

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022825\
 Data File : VY021381.D
 Acq On : 28 Feb 2025 15:31
 Operator : SY/MD
 Sample : Q1422-02
 Misc : 8.94g/5.0mL/MSVOA_Y/SOIL/B
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 B-154-SB02

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\82Y022525S.M

Title : SW846 8260

Signal : TIC: VY021381.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.610	465	474	483	rBV3	11650	35266	4.38%	0.847%
2	5.646	632	644	654	rVB6	4938	17162	2.13%	0.412%
3	7.634	958	970	976	rBV	120861	296244	36.79%	7.118%
4	7.707	976	982	994	rVB	182741	413606	51.36%	9.938%
5	8.054	1031	1039	1053	rBV	108052	247094	30.68%	5.937%
6	8.615	1122	1131	1142	rBV	295481	594714	73.85%	14.290%
7	10.109	1364	1376	1389	rBV	458408	805288	100.00%	19.350%
8	11.420	1583	1591	1604	rBV	405639	686233	85.22%	16.489%
9	12.407	1744	1753	1762	rBV	298576	479579	59.55%	11.523%
10	13.346	1900	1907	1916	rBV	330689	518132	64.34%	12.450%
11	13.889	1989	1996	2003	rVB2	42557	68454	8.50%	1.645%

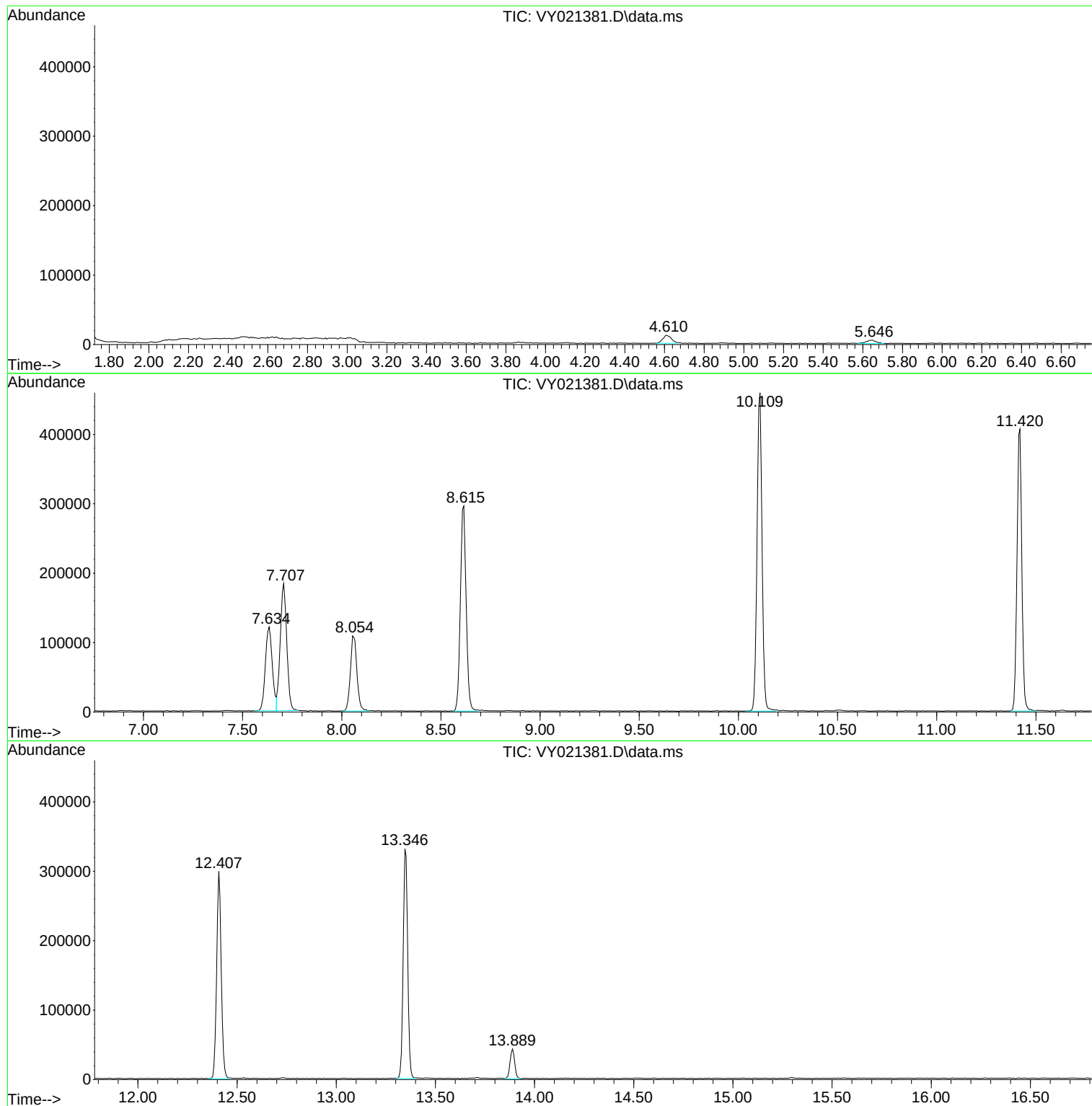
Sum of corrected areas: 4161772

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022525S.M
Quant Title : SW846 8260

