

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
 Data File : VY002984.D
 Acq On : 22 Jun 2020 19:21
 Operator : SY/MD
 Sample : L3082-03
 Misc : 3.56G/5ML/MSVOA Y/SOIL
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SB-3

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.724	475	490	510	rBV3	38976	198132	5.46%	0.553%
2	5.249	565	576	587	rBV3	10982	39221	1.08%	0.109%
3	5.755	648	659	670	rBV3	14062	43613	1.20%	0.122%
4	6.694	802	813	822	rBV4	13172	43265	1.19%	0.121%
5	6.797	822	830	838	rVB3	12936	37620	1.04%	0.105%
6	6.986	851	861	881	rBV	49080	144978	4.00%	0.405%
7	7.724	972	982	987	rBV2	190672	464706	12.82%	1.297%
8	7.797	987	994	1002	rVV	346401	845157	23.31%	2.358%
9	7.876	1002	1007	1017	rVB3	26577	67964	1.87%	0.190%
10	8.004	1017	1028	1037	rBV	44483	105400	2.91%	0.294%
11	8.145	1037	1051	1064	rVB2	187989	515143	14.21%	1.437%
12	8.327	1064	1081	1093	rBV	154802	421005	11.61%	1.175%
13	8.547	1107	1117	1129	rVB	28770	72748	2.01%	0.203%
14	8.693	1132	1141	1155	rBV	508366	1000250	27.58%	2.791%
15	9.187	1208	1222	1227	rBV2	72063	169135	4.66%	0.472%
16	9.242	1227	1231	1238	rVB	41709	78299	2.16%	0.218%
17	9.620	1285	1293	1303	rBV	139114	270695	7.47%	0.755%
18	9.748	1303	1314	1321	rBV	184565	411221	11.34%	1.147%
19	9.821	1321	1326	1329	rBV	34384	60575	1.67%	0.169%
20	9.955	1338	1348	1360	rBV2	46448	125309	3.46%	0.350%
21	10.114	1368	1374	1379	rBV	90422	154125	4.25%	0.430%
22	10.181	1379	1385	1390	rVV	774318	1301502	35.89%	3.631%
23	10.242	1390	1395	1402	rVB	393336	665221	18.35%	1.856%
24	10.370	1408	1416	1422	rVB4	45255	92337	2.55%	0.258%
25	10.577	1443	1450	1457	rBV	211863	404234	11.15%	1.128%
26	10.943	1506	1510	1516	rVB	69416	117014	3.23%	0.326%
27	11.278	1557	1565	1568	rBV2	47752	100799	2.78%	0.281%
28	11.315	1568	1571	1577	rVB	86458	140269	3.87%	0.391%
29	11.376	1577	1581	1590	rVB	40996	66096	1.82%	0.184%
30	11.492	1592	1600	1606	rBV	716938	1204137	33.21%	3.360%
31	11.589	1608	1616	1624	rVB	353738	572493	15.79%	1.597%
32	11.699	1625	1634	1645	rBV	1296908	2426306	66.91%	6.770%
33	12.028	1679	1688	1696	rVB	668345	1054136	29.07%	2.941%
34	12.327	1731	1737	1742	rBV	243955	368407	10.16%	1.028%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	12.479	1752	1762	1773	rVB3	719203	1391824	38.38%	3.883%
36	12.625	1781	1786	1788	rBV3	23071	38118	1.05%	0.106%
37	12.668	1788	1793	1798	rVV	465339	698594	19.27%	1.949%
38	12.735	1798	1804	1807	rVV	1136905	1880134	51.85%	5.246%
39	12.766	1807	1809	1812	rVV	590990	714180	19.70%	1.993%
40	12.808	1812	1816	1827	rVB	755030	1159051	31.96%	3.234%
41	12.973	1835	1843	1850	rBV	478688	782915	21.59%	2.184%
42	13.064	1855	1858	1861	rVV	52514	82606	2.28%	0.230%
43	13.119	1861	1867	1873	rVV	2437144	3626101	100.00%	10.117%
44	13.180	1873	1877	1881	rVV3	56567	95191	2.63%	0.266%
45	13.253	1881	1889	1891	rVV3	56667	151115	4.17%	0.422%
46	13.284	1891	1894	1896	rVV	69028	104257	2.88%	0.291%
47	13.314	1896	1899	1903	rVV	72784	121931	3.36%	0.340%
48	13.363	1903	1907	1910	rVV3	46036	87793	2.42%	0.245%
49	13.424	1910	1917	1920	rVV	830819	1408022	38.83%	3.929%
50	13.455	1920	1922	1937	rVV2	551676	859426	23.70%	2.398%
51	13.589	1939	1944	1946	rVV	122679	185475	5.11%	0.518%
52	13.625	1946	1950	1953	rVV2	576114	928821	25.61%	2.592%
53	13.662	1953	1956	1966	rVV3	453631	904718	24.95%	2.524%
54	13.753	1966	1971	1975	rVV2	65877	119448	3.29%	0.333%
55	13.820	1977	1982	1987	rVV	133137	223759	6.17%	0.624%
56	13.881	1987	1992	1995	rVV	216935	354092	9.77%	0.988%
57	13.912	1995	1997	2002	rVV	219700	294623	8.13%	0.822%
58	13.973	2002	2007	2012	rVV	409243	633448	17.47%	1.767%
59	14.028	2012	2016	2019	rVV	34934	58670	1.62%	0.164%
60	14.076	2019	2024	2029	rVV	137796	219191	6.04%	0.612%
61	14.211	2041	2046	2051	rBV	82842	132608	3.66%	0.370%
62	14.296	2053	2060	2063	rVV	272924	456128	12.58%	1.273%
63	14.332	2063	2066	2071	rVV	372566	534626	14.74%	1.492%
64	14.381	2071	2074	2077	rVV	97660	135931	3.75%	0.379%
65	14.418	2077	2080	2083	rVV	53549	81085	2.24%	0.226%
66	14.454	2083	2086	2093	rVV	60200	106625	2.94%	0.298%
67	14.521	2093	2097	2100	rVV	52800	76690	2.11%	0.214%
68	14.564	2100	2104	2109	rVV	170073	277352	7.65%	0.774%
69	14.613	2109	2112	2116	rVV	79042	106153	2.93%	0.296%
70	14.680	2116	2123	2129	rVV4	286364	726197	20.03%	2.026%
71	14.741	2129	2133	2139	rVB2	67128	116252	3.21%	0.324%

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72	14.832	2144	2148	2152	rBV2	29072	45505	1.25%	0.127%
73	14.942	2161	2166	2177	rVB3	95340	224514	6.19%	0.626%
74	15.052	2178	2184	2195	rVB3	70817	165904	4.58%	0.463%
75	15.235	2203	2214	2220	rBV	383516	668193	18.43%	1.864%
76	15.344	2227	2232	2237	rVB4	33561	54573	1.51%	0.152%
77	15.436	2243	2247	2255	rVB5	27643	52277	1.44%	0.146%
78	15.527	2255	2262	2271	rBV	56406	133010	3.67%	0.371%
79	15.692	2279	2289	2291	rBV5	32256	80682	2.23%	0.225%
80	15.814	2304	2309	2315	rBV	24954	45414	1.25%	0.127%
81	15.887	2316	2321	2327	rVB3	24063	50929	1.40%	0.142%
82	16.033	2338	2345	2349	rBV6	16165	37067	1.02%	0.103%
83	16.289	2381	2387	2405	rVB	204766	417694	11.52%	1.165%
84	16.491	2413	2420	2428	rBV	98724	209652	5.78%	0.585%

