

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070318\
 Data File : VU025192.D
 Acq On : 03 Jul 2018 15:41
 Operator : MD/SY
 Sample : J3790-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ZER-WC-L-01

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\82U061318W.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	1.088	136	155	174	rBV	538918	667338	58.13%	9.637%
2	1.387	241	248	259	rBV	22362	23950	2.09%	0.346%
3	2.323	527	539	553	rVB	17584	29050	2.53%	0.420%
4	4.680	1253	1272	1289	rBV	83691	198244	17.27%	2.863%
5	4.889	1318	1337	1353	rBV	176754	404186	35.21%	5.837%
6	4.989	1353	1368	1392	rVB2	264161	585921	51.04%	8.462%
7	5.313	1452	1469	1491	rBV	187735	401272	34.96%	5.795%
8	5.892	1630	1649	1672	rBV	365964	726185	63.26%	10.487%
9	7.570	2151	2171	2207	rBV	661844	1147952	100.00%	16.578%
10	9.094	2631	2645	2669	rBV	503739	844686	73.58%	12.198%
11	10.313	3012	3024	3046	rBV	493050	796087	69.35%	11.497%
12	11.487	3377	3389	3408	rBV	539684	869278	75.72%	12.554%
13	11.754	3460	3472	3491	rBV3	14951	31414	2.74%	0.454%
14	13.798	4100	4108	4115	rVV3	8051	11800	1.03%	0.170%
15	13.850	4115	4124	4137	rVB	38159	58871	5.13%	0.850%
16	13.943	4142	4153	4167	rBV2	55451	89230	7.77%	1.289%
17	16.567	4956	4969	4984	rBV3	20876	39086	3.40%	0.564%

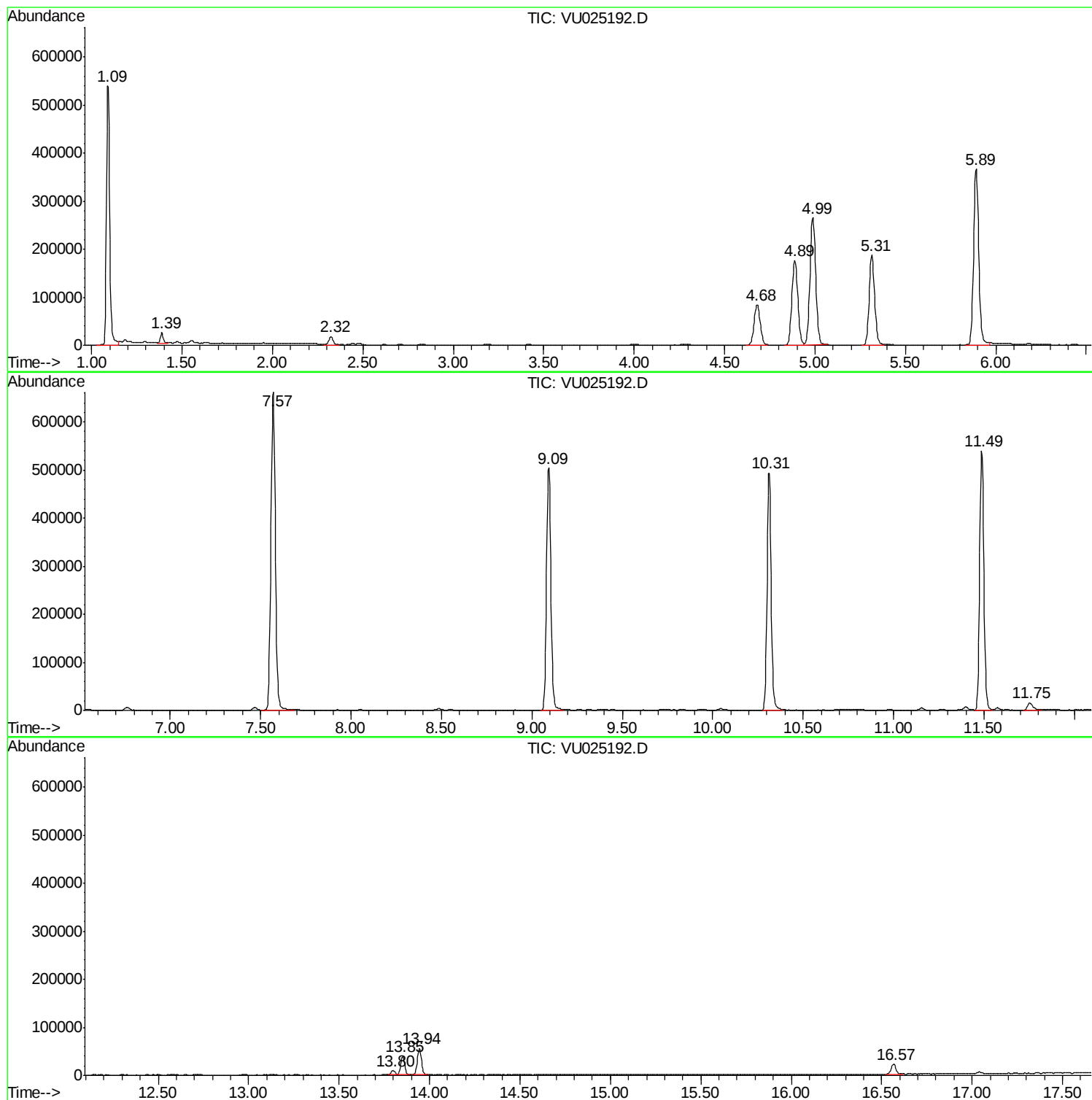
Sum of corrected areas: 6924550

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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



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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown13.94 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	5.13 ug/l	89230	1,4-Dichlorobenzene-d4	11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			dl-Homoserine	119	C4H9NO3	001927-25-9	27
2			1,3,4-Thiadiazol-2-amine, 5-methyl-	115	C3H5N3S	000108-33-8	12
3			Silanol, trimethyl-, nitrate	135	C3H9NO3Si	018026-82-9	12
4			1,2-Dithiane	120	C4H8S2	000505-20-4	12
5			1,3-Dithiane	120	C4H8S2	000505-23-7	12

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown13.94	13.94	5.1	ug/l	89230	4	11.49	869278	50.0