

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
 Data File : VY004262.D
 Acq On : 02 Apr 2021 15:02
 Operator : SY/MD
 Sample : M1911-02
 Misc : 5.05g/5mL/MSVOA_Y/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TP-6S1-(0-2)

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\82Y032921S.M
 Title : SW846 8260

Signal : TIC: VY004262.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.217	46	65	66	rBV5	10721	46348	2.93%	0.264%
2	2.564	113	122	123	rBV5	15208	38842	2.46%	0.222%
3	3.967	342	352	361	rBV	12054	33982	2.15%	0.194%
4	4.735	469	478	492	rVB3	10709	34118	2.16%	0.195%
5	7.728	959	969	975	rBV	146694	349012	22.10%	1.990%
6	7.801	975	981	995	rVB	252218	561009	35.52%	3.199%
7	8.155	1029	1039	1049	rVB	150934	328531	20.80%	1.874%
8	8.697	1117	1128	1138	rBV	393988	784548	49.67%	4.474%
9	10.185	1364	1372	1380	rBV	666086	1122927	71.09%	6.404%
10	10.374	1396	1403	1410	rVB2	12484	21004	1.33%	0.120%
11	10.581	1431	1437	1451	rBV	12820	26211	1.66%	0.149%
12	11.282	1544	1552	1563	rVB4	28630	67394	4.27%	0.384%
13	11.386	1563	1569	1574	rBV	25393	43795	2.77%	0.250%
14	11.496	1580	1587	1597	rVV	588392	952228	60.29%	5.430%
15	11.599	1598	1604	1608	rVV	16745	30944	1.96%	0.176%
16	11.715	1612	1623	1631	rVV3	204389	416900	26.39%	2.377%
17	11.782	1631	1634	1640	rVB2	18988	31894	2.02%	0.182%
18	12.032	1664	1675	1683	rBV3	58087	180232	11.41%	1.028%
19	12.130	1686	1691	1695	rVV	53293	77628	4.91%	0.443%
20	12.203	1699	1703	1706	rVV4	28751	51164	3.24%	0.292%
21	12.246	1706	1710	1719	rVV6	59181	132677	8.40%	0.757%
22	12.331	1720	1724	1727	rVV3	18092	33318	2.11%	0.190%
23	12.392	1728	1734	1735	rVV5	38994	82172	5.20%	0.469%
24	12.416	1735	1738	1740	rVV	85089	120767	7.65%	0.689%
25	12.441	1740	1742	1745	rVV	78695	107509	6.81%	0.613%
26	12.489	1745	1750	1754	rVV	459751	762835	48.30%	4.350%
27	12.526	1754	1756	1761	rVB	76896	89574	5.67%	0.511%
28	12.648	1772	1776	1777	rBV	29245	33163	2.10%	0.189%
29	12.739	1784	1791	1794	rBV	124686	216016	13.68%	1.232%
30	12.770	1794	1796	1797	rVV	62971	63452	4.02%	0.362%
31	12.806	1797	1802	1808	rVV	807812	1242551	78.67%	7.086%
32	12.861	1808	1811	1816	rVB	58719	90514	5.73%	0.516%
33	12.977	1822	1830	1833	rBV2	71258	185266	11.73%	1.057%
34	13.008	1833	1835	1837	rVV2	30944	36321	2.30%	0.207%
35	13.044	1837	1841	1847	rVB	200109	287296	18.19%	1.638%
36	13.123	1847	1854	1859	rBV	263794	420898	26.65%	2.400%

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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

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
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37	13.172	1859	1862	1867	rVB2	40387	52322	3.31%	0.298%
38	13.227	1867	1871	1873	rBV	50651	72739	4.61%	0.415%
39	13.325	1879	1887	1890	rBV3	163002	320704	20.30%	1.829%
40	13.367	1890	1894	1897	rVV2	156431	233182	14.76%	1.330%
41	13.428	1897	1904	1913	rVB3	701097	1350615	85.51%	7.702%
42	13.507	1913	1917	1923	rVB	134337	183644	11.63%	1.047%
43	13.575	1924	1928	1929	rBV3	20553	25332	1.60%	0.144%
44	13.629	1933	1937	1940	rVB	99112	129425	8.19%	0.738%
45	13.672	1940	1944	1952	rVB3	90053	173284	10.97%	0.988%
46	13.757	1952	1958	1962	rBV	1135729	1579486	100.00%	9.007%
47	13.800	1962	1965	1967	rVV2	26477	26982	1.71%	0.154%
48	13.831	1967	1970	1973	rVV	28936	34835	2.21%	0.199%
49	13.916	1981	1984	1990	rVB4	116893	175840	11.13%	1.003%
50	13.977	1990	1994	1998	rVB3	120021	173784	11.00%	0.991%
51	14.020	1998	2001	2006	rVB2	85456	123604	7.83%	0.705%
52	14.081	2006	2011	2016	rBV3	129891	204687	12.96%	1.167%
53	14.148	2016	2022	2028	rBV6	57496	125903	7.97%	0.718%
54	14.276	2028	2043	2047	rBV6	150659	624309	39.53%	3.560%
55	14.318	2047	2050	2057	rVB2	119593	252744	16.00%	1.441%
56	14.385	2057	2061	2064	rBV	109014	135460	8.58%	0.772%
57	14.422	2064	2067	2072	rVB3	61907	85078	5.39%	0.485%
58	14.568	2088	2091	2093	rBV2	28741	37702	2.39%	0.215%
59	14.617	2094	2099	2104	rVB	485620	668190	42.30%	3.810%
60	14.684	2104	2110	2115	rBV4	149080	342390	21.68%	1.953%
61	14.733	2115	2118	2125	rVB2	142749	234419	14.84%	1.337%
62	14.836	2132	2135	2143	rVB	90424	139157	8.81%	0.794%
63	14.946	2149	2153	2157	rBV5	37798	53921	3.41%	0.307%
64	15.056	2165	2171	2176	rBV4	48824	105397	6.67%	0.601%
65	15.214	2192	2197	2205	rVB2	75273	132970	8.42%	0.758%
66	15.300	2207	2211	2214	rVB3	27006	36404	2.30%	0.208%
67	15.349	2216	2219	2222	rBV3	17438	27893	1.77%	0.159%
68	15.440	2230	2234	2242	rVB	135952	210449	13.32%	1.200%
69	15.531	2244	2249	2256	rVV7	23054	64461	4.08%	0.368%
70	15.611	2259	2262	2267	rVB2	22249	33719	2.13%	0.192%
71	15.733	2278	2282	2289	rVB2	43412	79764	5.05%	0.455%
72	16.038	2326	2332	2338	rBV5	24928	55967	3.54%	0.319%
73	16.159	2349	2352	2357	rVB4	10695	15989	1.01%	0.091%
74	16.233	2359	2364	2370	rVV5	15996	31513	2.00%	0.180%
75	16.300	2370	2375	2379	rVV6	11012	22383	1.42%	0.128%
76	16.367	2382	2386	2392	rVB2	19506	32099	2.03%	0.183%

