

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD100418\
 Data File : VD060131.D
 Acq On : 4 Oct 2018 16:06
 Operator : VA/AP
 Sample : J5191-05
 Misc : 4.98g/5ml/MSVOA D/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 344-OILY-STONES

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	3.717	218	223	234	rVV	223856	631409	13.14%	1.894%
2	4.514	297	304	314	rBV	135242	387887	8.07%	1.164%
3	5.665	413	421	433	rBV	69669	228155	4.75%	0.684%
4	6.334	481	489	493	rBV	958485	2511451	52.27%	7.534%
5	6.413	493	497	518	rVB	1356572	3560958	74.12%	10.682%
6	6.797	527	536	555	rBV2	763880	2013318	41.90%	6.040%
7	7.456	597	603	618	rBV	1781215	4334430	90.21%	13.002%
8	8.834	737	743	750	rBV	42463	118702	2.47%	0.356%
9	9.464	802	807	821	rVB	2060758	4677500	97.36%	14.032%
10	10.211	875	883	889	rBV	185566	483623	10.07%	1.451%
11	10.792	938	942	949	rBV2	30480	56109	1.17%	0.168%
12	11.235	982	987	989	rBV	54536	105447	2.19%	0.316%
13	11.294	989	993	1001	rVB	2989027	4804569	100.00%	14.413%
14	12.445	1106	1110	1117	rBV	2545959	3744843	77.94%	11.234%
15	12.583	1119	1124	1134	rVB	333890	590168	12.28%	1.770%
16	13.390	1202	1206	1214	rBV	3295971	4707735	97.98%	14.122%
17	13.971	1261	1265	1273	rVB	195802	283589	5.90%	0.851%
18	15.152	1379	1385	1389	rBV	66453	95653	1.99%	0.287%

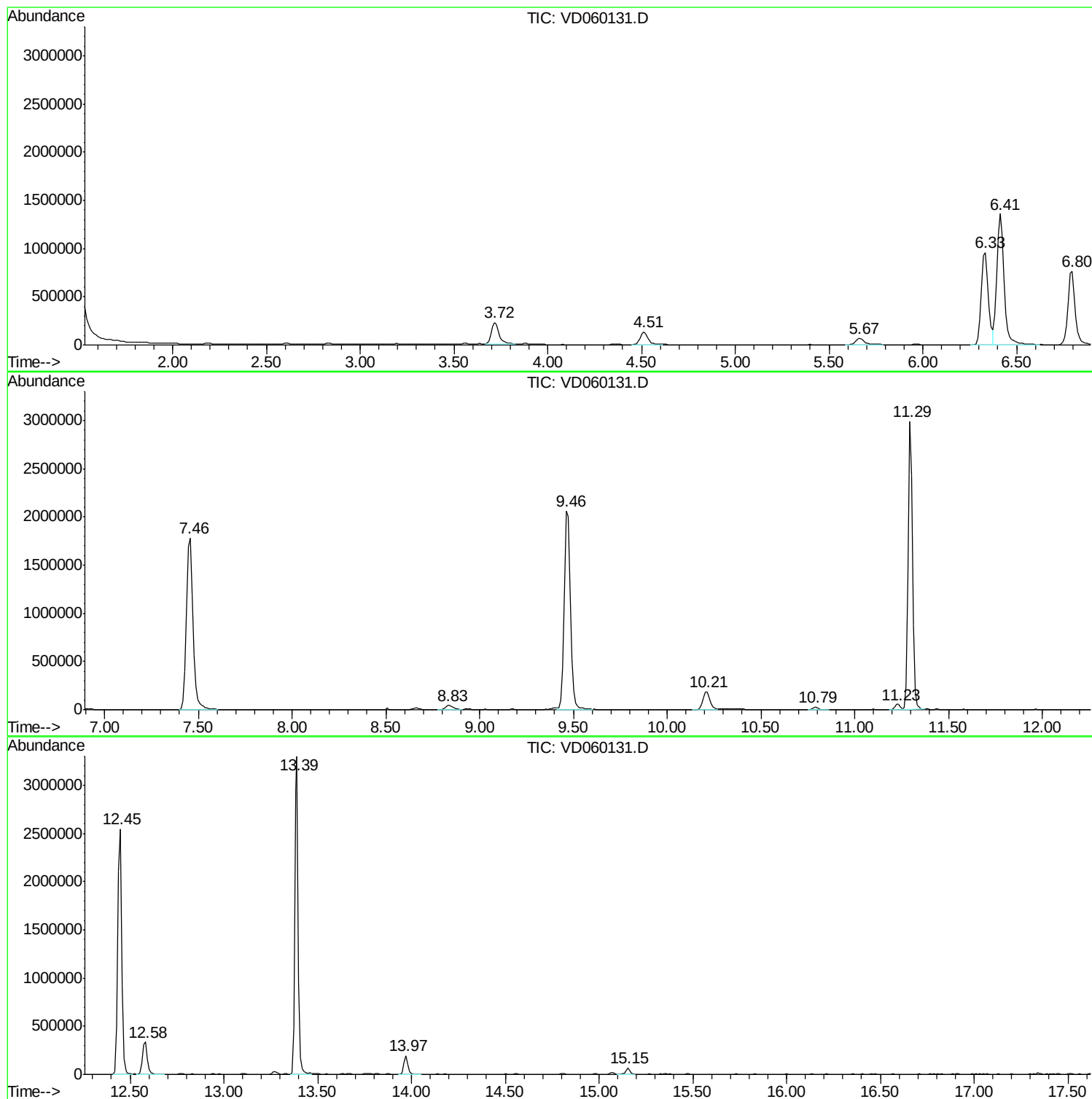
Sum of corrected areas: 33335546

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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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Peak Number 1 n-Hexane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.51	5.45 ug/l	387887	Pentafluorobenzene	6.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	91
2			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	37
3			Pentane, 3-methyl-	86	C6H14	000096-14-0	37
4			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	33
5			Butanal, 2-methyl-	86	C5H10O	000096-17-3	27

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexane	4.51	5.5	ug/l	387887	1	6.41	3560960	50.0