

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX072318\
 Data File : VX003612.D
 Acq On : 23 Jul 2018 18:43
 Operator : JC/SP
 Sample : J4048-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 YT-MW-6

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\82X071918W.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	1.075	76	80	94	rBV	1776543	2248439	100.00%	24.267%
2	1.301	114	117	121	rBV	95787	86752	3.86%	0.936%
3	1.386	121	131	132	rBV2	20883	59282	2.64%	0.640%
4	2.624	328	334	341	rVB	19548	36455	1.62%	0.393%
5	4.605	647	659	673	rBV2	56187	166904	7.42%	1.801%
6	5.507	795	807	820	rBV	160398	450661	20.04%	4.864%
7	5.672	822	834	850	rVV	231472	647539	28.80%	6.989%
8	6.074	887	900	917	rBV	159833	419676	18.67%	4.529%
9	6.867	1018	1030	1045	rBV	343329	801230	35.63%	8.647%
10	7.220	1079	1088	1095	rBV	22033	48216	2.14%	0.520%
11	8.720	1327	1334	1342	rBV	840375	1294071	57.55%	13.967%
12	9.336	1429	1435	1446	rVB	112204	167002	7.43%	1.802%
13	10.116	1557	1563	1575	rBV	719034	953191	42.39%	10.288%
14	11.140	1725	1731	1743	rBV	699028	859749	38.24%	9.279%
15	12.079	1880	1885	1893	rBV	817850	1026323	45.65%	11.077%

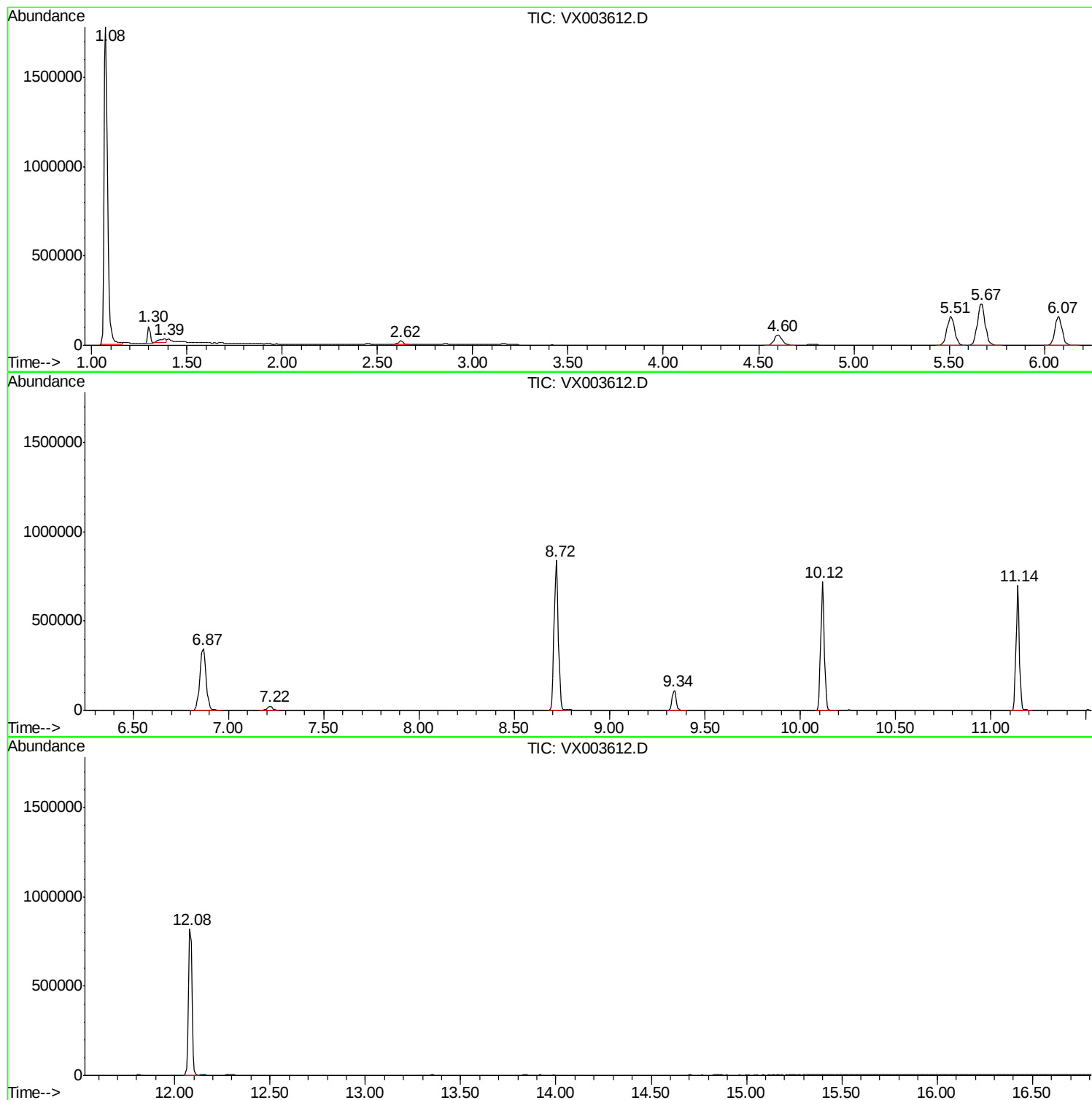
Sum of corrected areas: 9265490

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071918W.M
Quant Title : SW846 8260

