

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
 Data File : VN075883.D
 Acq On : 19 Dec 2022 19:58
 Operator : JC\MD
 Sample : N6096-02
 Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMP-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\82N121422W.M
 Title : SW846 8260

Signal : TIC: VN075883.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.224	1044	1049	1061	rVB	527410	1214122	1.74%	0.112%
2	9.100	1189	1198	1218	rBV	636027	1335019	1.91%	0.123%
3	9.600	1267	1283	1299	rBV	477836	1222437	1.75%	0.113%
4	10.206	1380	1386	1399	rVB2	484389	1118744	1.60%	0.103%
5	10.341	1399	1409	1423	rBV	442405	977187	1.40%	0.090%
6	10.571	1440	1448	1453	rBV	1932793	3822911	5.46%	0.352%
7	10.629	1453	1458	1464	rVB2	932064	1979025	2.83%	0.182%
8	10.747	1472	1478	1487	rVB	2027517	3641183	5.21%	0.336%
9	10.912	1501	1506	1514	rVB	692076	1254687	1.79%	0.116%
10	11.006	1514	1522	1531	rBV2	712553	1572845	2.25%	0.145%
11	11.177	1542	1551	1557	rBV	633578	1135220	1.62%	0.105%
12	11.276	1557	1568	1576	rBV2	532290	1445166	2.07%	0.133%
13	11.347	1576	1580	1584	rBV2	541857	958769	1.37%	0.088%
14	11.424	1584	1593	1597	rVV	1836172	3840939	5.49%	0.354%
15	11.476	1597	1602	1607	rVB	1658883	3037520	4.34%	0.280%
16	11.576	1614	1619	1623	rBV2	889992	1366479	1.95%	0.126%
17	11.647	1623	1631	1634	rBV2	2322528	4626432	6.61%	0.426%
18	11.747	1643	1648	1654	rVB	2256270	3478919	4.97%	0.321%
19	11.871	1663	1669	1673	rBV	712388	1227645	1.75%	0.113%
20	11.965	1679	1685	1689	rBV	2557092	4430214	6.33%	0.408%
21	12.006	1689	1692	1695	rVV	1254425	2032024	2.90%	0.187%
22	12.076	1695	1704	1708	rVV	18813954	35877972	51.29%	3.306%
23	12.112	1708	1710	1714	rVV	3588418	5387305	7.70%	0.496%
24	12.159	1714	1718	1723	rVB2	2475486	4349773	6.22%	0.401%
25	12.212	1723	1727	1734	rBV3	435573	912591	1.30%	0.084%
26	12.288	1734	1740	1747	rVB4	767982	1594839	2.28%	0.147%
27	12.400	1747	1759	1766	rBV3	6481272	20441915	29.22%	1.884%
28	12.488	1767	1774	1778	rVB2	3129141	6019876	8.61%	0.555%
29	12.541	1778	1783	1785	rBV2	1897104	3047828	4.36%	0.281%
30	12.594	1786	1792	1793	rVV	3870021	7764629	11.10%	0.715%
31	12.623	1794	1797	1805	rVV2	5495782	10892488	15.57%	1.004%
32	12.700	1805	1810	1814	rVV2	2399324	5087198	7.27%	0.469%
33	12.747	1814	1818	1819	rVV3	1689377	2673125	3.82%	0.246%
34	12.776	1819	1823	1825	rVV2	2850229	4736948	6.77%	0.436%
35	12.800	1825	1827	1832	rVB2	3344309	4260490	6.09%	0.393%
36	12.859	1832	1837	1838	rBV2	1651121	2380590	3.40%	0.219%

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37	12.882	1838	1841	1846	rVB	1938189	2693991	3.85%	0.248%
38	12.941	1846	1851	1855	rBV3	1424307	2382109	3.41%	0.220%
39	13.035	1860	1867	1872	rVB3	6939240	14572740	20.83%	1.343%
40	13.100	1872	1878	1882	rBV	18780830	32577025	46.57%	3.002%

41	13.170	1882	1890	1897	rVB2	32367136	69953846	100.00%	6.446%
42	13.229	1897	1900	1906	rVB2	3119433	5134939	7.34%	0.473%
43	13.341	1906	1919	1925	rBV	7408483	19199549	27.45%	1.769%
44	13.400	1925	1929	1938	rVB4	4171123	8550692	12.22%	0.788%
45	13.488	1938	1944	1954	rBV2	41371589	66009474	94.36%	6.083%

46	13.582	1956	1960	1962	rBV2	1207928	1835085	2.62%	0.169%
47	13.618	1962	1966	1971	rVB2	4386735	7401678	10.58%	0.682%
48	13.682	1971	1977	1982	rBV3	10744599	21157570	30.25%	1.950%
49	13.729	1982	1985	1992	rVB3	4118259	7264499	10.38%	0.669%
50	13.788	1992	1995	1997	rBV2	2318488	2985360	4.27%	0.275%

51	13.823	1997	2001	2006	rVB	12917908	18602685	26.59%	1.714%
52	13.865	2006	2008	2014	rVB3	1873867	2346444	3.35%	0.216%
53	13.953	2014	2023	2025	rBV4	5179747	10080104	14.41%	0.929%
54	13.988	2025	2029	2032	rVV2	13239979	18664554	26.68%	1.720%
55	14.029	2032	2036	2046	rVB3	15502296	34562125	49.41%	3.185%

56	14.112	2046	2050	2058	rBV3	22644526	41144465	58.82%	3.791%
57	14.188	2059	2063	2067	rVB	4239194	5488938	7.85%	0.506%
58	14.253	2069	2074	2076	rBV	7464796	12003735	17.16%	1.106%
59	14.276	2076	2078	2083	rVB	9179277	12506585	17.88%	1.152%
60	14.335	2083	2088	2093	rVB	14117606	20774335	29.70%	1.914%

61	14.400	2097	2099	2101	rBV2	794308	865768	1.24%	0.080%
62	14.447	2101	2107	2112	rVB2	14704070	24288076	34.72%	2.238%
63	14.523	2112	2120	2124	rVB3	3252496	6853141	9.80%	0.632%
64	14.582	2124	2130	2134	rBV3	6958852	12133019	17.34%	1.118%
65	14.635	2134	2139	2141	rBV	10550539	16333900	23.35%	1.505%

66	14.665	2141	2144	2147	rVB2	5460657	6445688	9.21%	0.594%
67	14.700	2147	2150	2153	rVB	9345697	11387277	16.28%	1.049%
68	14.741	2153	2157	2160	rVB	8273157	10231054	14.63%	0.943%
69	14.800	2160	2167	2172	rVB3	6763208	16503873	23.59%	1.521%
70	14.882	2172	2181	2185	rBV2	7719057	16059693	22.96%	1.480%

