

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
 Data File : VY020034.D
 Acq On : 28 Oct 2024 12:19
 Operator : SY/MD
 Sample : P4578-04
 Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WB-305-TOP-1

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\82Y100924S.M
 Title : SW846 8260

Signal : TIC: VY020034.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.873	343	353	362	rBV2	13876	41343	1.34%	0.110%
2	4.391	428	438	448	rBV	49013	144029	4.67%	0.383%
3	4.628	465	477	486	rBV2	17525	57380	1.86%	0.152%
4	7.640	961	971	977	rBV	185132	445648	14.46%	1.184%
5	7.713	977	983	994	rVB	267843	620236	20.12%	1.648%
6	8.067	1031	1041	1052	rBV	185887	405743	13.16%	1.078%
7	8.616	1122	1131	1142	rBV	486181	971182	31.51%	2.580%
8	9.115	1200	1213	1218	rBV5	21885	57414	1.86%	0.153%
9	9.890	1330	1340	1349	rBV6	16417	47345	1.54%	0.126%
10	10.109	1368	1376	1382	rBV	810445	1411187	45.78%	3.749%
11	10.170	1382	1386	1393	rVB	62901	115237	3.74%	0.306%
12	10.451	1428	1432	1437	rVB	18957	31292	1.02%	0.083%
13	10.542	1437	1447	1454	rBV7	13949	44557	1.45%	0.118%
14	10.731	1473	1478	1484	rVB	34777	55568	1.80%	0.148%
15	10.835	1491	1495	1501	rVB2	19418	37485	1.22%	0.100%
16	10.963	1513	1516	1520	rVB2	24551	32632	1.06%	0.087%
17	11.018	1520	1525	1531	rVB	49623	80149	2.60%	0.213%
18	11.127	1538	1543	1548	rBV3	23956	41601	1.35%	0.111%
19	11.219	1552	1558	1567	rVB6	36937	108034	3.50%	0.287%
20	11.310	1567	1573	1582	rBV	39147	78242	2.54%	0.208%
21	11.414	1582	1590	1603	rBV	736794	1211472	39.30%	3.219%
22	11.615	1616	1623	1624	rBV6	24793	49503	1.61%	0.132%
23	11.664	1628	1631	1635	rVV	38428	64030	2.08%	0.170%
24	11.706	1635	1638	1643	rVB4	27257	41398	1.34%	0.110%
25	11.950	1667	1678	1686	rBV5	74435	248276	8.05%	0.660%
26	12.048	1690	1694	1698	rVB	48274	67323	2.18%	0.179%
27	12.127	1702	1707	1709	rVV3	42759	69840	2.27%	0.186%
28	12.164	1709	1713	1719	rVV3	84395	157581	5.11%	0.419%
29	12.255	1723	1728	1732	rVV	85390	140131	4.55%	0.372%
30	12.298	1732	1735	1747	rVB	39307	120054	3.89%	0.319%
31	12.408	1747	1753	1762	rBV2	559242	954852	30.98%	2.537%
32	12.566	1775	1779	1787	rVB3	54536	116092	3.77%	0.308%
33	12.651	1787	1793	1796	rBV5	43482	85815	2.78%	0.228%
34	12.694	1796	1800	1802	rVV	74681	109505	3.55%	0.291%
35	12.737	1802	1807	1813	rVB2	140769	238740	7.75%	0.634%
36	12.895	1826	1833	1841	rBV	187387	316411	10.27%	0.841%

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37	12.962	1841	1844	1851	rVV2	54361	79887	2.59%	0.212%
38	13.042	1852	1857	1862	rVV	372917	556027	18.04%	1.477%
39	13.090	1862	1865	1871	rVV6	20984	30981	1.01%	0.082%
40	13.170	1871	1878	1882	rVV6	31592	69636	2.26%	0.185%
41	13.237	1883	1889	1894	rVV2	93994	182528	5.92%	0.485%
42	13.292	1894	1898	1901	rVV3	58315	90986	2.95%	0.242%
43	13.346	1901	1907	1910	rVV	594453	951084	30.86%	2.527%
44	13.377	1910	1912	1917	rVB	318273	398892	12.94%	1.060%
45	13.517	1928	1935	1936	rBV	145577	212416	6.89%	0.564%
46	13.548	1936	1940	1943	rVV	537754	812656	26.36%	2.159%
47	13.584	1943	1946	1955	rVB3	546456	1001985	32.51%	2.662%
48	13.676	1955	1961	1965	rBV3	127760	205989	6.68%	0.547%
49	13.743	1965	1972	1977	rBV	158449	263646	8.55%	0.700%
50	13.804	1977	1982	1985	rBV	463429	763041	24.75%	2.027%
51	13.834	1985	1987	1992	rVB	577070	690541	22.40%	1.835%
52	13.889	1992	1996	2002	rVB	1302836	1968530	63.86%	5.230%
53	13.999	2008	2014	2019	rBV2	616054	961247	31.19%	2.554%
54	14.054	2019	2023	2024	rBV	66581	91067	2.95%	0.242%
55	14.133	2030	2036	2040	rBV	427529	682650	22.15%	1.814%
56	14.218	2040	2050	2052	rVV	1044629	1836865	59.59%	4.880%
57	14.255	2052	2056	2060	rVV	1692804	2426610	78.72%	6.447%
58	14.304	2060	2064	2067	rVV2	502680	751564	24.38%	1.997%
59	14.340	2067	2070	2074	rVV	322300	477387	15.49%	1.268%
60	14.401	2077	2080	2083	rVV	80987	109686	3.56%	0.291%
61	14.444	2083	2087	2089	rVV	205232	271017	8.79%	0.720%
62	14.480	2089	2093	2098	rVV	857944	1389999	45.09%	3.693%
63	14.529	2098	2101	2106	rVB	723602	966094	31.34%	2.567%
64	14.596	2106	2112	2118	rBV3	1452966	3082391	100.00%	8.189%
65	14.657	2118	2122	2128	rVB3	412368	734382	23.83%	1.951%
66	14.749	2133	2137	2142	rVB	161871	240219	7.79%	0.638%
67	14.858	2149	2155	2163	rBV2	299676	606012	19.66%	1.610%
68	14.962	2167	2172	2177	rVV4	173101	336637	10.92%	0.894%
69	15.011	2177	2180	2183	rVV2	69340	101403	3.29%	0.269%
70	15.053	2183	2187	2193	rVB4	158148	252102	8.18%	0.670%
71	15.145	2193	2202	2207	rBV2	572518	1282122	41.60%	3.406%
72	15.255	2215	2220	2225	rBV5	77705	163525	5.31%	0.434%
73	15.346	2230	2235	2241	rVB	616168	883536	28.66%	2.347%
74	15.450	2245	2252	2256	rBV5	61815	154945	5.03%	0.412%
75	15.858	2316	2319	2323	rVB4	40965	60043	1.95%	0.160%
76	15.913	2324	2328	2335	rVV5	75618	174816	5.67%	0.464%

