

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
 Data File : VY015025.D
 Acq On : 10 Aug 2023 14:18
 Operator : SY/MD
 Sample : 03945-02
 Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 NEVINS-SB01-Z78-80

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_Y\methods\82Y080723S.M

Title : SW846 8260

Signal : TIC: VY015025.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.265	226	237	256	rVB3	772913	2396931	1.27%	0.104%
2	4.765	464	483	492	rBV	522432	2260492	1.20%	0.098%
3	5.747	630	644	662	rBV	1848336	5875083	3.11%	0.255%
4	6.789	805	815	827	rBV	626430	1891638	1.00%	0.082%
5	7.777	959	977	986	rBV5	1682452	5821251	3.08%	0.253%
6	8.002	1005	1014	1022	rBV	1271224	3403316	1.80%	0.148%
7	8.155	1030	1039	1044	rBV4	39094683	116222051	61.45%	5.043%
8	8.295	1053	1062	1068	rBV4	1432033	3743911	1.98%	0.162%
9	8.411	1076	1081	1094	rVB4	1158548	3364735	1.78%	0.146%
10	8.551	1094	1104	1118	rVB	5197687	11112996	5.88%	0.482%
11	8.691	1118	1127	1131	rBV2	1360281	3136044	1.66%	0.136%
12	8.959	1161	1171	1188	rVB3	1044114	2942904	1.56%	0.128%
13	9.185	1192	1208	1214	rBV2	5551671	14274578	7.55%	0.619%
14	9.246	1214	1218	1227	rVB2	1025066	2171514	1.15%	0.094%
15	9.563	1259	1270	1273	rBV	1007358	3053696	1.61%	0.133%
16	9.612	1273	1278	1286	rVB2	3181243	6358073	3.36%	0.276%
17	9.764	1298	1303	1308	rVV2	1877386	3413002	1.80%	0.148%
18	9.825	1308	1313	1326	rVB3	1818393	4771345	2.52%	0.207%
19	9.959	1326	1335	1338	rBV2	1056496	2636365	1.39%	0.114%
20	10.063	1345	1352	1359	rVV	2609378	5159144	2.73%	0.224%
21	10.173	1363	1370	1375	rVV4	3061093	7290321	3.85%	0.316%
22	10.246	1375	1382	1389	rVB2	9206566	20044368	10.60%	0.870%
23	10.374	1394	1403	1410	rVB2	5826014	12292723	6.50%	0.533%
24	10.526	1423	1428	1435	rVB2	1665422	3448201	1.82%	0.150%
25	10.612	1435	1442	1454	rBV5	3090543	9823102	5.19%	0.426%
26	10.807	1464	1474	1481	rBV2	3129312	7868194	4.16%	0.341%
27	10.886	1481	1487	1488	rBV	1697296	2751335	1.45%	0.119%
28	10.916	1488	1492	1497	rVB2	1969124	3616836	1.91%	0.157%
29	11.038	1509	1512	1517	rVB	1370672	2159707	1.14%	0.094%
30	11.093	1517	1521	1526	rBV	2350961	4162689	2.20%	0.181%
31	11.148	1526	1530	1536	rBV2	2326493	4701826	2.49%	0.204%
32	11.307	1548	1556	1566	rBV8	3076640	9361089	4.95%	0.406%
33	11.416	1566	1574	1582	rVB3	2701181	7721003	4.08%	0.335%
34	11.532	1582	1593	1597	rBV5	1861993	6594697	3.49%	0.286%
35	11.599	1597	1604	1608	rBV5	45579044	135809737	71.80%	5.893%
36	11.715	1617	1623	1626	rBV4	46274192	109146017	57.70%	4.736%

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Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
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37	11.849	1641	1645	1651	rBV3	4887390	8521134	4.51%	0.370%
38	11.910	1651	1655	1659	rBV4	1240840	2116804	1.12%	0.092%
39	12.044	1665	1677	1685	rBV4	50144942	174924155	92.48%	7.590%
40	12.130	1685	1691	1697	rVB2	6970624	13955288	7.38%	0.606%
41	12.209	1698	1704	1707	rBV5	4580516	9609692	5.08%	0.417%
42	12.252	1707	1711	1716	rVB3	6385798	11559876	6.11%	0.502%
43	12.337	1718	1725	1732	rBV2	40425057	109434015	57.86%	4.749%
44	12.538	1753	1758	1765	rBV4	3000530	7048731	3.73%	0.306%
45	12.617	1766	1771	1774	rBV	5501779	11117678	5.88%	0.482%
46	12.684	1775	1782	1786	rBV2	23314132	54011235	28.56%	2.344%
47	12.745	1786	1792	1795	rBV4	40775117	100196418	52.97%	4.348%
48	12.983	1824	1831	1839	rBV2	40199077	99136179	52.41%	4.302%
49	13.050	1839	1842	1846	rVV	7751185	12151222	6.42%	0.527%
50	13.129	1846	1855	1861	rVV4	51388505	189144905	100.00%	8.207%
51	13.184	1861	1864	1869	rVV	41346921	73317693	38.76%	3.181%
52	13.245	1869	1874	1880	rVV4	7325665	16279252	8.61%	0.706%
53	13.325	1880	1887	1892	rVV	31078263	68987210	36.47%	2.993%
54	13.385	1892	1897	1904	rVV	21897086	44554650	23.56%	1.933%
55	13.471	1904	1911	1916	rVV2	37712672	96443952	50.99%	4.185%
56	13.526	1916	1920	1929	rVB	26423008	51640263	27.30%	2.241%
57	13.635	1929	1938	1941	rBV4	53745274	169133651	89.42%	7.339%
58	13.678	1942	1945	1953	rVV4	55533511	93401422	49.38%	4.053%
59	13.751	1953	1957	1962	rVV	12913260	20425932	10.80%	0.886%
60	13.818	1963	1968	1973	rVB2	7352382	13161318	6.96%	0.571%
61	13.885	1975	1979	1981	rBV	7647937	11214171	5.93%	0.487%
62	13.916	1981	1984	1988	rVB	13497269	18609545	9.84%	0.807%
63	13.971	1988	1993	1998	rBV	29183125	45810907	24.22%	1.988%
64	14.080	2006	2011	2021	rBV	15864647	33736585	17.84%	1.464%
65	14.160	2021	2024	2027	rVB	2773676	3644661	1.93%	0.158%
66	14.215	2027	2033	2037	rBV3	1581767	2580982	1.36%	0.112%
67	14.263	2037	2041	2043	rBV3	1267392	1995512	1.06%	0.087%
68	14.294	2043	2046	2049	rVV2	2875063	3814295	2.02%	0.166%
69	14.330	2049	2052	2056	rVB	5909836	8120631	4.29%	0.352%
70	14.379	2056	2060	2064	rVB	4520185	6207868	3.28%	0.269%
71	14.440	2064	2070	2074	rBV4	1602720	3261817	1.72%	0.142%
72	14.562	2085	2090	2095	rBV	14349821	20714672	10.95%	0.899%
73	14.605	2095	2097	2102	rVB2	1878287	2492765	1.32%	0.108%
74	14.672	2102	2108	2114	rBV2	4515195	9310484	4.92%	0.404%
75	14.745	2114	2120	2128	rVV	11596346	18063039	9.55%	0.784%
76	14.830	2128	2134	2143	rVB	15111721	23783592	12.57%	1.032%