71	14.964	2185	2195	2202	rBV5	20105955	49684107	71.02%	4.578%
72	15.053	2202	2210	2216	rBV4	16806280	53108449	75.92%	4.894%
73	15.217	2233	2238	2244	rVB	12768892	21980863	31.42%	2.025%
74	15.282	2244	2249	2251	rBV4	3512455	5468688	7.82%	0.504%
75	15.323	2251	2256	2261	rBV2	8277156	14632848	20.92%	1.348%

76	15.447	2270	2277	2284	rVB5	9660227	23198980	33.16%	2.138%
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77	15.523	2286	2290	2291	rBV3	1272059	1851101	2.65%	0.171%
78	15.600	2298	2303	2307	rBV2	5325368	9998752	14.29%	0.921%
79	15.706	2314	2321	2325	rBV	11366429	22869112	32.69%	2.107%
80	15.753	2325	2329	2332	rVV2	7332829	14321598	20.47%	1.320%
81	15.788	2332	2335	2340	rVV3	7424281	16609163	23.74%	1.530%
82	15.841	2340	2344	2354	rVB3	8143998	18237028	26.07%	1.681%
83	15.953	2354	2363	2370	rBV4	6283397	16130139	23.06%	1.486%
84	16.029	2371	2376	2381	rBV	1592481	3454211	4.94%	0.318%
85	16.129	2386	2393	2394	rVV4	2903241	5972435	8.54%	0.550%
86	16.176	2395	2401	2410	rVB	18204599	37931722	54.22%	3.495%
87	16.264	2411	2416	2422	rBV6	2121240	4371781	6.25%	0.403%
88	16.347	2424	2430	2436	rVB2	1792377	3905410	5.58%	0.360%
89	16.511	2448	2458	2469	rVB2	8344985	20878301	29.85%	1.924%
90	16.717	2483	2493	2500	rBV2	8617197	20468771	29.26%	1.886%

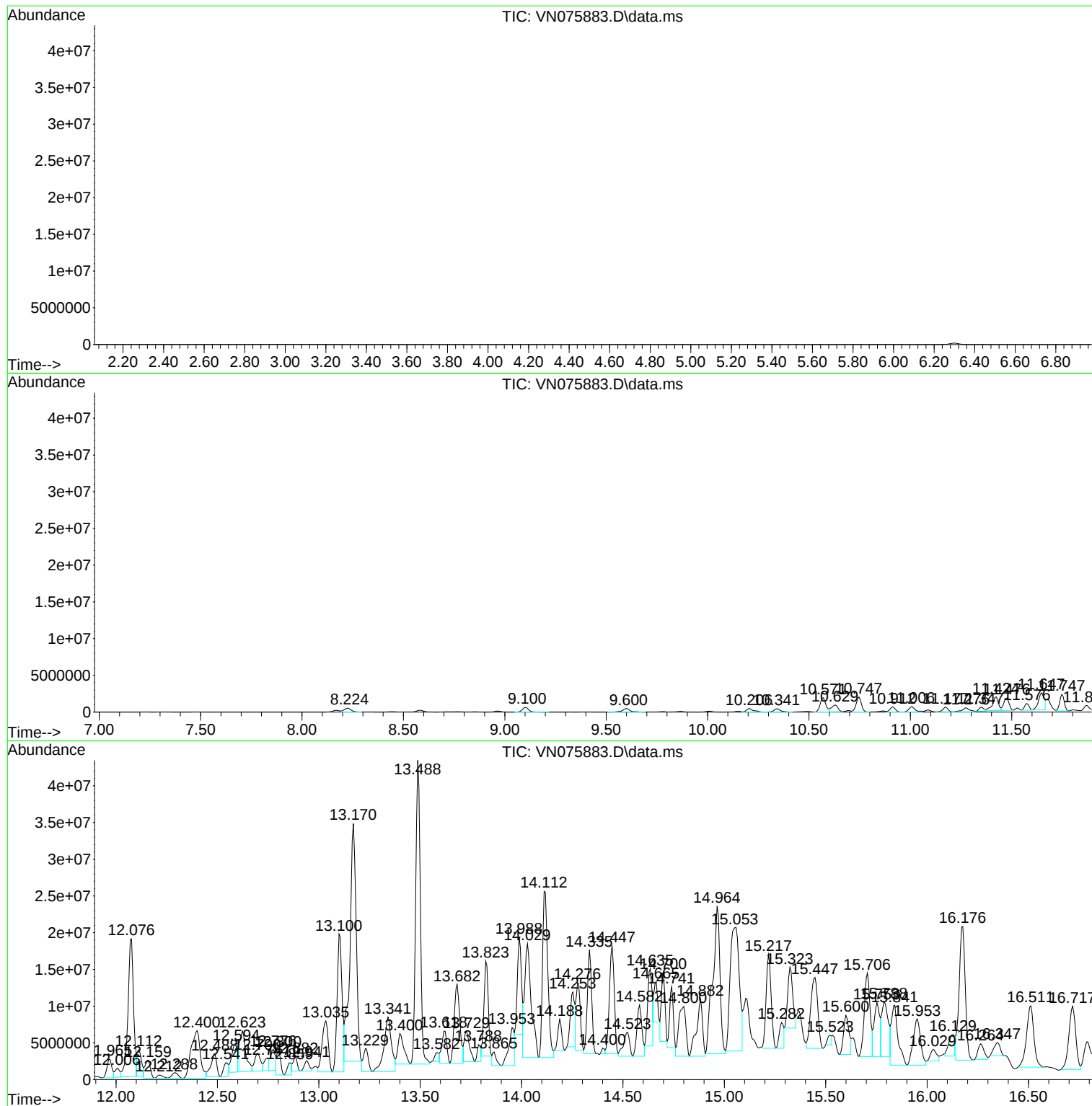
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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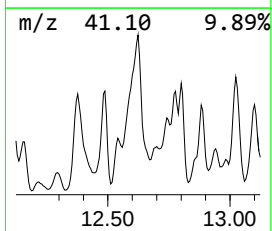
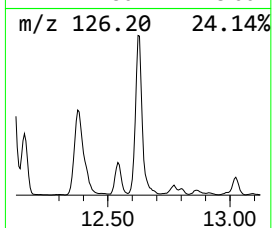
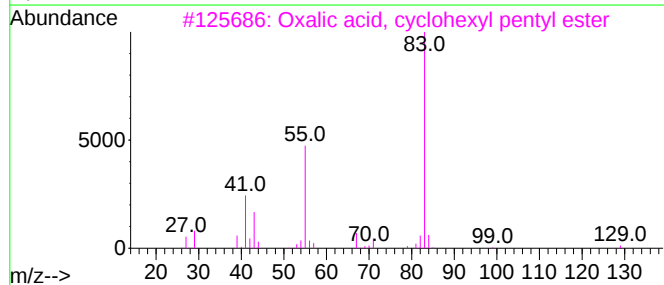
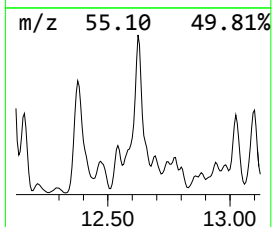
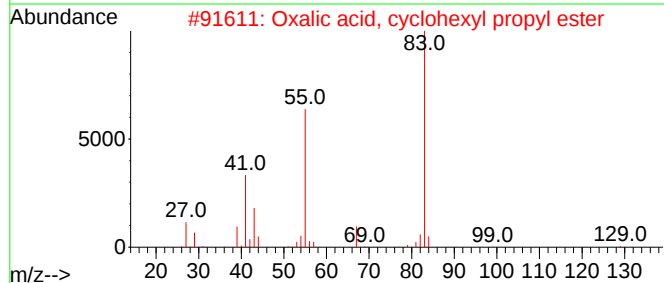
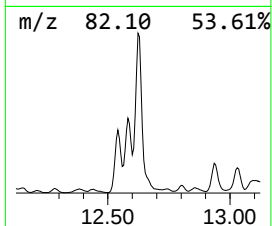
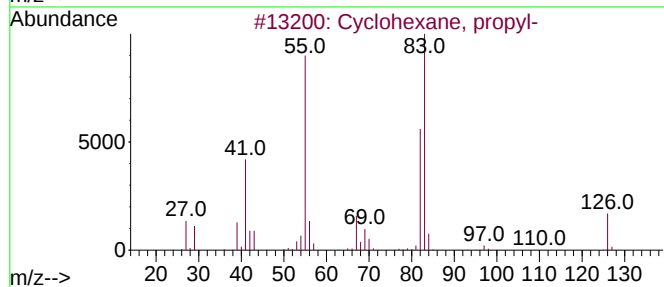
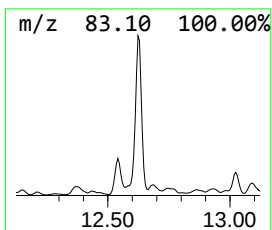
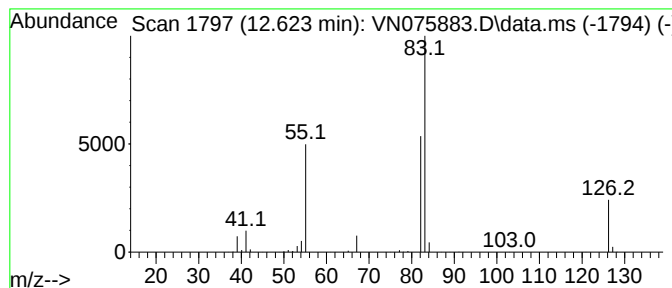
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.624	443.63 ug/l	10892500	Chlorobenzene-d5	11.871