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TP-6S1-(0-2)

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

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Peak separation: 5

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77	16.489	2401	2406	2412	rBV9	8863	23772	1.51%	0.136%
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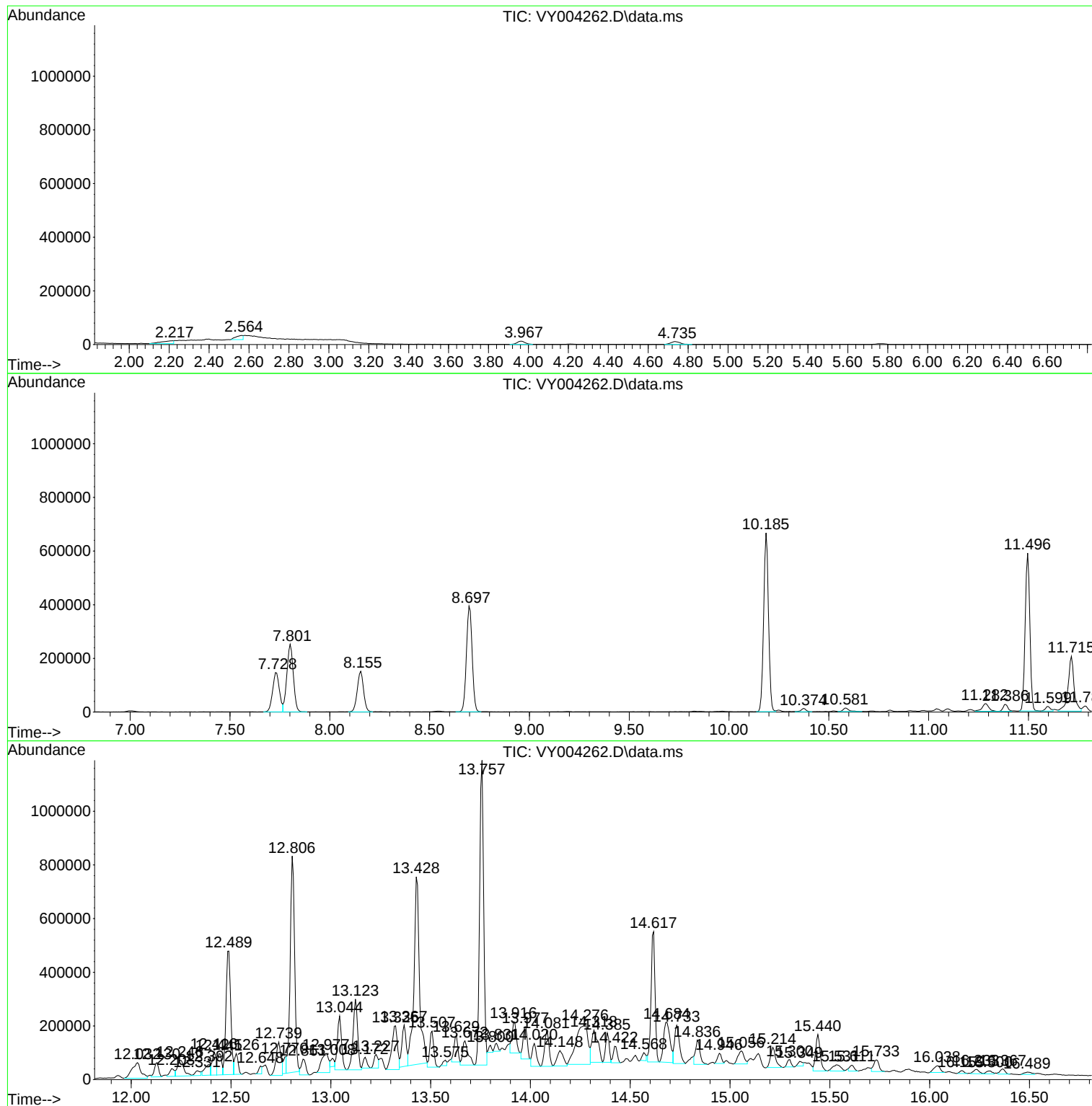
Sum of corrected areas: 17535558

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Sample : M1911-02
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Instrument :
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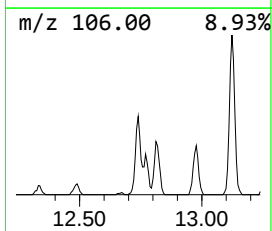
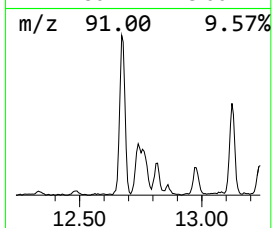
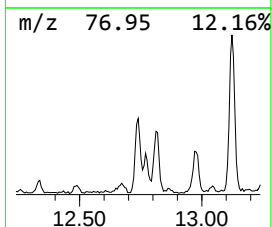
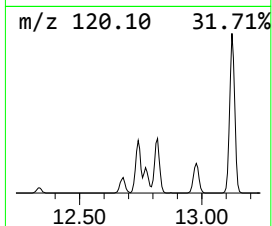
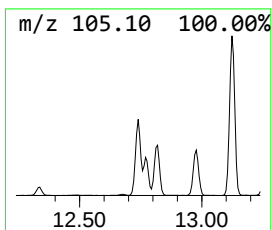
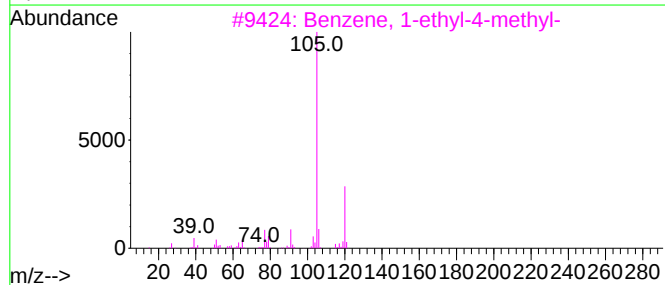
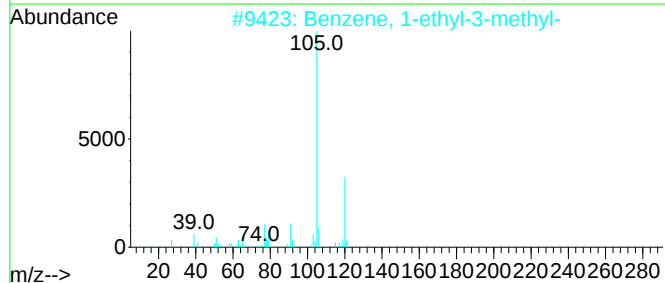
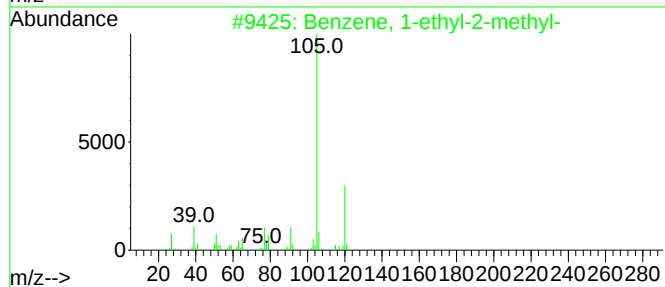
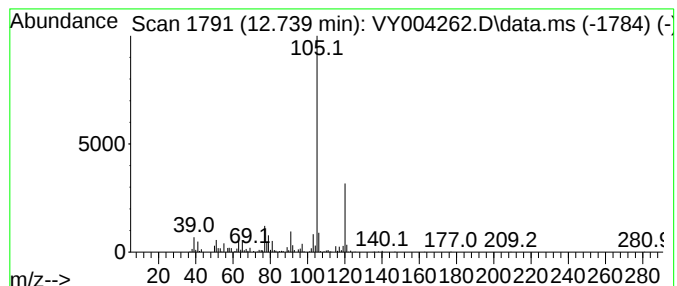
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.739	8.00 ug/l	216016	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5			Mesitylene	120	C9H12	000108-67-8	90



