

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101520\
 Data File : VN064148.D
 Acq On : 15 Oct 2020 17:07
 Operator : JC/MD
 Sample : L4417-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-7

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_N\METHODS\82N101420W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.428	198	214	224	rBV3	50233	128920	7.10%	1.332%
2	7.550	1786	1807	1819	rBV2	243722	616181	33.93%	6.368%
3	7.627	1819	1831	1858	rVB3	359786	864815	47.63%	8.938%
4	7.988	1925	1943	1963	rBV	259884	607824	33.47%	6.282%
5	8.550	2100	2118	2142	rBV2	599988	1271723	70.04%	13.143%
6	10.061	2573	2588	2603	rBV	982911	1815836	100.00%	18.767%
7	10.605	2745	2757	2772	rVB2	57083	105324	5.80%	1.089%
8	11.380	2984	2998	3021	rBV	970085	1719476	94.69%	17.771%
9	12.376	3295	3308	3328	rBV	708923	1211617	66.73%	12.522%
10	13.319	3588	3601	3626	rBV	772148	1334227	73.48%	13.789%

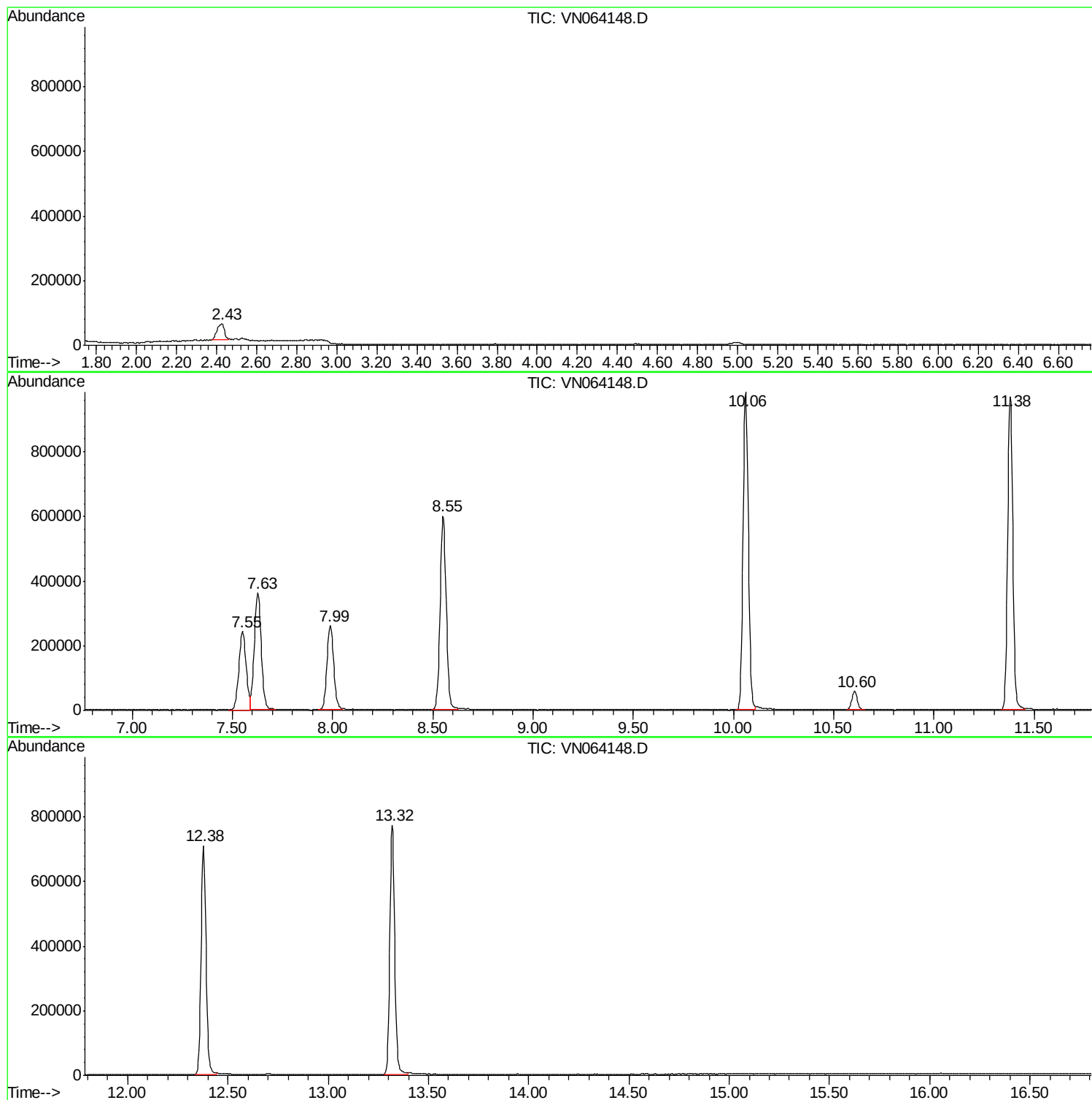
Sum of corrected areas: 9675943

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101420W.M
Quant Title : SW846 8260

