

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
 Data File : VY004226.D
 Acq On : 30 Mar 2021 11:45
 Operator : SY/MD
 Sample : M1836-08
 Misc : 9.12g/5mL/MSVOA_Y/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 B4-0-2

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: VY004226.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.083	36	43	54	rBV	155846	255302	2.60%	0.274%
2	2.253	61	71	82	rBV	976703	1626933	16.56%	1.748%
3	2.912	170	179	219	rVB	3515213	7751291	78.90%	8.327%
4	3.265	228	237	260	rVB	1147238	2856972	29.08%	3.069%
5	3.454	260	268	279	rVB	56281	129222	1.32%	0.139%
6	3.710	300	310	329	rVB	792393	2062991	21.00%	2.216%
7	3.954	336	350	371	rVB	137486	448778	4.57%	0.482%
8	4.716	463	475	479	rVV2	474936	1558205	15.86%	1.674%
9	4.783	480	486	526	rVB3	846823	3566586	36.31%	3.831%
10	5.253	546	563	591	rBV	520804	1879228	19.13%	2.019%
11	5.570	602	615	632	rBV3	46651	156654	1.59%	0.168%
12	5.765	632	647	666	rBV	452576	1413812	14.39%	1.519%
13	6.112	684	704	717	rBV	240626	786188	8.00%	0.845%
14	6.253	717	727	737	rBV	113009	348188	3.54%	0.374%
15	6.551	765	776	789	rBV2	231309	705295	7.18%	0.758%
16	6.704	789	801	809	rVV	358870	1146014	11.67%	1.231%
17	6.807	809	818	827	rVV	670142	2066537	21.04%	2.220%
18	6.887	827	831	840	rVV2	67431	177314	1.80%	0.190%
19	7.539	930	938	948	rVB	292513	729379	7.42%	0.784%
20	7.783	958	978	987	rBV6	738739	2605577	26.52%	2.799%
21	7.880	987	994	1006	rVB	709961	1823920	18.57%	1.959%
22	8.008	1006	1015	1023	rBV	585906	1362480	13.87%	1.464%
23	8.173	1031	1042	1052	rBV	2005927	4522408	46.04%	4.858%
24	8.338	1052	1069	1079	rBV	3756736	9823582	100.00%	10.553%
25	8.423	1079	1083	1094	rVB	171611	398226	4.05%	0.428%
26	8.551	1094	1104	1116	rVB	372509	822308	8.37%	0.883%
27	8.697	1118	1128	1133	rBV2	504120	1158147	11.79%	1.244%
28	8.856	1149	1154	1161	rVB	75908	137108	1.40%	0.147%
29	8.978	1169	1174	1187	rVB2	98395	241353	2.46%	0.259%
30	9.191	1195	1209	1214	rBV2	1136053	2630827	26.78%	2.826%
31	9.246	1214	1218	1225	rVV	509820	985785	10.03%	1.059%
32	9.325	1225	1231	1236	rVV	232872	486826	4.96%	0.523%
33	9.374	1236	1239	1245	rVV	154853	261108	2.66%	0.281%
34	9.465	1245	1254	1261	rVV3	77955	194998	1.98%	0.209%
35	9.624	1272	1280	1289	rVB	2600978	4999636	50.89%	5.371%
36	9.758	1289	1302	1309	rBV	3112736	6880492	70.04%	7.392%

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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	9.825	1309	1313	1316	rBV	194295	290809	2.96%	0.312%
38	9.965	1326	1336	1347	rBV	309913	869659	8.85%	0.934%
39	10.063	1347	1352	1356	rBV	98633	181910	1.85%	0.195%
40	10.118	1356	1361	1366	rVV	386762	674342	6.86%	0.724%

41	10.185	1366	1372	1380	rVV2	491678	1009367	10.27%	1.084%
42	10.246	1380	1382	1387	rVV	76590	126190	1.28%	0.136%
43	10.307	1387	1392	1395	rVV5	69250	130361	1.33%	0.140%
44	10.374	1395	1403	1409	rVB4	262451	585405	5.96%	0.629%
45	10.526	1424	1428	1432	rBV2	69852	117274	1.19%	0.126%

46	10.612	1437	1442	1453	rVB4	172820	483576	4.92%	0.519%
47	10.715	1454	1459	1464	rVB3	54559	101167	1.03%	0.109%
48	10.800	1464	1473	1478	rBV4	43075	104178	1.06%	0.112%
49	10.910	1487	1491	1497	rVV	62241	124097	1.26%	0.133%
50	11.008	1503	1507	1509	rVV	69840	111013	1.13%	0.119%

51	11.038	1509	1512	1517	rVV	76534	130154	1.32%	0.140%
52	11.282	1544	1552	1563	rVB3	123323	349845	3.56%	0.376%
53	11.386	1563	1569	1578	rBV	97575	204216	2.08%	0.219%
54	11.495	1581	1587	1592	rBV	330457	545720	5.56%	0.586%
55	11.599	1596	1604	1613	rVB	3751897	5793729	58.98%	6.224%

56	11.703	1613	1621	1632	rBV	564422	1180091	12.01%	1.268%
57	12.032	1665	1675	1683	rBV2	92593	229899	2.34%	0.247%
58	12.331	1719	1724	1730	rBV	475490	727991	7.41%	0.782%
59	12.489	1741	1750	1759	rVB4	354940	693433	7.06%	0.745%
60	12.672	1769	1780	1786	rBV	1131133	1703706	17.34%	1.830%

61	12.739	1786	1791	1793	rVV	95387	145632	1.48%	0.156%
62	12.770	1793	1796	1800	rVV	155664	235287	2.40%	0.253%
63	12.812	1800	1803	1815	rVB	116868	195241	1.99%	0.210%
64	12.977	1823	1830	1837	rBV	241086	378368	3.85%	0.406%
65	13.123	1849	1854	1860	rVB	342533	495617	5.05%	0.532%

