

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091522\
Data File : VX031368.D
Acq On : 15 Sep 2022 20:53
Operator : JC/MD
Sample : N4508-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EB01-090222

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091422W.M

Title : SW846 8260

Signal : TIC: VX031368.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.618	77	87	98	rBV7	6857	28602	3.55%	0.858%
2	5.385	693	705	722	rBV	94312	273073	33.89%	8.193%
3	5.556	722	733	748	rVB2	94279	268483	33.32%	8.055%
4	5.958	789	799	812	rBV	105318	277824	34.48%	8.335%
5	6.763	922	931	941	rBV	150092	353615	43.89%	10.609%
6	8.653	1233	1241	1255	rBV	509311	805741	100.00%	24.174%
7	10.055	1465	1471	1481	rBV	318229	424111	52.64%	12.724%
8	11.085	1634	1640	1652	rBV	393357	516605	64.12%	15.499%
9	12.024	1788	1794	1802	rBV	318820	385049	47.79%	11.552%

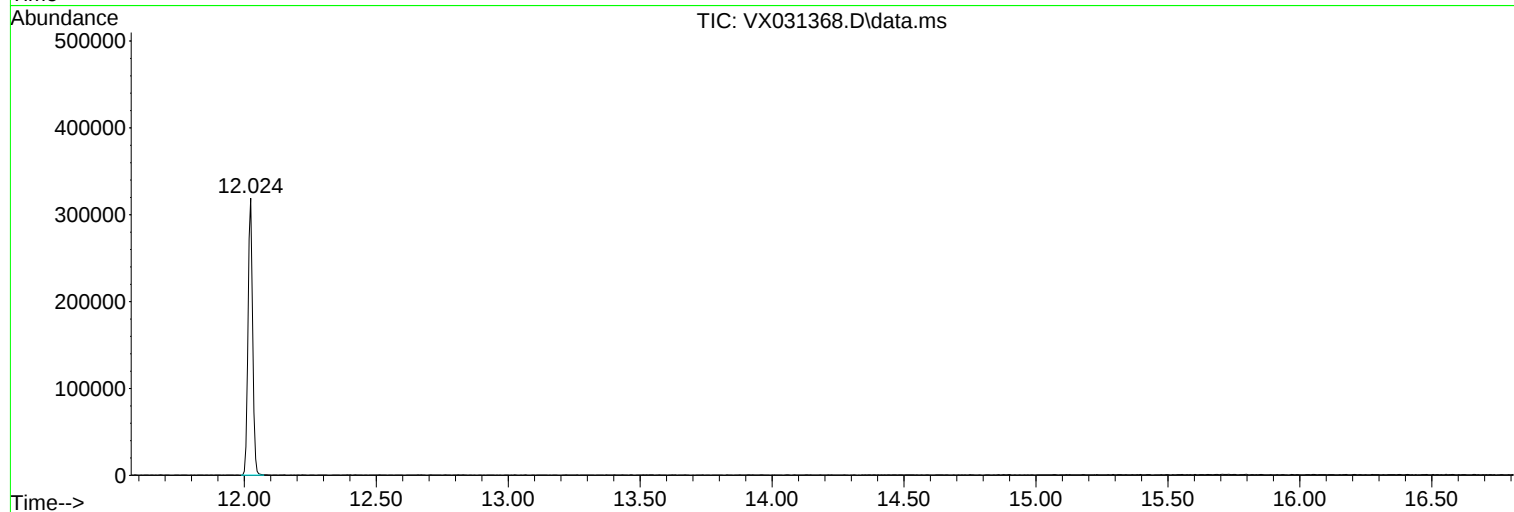
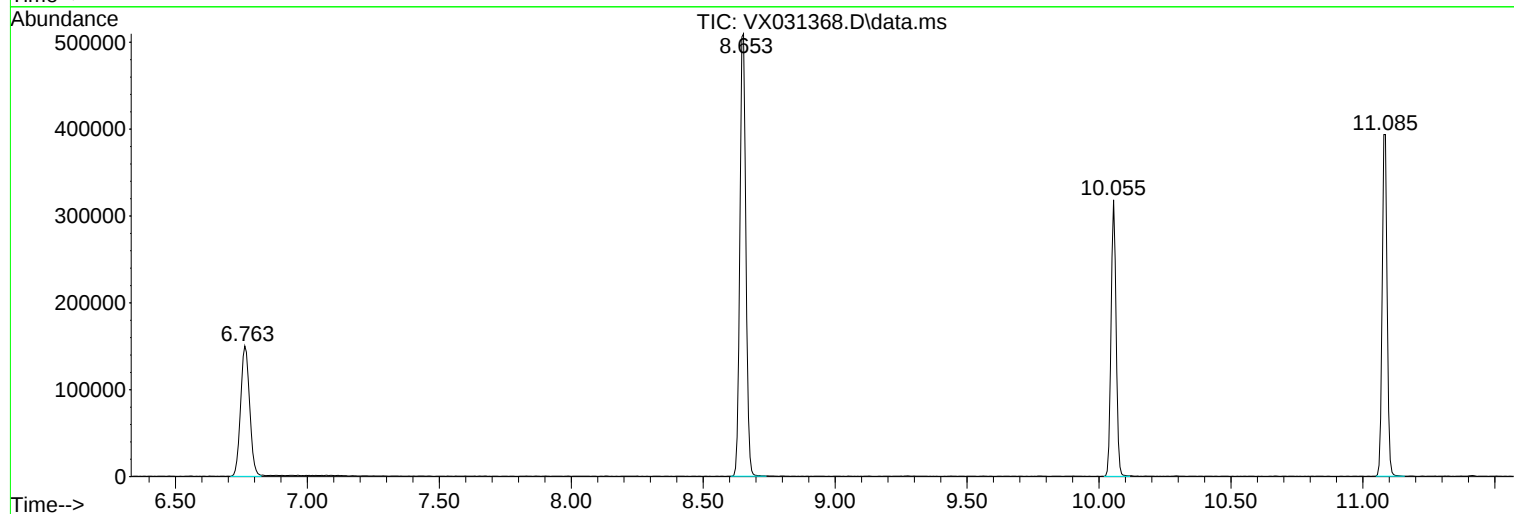
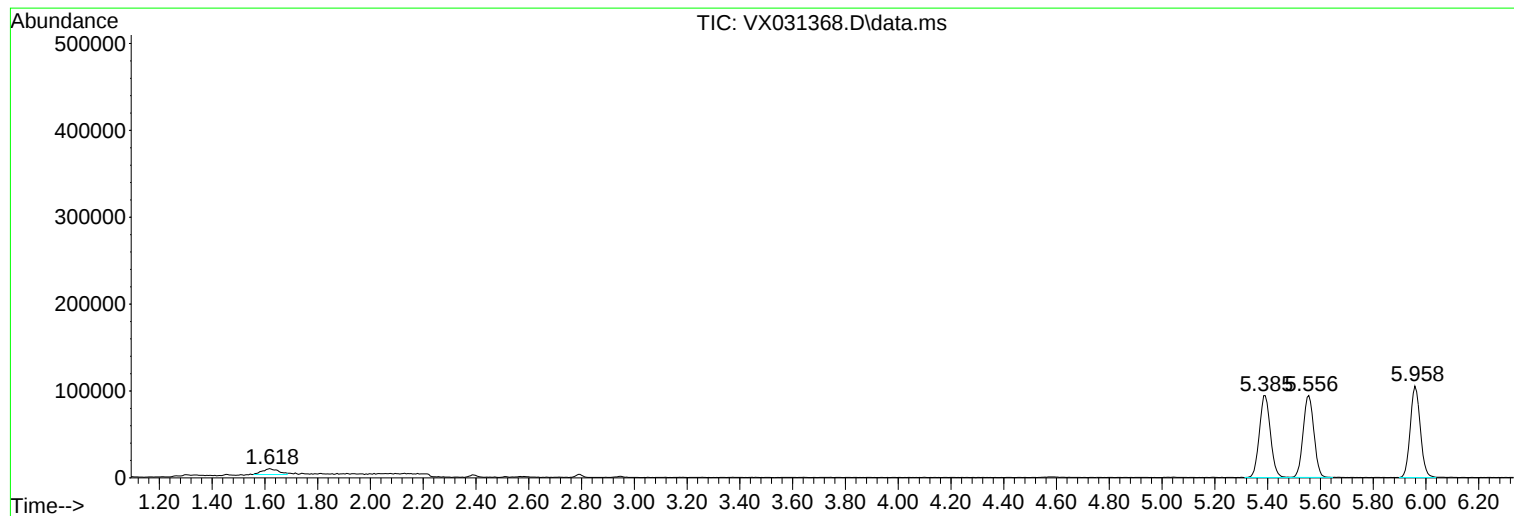
Sum of corrected areas: 3333103

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091422W.M
Quant Title : SW846 8260

