

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
 Data File : VY020696.D
 Acq On : 23 Dec 2024 17:04
 Operator : SY/MD
 Sample : P5342-01
 Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 CHRT26634

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\82Y122324S.M

Title : SW846 8260

Signal : TIC: VY020696.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.178	71	75	81	rVB2	1522	2201	3.23%	0.216%
2	3.872	346	353	358	rBV2	981	2256	3.31%	0.222%
3	4.122	387	394	399	rBV3	612	1424	2.09%	0.140%
4	6.890	837	848	859	rBV	5463	16117	23.67%	1.583%
5	7.433	934	937	941	rBV2	506	797	1.17%	0.078%
6	7.646	964	972	977	rBV3	10615	25783	37.87%	2.533%
7	7.719	977	984	993	rVB2	23072	51914	76.26%	5.100%
8	8.067	1030	1041	1050	rBV4	9638	27547	40.46%	2.706%
9	8.622	1125	1132	1141	rBV2	27601	54287	79.74%	5.334%
10	10.115	1370	1377	1383	rBV	38979	68077	100.00%	6.688%
11	10.176	1383	1387	1395	rVB	9401	15589	22.90%	1.532%
12	10.310	1405	1409	1411	rBV2	730	739	1.09%	0.073%
13	10.518	1437	1443	1450	rBV2	2094	4358	6.40%	0.428%
14	10.890	1497	1504	1508	rVB3	1070	1875	2.75%	0.184%
15	10.975	1511	1518	1522	rVV2	2742	4642	6.82%	0.456%
16	11.030	1522	1527	1532	rVB4	1211	1797	2.64%	0.177%
17	11.219	1548	1558	1560	rBV4	1672	3031	4.45%	0.298%
18	11.255	1560	1564	1570	rVV2	2590	4302	6.32%	0.423%
19	11.316	1570	1574	1578	rVB2	1303	1688	2.48%	0.166%
20	11.426	1584	1592	1598	rBV	27479	45423	66.72%	4.463%
21	11.524	1603	1608	1616	rVB	4581	8008	11.76%	0.787%
22	11.633	1616	1626	1636	rBV3	13266	31338	46.03%	3.079%
23	11.713	1636	1639	1645	rVB5	1309	1956	2.87%	0.192%
24	11.963	1669	1680	1684	rBV3	8359	18505	27.18%	1.818%
25	12.054	1692	1695	1698	rBV4	2286	3110	4.57%	0.306%
26	12.097	1698	1702	1704	rVV3	1665	1898	2.79%	0.186%
27	12.133	1704	1708	1711	rVV6	2295	3032	4.45%	0.298%
28	12.176	1711	1715	1723	rVB6	4951	8683	12.75%	0.853%
29	12.267	1723	1730	1735	rBV	35523	55949	82.18%	5.497%
30	12.347	1741	1743	1749	rVB5	2258	3647	5.36%	0.358%
31	12.414	1749	1754	1760	rBV4	19677	35231	51.75%	3.461%
32	12.505	1764	1769	1773	rBV7	1848	3573	5.25%	0.351%
33	12.572	1776	1780	1788	rVB6	4766	11189	16.44%	1.099%
34	12.664	1788	1795	1799	rBV3	8112	17276	25.38%	1.697%
35	12.694	1799	1800	1801	rVV	1851	1263	1.86%	0.124%
36	12.743	1802	1808	1813	rVB4	15281	26888	39.50%	2.642%

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37	12.786	1813	1815	1822	rVB5	3373	5350	7.86%	0.526%
38	12.877	1825	1830	1831	rBV4	2964	2887	4.24%	0.284%
39	12.944	1838	1841	1842	rBV3	1007	1057	1.55%	0.104%
40	12.968	1842	1845	1853	rVB6	7863	14179	20.83%	1.393%
41	13.048	1853	1858	1863	rBV	15862	26412	38.80%	2.595%
42	13.109	1863	1868	1872	rBV7	1966	4006	5.88%	0.394%
43	13.182	1872	1880	1884	rBV	8896	13500	19.83%	1.326%
44	13.249	1884	1891	1896	rBV8	9020	21813	32.04%	2.143%
45	13.353	1904	1908	1912	rBV	8004	13553	19.91%	1.332%
46	13.432	1918	1921	1924	rVB5	1656	2027	2.98%	0.199%
47	13.493	1924	1931	1936	rBV9	3207	8460	12.43%	0.831%
48	13.554	1936	1941	1945	rVV	26935	40908	60.09%	4.019%
49	13.602	1945	1949	1955	rVB5	5511	10262	15.07%	1.008%
50	13.676	1955	1961	1966	rBV5	14618	26746	39.29%	2.628%
51	13.737	1966	1971	1976	rBV4	8714	16152	23.73%	1.587%
52	13.804	1981	1982	1985	rVV3	2833	3502	5.14%	0.344%
53	13.846	1986	1989	1993	rVV5	5126	6801	9.99%	0.668%
54	13.895	1994	1997	2002	rVB3	6933	11057	16.24%	1.086%
55	13.944	2002	2005	2010	rVB7	2332	3944	5.79%	0.387%
56	14.011	2010	2016	2020	rBV	15197	26178	38.45%	2.572%
57	14.060	2020	2024	2025	rBV3	3554	4168	6.12%	0.410%
58	14.121	2032	2034	2037	rBV4	1261	1519	2.23%	0.149%
59	14.182	2039	2044	2061	rVB8	10199	31463	46.22%	3.091%
60	14.310	2061	2065	2068	rBV5	4184	6753	9.92%	0.663%
61	14.346	2068	2071	2075	rVB3	7591	10439	15.33%	1.026%
62	14.401	2075	2080	2089	rVB9	3938	7170	10.53%	0.704%
63	14.493	2093	2095	2098	rBV4	1448	1871	2.75%	0.184%
64	14.541	2100	2103	2107	rVB4	2965	4211	6.19%	0.414%
65	14.608	2107	2114	2119	rBV6	7424	17620	25.88%	1.731%
66	14.657	2119	2122	2129	rVB7	6185	11089	16.29%	1.089%
67	14.712	2129	2131	2132	rBV2	1367	1053	1.55%	0.103%
68	14.761	2133	2139	2143	rBV3	8582	14459	21.24%	1.421%
69	14.803	2143	2146	2153	rVB9	4190	6163	9.05%	0.606%
70	14.864	2153	2156	2161	rVB5	4153	6845	10.05%	0.673%
71	14.925	2161	2166	2167	rBV4	2281	3720	5.46%	0.365%
72	15.005	2177	2179	2180	rBV	1103	765	1.12%	0.075%
73	15.102	2192	2195	2196	rBV3	748	946	1.39%	0.093%
74	15.120	2196	2198	2199	rBV2	1068	882	1.30%	0.087%
75	15.151	2199	2203	2208	rBV	10051	18451	27.10%	1.813%
76	15.291	2224	2226	2231	rVB6	4239	7337	10.78%	0.721%