66	13.251	1868	1875	1879	rBV3	93493	242039	2.46%	0.260%
67	13.428	1898	1904	1906	rBV	255624	385437	3.92%	0.414%
68	13.459	1906	1909	1918	rVB	316304	482896	4.92%	0.519%
69	13.593	1925	1931	1933	rBV	266850	382662	3.90%	0.411%
70	13.629	1933	1937	1941	rVV	843431	1304217	13.28%	1.401%

71	13.672	1941	1944	1953	rVB3	163680	375809	3.83%	0.404%
72	13.757	1953	1958	1963	rVB2	70156	108867	1.11%	0.117%
73	13.824	1963	1969	1975	rBV	187606	279293	2.84%	0.300%
74	13.977	1988	1994	1999	rVB	576696	839048	8.54%	0.901%
75	14.032	1999	2003	2007	rBV	67202	103113	1.05%	0.111%

76	14.080	2007	2011	2017	rVB2	276667	432941	4.41%	0.465%
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Instrument :
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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	14.300	2038	2047	2051	rBV	253118	399618	4.07%	0.429%
78	14.568	2086	2091	2096	rBV	134779	207174	2.11%	0.223%
79	14.684	2103	2110	2116	rBV3	263641	522133	5.32%	0.561%
80	15.056	2165	2171	2176	rVB2	50092	102876	1.05%	0.111%

Sum of corrected areas: 93086070

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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

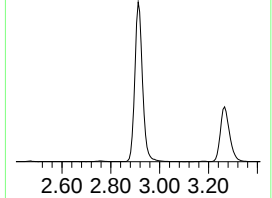
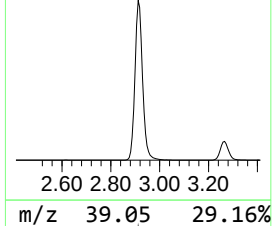
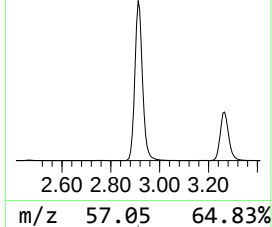
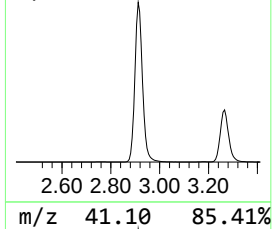
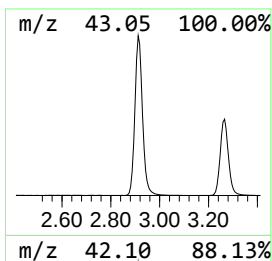
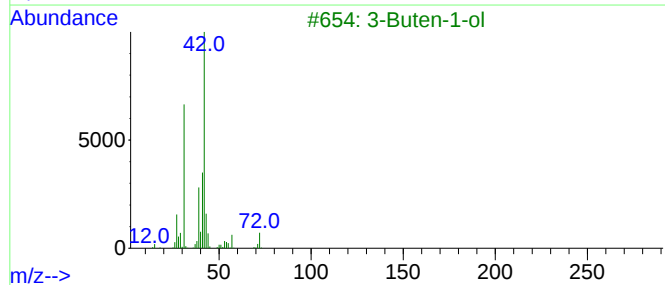
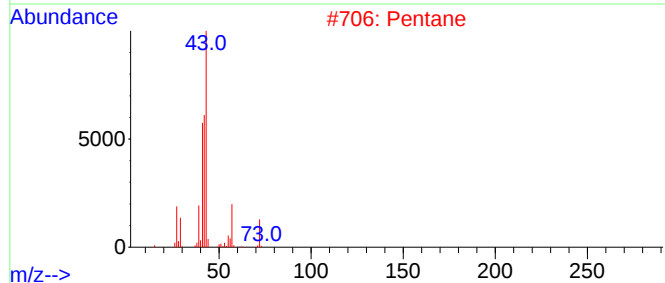
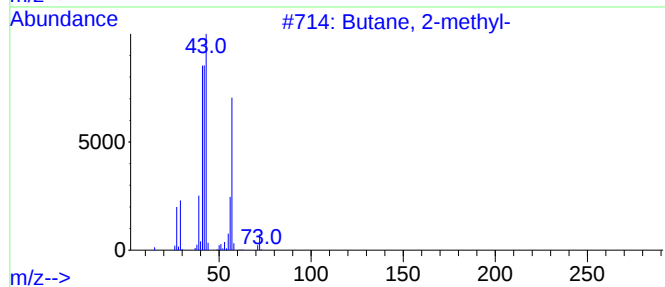
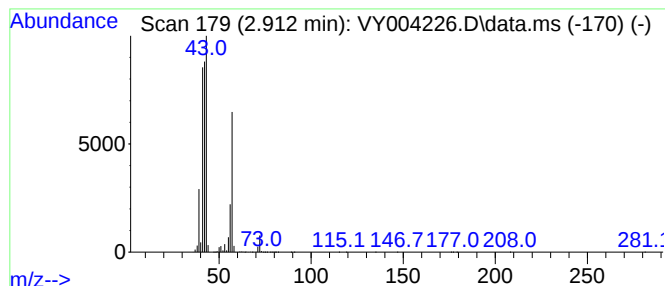
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.912	148.74 ug/l	7751290	Pentafluorobenzene	7.801

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	94
2			Pentane	72	C5H12	000109-66-0	36
3			3-Buten-1-ol	72	C4H8O	000627-27-0	28
4			1-Butene	56	C4H8	000106-98-9	9
5			1-Propene, 2-methyl-	56	C4H8	000115-11-7	9