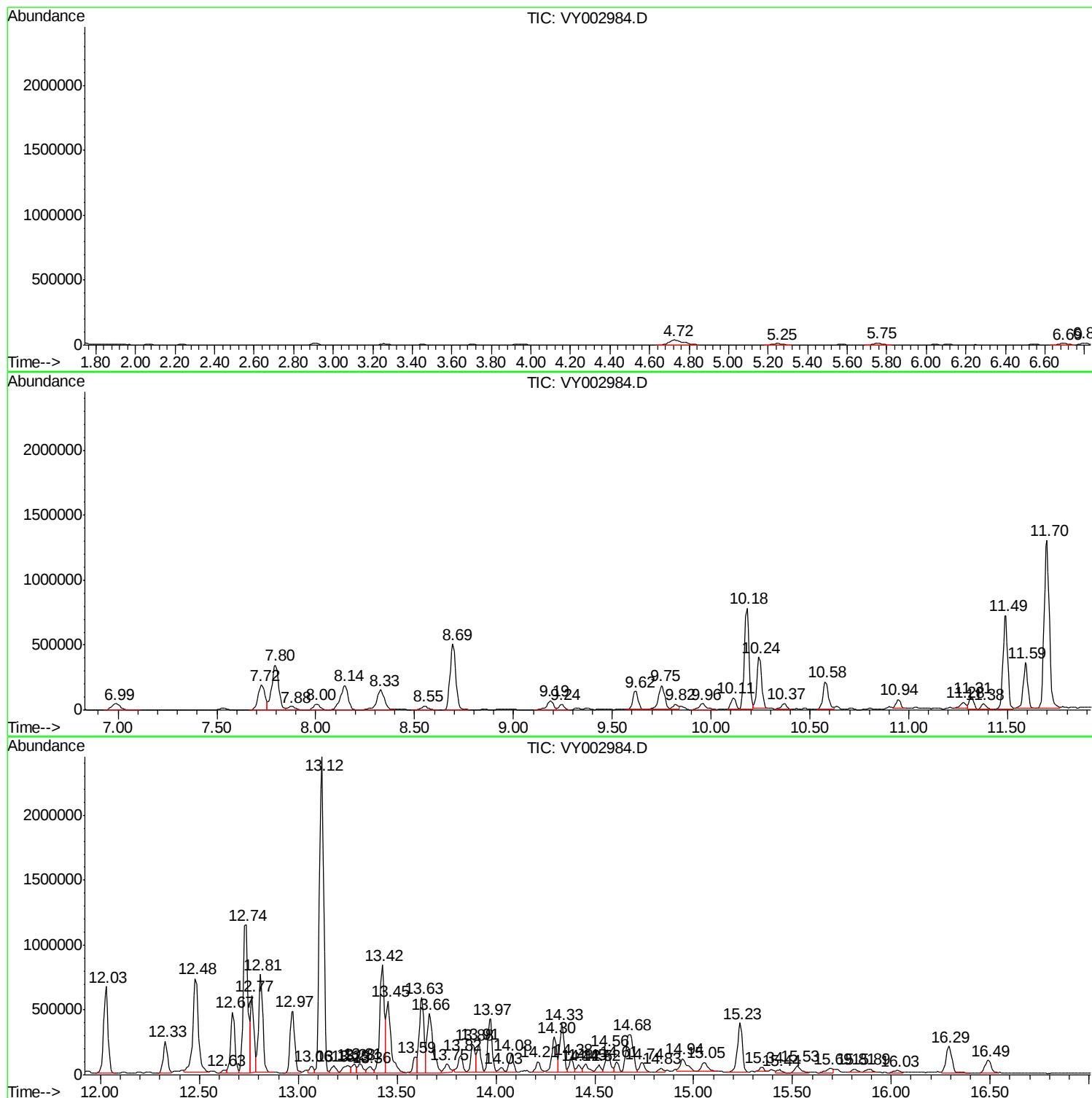
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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



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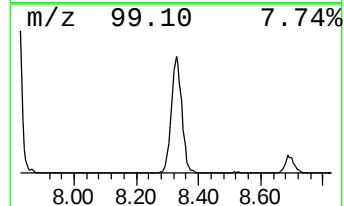
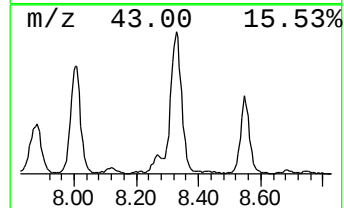
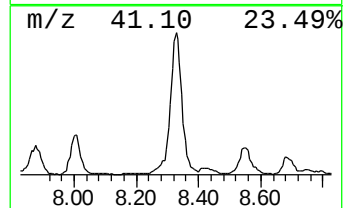
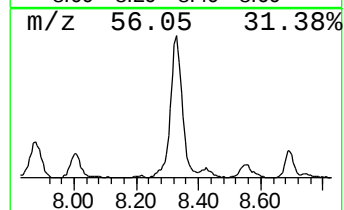
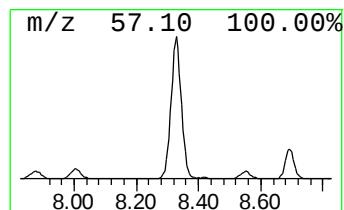
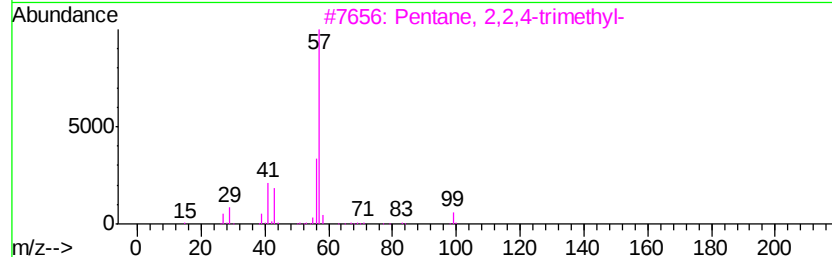
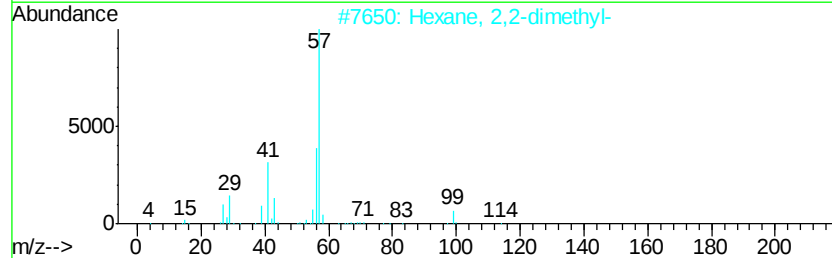
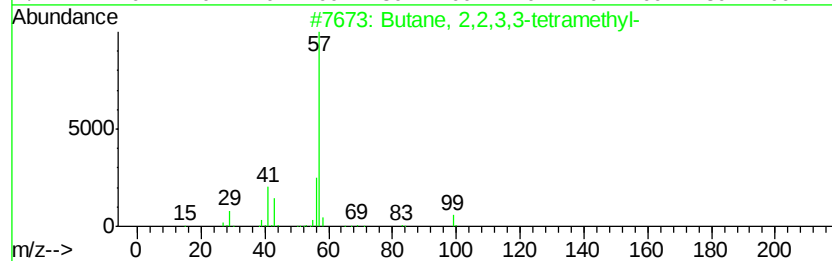
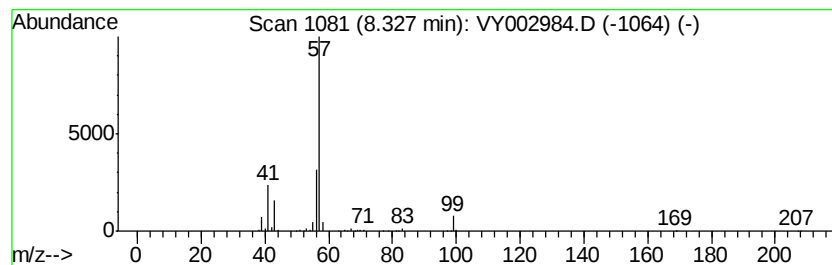
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.33	21.05 ug/l	421005	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



