

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
 Data File : VY008385.D
 Acq On : 18 Apr 2022 19:50
 Operator : KP/MD
 Sample : N2452-07
 Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WC-15-1-4-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_Y\methods\82Y040822S.M
 Title : SW846 8260

Signal : TIC: VY008385.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.771	460	484	511	rVB3	131766	672083	1.57%	0.121%
2	5.240	546	561	581	rBV	125189	440928	1.03%	0.079%
3	6.795	806	816	837	rVB	228851	734867	1.72%	0.132%
4	7.783	971	978	986	rVB3	437060	1102933	2.58%	0.198%
5	7.874	986	993	1004	rVB	193335	497306	1.16%	0.089%
6	8.002	1004	1014	1020	rBV	291483	706918	1.65%	0.127%
7	8.325	1048	1067	1076	rBV	1032140	3300357	7.71%	0.593%
8	8.417	1076	1082	1093	rVB	370293	854593	2.00%	0.154%
9	8.691	1118	1127	1139	rBV	219751	476799	1.11%	0.086%
10	9.185	1192	1208	1214	rBV	2531217	5992174	14.01%	1.077%
11	9.246	1214	1218	1225	rVB	323536	604265	1.41%	0.109%
12	9.465	1244	1254	1261	rVB2	437646	1039695	2.43%	0.187%
13	9.612	1270	1278	1287	rVB3	919051	1954630	4.57%	0.351%
14	9.752	1288	1301	1306	rBV2	636235	1491852	3.49%	0.268%
15	9.941	1324	1332	1337	rBV3	522570	1296423	3.03%	0.233%
16	9.990	1337	1340	1347	rVB	323637	574546	1.34%	0.103%
17	10.112	1356	1360	1365	rVV2	346910	651653	1.52%	0.117%
18	10.179	1365	1371	1375	rVV	2681529	4955553	11.58%	0.890%
19	10.215	1375	1377	1385	rVB	991615	1462221	3.42%	0.263%
20	10.349	1394	1399	1410	rVB3	992934	2800793	6.55%	0.503%
21	10.483	1410	1421	1423	rBV2	521576	1055294	2.47%	0.190%
22	10.520	1423	1427	1436	rVB	1953021	3448409	8.06%	0.620%
23	10.618	1436	1443	1449	rBV3	961403	2199339	5.14%	0.395%
24	10.666	1449	1451	1455	rVV2	338061	477124	1.12%	0.086%
25	10.715	1455	1459	1463	rVB	627350	946499	2.21%	0.170%
26	10.807	1469	1474	1479	rVB	1286851	2133222	4.99%	0.383%
27	10.910	1479	1491	1497	rBV	876771	2203924	5.15%	0.396%
28	10.971	1497	1501	1504	rBV2	661463	1069108	2.50%	0.192%
29	11.038	1508	1512	1516	rVB	1830517	2598968	6.07%	0.467%
30	11.093	1516	1521	1527	rVB	7004625	10862014	25.39%	1.952%
31	11.148	1527	1530	1534	rVB	597794	748517	1.75%	0.134%
32	11.203	1534	1539	1544	rBV3	1394025	2626880	6.14%	0.472%
33	11.300	1550	1555	1563	rVB	2099648	3840952	8.98%	0.690%
34	11.380	1563	1568	1572	rBV2	713312	1274919	2.98%	0.229%
35	11.483	1581	1585	1589	rBV2	292710	495104	1.16%	0.089%
36	11.581	1596	1601	1605	rBV4	315926	545167	1.27%	0.098%

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
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37	11.630	1605	1609	1613	rBV	1199736	1709009	3.99%	0.307%
38	11.691	1613	1619	1623	rBV5	2035282	4277840	10.00%	0.769%
39	11.739	1623	1627	1632	rBV	2855638	5234947	12.24%	0.941%
40	11.843	1639	1644	1648	rBV3	802631	1583479	3.70%	0.285%
41	11.898	1648	1653	1656	rVV2	635476	1088945	2.55%	0.196%
42	11.941	1656	1660	1664	rVB2	996324	1532040	3.58%	0.275%
43	12.008	1664	1671	1678	rBV5	5249557	12761557	29.83%	2.293%
44	12.062	1678	1680	1682	rVV2	584479	622143	1.45%	0.112%
45	12.093	1682	1685	1687	rVV	647932	773297	1.81%	0.139%
46	12.130	1687	1691	1695	rVB	5490448	6826880	15.96%	1.227%
47	12.172	1695	1698	1699	rBV2	635340	800740	1.87%	0.144%
48	12.209	1699	1704	1706	rVV2	4614980	7370865	17.23%	1.324%
49	12.245	1706	1710	1715	rVB2	7324127	13143988	30.72%	2.362%
50	12.379	1727	1732	1738	rVB5	2538483	4988173	11.66%	0.896%
51	12.471	1743	1747	1748	rBV	1933772	2263861	5.29%	0.407%
52	12.575	1756	1764	1768	rBV7	2463452	6968822	16.29%	1.252%
53	12.648	1769	1776	1784	rVB3	9749530	20200747	47.21%	3.630%
54	12.733	1784	1790	1796	rBV4	3412782	8704262	20.34%	1.564%
55	12.818	1796	1804	1809	rBV5	12367758	25274088	59.07%	4.541%
56	12.959	1822	1827	1831	rBV5	2486135	4137188	9.67%	0.743%
57	13.014	1834	1836	1837	rBV	637436	528257	1.23%	0.095%
58	13.050	1838	1842	1843	rBV2	4692805	6066390	14.18%	1.090%
59	13.074	1843	1846	1851	rVB	11605230	16445919	38.44%	2.955%
60	13.123	1851	1854	1859	rVB	7019919	9882704	23.10%	1.776%
61	13.184	1859	1864	1869	rBV	6908390	11317690	26.45%	2.033%
62	13.288	1869	1881	1882	rBV3	19307108	31547414	73.73%	5.668%
63	13.361	1891	1893	1897	rVB4	1530406	1756651	4.11%	0.316%
64	13.404	1897	1900	1904	rVB	4930208	6815014	15.93%	1.224%
65	13.477	1904	1912	1914	rBV3	17717608	30510888	71.31%	5.482%
66	13.507	1915	1917	1924	rVB	17412070	22766087	53.21%	4.090%
67	13.574	1924	1928	1930	rBV4	2129261	3421792	8.00%	0.615%
68	13.599	1930	1932	1934	rBV	850475	877410	2.05%	0.158%
69	13.654	1934	1941	1949	rVB5	16838010	42784920	100.00%	7.687%
70	13.727	1949	1953	1955	rBV2	1916423	2675015	6.25%	0.481%
71	13.757	1955	1958	1961	rBV2	2409089	2822108	6.60%	0.507%
72	13.794	1961	1964	1966	rBV3	3657234	4872672	11.39%	0.875%
73	13.830	1966	1970	1976	rVB2	8005038	16019527	37.44%	2.878%
74	13.885	1976	1979	1982	rVB3	2771182	3563944	8.33%	0.640%
75	13.928	1982	1986	1987	rBV2	4343057	5527425	12.92%	0.993%
76	13.958	1987	1991	1995	rBV2	6939302	13500061	31.55%	2.426%