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	64
2			Oxalic acid, cyclohexyl propyl e...	214	C11H18O4	1000309-30-3	40
3			Oxalic acid, cyclohexyl pentyl e...	242	C13H22O4	1000309-30-6	40
4			Oxalic acid, cyclohexyl butyl ester	228	C12H20O4	1000309-30-5	40
5			Oxalic acid, cyclohexyl ethyl ester	200	C10H16O4	1000309-30-2	40



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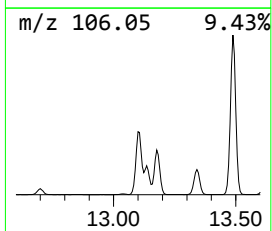
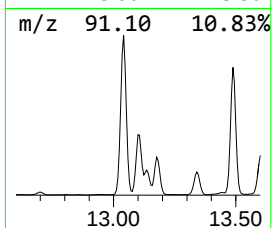
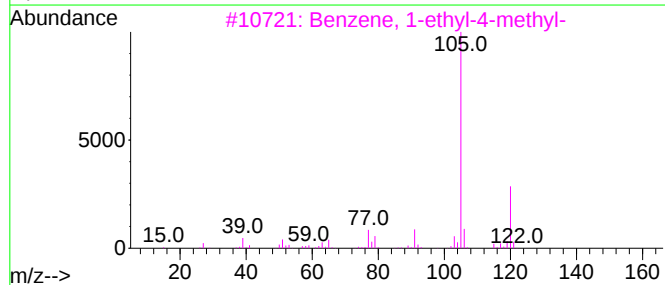
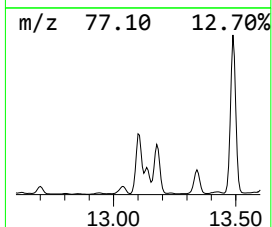
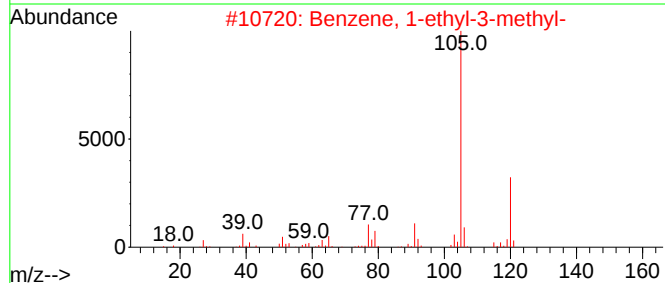
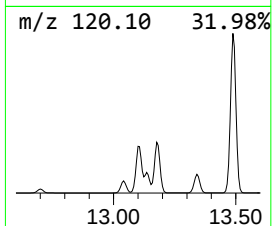
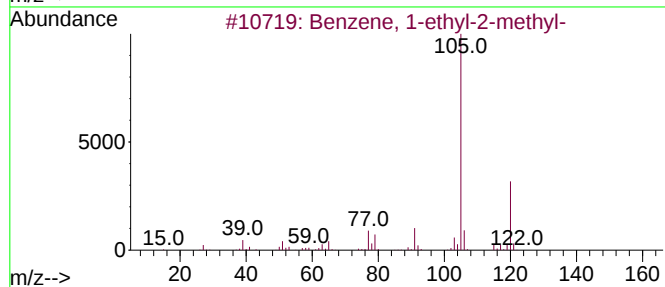
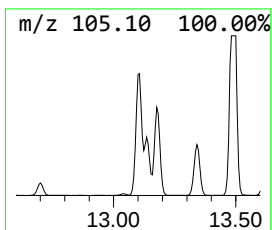
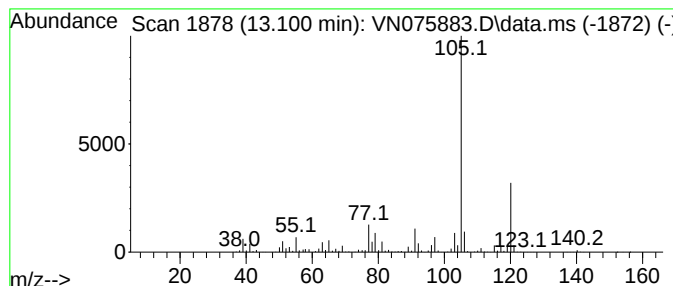
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	545.61 ug/l	32577000	1,4-Dichlorobenzene-d4	13.794