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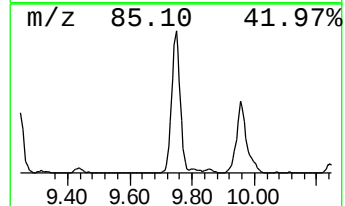
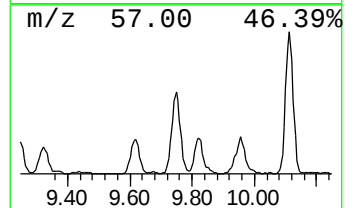
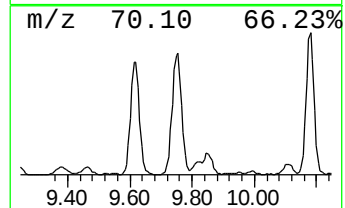
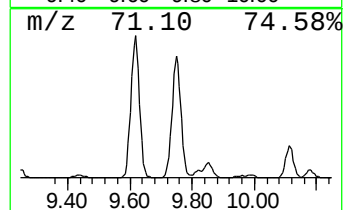
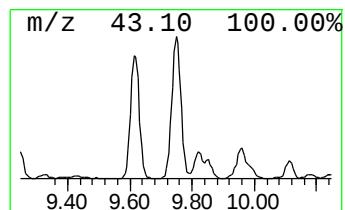
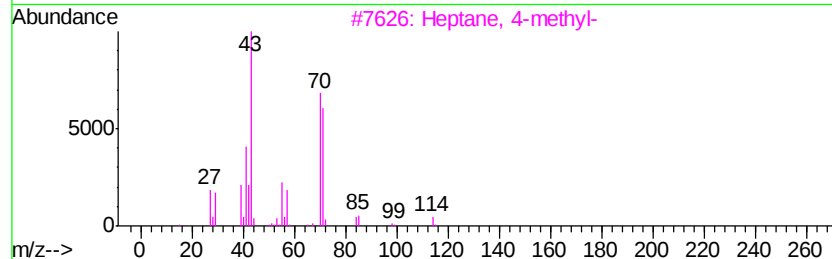
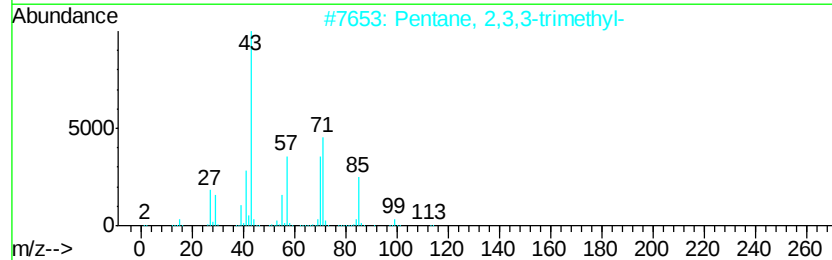
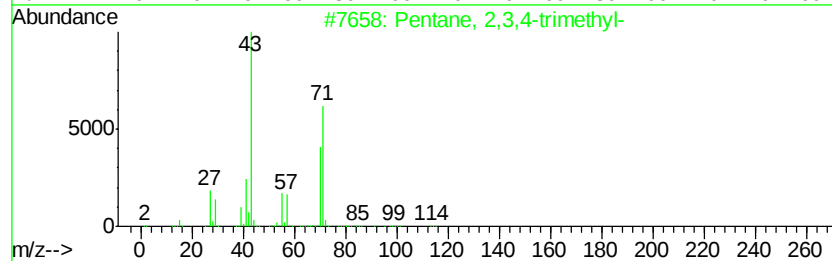
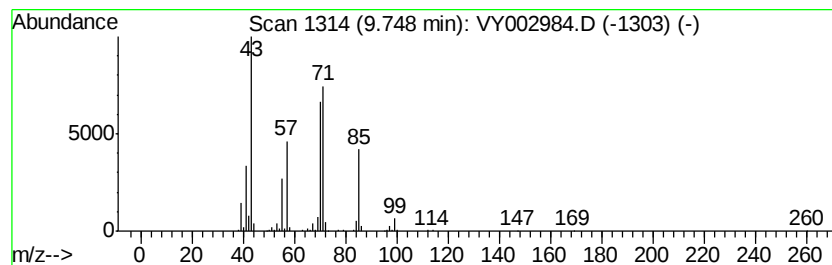
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.75	20.56 ug/l	411221	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	64
2		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	64
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4		Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	59
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	58