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B4-0-2

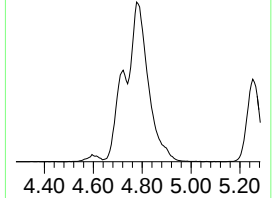
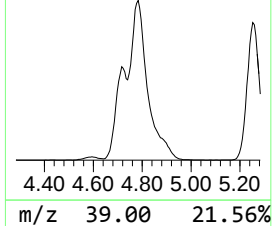
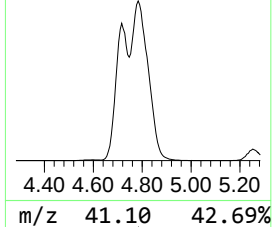
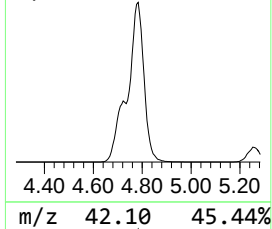
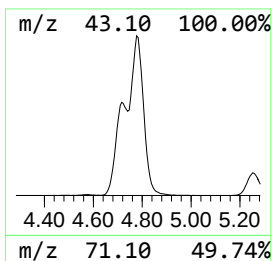
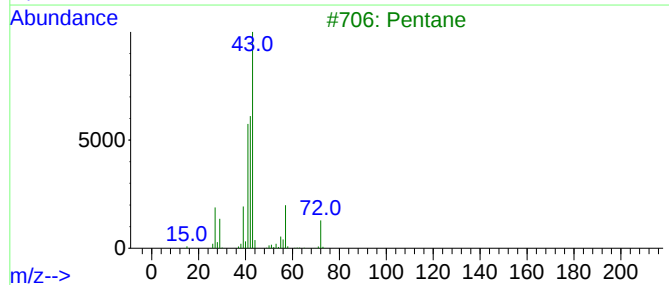
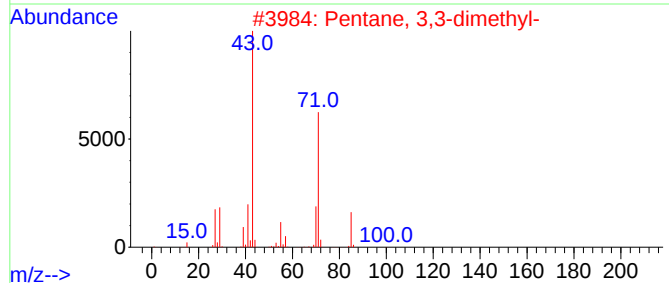
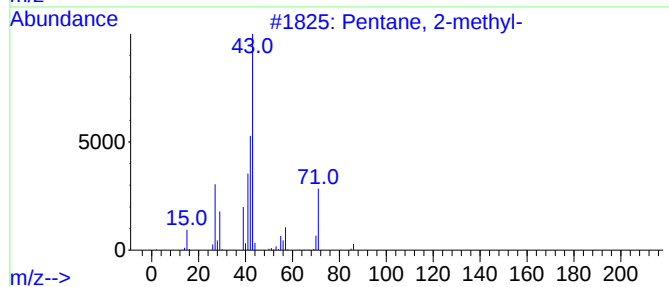
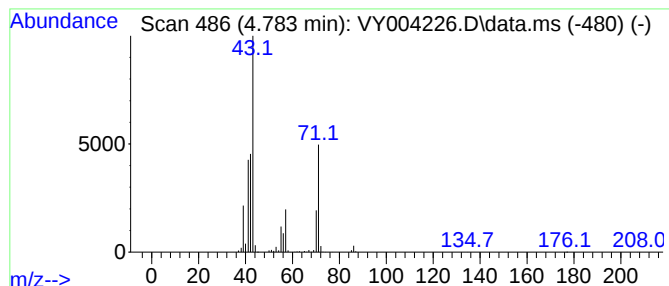
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032921S.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.783	68.44 ug/l	3566590	Pentafluorobenzene	7.801

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	59
2			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	50
3			Pentane	72	C5H12	000109-66-0	43
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	25
5			Heptane, 4-methyl-	114	C8H18	000589-53-7	25



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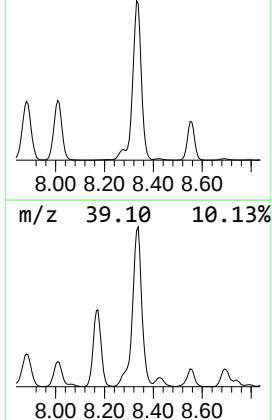
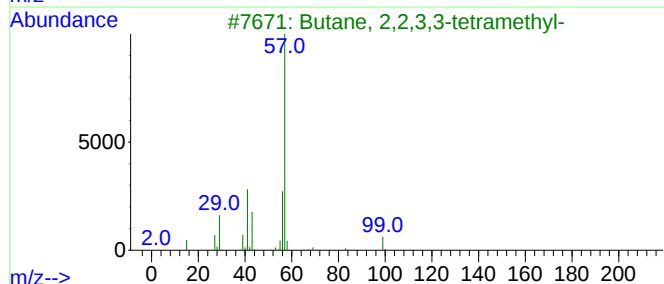
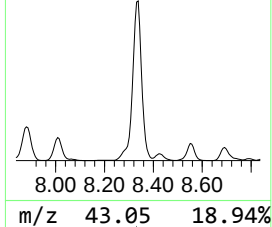
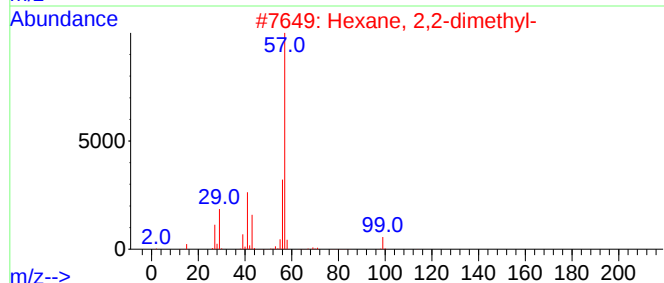
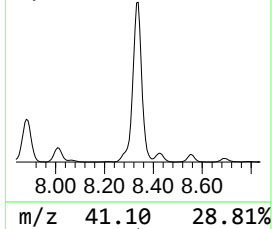
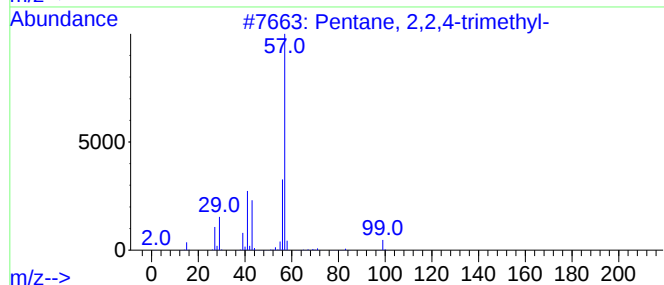
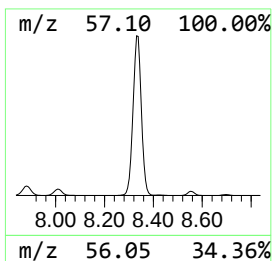
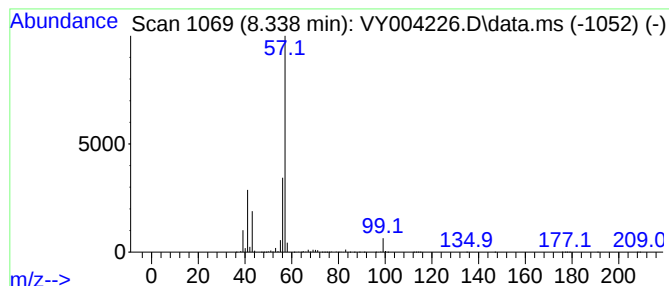
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.338	424.11 ug/l	9823580	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
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			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59