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77	14.001	1995	1998	2007	rVB3	15490603	31903488	74.57%	5.732%
78	14.080	2007	2011	2015	rBV4	2241502	3862396	9.03%	0.694%
79	14.147	2015	2022	2023	rBV	9135041	14942383	34.92%	2.685%
80	14.263	2037	2041	2043	rBV2	2117659	3273522	7.65%	0.588%
81	14.306	2044	2048	2052	rBV4	2134202	4024793	9.41%	0.723%
82	14.385	2056	2061	2063	rBV2	17572571	29787133	69.62%	5.352%
83	14.483	2073	2077	2080	rBV	2159070	3119063	7.29%	0.560%
84	14.525	2080	2084	2088	rVV	7213609	11151855	26.06%	2.004%
85	14.568	2088	2091	2096	rVB2	3497990	5489767	12.83%	0.986%
86	14.635	2096	2102	2105	rBV5	1304223	2885625	6.74%	0.518%
87	14.678	2105	2109	2114	rVV3	1705922	4415296	10.32%	0.793%
88	14.751	2118	2121	2125	rVV4	1742075	2530555	5.91%	0.455%
89	14.818	2129	2132	2136	rVV	931297	1450975	3.39%	0.261%
90	14.861	2136	2139	2149	rVB3	1247742	2936912	6.86%	0.528%
91	14.946	2149	2153	2157	rVB5	470560	661271	1.55%	0.119%
92	14.995	2157	2161	2168	rBV5	340847	917058	2.14%	0.165%
93	15.227	2195	2199	2205	rVB	682164	1111135	2.60%	0.200%

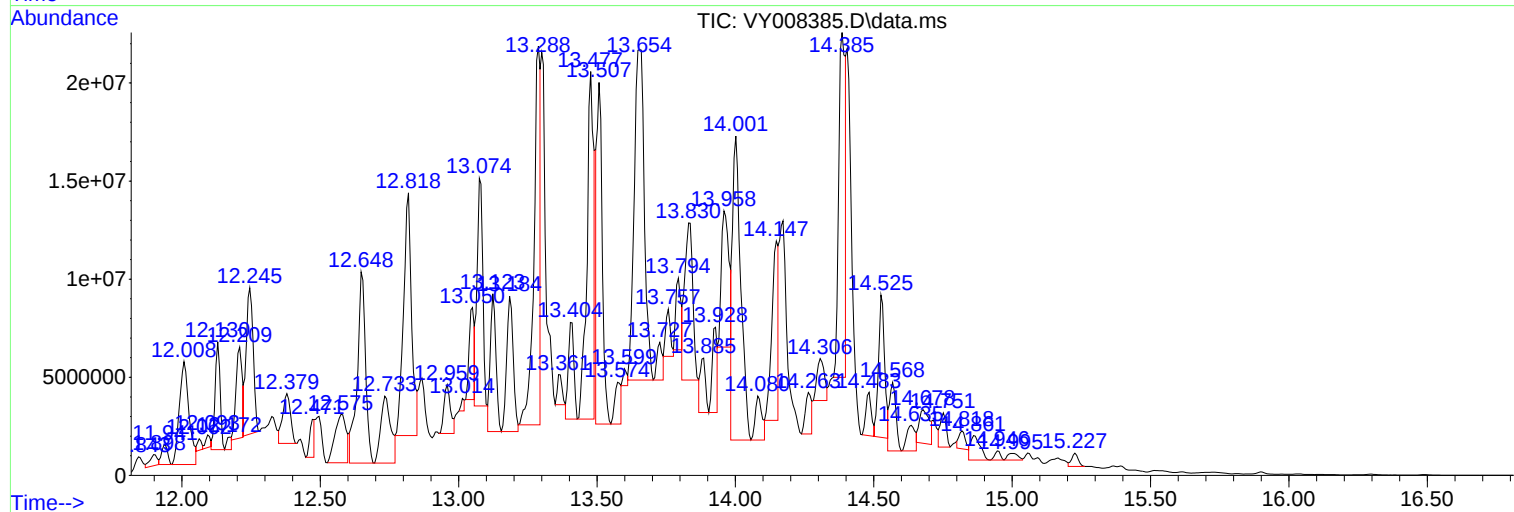
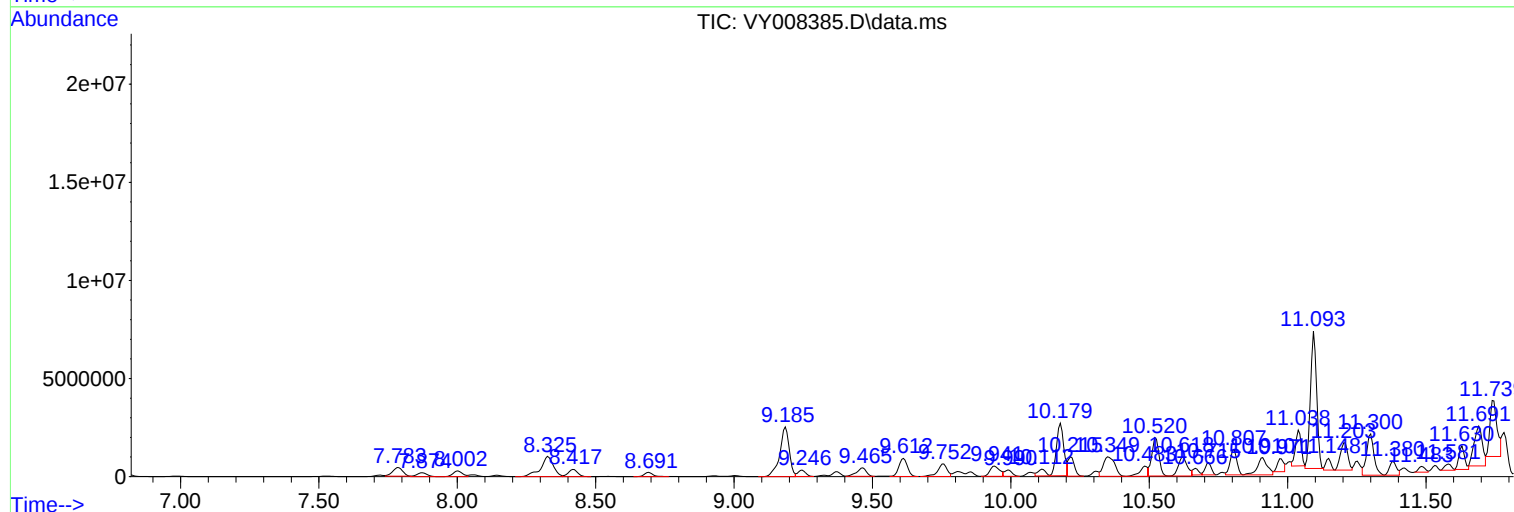
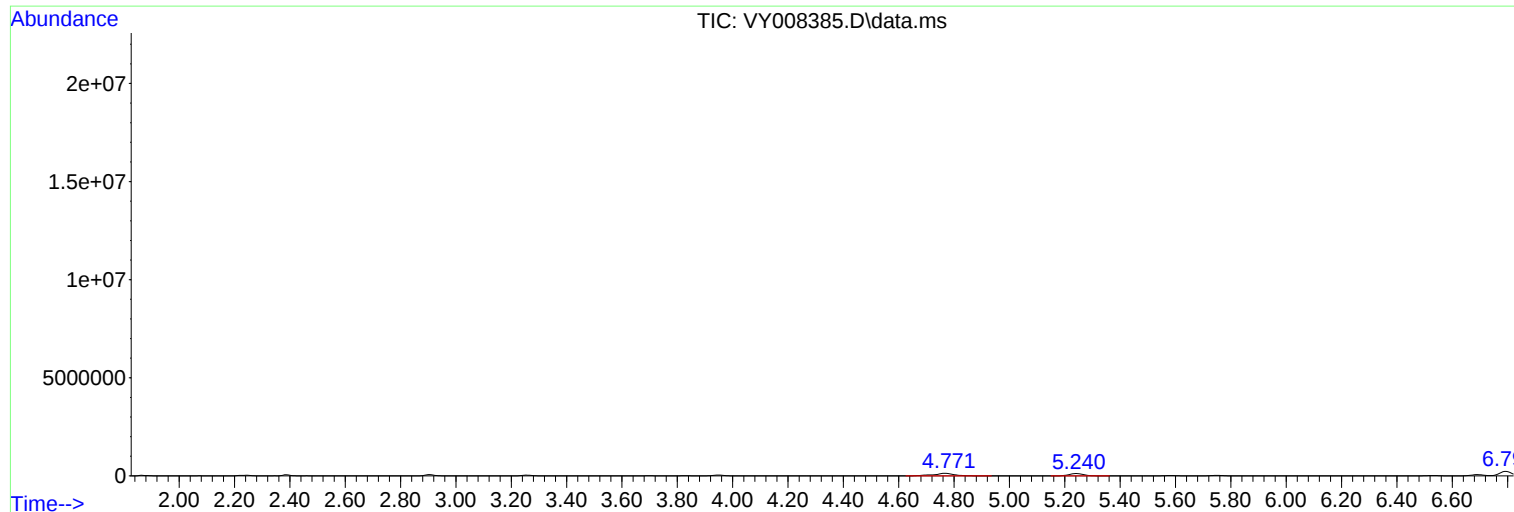
Sum of corrected areas: 556564015