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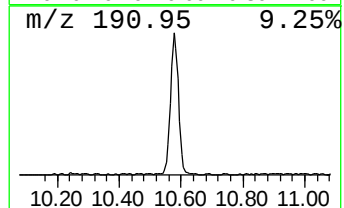
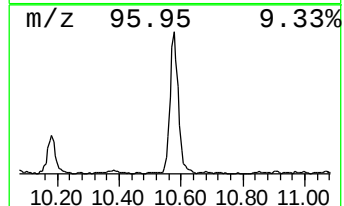
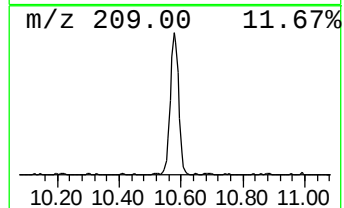
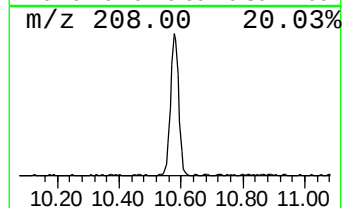
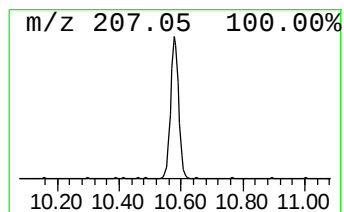
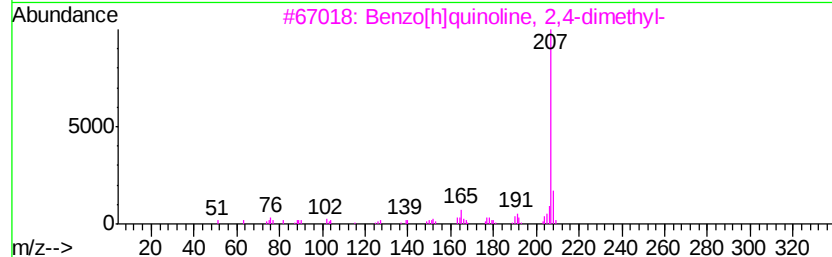
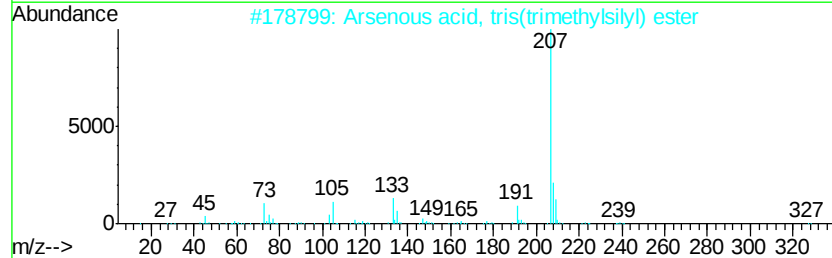
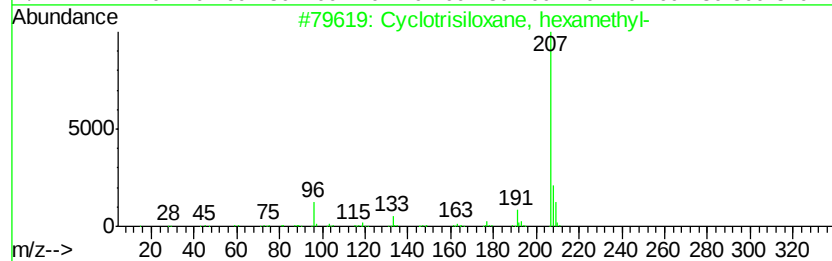
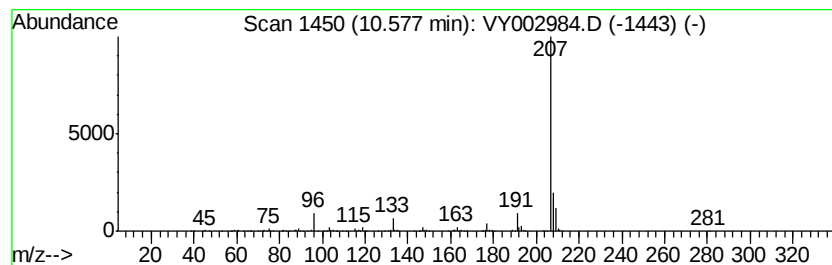
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclotrisiloxane, hexamethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	16.79 ug/l	404234	Chlorobenzene-d5	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	91
2		Arsenous acid, tris(trimethylsil...	342	C9H27AsO3Si3	055429-29-3	56
3		Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	42
4		4-Cyclohexene-1,2-dicarboximide,...	207	C12H17N02	028916-00-9	42
5		1H-Indole, 5-methyl-2-phenyl-	207	C15H13N	013228-36-9	38



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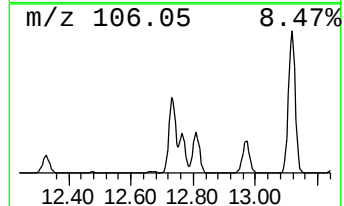
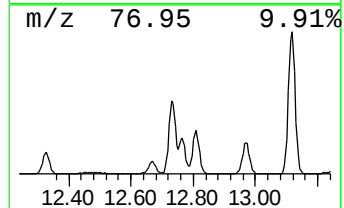
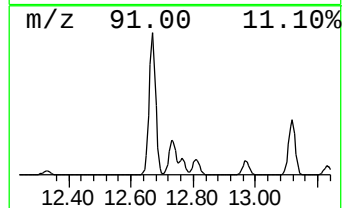
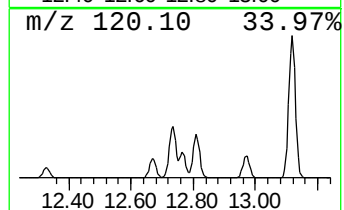
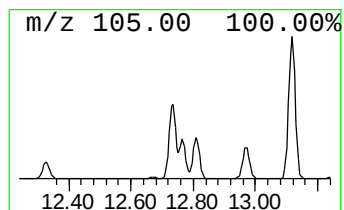
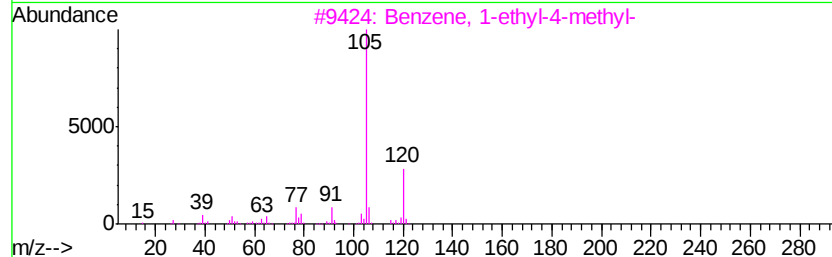
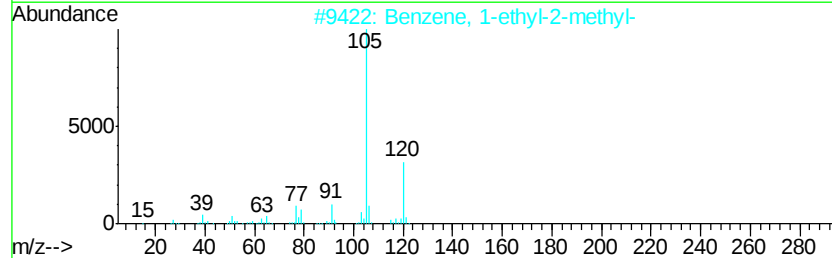
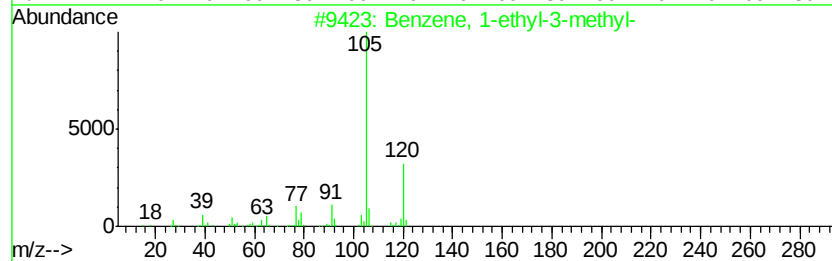
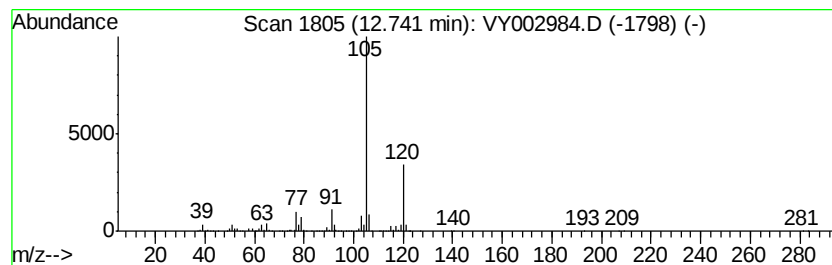
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	66.77 ug/l	1880130	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

