

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011022\
 Data File : VY007108.D
 Acq On : 10 Jan 2022 17:46
 Operator : SY/MD
 Sample : VIBLK
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VIBLK

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\82Y010822S.M
 Title : SW846 8260

Signal : TIC: VY007108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.058	34	39	48	rBV	5889	16470	1.97%	0.263%
2	4.716	466	475	493	rBV4	7651	27734	3.31%	0.442%
3	5.759	637	646	654	rBV2	3164	8730	1.04%	0.139%
4	6.984	834	847	862	rBV	15689	51438	6.15%	0.820%
5	7.722	959	968	974	rBV	116881	273333	32.66%	4.359%
6	7.795	974	980	992	rVB	184815	426898	51.01%	6.808%
7	8.142	1022	1037	1049	rBV2	114424	328468	39.25%	5.238%
8	8.697	1119	1128	1140	rBV	299659	589435	70.44%	9.399%
9	10.185	1364	1372	1384	rVB	483334	822062	98.23%	13.109%
10	10.581	1430	1437	1447	rVB	98093	173303	20.71%	2.764%
11	10.947	1492	1497	1505	rBV	25869	44239	5.29%	0.705%
12	11.044	1508	1513	1517	rBV3	5103	8782	1.05%	0.140%
13	11.099	1517	1522	1526	rVB3	7453	12129	1.45%	0.193%
14	11.319	1551	1558	1564	rVB	65703	116999	13.98%	1.866%
15	11.496	1580	1587	1597	rVB	438835	731405	87.40%	11.663%
16	11.739	1622	1627	1631	rBV3	7480	14542	1.74%	0.232%
17	12.008	1664	1671	1677	rBV5	12390	30275	3.62%	0.483%
18	12.123	1686	1690	1694	rBV	18738	26001	3.11%	0.415%
19	12.203	1699	1703	1706	rVV5	15191	29671	3.55%	0.473%
20	12.239	1706	1709	1718	rVB6	20705	45088	5.39%	0.719%
21	12.373	1728	1731	1738	rVB5	11182	21395	2.56%	0.341%
22	12.489	1743	1750	1757	rBV3	482332	836839	100.00%	13.345%
23	12.575	1758	1764	1768	rVV9	6226	14700	1.76%	0.234%
24	12.648	1772	1776	1784	rVB4	23122	46642	5.57%	0.744%
25	12.733	1784	1790	1796	rBV7	11149	28615	3.42%	0.456%
26	12.818	1796	1804	1808	rVV5	29897	80947	9.67%	1.291%
27	12.861	1808	1811	1817	rVB4	15878	26586	3.18%	0.424%
28	12.953	1822	1826	1830	rBV4	13046	19287	2.30%	0.308%
29	13.001	1830	1834	1837	rVV4	6025	12406	1.48%	0.198%
30	13.038	1837	1840	1848	rVB6	13310	28110	3.36%	0.448%
31	13.184	1858	1864	1868	rBV2	23497	43195	5.16%	0.689%
32	13.324	1879	1887	1890	rBV5	25427	52825	6.31%	0.842%
33	13.428	1898	1904	1918	rVB	437425	697280	83.32%	11.119%
34	13.568	1923	1927	1930	rBV2	8399	12384	1.48%	0.197%
35	13.702	1943	1949	1951	rBV6	6801	11375	1.36%	0.181%
36	13.751	1952	1957	1961	rVV3	35470	62895	7.52%	1.003%

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37	13.794	1962	1964	1969	rVV3	11189	18095	2.16%	0.289%
38	13.861	1971	1975	1982	rVB8	9163	19621	2.34%	0.313%
39	13.971	1987	1993	1999	rVB2	43420	82443	9.85%	1.315%
40	14.080	2006	2011	2016	rBV6	9915	19020	2.27%	0.303%
41	14.147	2018	2022	2028	rVV8	9098	17155	2.05%	0.274%
42	14.202	2028	2031	2034	rVV4	6519	8839	1.06%	0.141%
43	14.269	2035	2042	2055	rVB6	33133	85405	10.21%	1.362%
44	14.428	2063	2068	2072	rBV5	17363	27278	3.26%	0.435%
45	14.623	2096	2100	2106	rVV8	8331	18048	2.16%	0.288%
46	14.684	2106	2110	2115	rVV5	16086	32103	3.84%	0.512%
47	14.739	2115	2119	2125	rVB5	19956	33920	4.05%	0.541%
48	14.946	2150	2153	2157	rVB4	12212	16630	1.99%	0.265%
49	15.214	2193	2197	2204	rBV	22709	35503	4.24%	0.566%
50	15.324	2211	2215	2221	rVB3	11592	19010	2.27%	0.303%
51	15.611	2257	2262	2267	rVB7	10029	18422	2.20%	0.294%
52	16.031	2325	2331	2337	rBV7	8030	19582	2.34%	0.312%
53	16.159	2348	2352	2357	rVB3	11549	18940	2.26%	0.302%
54	16.482	2401	2405	2408	rBV5	5077	8483	1.01%	0.135%