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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



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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

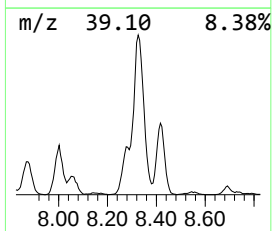
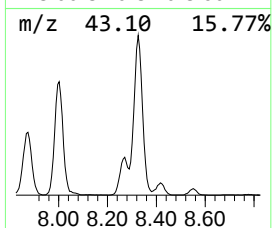
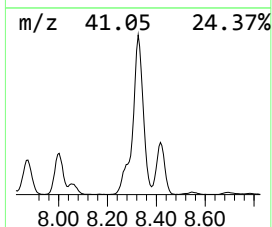
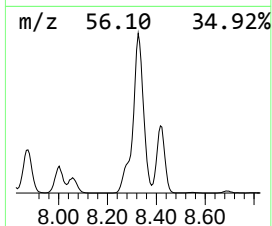
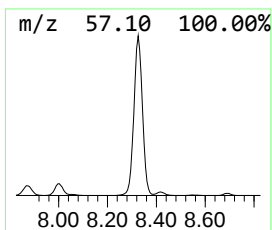
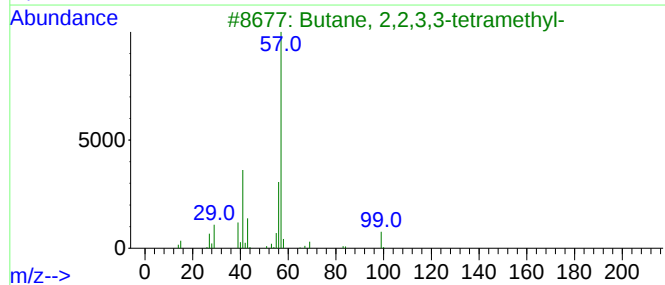
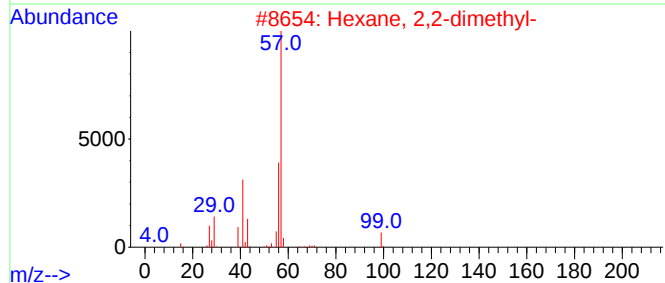
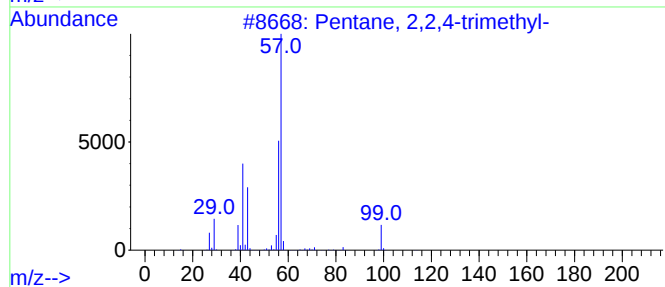
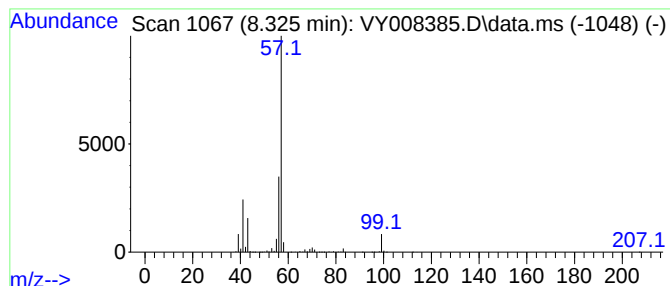
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.325	346.10 ug/l	3300360	1,4-Difluorobenzene	8.691

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
4			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	59
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