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-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5			Mesitylene	120	C9H12	000108-67-8	90



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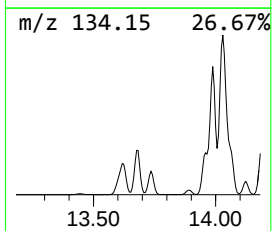
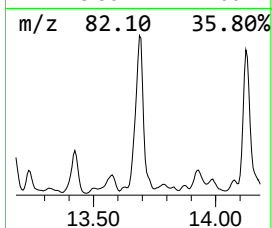
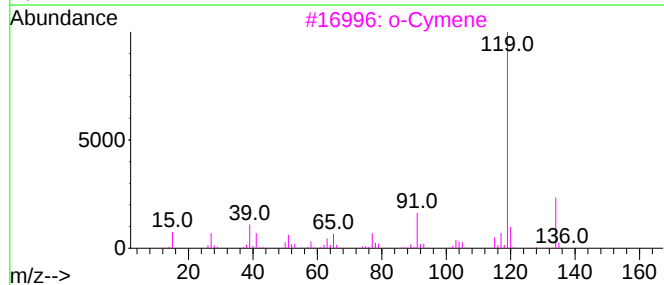
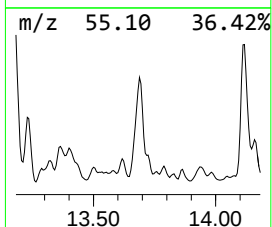
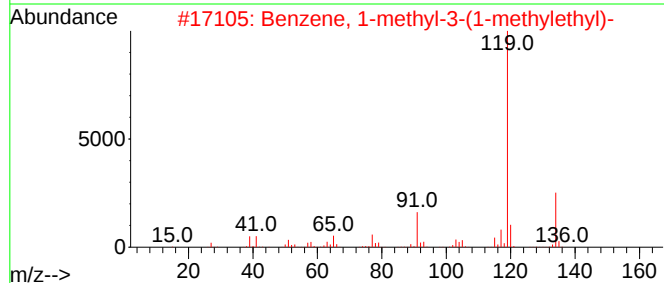
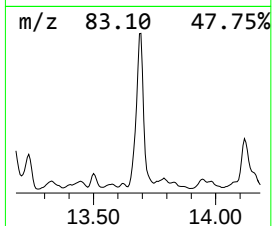
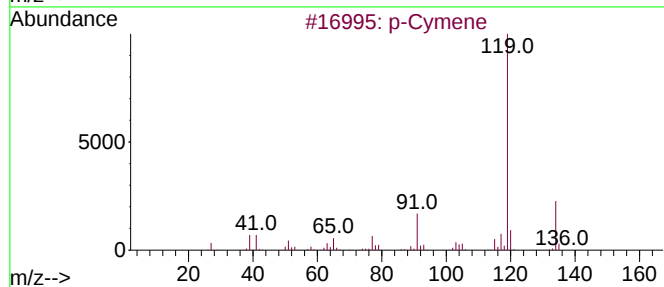
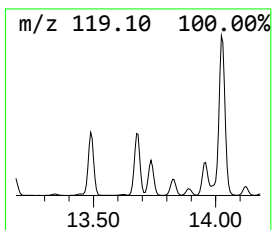
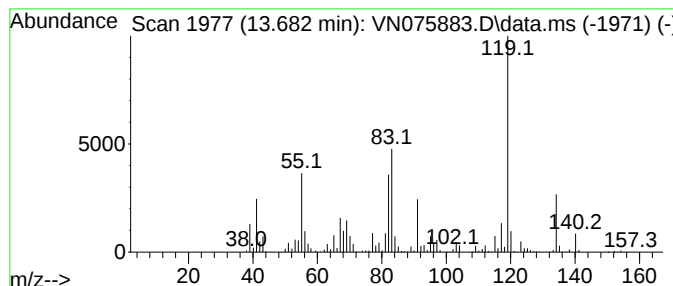
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.682	354.36 ug/l	21157600	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	92
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	91
3			o-Cymene	134	C10H14	000527-84-4	90
4			1,3,5-Cycloheptatriene, 3,7,7-trimethyl-	134	C10H14	003479-89-8	78
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