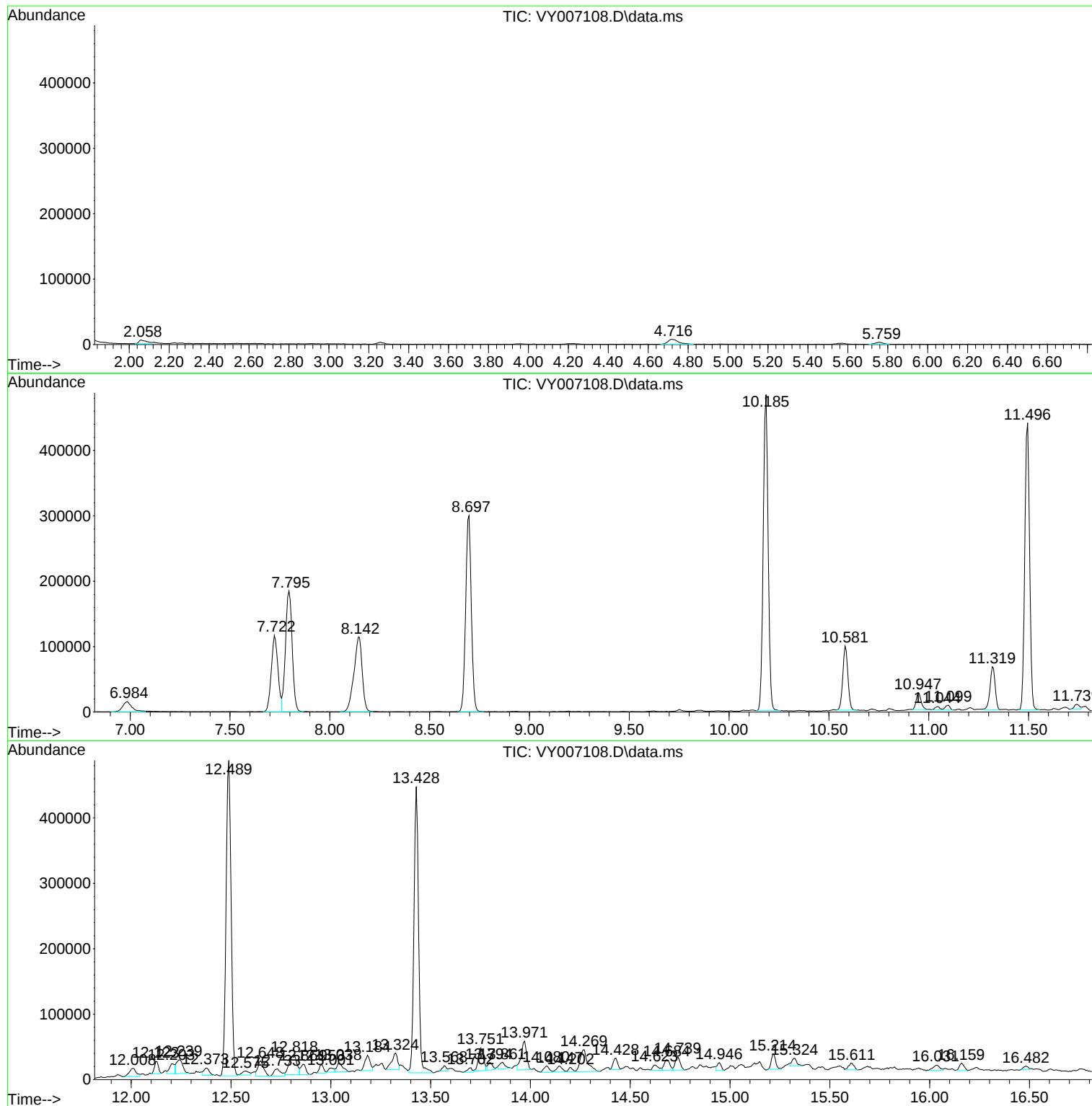
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y010822S.M
Quant Title : SW846 8260