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77	15.986	2335	2340	2346	rVB7	78842	167004	5.42%	0.444%
78	16.053	2346	2351	2357	rVB3	78880	137995	4.48%	0.367%
79	16.181	2366	2372	2378	rVB	416372	774786	25.14%	2.058%
80	16.248	2378	2383	2389	rBV	234676	386642	12.54%	1.027%
81	16.370	2396	2403	2417	rVB	456840	1040176	33.75%	2.764%

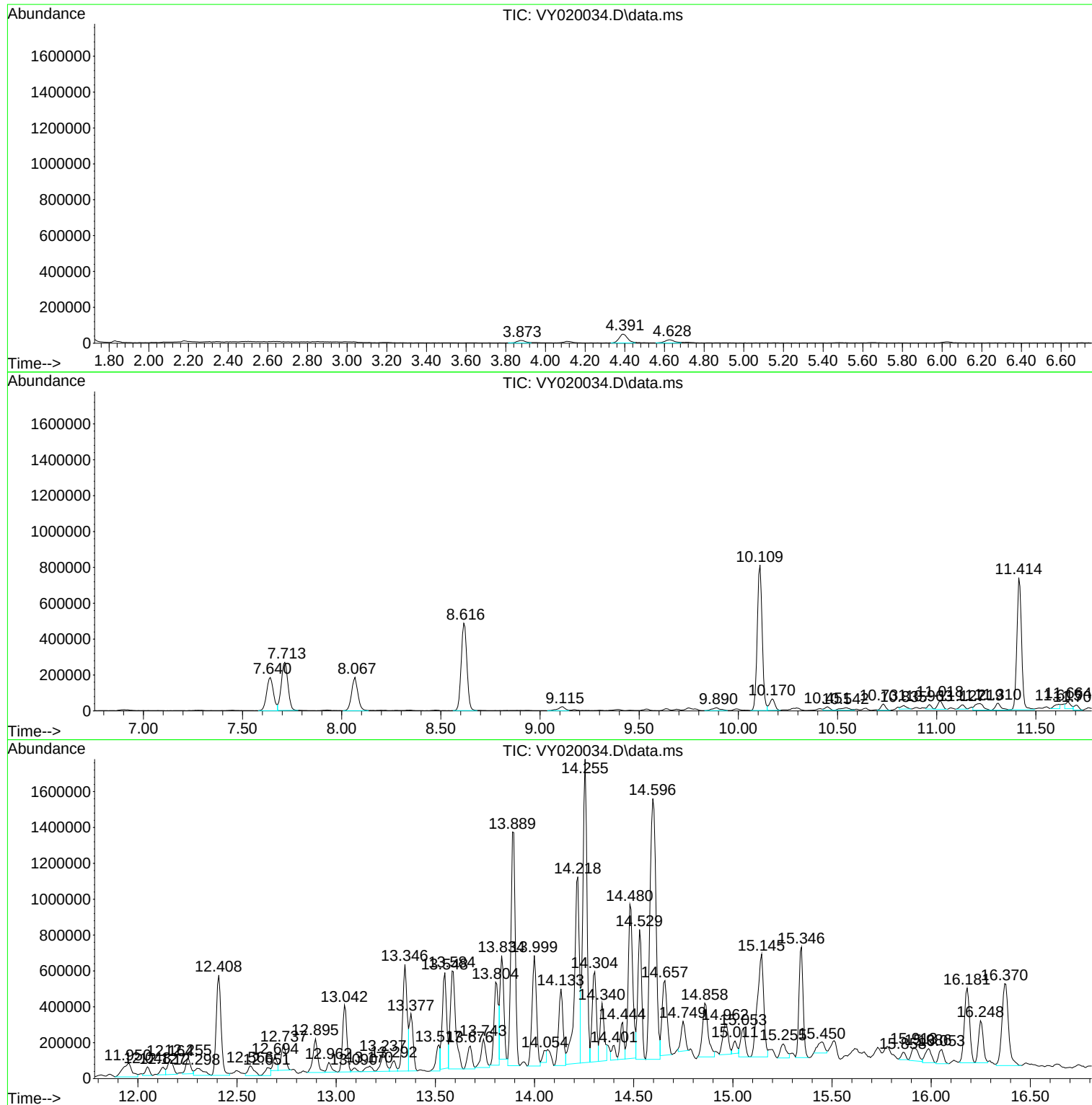
Sum of corrected areas: 37639074

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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



