

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011422\
 Data File : VY007169.D
 Acq On : 14 Jan 2022 15:20
 Operator : SY/MD
 Sample : N1143-02
 Misc : 4.99g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WEST-BOILER-ROOM-WEST-SIDEWALL

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\82Y011322S.M

Title : SW846 8260

Signal : TIC: VY007169.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.685	1115	1126	1140	rBV	220058	439993	1.15%	0.075%
2	10.179	1364	1371	1385	rVB	418454	809983	2.11%	0.138%
3	10.368	1393	1402	1410	rVB	239665	412147	1.08%	0.070%
4	10.801	1464	1473	1481	rBV	329189	573166	1.50%	0.098%
5	10.904	1483	1490	1497	rBV	437607	926966	2.42%	0.158%
6	11.032	1507	1511	1516	rVB	267952	410042	1.07%	0.070%
7	11.093	1516	1521	1526	rVB	356447	599326	1.56%	0.102%
8	11.209	1534	1540	1543	rBV	611858	1086456	2.84%	0.185%
9	11.282	1543	1552	1562	rVB2	4258285	8894901	23.22%	1.518%
10	11.380	1562	1568	1578	rBV	3995898	6487627	16.94%	1.107%
11	11.477	1578	1584	1591	rBV2	591281	1240424	3.24%	0.212%
12	11.569	1594	1599	1605	rBV3	444898	952842	2.49%	0.163%
13	11.715	1612	1623	1631	rBV3	8174540	18346155	47.89%	3.130%
14	11.782	1631	1634	1639	rVB	1353876	2052916	5.36%	0.350%
15	11.880	1639	1650	1654	rBV3	832424	2368984	6.18%	0.404%
16	11.934	1654	1659	1664	rBV2	3746391	6178170	16.13%	1.054%
17	12.008	1664	1671	1677	rBV3	7029719	15140265	39.53%	2.583%
18	12.062	1677	1680	1683	rVV	1935542	2437559	6.36%	0.416%
19	12.130	1686	1691	1695	rVB	8124808	12419346	32.42%	2.119%
20	12.203	1695	1703	1706	rBV3	4916212	10208782	26.65%	1.742%
21	12.245	1706	1710	1715	rVB3	5313092	9698278	25.32%	1.655%
22	12.325	1721	1723	1725	rBV	448964	517631	1.35%	0.088%
23	12.392	1725	1734	1735	rVV5	6862080	15893842	41.49%	2.712%
24	12.422	1735	1739	1741	rVV	10578455	19826153	51.76%	3.383%
25	12.447	1741	1743	1748	rVB	11193455	13960082	36.44%	2.382%
26	12.532	1752	1757	1761	rVB	10334184	15849863	41.38%	2.705%
27	12.581	1761	1765	1768	rVB2	1325787	1835937	4.79%	0.313%
28	12.648	1769	1776	1784	rVB3	4192599	12815437	33.46%	2.187%
29	12.739	1784	1791	1794	rBV	10251526	18410701	48.06%	3.141%
30	12.770	1794	1796	1798	rVV	2267682	2685022	7.01%	0.458%
31	12.812	1798	1803	1808	rVB2	21026382	38305027	100.00%	6.536%
32	12.867	1808	1812	1818	rVB2	3907039	6079134	15.87%	1.037%
33	12.959	1822	1827	1832	rBV2	6443005	15465052	40.37%	2.639%
34	13.007	1832	1835	1838	rVV3	4051423	6187966	16.15%	1.056%
35	13.044	1838	1841	1847	rVB2	8126449	11077113	28.92%	1.890%
36	13.123	1847	1854	1859	rVB	16619187	31169334	81.37%	5.318%

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37	13.172	1859	1862	1866	rVB	4420732	5698995	14.88%	0.972%
38	13.227	1866	1871	1874	rBV3	3409838	6054656	15.81%	1.033%
39	13.318	1879	1886	1890	rVV4	11491373	23201413	60.57%	3.959%
40	13.367	1890	1894	1897	rVV2	10480686	17267016	45.08%	2.946%
41	13.404	1897	1900	1902	rVV	6960589	9921433	25.90%	1.693%
42	13.434	1902	1905	1907	rVV	8646474	12907383	33.70%	2.202%
43	13.459	1907	1909	1913	rVV	9432037	13294345	34.71%	2.268%
44	13.507	1913	1917	1923	rVV2	7349299	11763435	30.71%	2.007%
45	13.593	1923	1931	1933	rVV3	3787875	7452526	19.46%	1.272%
46	13.623	1933	1936	1940	rVV	9200329	14327145	37.40%	2.445%
47	13.666	1940	1943	1952	rVV3	10812879	21165482	55.26%	3.612%
48	13.757	1952	1958	1962	rVV3	10826470	18159421	47.41%	3.099%
49	13.824	1962	1969	1973	rVV	5194485	11220200	29.29%	1.915%
50	13.885	1975	1979	1981	rVV	4977049	8293952	21.65%	1.415%
51	13.916	1981	1984	1989	rVV2	8031112	13747149	35.89%	2.346%
52	13.971	1989	1993	1998	rVV	7348178	11869055	30.99%	2.025%
53	14.013	1998	2000	2005	rVV4	1176272	1965583	5.13%	0.335%
54	14.080	2005	2011	2015	rVV2	5791365	9663542	25.23%	1.649%
55	14.135	2015	2020	2027	rVV4	2062937	5165924	13.49%	0.881%
56	14.215	2027	2033	2037	rVV4	2721902	5757259	15.03%	0.982%
57	14.294	2038	2046	2050	rVV3	4004758	12389232	32.34%	2.114%
58	14.336	2050	2053	2057	rVV	4388612	5972365	15.59%	1.019%
59	14.379	2057	2060	2064	rVV2	3155260	4501008	11.75%	0.768%
60	14.422	2064	2067	2070	rVV	1609254	2257995	5.89%	0.385%
61	14.452	2070	2072	2075	rVV	727127	902018	2.35%	0.154%
62	14.519	2077	2083	2087	rVB2	1055349	2382093	6.22%	0.406%
63	14.568	2087	2091	2094	rBV2	596177	1087094	2.84%	0.185%
64	14.611	2094	2098	2103	rVB	1910609	2791646	7.29%	0.476%
65	14.672	2103	2108	2115	rBV3	3025865	5912104	15.43%	1.009%
66	14.733	2115	2118	2124	rVV2	692730	1163152	3.04%	0.198%
67	14.788	2124	2127	2131	rVV2	334484	471360	1.23%	0.080%
68	14.830	2131	2134	2138	rVB	637657	810593	2.12%	0.138%
69	14.946	2148	2153	2161	rBV2	767232	1761690	4.60%	0.301%
70	15.050	2165	2170	2176	rVB3	487868	967710	2.53%	0.165%
71	15.141	2180	2185	2190	rVB3	206880	384112	1.00%	0.066%
72	15.233	2190	2200	2206	rVB	875811	1642279	4.29%	0.280%