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ClientSampleId :
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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

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Title : SW846 8260

77	14.928	2144	2150	2157	rVB3	2461632	4798142	2.54%	0.208%
78	15.233	2192	2200	2211	rVV3	39088692	83224198	44.00%	3.611%
79	15.330	2211	2216	2229	rVB	1017588	2130736	1.13%	0.092%

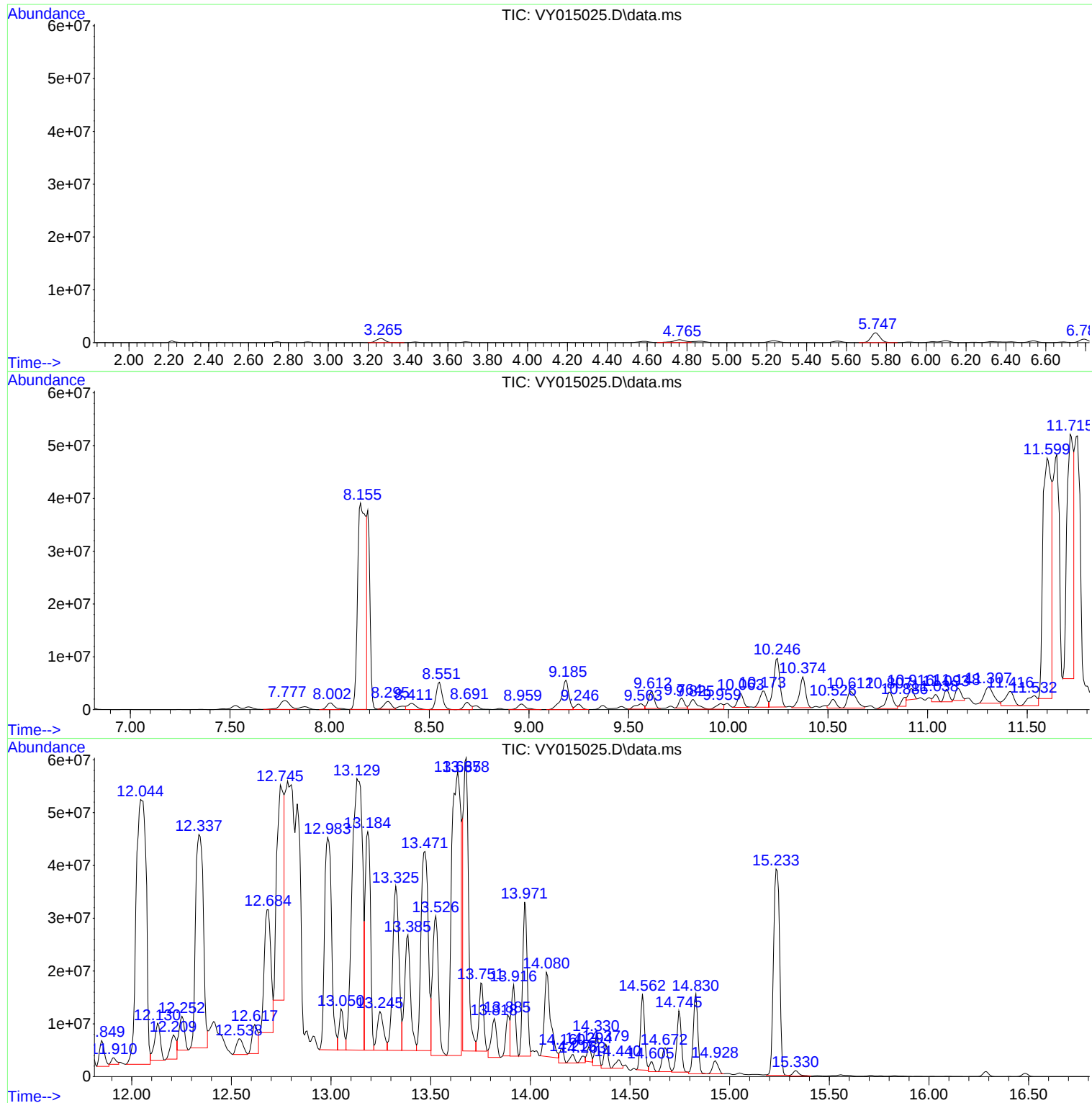
Sum of corrected areas: 2304588191

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
Quant Title : SW846 8260

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



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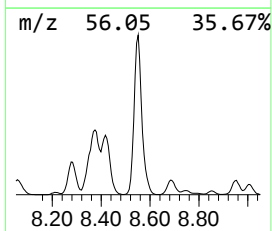
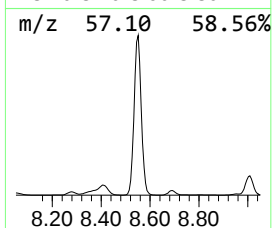
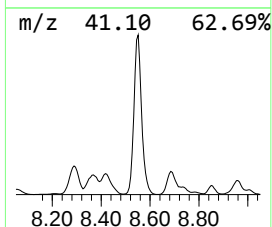
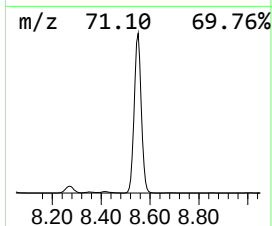
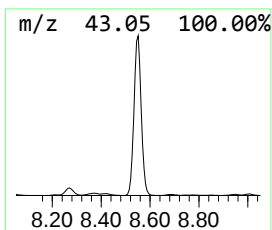
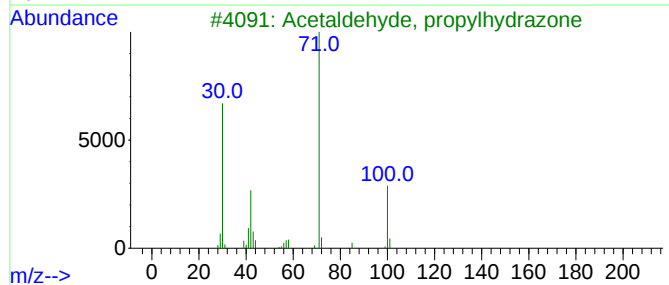
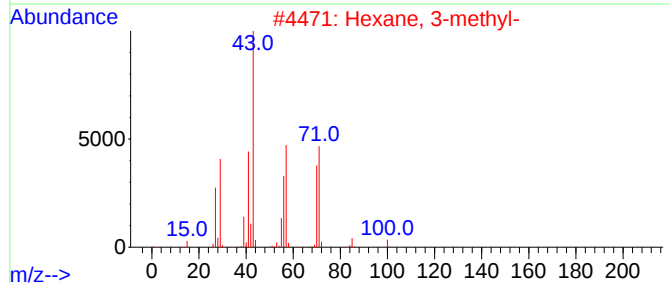
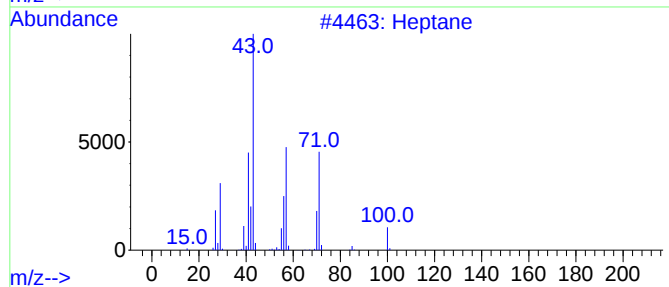
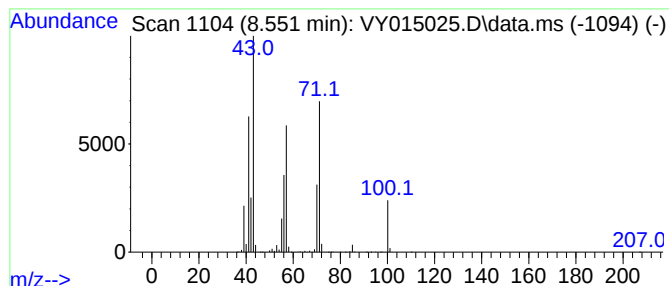
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.551	177.18 ug/l	11113000	1,4-Difluorobenzene	8.691