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77	15.352	2233	2236	2240	rBV6	2061	3273	4.81%	0.322%
78	15.407	2242	2245	2250	rBV7	1177	2355	3.46%	0.231%
79	15.468	2252	2255	2259	rVV6	1391	2385	3.50%	0.234%
80	15.511	2259	2262	2269	rVB8	3753	8217	12.07%	0.807%
81	15.596	2274	2276	2278	rBV3	720	782	1.15%	0.077%
82	15.651	2279	2285	2289	rVV9	2160	3948	5.80%	0.388%
83	15.846	2315	2317	2320	rVB4	1172	1097	1.61%	0.108%
84	16.053	2348	2351	2352	rBV3	662	685	1.01%	0.067%
85	16.181	2368	2372	2374	rBV5	1840	2480	3.64%	0.244%
86	16.376	2399	2404	2406	rBV5	2307	3837	5.64%	0.377%
87	16.437	2412	2414	2418	rBV5	924	1022	1.50%	0.100%
88	16.510	2424	2426	2428	rVB3	1262	1014	1.49%	0.100%
89	16.535	2428	2430	2432	rBV3	946	917	1.35%	0.090%
90	16.651	2441	2449	2452	rBV8	1840	4702	6.91%	0.462%

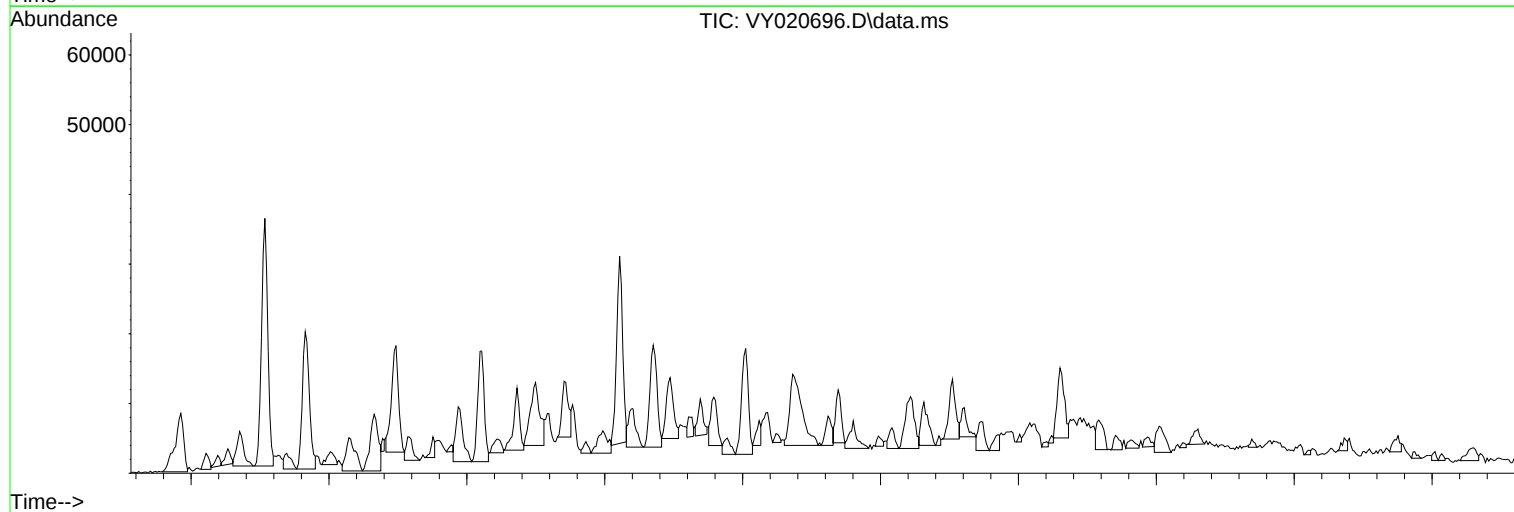
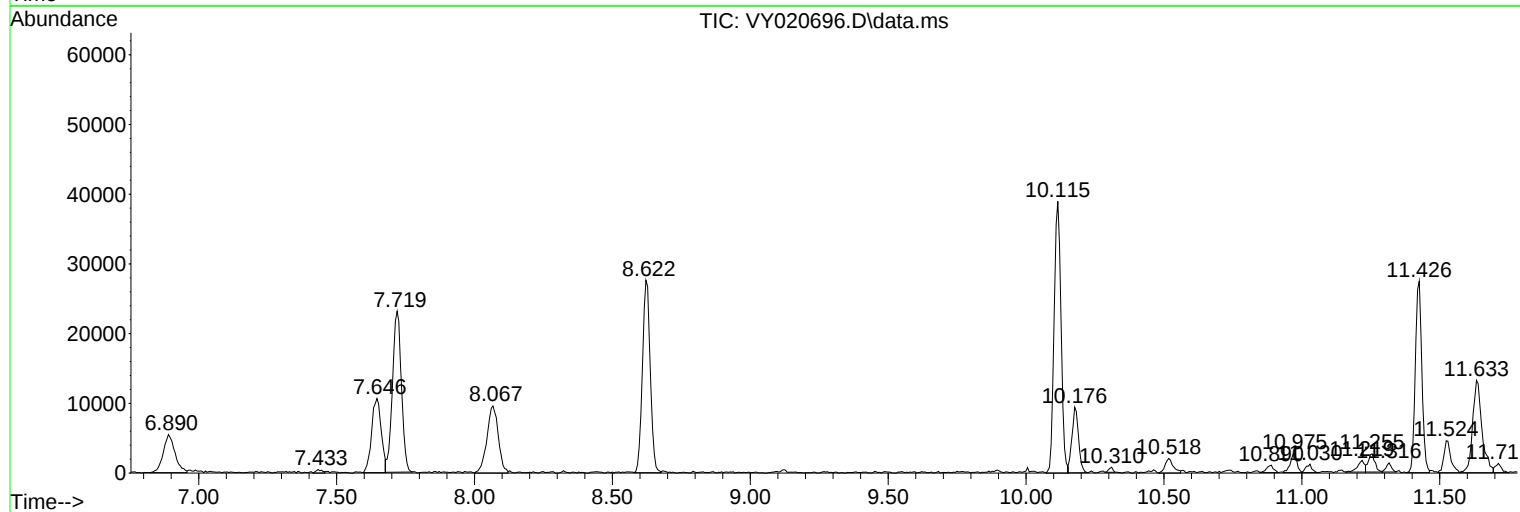
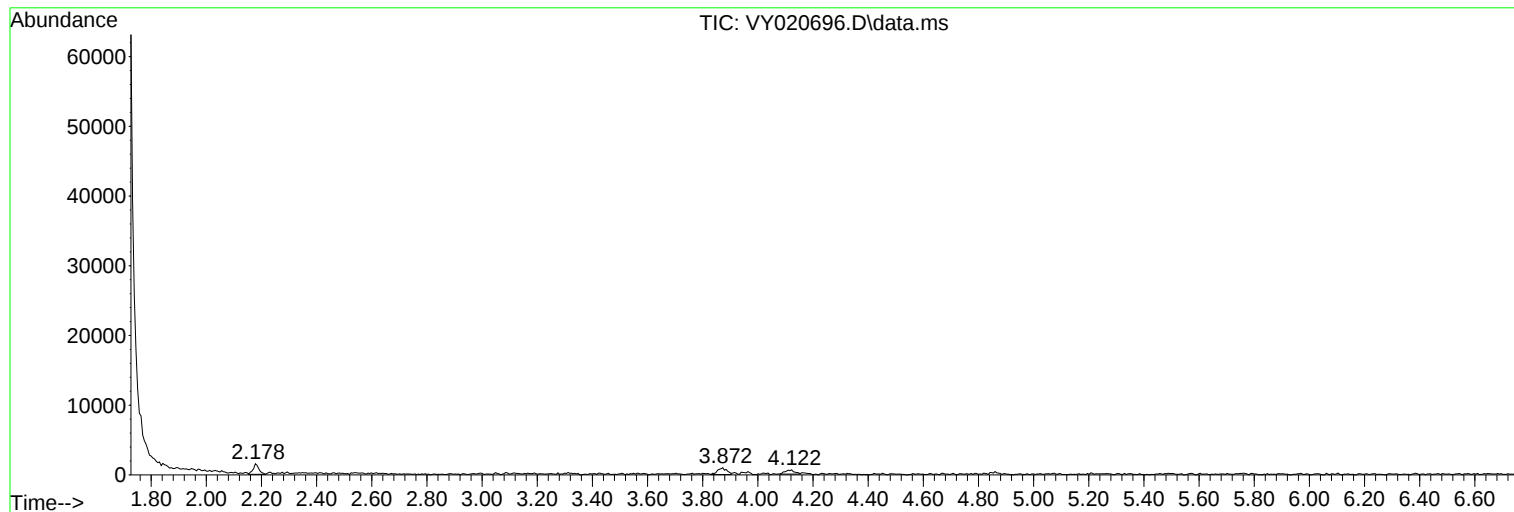
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y122324S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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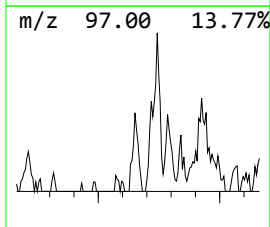
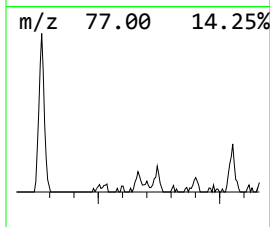
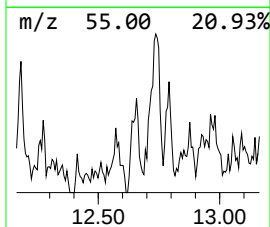
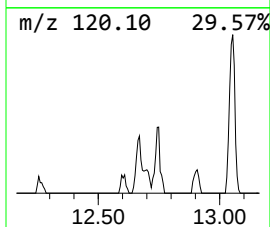
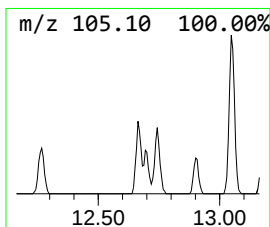
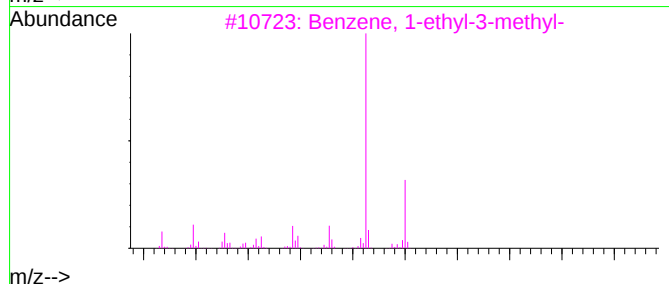
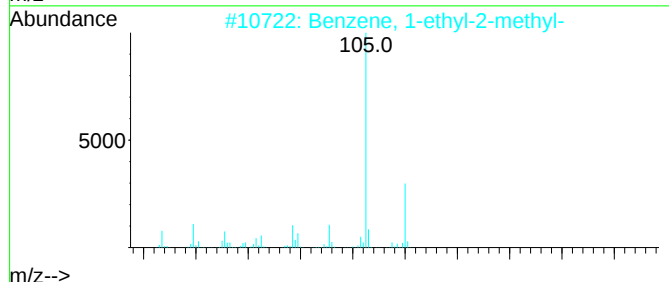
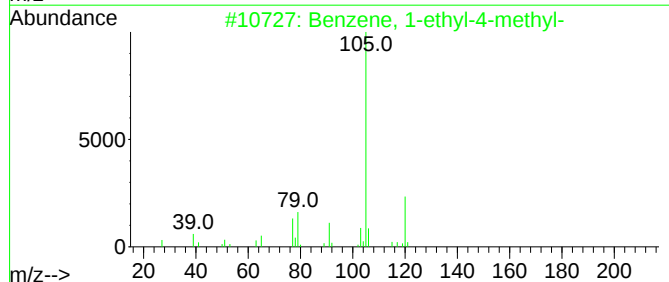
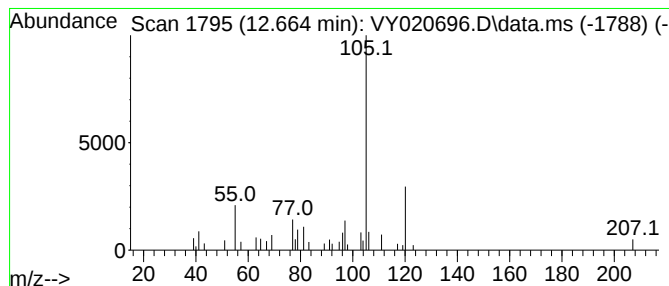
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.664	63.73 ug/l	17276	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	64
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	64
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	52
5	Mesitylene	120	C9H12	000108-67-8	52



