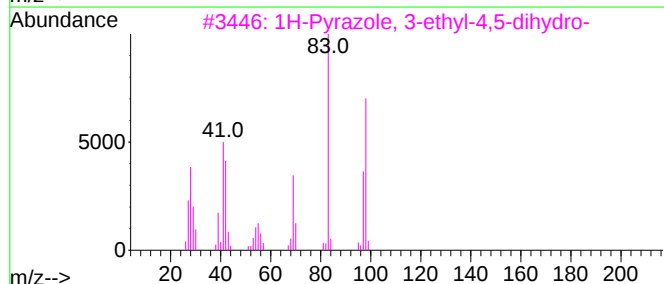
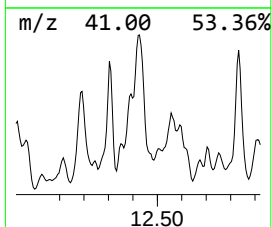
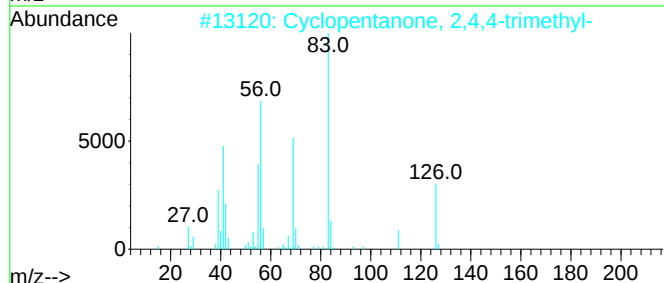
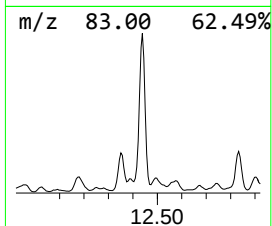
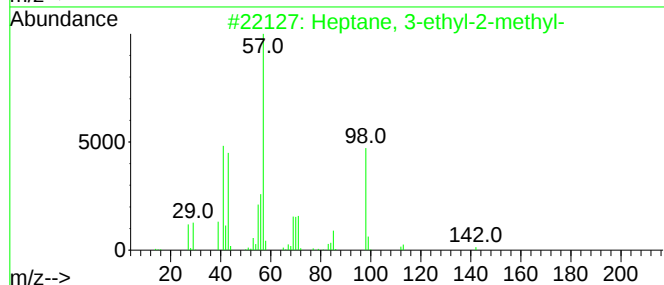
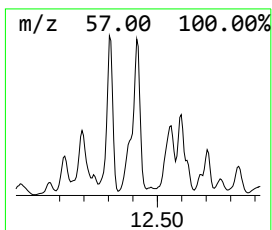
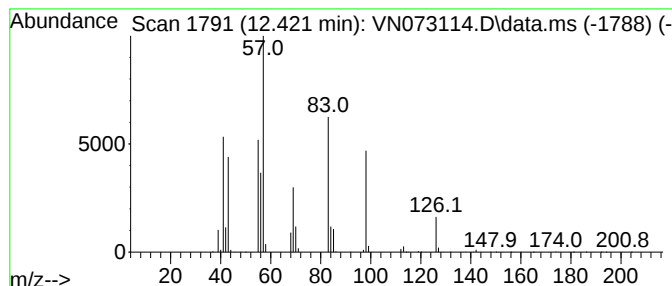
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Quant Title : SW846 8260

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

Peak Number 7 unknown12.421 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.421	176.57 ug/l	5765970	Chlorobenzene-d5	11.686

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	46
2			Cyclopentanone, 2,4,4-trimethyl-	126	C8H14O	004694-12-6	43
3			1H-Pyrazole, 3-ethyl-4,5-dihydro-	98	C5H10N2	005920-29-6	43
4			Furan, 2-methoxy-	98	C5H6O2	025414-22-6	38
5			Heptane, 4-propyl-	142	C10H22	003178-29-8	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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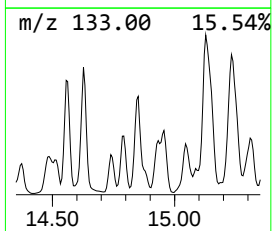
