

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012522\
 Data File : VY007296.D
 Acq On : 25 Jan 2022 17:42
 Operator : SY/MD
 Sample : N1266-06
 Misc : 7.16g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 RB-08-(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\82Y011322S.M

Title : SW846 8260

Signal : TIC: VY007296.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.716	957	967	973	rBV2	69887	168075	9.62%	4.102%
2	7.789	973	979	990	rVB	112546	249862	14.30%	6.099%
3	8.142	1026	1037	1048	rVB	70198	160642	9.19%	3.921%
4	8.685	1118	1126	1139	rVB	172437	346605	19.84%	8.460%
5	10.179	1363	1371	1380	rBV	289306	480051	27.48%	11.717%
6	11.489	1577	1586	1595	rBV	270698	437209	25.02%	10.671%
7	12.489	1741	1750	1760	rBV2	1010493	1747124	100.00%	42.644%
8	13.422	1893	1903	1910	rBV	293592	454317	26.00%	11.089%
9	13.964	1986	1992	1998	rVB	33662	53116	3.04%	1.296%

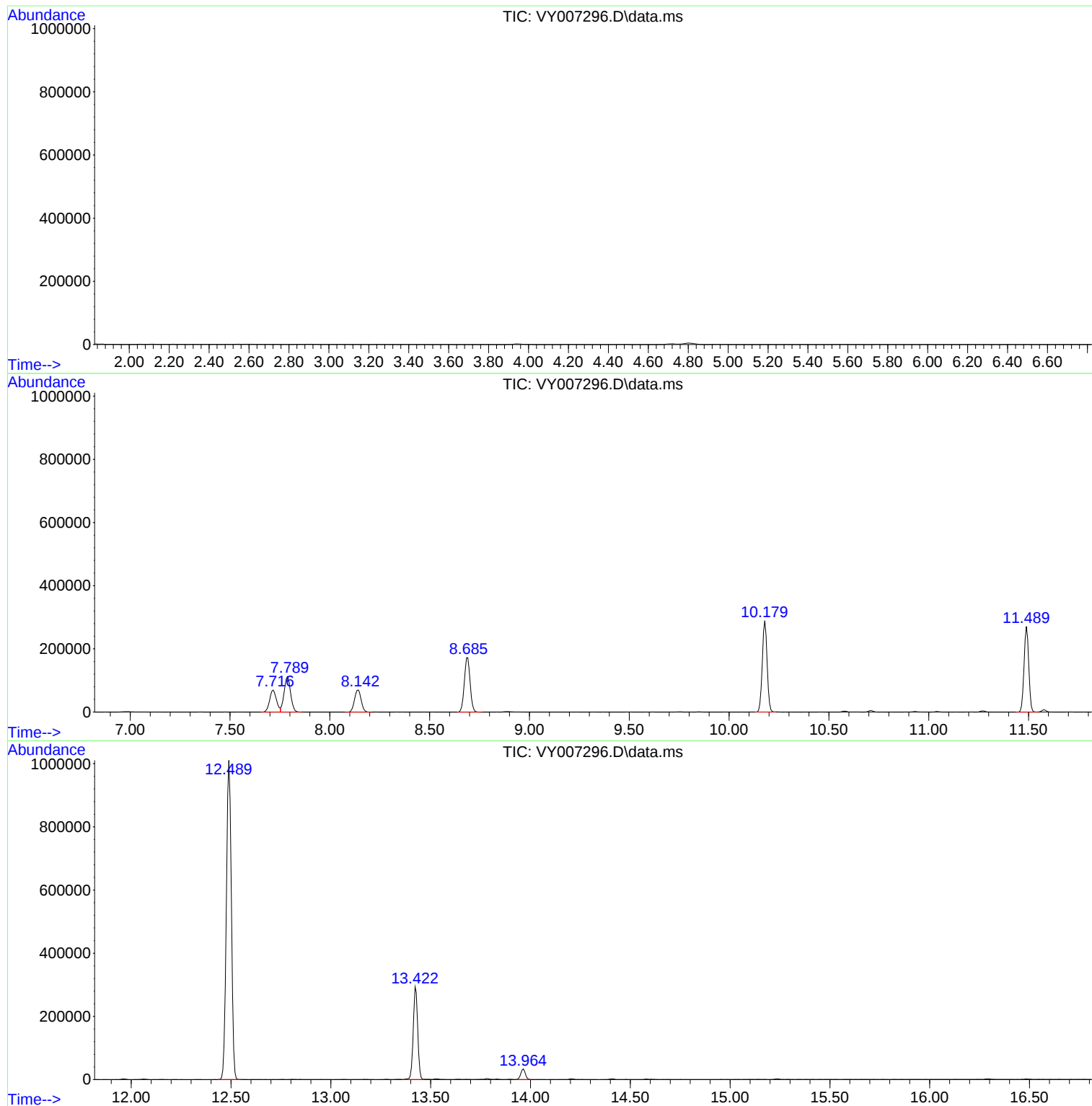
Sum of corrected areas: 4097001

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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



