

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX012920\
 Data File : VX014765.D
 Acq On : 29 Jan 2020 15:45
 Operator : JC/SP
 Sample : VX0129MBL02
 Misc : 5.0g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX0129MBL02

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\82X012220W.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	1.070	26	30	42	rBV	378791	484120	22.87%	4.066%
2	1.283	63	65	72	rVB	20336	22314	1.05%	0.187%
3	1.637	103	123	124	rBV	57815	216551	10.23%	1.819%
4	5.490	745	755	771	rBV	198770	665474	31.44%	5.589%
5	5.648	771	781	796	rVB2	393658	1181558	55.82%	9.924%
6	6.051	835	847	863	rBV2	238715	713279	33.69%	5.991%
7	6.849	968	978	991	rBV	589775	1536846	72.60%	12.908%
8	8.709	1276	1283	1291	rBV	1297374	2116912	100.00%	17.780%
9	10.111	1503	1513	1521	rBV	1264411	1827168	86.31%	15.347%
10	11.135	1675	1681	1692	rBV	1105806	1414105	66.80%	11.877%
11	12.074	1830	1835	1843	rBV	1420870	1727702	81.61%	14.511%

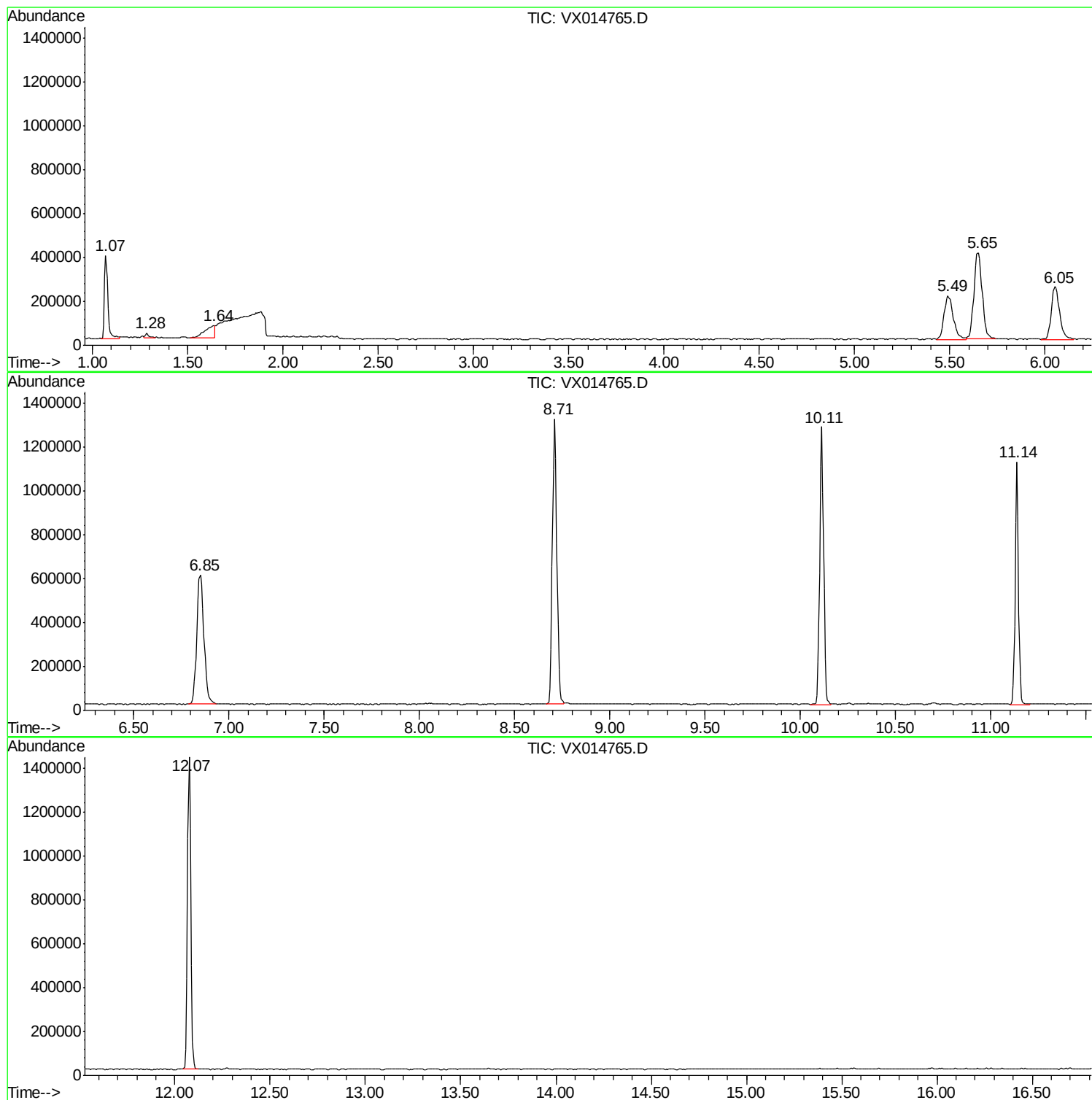
Sum of corrected areas: 11906029

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X012220W.M
Quant Title : SW846 8260

