

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081922\
 Data File : VY010131.D
 Acq On : 20 Aug 2022 03:12
 Operator : KP/MD
 Sample : N4142-05
 Misc : 6.84g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 UST-BOT

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\82Y081922S.M

Title : SW846 8260

Signal : TIC: VY010131.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.186	380	388	403	rVB	17366	47398	5.24%	0.977%
2	4.710	464	474	481	rBV4	12168	39033	4.31%	0.804%
3	4.795	483	488	500	rVB3	7283	22284	2.46%	0.459%
4	7.710	956	966	972	rBV	76152	184956	20.43%	3.811%
5	7.783	972	978	994	rVB	240397	533984	58.99%	11.004%
6	8.136	1027	1036	1049	rBV	72299	156685	17.31%	3.229%
7	8.685	1117	1126	1142	rBV	378663	730990	80.75%	15.063%
8	10.179	1363	1371	1380	rBV	534909	905230	100.00%	18.654%
9	10.575	1428	1436	1443	rBV2	9340	17459	1.93%	0.360%
10	11.489	1578	1586	1596	rBV	512185	835334	92.28%	17.213%
11	12.483	1742	1749	1758	rBV2	393802	708535	78.27%	14.601%
12	13.422	1896	1903	1912	rVB	384825	594958	65.72%	12.260%
13	13.843	1966	1972	1977	rBV3	8012	14257	1.57%	0.294%
14	13.958	1985	1991	2000	rVB	36688	61702	6.82%	1.271%

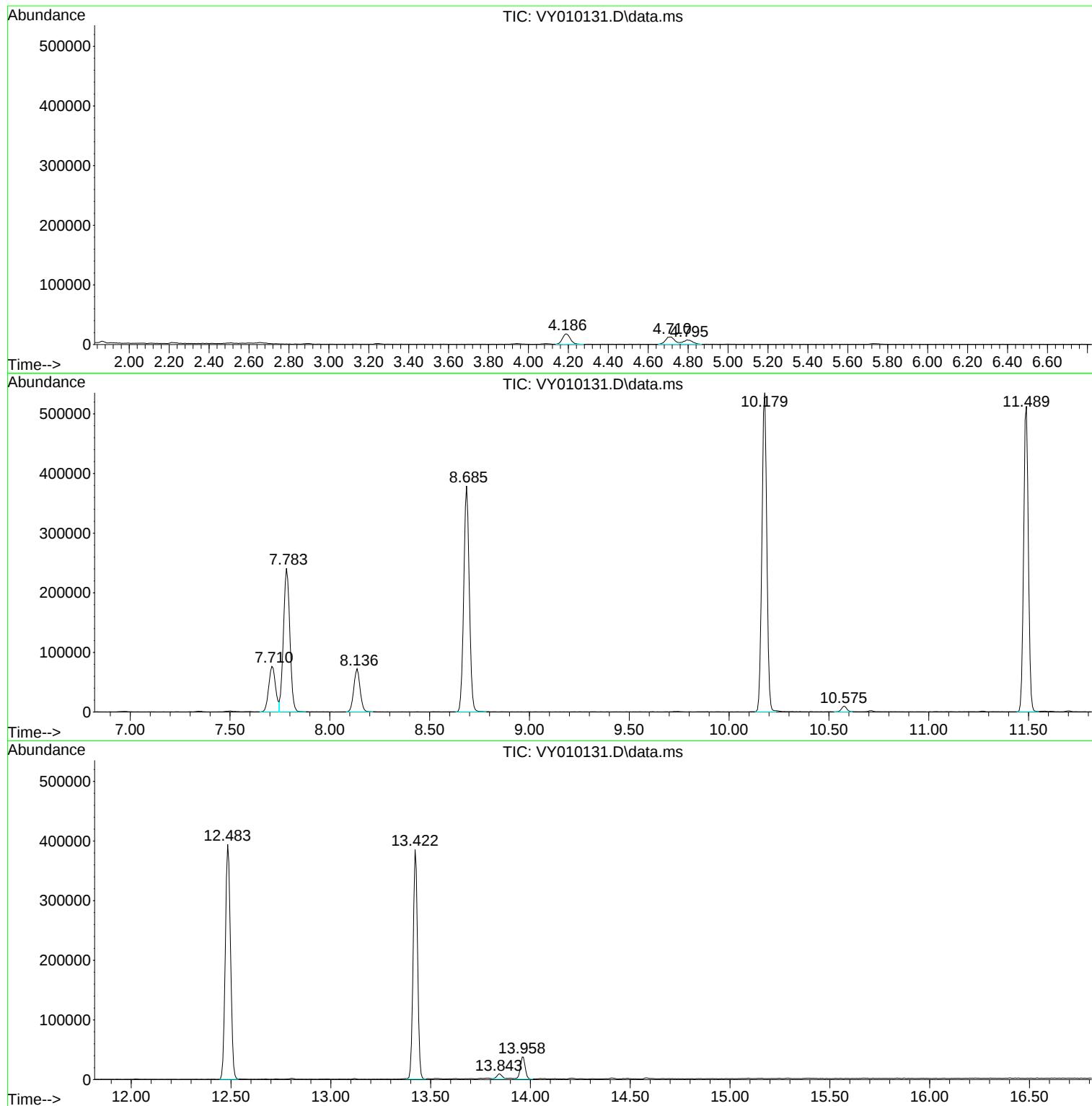
Sum of corrected areas: 4852805

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y081922S.M
Quant Title : SW846 8260

