

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041621\
 Data File : VY004470.D
 Acq On : 16 Apr 2021 15:22
 Operator : SY/MD
 Sample : M2050-07
 Misc : 5.57G/5ML/MSVOA_Y/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TP-5

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\82Y041221S.M

Title : SW846 8260

Signal : TIC: VY004470.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.229	53	67	69	rBV6	8497	27401	1.64%	0.496%
2	2.589	113	126	127	rBV6	37308	122779	7.35%	2.224%
3	4.735	466	478	499	rVB4	8043	28520	1.71%	0.517%
4	7.728	958	969	975	rBV2	112566	278218	16.65%	5.039%
5	7.801	975	981	992	rVB	196100	444649	26.61%	8.053%
6	8.155	1026	1039	1048	rBV	124190	279284	16.72%	5.058%
7	8.697	1120	1128	1139	rVB	314861	611257	36.59%	11.071%
8	10.185	1362	1372	1380	rBV	513604	869261	52.03%	15.743%
9	11.496	1577	1587	1595	rBV	412621	673219	40.30%	12.193%
10	12.489	1741	1750	1759	rBV2	962296	1670714	100.00%	30.259%
11	13.428	1898	1904	1912	rVB	295381	447865	26.81%	8.111%
12	13.965	1986	1992	1998	rBV	41474	68231	4.08%	1.236%