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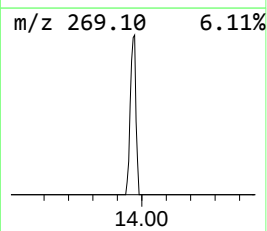
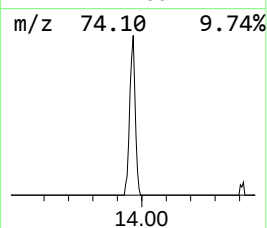
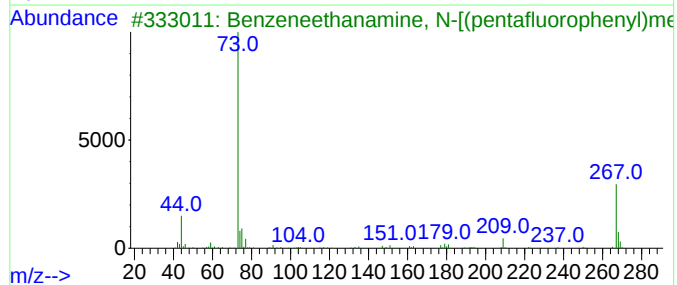
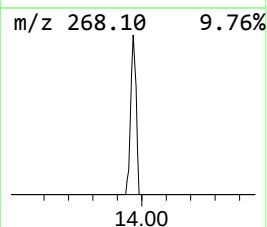
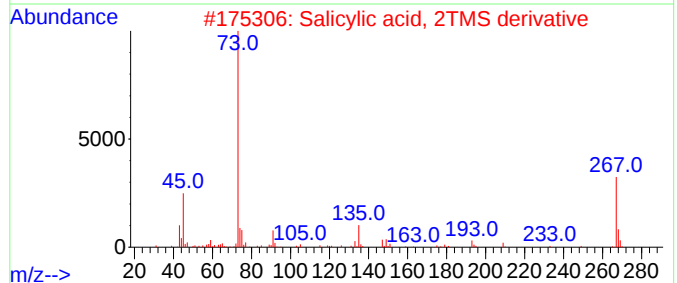
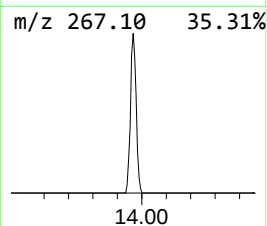
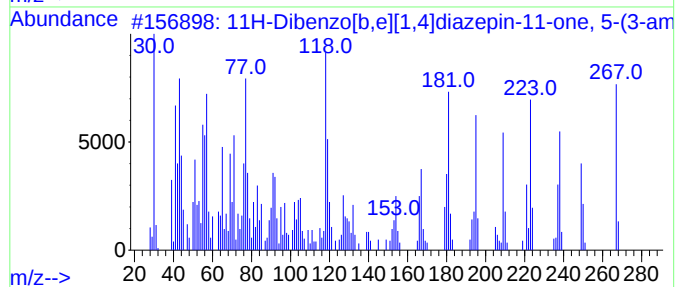
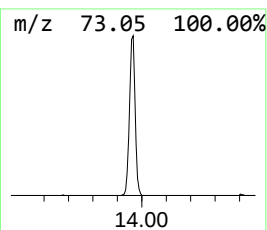
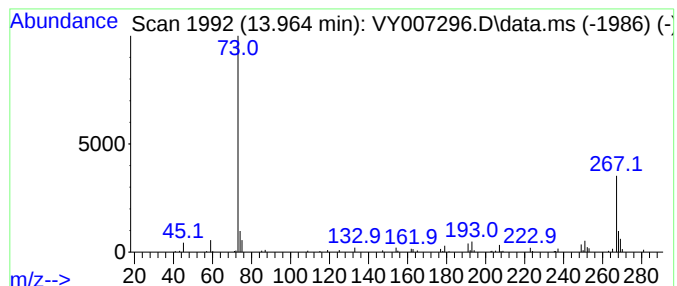
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011322S.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown13.964 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.964	5.85 ug/l	53116	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	43
2			Salicylic acid, 2TMS derivative	282	C13H22O3Si2	003789-85-3	42
3			Benzeneethanamine, N-[(pentafluo...	475	C21H26F5NO2Si2	055429-85-1	42
4			Acridone, TMS derivative	267	C16H17NO5i	1000478-16-0	42
5			Benzeneethanamine, N-methyl-.bet...	311	C15H29NO2Si2	029522-18-7	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown13.964	13.964	5.8	ug/l	53116	4	13.422	454317	50.0