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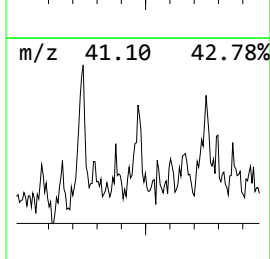
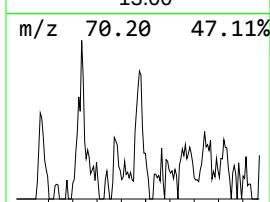
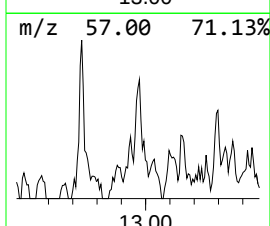
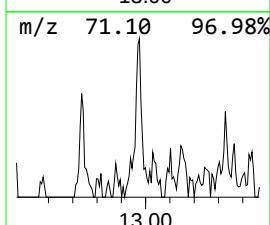
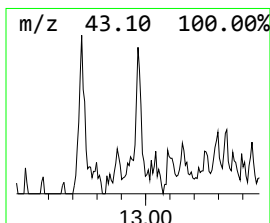
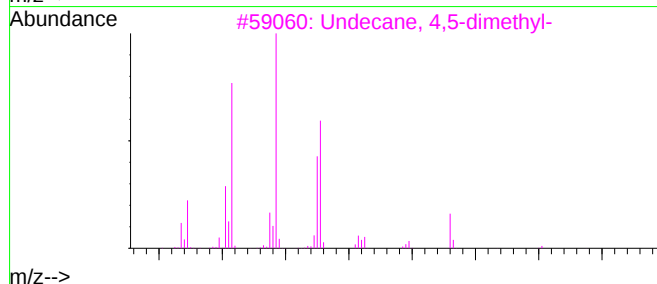
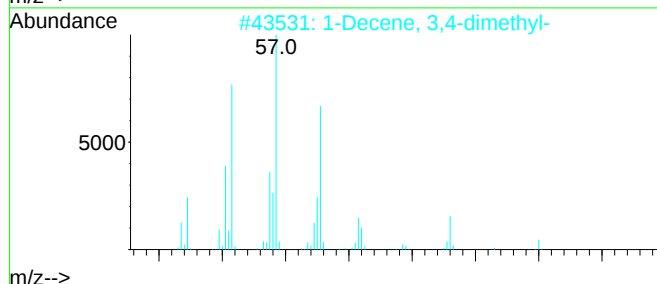
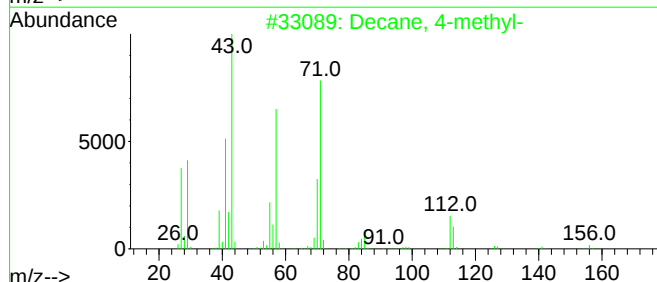
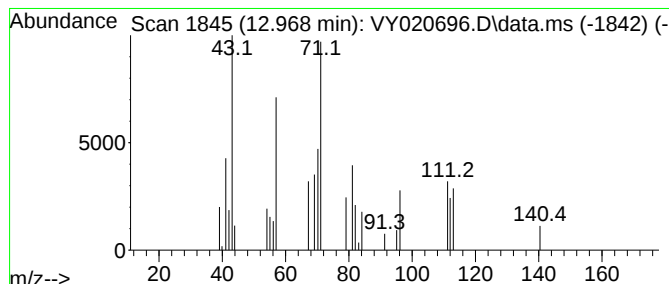
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y122324S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown12.968 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.968	52.31 ug/l	14179	1,4-Dichlorobenzene-d4	13.359