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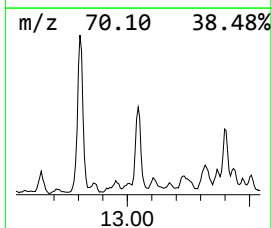
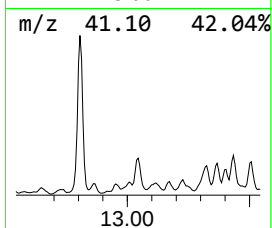
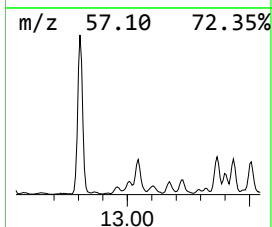
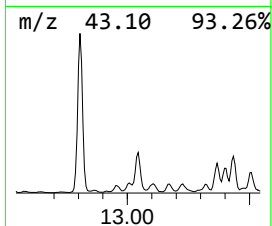
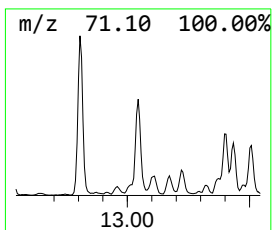
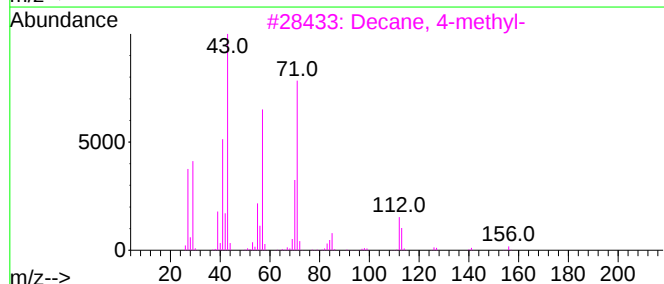
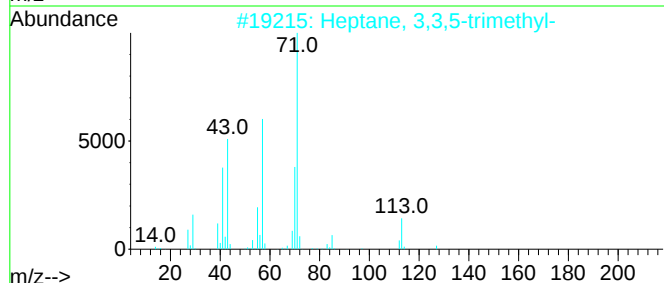
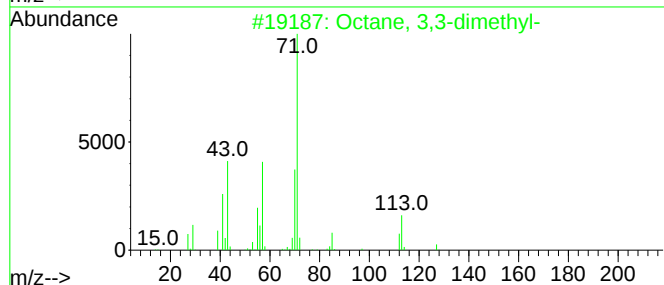
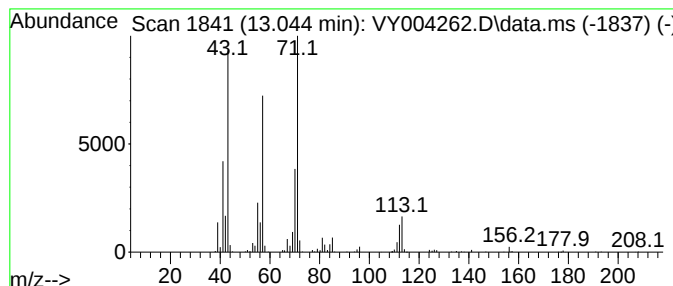
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.044	10.64 ug/l	287296	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	78
2			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	78
3			Decane, 4-methyl-	156	C11H24	002847-72-5	60
4			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	59
5			Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	59