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	59
3			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	49
4			Pentane, 2-methyl-	86	C6H14	000107-83-5	43
5			3-Hexanone	100	C6H12O	000589-38-8	38



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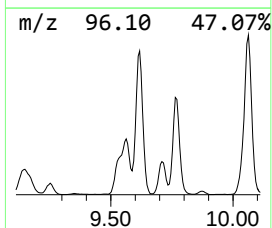
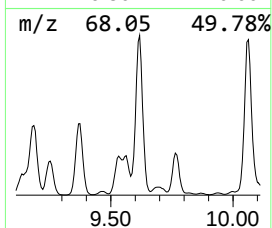
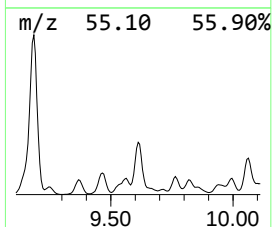
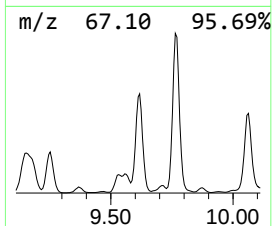
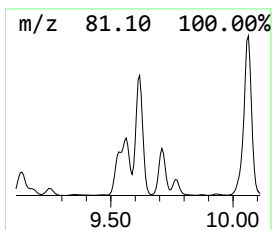
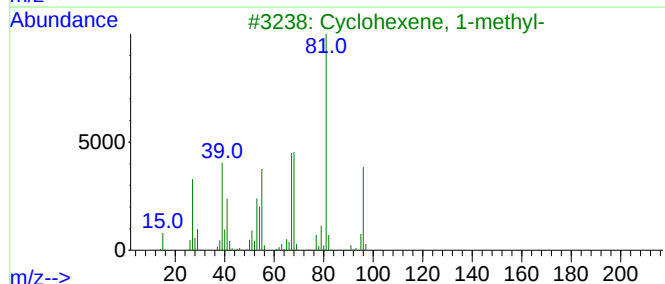
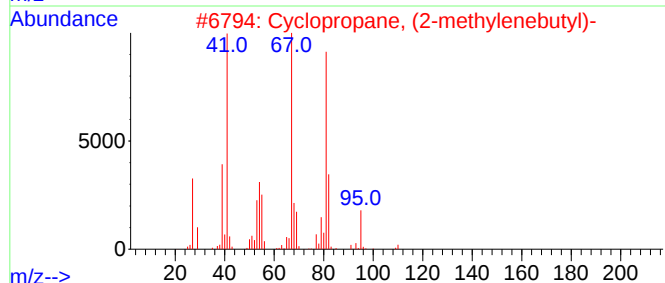
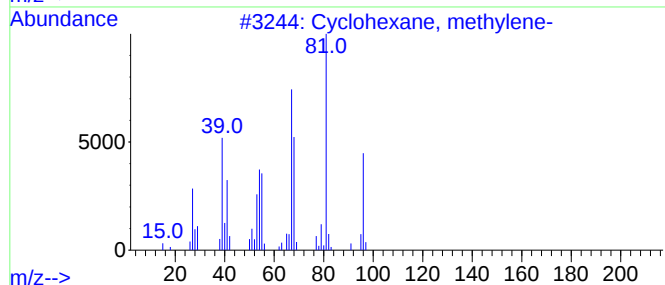
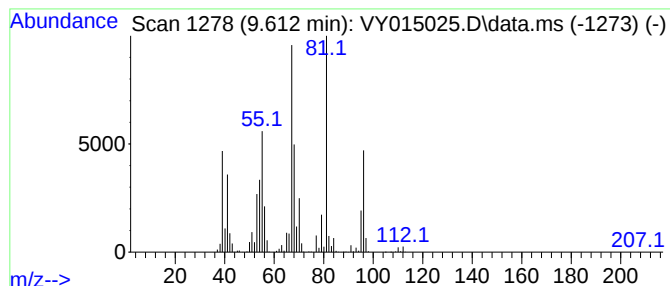
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, methylene- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.612	101.37 ug/l	6358070	1,4-Difluorobenzene	8.691

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, methylene-	96	C7H12	001192-37-6	95
2			Cyclopropane, (2-methylenebutyl)-	110	C8H14	074685-56-6	70
3			Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	68
4			Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	50
5			1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	43