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	47
2			1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	40
3			Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	38
4			Sulfurous acid, octyl 2-pentyl e...	264	C13H28O3S	1000309-15-7	32
5			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	25



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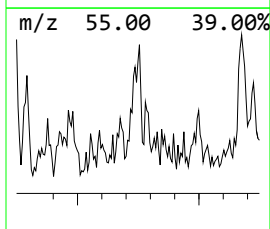
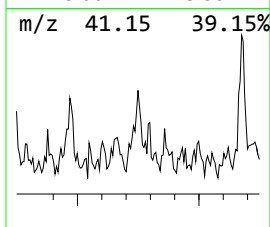
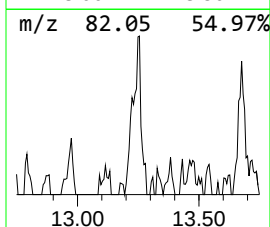
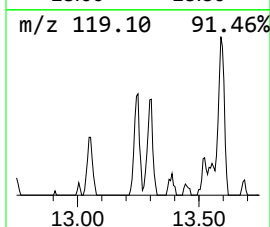
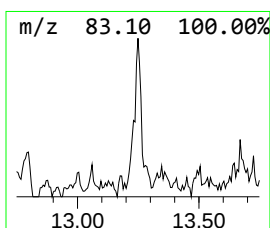
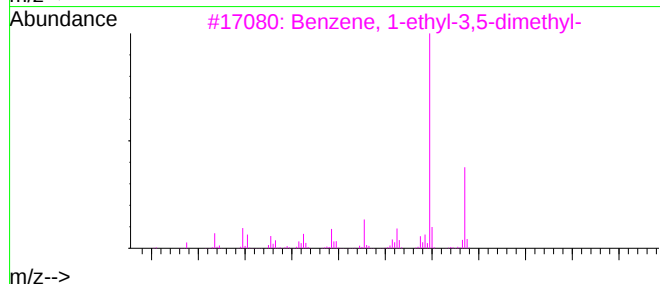
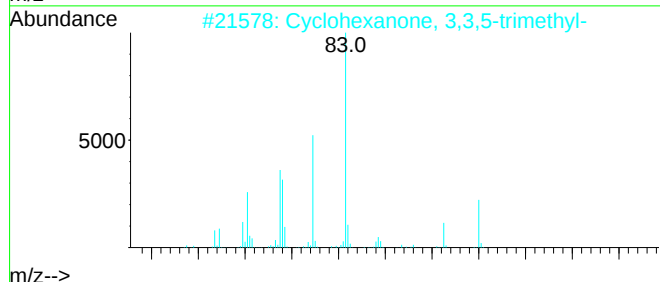
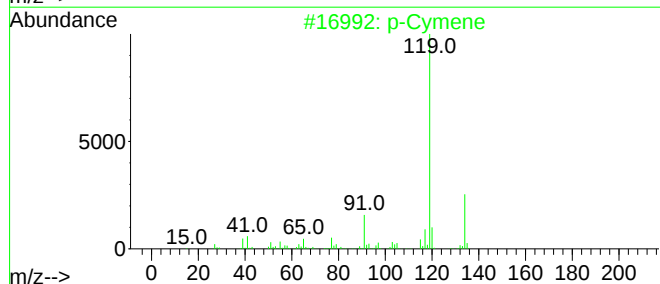
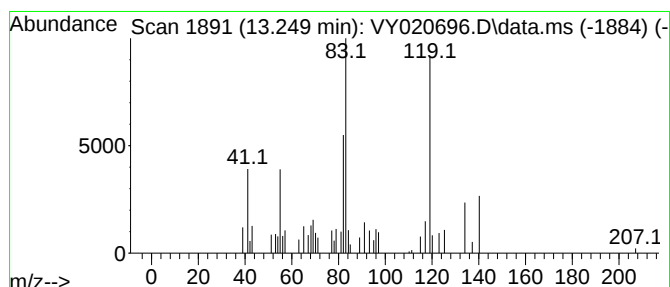
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y122324S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown13.249 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.249	80.47 ug/l	21813	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	30
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	30
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	25
4	o-Cymene	134	C10H14	000527-84-4	25
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CHRT26634