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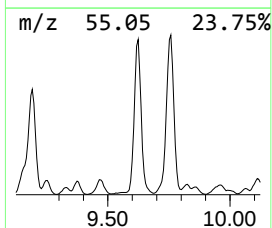
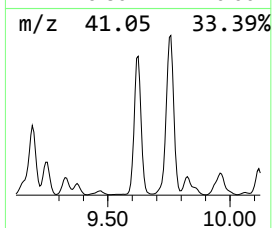
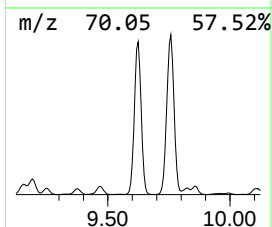
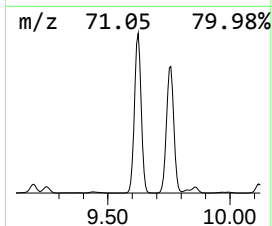
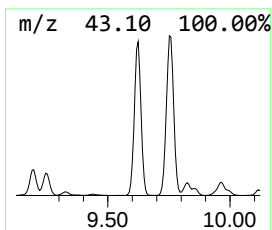
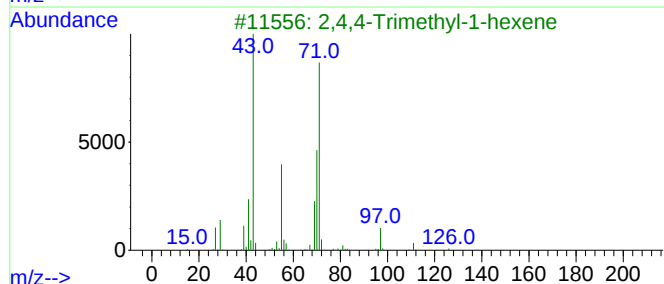
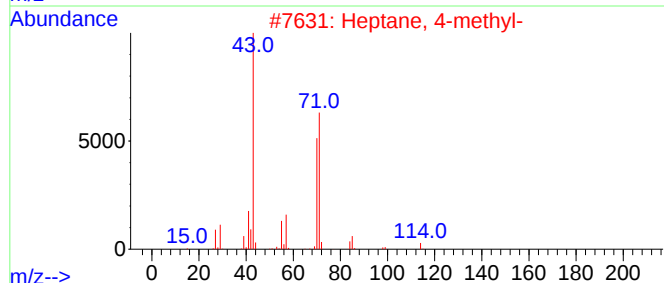
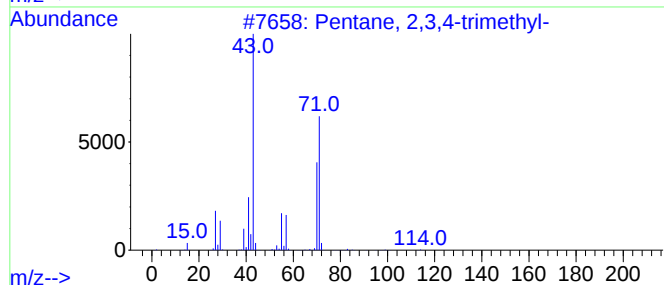
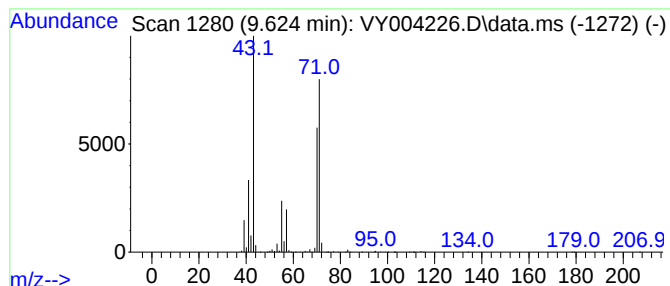
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.624	215.85 ug/l	4999640	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
4			Pentane, 3-ethyl-	100	C7H16	000617-78-7	74
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

