

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_R\DATA\VR112118\
 Data File : VR026262.D
 Acq On : 21 Nov 2018 15:45
 Operator : SY/MD
 Sample : VR1121WBL01
 Misc : 25mL/MSVOA_R/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_R
 ClientSampleId :
 VBLK99

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 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_R\METHODS\SOMRTR112118WMA.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	2.160	70	79	93	rBV2	92580	323942	15.41%	2.256%
2	2.628	148	156	168	rVB2	139518	349905	16.64%	2.437%
3	3.614	306	318	334	rBV2	276605	1138802	54.17%	7.932%
4	6.588	798	807	816	rBV2	98077	292416	13.91%	2.037%
5	7.203	897	908	919	rBV	353553	959198	45.62%	6.681%
6	7.854	1005	1015	1031	rBV2	769921	2102358	100.00%	14.644%
7	8.462	1105	1115	1124	rBV	778044	1436017	68.31%	10.002%
8	8.900	1179	1187	1194	rBV	469789	887663	42.22%	6.183%
9	9.685	1310	1316	1327	rVB	176856	305141	14.51%	2.125%
10	9.971	1354	1363	1370	rBV	1280759	1978078	94.09%	13.778%
11	10.032	1370	1373	1381	rVB	27214	48409	2.30%	0.337%
12	10.239	1400	1407	1414	rVB	99762	160081	7.61%	1.115%
13	10.591	1459	1465	1475	rBV	449086	693253	32.98%	4.829%
14	11.291	1573	1580	1589	rBV	999035	1524057	72.49%	10.616%
15	12.362	1750	1756	1763	rBV	170594	261308	12.43%	1.820%
16	13.226	1891	1898	1905	rBV	700261	985083	46.86%	6.862%
17	13.518	1939	1946	1956	rBV	605288	910936	43.33%	6.345%

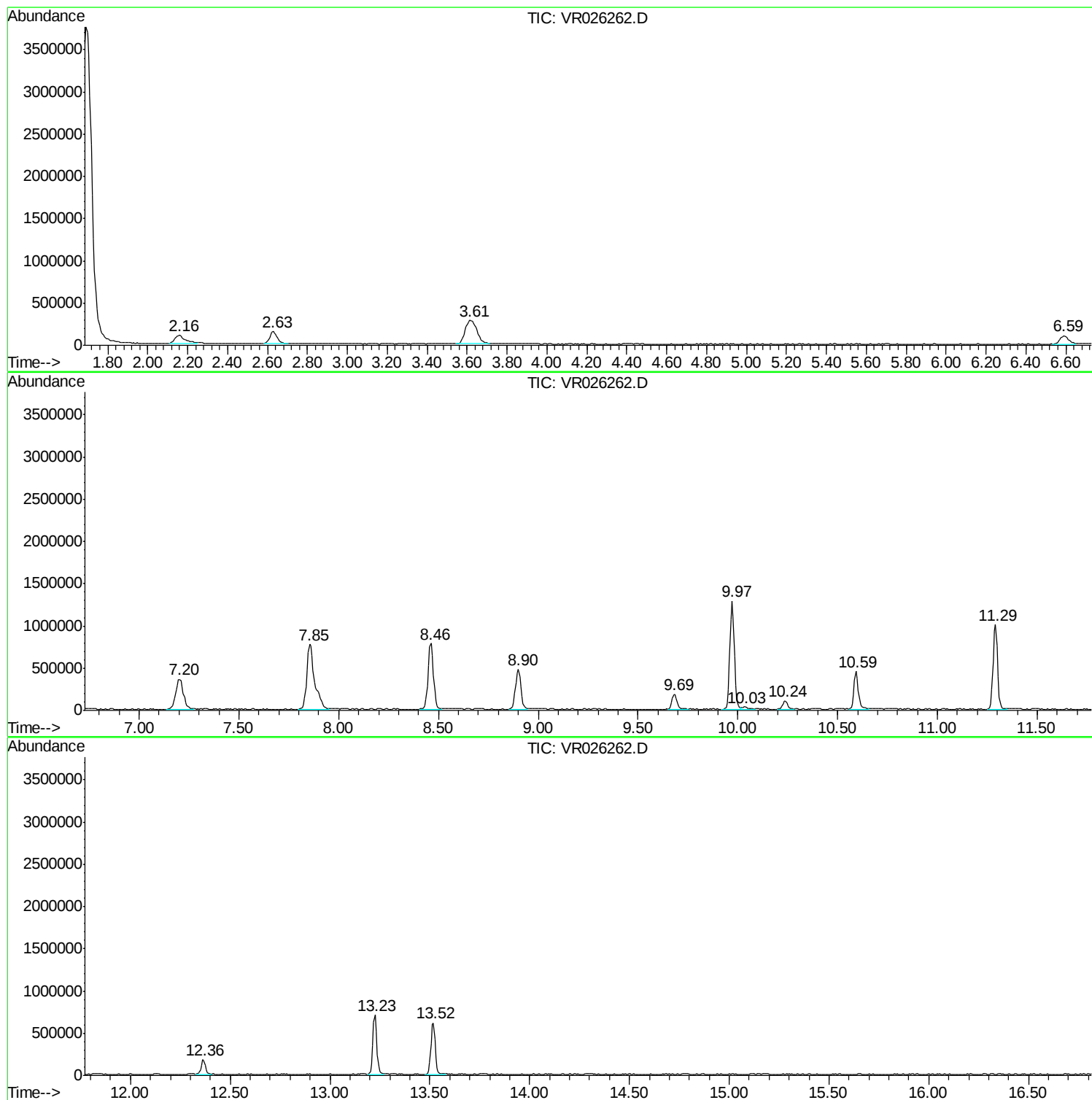
Sum of corrected areas: 14356647

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR112118WMA.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.
9.69	1.06 ug/L	305141	1,4-Difluorobenzene	8.46

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1,3-dichloro-	128	C3H6Cl2O	000096-23-1	17
2			Cyclopropane	42	C3H6	000075-19-4	9
3			n-Propyl chloride	78	C3H7Cl	000540-54-5	9
4			n-Propyl chloride	78	C3H7Cl	000540-54-5	9
5			Methane, oxybis[chloro-	114	C2H4Cl2O	000542-88-1	9

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	9.69	1.1	ug/L	305141	1	8.46	1436020	5.0