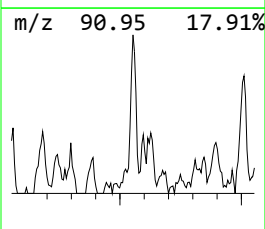
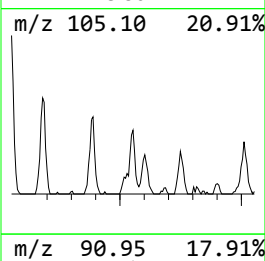
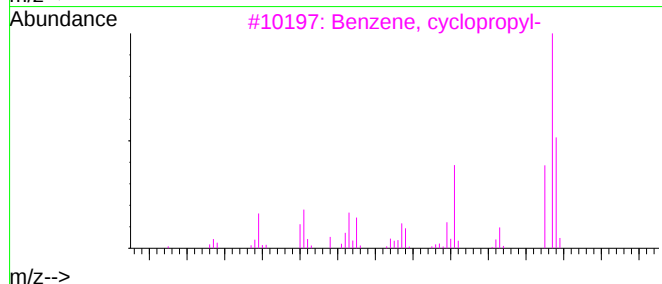
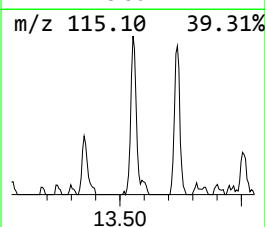
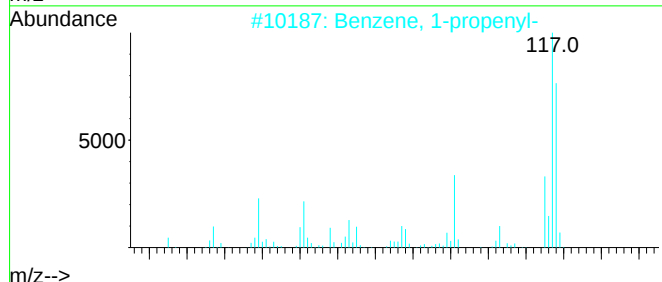
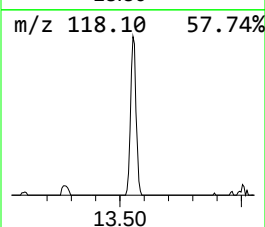
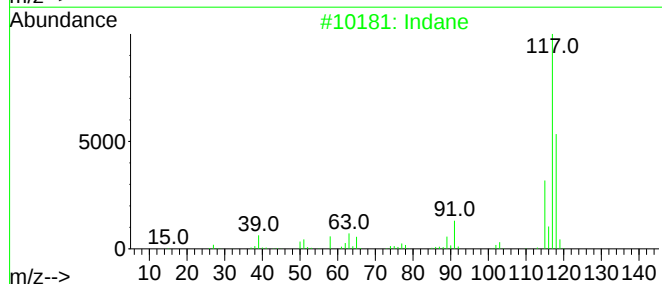
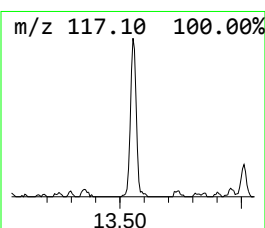
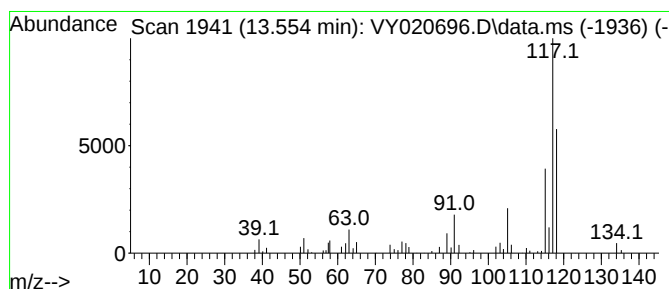
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.554	150.92 ug/l	40908	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	58
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	53
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	52
5	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CHRT26634

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y122324S.M
Quant Title : SW846 8260

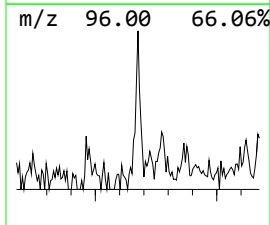
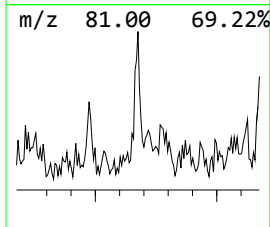
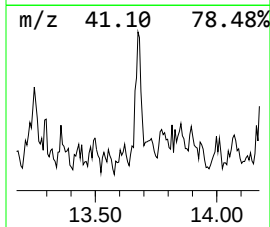
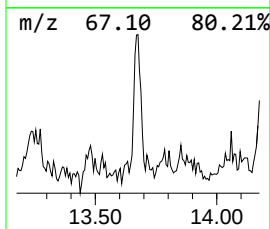
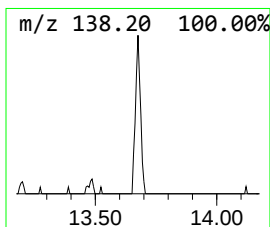
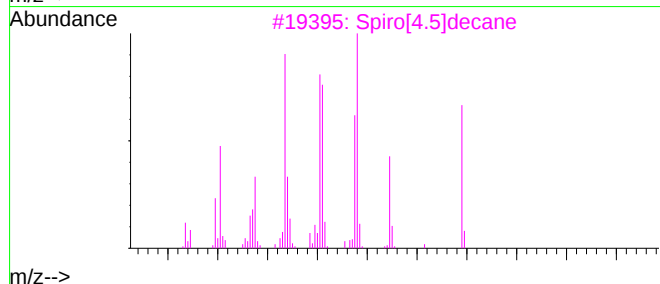
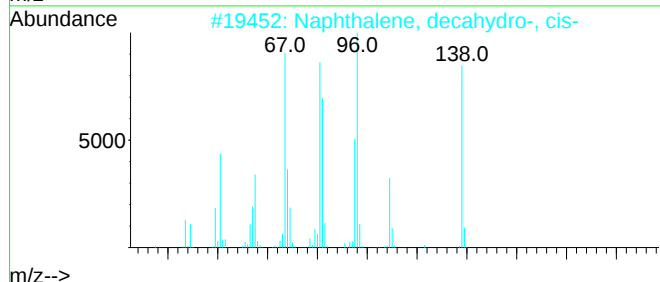
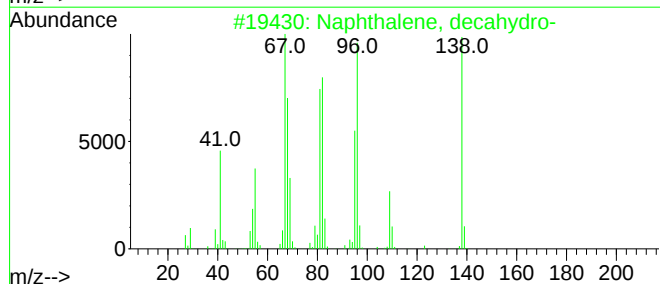
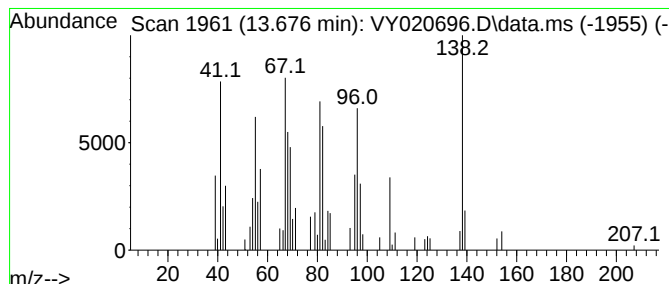
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, decahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.676	98.67 ug/l	26746	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-	138	C10H18	000091-17-8	87
2	Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	87
3	Spiro[4.5]decane	138	C10H18	000176-63-6	83
4	Cyclodecene, (Z)-	138	C10H18	000935-31-9	81
5	Cyclohexene, 1-butyl-	138	C10H18	003282-53-9	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CHRT26634