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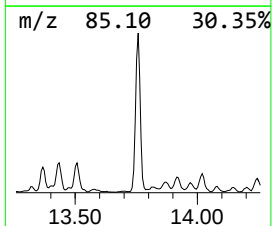
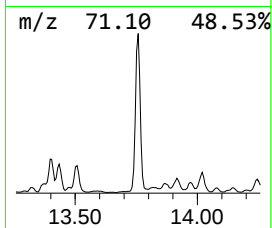
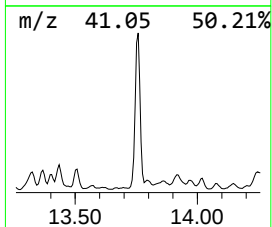
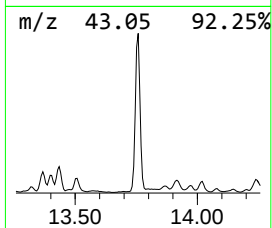
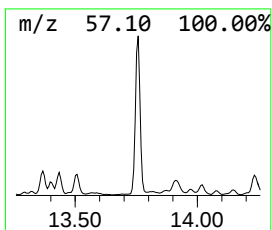
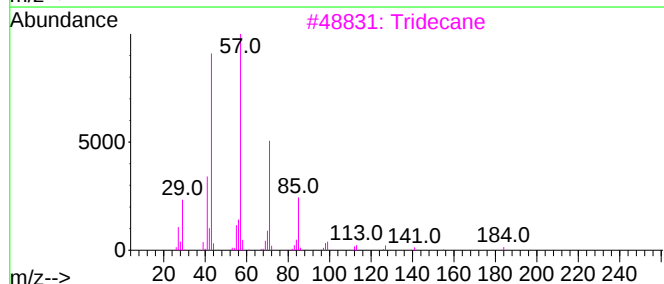
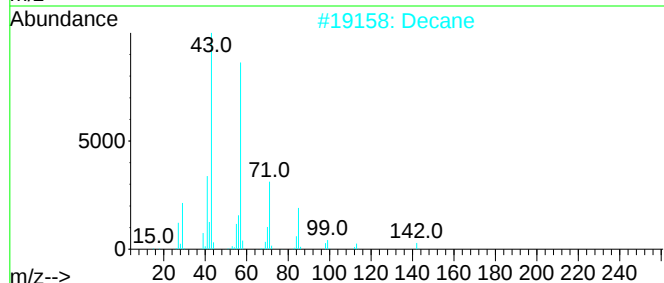
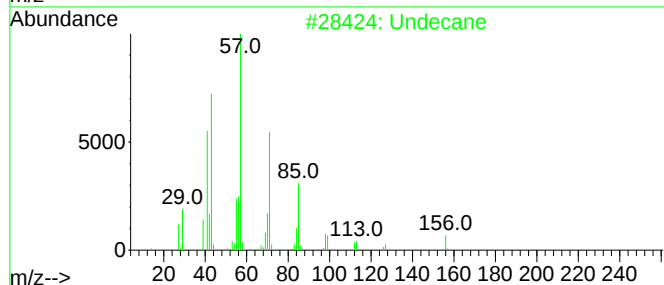
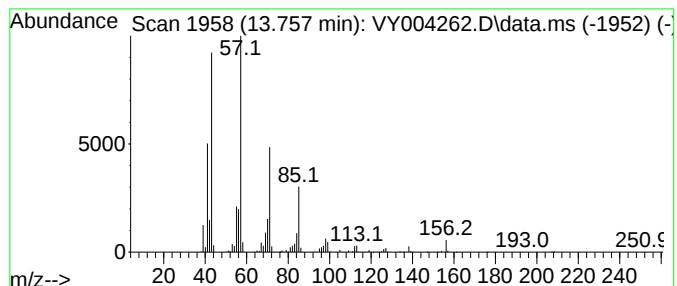
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Undecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.757	58.47 ug/l	1579490	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	96
2	Decane			142	C10H22	000124-18-5	90
3	Tridecane			184	C13H28	000629-50-5	72
4	Hexadecane			226	C16H34	000544-76-3	72
5	1-Iodo-2-methylnonane			268	C10H21I	1000101-47-9	64



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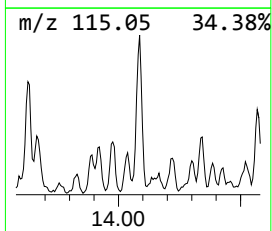
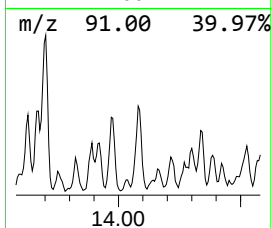
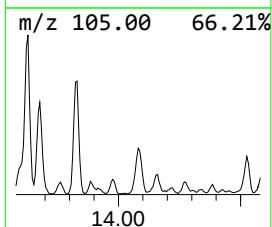
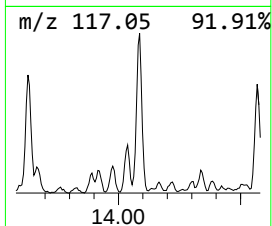
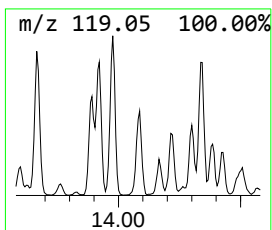
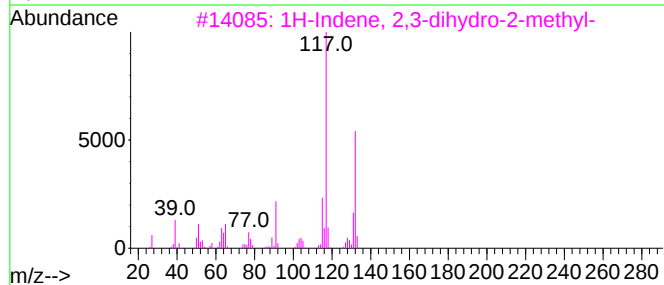
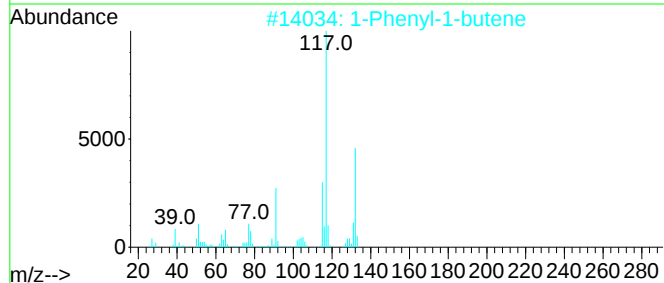
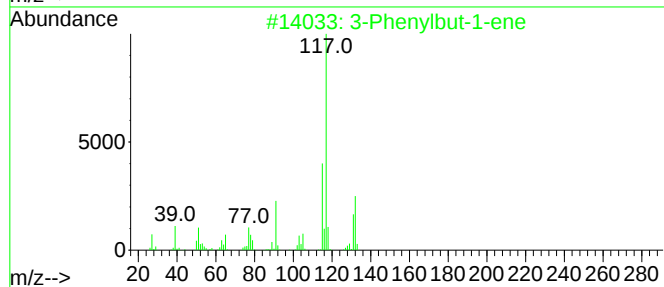
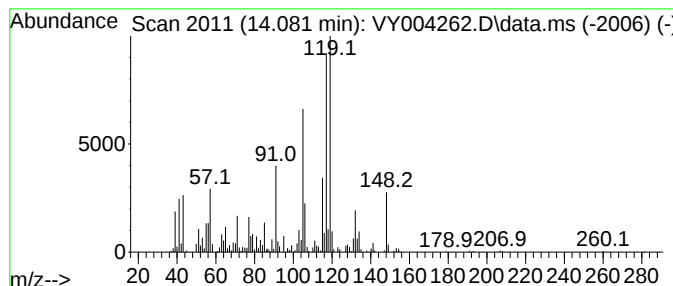
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 3-Phenylbut-1-ene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	7.58 ug/l	204687	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	46
3			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	46
4			o-Cymene	134	C10H14	000527-84-4	46
5			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY04262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