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Instrument :
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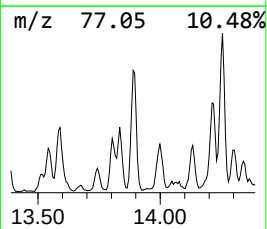
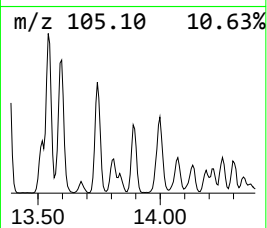
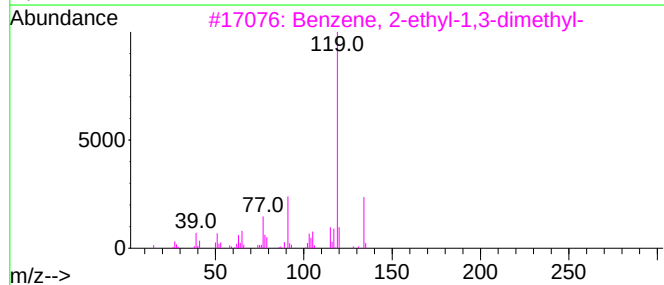
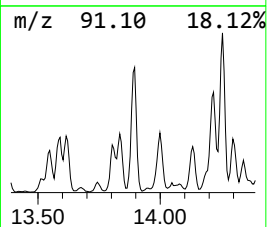
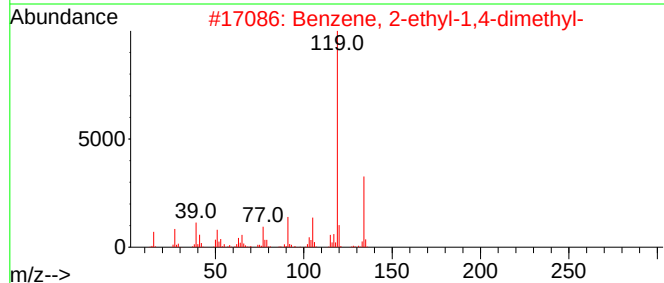
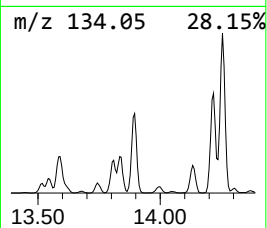
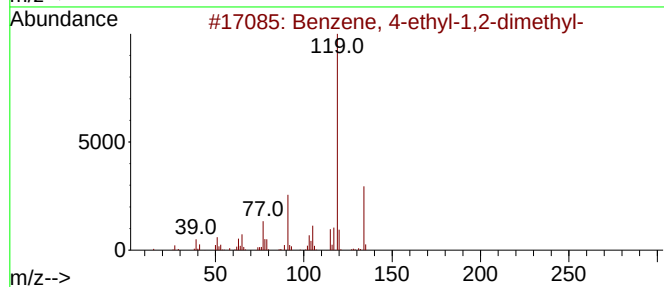
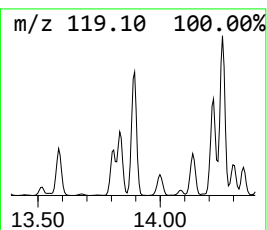
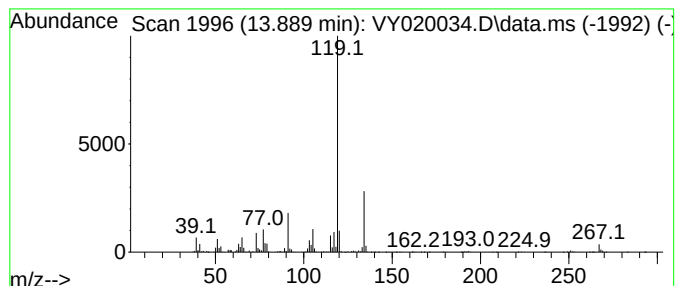
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.889	103.49 ug/l	1968530	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