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



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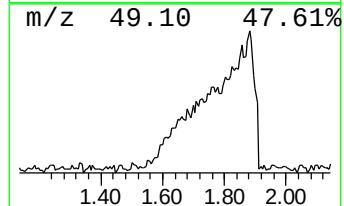
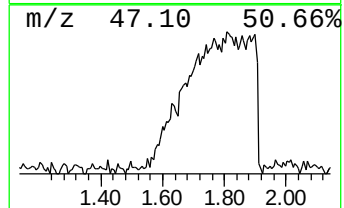
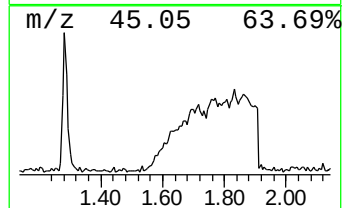
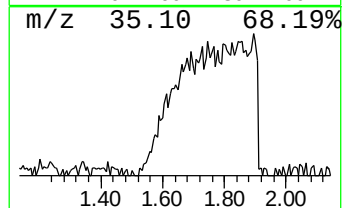
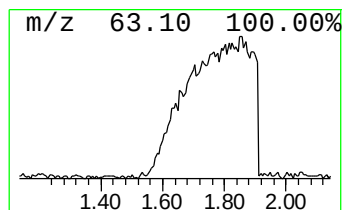
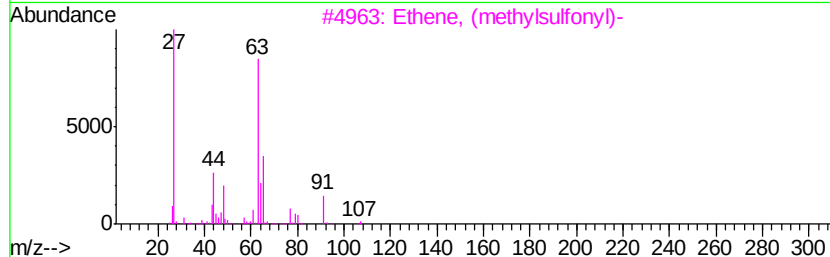
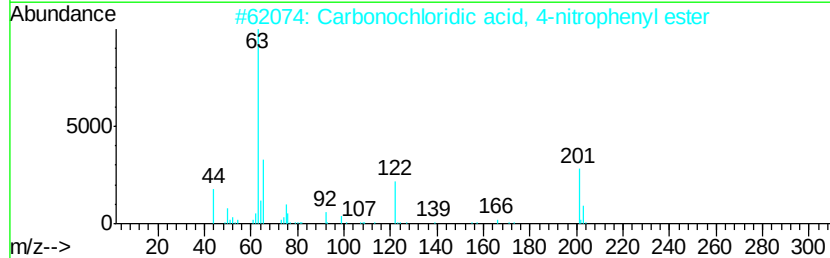
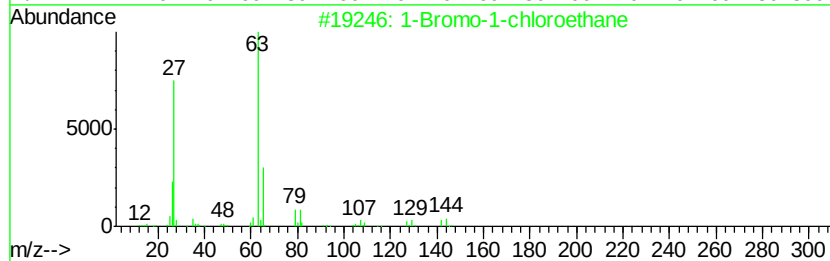
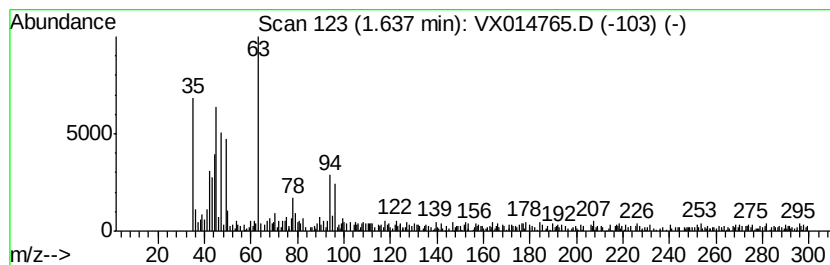
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X012220W.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown1.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
1.64	9.16 ug/l	216551	Pentafluorobenzene	5.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Bromo-1-chloroethane	142	C2H4BrCl	000593-96-4	30
2		Carbonochloridic acid, 4-nitroph...	201	C7H4ClN04	007693-46-1	20
3		Ethene, (methylsulfonyl)-	106	C3H6O2S	003680-02-2	17
4		Acetaldehyde, bis(2-chloroethyl)...	186	C6H12Cl2O2	014689-97-5	17
5		Benzaldehyde, 5-bromo-2-hydroxy-...	359	C13H9BrF3N3O	1000276-33-1	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.64	1.64	9.2	ug/l	216551	1	5.65	1181560	50.0