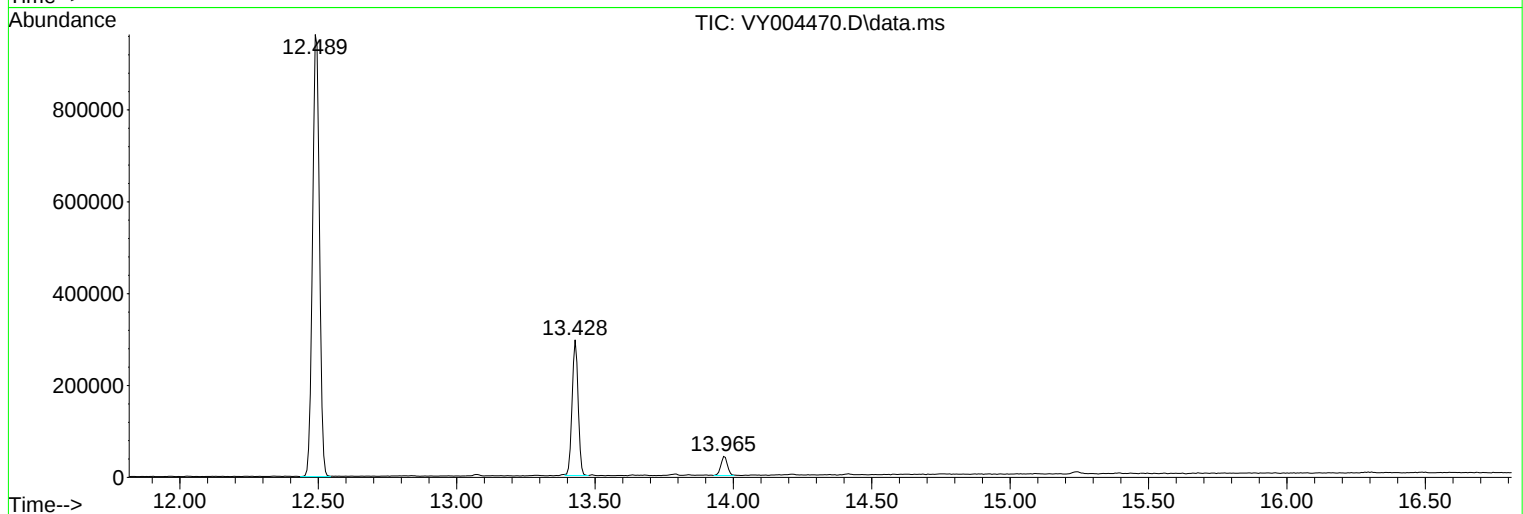
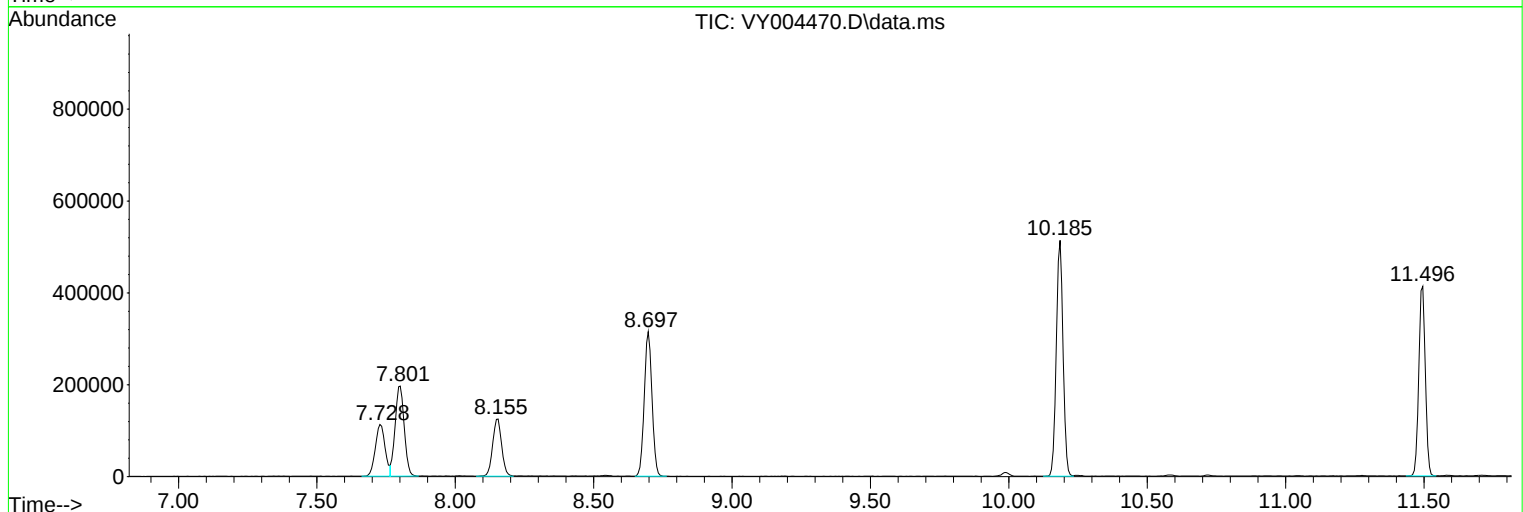
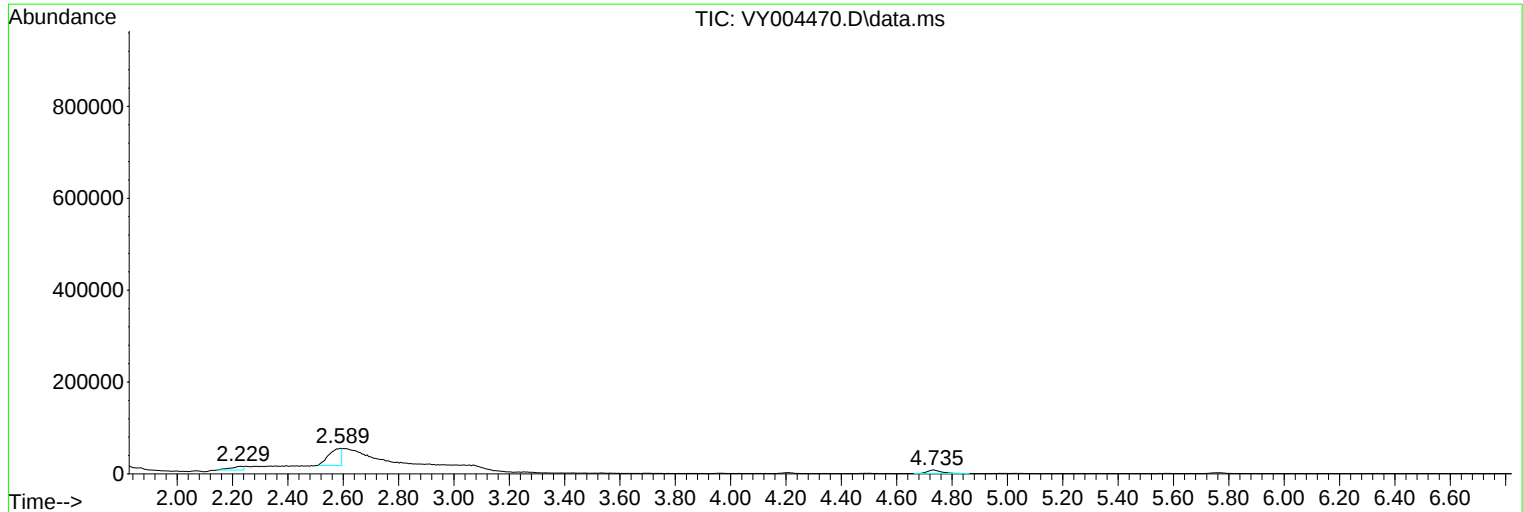
Sum of corrected areas: 5521398