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Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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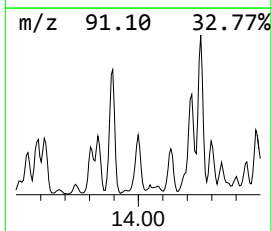
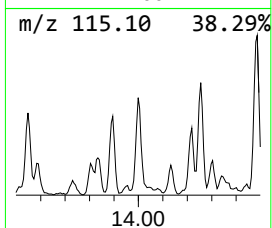
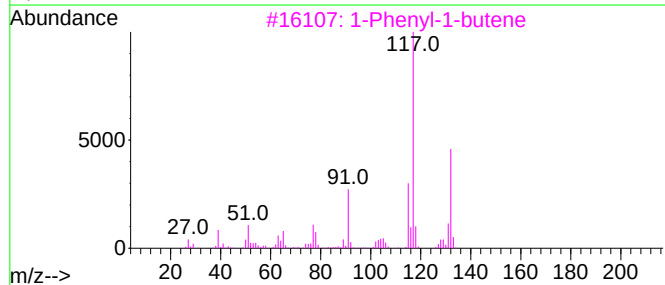
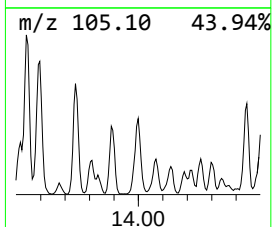
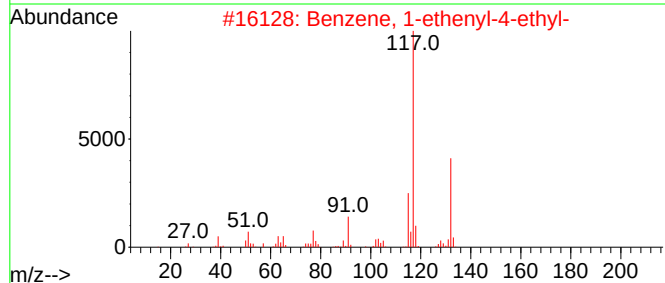
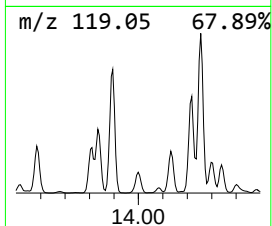
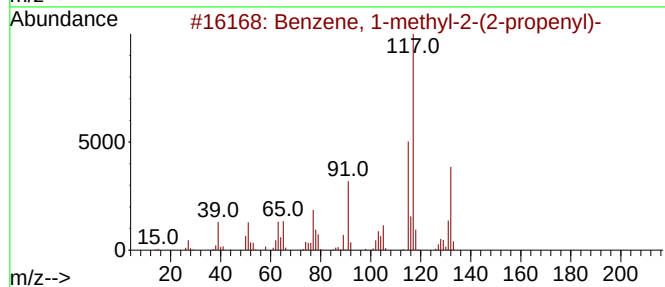
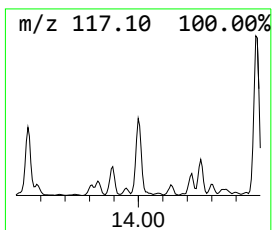
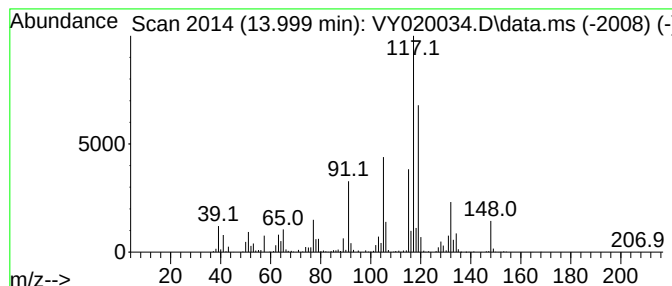
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-methyl-2-(2-prop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.999	50.53 ug/l	961247	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	70
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70



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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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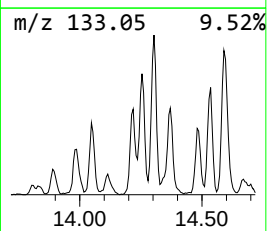
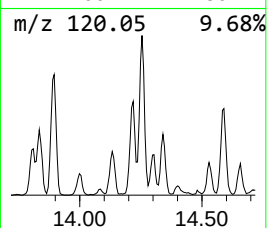
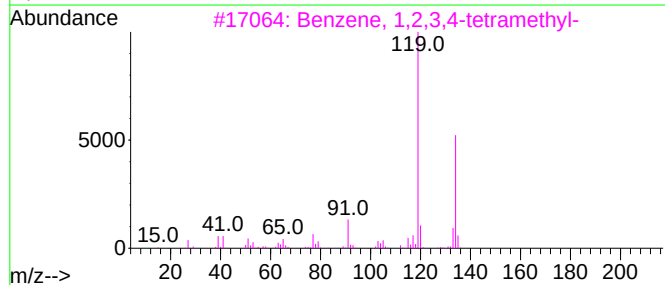
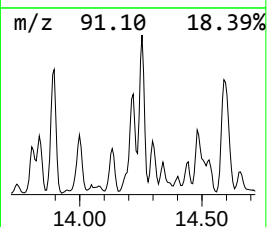
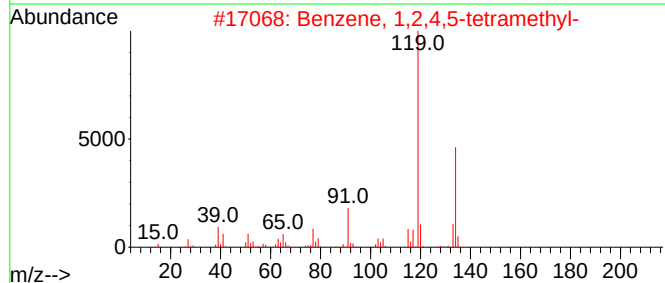
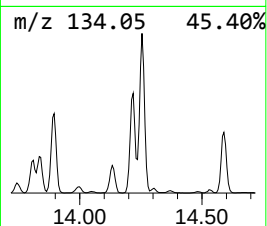
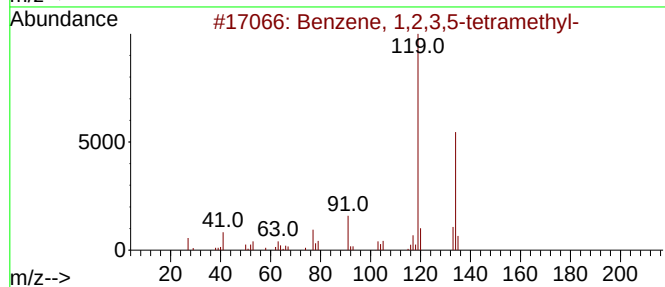
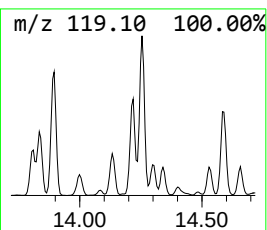
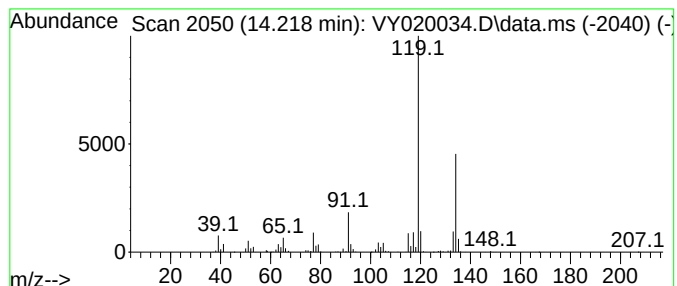
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.218	96.57 ug/l	1836870	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5			o-Cymene	134	C10H14	000527-84-4	93



