

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY031122\
 Data File : VY007864.D
 Acq On : 11 Mar 2022 18:57
 Operator : SY/MD
 Sample : N1864-14
 Misc : 6.46g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 MB-B1-030922-5-6

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\82Y022822S.M

Title : SW846 8260

Signal : TIC: VY007864.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.516	108	114	125	rBV3	23577	62524	5.37%	1.998%
2	3.046	199	201	209	rVB6	5755	12089	1.04%	0.386%
3	4.710	465	474	486	rVB5	12002	39095	3.36%	1.249%
4	7.716	958	967	973	rBV2	61711	145710	12.51%	4.655%
5	7.789	973	979	993	rVB	96596	211652	18.17%	6.762%
6	8.143	1028	1037	1048	rBV2	61247	145086	12.45%	4.635%
7	8.691	1118	1127	1136	rBV	134571	275208	23.62%	8.792%
8	10.179	1364	1371	1378	rBV	222339	378693	32.50%	12.098%
9	11.490	1577	1586	1594	rBV	206674	336355	28.87%	10.746%
10	12.489	1741	1750	1759	rBV2	676427	1165091	100.00%	37.222%
11	13.422	1897	1903	1911	rBV	182595	289099	24.81%	9.236%
12	13.959	1985	1991	1999	rBV2	41140	69501	5.97%	2.220%

