

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD052323\
 Data File : VD076199.D
 Acq On : 23 May 2023 18:00
 Operator : KP/SY
 Sample : 02864-05
 Misc : 5.05G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 ETGI-359

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_D\Method\82D051523S.M
 Title : SW846 8260

Signal : TIC: VD076199.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.734	18	25	38	rVV	303305	571568	100.00%	49.280%
2	4.799	534	546	558	rBV	25514	75117	13.14%	6.476%
3	7.063	920	931	943	rBV2	10225	30471	5.33%	2.627%
4	7.805	1047	1057	1063	rBV2	40818	95870	16.77%	8.266%
5	7.875	1063	1069	1079	rVB	42225	94922	16.61%	8.184%
6	8.228	1119	1129	1139	rVB2	29528	67112	11.74%	5.786%
7	8.775	1215	1222	1232	rBV	50183	99801	17.46%	8.605%
8	10.269	1468	1476	1483	rBV	35430	62260	10.89%	5.368%
9	10.663	1537	1543	1549	rVB3	3834	6118	1.07%	0.527%
10	11.028	1601	1605	1611	rBV2	5548	8362	1.46%	0.721%
11	11.581	1692	1699	1706	rBV	18076	30836	5.39%	2.659%
12	12.575	1863	1868	1872	rVB5	6991	9738	1.70%	0.840%
13	13.516	2023	2028	2034	rVB2	4885	7673	1.34%	0.662%