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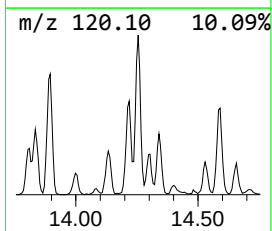
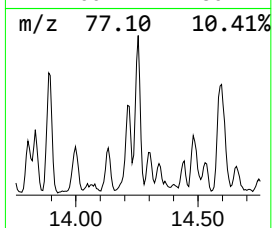
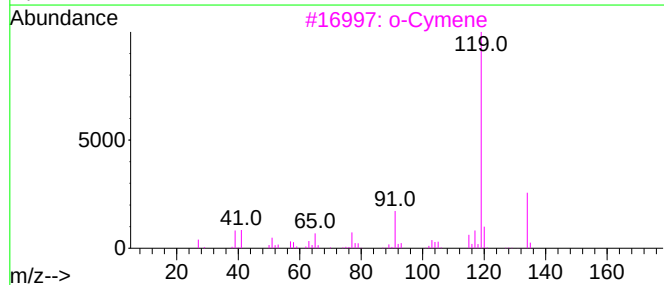
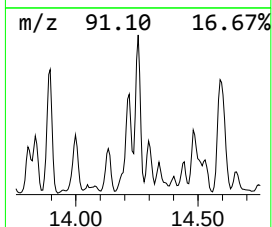
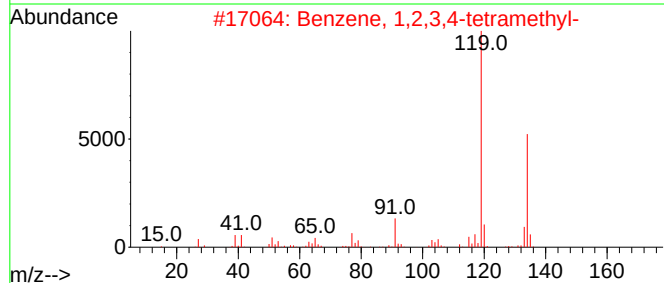
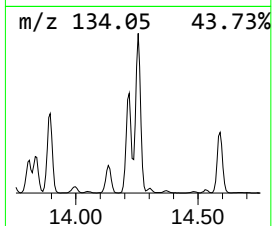
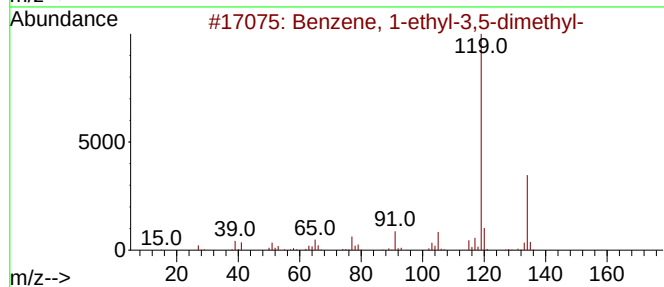
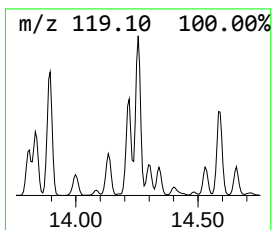
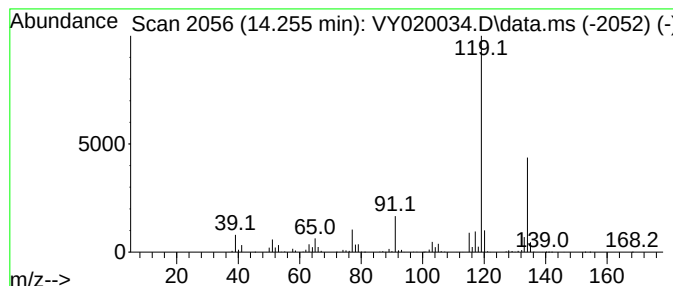
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.255	127.57 ug/l	2426610	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3			o-Cymene	134	C10H14	000527-84-4	94
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020034.D
Acq On : 28 Oct 2024 12:19
Operator : SY/MD
Sample : P4578-04
Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP-1