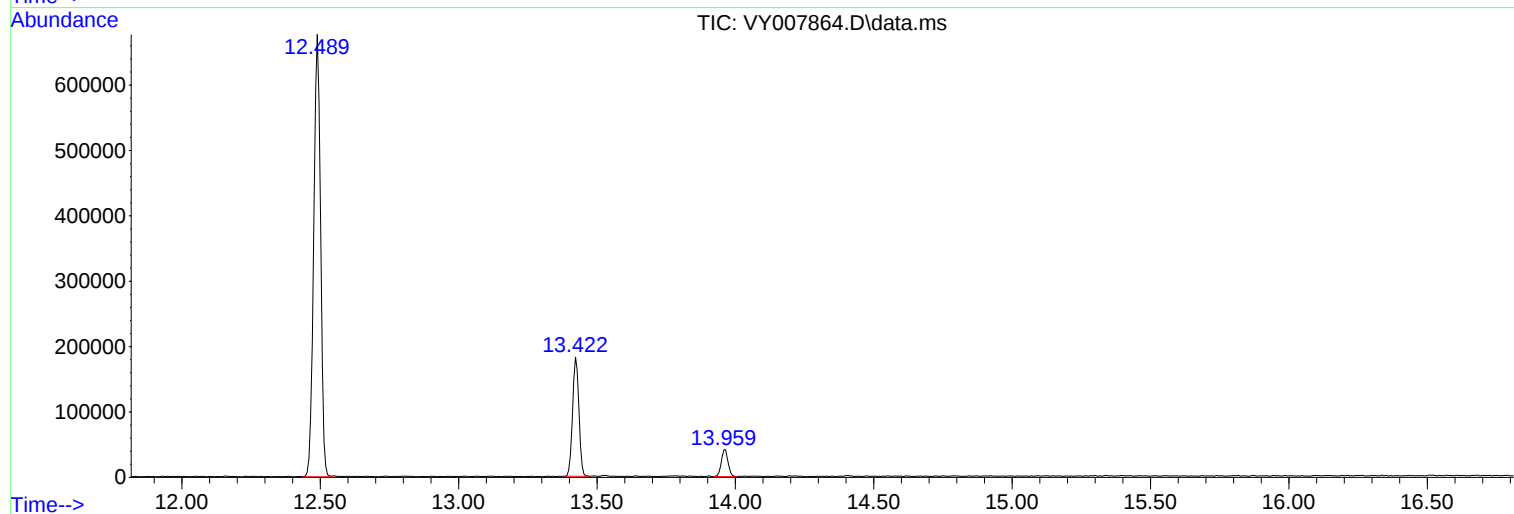
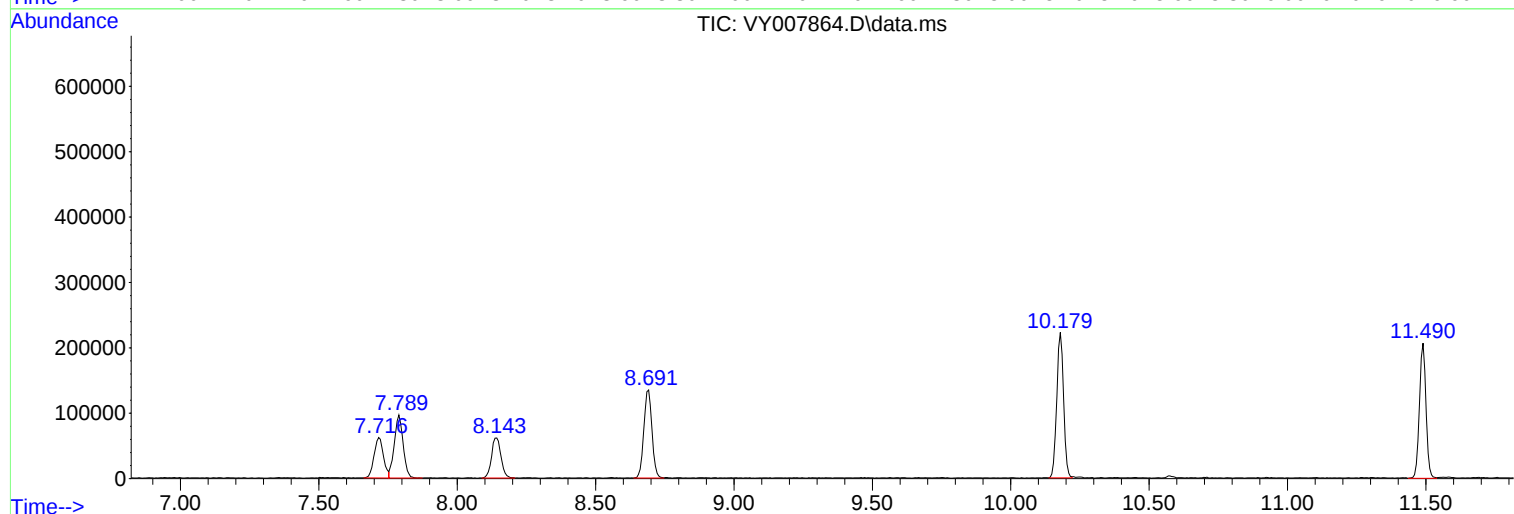
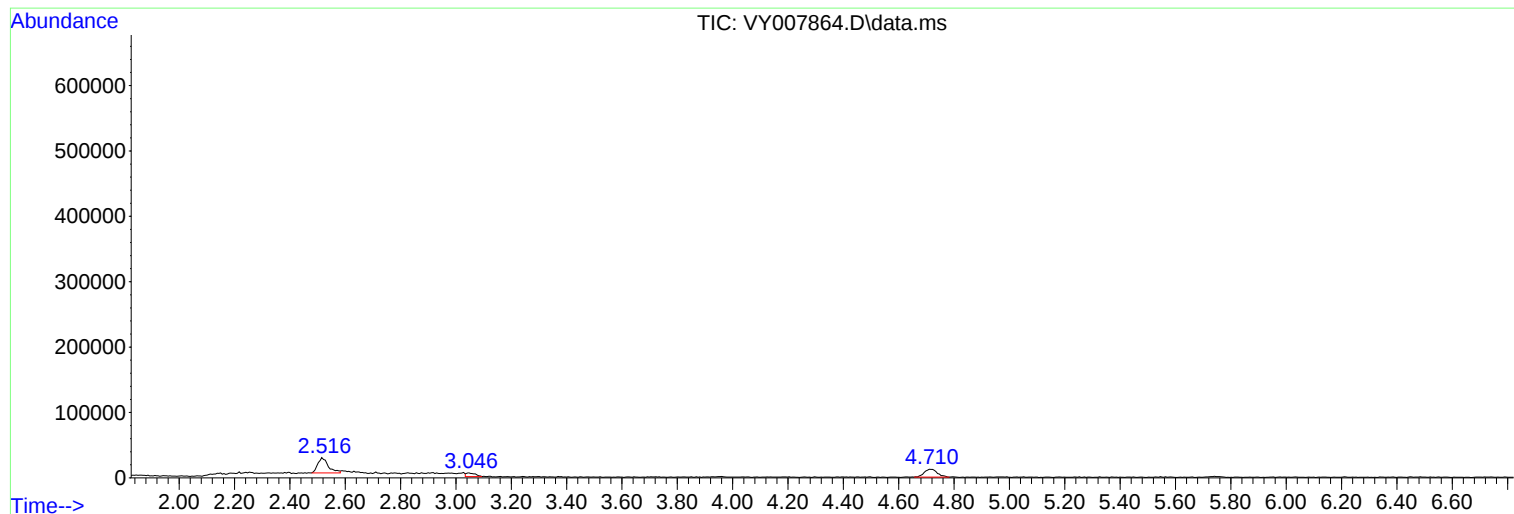
Sum of corrected areas: 3130103