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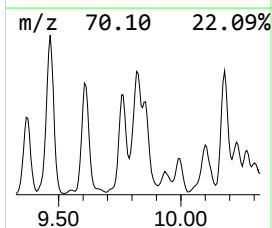
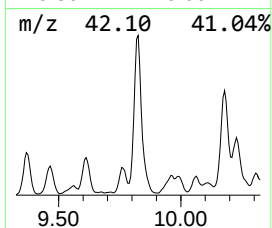
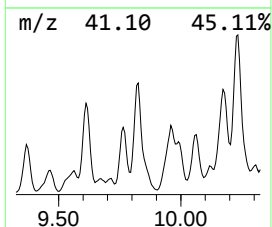
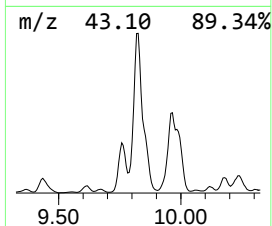
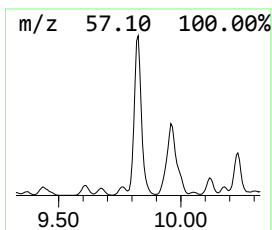
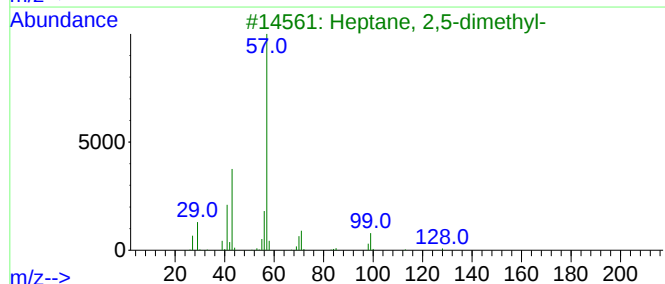
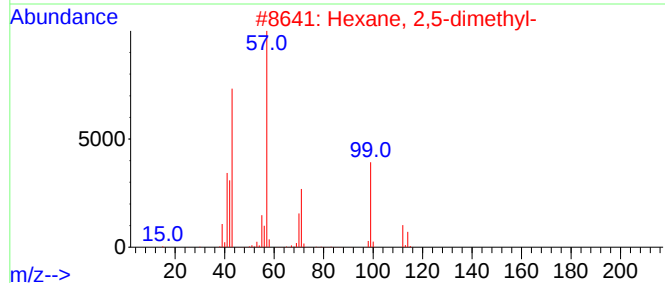
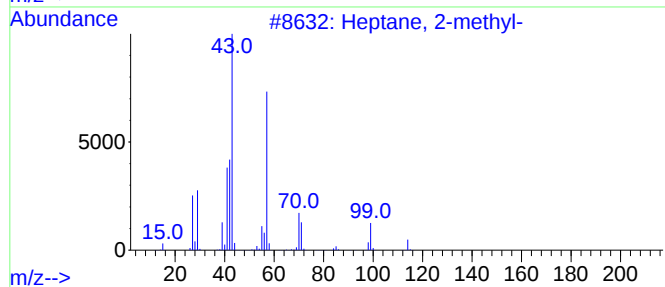
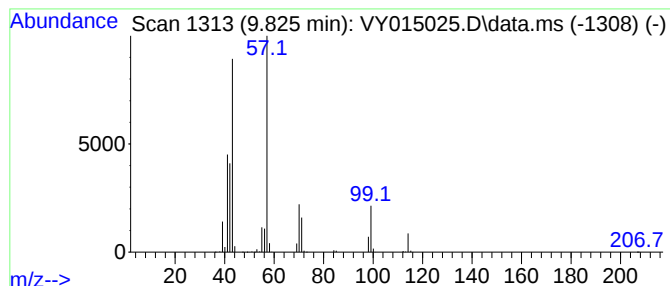
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
Quant Title : SW846 8260

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

Peak Number 3 Heptane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.825	76.07 ug/l	4771350	1,4-Difluorobenzene	8.691

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	64
3			Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	38
4			Hexane, 4-ethyl-2-methyl-	128	C9H20	003074-75-7	38
5			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	37



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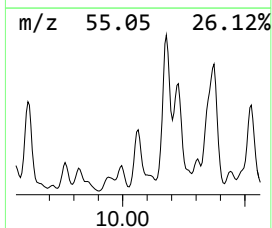
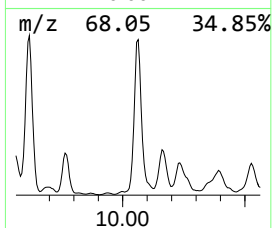
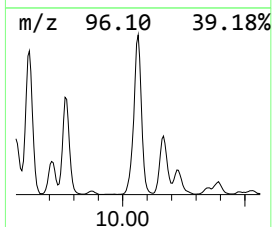
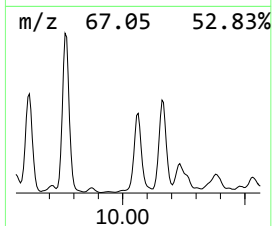
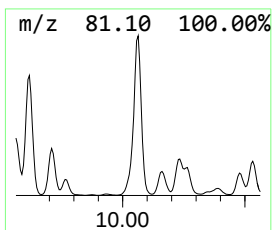
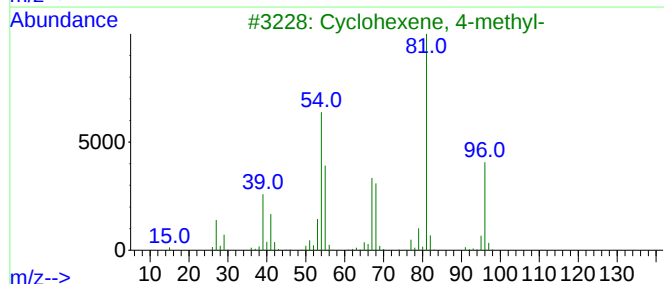
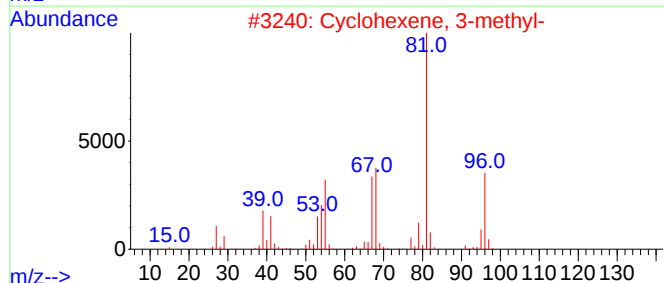
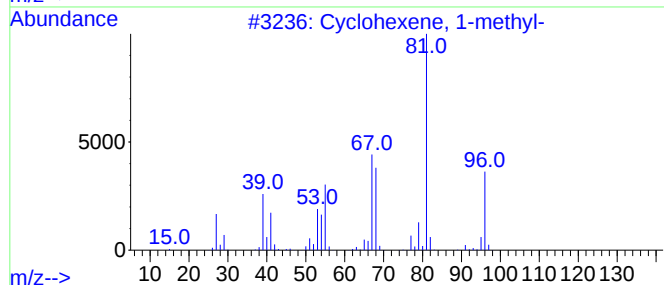
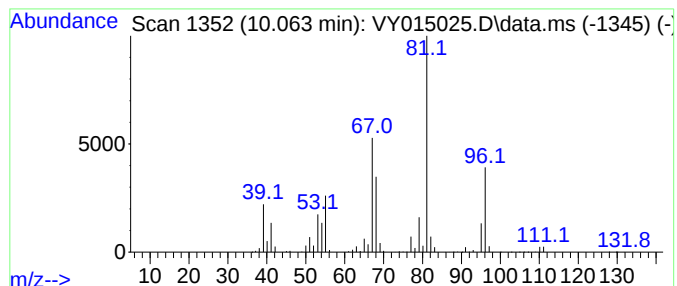
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.063	82.26 ug/l	5159140	1,4-Difluorobenzene	8.691