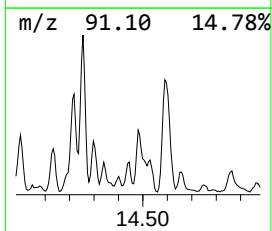
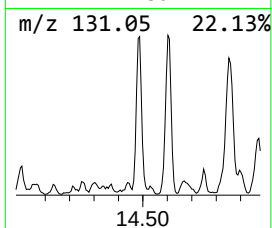
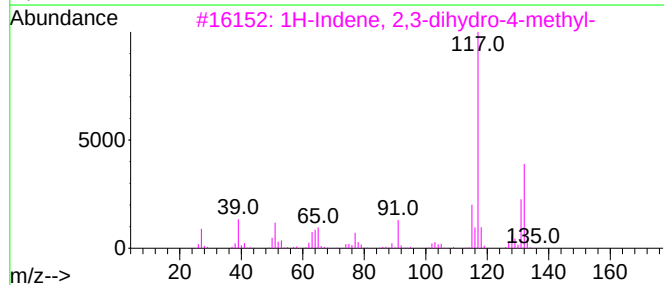
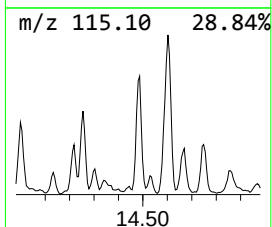
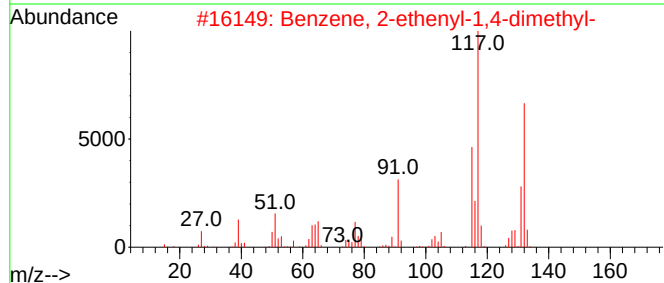
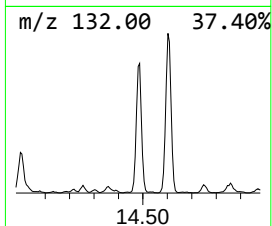
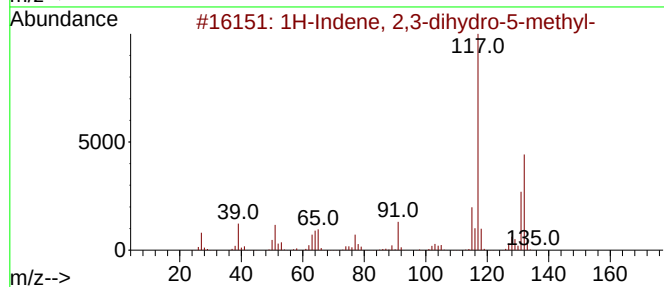
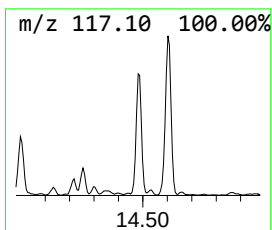
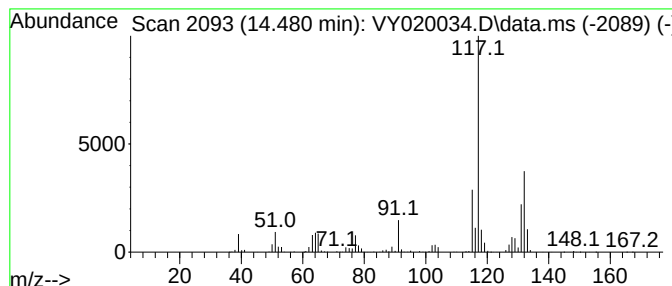
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.480	73.07 ug/l	1390000	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020034.D
Acq On : 28 Oct 2024 12:19
Operator : SY/MD
Sample : P4578-04
Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP-1

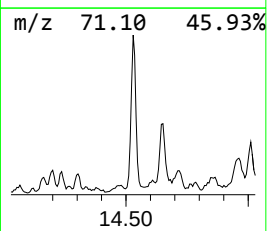
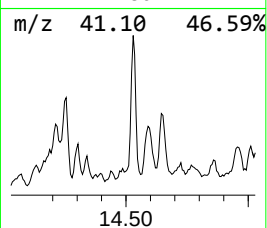
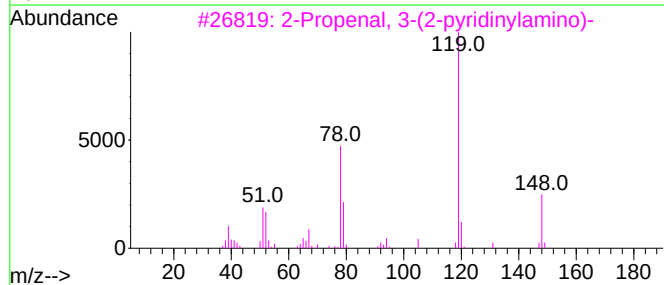
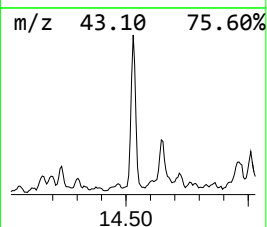
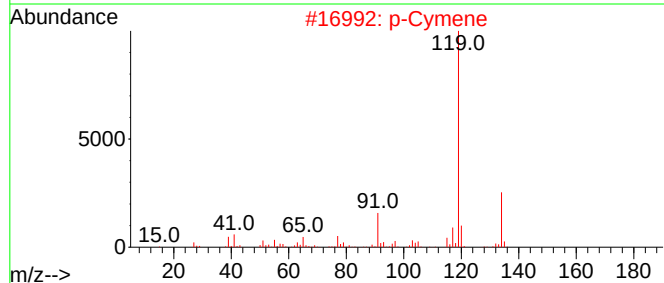
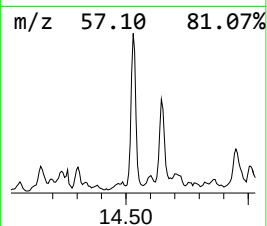
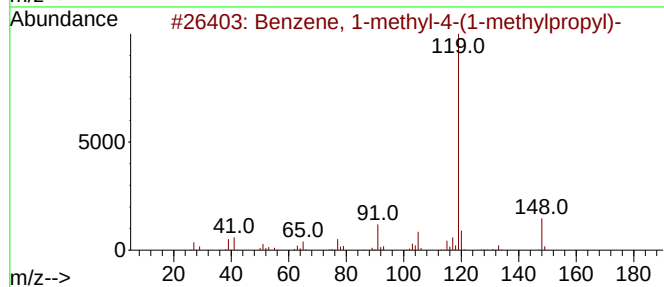
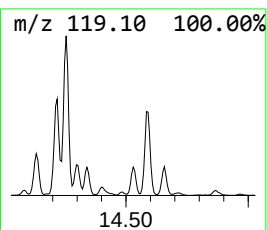
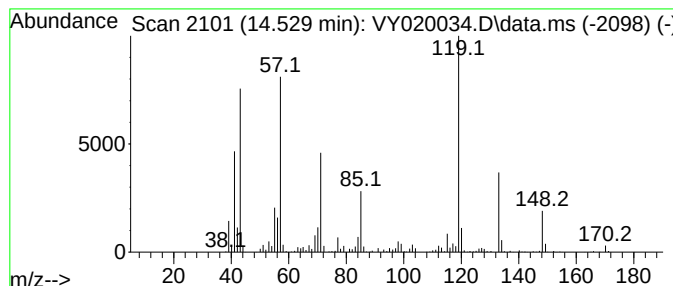
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Quant Title : SW846 8260