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TP-5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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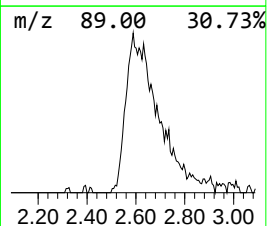
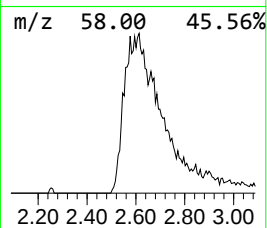
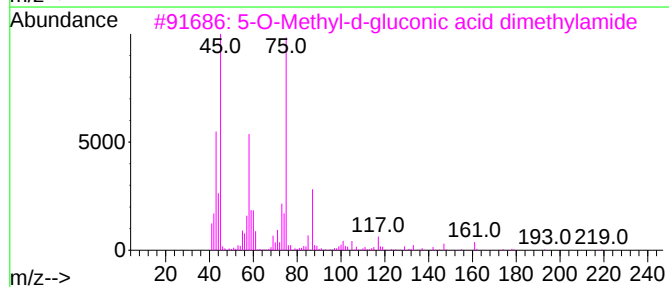
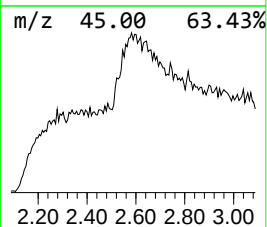
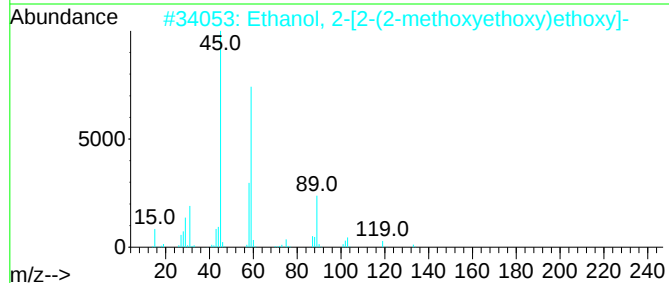
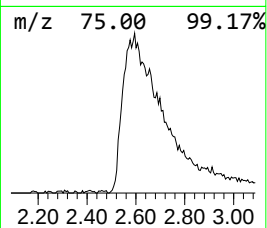
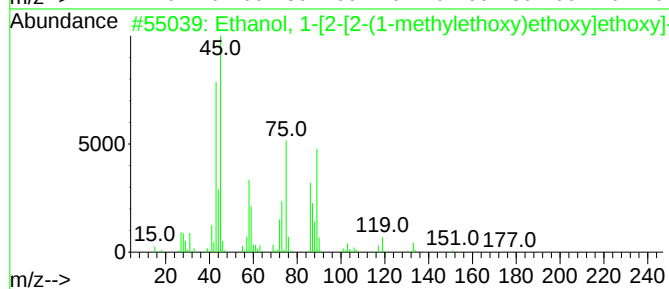
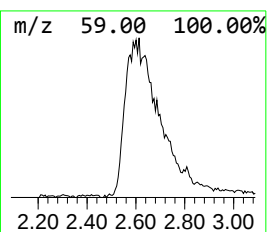
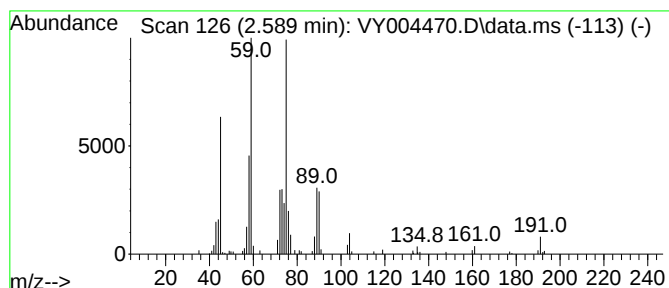
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
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Peak Number 1 unknown2.589 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.589	13.81 ug/l	122779	Pentafluorobenzene	7.801

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 1-[2-[2-(1-methylethoxy...	192	C9H20O4	029681-21-8	43
2			Ethanol, 2-[2-(2-methoxyethoxy)e...	164	C7H16O4	000112-35-6	38
3			5-O-Methyl-d-gluconic acid dimet...	237	C9H19NO6	013096-67-8	38
4			Ethane, 1-ethoxy-1-methoxy-	104	C5H12O2	010471-14-4	35
5			Tetraethyleneglycol monomethylether	208	C9H20O5	023783-42-8	32