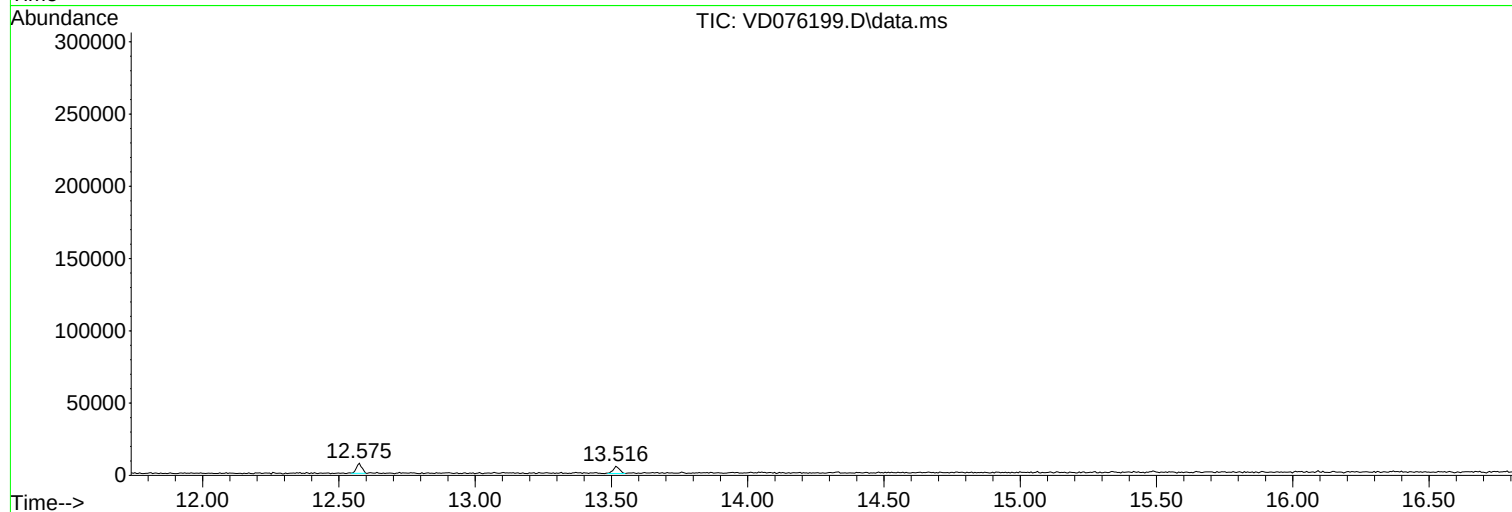
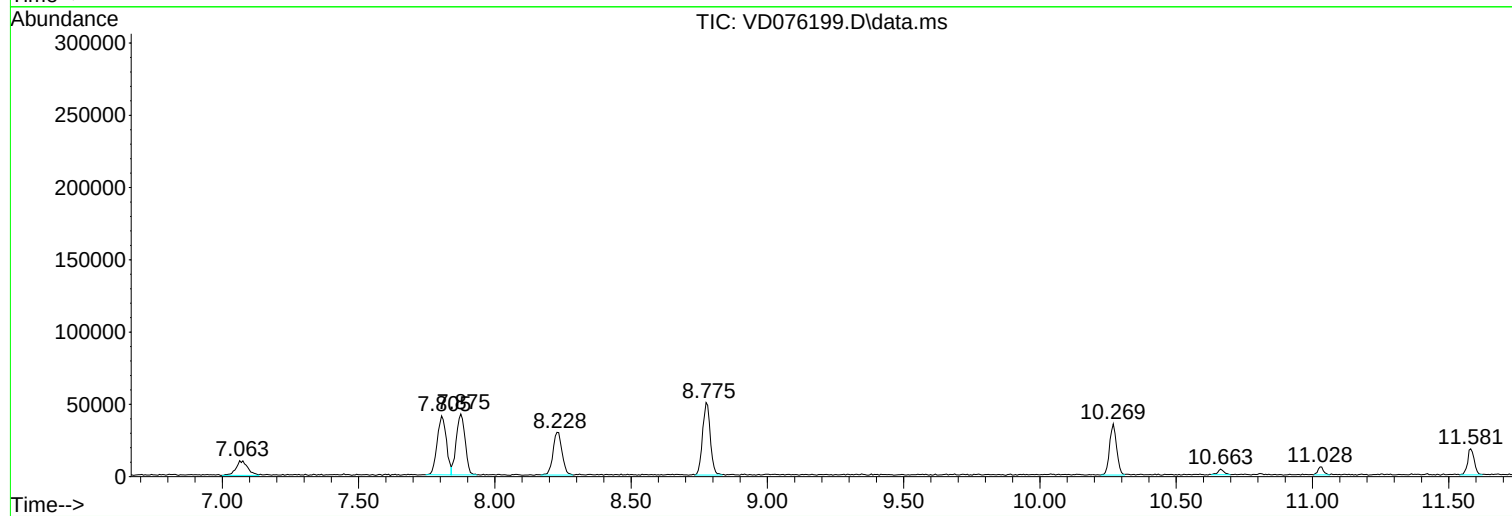
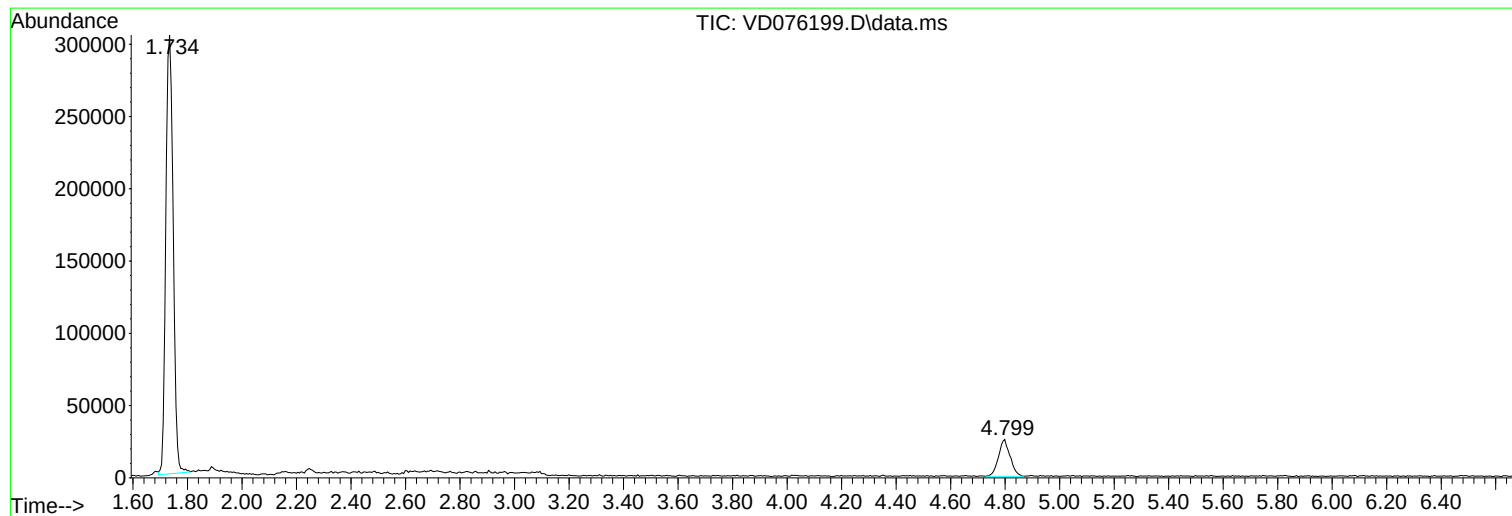
Sum of corrected areas: 1159848