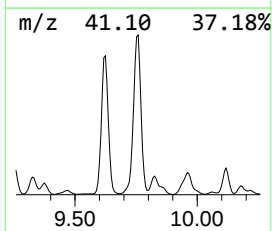
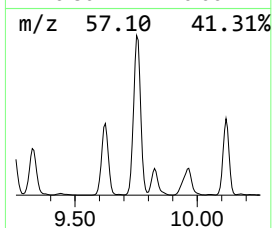
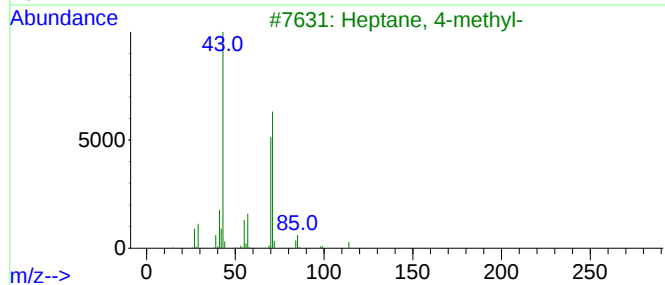
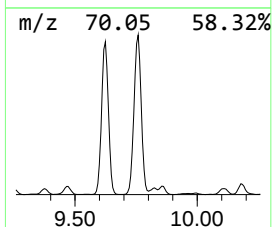
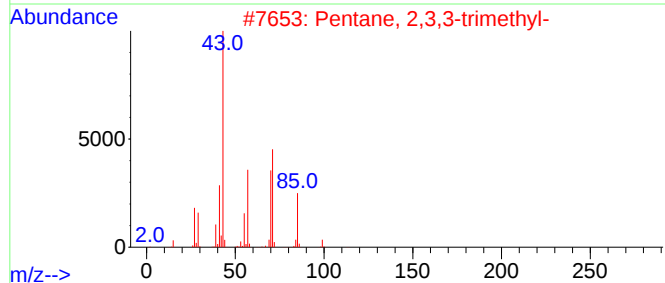
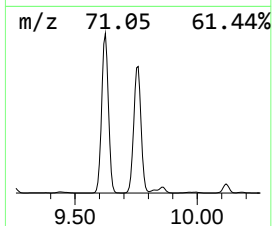
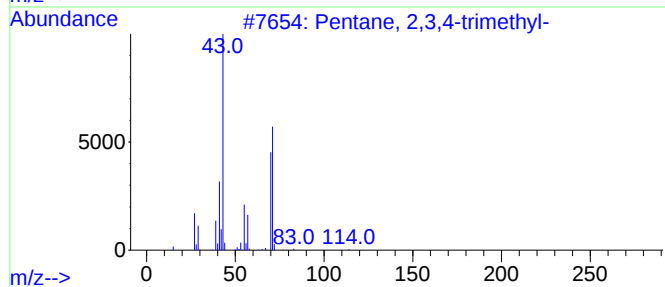
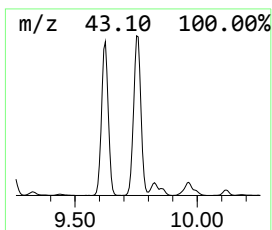
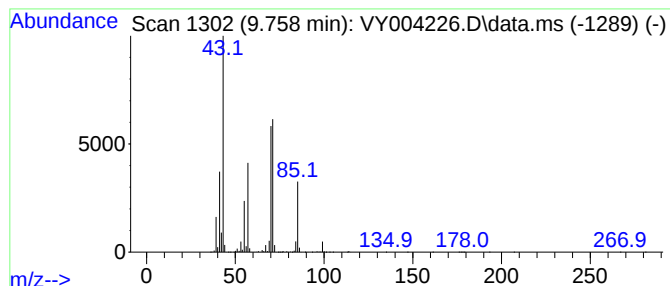
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.758	297.05 ug/l	6880490	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5			Hexane, 3-methyl-	100	C7H16	000589-34-4	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

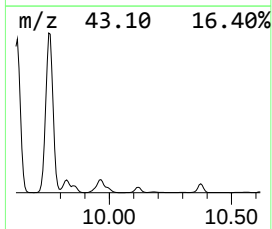
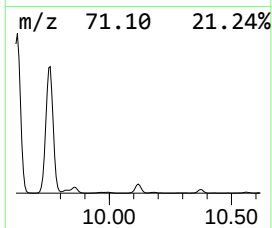
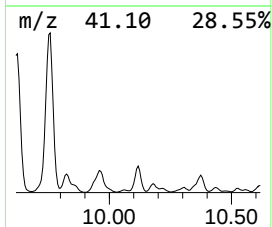
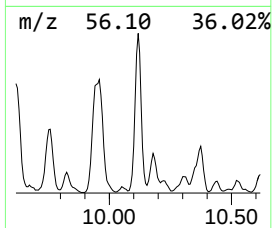
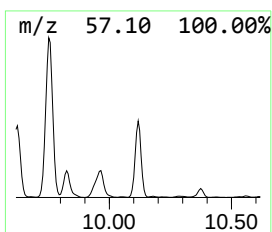
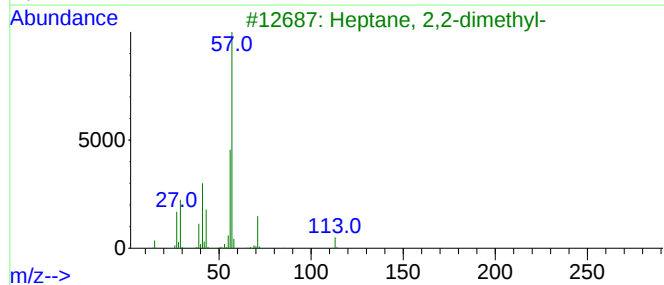
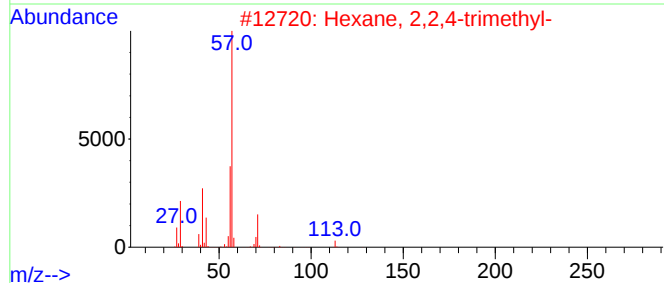
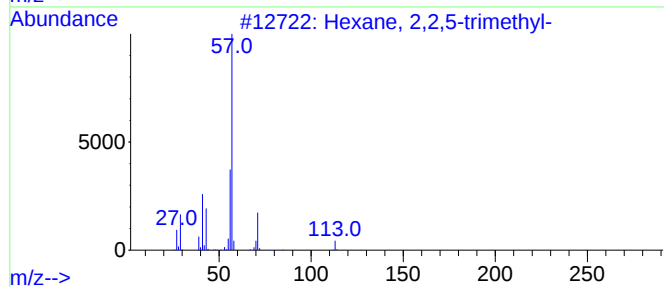
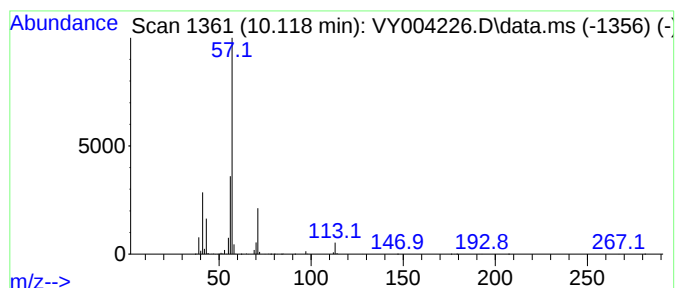
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 2,2,5-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.118	61.78 ug/l	674342	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	83
2			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	83
3			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64
4			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	53
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