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

Peak Number 6 unknown14.529 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.529	50.79 ug/l	966094	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	43
2			p-Cymene	134	C10H14	000099-87-6	42
3			2-Propenal, 3-(2-pyridinylamino)-	148	C8H8N2O	068970-82-1	38
4			Dodecane	170	C12H26	000112-40-3	38
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020034.D
Acq On : 28 Oct 2024 12:19
Operator : SY/MD
Sample : P4578-04
Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP-1

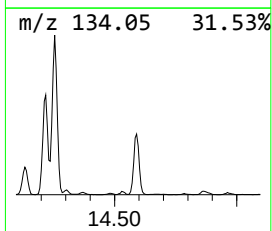
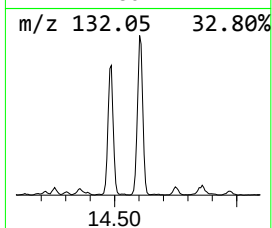
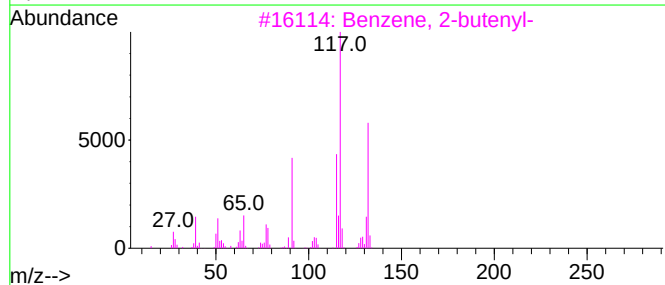
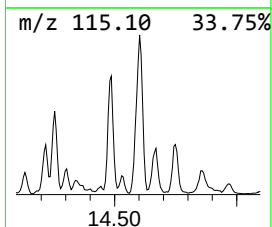
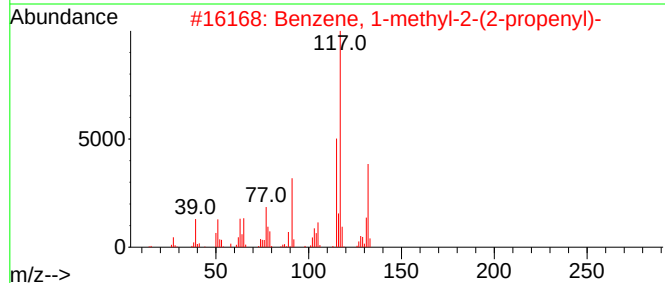
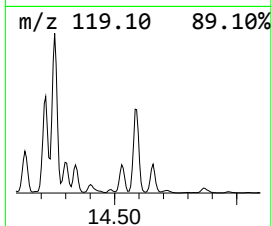
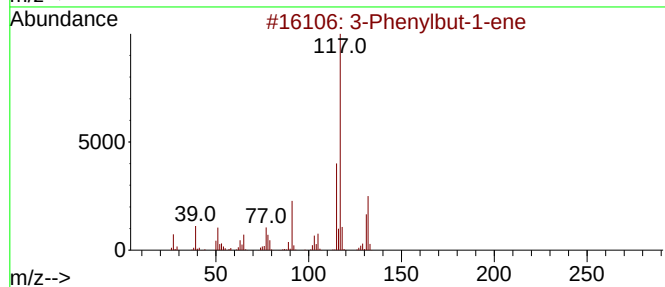
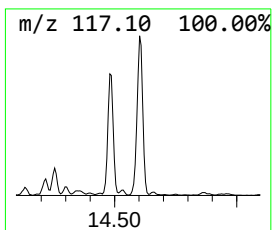
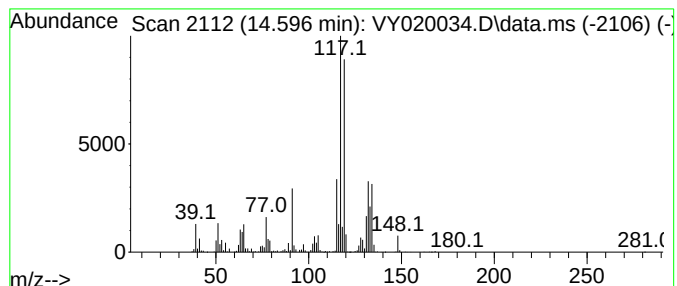
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Quant Title : SW846 8260