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.M
Quant Title : SW846 8260

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



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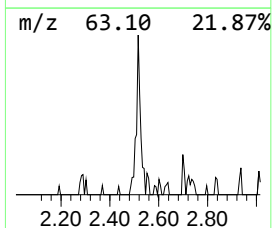
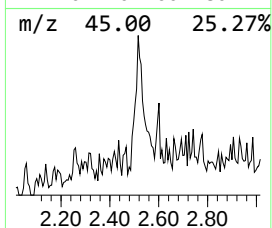
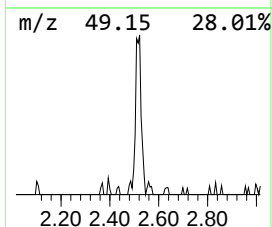
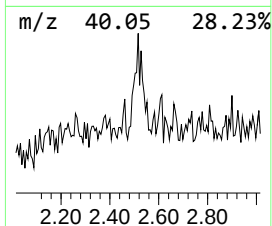
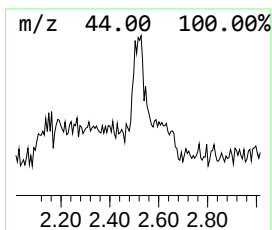
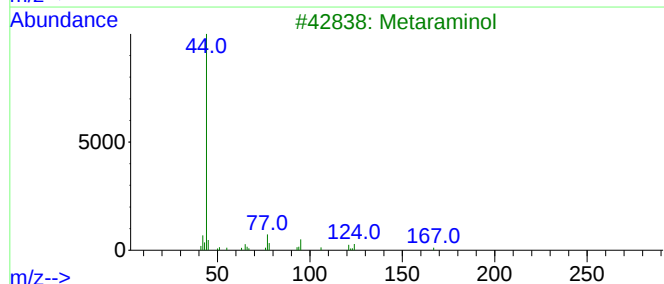
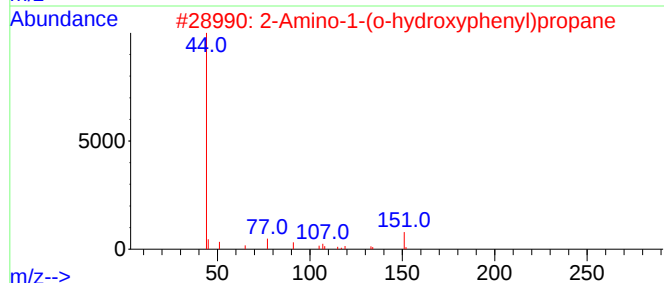
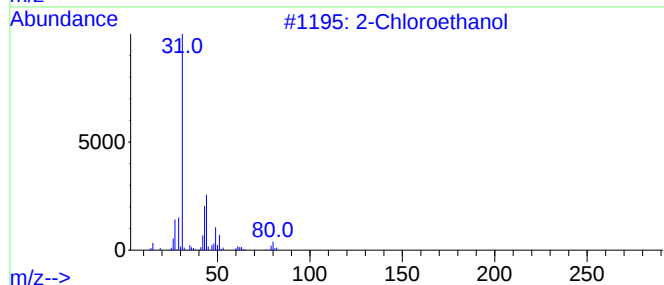
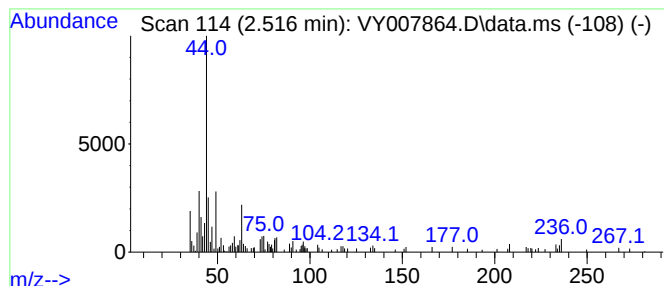
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022822S.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.516 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.516	14.77 ug/l	62524	Pentafluorobenzene	7.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloroethanol	80	C2H5ClO	000107-07-3	47
2			2-Amino-1-(o-hydroxyphenyl)propane	151	C9H13NO	073470-00-5	25
3			Metaraminol	167	C9H13NO2	000054-49-9	22
4			Ethyl oxamate	117	C4H7NO3	000617-36-7	22
5			Phenethylamine, 4-iodo-.alpha.-m...	261	C9H12IN	021894-58-6	16