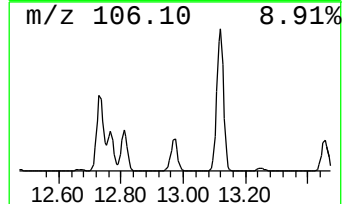
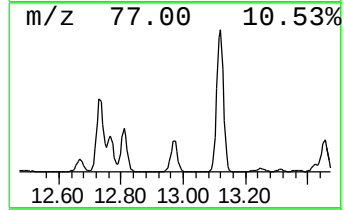
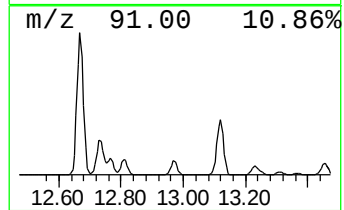
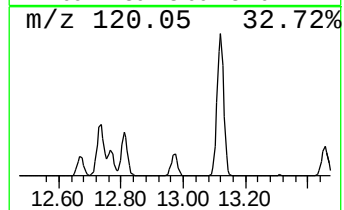
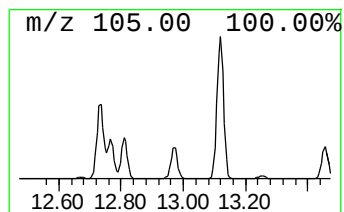
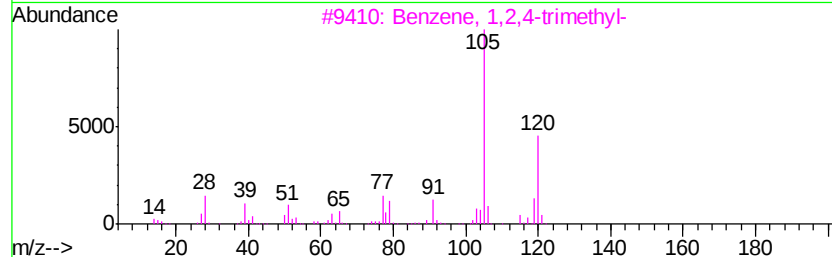
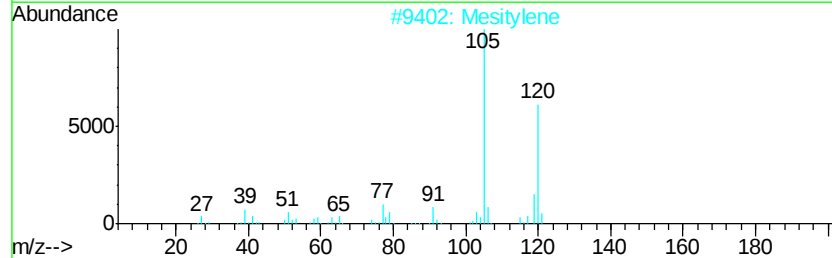
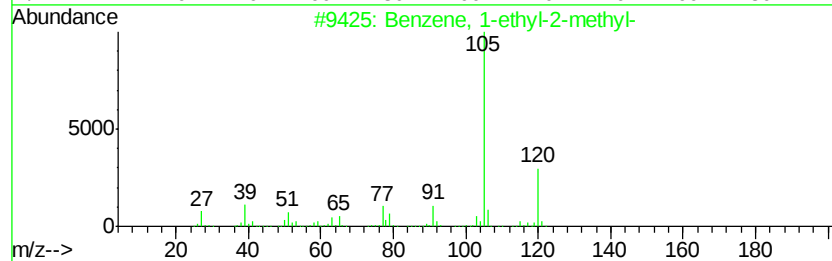
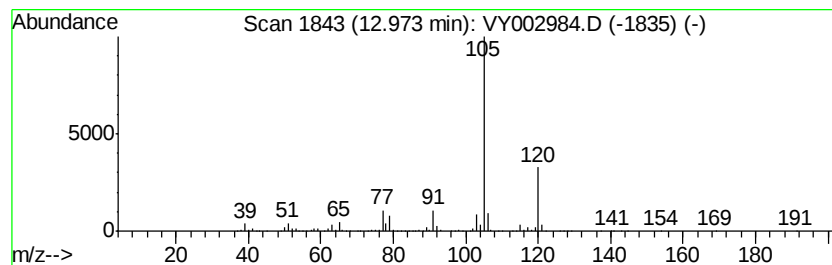
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	27.80 ug/l	782915	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Mesitylene	120	C9H12	000108-67-8	90
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