Sum of corrected areas: 586054987

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MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

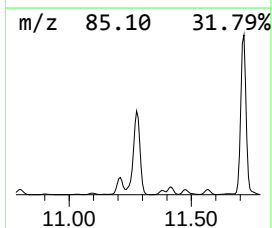
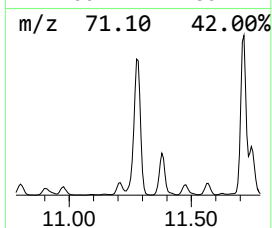
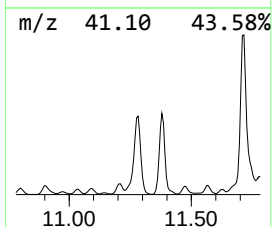
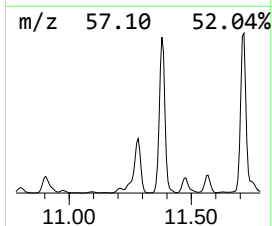
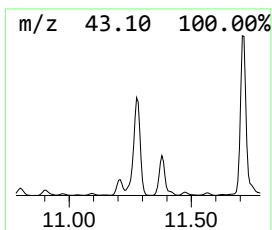
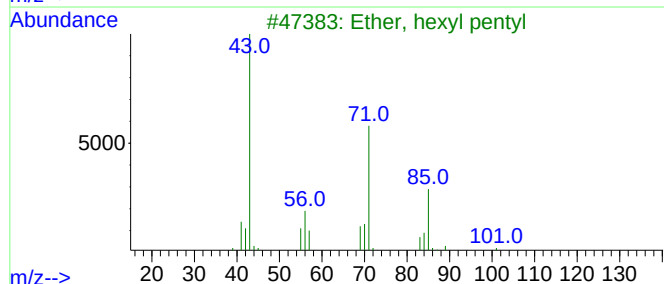
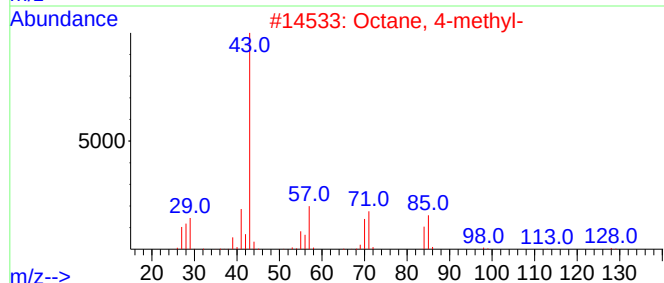
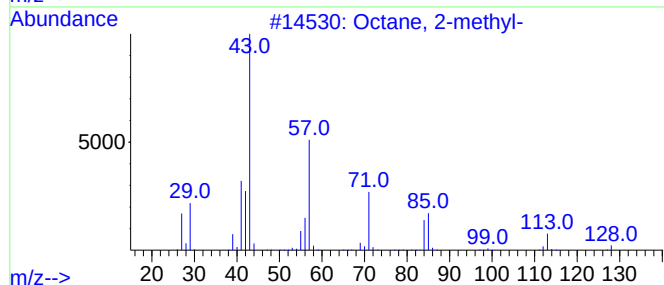
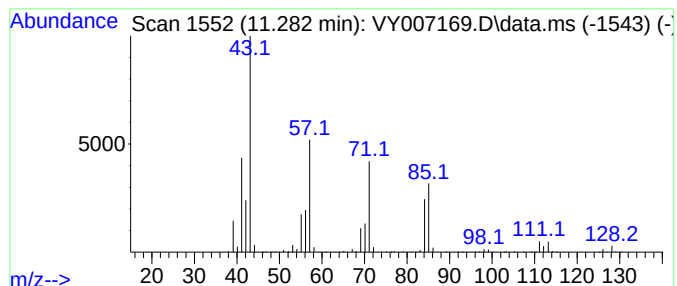
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.282	358.54 ug/l	8894900	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2-methyl-	128	C9H20	003221-61-2	83
2			Octane, 4-methyl-	128	C9H20	002216-34-4	59
3			Ether, hexyl pentyl	172	C11H24O	032357-83-8	53
4			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	52
5			Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	50



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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