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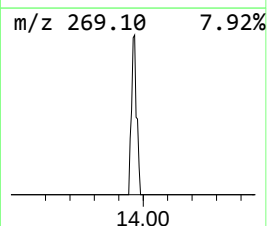
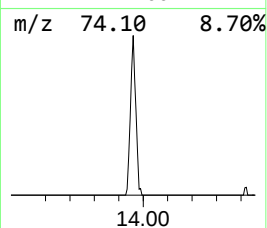
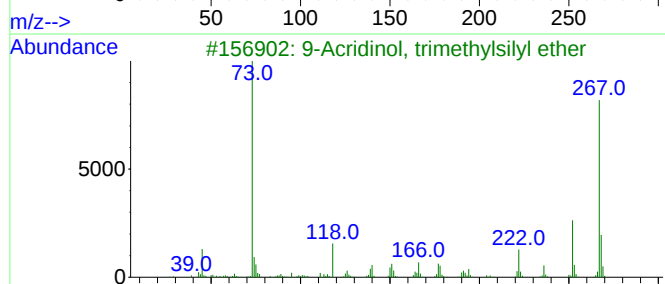
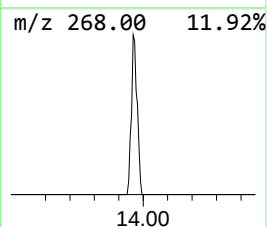
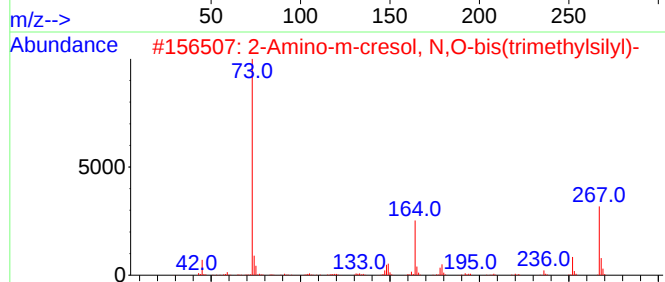
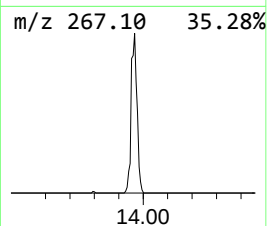
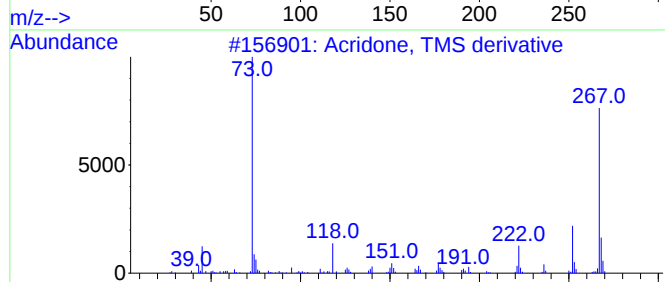
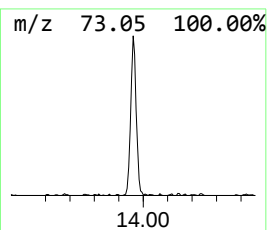
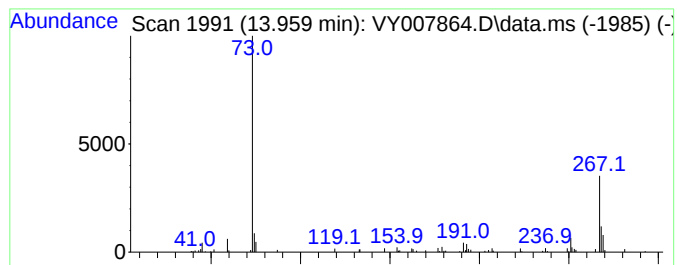
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Peak Number 2 unknown13.959 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.959	12.02 ug/l	69501	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acridone, TMS derivative	267	C16H17NOSi	1000478-16-0	45
2			2-Amino-m-cresol, N,O-bis(trimet...	267	C13H25NOSi2	1000449-77-1	42
3			9-Acridinol, trimethylsilyl ether	267	C16H17NOSi	1000463-65-5	42
4			Salicylic acid, 2TMS derivative	282	C13H22O3Si2	003789-85-3	42
5			Benzeneethanamine, N-methyl-.bet...	311	C15H29NO2Si2	029522-18-7	39



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.516	2.516	14.8	ug/l	62524	1	7.789	211652	50.0
unknown13.959	13.959	12.0	ug/l	69501	4	13.422	289099	50.0