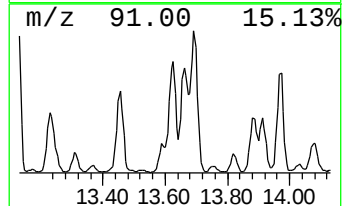
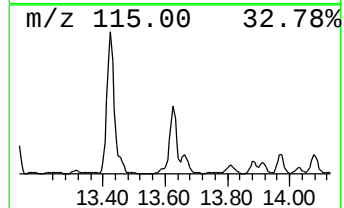
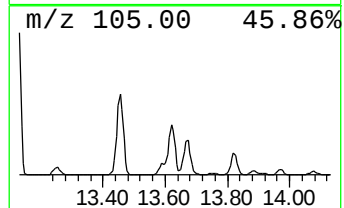
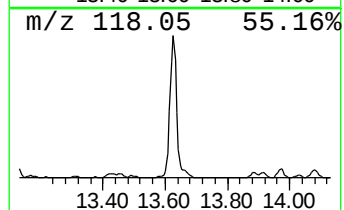
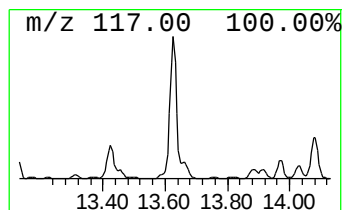
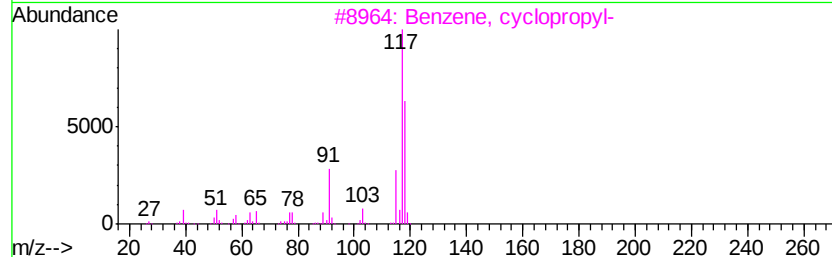
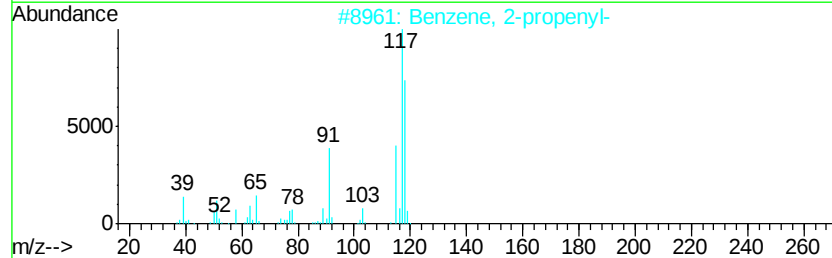
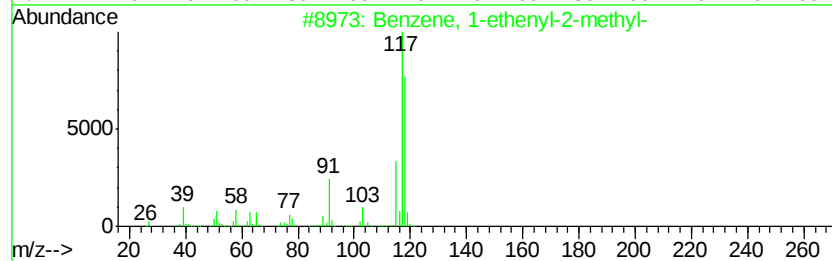
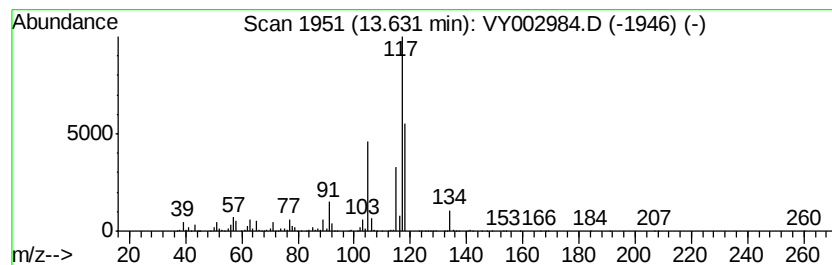
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	32.98 ug/l	928821	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5		Indane	118	C9H10	000496-11-7	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

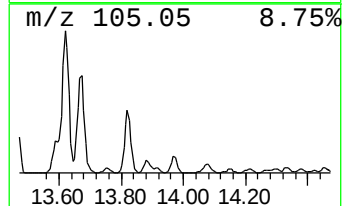
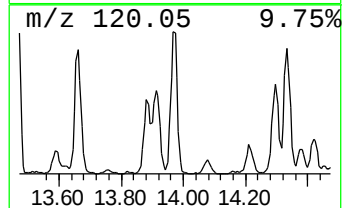
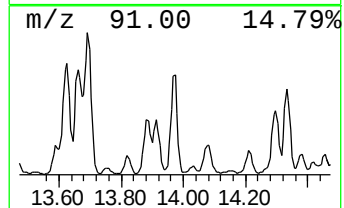
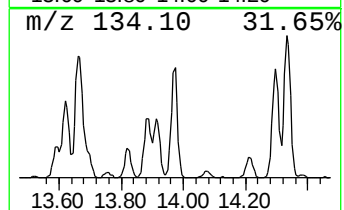
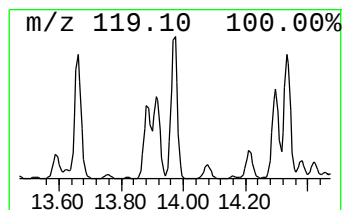
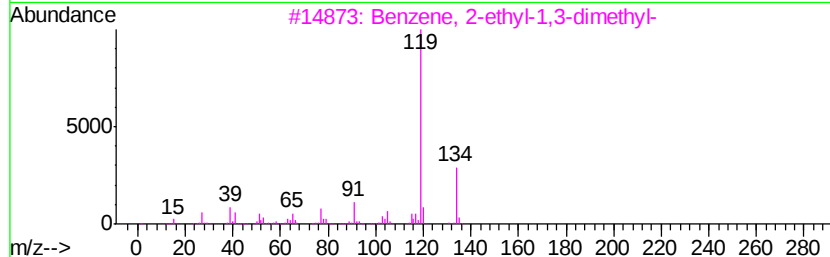
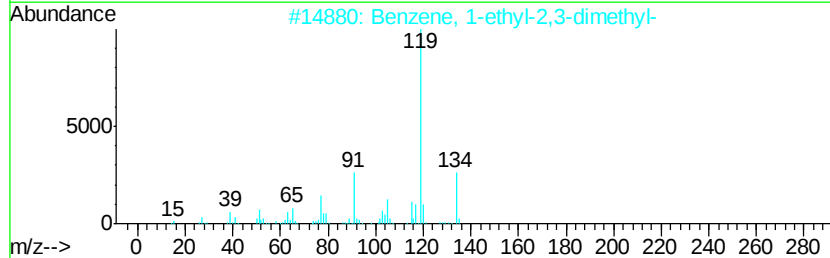
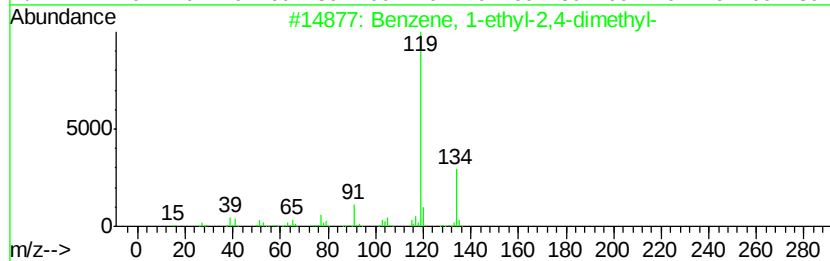
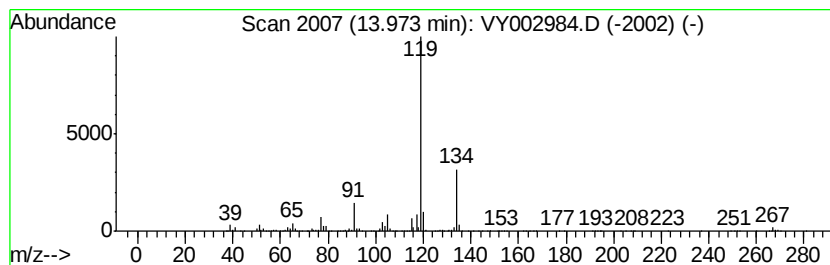
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	22.49 ug/l	633448	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