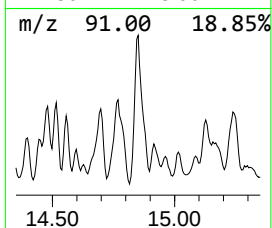
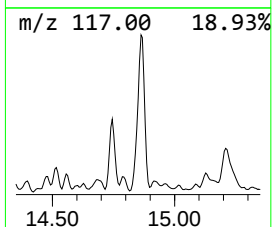
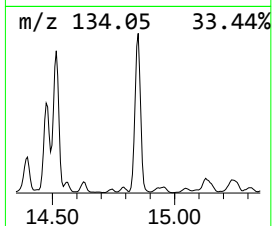
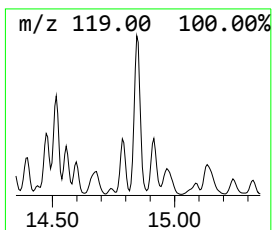
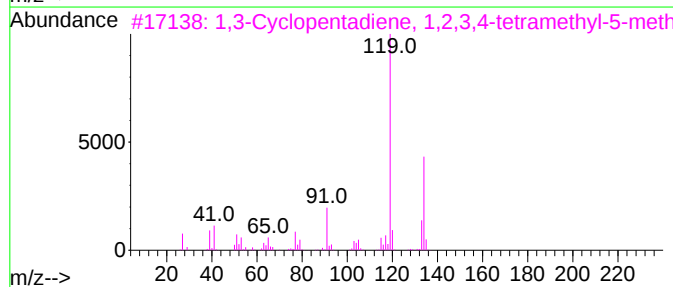
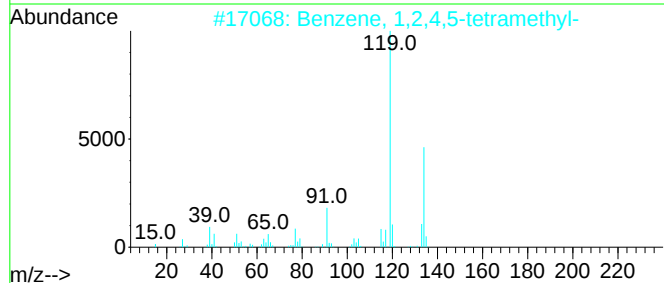
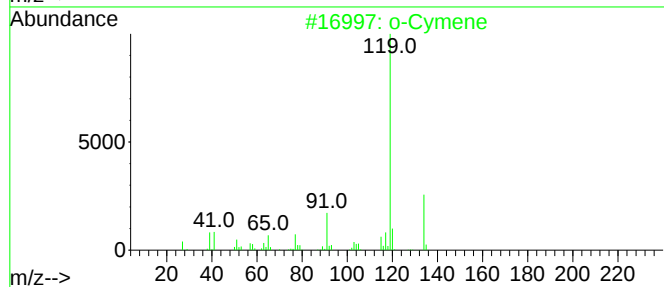
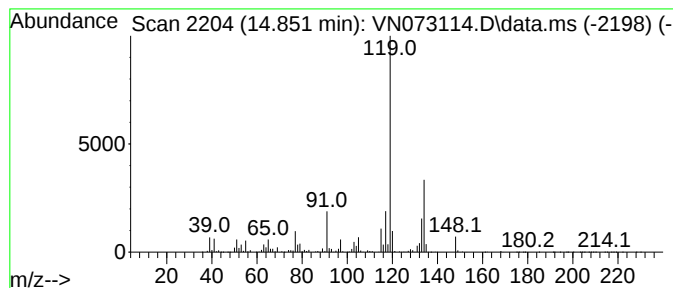
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Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.851	96.78 ug/l	13841000	1,4-Dichlorobenzene-d4	13.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Cymene	134	C10H14	000527-84-4	94
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
3		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	90
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	87
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
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Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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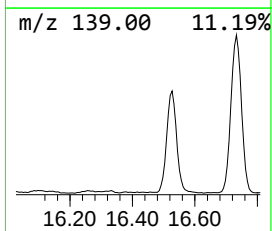