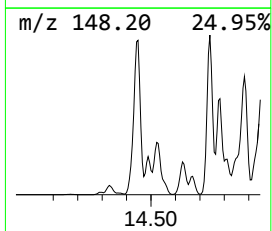
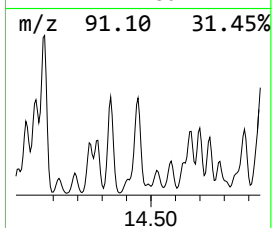
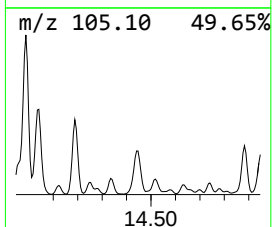
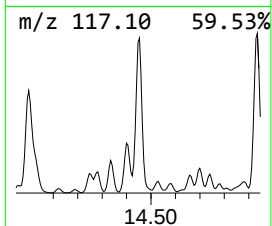
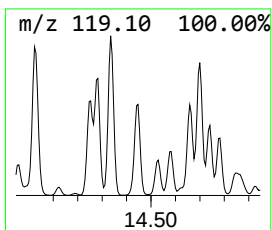
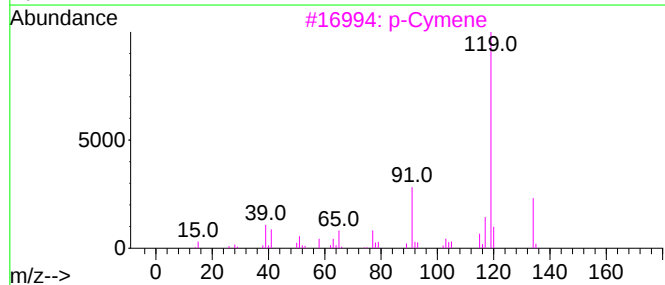
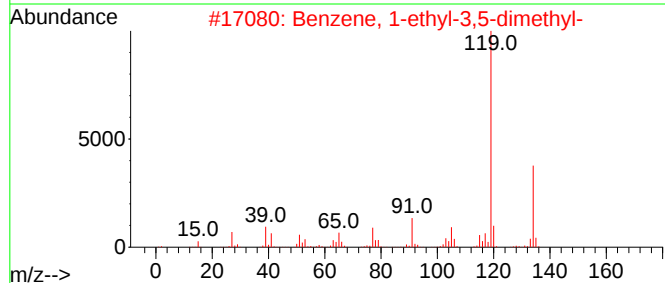
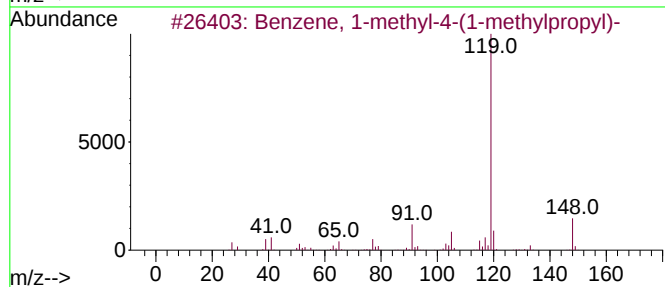
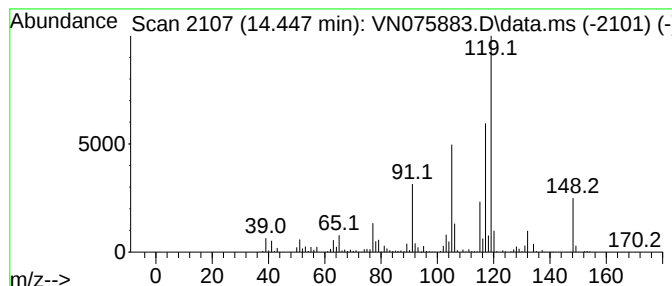
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-4-(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.447	406.79 ug/l	24288100	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	53
2			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	49
3			p-Cymene	134	C10H14	000099-87-6	46
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	46
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

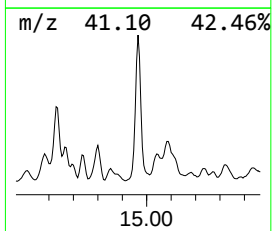
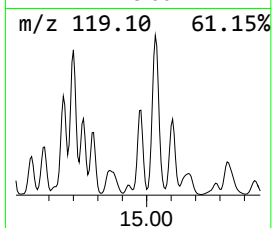
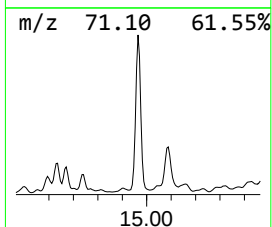
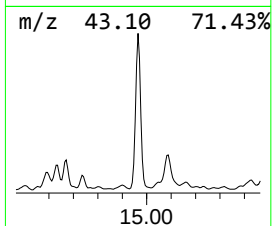
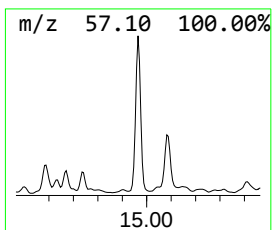
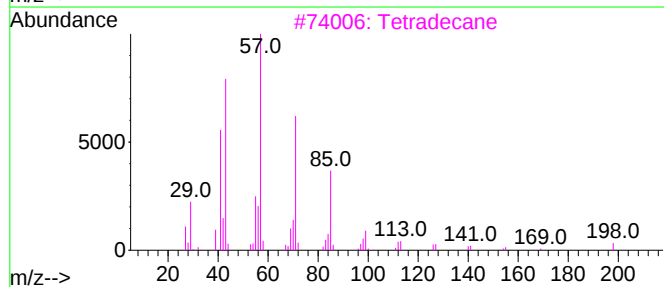
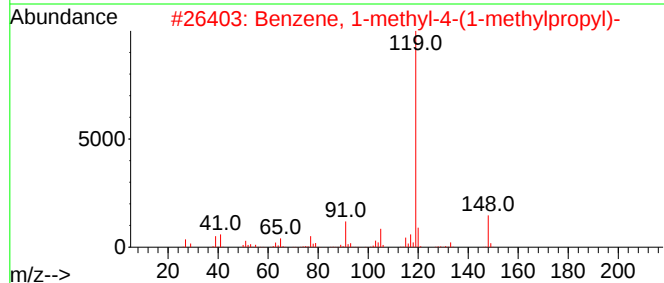
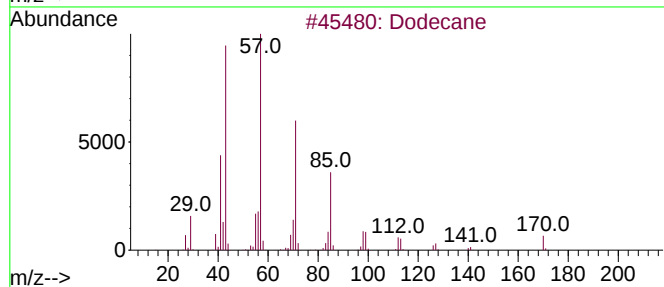
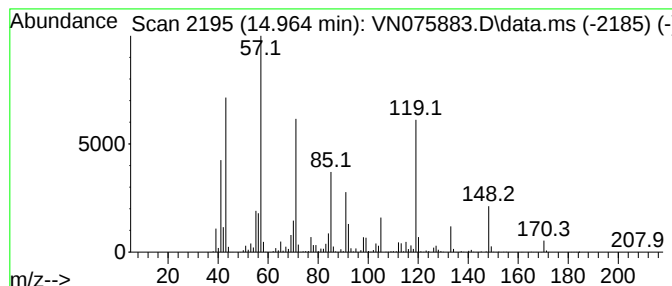
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Quant Title : SW846 8260

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

Peak Number 5 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.964	832.13 ug/l	49684100	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	95
2			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	42
3			Tetradecane	198	C14H30	000629-59-4	41
4			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38
5			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