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D051523S.M
Quant Title : SW846 8260

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



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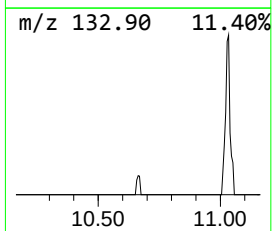
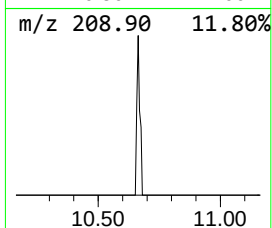
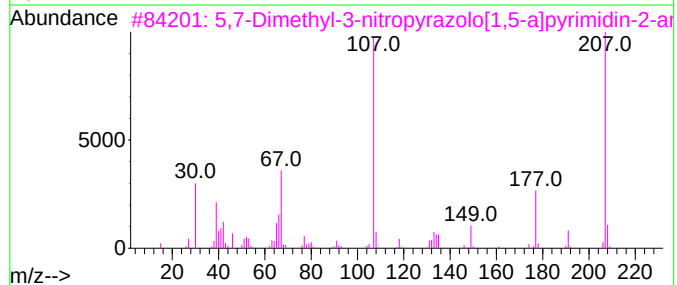
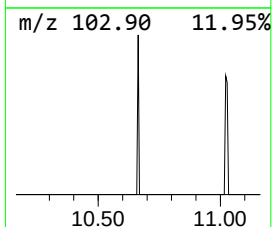
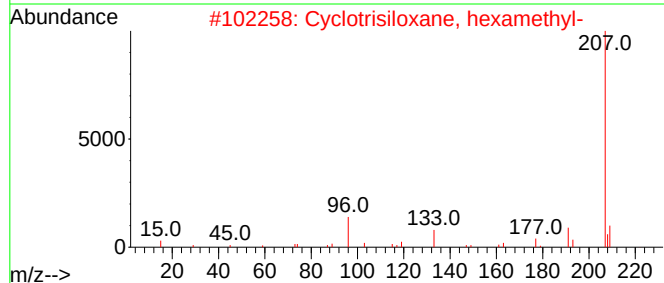
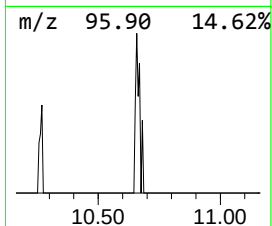
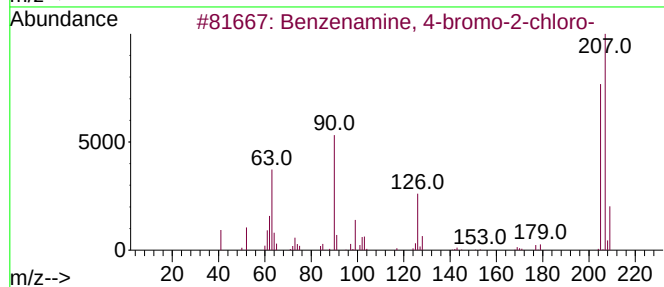
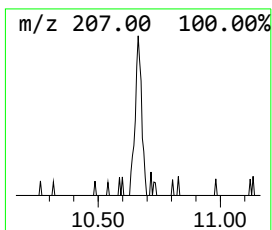
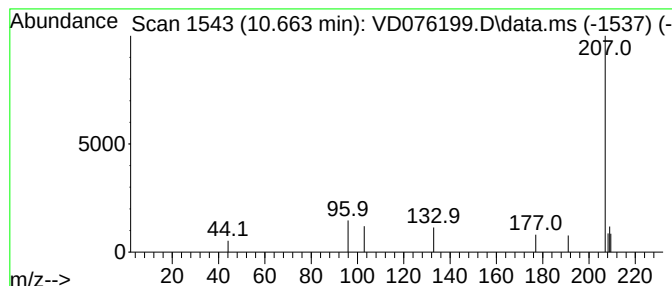
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Quant Title : SW846 8260

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Peak Number 3 Benzenamine, 4-bromo-2-chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.663	9.92 ug/l	6118	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 4-bromo-2-chloro-	205	C6H5BrClN	038762-41-3	56
2			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	39
3			5,7-Dimethyl-3-nitropyrzolo[1,5...	207	C8H9N5O2	1000489-27-6	36
4			2-Bromo-4-chloroaniline	205	C6H5BrClN	000873-38-1	9
5			Toloxatone	207	C11H13NO3	029218-27-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzenamine, 4-...	10.663	9.9	ug/l	6118	3	11.581	30836	50.0