

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112118\
 Data File : VX006118.D
 Acq On : 21 Nov 2018 12:33
 Operator : JC/MD
 Sample : J5967-06
 Misc : 5.0g/10ml/100ul/5.00mL/MSVOA_X/MEOH
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FDSB-13(16-18)

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\82X112018W.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	1.672	63	85	87	rBV10	60237	265813	16.10%	2.849%
2	5.501	702	713	727	rBV	181587	548623	33.24%	5.881%
3	5.653	727	738	752	rVB	333373	951957	57.67%	10.204%
4	6.062	795	805	823	rBV	210449	581567	35.23%	6.234%
5	6.854	923	935	947	rBV	489759	1197680	72.56%	12.838%
6	8.714	1233	1240	1256	rBV	1020994	1650554	100.00%	17.692%
7	10.116	1464	1470	1482	rBV	1017876	1448700	87.77%	15.528%
8	11.140	1633	1638	1650	rBV	887127	1166647	70.68%	12.505%
9	12.079	1787	1792	1802	rBV	1214679	1517854	91.96%	16.270%

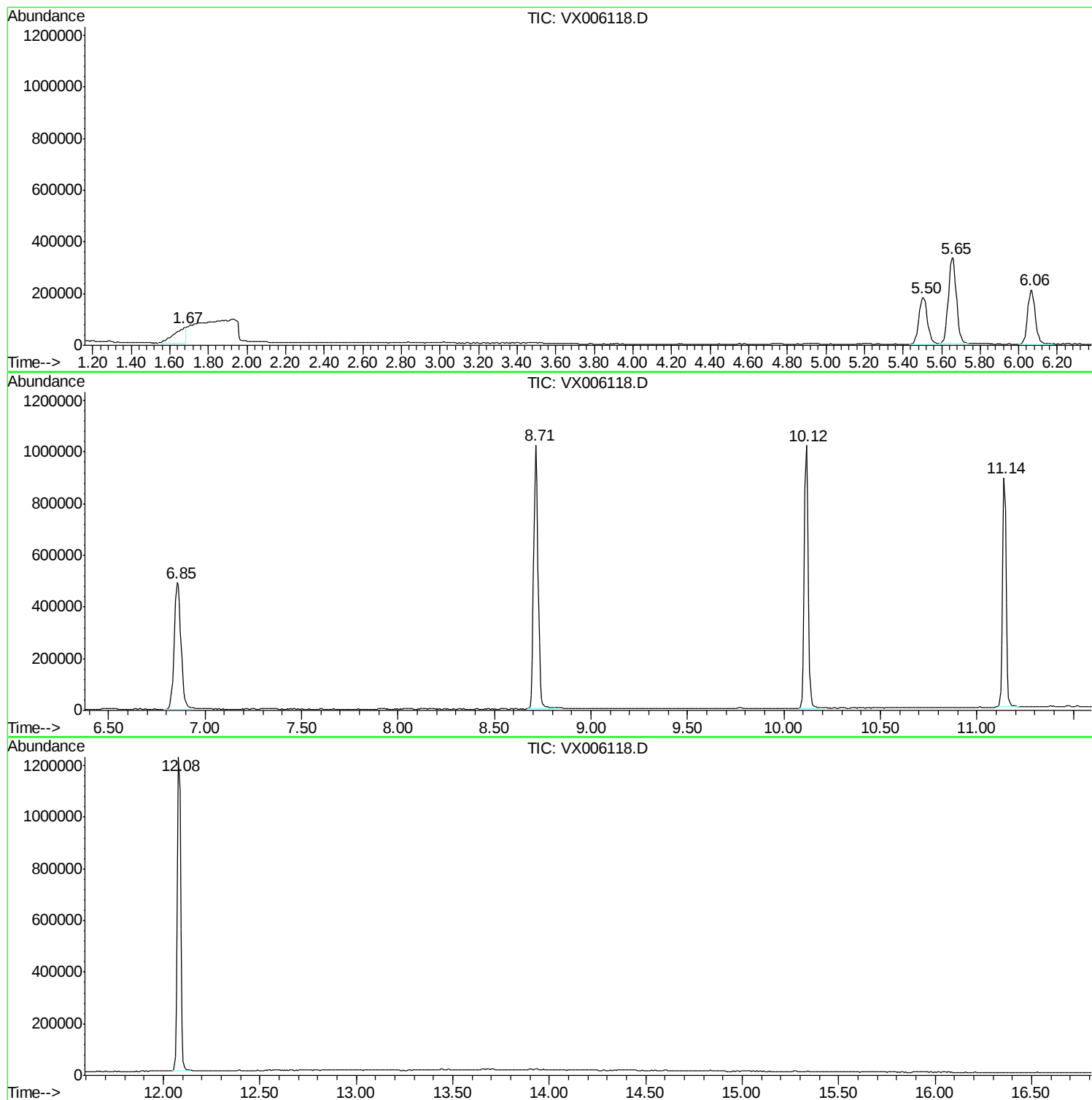
Sum of corrected areas: 9329395

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112018W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampled :
FDSB-13(16-18)