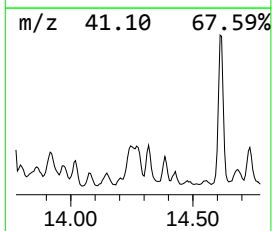
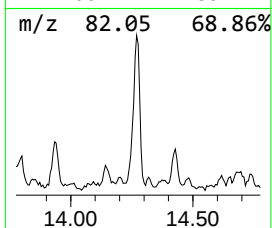
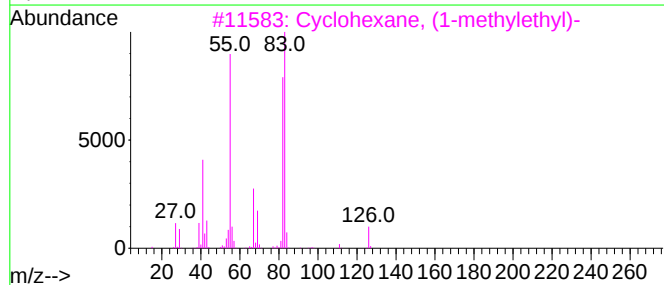
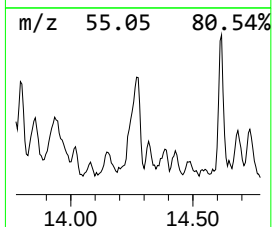
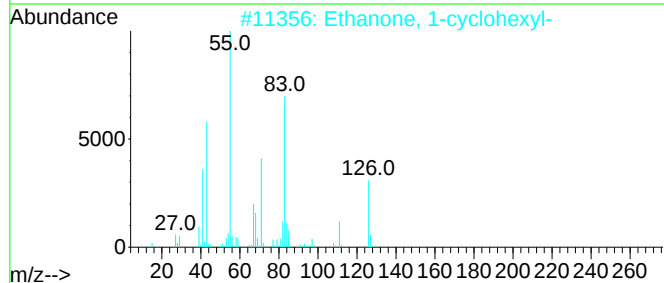
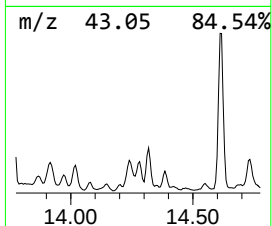
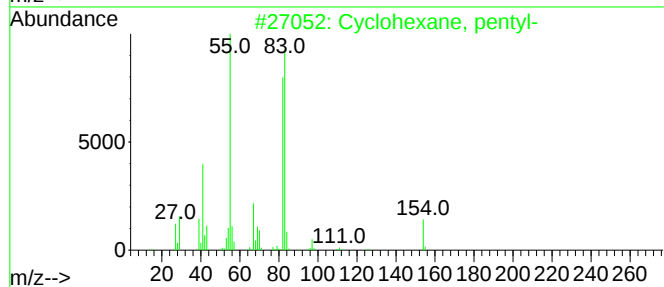
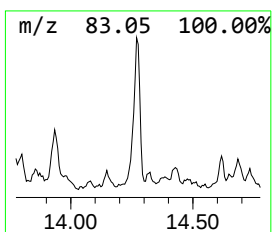
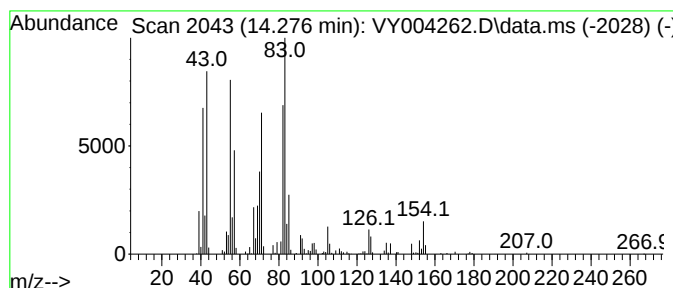
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclohexane, pentyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.276	23.11 ug/l	624309	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	60
2			Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	52
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	49
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	49
5			Cyclohexane, butyl-	140	C10H20	001678-93-9	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY04262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