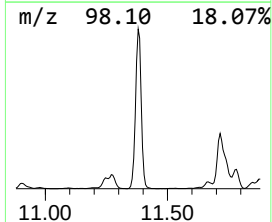
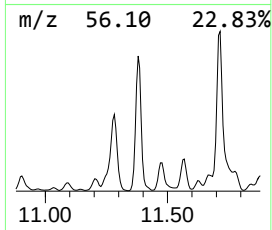
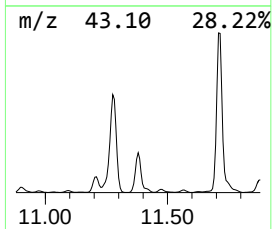
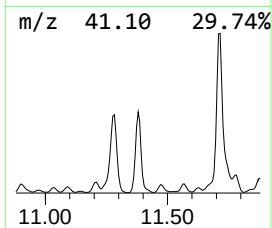
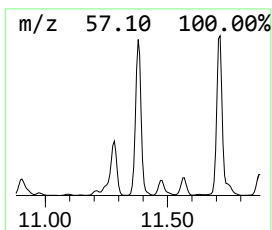
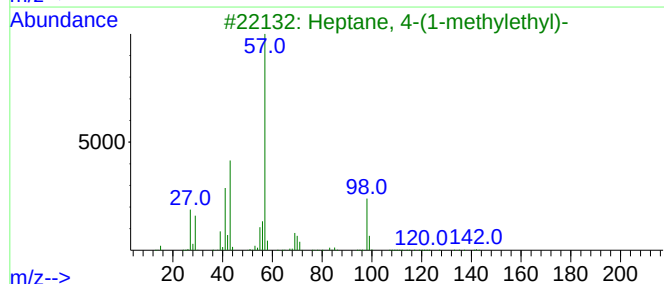
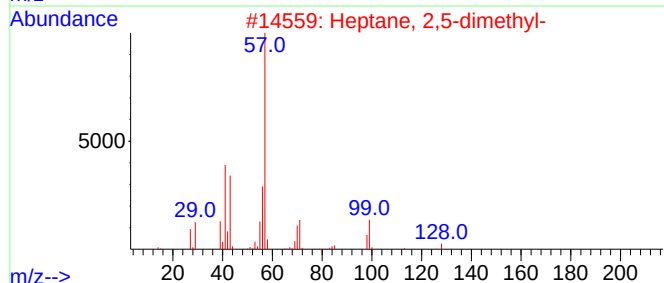
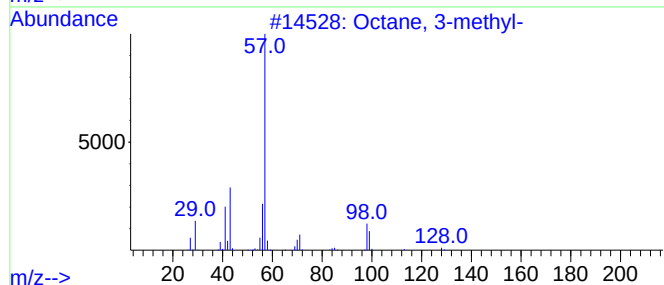
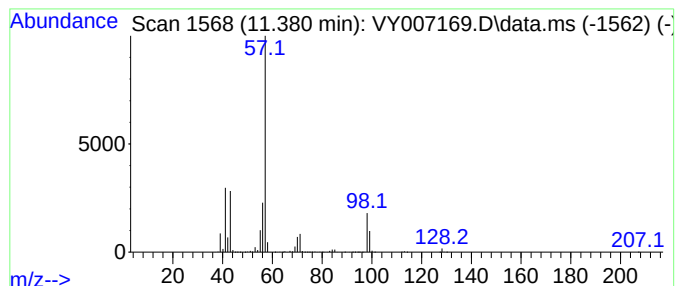
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 2 Octane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.380	261.51 ug/l	6487630	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3-methyl-	128	C9H20	002216-33-3	91
2			Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	81
3			Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	64
4			Heptane, 3-ethyl-	128	C9H20	015869-80-4	50
5			Butane, 1-chloro-2-methyl-, (S)-	106	C5H11Cl	040560-29-0	47



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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