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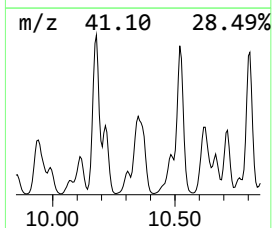
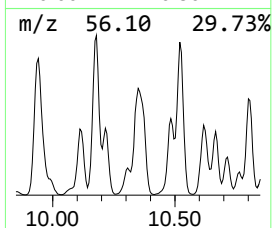
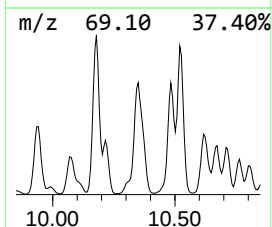
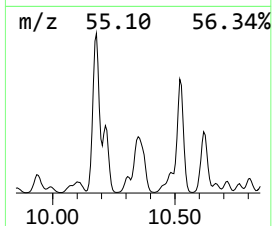
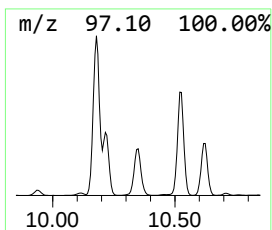
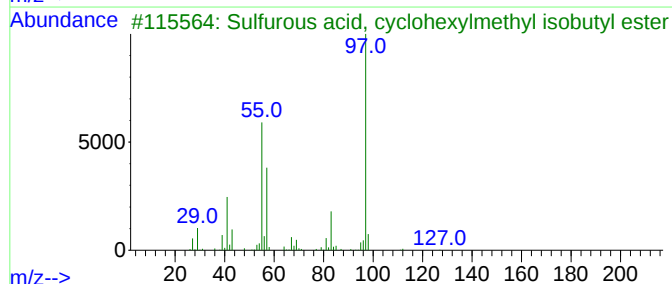
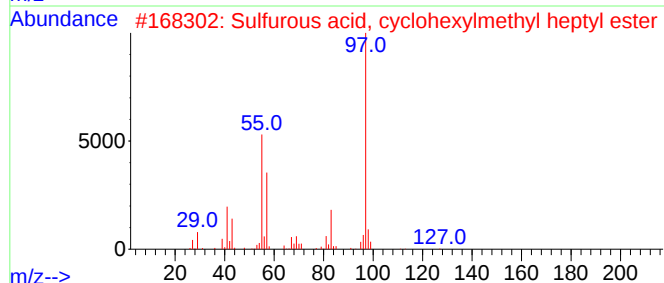
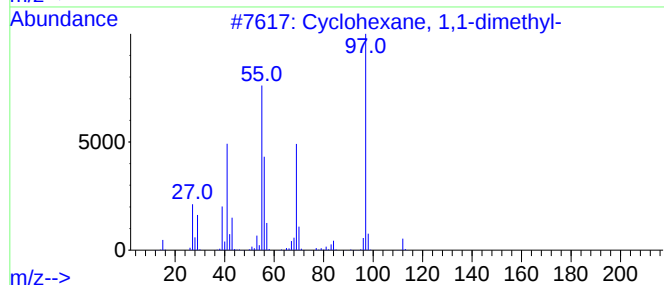
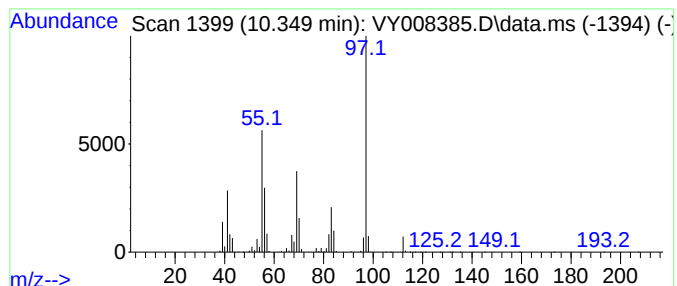
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, 1,1-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.349	282.85 ug/l	2800790	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	81
2			Sulfurous acid, cyclohexylmethyl...	276	C14H28O3S	1000309-21-7	59
3			Sulfurous acid, cyclohexylmethyl...	234	C11H22O3S	1000309-21-3	59
4			1H-Pyrazole, 4,5-dihydro-3,5,5-t...	112	C6H12N2	003975-85-7	53
5			Cycloheptane, methyl-	112	C8H16	004126-78-7	50



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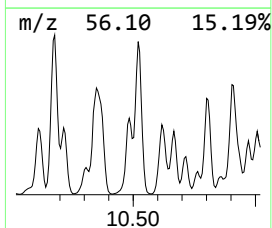
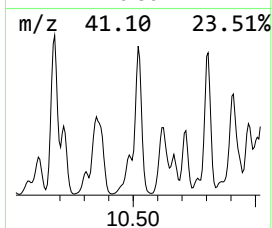
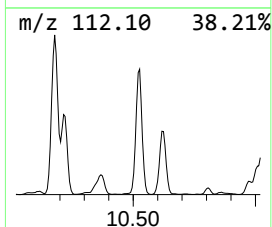
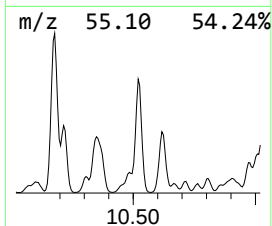
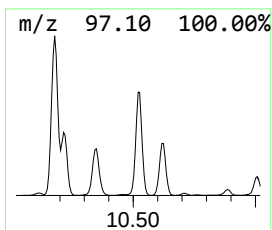
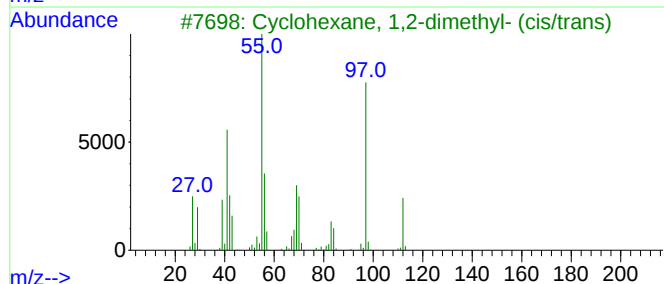
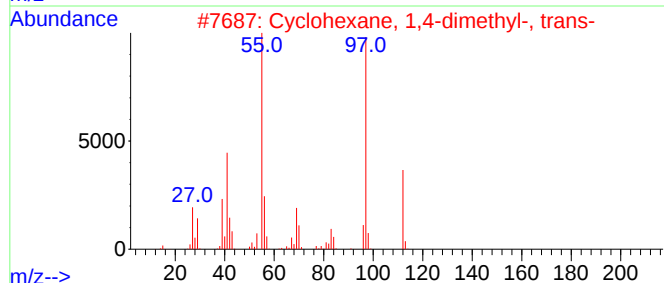
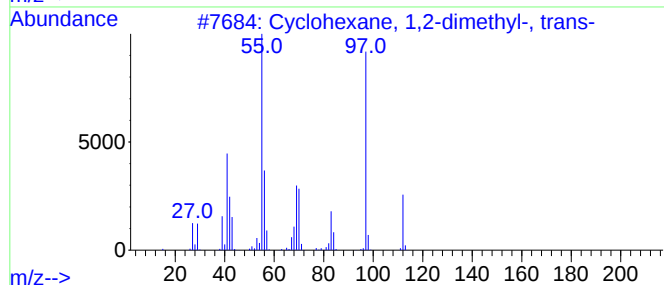
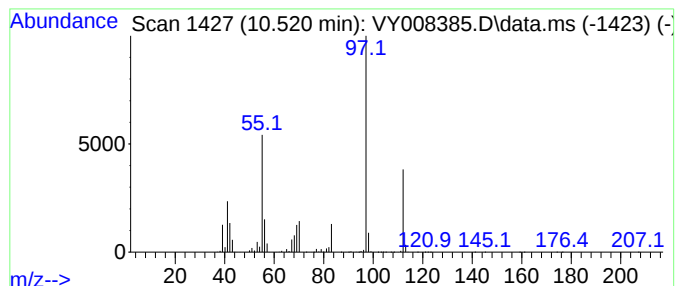
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.520	348.25 ug/l	3448410	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	91
2			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	91
3			Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	91
4			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	91
5			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

