

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU040819\
 Data File : VU030878.D
 Acq On : 08 Apr 2019 00:12
 Operator : JC/SP
 Sample : K2284-11
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BEZA9

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR032919WMA.M
 Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.396	88	95	105	rBV	171566	179947	13.94%	2.092%
2	1.679	175	183	198	rVB	128641	165442	12.82%	1.924%
3	2.270	356	367	378	rBV	269806	417845	32.37%	4.858%
4	4.186	949	963	992	rBV	180387	508686	39.41%	5.914%
5	4.392	1015	1027	1058	rVB2	50476	128081	9.92%	1.489%
6	4.643	1088	1105	1130	rBV3	164487	447381	34.66%	5.202%
7	5.331	1296	1319	1340	rBV2	371214	1039375	80.52%	12.084%
8	5.881	1477	1490	1509	rBV	285430	572257	44.33%	6.653%
9	6.328	1612	1629	1650	rBV	250390	501329	38.84%	5.829%
10	7.225	1896	1908	1935	rBV2	98717	194746	15.09%	2.264%
11	7.563	1999	2013	2032	rBV	419476	754946	58.48%	8.777%
12	7.852	2092	2103	2120	rBV	61224	117477	9.10%	1.366%
13	8.312	2235	2246	2288	rBV2	580546	1290892	100.00%	15.009%
14	9.087	2475	2487	2523	rBV	429171	767090	59.42%	8.919%
15	10.431	2893	2905	2921	rBV	183578	302467	23.43%	3.517%
16	11.482	3221	3232	3257	rBV	338543	590319	45.73%	6.863%
17	11.858	3335	3349	3374	rBV	351389	622646	48.23%	7.239%

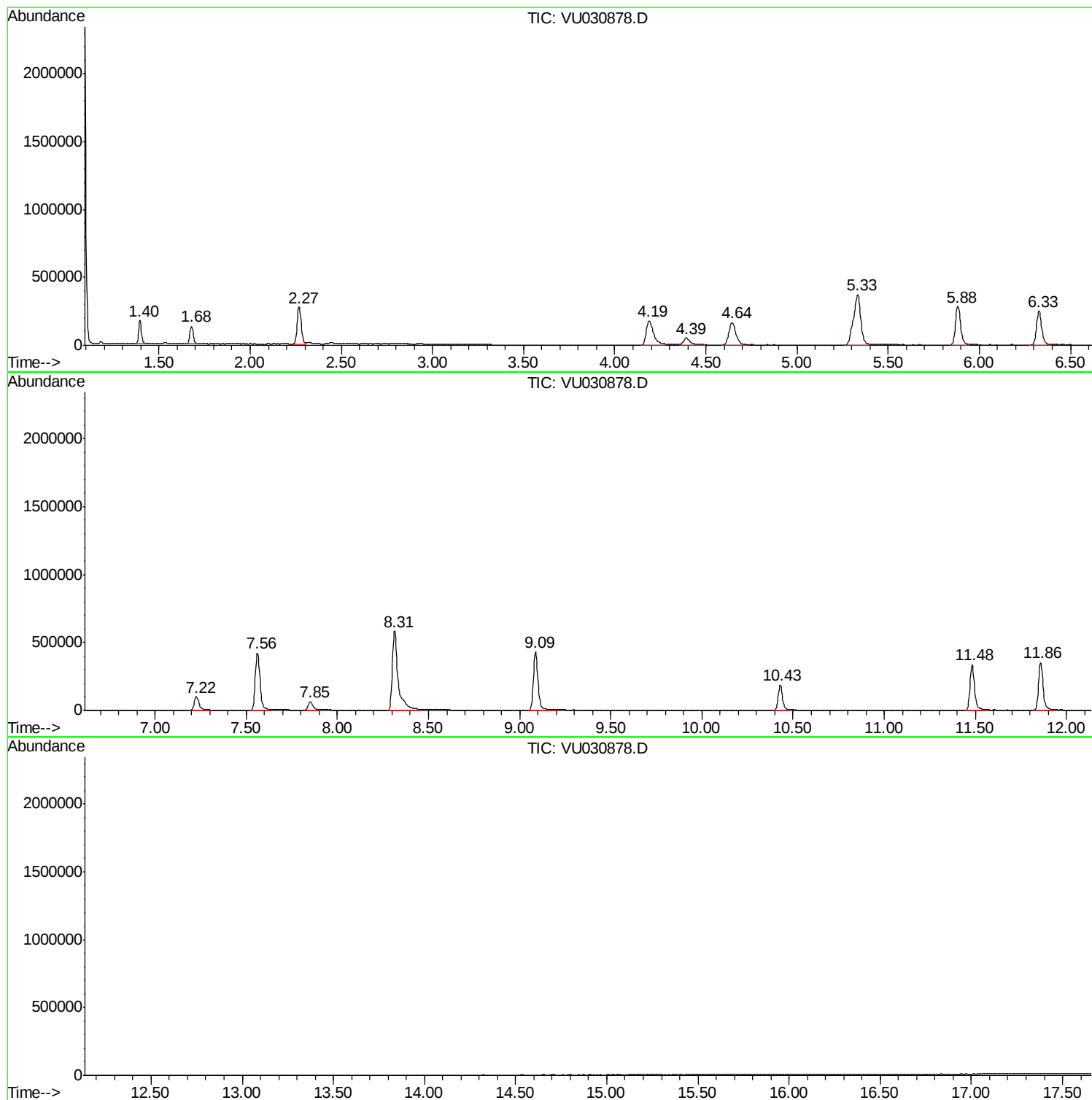
Sum of corrected areas: 8600926

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR032919WMA.M
Quant Title : TRACE VOA SOM01.0

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



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TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethyl Acetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	1.12 ug/L	128081	1,4-Difluorobenzene	5.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl Acetate	88	C4H8O2	000141-78-6	86
2			Ethyl Acetate	88	C4H8O2	000141-78-6	86
3			Ethyl Acetate	88	C4H8O2	000141-78-6	86
4			Ethyl Acetate	88	C4H8O2	000141-78-6	78
5			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	23

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethyl Acetate	4.39	1.1	ug/L	128081	1	5.88	572257	5.0