

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW091019\
 Data File : VW012893.D
 Acq On : 11 Sep 2019 02:46
 Operator : SY/VA
 Sample : K4696-05
 Misc : 5.08G/5ML/MSVOA W/SOIL
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WS-SB06-4.75-5.0

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_W\METHOD\82W090719S.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	2.148	35	40	49	rBV	24184	43156	5.71%	1.056%
2	2.709	124	132	133	rBV8	3659	7733	1.02%	0.189%
3	4.903	481	492	505	rBV2	15832	52376	6.93%	1.281%
4	5.934	650	661	674	rBV3	16059	50370	6.66%	1.232%
5	7.141	850	859	875	rVB	24465	71985	9.52%	1.761%
6	7.884	969	981	985	rBV	100438	241750	31.99%	5.913%
7	7.945	985	991	1002	rVB	209888	458652	60.69%	11.219%
8	8.299	1037	1049	1059	rBV	105408	238863	31.61%	5.843%
9	8.842	1129	1138	1152	rBV	282759	579767	76.71%	14.181%
10	10.323	1370	1381	1389	rBV	425850	755766	100.00%	18.486%
11	10.701	1437	1443	1449	rBV2	18533	31678	4.19%	0.775%
12	11.061	1496	1502	1511	rVB	18975	31805	4.21%	0.778%
13	11.628	1588	1595	1607	rVB	359086	602659	79.74%	14.741%
14	12.615	1749	1757	1767	rVB	269128	440457	58.28%	10.774%
15	12.932	1802	1809	1814	rBV2	6065	8644	1.14%	0.211%
16	13.554	1904	1911	1920	rBV	275061	450445	59.60%	11.018%
17	14.072	1988	1996	2001	rVB2	4910	8624	1.14%	0.211%
18	15.438	2213	2220	2227	rBV3	7977	13595	1.80%	0.333%

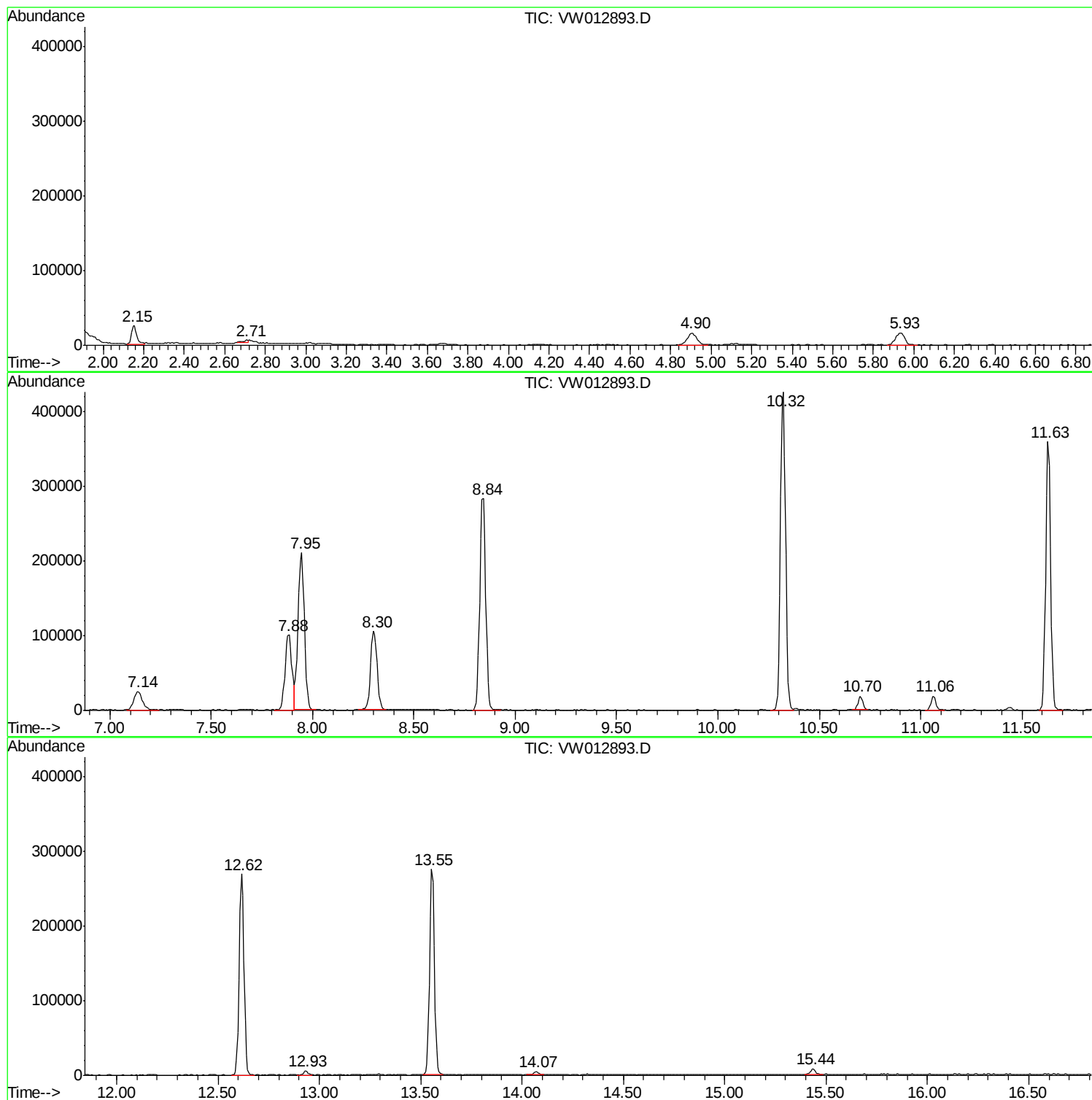
Sum of corrected areas: 4088325

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W090719S.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.
5.93	5.49 ug/l	50370	Pentafluorobenzene	7.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	86
2			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	42
3			Butane, 2-nitro-	103	C4H9NO2	000600-24-8	36
4			Cyclobutane	56	C4H8	000287-23-0	35
5			1-Propene, 2-methyl-	56	C4H8	000115-11-7	35

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexane	5.93	5.5	ug/l	50370	1	7.95	458652	50.0