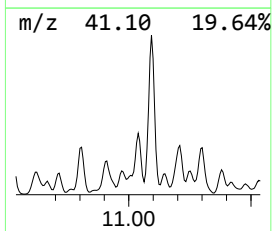
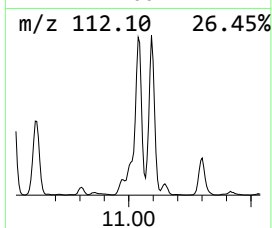
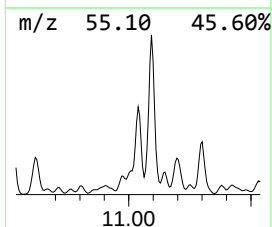
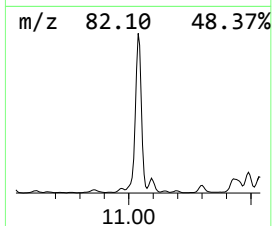
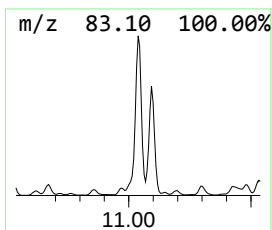
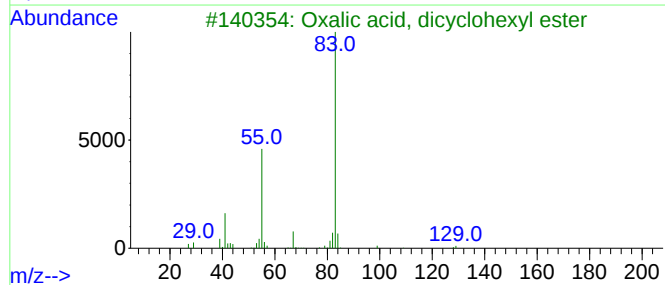
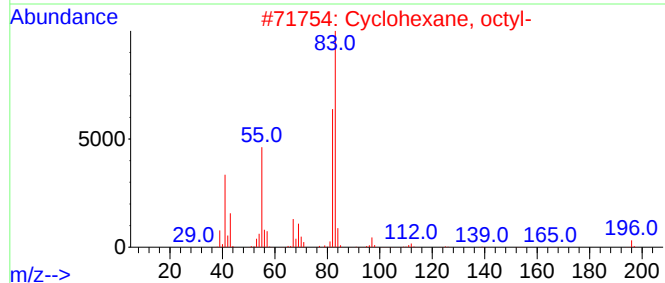
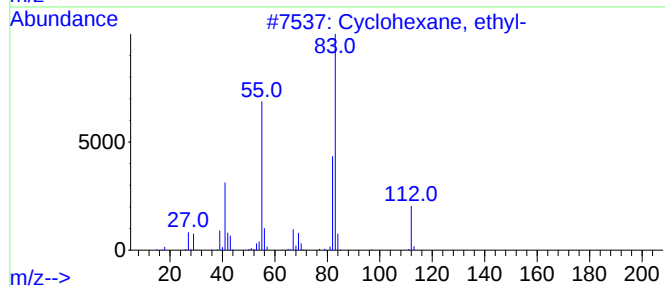
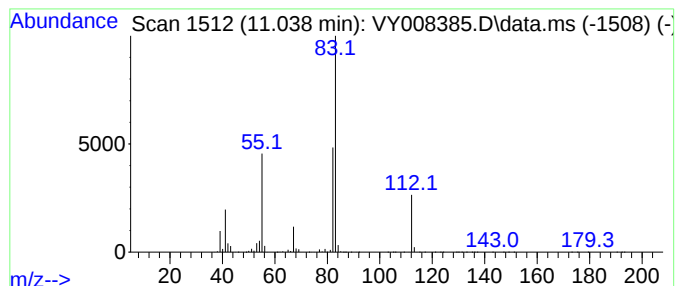
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.038	262.47 ug/l	2598970	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	83
2			Cyclohexane, octyl-	196	C14H28	001795-15-9	56
3			Oxalic acid, dicyclohexyl ester	254	C14H22O4	1000309-30-9	50
4			Cyclohexane, bromo-	162	C6H11Br	000108-85-0	47
5			Cyclopropane, 2-bromo-1,1,3-trim...	162	C6H11Br	036617-00-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

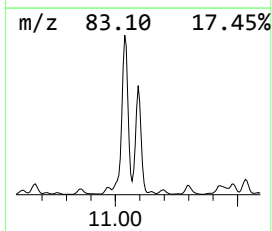
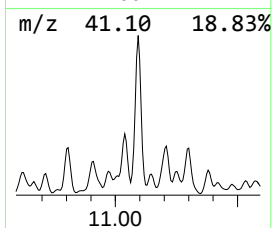
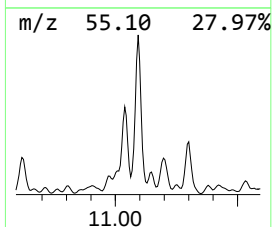
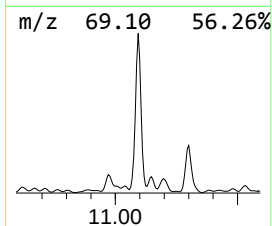
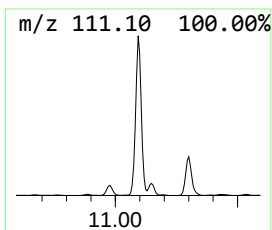
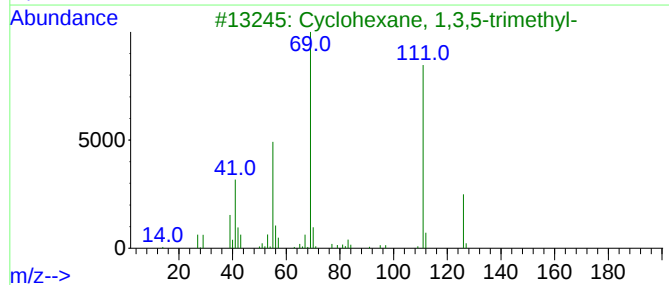
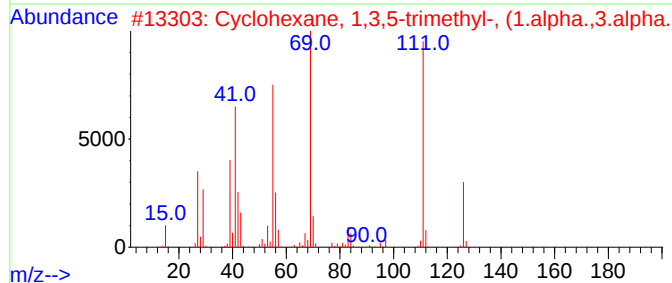
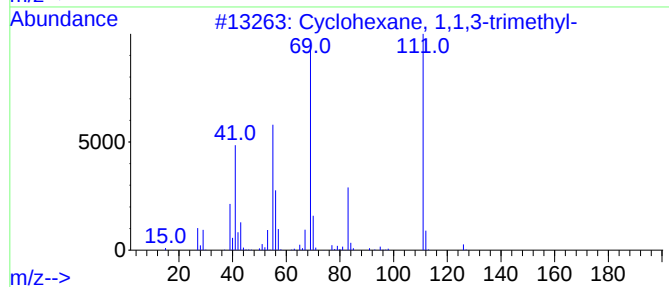
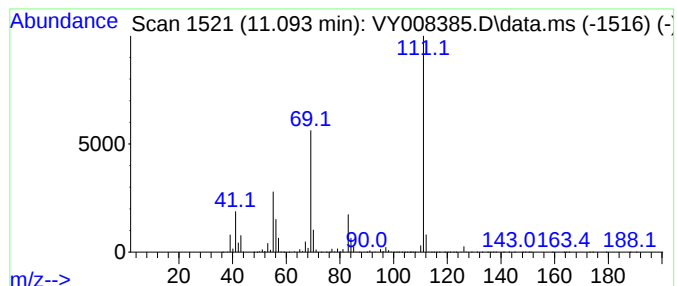
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.093	1096.94 ug/l	10862000	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	91
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	78
3			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
4			Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	72
5			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