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



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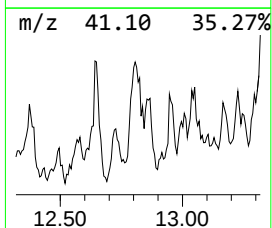
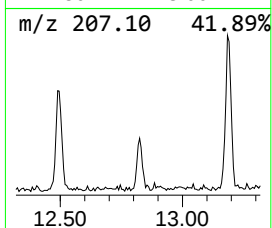
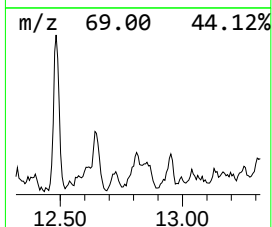
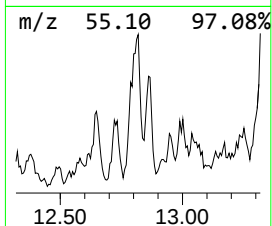
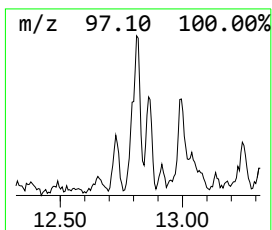
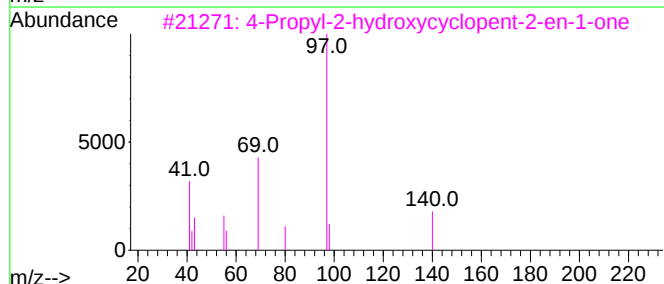
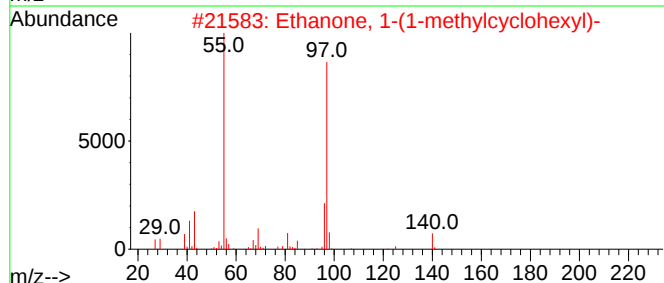
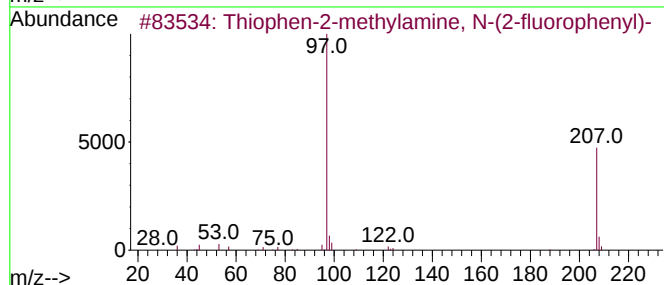
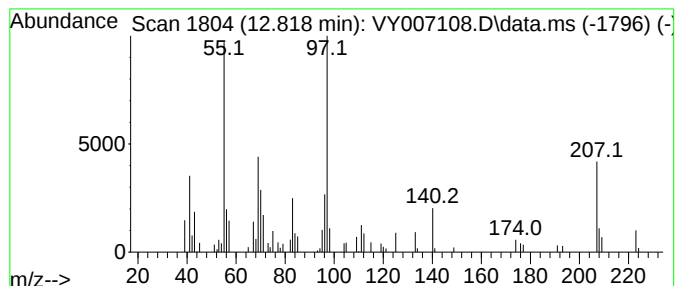
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y010822S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Thiophen-2-methylamine, N-(... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.819	5.80 ug/l	80947	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiophen-2-methylamine, N-(2-flu...	207	C11H10FNS	1000310-35-2	50
2			Ethanone, 1-(1-methylcyclohexyl)-	140	C9H16O	002890-62-2	47
3			4-Propyl-2-hydroxycyclopent-2-en...	140	C8H12O2	082147-26-0	43
4			1-Methyl-4-(1-methylethyl)-cyclo...	140	C10H20	000099-82-1	38
5			Semustine	247	C10H18ClN3O2	013909-09-6	38