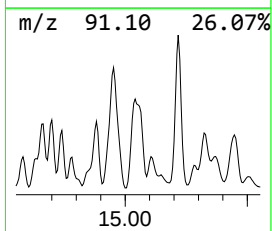
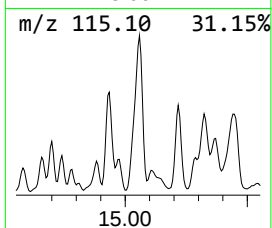
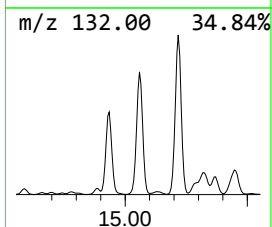
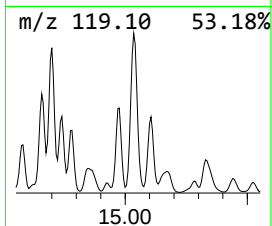
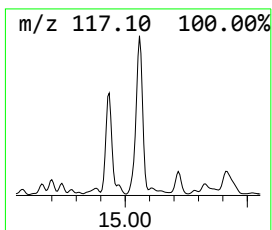
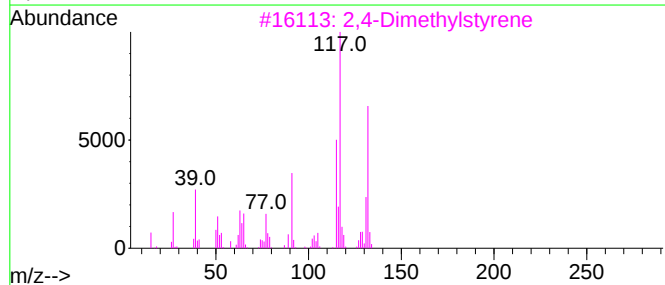
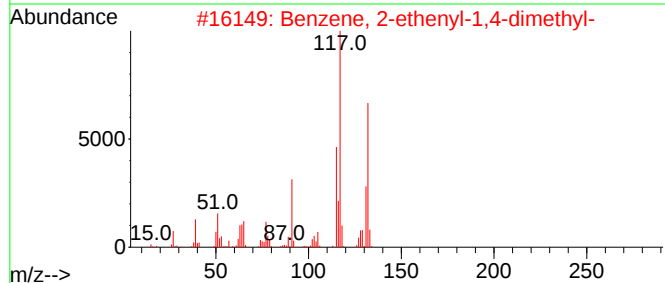
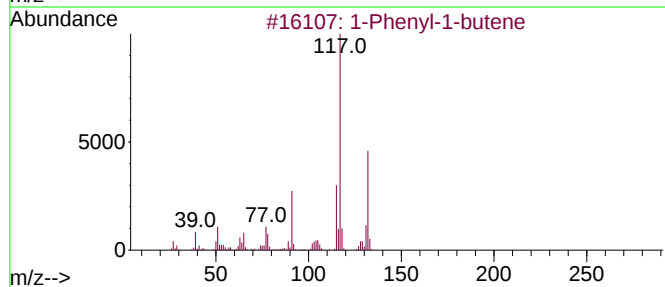
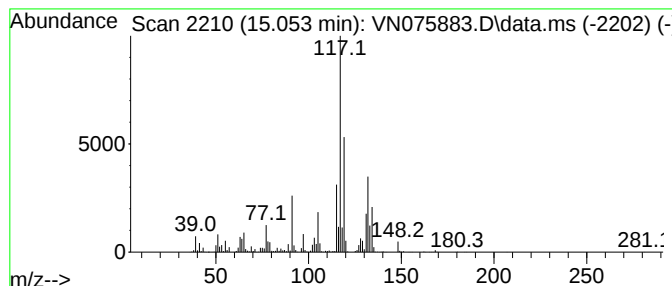
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Quant Title : SW846 8260

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

Peak Number 6 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	889.48 ug/l	53108400	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	91
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	89
3			2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

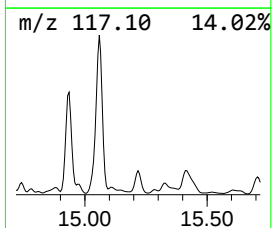
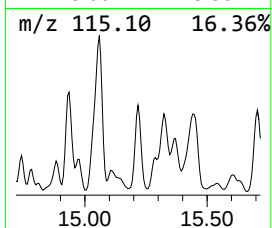
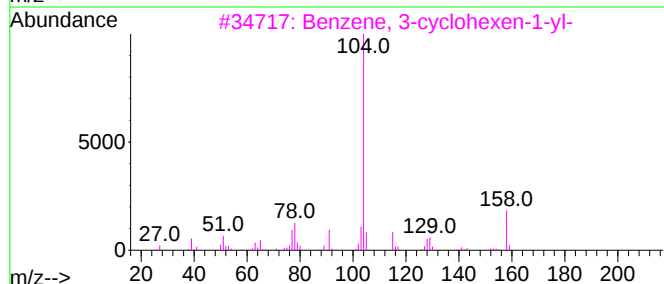
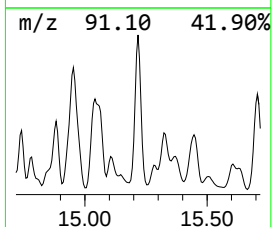
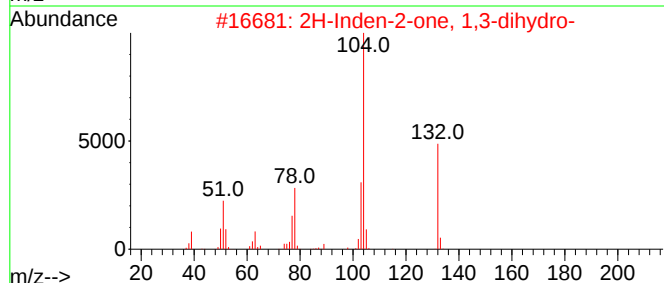
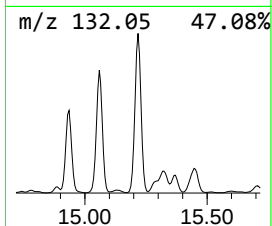
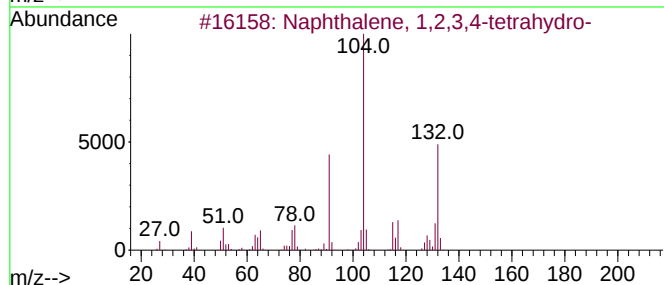
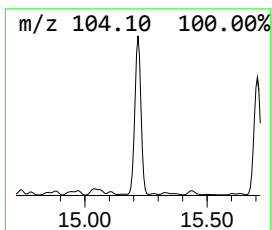
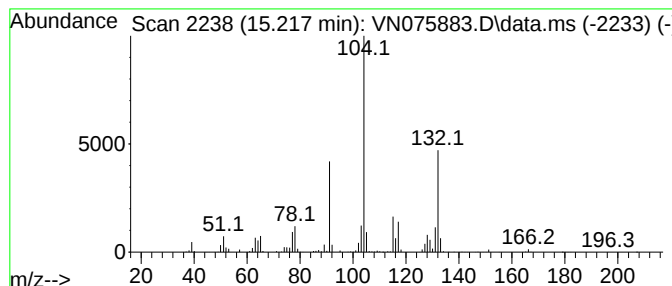
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.217	368.14 ug/l	21980900	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
3			Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47
4			Benzenepropanoic acid, methyl ester	164	C10H12O2	000103-25-3	43
5			Acetic acid, trifluoro-, 2-phenyl-	218	C10H9F3O2	055419-66-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

