

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU111519\
 Data File : VU035723.D
 Acq On : 15 Nov 2019 13:50
 Operator : JC/SP
 Sample : K5759-01 50X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 F2J21

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\SOMUTR110419WMA.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	1.613	78	85	100	rBV	120706	135602	12.89%	1.786%
2	1.932	177	184	197	rVB	112117	146744	13.95%	1.933%
3	2.594	379	390	406	rBV	185748	304961	29.00%	4.018%
4	3.083	533	542	554	rVB3	8066	15862	1.51%	0.209%
5	4.713	1032	1049	1083	rBV2	147638	465529	44.26%	6.133%
6	5.115	1156	1174	1200	rBV2	175561	447286	42.53%	5.893%
7	5.771	1358	1378	1412	rBV2	356612	899418	85.52%	11.849%
8	6.292	1526	1540	1571	rBV	262705	519255	49.37%	6.841%
9	6.735	1664	1678	1697	rBV	222140	439327	41.77%	5.788%
10	7.607	1931	1949	1972	rBV	97029	190550	18.12%	2.510%
11	7.935	2037	2051	2071	rBV	365582	658231	62.59%	8.672%
12	8.218	2127	2139	2159	rBV2	60544	117546	11.18%	1.549%
13	8.687	2267	2285	2348	rBV	382059	1051725	100.00%	13.856%
14	9.449	2506	2522	2544	rBV	394646	715052	67.99%	9.420%
15	10.787	2924	2938	2961	rVB	187533	326722	31.07%	4.304%
16	11.841	3253	3266	3286	rBV2	313085	598923	56.95%	7.890%
17	12.221	3367	3384	3401	rBV	303657	557842	53.04%	7.349%

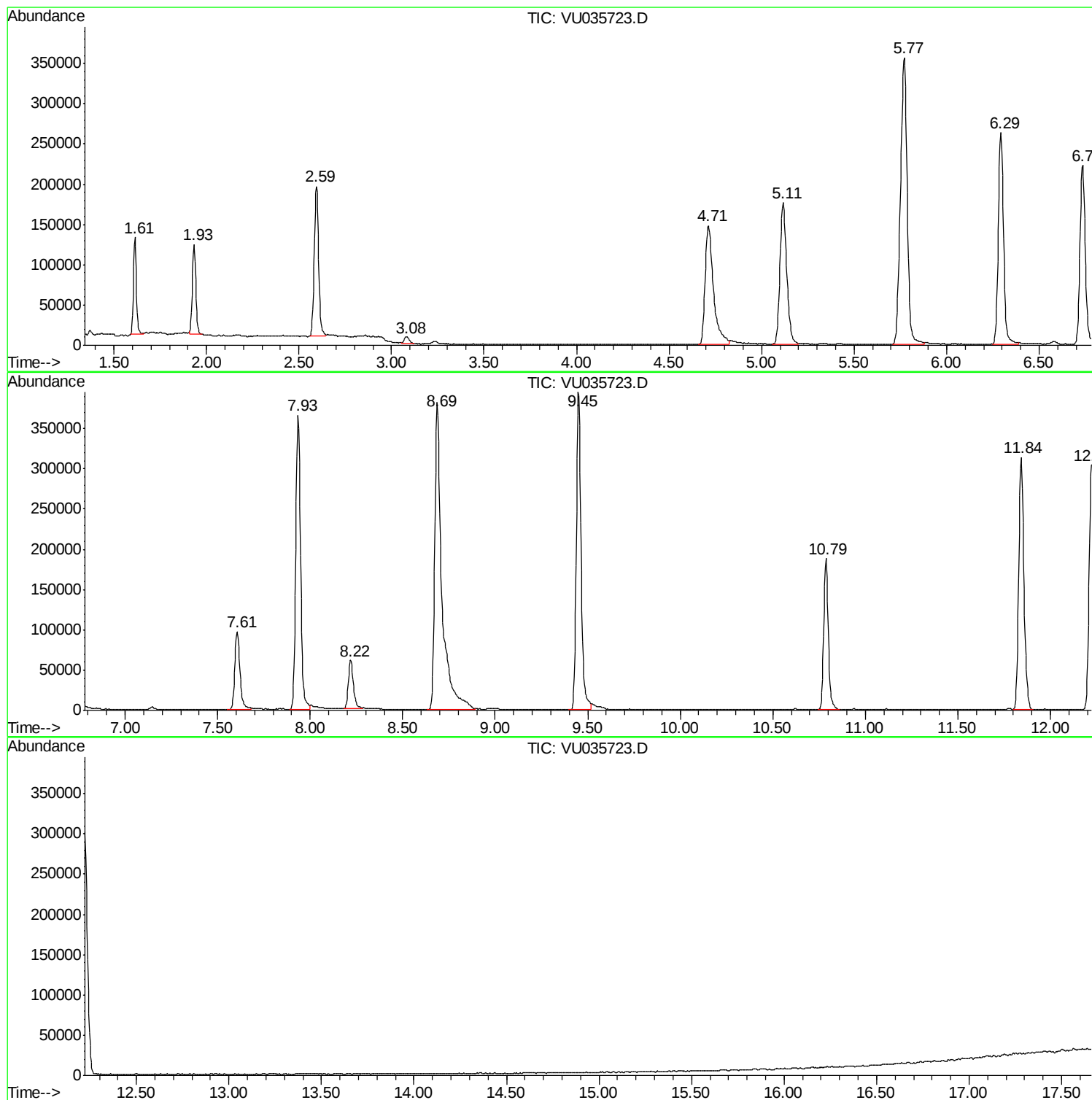
Sum of corrected areas: 7590575

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR110419WMA.M
Quant Title : TRACE VOA SOM01.0

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



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Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	1.83 ug/L	190550	1,4-Difluorobenzene	6.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclohexadiene	80	C6H8	000592-57-4	28
2			2-Propanol, 1,3-dichloro-	128	C3H6Cl2O	000096-23-1	25
3			Methane, oxybis[chloro-	114	C2H4Cl2O	000542-88-1	25
4			Thiophosgene	114	CCl2S	000463-71-8	9
5			2-Propanol, 1,3-dichloro-	128	C3H6Cl2O	000096-23-1	9

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	7.61	1.8	ug/L	190550	1	6.29	519255	5.0