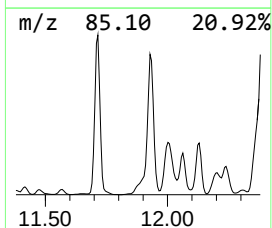
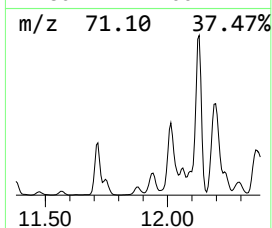
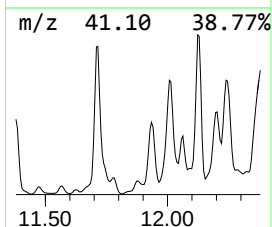
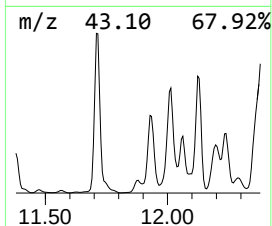
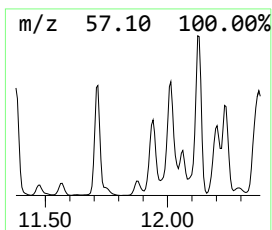
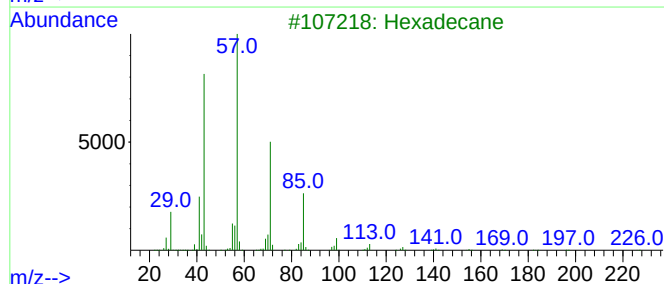
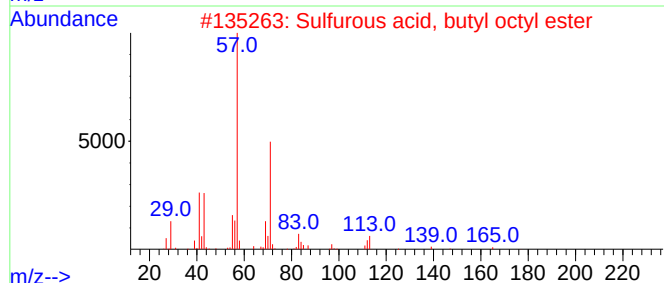
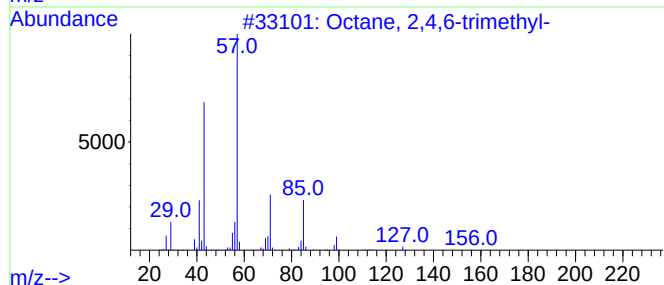
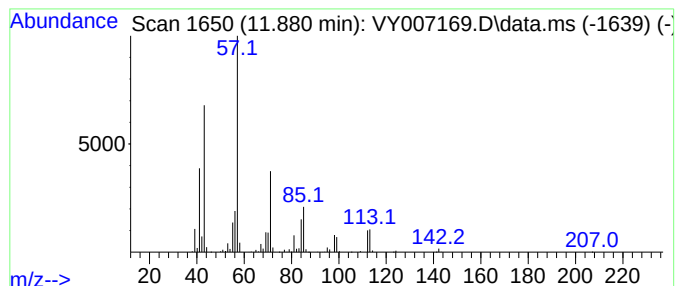
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 3 Octane, 2,4,6-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	95.49 ug/l	2368980	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	59
2			Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	53
3			Hexadecane	226	C16H34	000544-76-3	53
4			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	52
5			Tridecane, 6-methyl-	198	C14H30	013287-21-3	50



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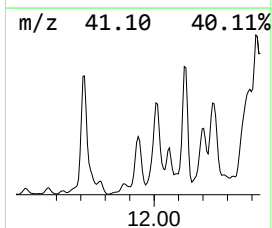
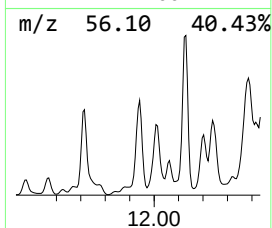
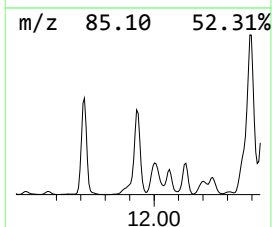
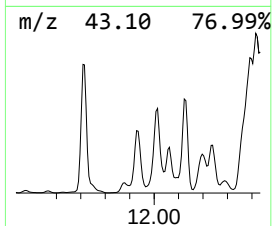
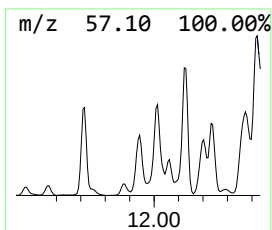
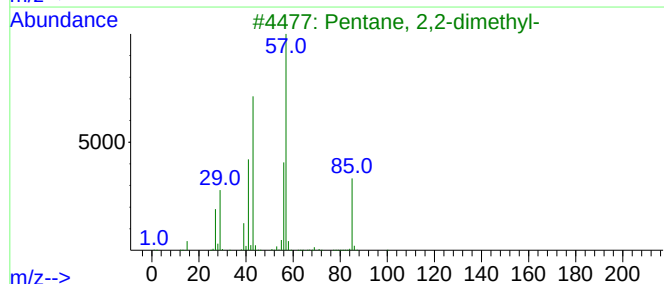
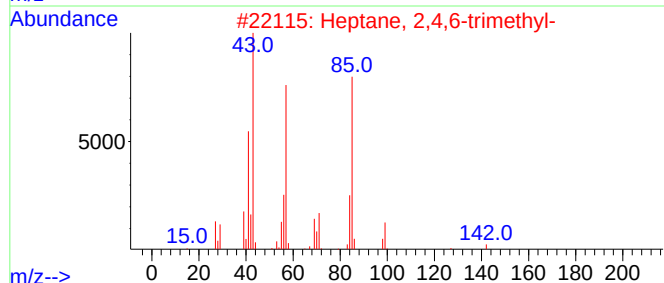
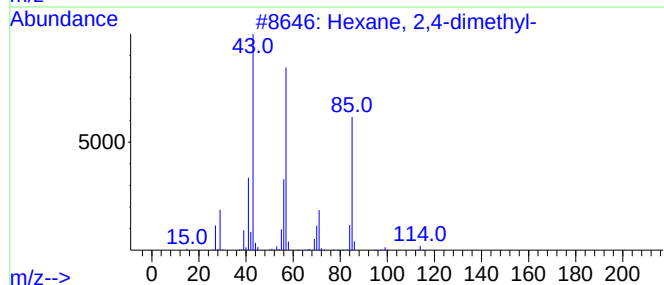
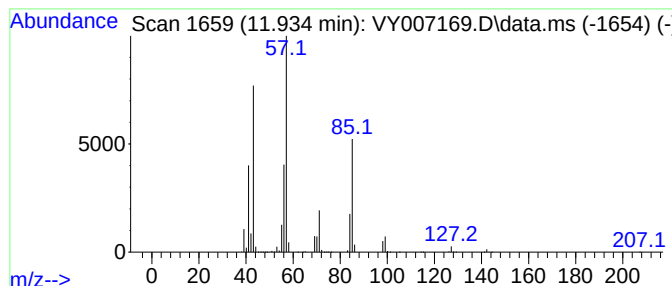
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 4 Hexane, 2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.935	249.04 ug/l	6178170	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	78
2			Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	60
3			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	59
4			Heptane, 4,4-dimethyl-	128	C9H20	001068-19-5	47
5			Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	47