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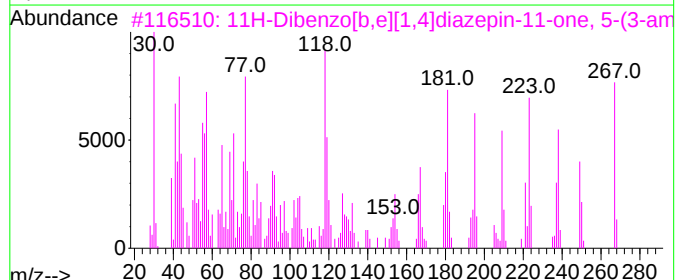
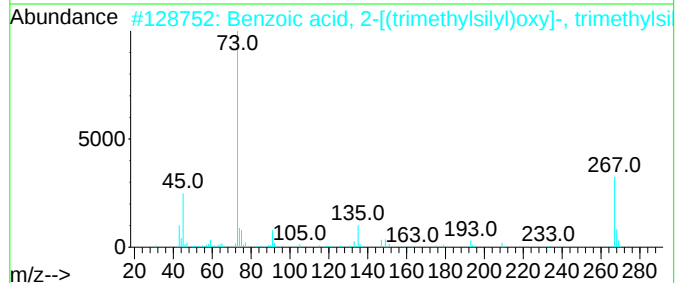
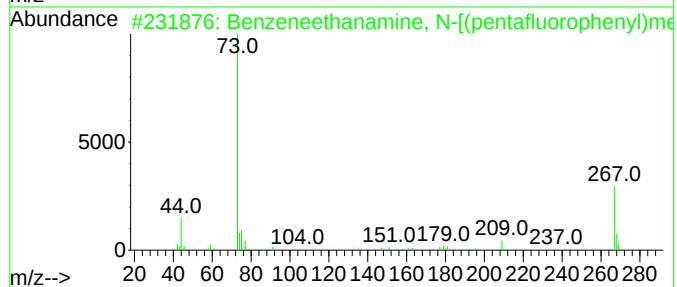
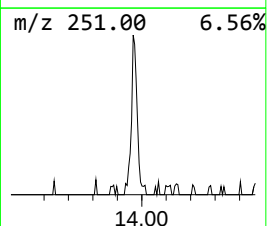
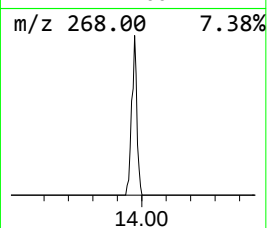
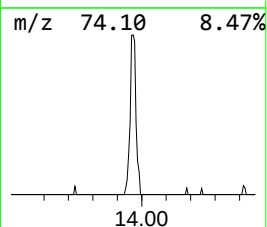
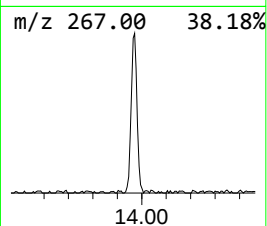
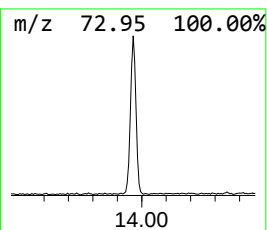
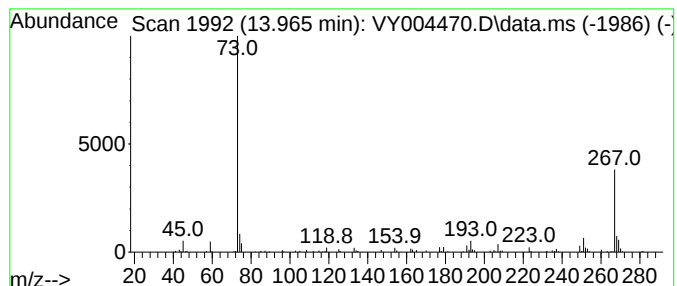
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Peak Number 2 Benzeneethanamine, N-[(pent... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.965	7.62 ug/l	68231	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5NO2Si2	055429-85-1	59
2			Benzoic acid, 2-[(trimethylsilyl)...	282	C13H22O3Si2	003789-85-3	45
3			11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	43
4			4-Methylcatechol, bis(trimethyls...	268	C13H24O2Si2	1000332-79-9	14
5			Monobenzylidene-d-glucose	268	C13H16O6	1000127-04-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.589	2.589	13.8	ug/l	122779	1	7.801	444649	50.0
Benzeneethanami...	13.965	7.6	ug/l	68231	4	13.428	447865	50.0