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



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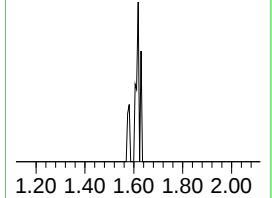
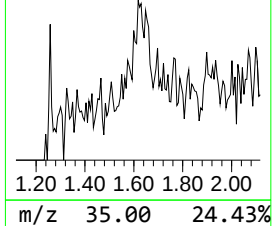
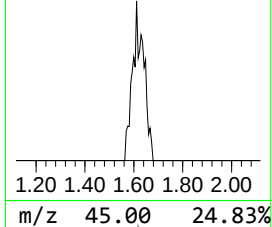
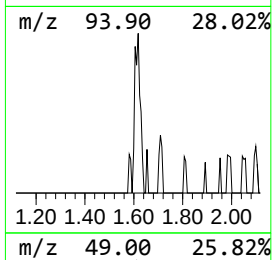
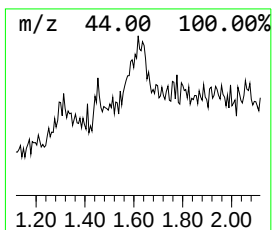
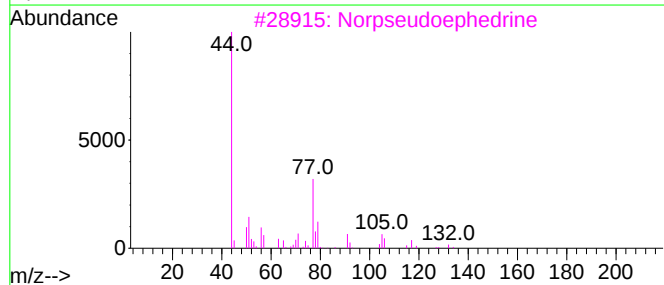
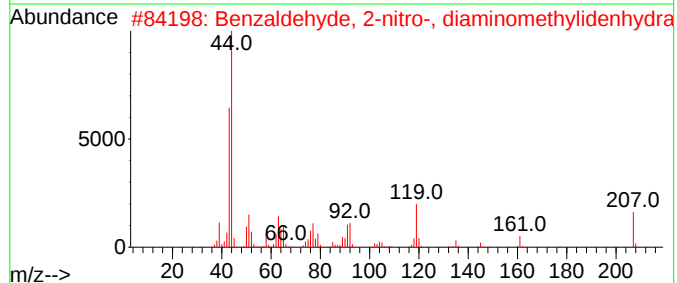
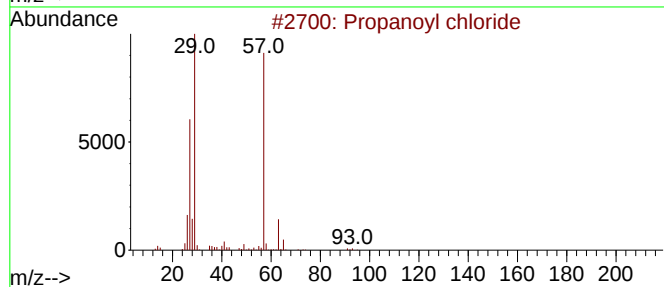
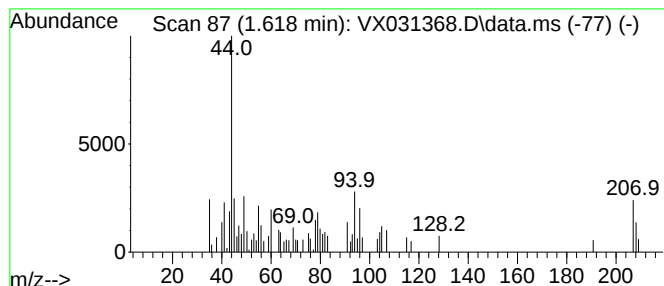
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091422W.M
Quant Title : SW846 8260

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

Peak Number 1 unknown1.618 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.618	5.33 ug/l	28602	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoyl chloride	92	C3H5ClO	000079-03-8	38
2			Benzaldehyde, 2-nitro-, diaminom...	207	C8H9N5O2	102632-31-5	25
3			Norpseudoephedrine	151	C9H13NO	036393-56-3	14
4			1,3-Bis(2-chloroethyl)urea	184	C5H10Cl2N2O	002214-72-4	13
5			Phosphonoacetic acid	140	C2H5O5P	004408-78-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.618	1.618	5.3	ug/l	28602	1	5.556	268483	50.0