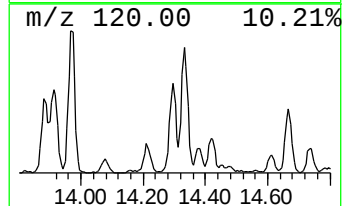
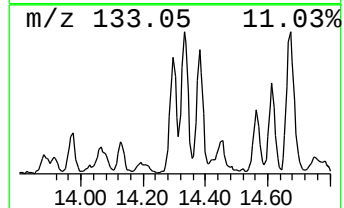
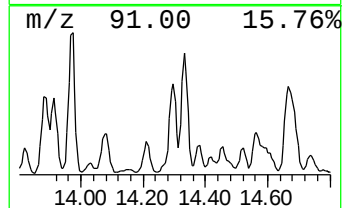
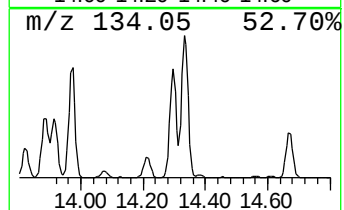
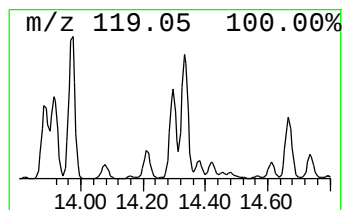
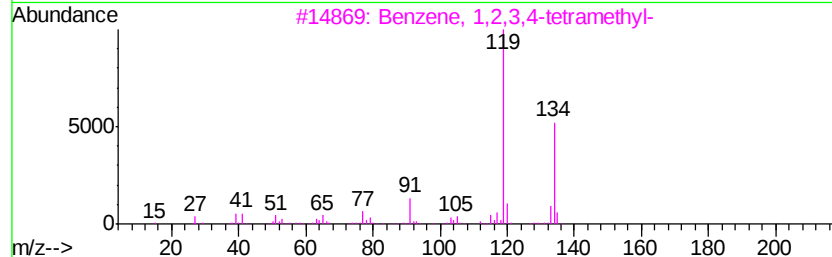
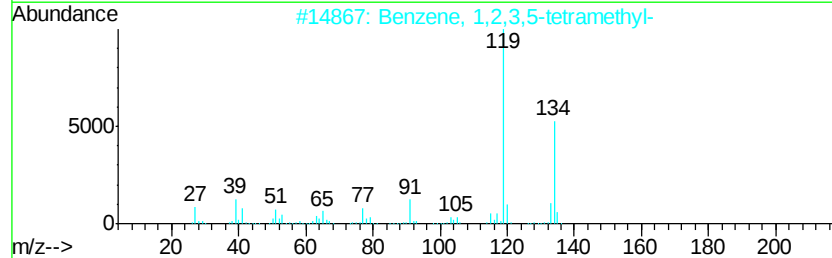
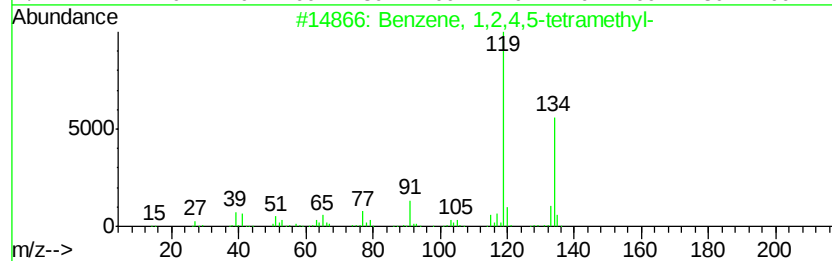
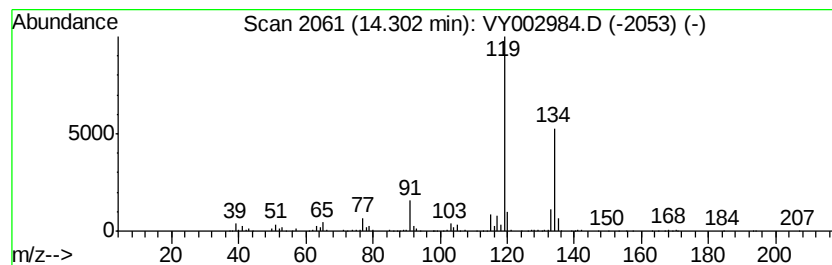
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	16.20 ug/l	456128	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

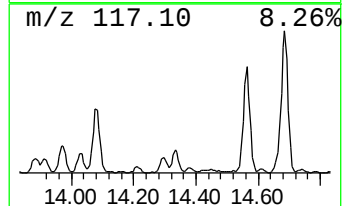
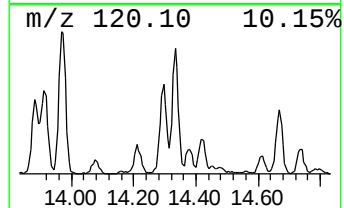
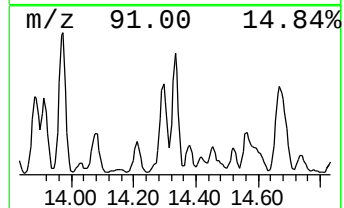
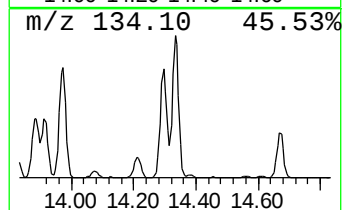
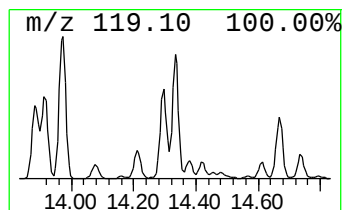
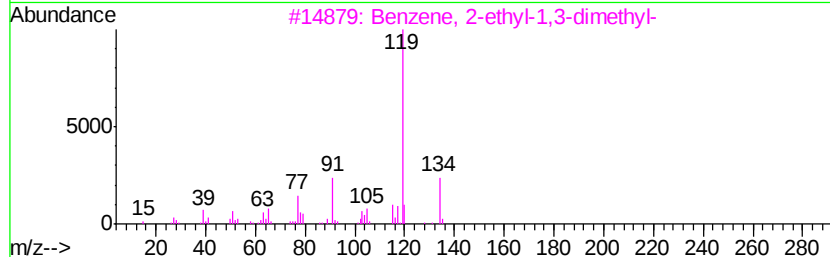
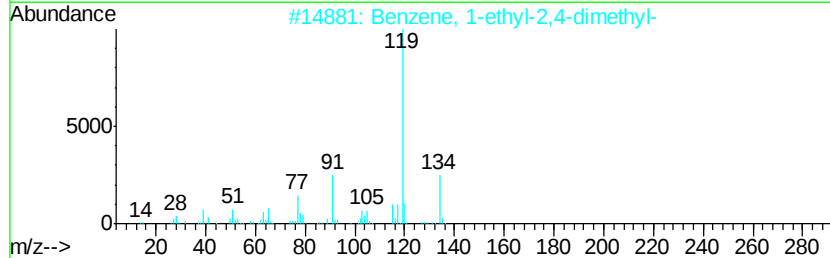
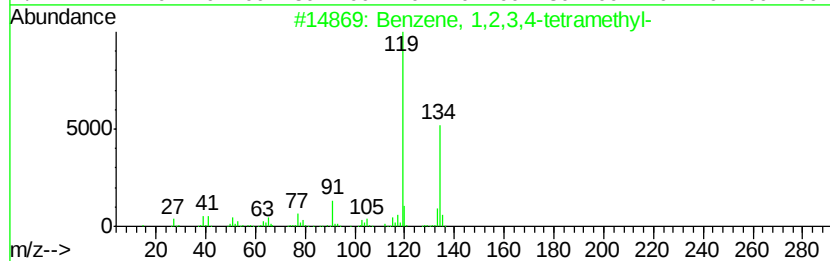
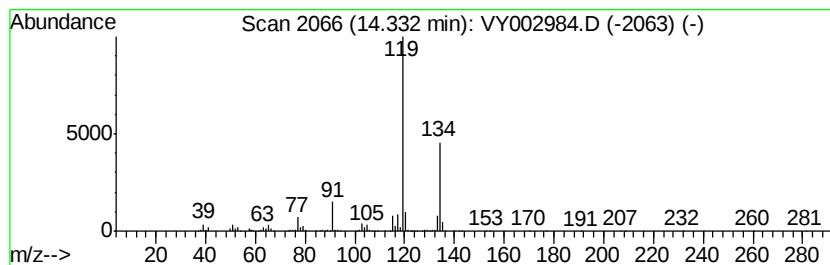
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	18.99 ug/l	534626	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

