

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD092820\
 Data File : VD066919.D
 Acq On : 28 Sep 2020 17:55
 Operator : VA/SY
 Sample : L4136-01
 Misc : 5.60G/5.00ml/MSVOA D/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 PAV-B26-092420-0-1

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_D\METHOD\82D092220S.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	1.944	3	4	18	rVB	171364	178075	8.08%	1.442%
2	2.314	61	67	77	rBV2	69306	145225	6.59%	1.176%
3	3.314	229	237	244	rBV3	16815	35745	1.62%	0.289%
4	3.426	247	256	267	rVB7	18731	61558	2.79%	0.498%
5	3.767	305	314	324	rVB7	7533	23081	1.05%	0.187%
6	4.967	507	518	532	rBV2	48569	155887	7.07%	1.262%
7	7.914	1008	1019	1024	rBV	299341	734684	33.34%	5.949%
8	7.979	1024	1030	1043	rVB	527832	1170311	53.11%	9.476%
9	8.326	1079	1089	1101	rBV	269254	614864	27.90%	4.979%
10	8.861	1171	1180	1194	rBV	768836	1574648	71.46%	12.751%
11	10.338	1421	1431	1439	rBV	1224153	2203544	100.00%	17.843%
12	11.643	1644	1653	1665	rBV	1180014	2053495	93.19%	16.628%
13	12.632	1814	1821	1835	rVB	912533	1494113	67.80%	12.098%
14	13.579	1974	1982	1990	rBV	1175396	1876174	85.14%	15.192%
15	15.402	2286	2292	2301	rBV2	14529	28231	1.28%	0.229%

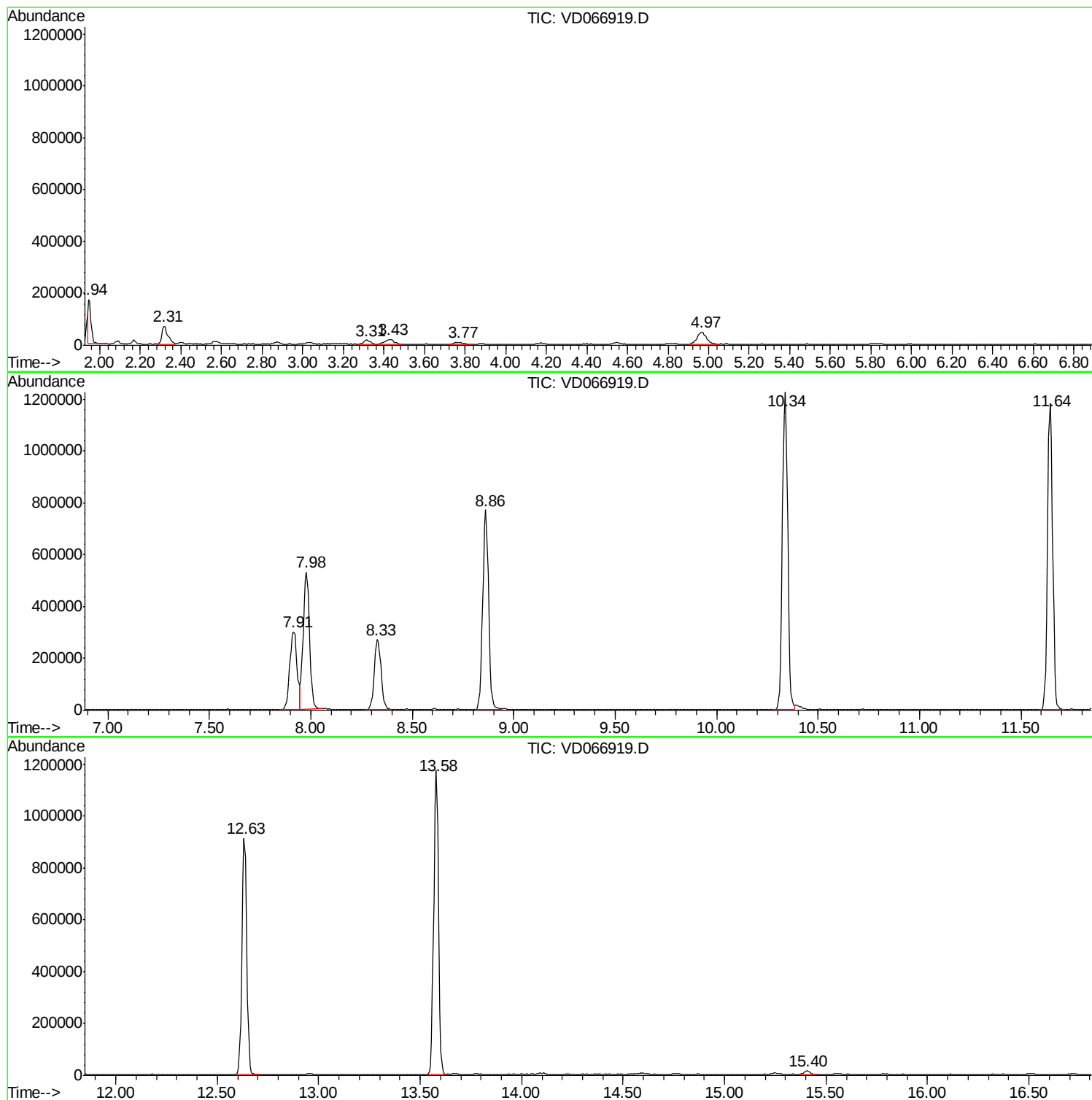
Sum of corrected areas: 12349635

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.M
Quant Title : SW846 8260

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



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Peak Number 1 3-Butenoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.94	7.61 ug/l	178075	Pentafluorobenzene	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Butenoic acid	86	C4H6O2	000625-38-7	64
2		Bis(3-methylbutyl) fluorene-2,7-...	466	C23H30O6S2	253664-95-8	40
3		2-Butenal, (E)-	70	C4H6O	000123-73-9	40
4		Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	39
5		Propene	42	C3H6	000115-07-1	22