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Data File : VY007169.D
Acq On : 14 Jan 2022 15:20
Operator : SY/MD
Sample : N1143-02
Misc : 4.99g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

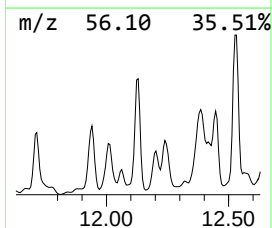
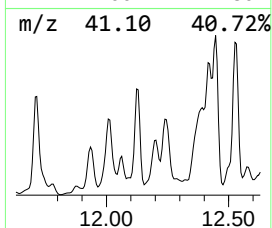
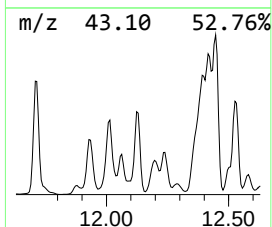
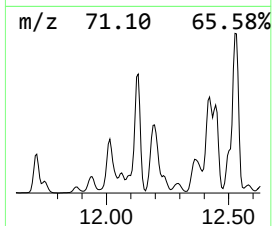
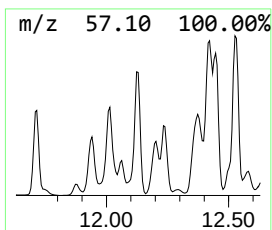
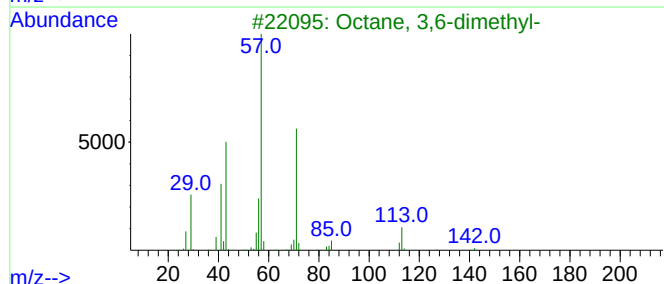
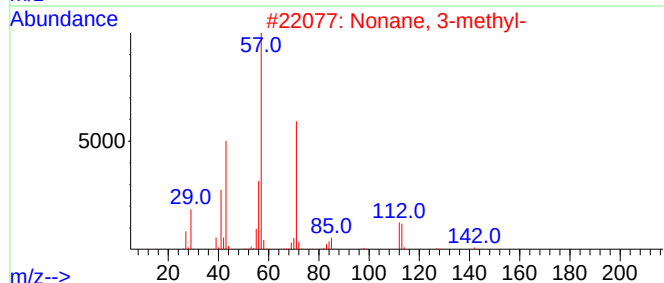
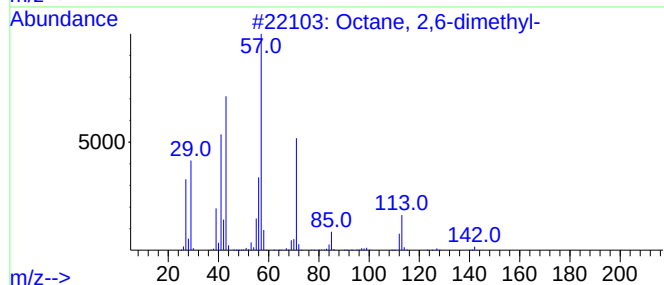
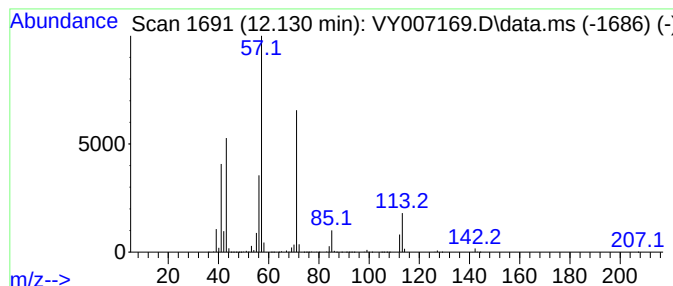
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Quant Title : SW846 8260

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

Peak Number 5 Octane, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.130	500.61 ug/l	12419300	Chlorobenzene-d5	11.489

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	90
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	86
4			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	64
5			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011422\
Data File : VY007169.D
Acq On : 14 Jan 2022 15:20
Operator : SY/MD
Sample : N1143-02
Misc : 4.99g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