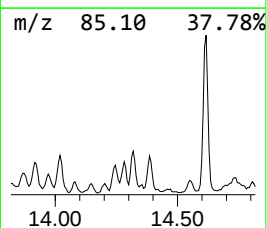
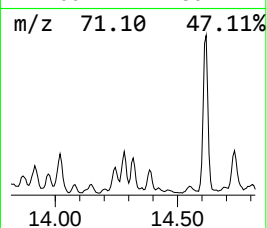
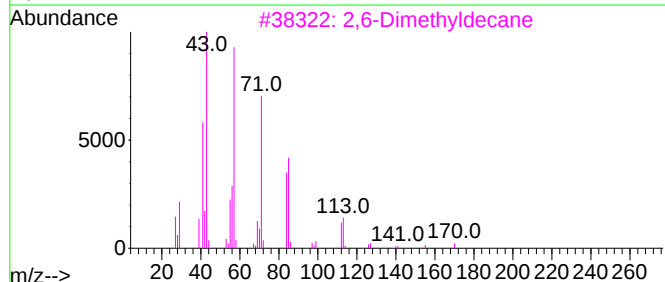
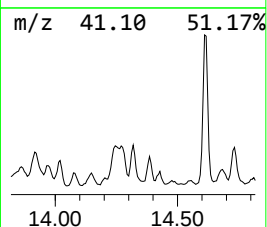
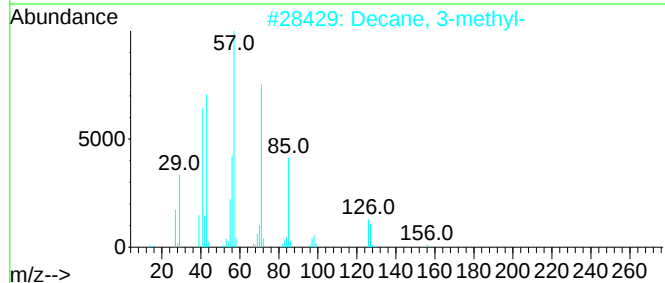
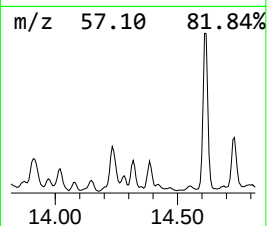
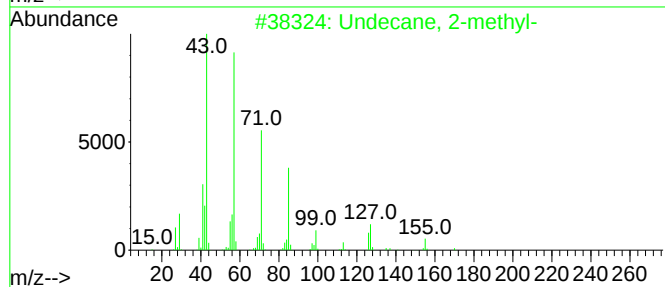
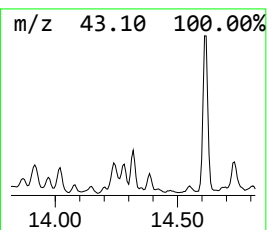
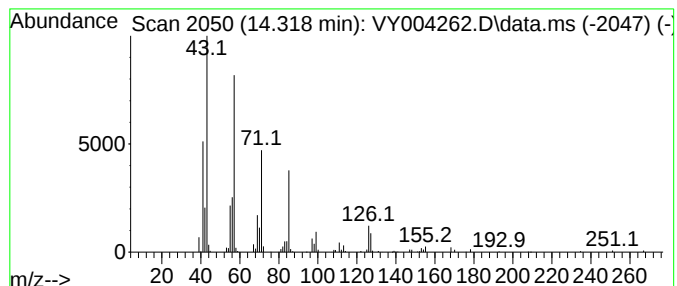
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Undecane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.318	9.36 ug/l	252744	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2-methyl-	170	C12H26	007045-71-8	90
2			Decane, 3-methyl-	156	C11H24	013151-34-3	64
3			2,6-Dimethyldecane	170	C12H26	013150-81-7	59
4			Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	59
5			Hexadecane	226	C16H34	000544-76-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY040262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

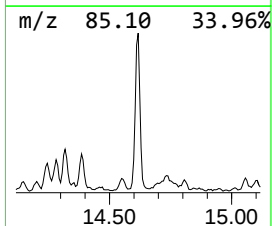
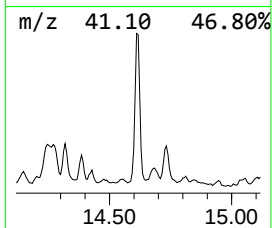
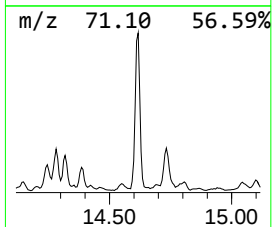
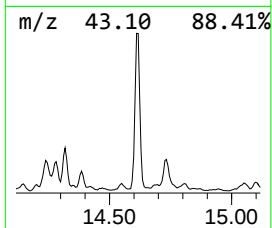
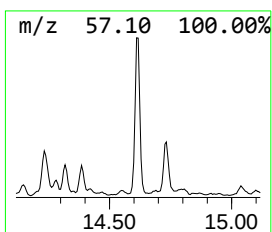
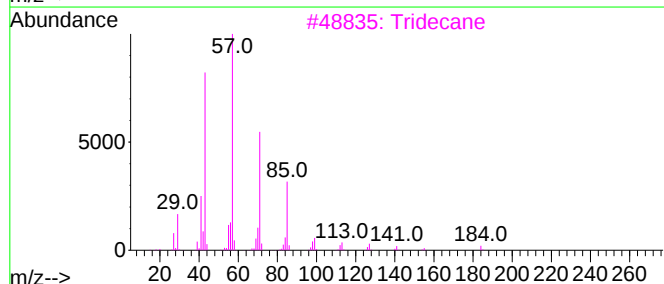
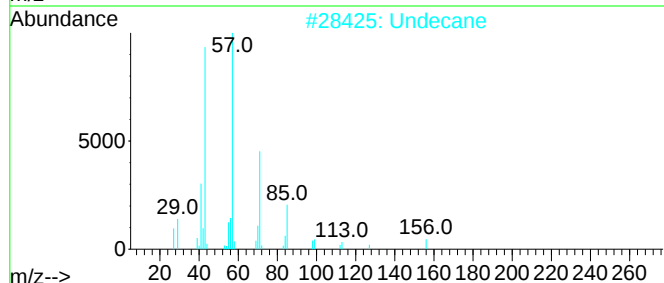
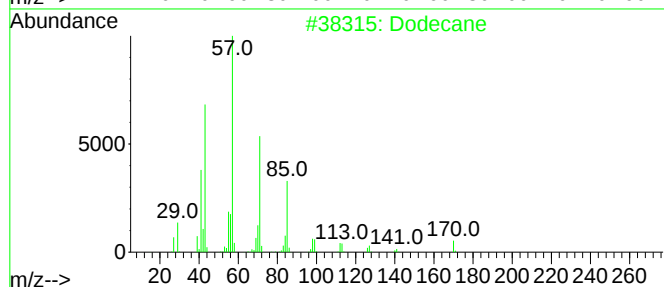
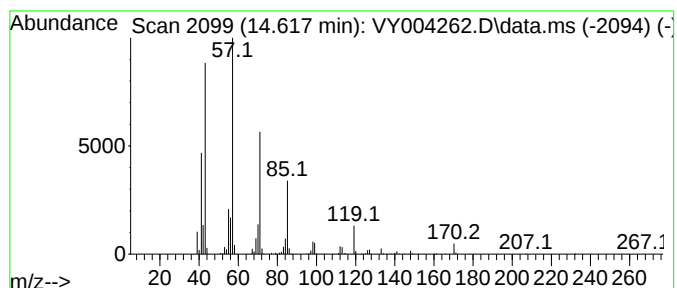
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.617	24.74 ug/l	668190	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	93
2			Undecane	156	C11H24	001120-21-4	80
3			Tridecane	184	C13H28	000629-50-5	80
4			Tetradecane	198	C14H30	000629-59-4	80
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY004262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