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	94
2			Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	87
3			Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	83
4			Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	50
5			Cyclohexane, methylene-	96	C7H12	001192-37-6	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

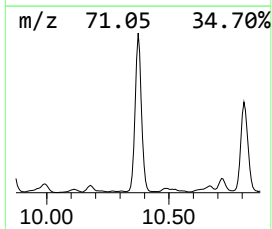
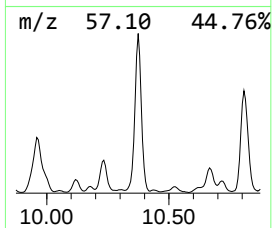
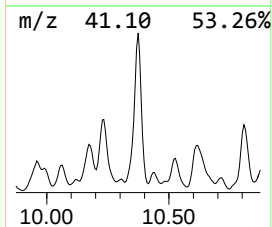
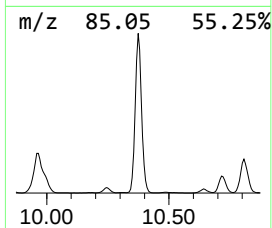
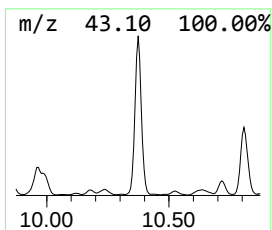
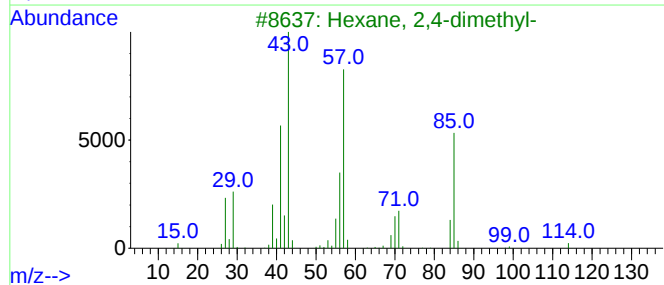
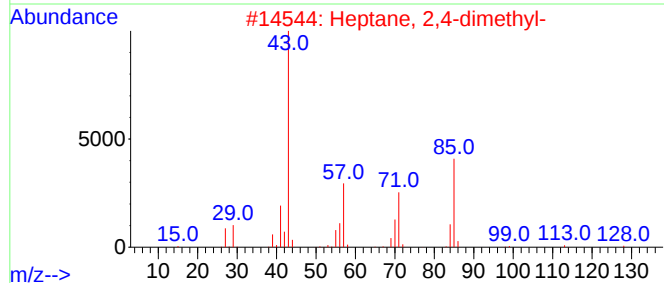
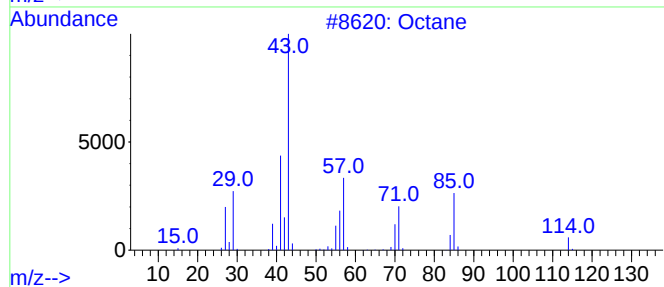
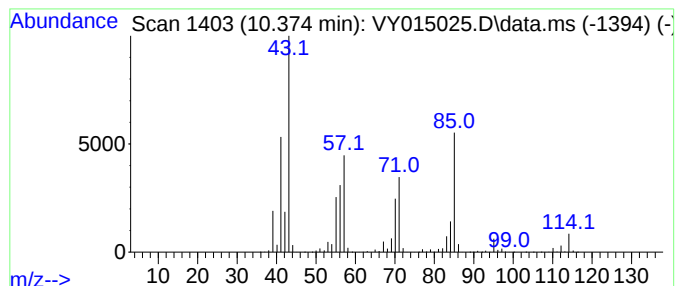
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Quant Title : SW846 8260

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

Peak Number 5 Octane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.374	93.20 ug/l	12292700	Chlorobenzene-d5	11.502

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane	114	C8H18	000111-65-9	91
2			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	72
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4			Oxalic acid, diisohexyl ester	258	C14H26O4	1010309-32-9	53
5			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SoIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