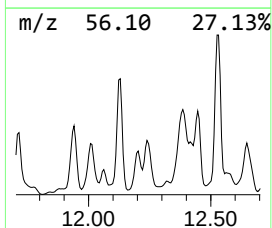
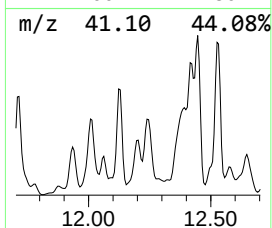
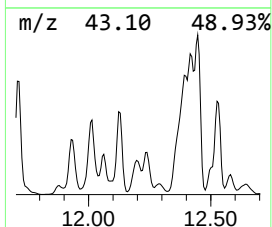
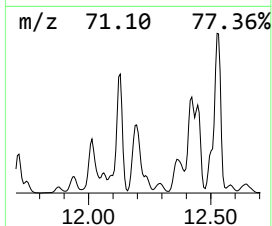
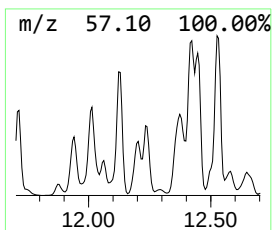
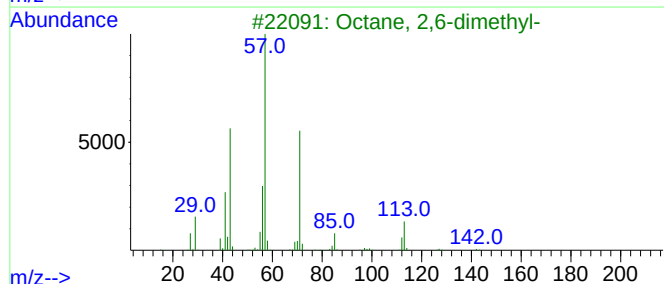
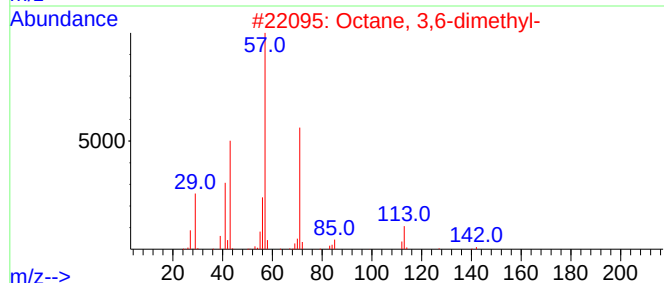
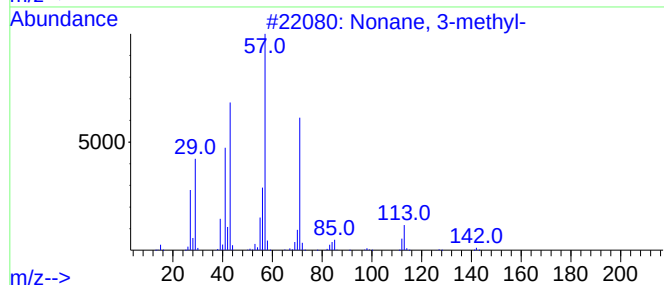
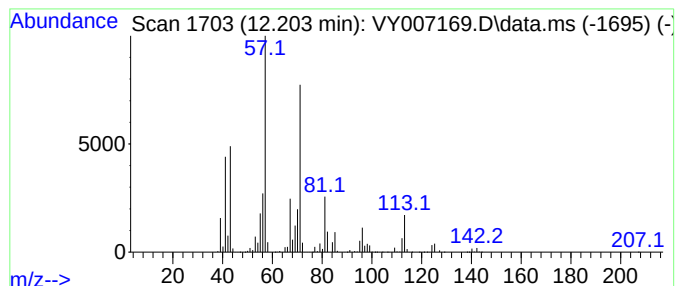
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.203	411.50 ug/l	10208800	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	76
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	62
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	59
4			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	52
5			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011422\
Data File : VY007169.D
Acq On : 14 Jan 2022 15:20
Operator : SY/MD
Sample : N1143-02
Misc : 4.99g/5.0mL/MSVOA_Y/soil
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

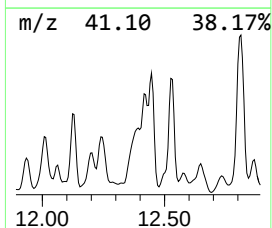
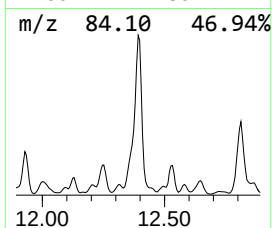
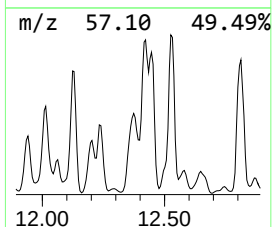
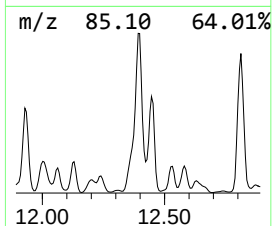
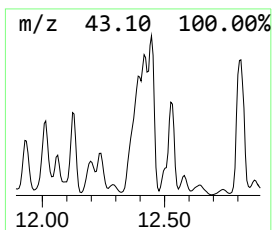
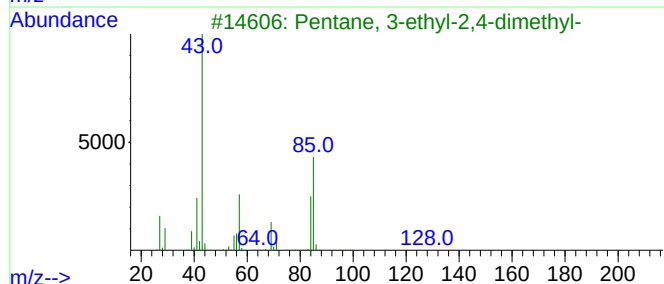
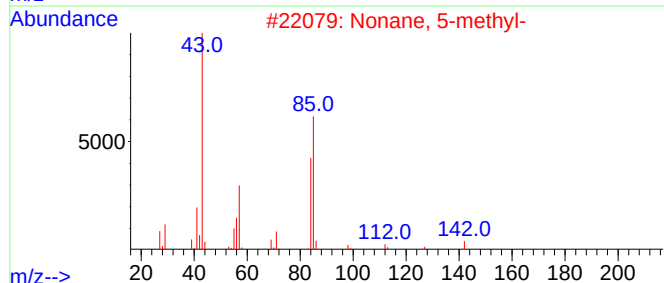
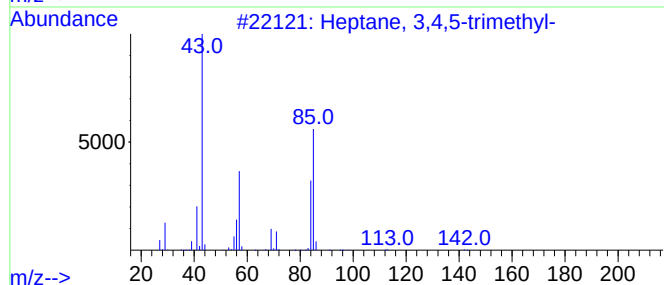
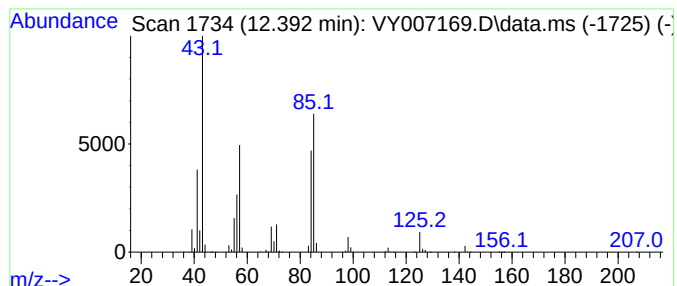
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 8 Heptane, 3,4,5-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.392	640.66 ug/l	15893800	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	72
2			Nonane, 5-methyl-	142	C10H22	015869-85-9	64
3			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	64
4			Hexane, 3-ethyl-	114	C8H18	000619-99-8	53
5			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011422\
Data File : VY007169.D
Acq On : 14 Jan 2022 15:20
Operator : SY/MD
Sample : N1143-02
Misc : 4.99g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

