

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010519\
 Data File : VU029041.D
 Acq On : 04 Jan 2019 20:43
 Operator : MD/SY
 Sample : K1053-07 10X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 COMP-3-DRUM

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_U\METHOD\82U010319W.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	1.283	27	34	44	rBV	183993	175065	28.81%	5.200%
2	1.472	85	93	108	rBV	78240	88061	14.49%	2.616%
3	1.569	112	123	133	rVB8	6699	14520	2.39%	0.431%
4	2.013	252	261	278	rVB	37082	54816	9.02%	1.628%
5	4.267	951	962	992	rBV2	12762	32306	5.32%	0.960%
6	4.881	1138	1153	1170	rBV	92715	206640	34.00%	6.138%
7	4.980	1170	1184	1205	rVB2	136088	299176	49.23%	8.887%
8	5.305	1271	1285	1307	rBV	97451	206842	34.04%	6.144%
9	5.884	1450	1465	1490	rBV	218202	429534	70.68%	12.759%
10	7.562	1974	1987	2025	rBV	341926	607693	100.00%	18.051%
11	9.090	2449	2462	2486	rBV	295985	501808	82.58%	14.905%
12	10.305	2828	2840	2866	rBV	220061	350575	57.69%	10.413%
13	11.482	3195	3206	3228	rBV	250425	399588	65.75%	11.869%

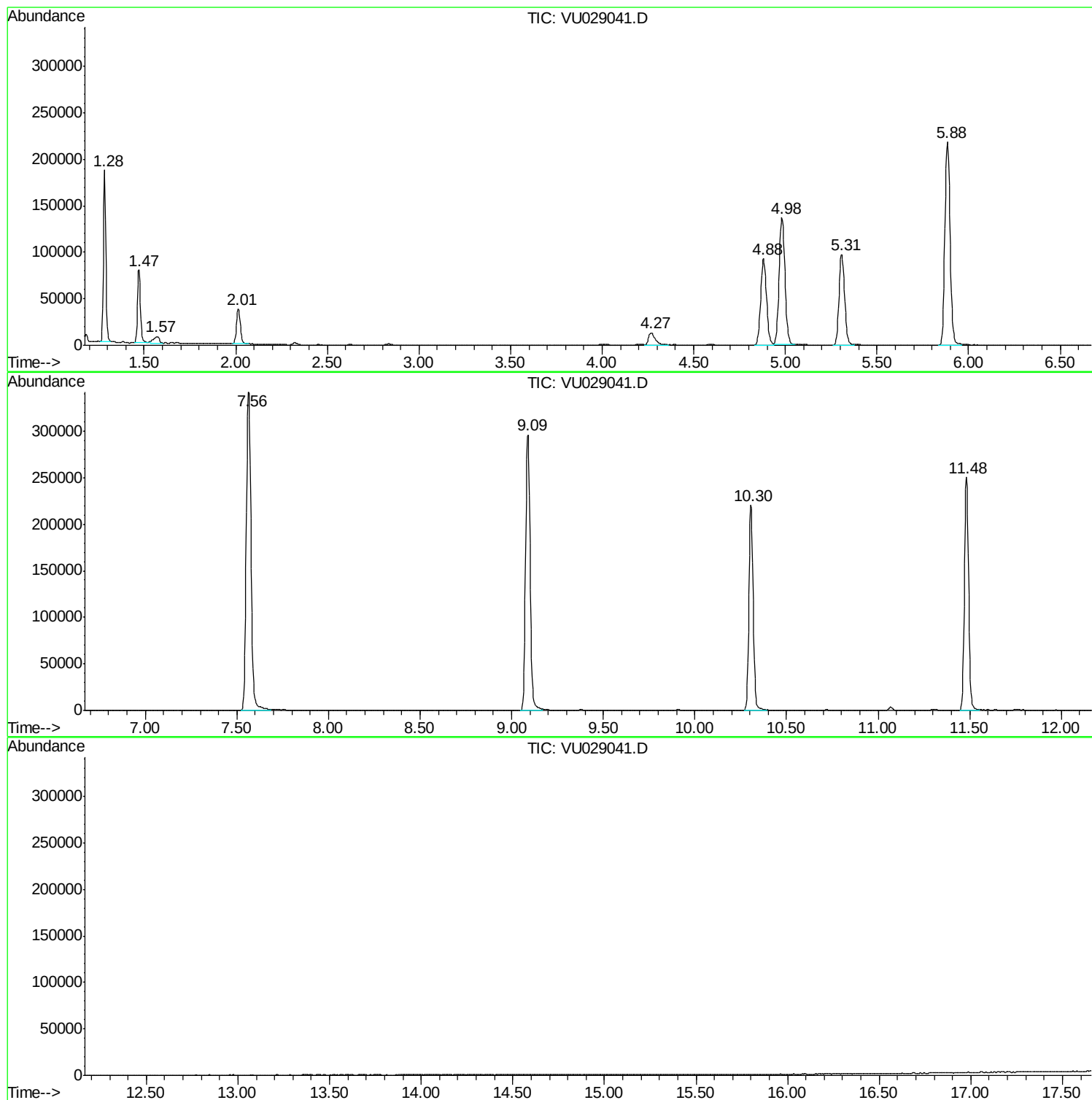
Sum of corrected areas: 3366624

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
Quant Title : SW846 8260

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



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

Peak Number 1 Acetaldehyde Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.47	14.72 ug/l	88061	Pentafluorobenzene	4.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Acetaldehyde	44	C2H4O	000075-07-0 74
2	Ethylene oxide	44	C2H4O	000075-21-8 5
3	Propane	44	C3H8	000074-98-6 4
4	Alanine	89	C3H7NO2	000056-41-7 4
5	Carbon dioxide	44	CO2	000124-38-9 3

Peak Number 2 Ethanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.01	9.16 ug/l	54816	Pentafluorobenzene	4.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Ethanol	46	C2H6O	000064-17-5 74
2	Dimethyl ether	46	C2H6O	000115-10-6 9
3	Methane, nitroso-	45	CH3NO	000865-40-7 4
4	Formic acid	46	CH2O2	000064-18-6 4
5	Oxalic acid	90	C2H2O4	000144-62-7 4

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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Acetaldehyde	1.47	14.7	ug/l	88061	1	4.98	299176	50.0
Ethanol	2.01	9.2	ug/l	54816	1	4.98	299176	50.0