

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD100318\
 Data File : VD060118.D
 Acq On : 3 Oct 2018 15:10
 Operator : VA/AP
 Sample : J5219-02
 Misc : 10.11g/5ml/MSVOA D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 FK-BPM-01-E10-F10-E10A-F10A

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_D\METHOD\82D100118S.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	1.700	15	18	28	rVB	105719	182580	3.97%	0.615%
2	3.147	160	165	177	rBV	35383	108934	2.37%	0.367%
3	3.698	216	221	232	rBV	319508	921651	20.04%	3.104%
4	3.856	234	237	244	rVB3	19770	56565	1.23%	0.190%
5	4.495	291	302	311	rVB3	18462	57533	1.25%	0.194%
6	5.647	411	419	433	rBV	96546	319980	6.96%	1.078%
7	6.316	480	487	491	rBV	826313	2171504	47.22%	7.313%
8	6.394	491	495	512	rVB	1222198	3146042	68.41%	10.595%
9	6.778	527	534	550	rBV	622090	1593347	34.65%	5.366%
10	7.428	594	600	617	rBV	1504403	3665696	79.71%	12.345%
11	9.445	800	805	813	rBV	2110201	4598651	100.00%	15.487%
12	10.183	874	880	889	rBV	123355	334217	7.27%	1.126%
13	11.276	987	991	1002	rBV	2546854	4069369	88.49%	13.705%
14	12.427	1104	1108	1117	rBV	2286680	3504647	76.21%	11.803%
15	12.565	1117	1122	1127	rVB	279831	494370	10.75%	1.665%
16	13.372	1200	1204	1213	rBV	2933026	4271062	92.88%	14.384%
17	13.952	1259	1263	1271	rBV	136708	197258	4.29%	0.664%

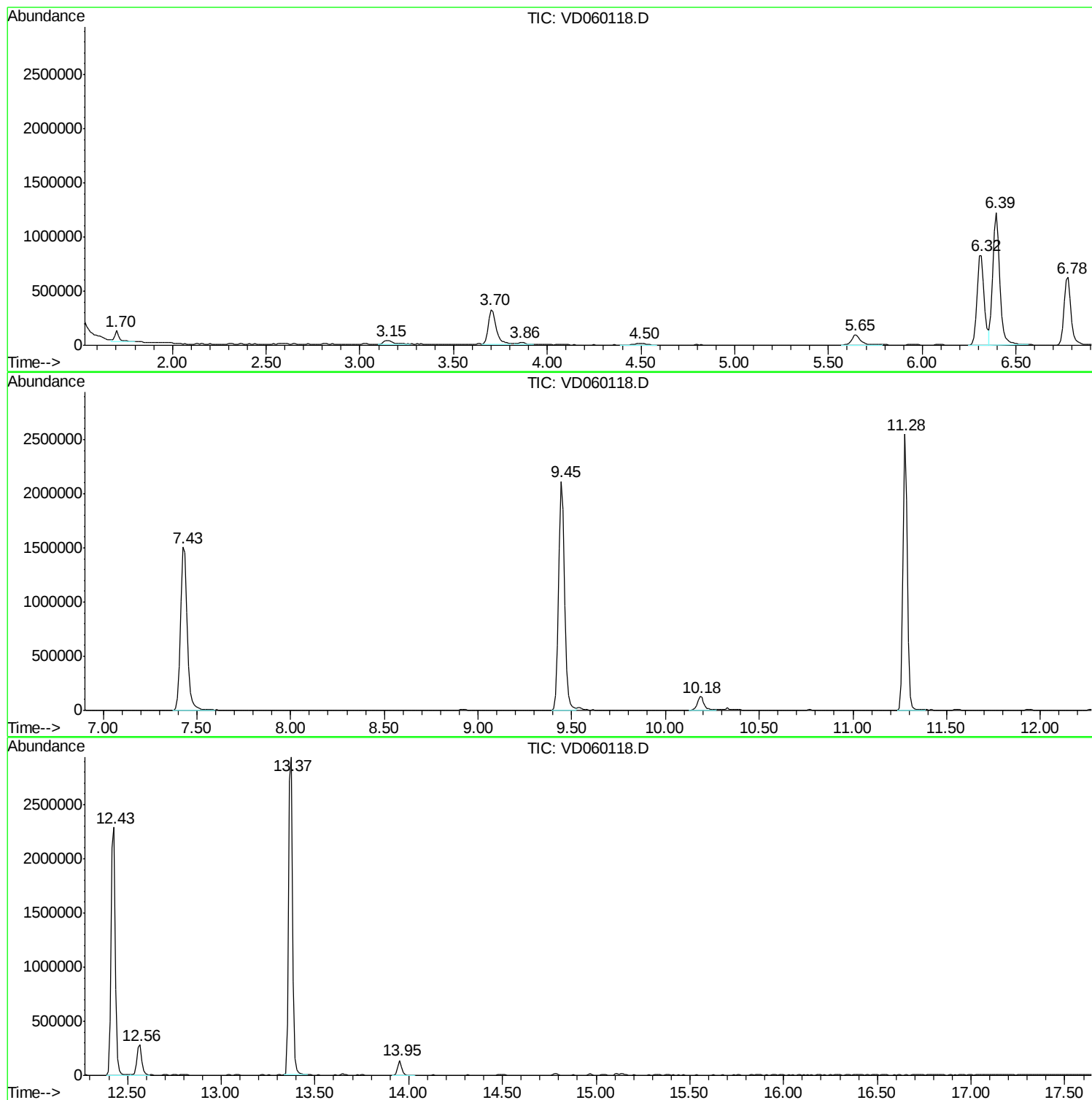
Sum of corrected areas: 29693406

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.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 7H-Dibenzo[b,g]carbazole, 7... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	5.79 ug/l	494370	1,4-Dichlorobenzene-d4	13.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7H-Dibenzo[b,g]carbazole, 7-methyl-	281	C21H15N	003557-49-1	59
2			4H-1,2,4-Triazole-3-thiol, 4-all...	281	C16H15N3S	031803-13-1	59
3			Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	53
4			5H-Naphtho[2,3-c]carbazole, 5-me...	281	C21H15N	100025-44-3	45
5			2',6'-Dihydroxyacetophenone, bis...	296	C14H24O3Si2	1000352-81-3	45

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
7H-Dibenzo[b,g]ca...	12.56	5.8	ug/l	494370	4	13.37	4271060	50.0