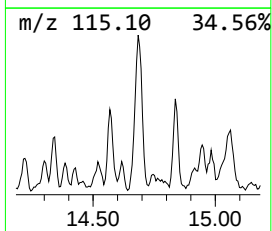
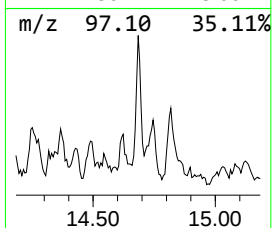
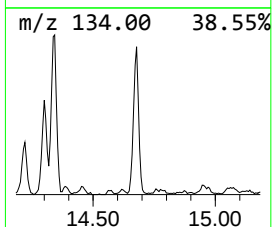
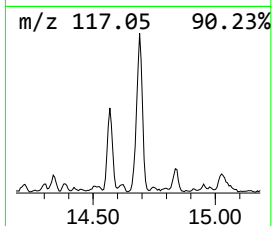
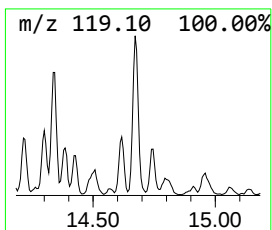
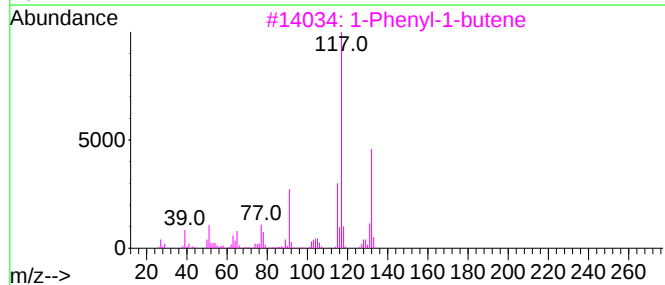
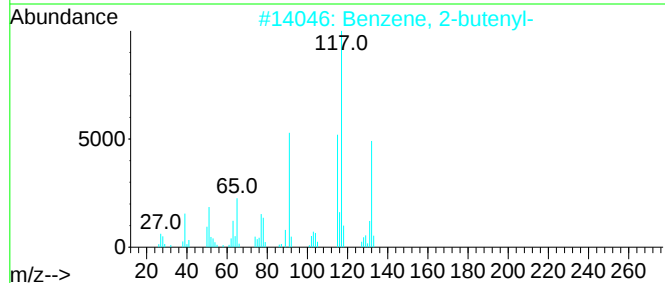
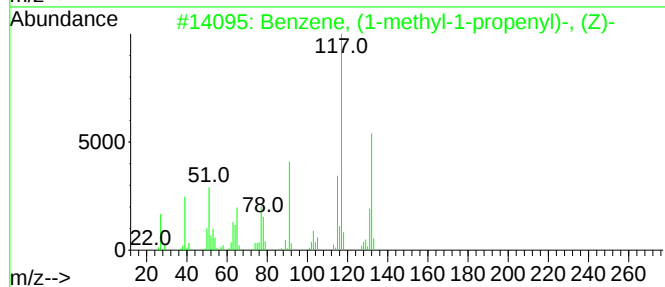
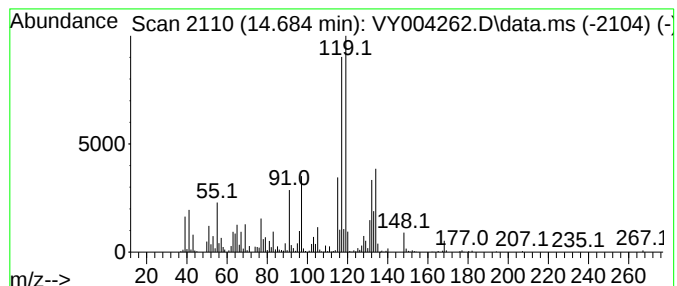
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, (1-methyl-1-propen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	12.68 ug/l	342390	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	86
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	84
4			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	68
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY04262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

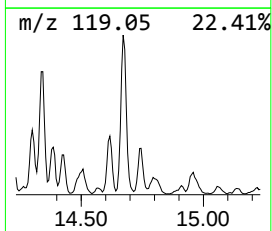
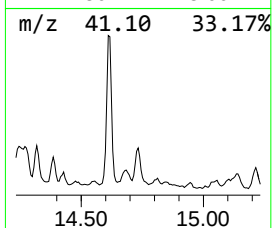
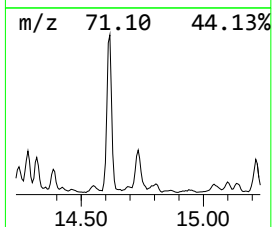
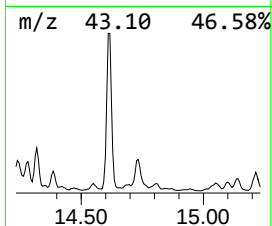
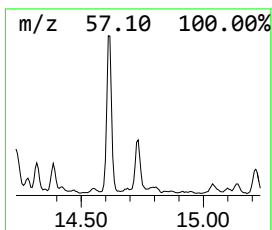
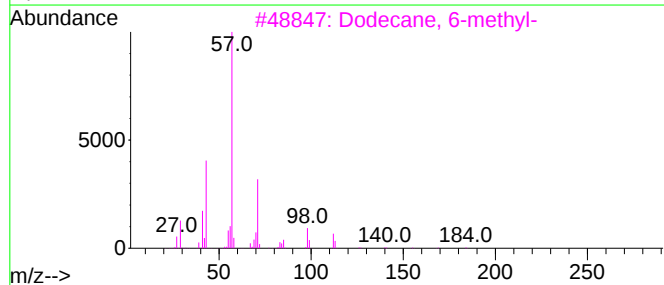
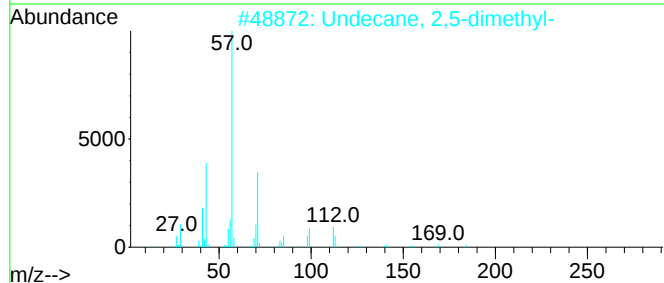
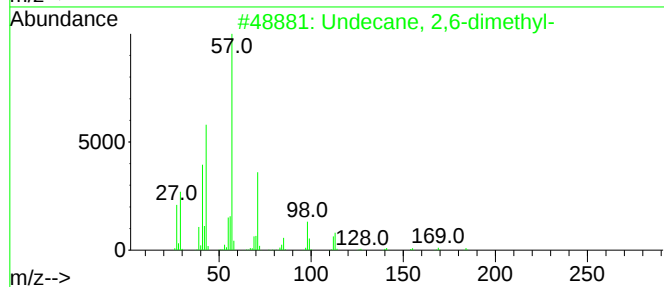
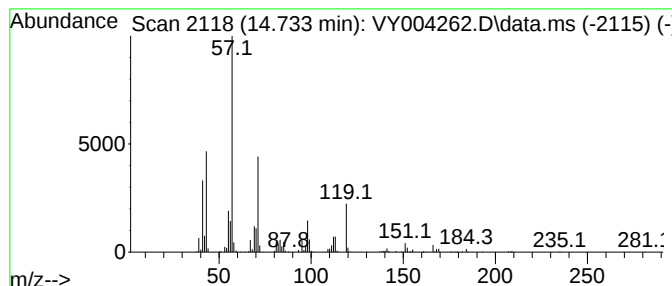
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Undecane, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	8.68 ug/l	234419	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	76
2			Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	60
3			Dodecane, 6-methyl-	184	C13H28	006044-71-9	53
4			Nonane, 3-methyl-	142	C10H22	005911-04-6	53
5			Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY04262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