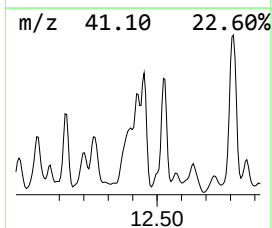
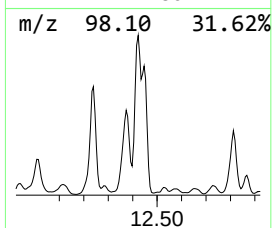
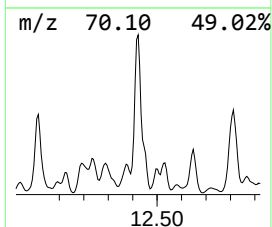
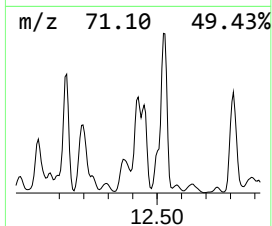
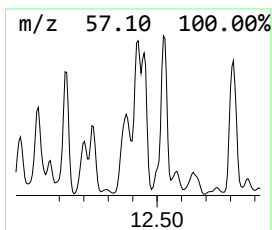
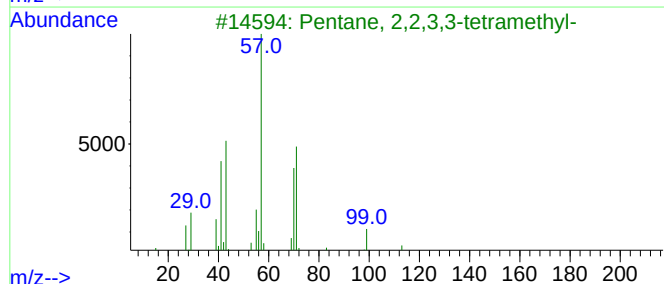
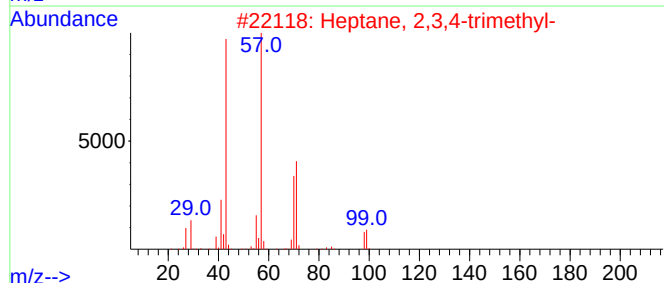
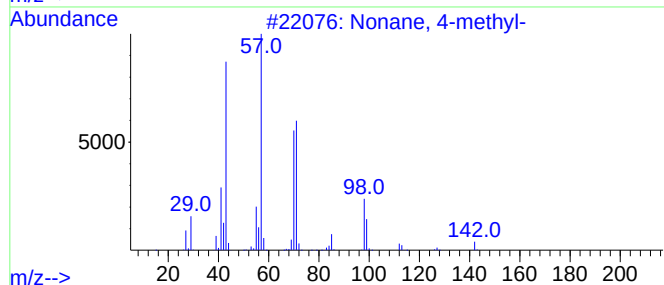
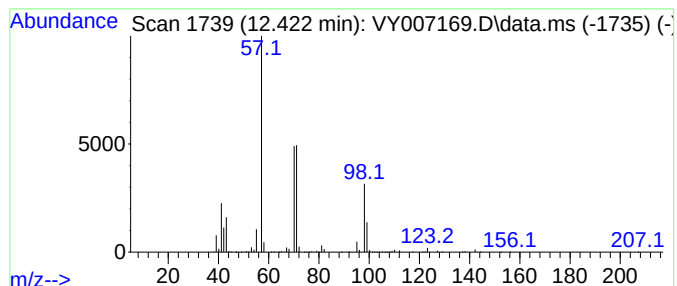
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 9 Nonane, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.422	799.17 ug/l	19826200	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	64
2			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	59
3			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	45
4			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	43
5			Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011422\
Data File : VY007169.D
Acq On : 14 Jan 2022 15:20
Operator : SY/MD
Sample : N1143-02
Misc : 4.99g/5.0mL/MSVOA_Y/soil
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