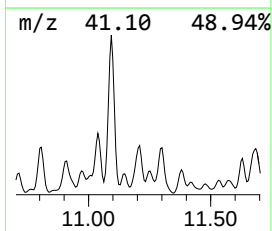
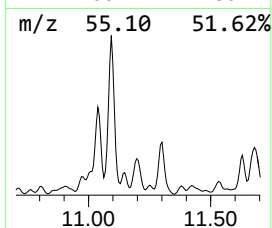
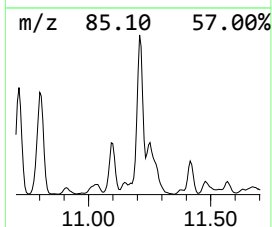
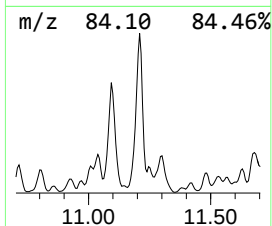
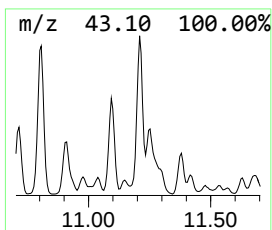
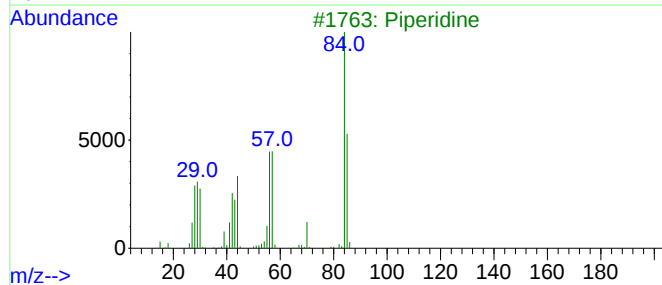
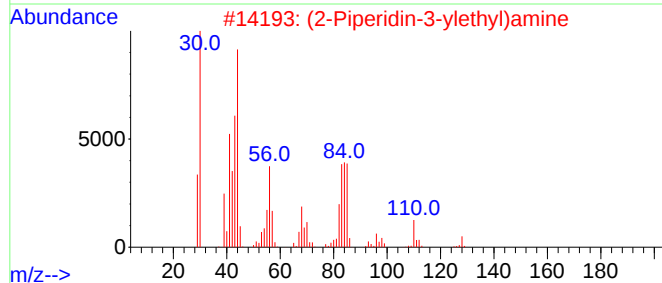
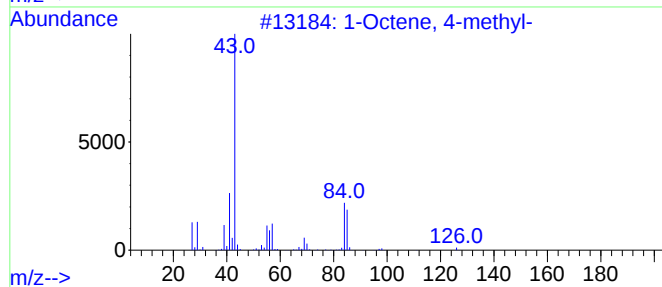
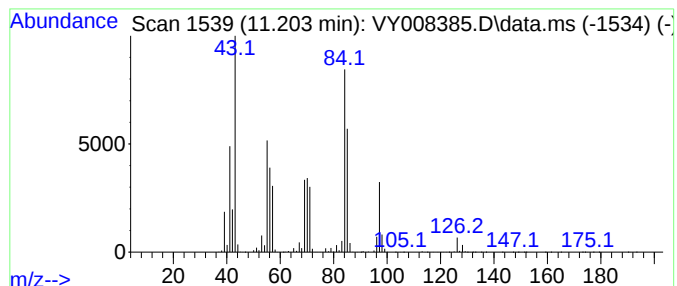
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Quant Title : SW846 8260

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

Peak Number 6 1-Octene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.203	265.29 ug/l	2626880	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octene, 4-methyl-	126	C9H18	013151-12-7	64
2			(2-Piperidin-3-ylethyl)amine	128	C7H16N2	089850-94-2	64
3			Piperidine	85	C5H11N	000110-89-4	59
4			1-Fluorononane	146	C9H19F	000463-18-3	43
5			Cyclopentanone	84	C5H8O	000120-92-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

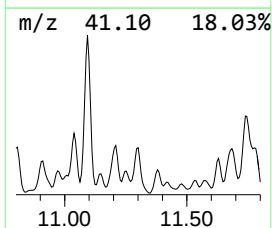
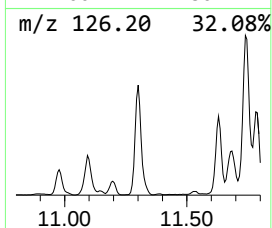
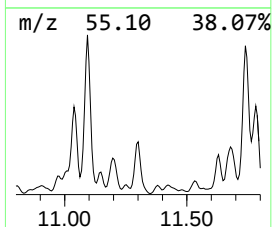
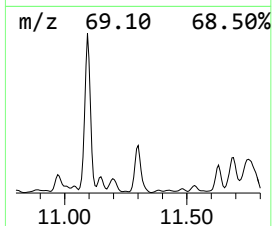
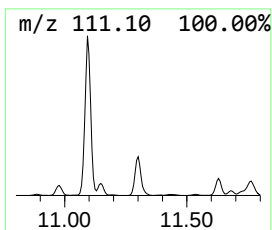
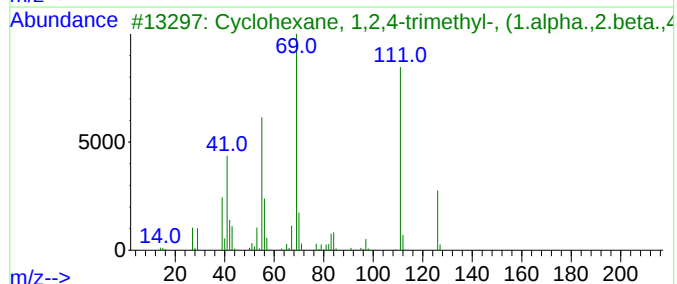
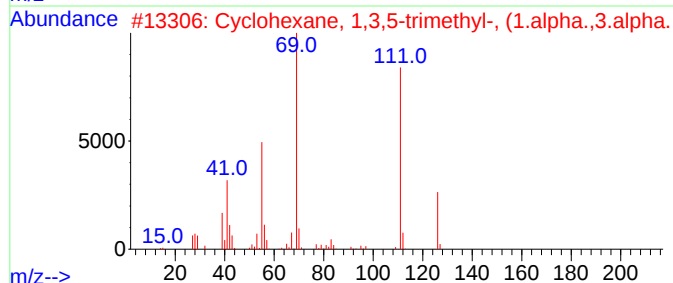
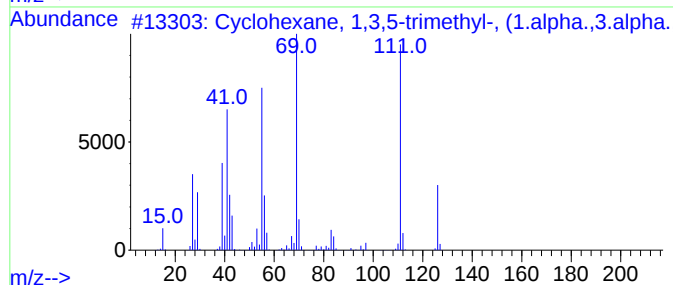
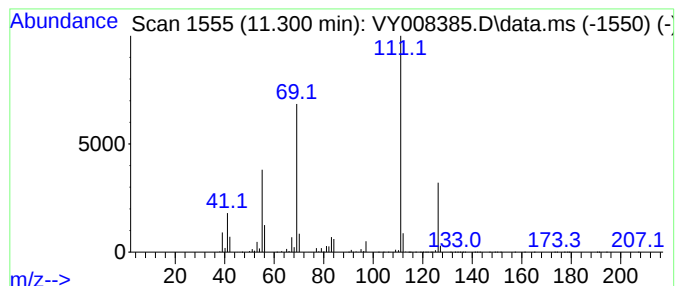
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, 1,3,5-trimethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.300	387.89 ug/l	3840950	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	91
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	91
3			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	91
4			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	90
5			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