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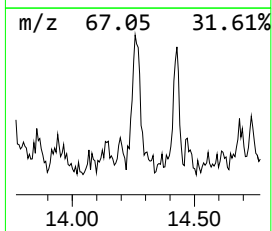
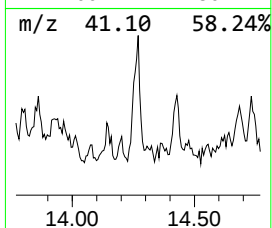
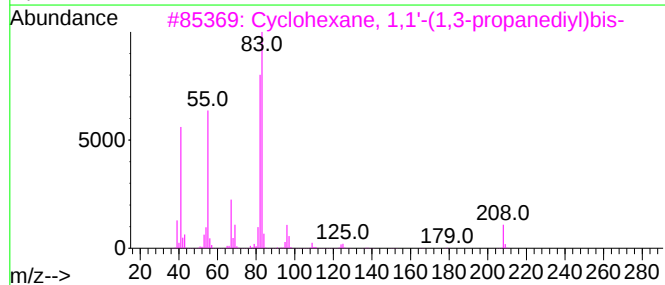
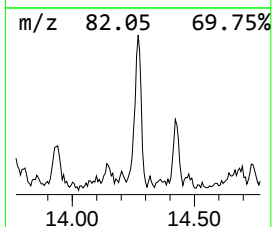
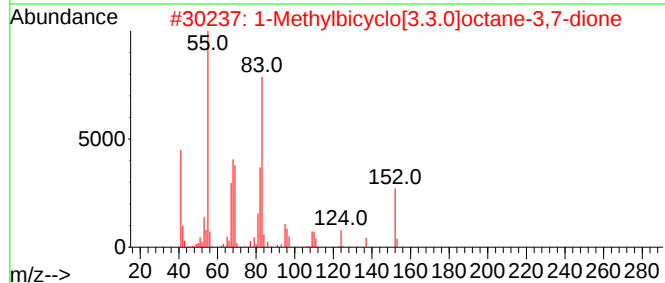
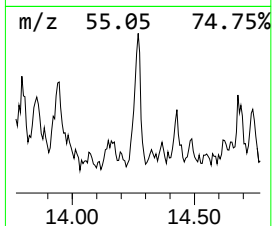
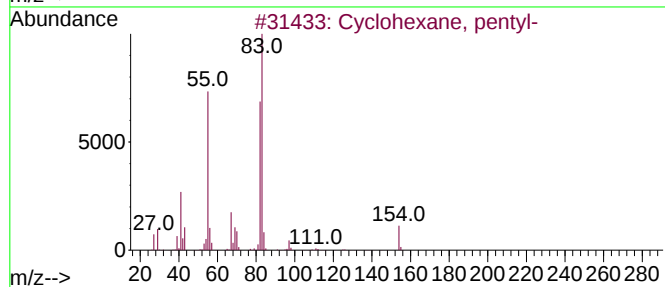
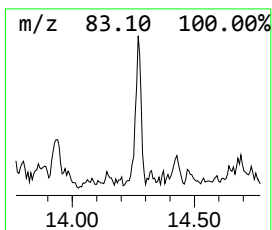
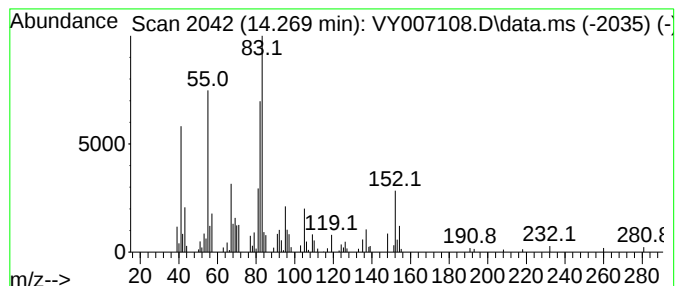
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclohexane, pentyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.269	6.12 ug/l	85405	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	78
2			1-Methylbicyclo[3.3.0]octane-3,7...	152	C9H12O2	021170-08-1	58
3			Cyclohexane, 1,1'-(1,3-propanedi...	208	C15H28	003178-24-3	49
4			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	47
5			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	46



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Thiophen-2-meth...	12.819	5.8	ug/l	80947	4	13.428	697280	50.0
Cyclohexane, pe...	14.269	6.1	ug/l	85405	4	13.428	697280	50.0