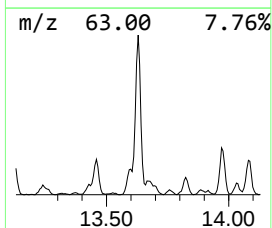
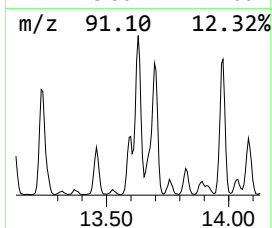
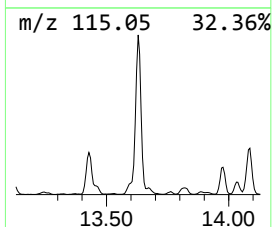
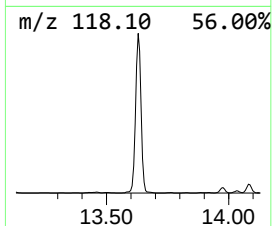
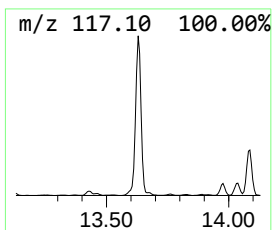
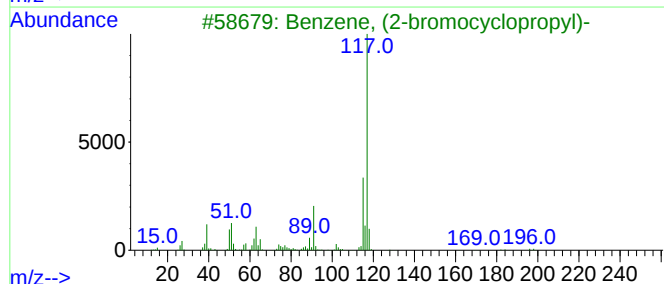
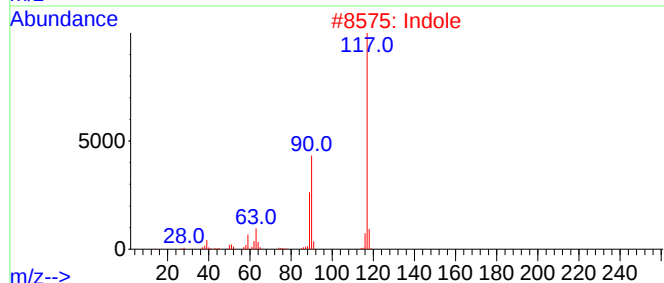
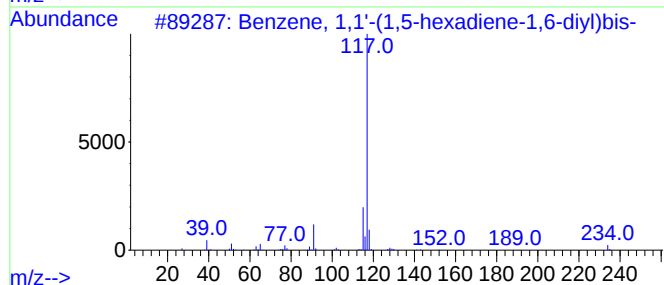
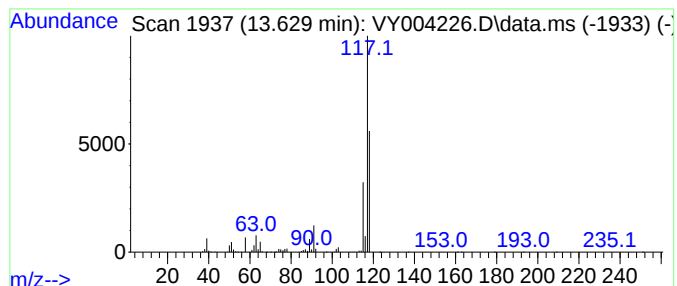
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.629	169.19 ug/l	1304220	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2			Indole	117	C8H7N	000120-72-9	46
3			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	42
4			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42
5			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