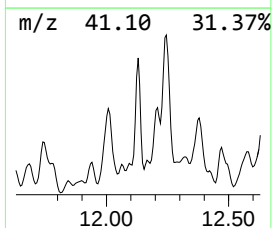
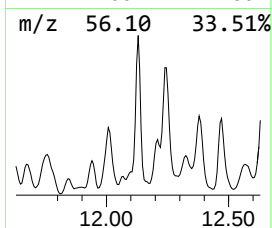
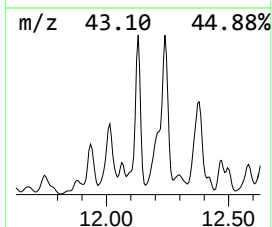
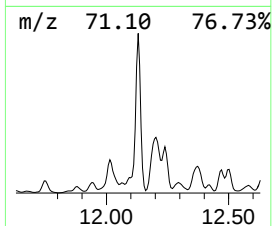
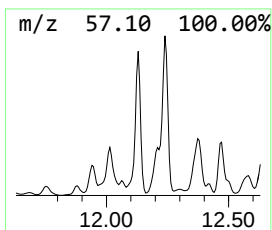
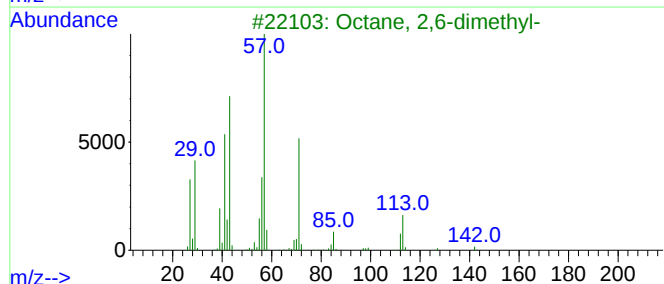
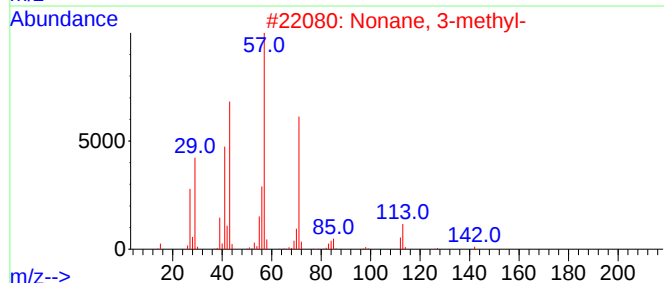
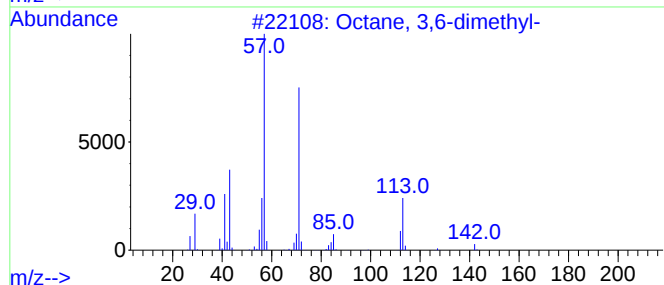
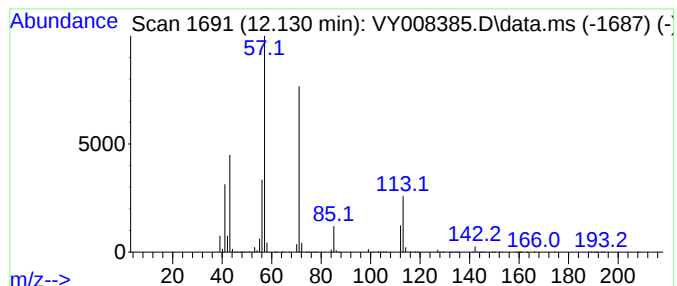
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Quant Title : SW846 8260

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

Peak Number 8 Octane, 3,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.130	689.44 ug/l	6826880	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	91
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	91
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	91
4			Octyl thioglycolate	204	C10H20O2S	007664-80-4	78
5			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WC-15-1-4-GR

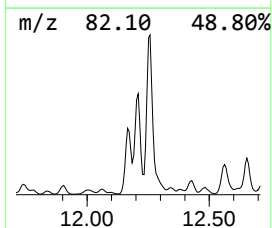
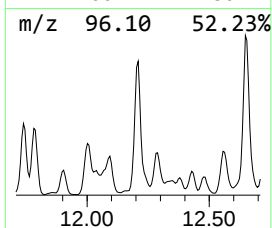
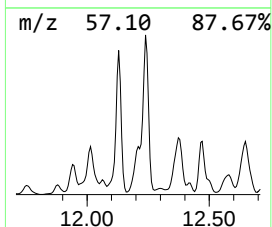
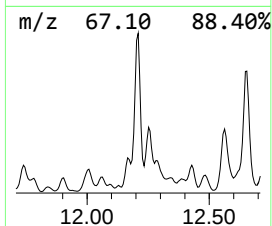
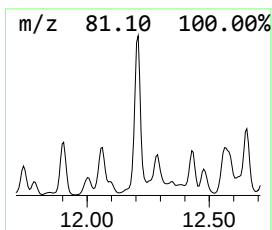
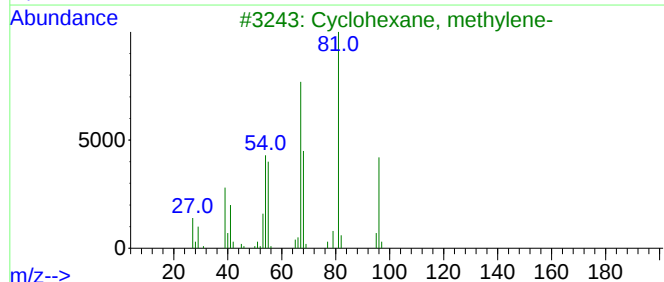
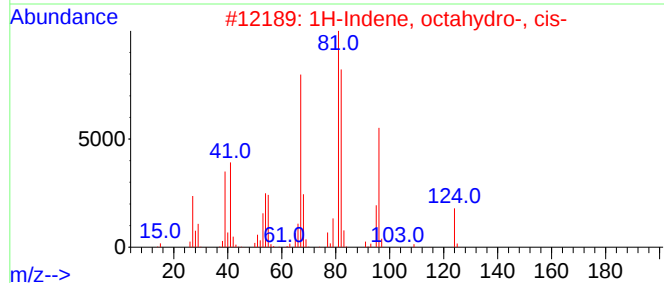
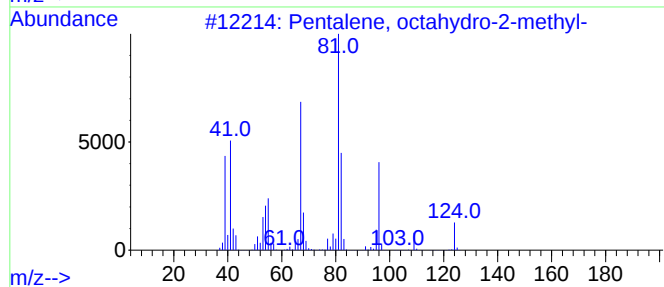
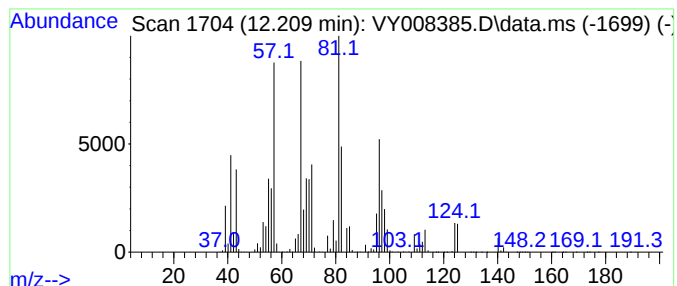
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Quant Title : SW846 8260