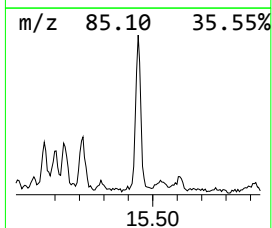
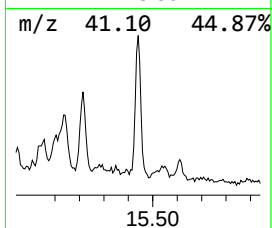
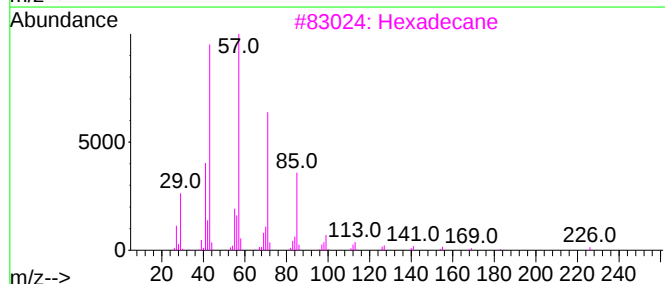
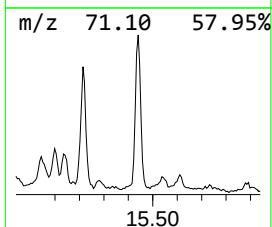
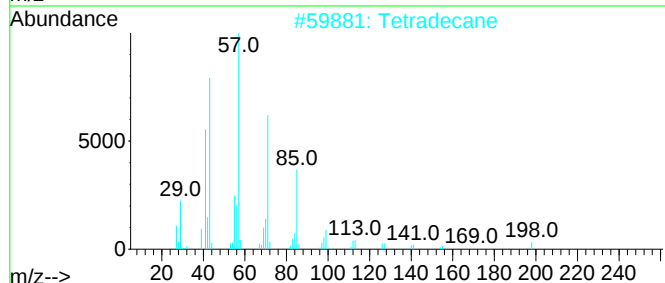
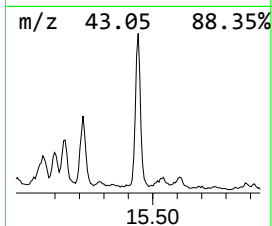
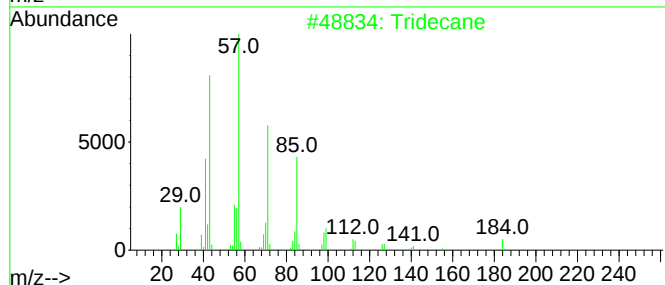
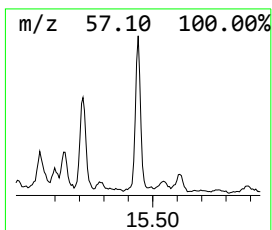
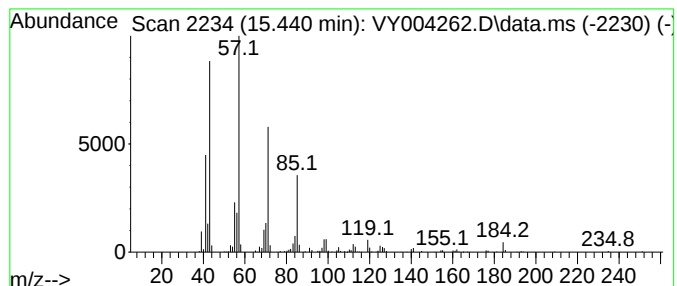
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.440	7.79 ug/l	210449	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	97
2			Tetradecane	198	C14H30	000629-59-4	86
3			Hexadecane	226	C16H34	000544-76-3	80
4			Dodecane	170	C12H26	000112-40-3	78
5			Undecane	156	C11H24	001120-21-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY040221\
Data File : VY004262.D
Acq On : 02 Apr 2021 15:02
Operator : SY/MD
Sample : M1911-02
Misc : 5.05g/5mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-6S1-(0-2)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	12.739	8.0	ug/l	216016	4	13.428	1350620	50.0
Octane, 3,3-dim...	13.044	10.6	ug/l	287296	4	13.428	1350620	50.0
Undecane	13.757	58.5	ug/l	1579490	4	13.428	1350620	50.0
3-Phenylbut-1-ene	14.081	7.6	ug/l	204687	4	13.428	1350620	50.0
Cyclohexane, pe...	14.276	23.1	ug/l	624309	4	13.428	1350620	50.0
Undecane, 2-met...	14.318	9.4	ug/l	252744	4	13.428	1350620	50.0
Dodecane	14.617	24.7	ug/l	668190	4	13.428	1350620	50.0
Benzene, (1-met...	14.684	12.7	ug/l	342390	4	13.428	1350620	50.0
Undecane, 2,6-d...	14.733	8.7	ug/l	234419	4	13.428	1350620	50.0
Tridecane	15.440	7.8	ug/l	210449	4	13.428	1350620	50.0