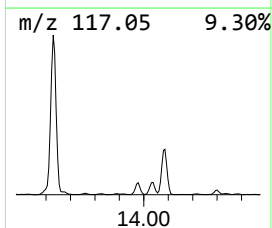
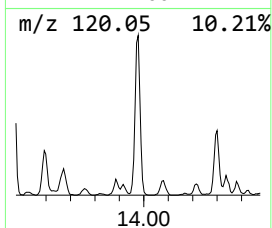
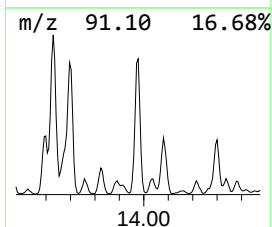
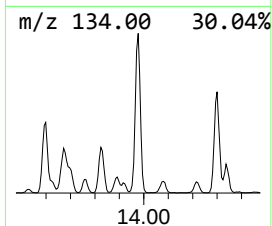
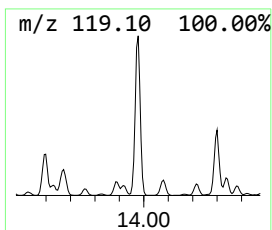
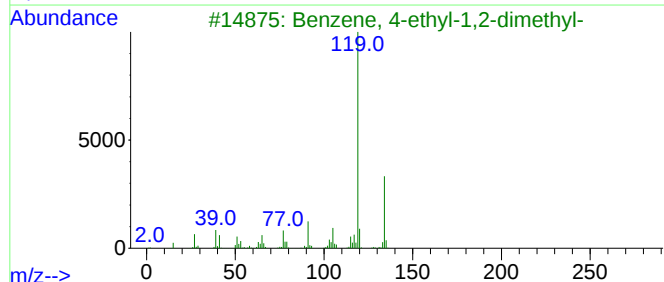
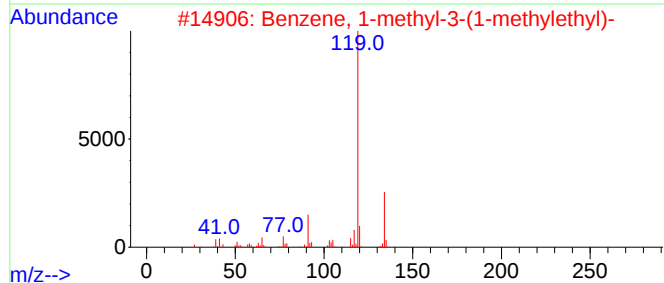
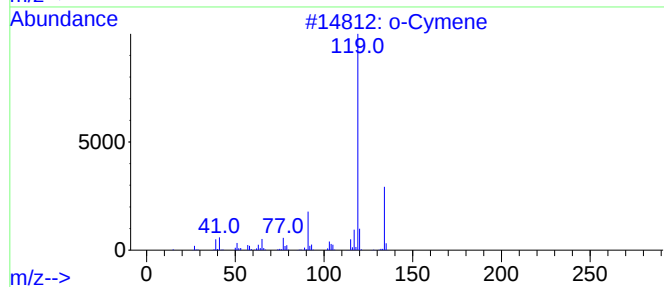
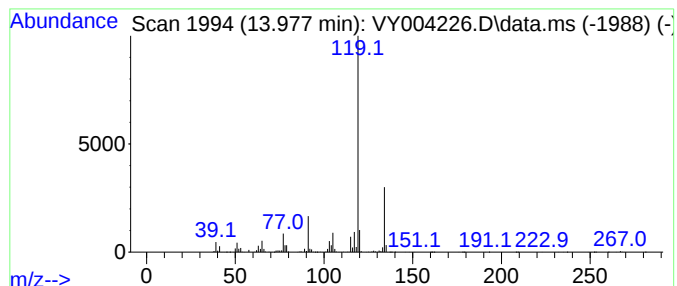
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Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.977	108.84 ug/l	839048	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

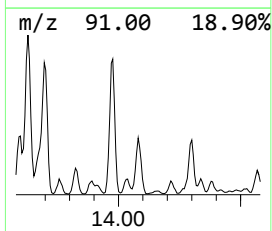
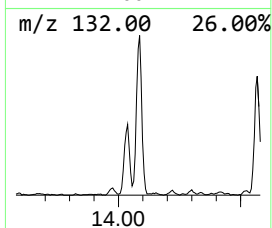
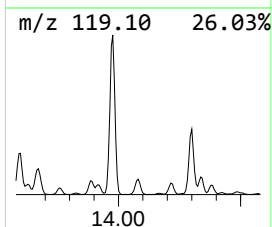
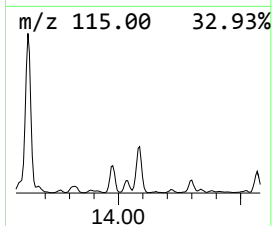
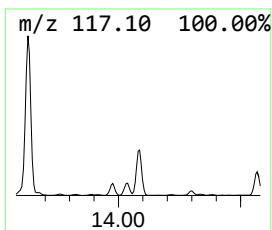
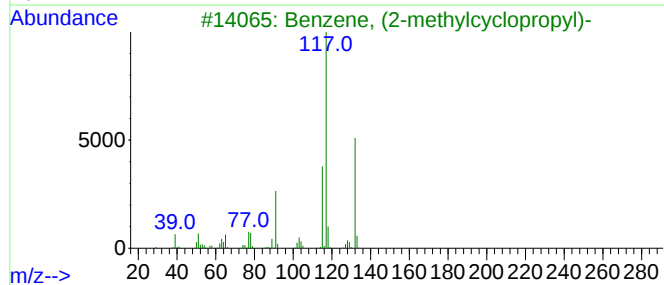
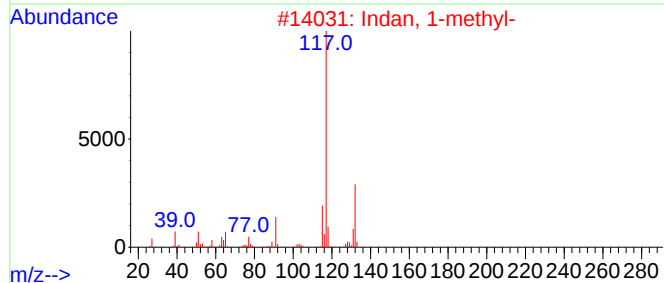
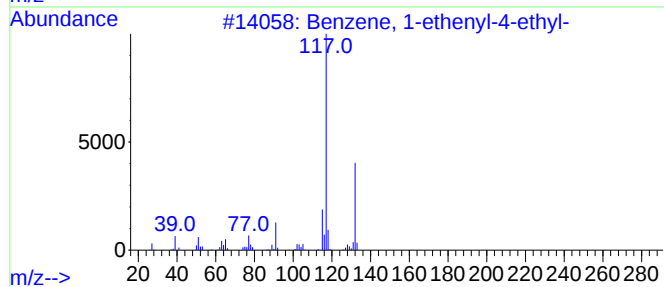
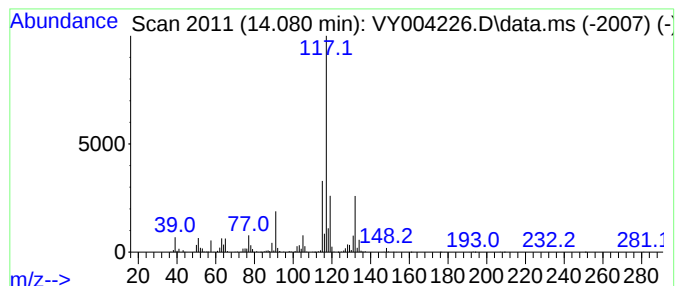
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	56.16 ug/l	432941	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
2			Indan, 1-methyl-	132	C10H12	000767-58-8	76
3			Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
4			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	74
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