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

Peak Number 9 Pentalene, octahydro-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.209	744.38 ug/l	7370870	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	86
2			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	53
3			Cyclohexane, methylene-	96	C7H12	001192-37-6	49
4			1H-Indene, octahydro-, trans-	124	C9H16	003296-50-2	43
5			1H-Indene, octahydro-	124	C9H16	000496-10-6	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

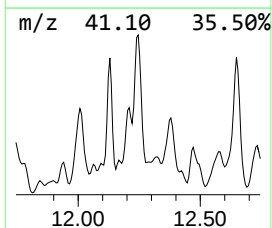
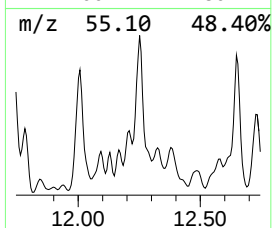
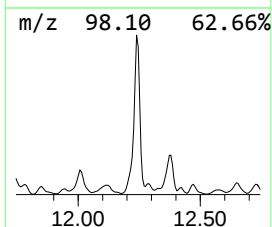
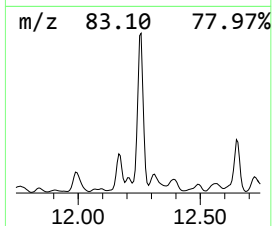
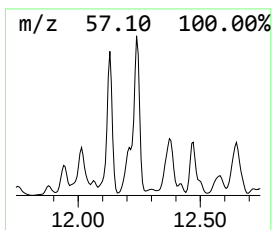
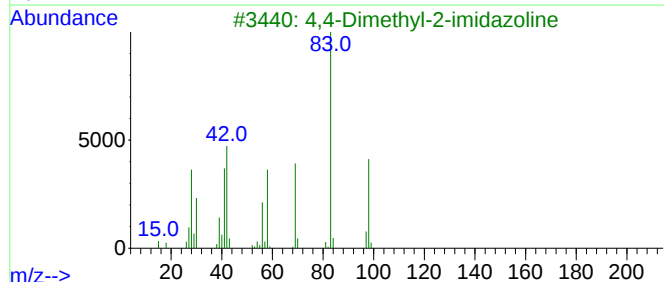
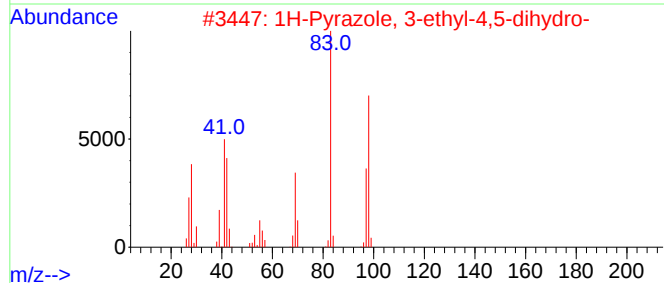
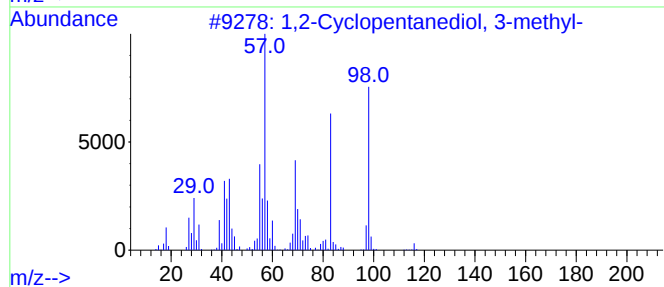
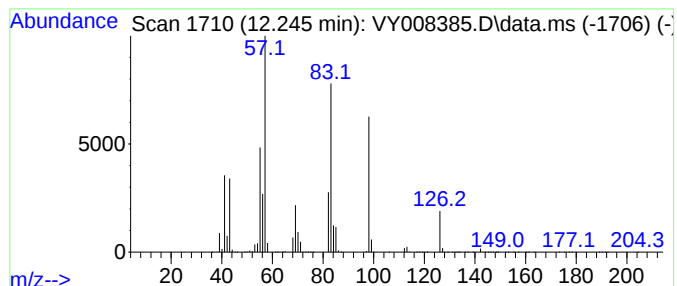
Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 10 1,2-Cyclopentanediol, 3-met... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.245	1327.40 ug/l	13144000	Chlorobenzene-d5	11.489	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Cyclopentanediol, 3-methyl-	116	C6H12O2	027583-37-5	53
2	1H-Pyrazole, 3-ethyl-4,5-dihydro-	98	C5H10N2	005920-29-6	49
3	4,4-Dimethyl-2-imidazoline	98	C5H10N2	002305-59-1	43
4	Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	43
5	Furan, 2-methoxy-	98	C5H6O2	025414-22-6	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041822\
Data File : VY008385.D
Acq On : 18 Apr 2022 19:50
Operator : KP/MD
Sample : N2452-07
Misc : 4.73g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WC-15-1-4-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.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
Pentane, 2,2,4-...	8.325	346.1	ug/l	3300360	2	8.691	476799	50.0
Cyclohexane, 1,...	10.349	282.9	ug/l	2800790	3	11.489	495104	50.0
Cyclohexane, 1,...	10.520	348.3	ug/l	3448410	3	11.489	495104	50.0
Cyclohexane, et...	11.038	262.5	ug/l	2598970	3	11.489	495104	50.0
Cyclohexane, 1,...	11.093	1096.9	ug/l	10862000	3	11.489	495104	50.0
1-Octene, 4-met...	11.203	265.3	ug/l	2626880	3	11.489	495104	50.0
Cyclohexane, 1,...	11.300	387.9	ug/l	3840950	3	11.489	495104	50.0
Octane, 3,6-dim...	12.130	689.4	ug/l	6826880	3	11.489	495104	50.0
Pentalene, octa...	12.209	744.4	ug/l	7370870	3	11.489	495104	50.0
1,2-Cyclopentan...	12.245	1327.4	ug/l	13144000	3	11.489	495104	50.0