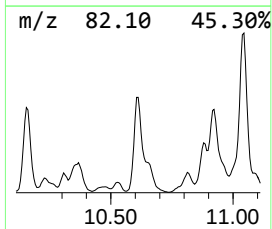
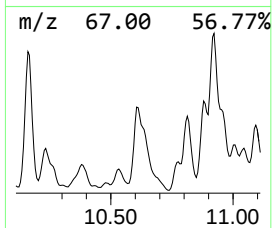
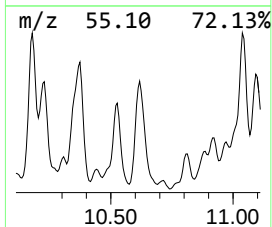
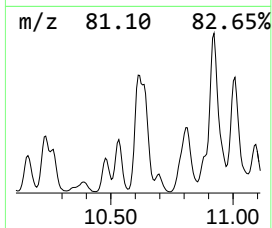
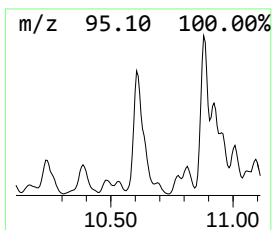
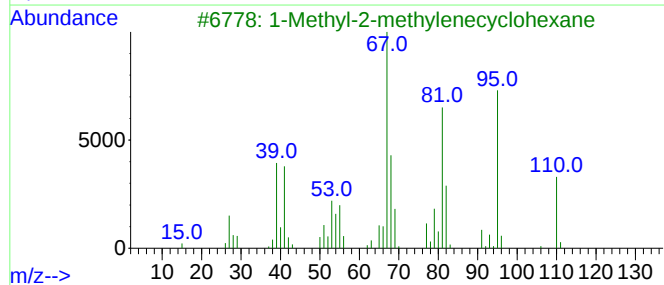
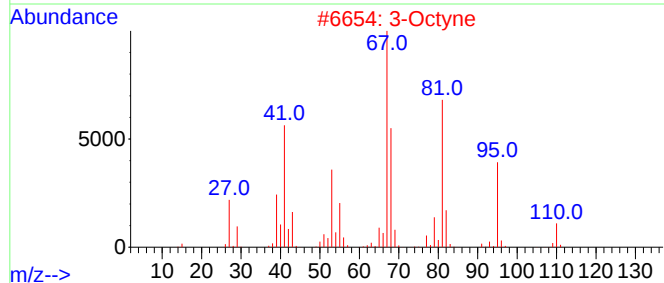
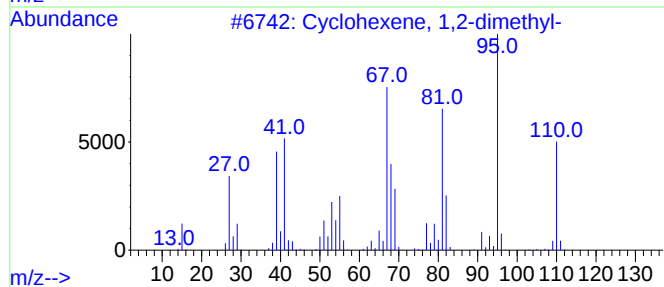
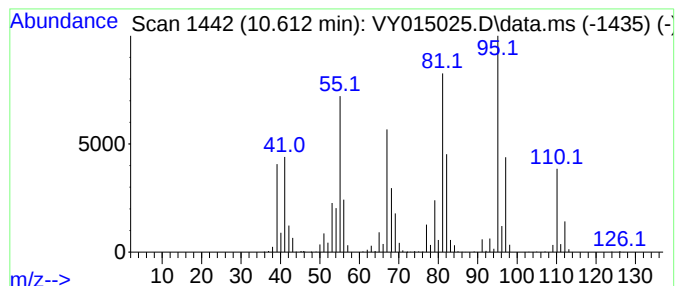
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Quant Title : SW846 8260

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

Peak Number 6 Cyclohexene, 1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.612	74.48 ug/l	9823100	Chlorobenzene-d5	11.502

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1,2-dimethyl-	110	C8H14	001674-10-8	90
2			3-Octyne	110	C8H14	015232-76-5	68
3			1-Methyl-2-methylenecyclohexane	110	C8H14	002808-75-5	58
4			Bicyclo[2.2.2]octane	110	C8H14	000280-33-1	53
5			Cyclopentane, 1,2-dimethyl-3-met...	110	C8H14	1000150-99-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

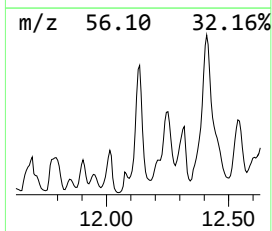
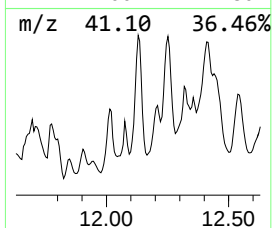
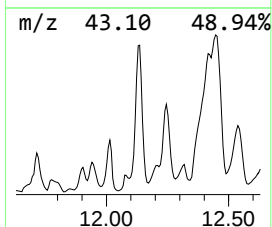
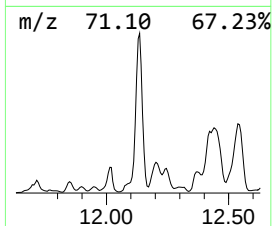
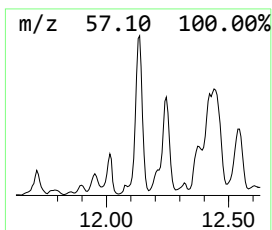
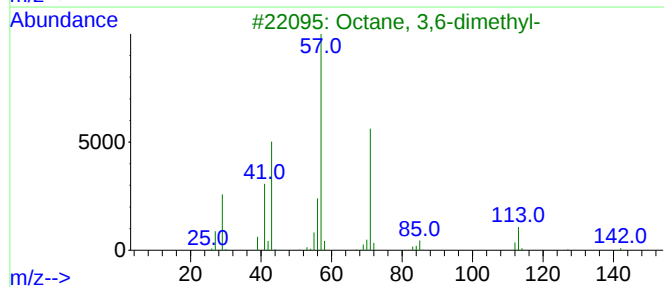
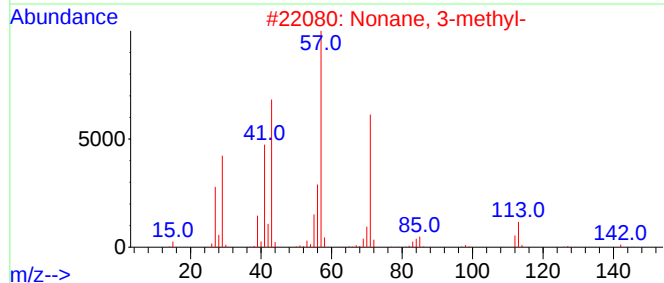
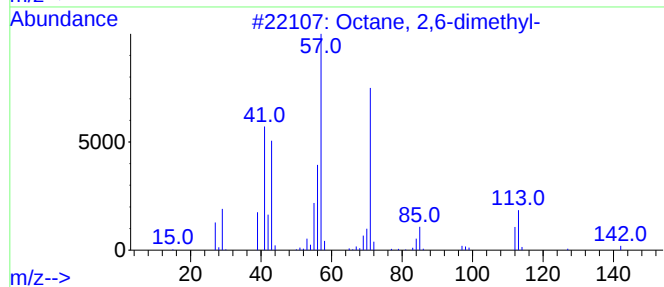
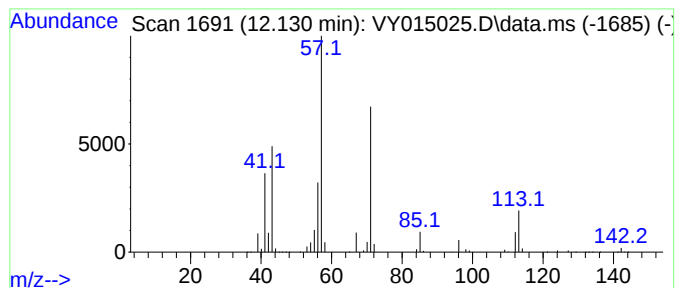
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Quant Title : SW846 8260

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

Peak Number 7 Octane, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.130	105.81 ug/l	13955300	Chlorobenzene-d5	11.502

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	83
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	76
4			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	59
5			Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