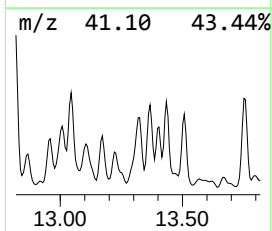
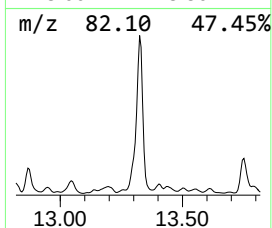
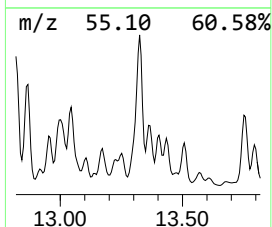
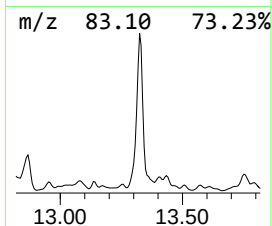
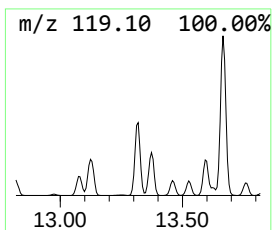
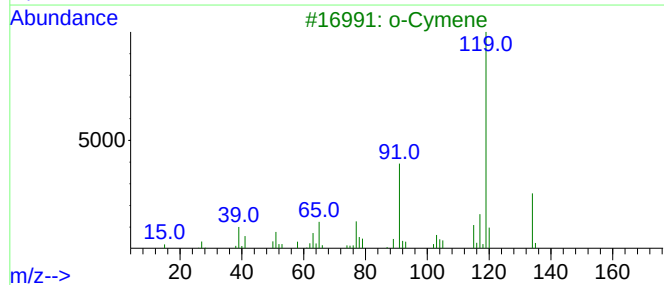
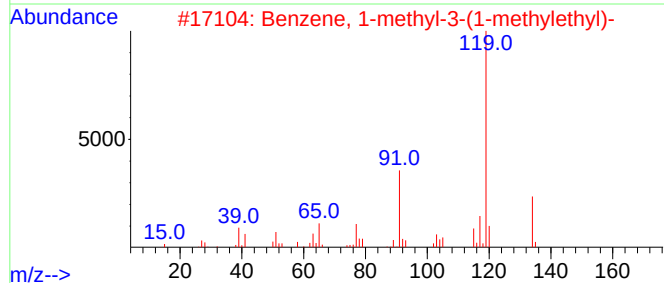
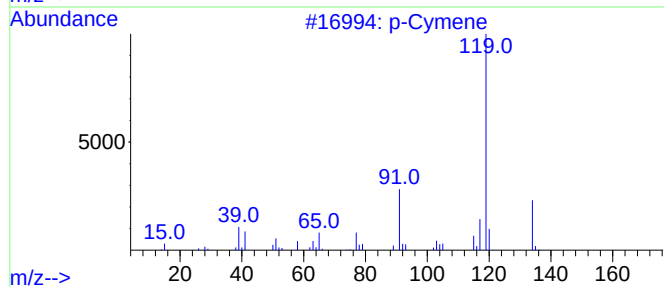
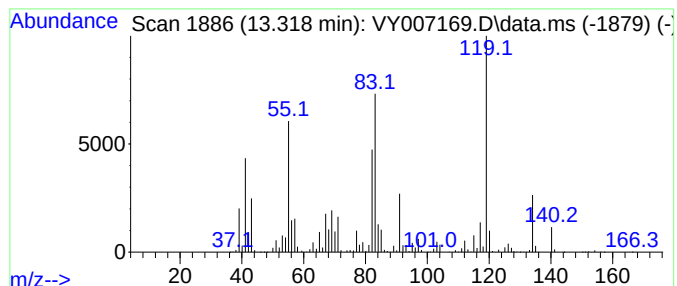
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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

Peak Number 10 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.318	89.88 ug/l	23201400	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	86
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	86
3			o-Cymene	134	C10H14	000527-84-4	81
4			1,3,5-Cycloheptatriene, 3,7,7-trimethyl-	134	C10H14	003479-89-8	53
5			Cyclohexane, pentyl-	154	C11H22	004292-92-6	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011422\
Data File : VY007169.D
Acq On : 14 Jan 2022 15:20
Operator : SY/MD
Sample : N1143-02
Misc : 4.99g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WEST-BOILER-ROOM-WEST-SIDEWALL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.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
Octane, 2-methyl-	11.282	358.5	ug/l	8894900	3	11.489	1240420	50.0
Octane, 3-methyl-	11.380	261.5	ug/l	6487630	3	11.489	1240420	50.0
Octane, 2,4,6-t...	11.880	95.5	ug/l	2368980	3	11.489	1240420	50.0
Hexane, 2,4-dim...	11.935	249.0	ug/l	6178170	3	11.489	1240420	50.0
Octane, 2,6-dim...	12.130	500.6	ug/l	12419300	3	11.489	1240420	50.0
Nonane, 3-methyl-	12.203	411.5	ug/l	10208800	3	11.489	1240420	50.0
Heptane, 3,4,5-...	12.392	640.7	ug/l	15893800	3	11.489	1240420	50.0
Nonane, 4-methyl-	12.422	799.2	ug/l	19826200	3	11.489	1240420	50.0
p-Cymene	13.318	89.9	ug/l	23201400	4	13.428	12907400	50.0