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



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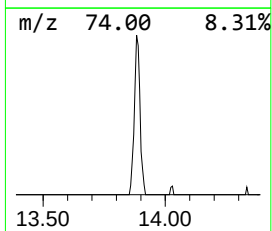
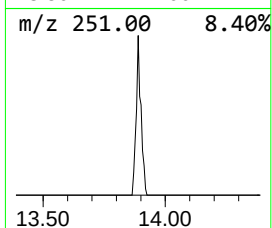
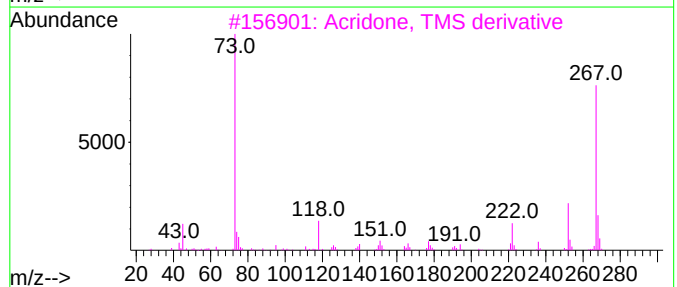
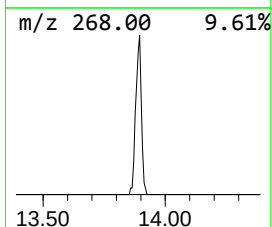
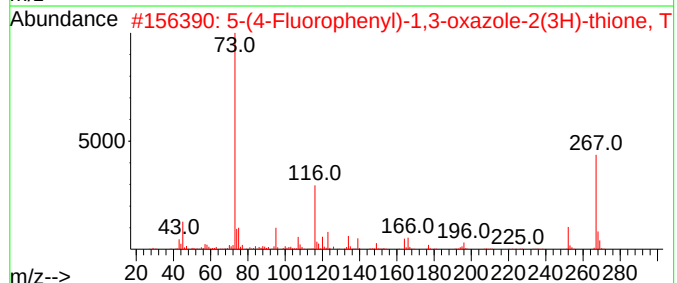
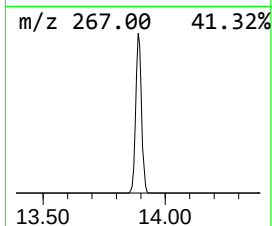
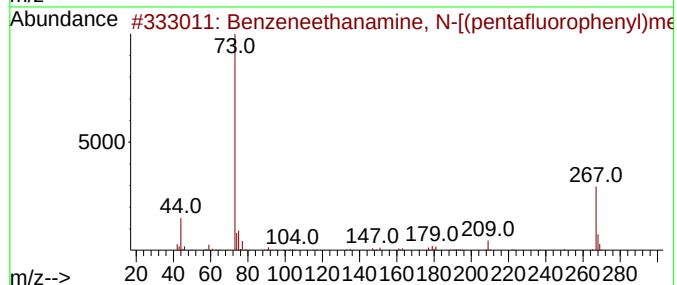
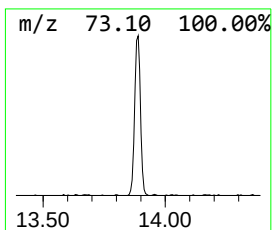
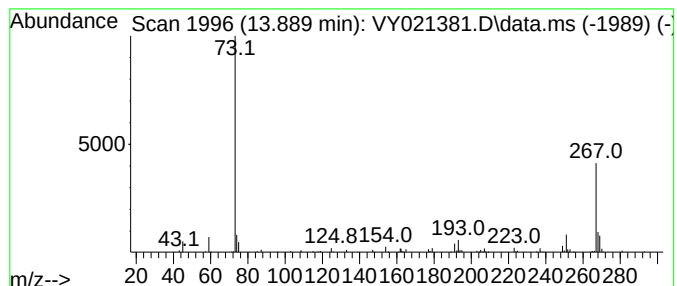
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022525S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzeneethanamine, N-[(pent... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.889	6.61 ug/l	68454	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5NO2Si2	055429-85-1	64
2			5-(4-Fluorophenyl)-1,3-oxazole-2...	267	C12H14FNOSSi	1000497-07-6	64
3			Acridone, TMS derivative	267	C16H17NOSi	1000478-16-0	59
4			Salicylic acid, 2TMS derivative	282	C13H22O3Si2	003789-85-3	50
5			4-(Methylamino)phenol, N,O-bis(t...	267	C13H25NOSi2	1000471-13-0	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanami...	13.889	6.6	ug/l	68454	4	13.352	518132	50.0