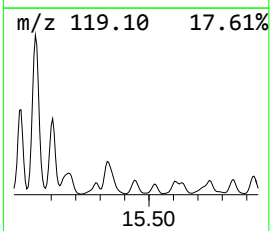
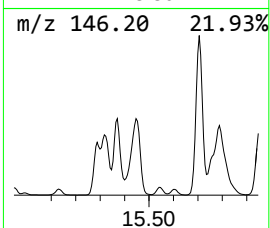
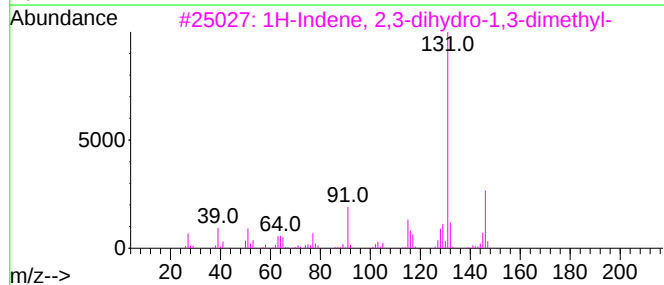
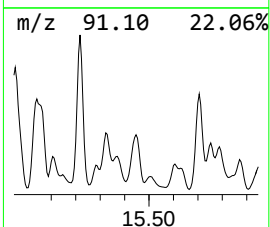
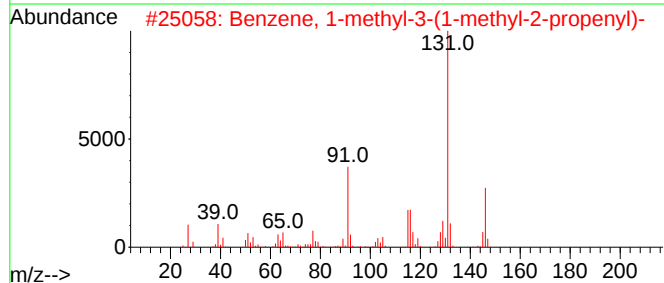
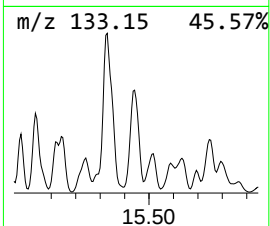
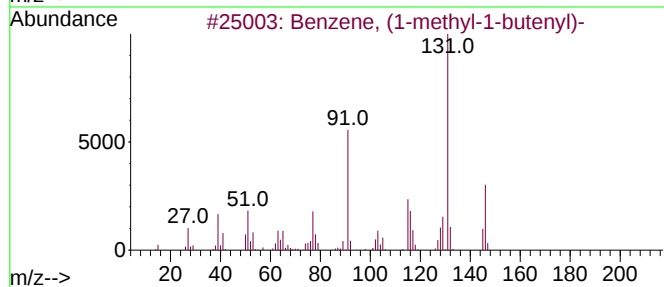
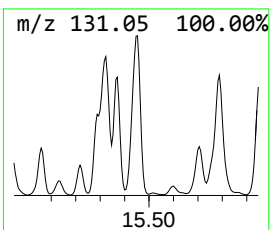
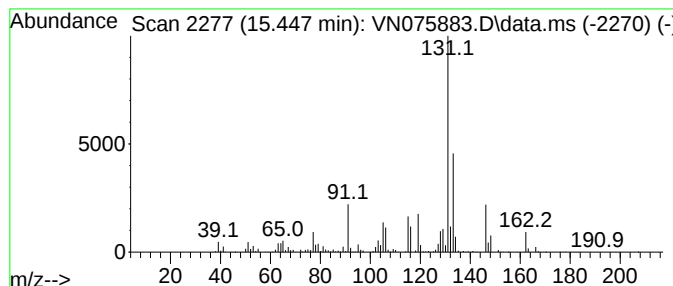
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.447	388.55 ug/l	23199000	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
2			Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	64
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
4			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	64
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

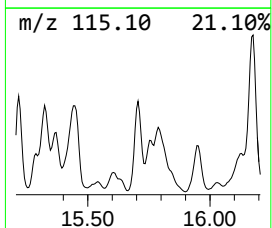
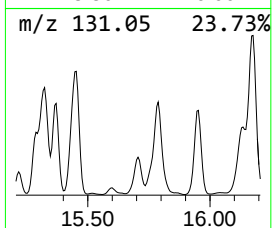
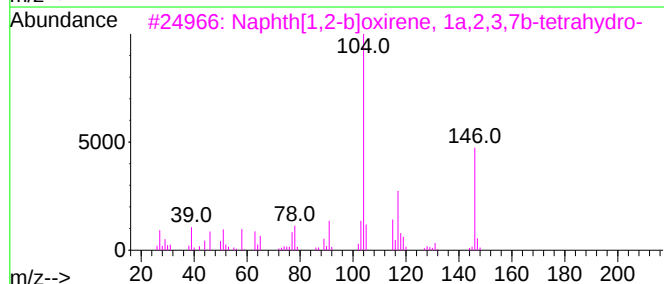
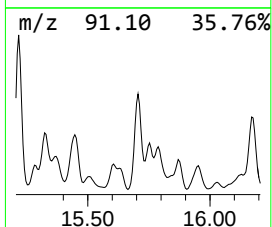
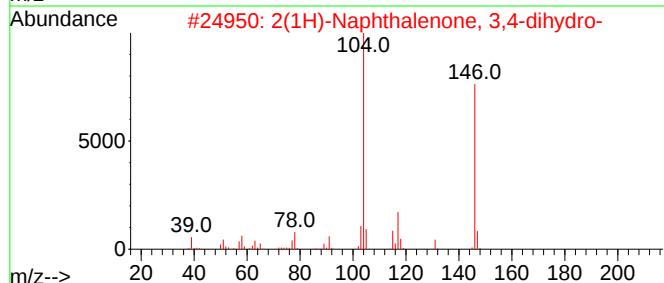
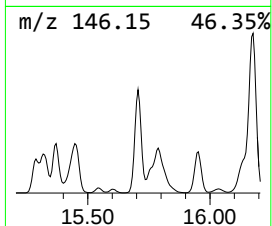
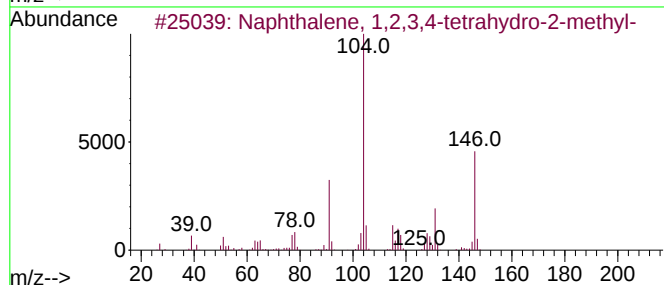
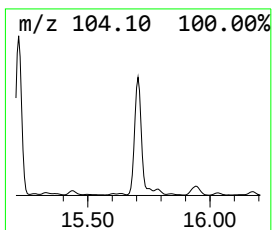
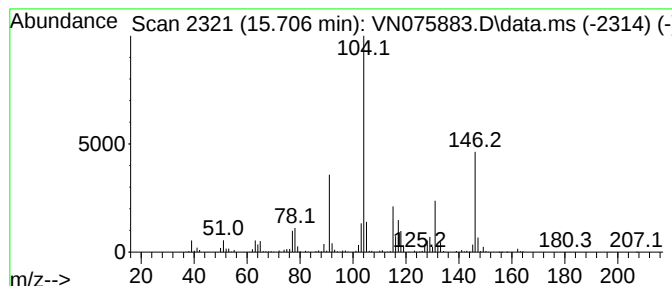
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.706	383.02 ug/l	22869100	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	95
2			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	64
3			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	64
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	38
5			Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