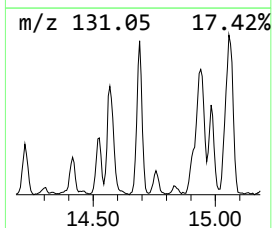
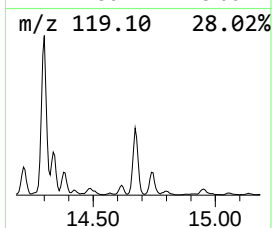
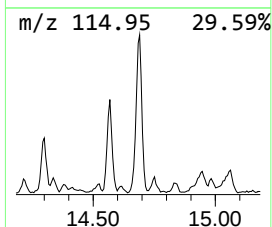
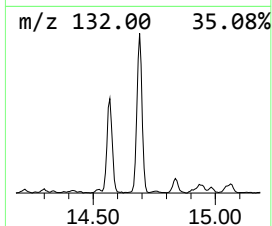
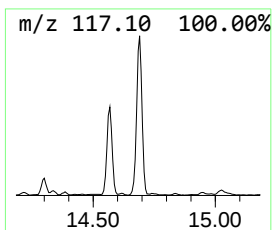
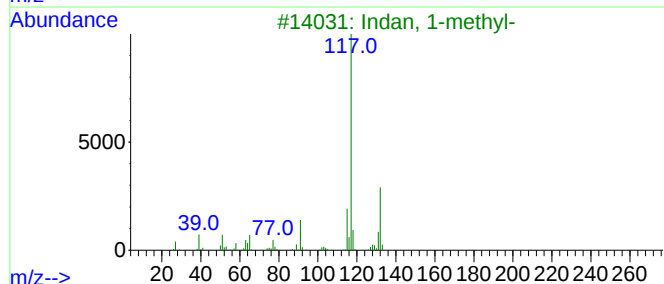
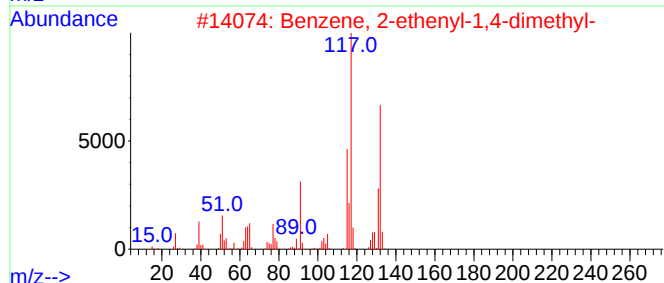
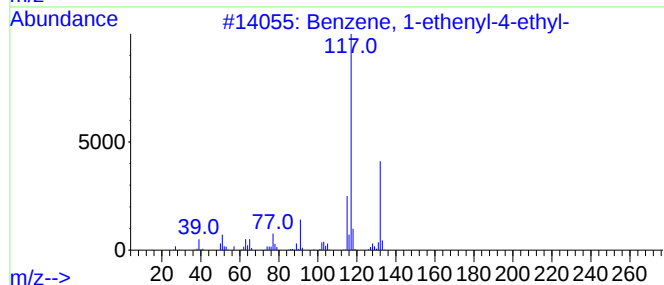
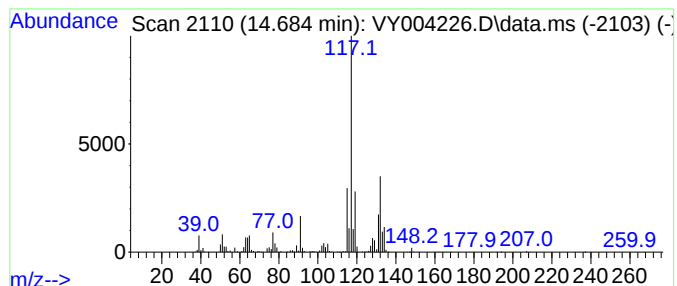
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	67.73 ug/l	522133	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3			Indan, 1-methyl-	132	C10H12	000767-58-8	81
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
5			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033021\
Data File : VY004226.D
Acq On : 30 Mar 2021 11:45
Operator : SY/MD
Sample : M1836-08
Misc : 9.12g/5mL/MSVOA_Y/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B4-0-2

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
Butane, 2-methyl-	2.912	148.7	ug/l	7751290	1	7.801	2605580	50.0
Pentane, 2-methyl-	4.783	68.4	ug/l	3566590	1	7.801	2605580	50.0
Pentane, 2,2,4-...	8.338	424.1	ug/l	9823580	2	8.697	1158150	50.0
Pentane, 2,3,4-...	9.624	215.8	ug/l	4999640	2	8.697	1158150	50.0
Pentane, 2,3,3-...	9.758	297.1	ug/l	6880490	2	8.697	1158150	50.0
Hexane, 2,2,5-t...	10.118	61.8	ug/l	674342	3	11.496	545720	50.0
Benzene, 1,1'-(...	13.629	169.2	ug/l	1304220	4	13.428	385437	50.0
o-Cymene	13.977	108.8	ug/l	839048	4	13.428	385437	50.0
Benzene, 1-ethe...	14.080	56.2	ug/l	432941	4	13.428	385437	50.0
Benzene, 2-ethe...	14.684	67.7	ug/l	522133	4	13.428	385437	50.0