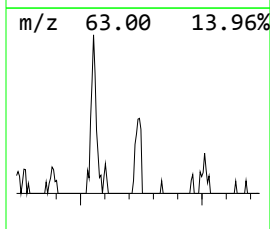
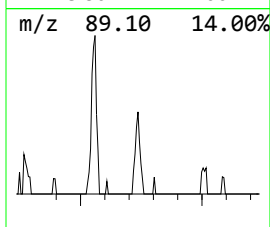
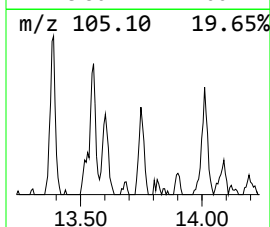
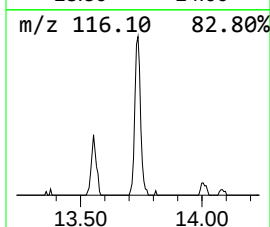
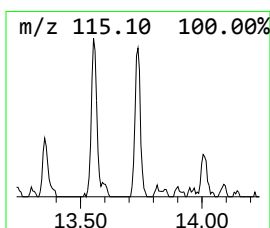
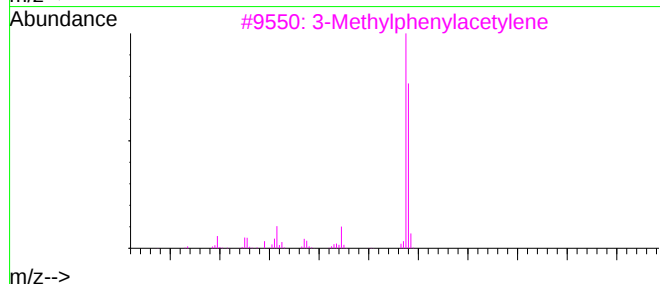
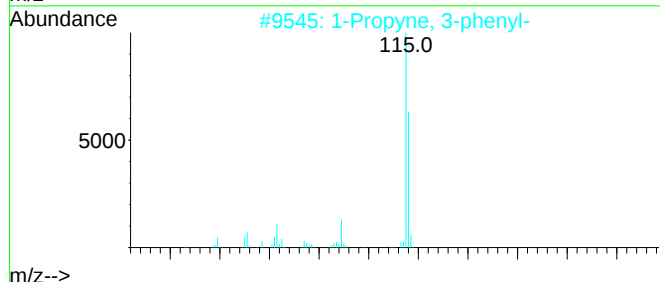
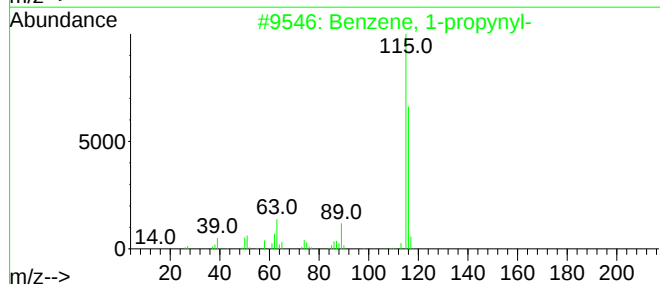
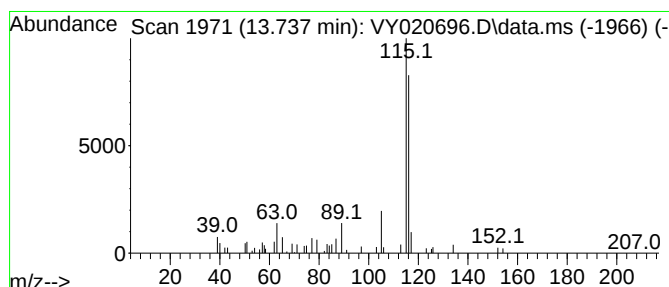
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-propynyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.737	59.59 ug/l	16152	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	80
2	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	72
3	3-Methylphenylacetylene	116	C9H8	000766-82-5	72
4	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	64
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CHRT26634

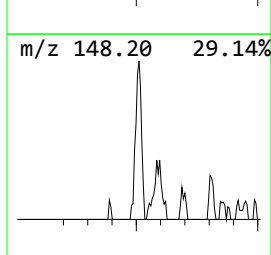
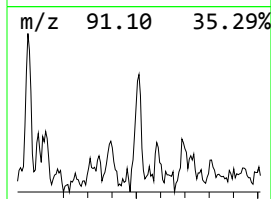
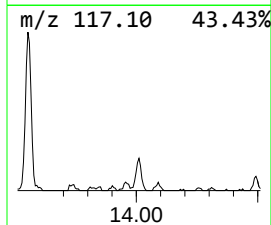
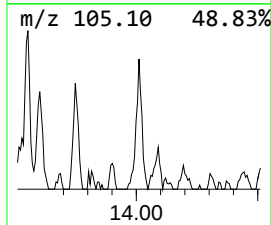
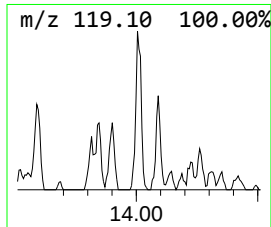
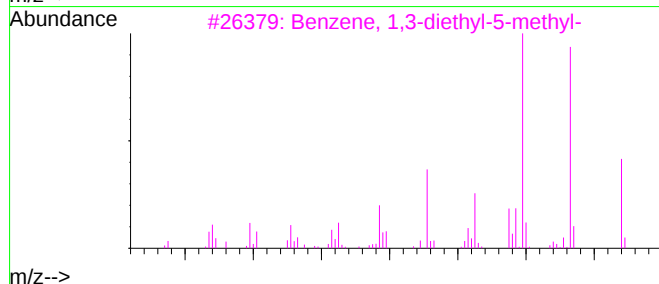
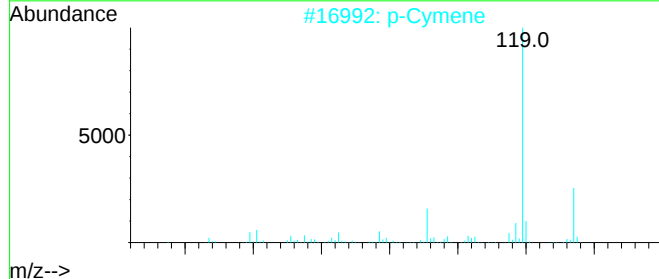
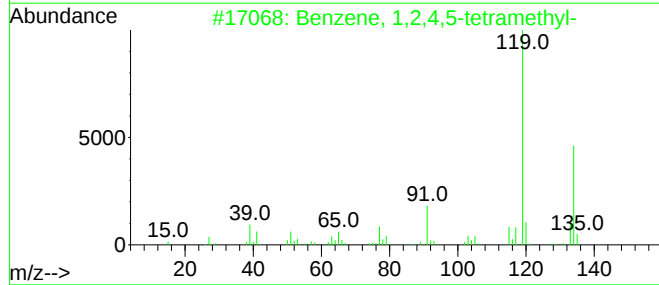
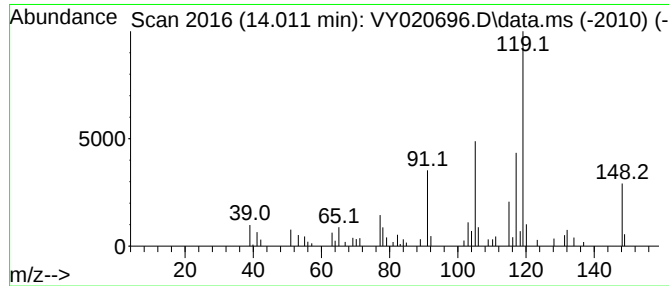
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Quant Title : SW846 8260