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



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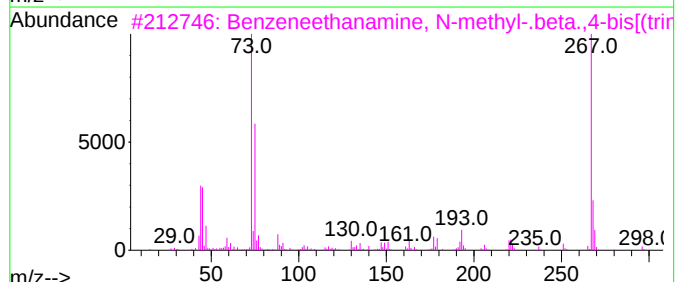
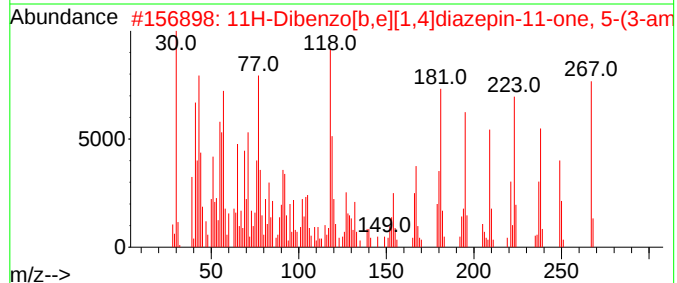
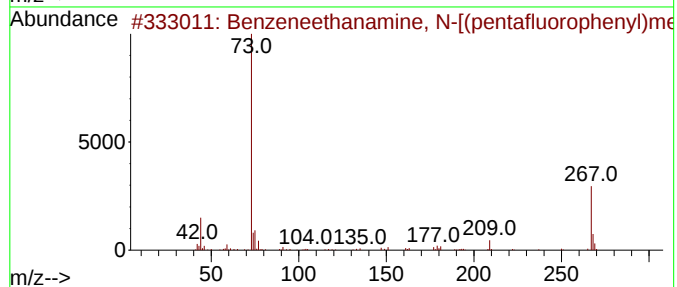
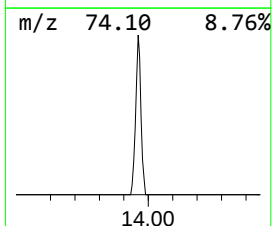
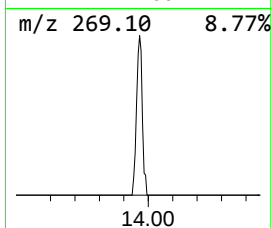
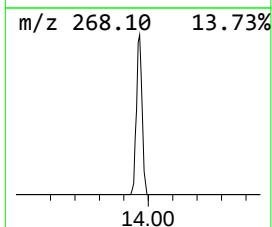
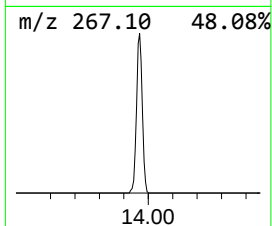
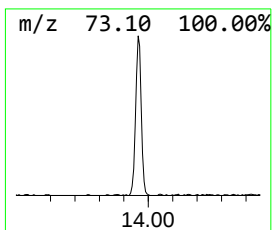
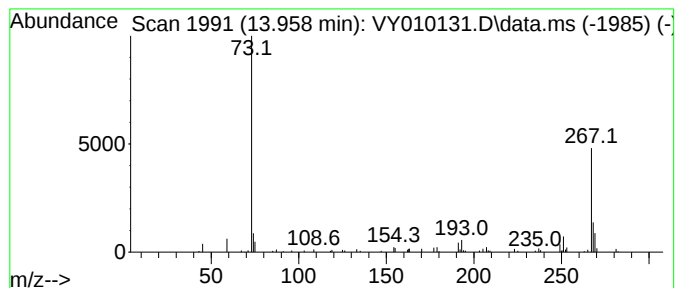
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Peak Number 1 Benzeneethanamine, N-[(pent... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.958	5.19 ug/l	61702	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5NO2Si2	055429-85-1	56
2			11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	53
3			Benzeneethanamine, N-methyl-.bet...	311	C15H29NO2Si2	029522-18-7	45
4			9-Acridinol, trimethylsilyl ether	267	C16H17NOSi	1000463-65-5	45
5			Salicylic acid, 2TMS derivative	282	C13H22O3Si2	003789-85-3	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanami...	13.958	5.2	ug/l	61702	4	13.422	594958	50.0