

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112224\
 Data File : VX043964.D
 Acq On : 22 Nov 2024 16:52
 Operator : JC/MD
 Sample : P4846-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW-15-20241113

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_X\Method\82X112124W.M
 Title : SW846 8260

Signal : TIC: VX043964.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.240	22	25	39	rVB2	8803	14262	1.14%	0.286%
2	1.611	67	86	87	rBV5	16513	68165	5.47%	1.366%
3	5.385	693	705	721	rBV	87558	270381	21.69%	5.417%
4	5.544	721	731	745	rBV	124812	359858	28.87%	7.210%
5	5.946	789	797	812	rBV	101415	270468	21.70%	5.419%
6	6.757	921	930	946	rBV	210218	522450	41.92%	10.468%
7	7.117	980	989	1005	rBV	554799	1246312	100.00%	24.971%
8	8.647	1233	1240	1253	rBV	431444	706812	56.71%	14.162%
9	10.049	1465	1470	1483	rBV	424273	605003	48.54%	12.122%
10	11.079	1634	1639	1650	rBV	325301	428001	34.34%	8.575%
11	12.018	1788	1793	1804	rBV	387574	499321	40.06%	10.004%

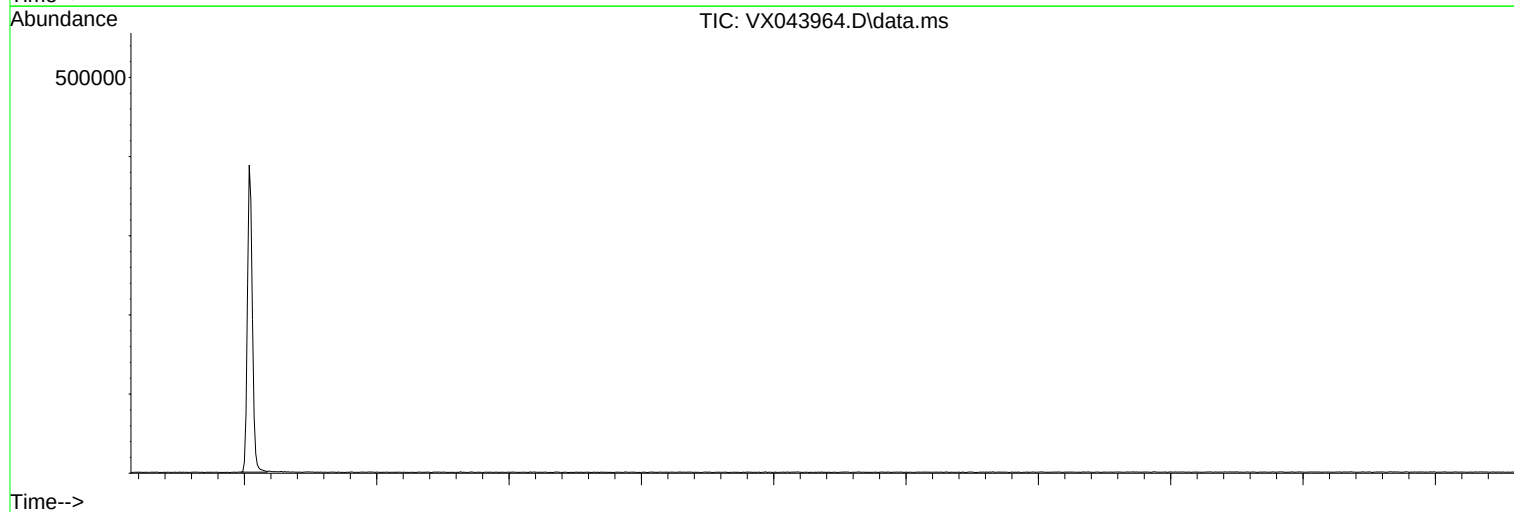
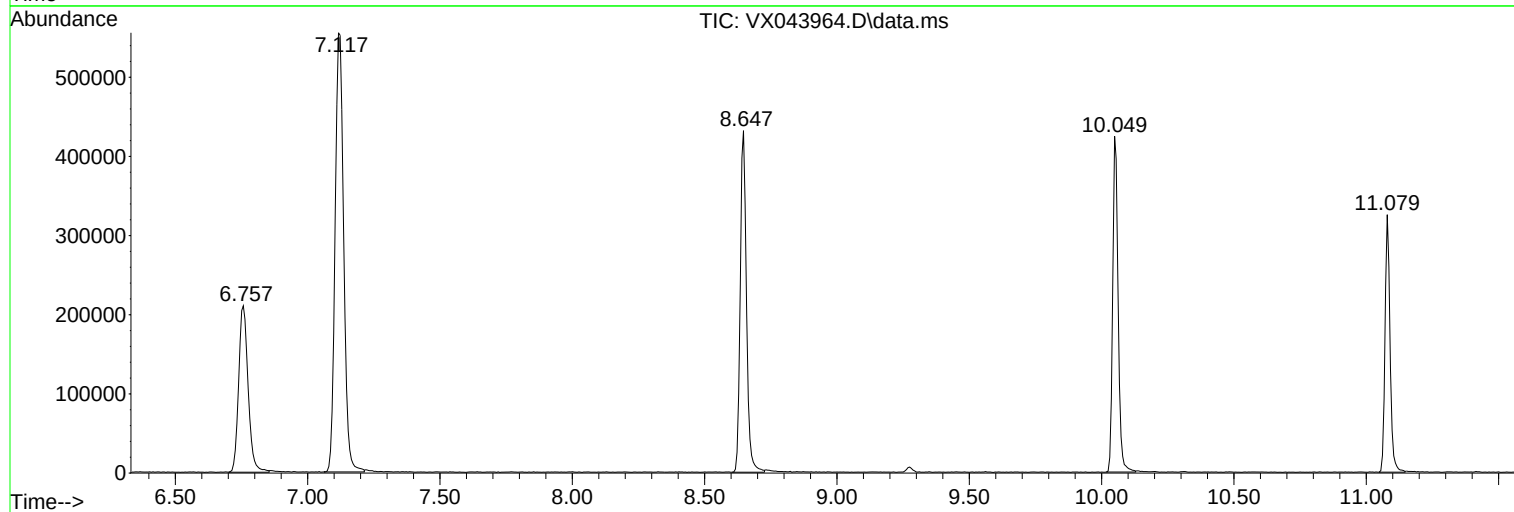
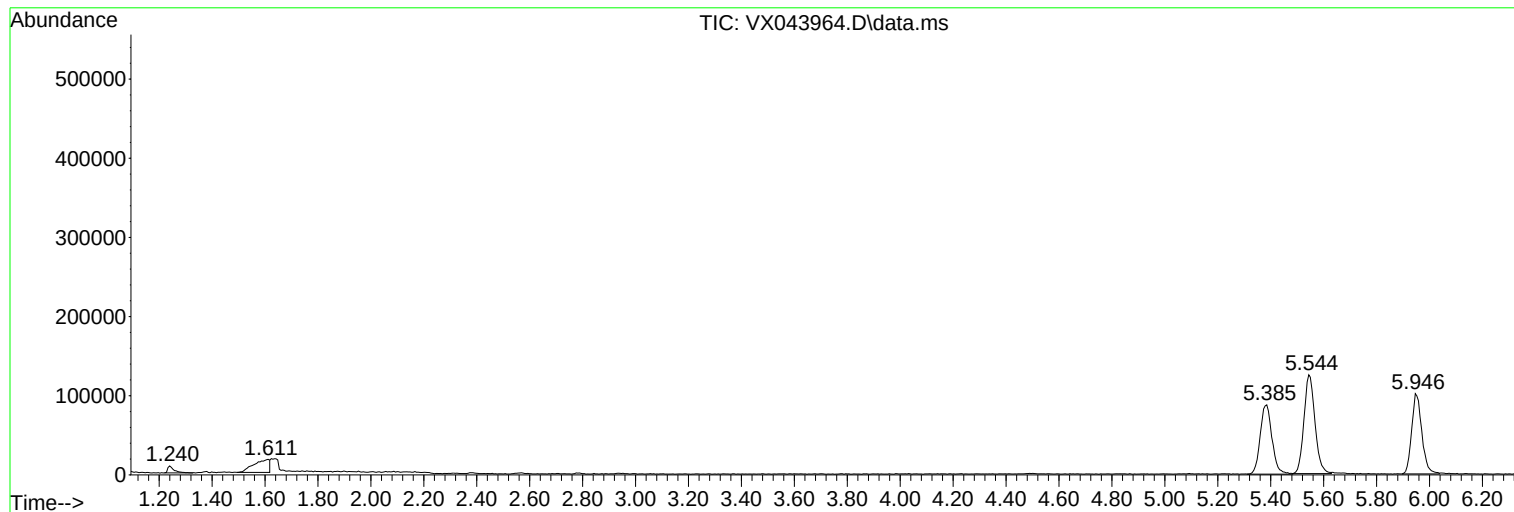
Sum of corrected areas: 4991033