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

Peak Number 7 unknown14.011 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.011	96.58 ug/l	26178	1,4-Dichlorobenzene-d4	13.359

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	47
2			p-Cymene	134	C10H14	000099-87-6	47
3			Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	46
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	46
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/soil/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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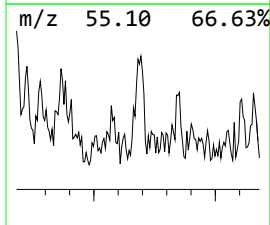
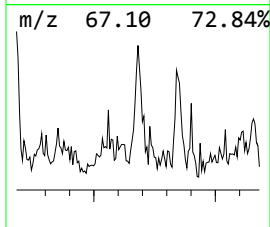
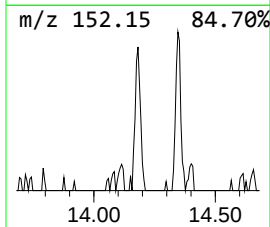
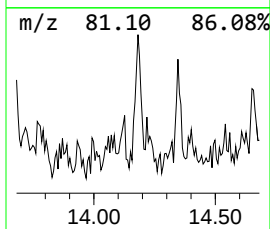
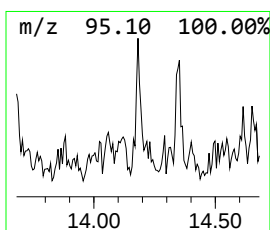
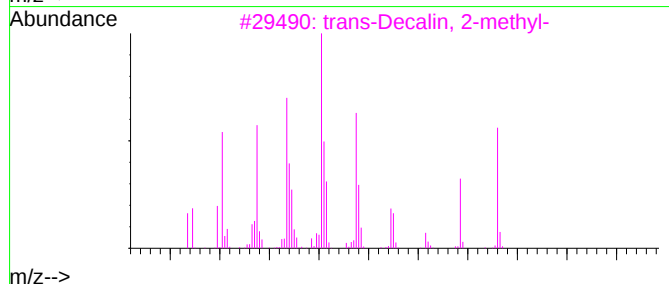
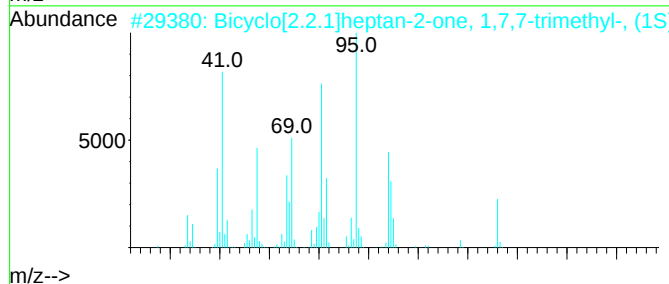
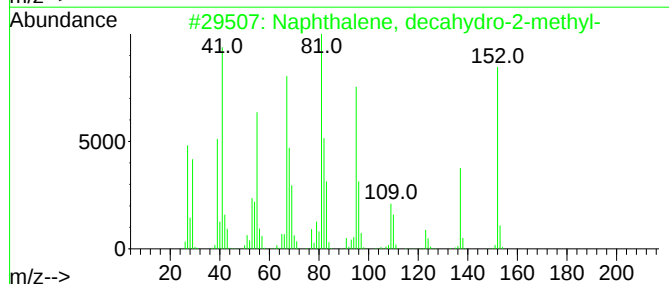
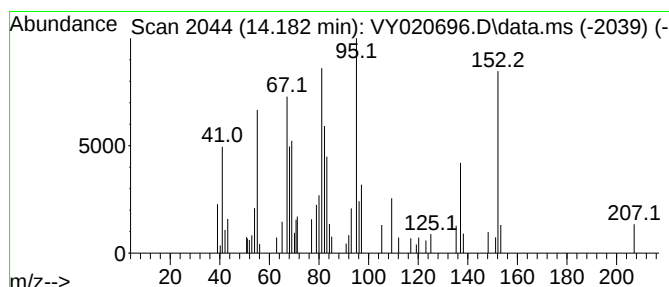
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TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, decahydro-2-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.182	116.07 ug/l	31463	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	89
2	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	76
3	trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	70
4	cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	62
5	Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
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ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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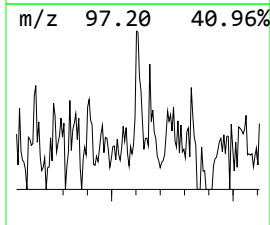
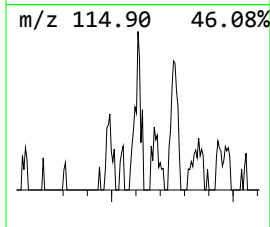
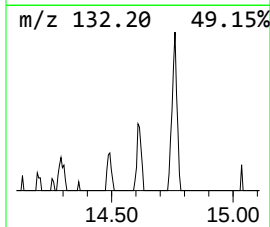
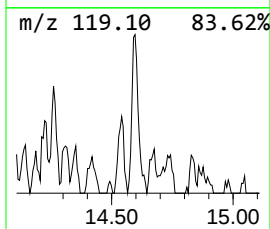
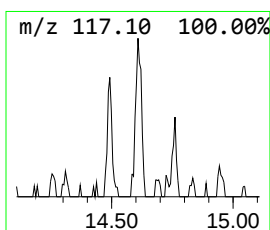
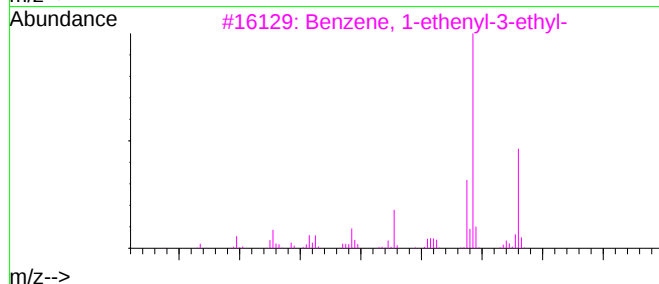
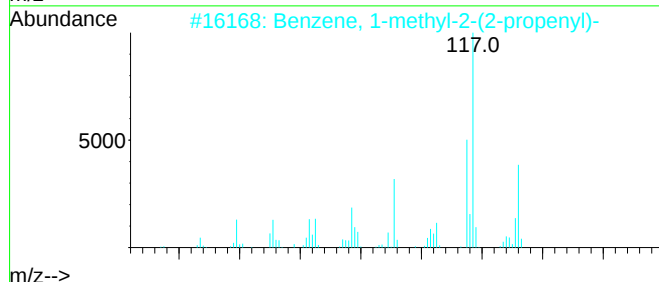
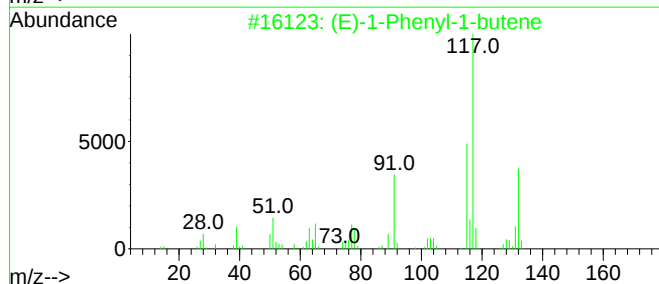
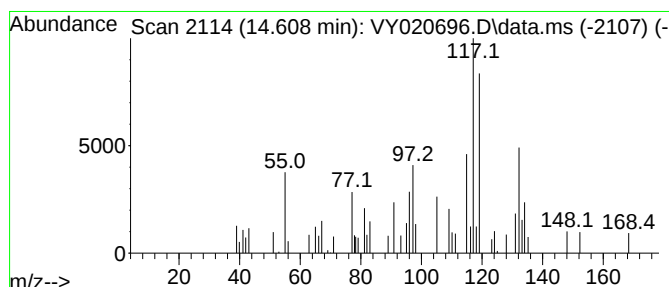
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Quant Title : SW846 8260