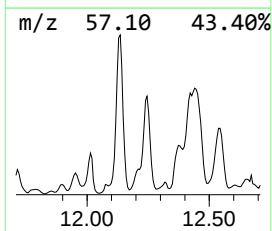
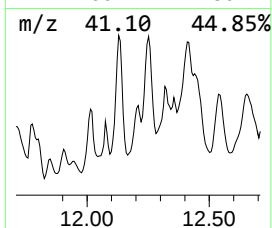
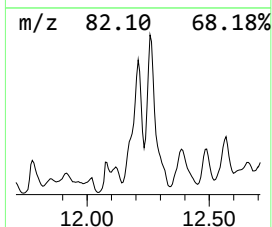
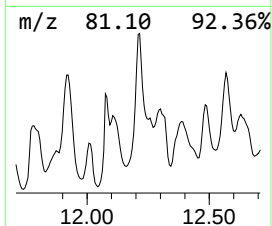
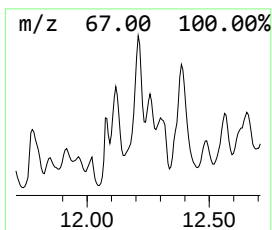
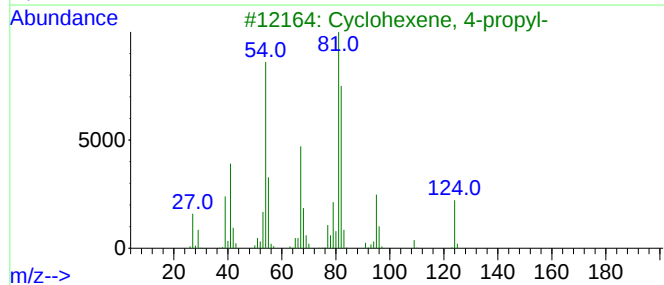
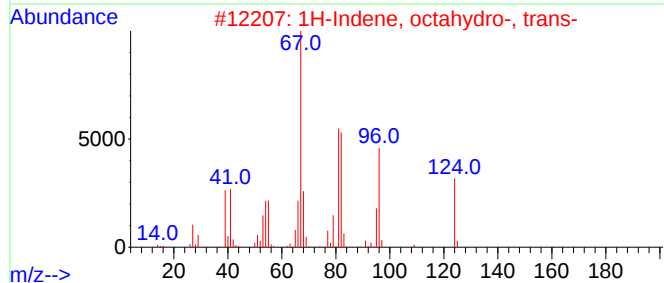
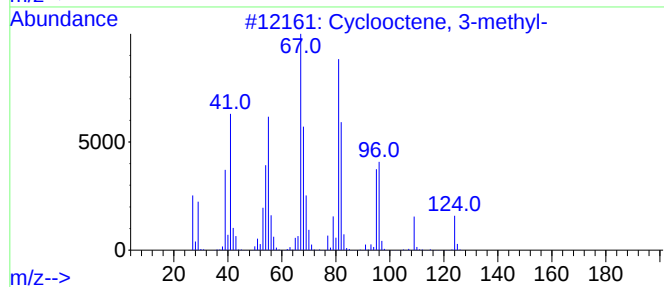
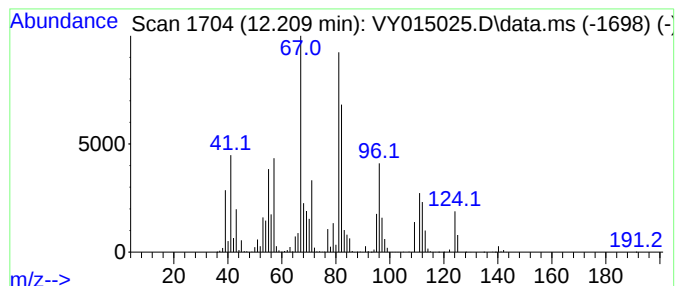
Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
Quant Title : SW846 8260

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

Peak Number 8 Cyclooctene, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.209	72.86 ug/l	9609690	Chlorobenzene-d5		11.502	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclooctene, 3-methyl-		124	C9H16	013152-05-1	72
2	1H-Indene, octahydro-, trans-		124	C9H16	003296-50-2	49
3	Cyclohexene, 4-propyl-		124	C9H16	013487-64-4	43
4	7-Methyl-8-oxabicyclo[4.3.0]nonane		140	C9H16O	1000216-87-6	43
5	1-Methylcyclooctene		124	C9H16	000933-11-9	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

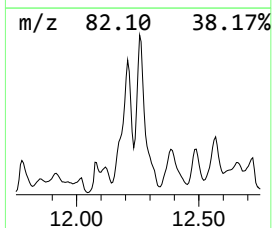
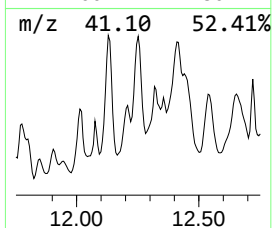
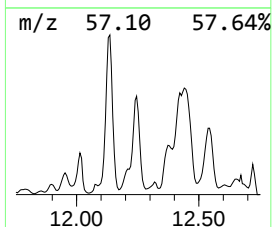
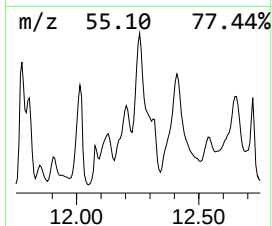
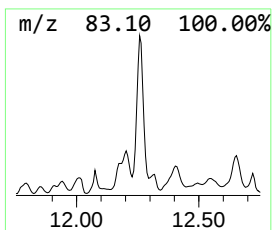
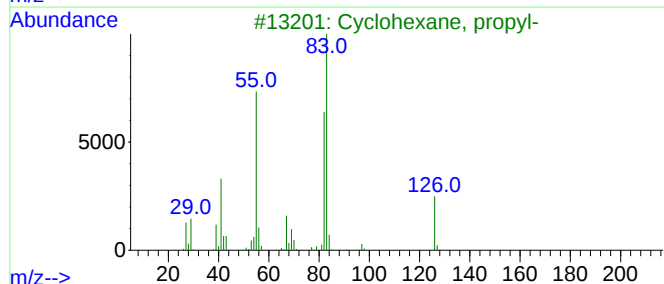
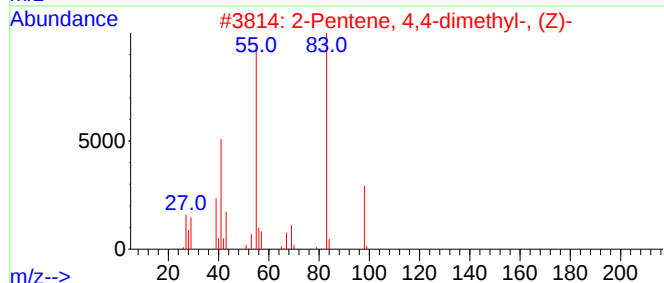
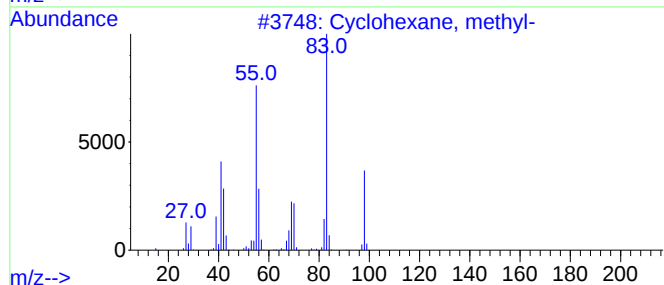
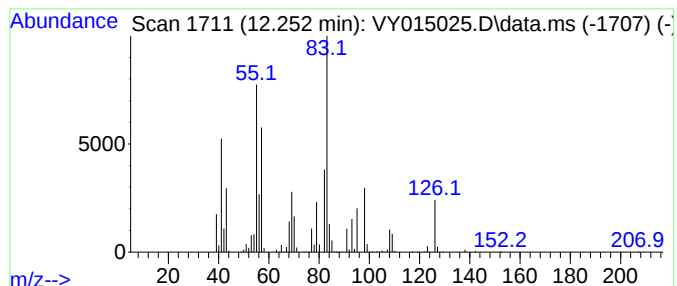
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Quant Title : SW846 8260