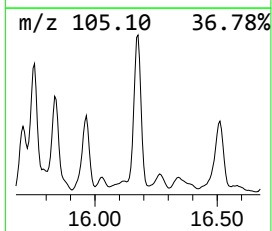
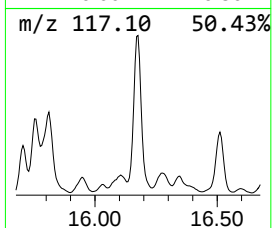
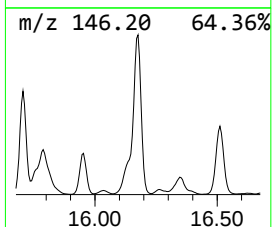
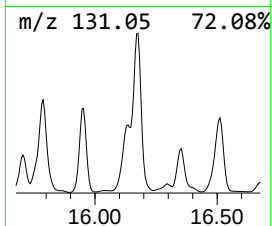
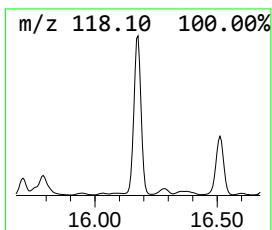
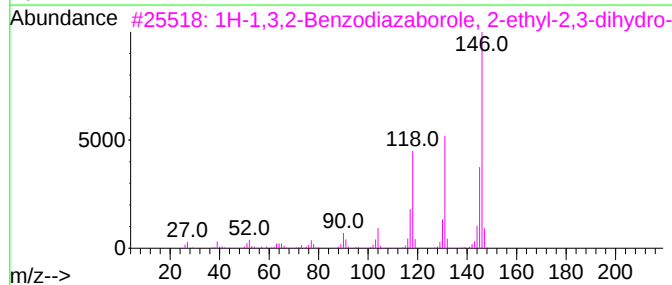
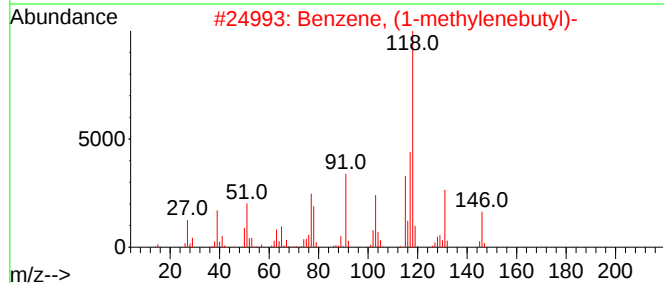
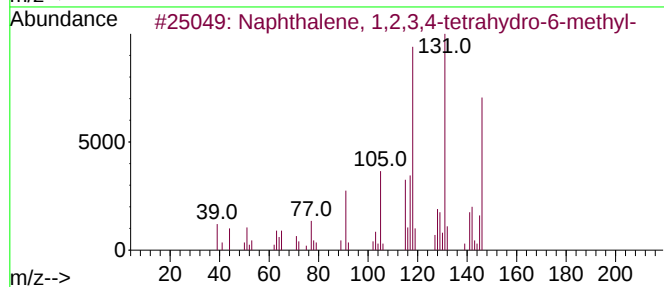
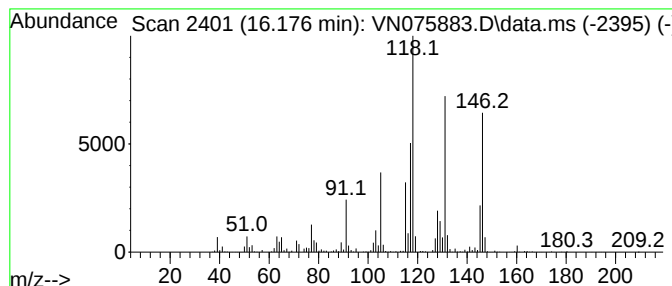
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.176	635.30 ug/l	37931700	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
2			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	72
3			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	70
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	70
5			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075883.D
Acq On : 19 Dec 2022 19:58
Operator : JC\MD
Sample : N6096-02
Misc : 5.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMP-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.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
Cyclohexane, pr...	12.624	443.6	ug/l	10892500	3	11.871	1227650	50.0
Benzene, 1-ethy...	13.100	545.6	ug/l	32577000	4	13.794	2985360	50.0
p-Cymene	13.682	354.4	ug/l	21157600	4	13.794	2985360	50.0
Benzene, 1-meth...	14.447	406.8	ug/l	24288100	4	13.794	2985360	50.0
Dodecane	14.964	832.1	ug/l	49684100	4	13.794	2985360	50.0
1-Phenyl-1-butene	15.053	889.5	ug/l	53108400	4	13.794	2985360	50.0
Naphthalene, 1,...	15.217	368.1	ug/l	21980900	4	13.794	2985360	50.0
Benzene, (1-met...	15.447	388.6	ug/l	23199000	4	13.794	2985360	50.0
Naphthalene, 1,...	15.706	383.0	ug/l	22869100	4	13.794	2985360	50.0
Naphthalene, 1,...	16.176	635.3	ug/l	37931700	4	13.794	2985360	50.0