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

Peak Number 9 unknown14.608 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.608	65.00 ug/l	17620	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	41
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	41
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	38
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	38
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CHRT26634

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Quant Title : SW846 8260

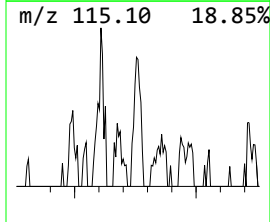
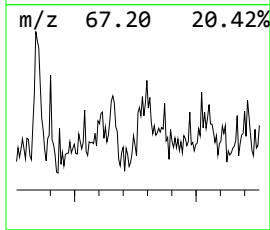
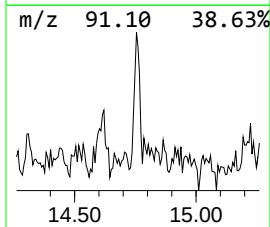
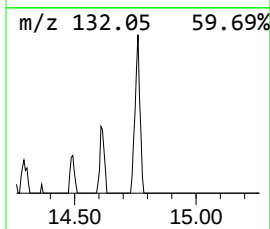
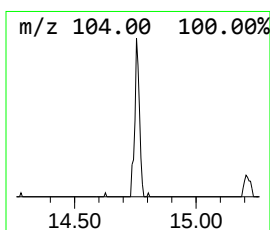
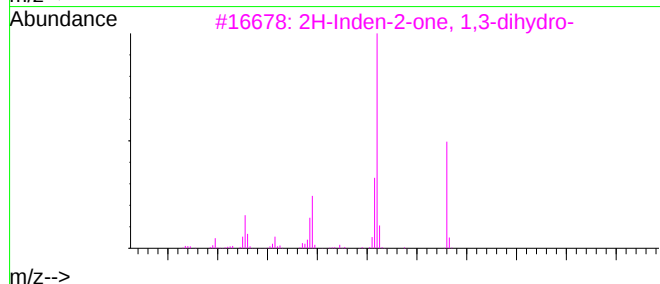
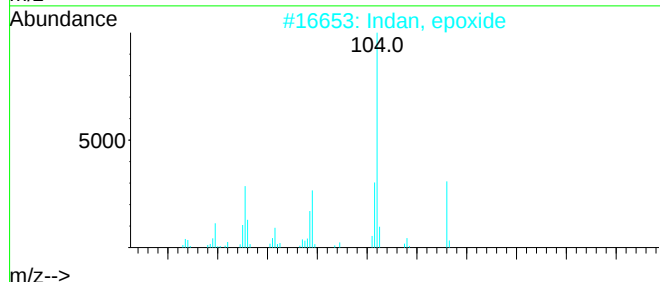
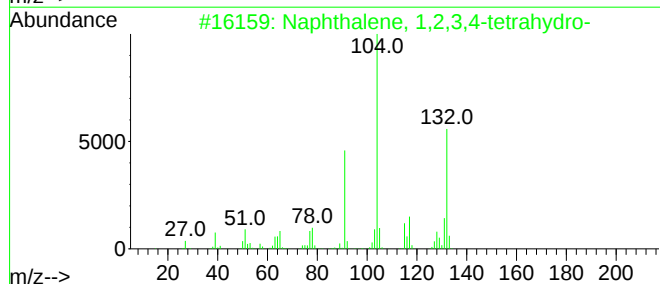
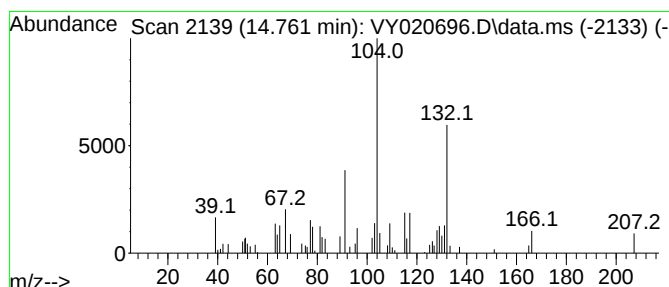
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.761	53.34 ug/l	14459	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	89
2	Indan, epoxide	132	C9H8O	000768-22-9	49
3	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	47
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5	Butanedioic acid, phenyl-	194	C10H10O4	000635-51-8	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY122324\
Data File : VY020696.D
Acq On : 23 Dec 2024 17:04
Operator : SY/MD
Sample : P5342-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CHRT26634

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y122324S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	12.664	63.7	ug/l	17276	4	13.359	13553	50.0
unknown12.968	12.968	52.3	ug/l	14179	4	13.359	13553	50.0
unknown13.249	13.249	80.5	ug/l	21813	4	13.359	13553	50.0
Indane	13.554	150.9	ug/l	40908	4	13.359	13553	50.0
Naphthalene, de...	13.676	98.7	ug/l	26746	4	13.359	13553	50.0
Benzene, 1-prop...	13.737	59.6	ug/l	16152	4	13.359	13553	50.0
unknown14.011	14.011	96.6	ug/l	26178	4	13.359	13553	50.0
Naphthalene, de...	14.182	116.1	ug/l	31463	4	13.359	13553	50.0
unknown14.608	14.608	65.0	ug/l	17620	4	13.359	13553	50.0
Naphthalene, 1,...	14.761	53.3	ug/l	14459	4	13.359	13553	50.0