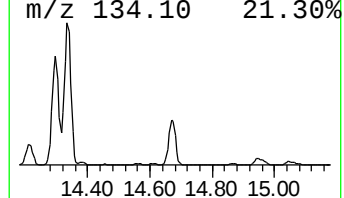
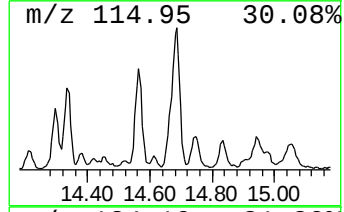
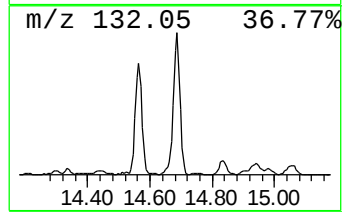
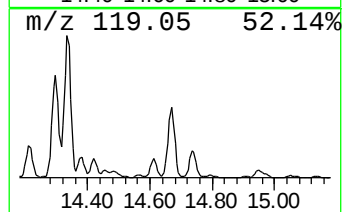
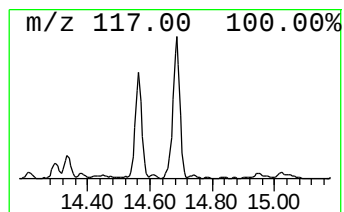
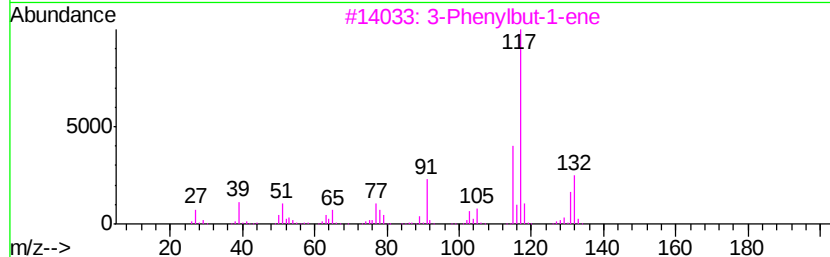
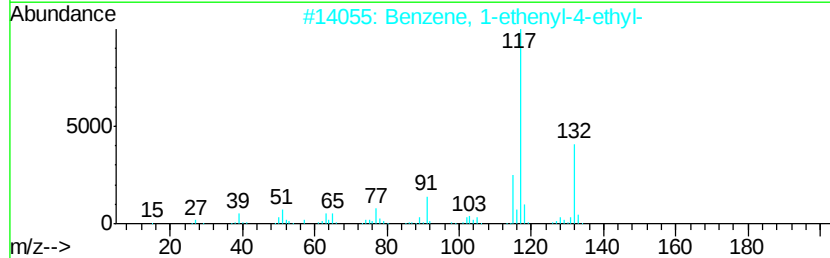
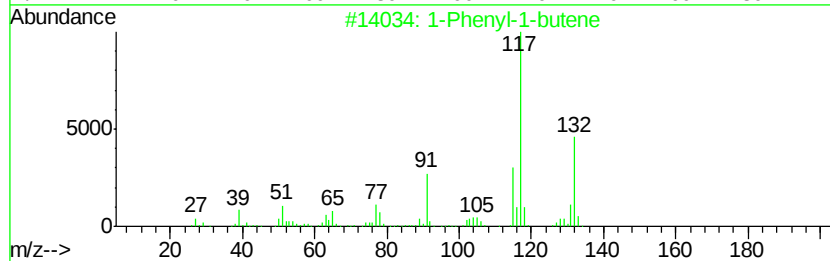
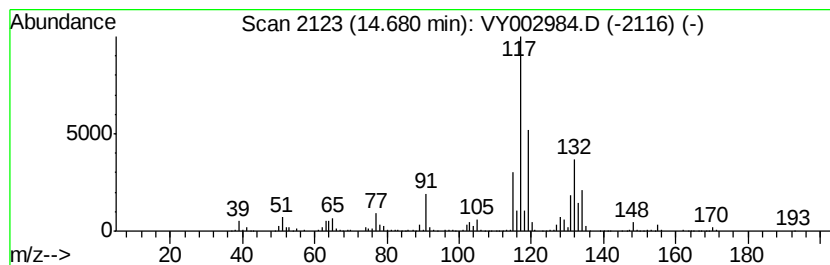
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Quant Title : SW846 8260

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

Peak Number 10 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	25.79 ug/l	726197	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	70
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	62
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_Y\DATA\VY062220\
Data File : VY002984.D
Acq On : 22 Jun 2020 19:21
Operator : SY/MD
Sample : L3082-03
Misc : 3.56G/5ML/MSVOA_Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.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, 2,2,3,3-t...	8.33	21.1	ug/l	421005	2	8.69	1000250	50.0
Pentane, 2,3,4-tr...	9.75	20.6	ug/l	411221	2	8.69	1000250	50.0
Cyclotrisiloxane,...	10.58	16.8	ug/l	404234	3	11.49	1204140	50.0
Benzene, 1-ethyl-...	12.74	66.8	ug/l	1880130	4	13.42	1408020	50.0
Benzene, 1-ethyl-...	12.97	27.8	ug/l	782915	4	13.42	1408020	50.0
Benzene, 1-etheny...	13.63	33.0	ug/l	928821	4	13.42	1408020	50.0
Benzene, 1-ethyl-...	13.97	22.5	ug/l	633448	4	13.42	1408020	50.0
Benzene, 1,2,4,5-...	14.30	16.2	ug/l	456128	4	13.42	1408020	50.0
Benzene, 1,2,3,4-...	14.33	19.0	ug/l	534626	4	13.42	1408020	50.0
1-Phenyl-1-butene	14.68	25.8	ug/l	726197	4	13.42	1408020	50.0