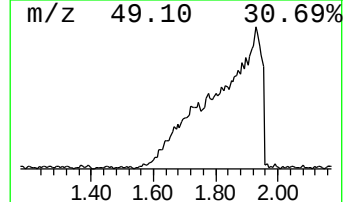
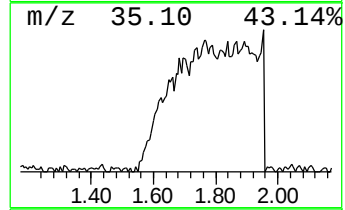
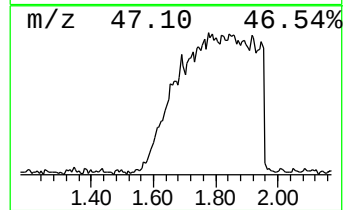
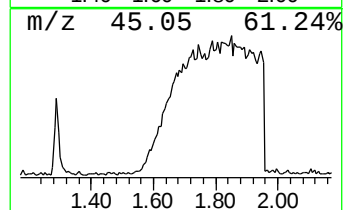
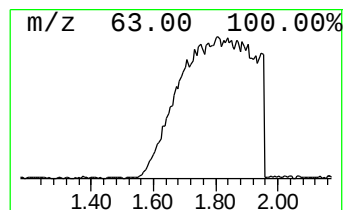
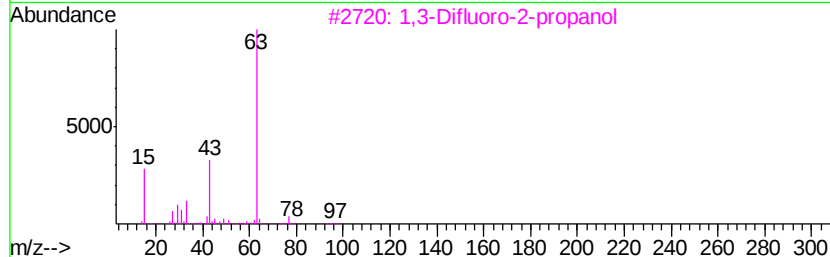
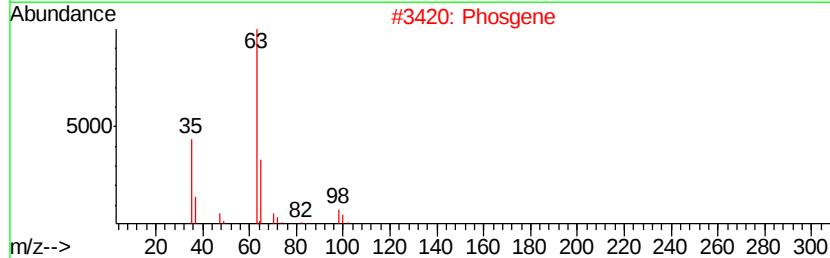
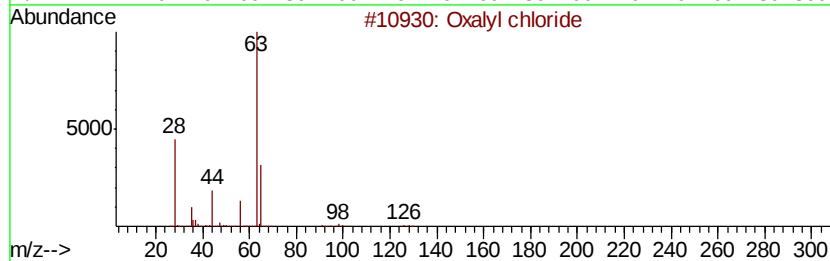
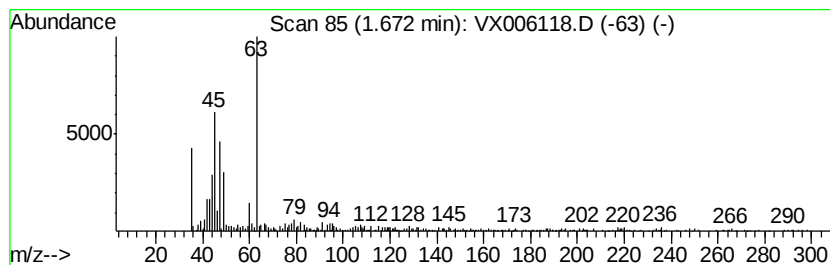
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Quant Title : SW846 8260

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

Peak Number 1 unknown1.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.67	13.96 ug/l	265813	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxalyl chloride	126	C2Cl2O2	000079-37-8	42
2		Phosgene	98	CCl2O	000075-44-5	36
3		1,3-Difluoro-2-propanol	96	C3H6F2O	000453-13-4	17
4		Bis(2-chloroethyl) sulfone	190	C4H8Cl2O2S	000471-03-4	17
5		1-Chloroethyl sulfone	190	C4H8Cl2O2S	010038-07-0	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.67	1.67	14.0	ug/l	265813	1	5.66	951957	50.0