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



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ClientSampled :
MW-7

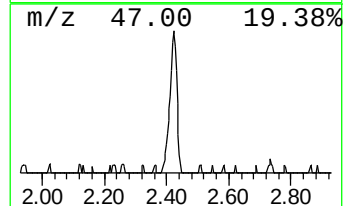
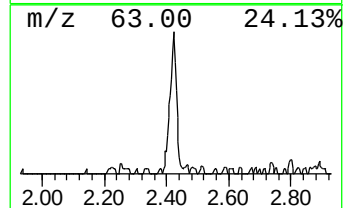
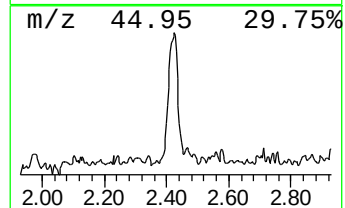
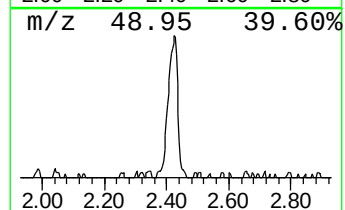
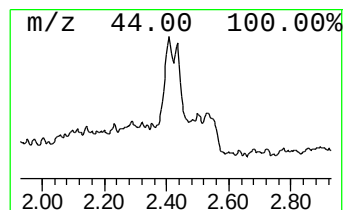
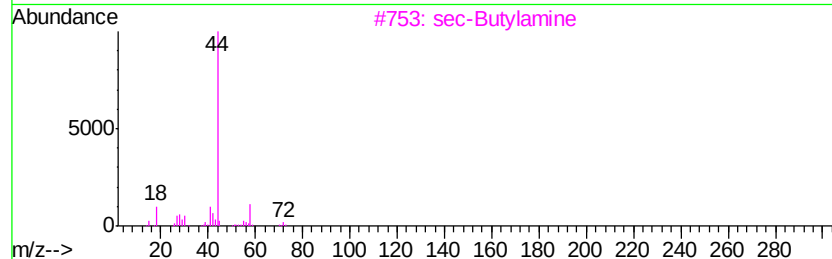
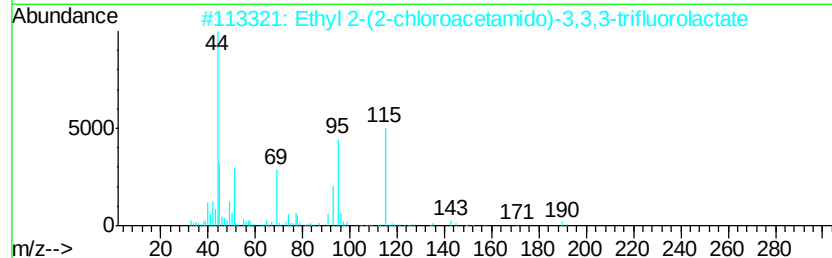
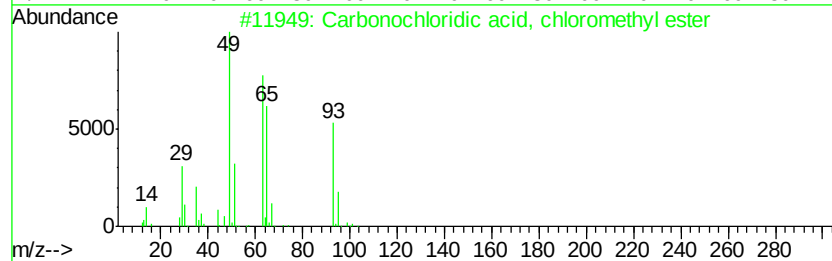
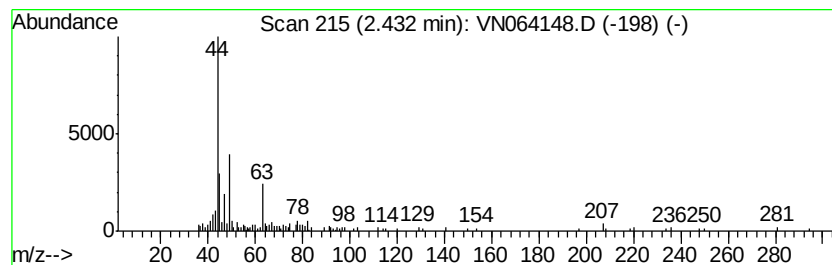
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Quant Title : SW846 8260

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

Peak Number 1 unknown2.43 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.43	7.45 ug/l	128920	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbonochloridic acid, chloromet...	128	C2H2Cl2O2	022128-62-7	17
2		Ethyl 2-(2-chloroacetamido)-3,3,...	263	C7H9ClF3N04	1000224-74-2	14
3		sec-Butylamine	73	C4H11N	013952-84-6	10
4		2-Butanamine, (S)-	73	C4H11N	000513-49-5	10
5		1-Propanamine, N,2-dimethyl-	87	C5H13N	000625-43-4	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.43	2.43	7.5	ug/l	128920	1	7.63	864815	50.0