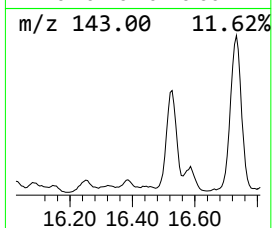
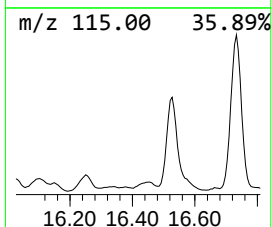
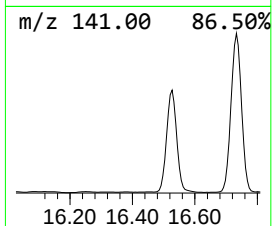
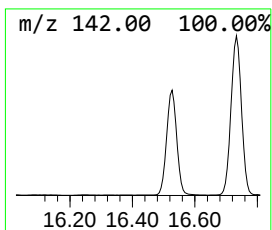
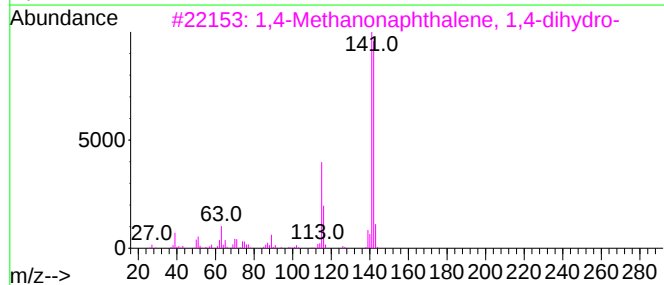
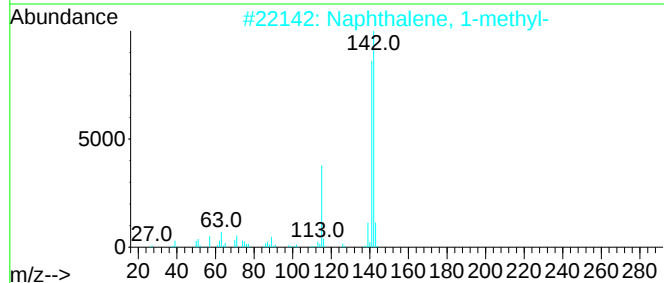
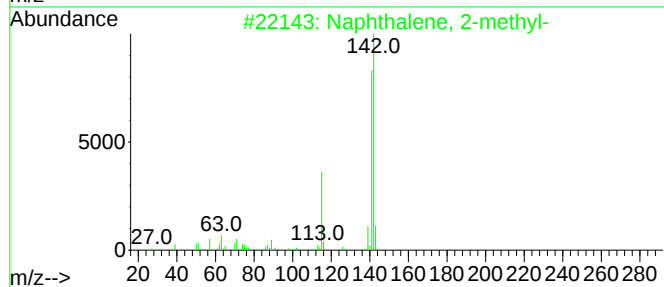
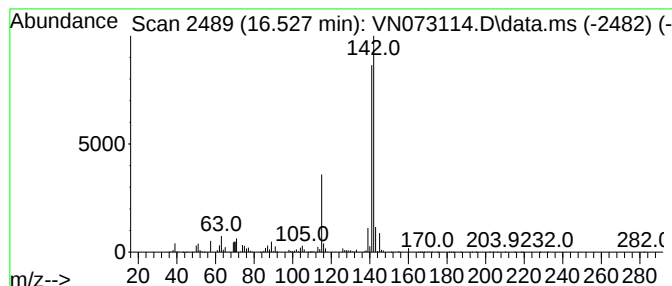
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.527	114.97 ug/l	16443100	1,4-Dichlorobenzene-d4	13.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14A-1A-1-GR

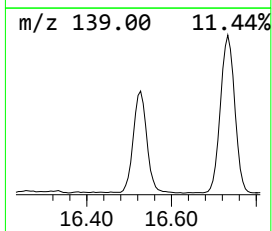
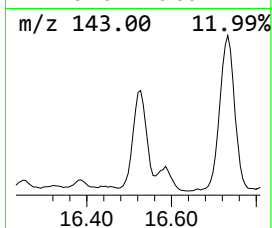
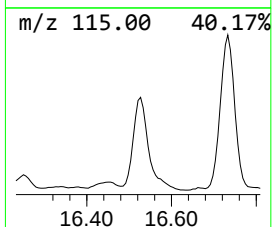
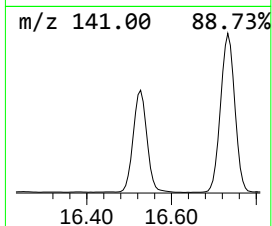
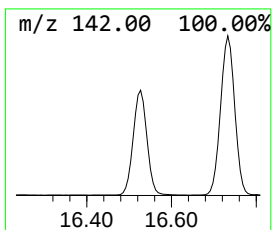
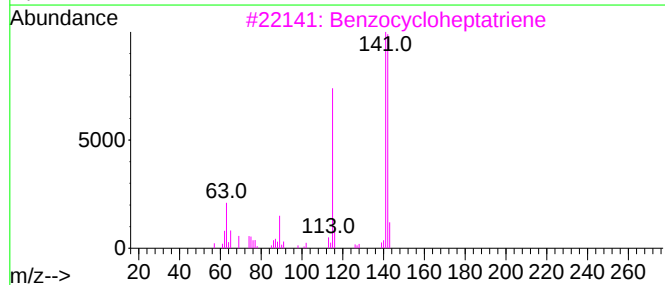
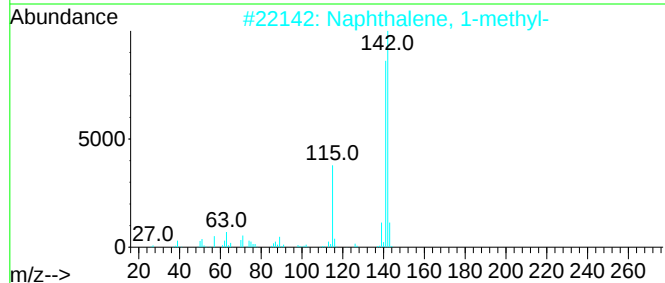
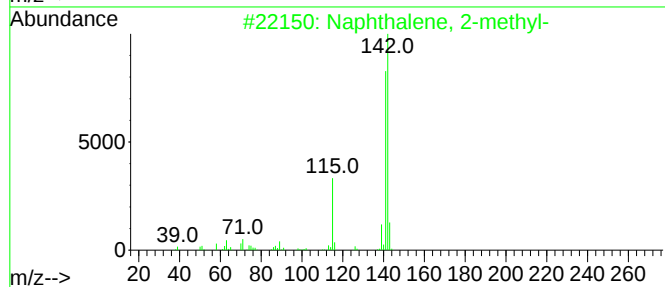
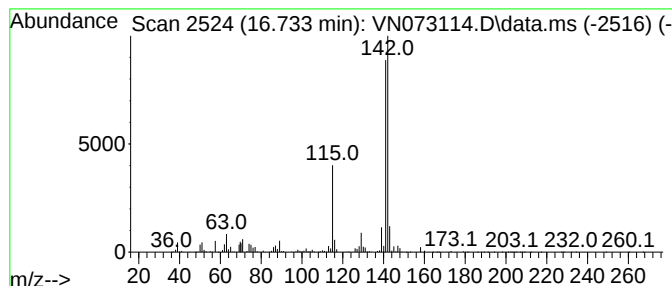
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.733	142.03 ug/l	20313100	1,4-Dichlorobenzene-d4	13.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	93
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073114.D
Acq On : 25 Jun 2022 16:43
Operator : JC\MD
Sample : N3454-07
Misc : 5.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14A-1A-1-GR

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
Heptane, 3-methyl-	10.145	67.3	ug/l	1993900	2	8.898	1480610	50.0
Cyclohexane, et...	11.227	99.1	ug/l	3237440	3	11.686	1632810	50.0
Cyclohexane, 1,...	11.280	134.6	ug/l	4396040	3	11.686	1632810	50.0
Cyclohexane, 1,...	11.480	119.8	ug/l	3911010	3	11.686	1632810	50.0
Nonane, 3-methyl-	12.304	113.1	ug/l	3694470	3	11.686	1632810	50.0
Bicyclo[3.3.1]n...	12.392	109.6	ug/l	3577360	3	11.686	1632810	50.0
unknown12.421	12.421	176.6	ug/l	5765970	3	11.686	1632810	50.0
o-Cymene	14.851	96.8	ug/l	13841000	4	13.610	7151080	50.0
Naphthalene, 1-...	16.527	115.0	ug/l	16443100	4	13.610	7151080	50.0
Benzocyclohepta...	16.733	142.0	ug/l	20313100	4	13.610	7151080	50.0