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



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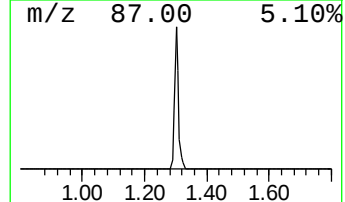
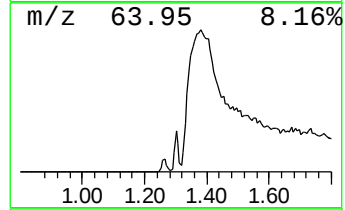
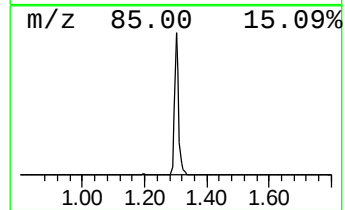
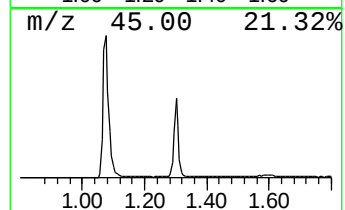
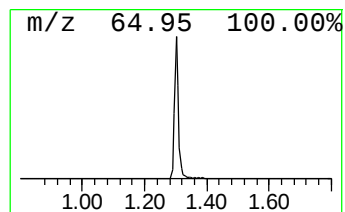
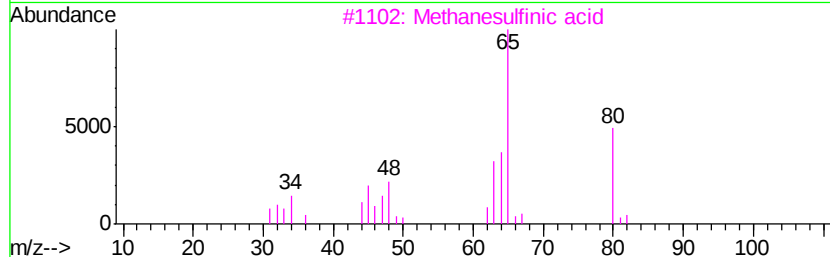
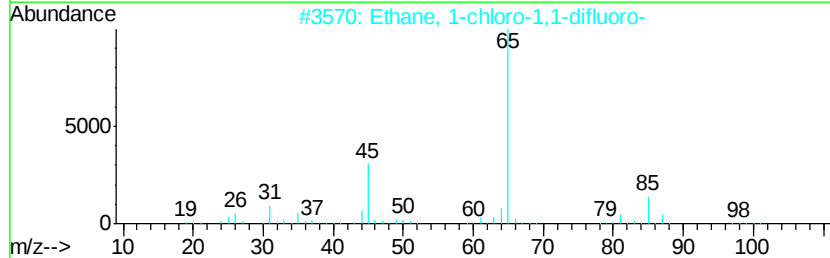
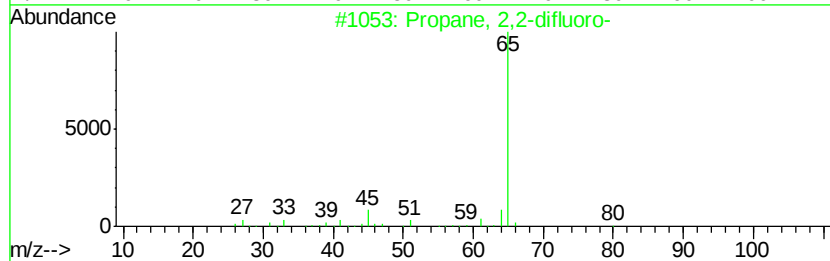
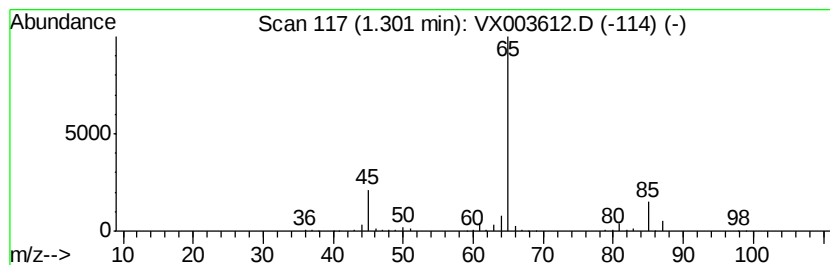
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Peak Number 2 unknown1.30 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.30	6.70 ug/l	86752	Pentafluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	43
2			Ethane, 1-chloro-1,1-difluoro-	100	C2H3ClF2	000075-68-3	37
3			Methanesulfinic acid	80	CH4O2S	017696-73-0	4
4			Benzene, 1-nitro-2-(p-methylphen...	247	C13H10FN03	028987-57-7	1
5			Propanoyl chloride	92	C3H5ClO	000079-03-8	1



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.30	1.30	6.7	ug/l	86752	1	5.67	647539	50.0