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

Peak Number 9 unknown12.252 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.252	87.65 ug/l	11559900	Chlorobenzene-d5	11.502

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, methyl-	98	C7H14	000108-87-2	49
2			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	43
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	43
4			2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	43
5			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

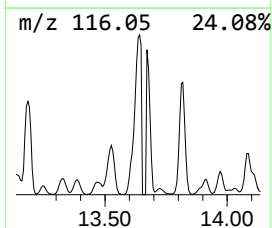
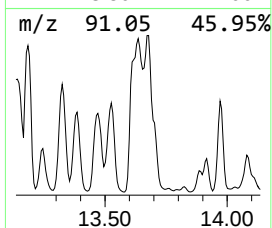
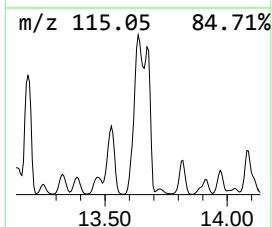
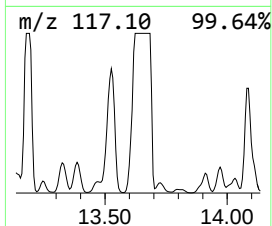
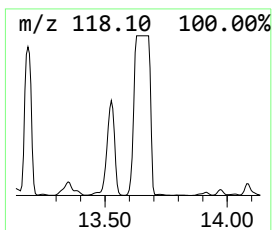
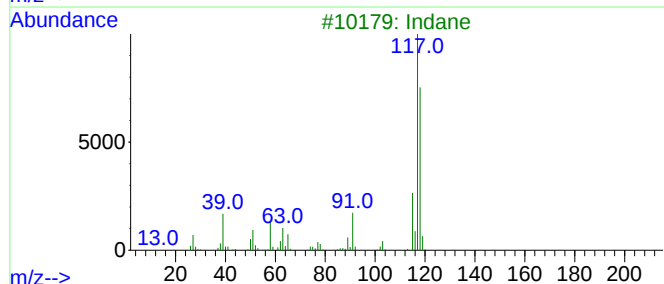
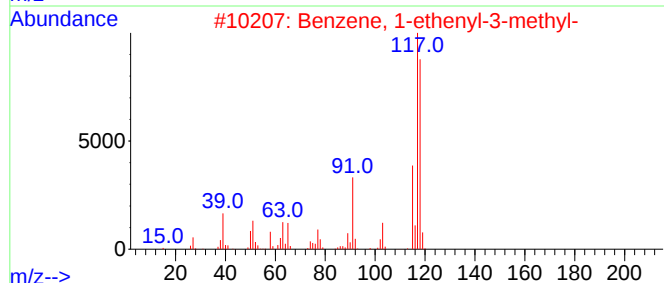
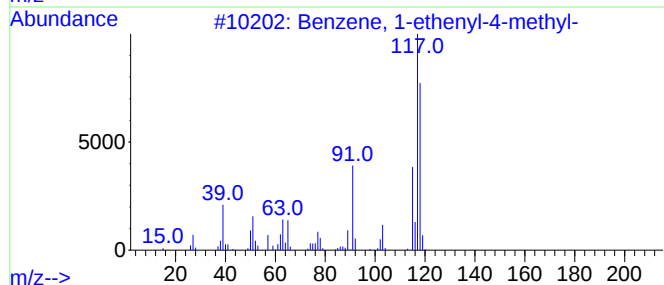
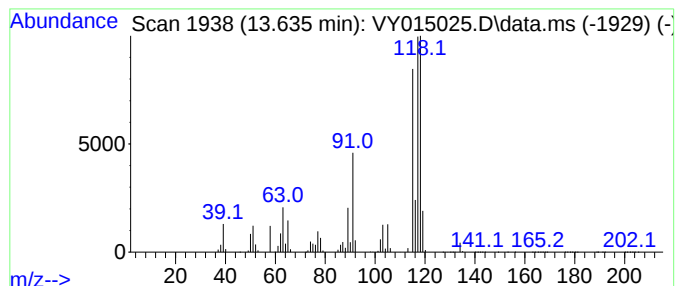
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethenyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.635	87.68 ug/l	169134000	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	55
2			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	55
3			Indane	118	C9H10	000496-11-7	53
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	50
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081023\
Data File : VY015025.D
Acq On : 10 Aug 2023 14:18
Operator : SY/MD
Sample : 03945-02
Misc : 6.78g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NEVINS-SB01-Z78-80

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y080723S.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	8.551	177.2	ug/l	11113000	2	8.691	3136040	50.0
Cyclohexane, me...	9.612	101.4	ug/l	6358070	2	8.691	3136040	50.0
Heptane, 2-methyl-	9.825	76.1	ug/l	4771350	2	8.691	3136040	50.0
Cyclohexene, 1-...	10.063	82.3	ug/l	5159140	2	8.691	3136040	50.0
Octane	10.374	93.2	ug/l	12292700	3	11.502	6594700	50.0
Cyclohexene, 1,...	10.612	74.5	ug/l	9823100	3	11.502	6594700	50.0
Octane, 2,6-dim...	12.130	105.8	ug/l	13955300	3	11.502	6594700	50.0
Cyclooctene, 3-...	12.209	72.9	ug/l	9609690	3	11.502	6594700	50.0
unknown12.252	12.252	87.7	ug/l	11559900	3	11.502	6594700	50.0
Benzene, 1-ethe...	13.635	87.7	ug/l	169134000	4	13.440	96444000	50.0