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

Peak Number 7 3-Phenylbut-1-ene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.596	162.05 ug/l	3082390	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020034.D
Acq On : 28 Oct 2024 12:19
Operator : SY/MD
Sample : P4578-04
Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

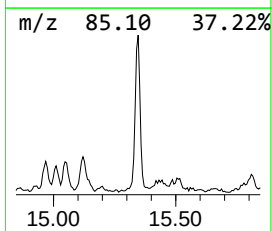
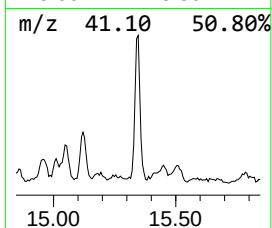
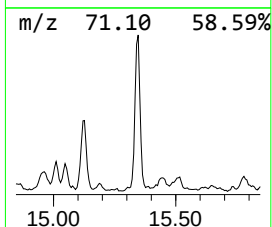
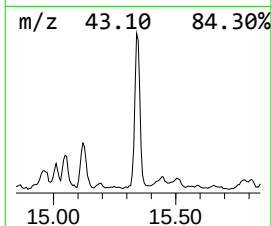
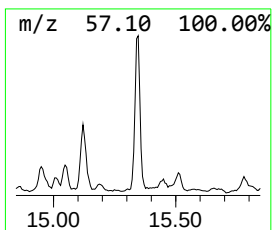
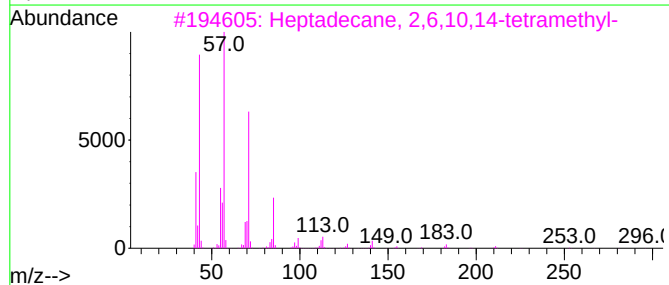
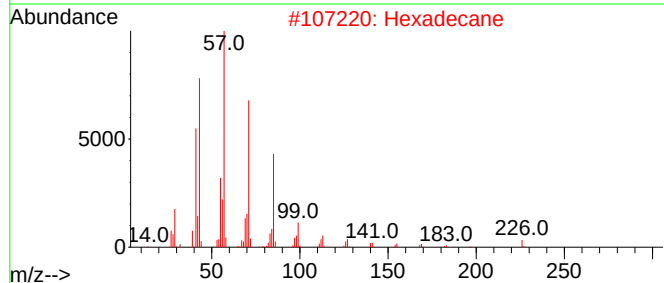
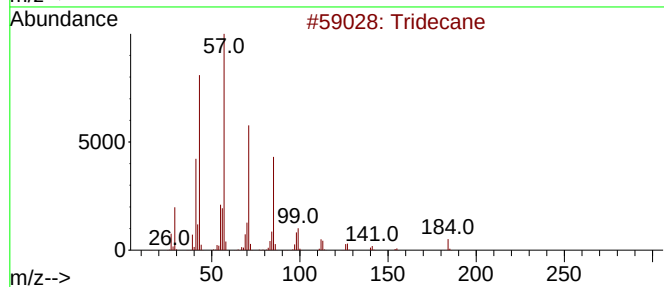
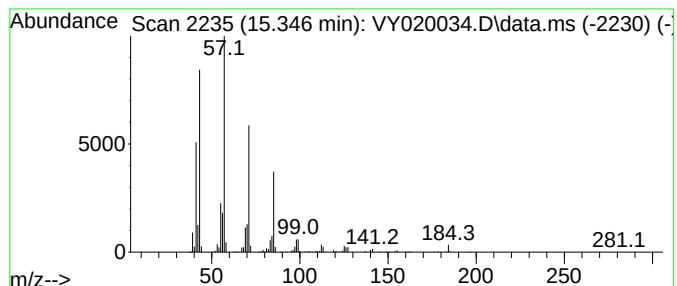
Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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

Peak Number 9 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.346	46.45 ug/l	883536	1,4-Dichlorobenzene-d4			13.346
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	95
2	Hexadecane		226	C16H34	000544-76-3	91
3	Heptadecane, 2,6,10,14-tetramethyl-		296	C21H44	018344-37-1	86
4	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	80
5	1-Iodo-2-methylnonane		268	C10H21I	1000101-47-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020034.D
Acq On : 28 Oct 2024 12:19
Operator : SY/MD
Sample : P4578-04
Misc : 7.19g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 4-ethy...	13.889	103.5	ug/l	1968530	4	13.346	951084	50.0
Benzene, 1-meth...	13.999	50.5	ug/l	961247	4	13.346	951084	50.0
Benzene, 1,2,3,...	14.218	96.6	ug/l	1836870	4	13.346	951084	50.0
Benzene, 1-ethy...	14.255	127.6	ug/l	2426610	4	13.346	951084	50.0
1H-Indene, 2,3-...	14.480	73.1	ug/l	1390000	4	13.346	951084	50.0
unknown14.529	14.529	50.8	ug/l	966094	4	13.346	951084	50.0
3-Phenylbut-1-ene	14.596	162.1	ug/l	3082390	4	13.346	951084	50.0
Tridecane	15.346	46.5	ug/l	883536	4	13.346	951084	50.0