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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



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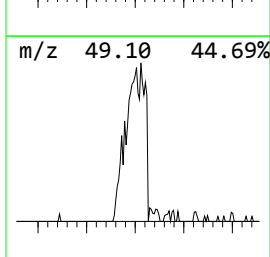
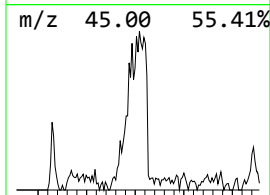
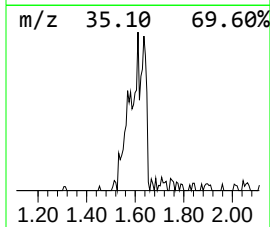
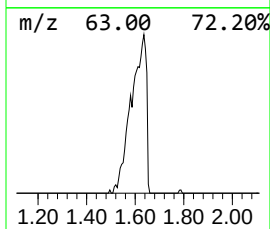
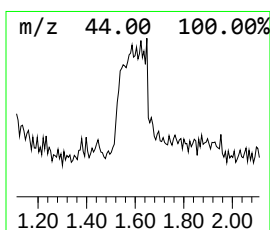
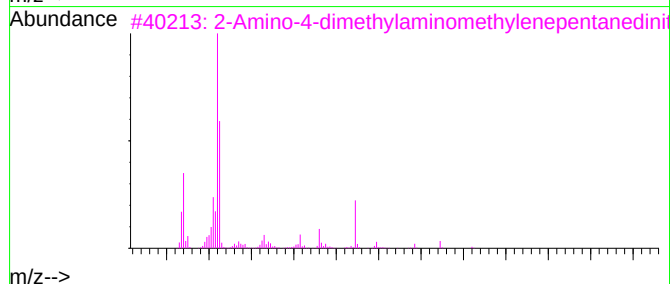
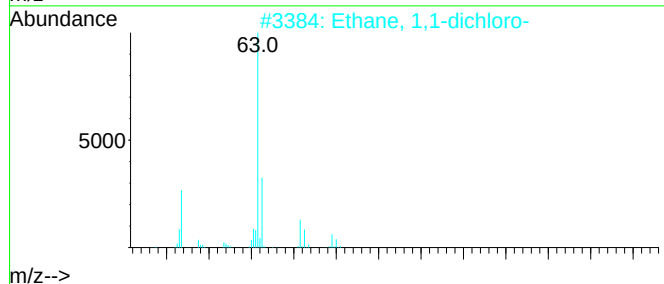
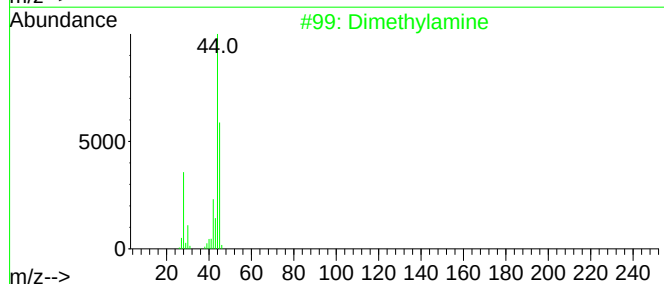
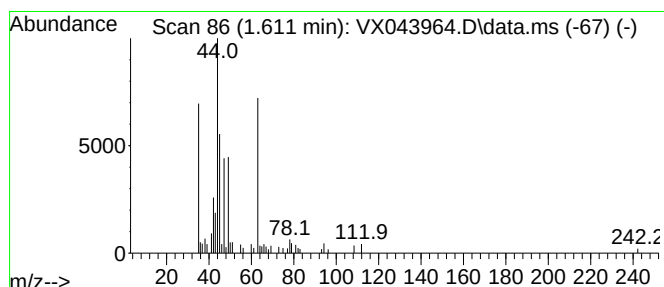
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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

Peak Number 1 unknown1.611 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.611	9.47 ug/l	68165	Pentafluorobenzene	5.544

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dimethylamine	45	C2H7N	000124-40-3	10
2	Ethane, 1,1-dichloro-	98	C2H4Cl2	000075-34-3	9
3	2-Amino-4-dimethylaminomethylene...	164	C8H12N4	1000191-40-0	9
4	2-(Oxan-3-yl)ethanamine	129	C7H15NO	1000411-07-5	9
5	Dihydroxymaleic acid	148	C4H4O6	000526-84-1	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.611	1.611	9.5	ug/l	68165	1	5.544	359858	50.0