

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091019\
 Data File : VN057850.D
 Acq On : 10 Sep 2019 10:32
 Operator : JC/SP
 Sample : K4781-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GW801-8.0-13.0-090719

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_N\METHODS\82N091019W.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	1.799	3	8	21	rVB2	73927	93856	3.92%	0.805%
2	2.137	104	113	130	rVB2	37158	73539	3.07%	0.631%
3	2.426	194	203	211	rBV3	20925	35275	1.47%	0.303%
4	3.805	619	632	651	rBV2	10691	28050	1.17%	0.241%
5	4.050	690	708	728	rBV3	27663	83480	3.49%	0.716%
6	6.812	1550	1567	1591	rBV4	8859	29298	1.22%	0.251%
7	7.355	1716	1736	1737	rBV3	20214	30720	1.28%	0.263%
8	7.577	1786	1805	1816	rBV	289696	716662	29.94%	6.147%
9	7.651	1816	1828	1850	rVB	377207	907131	37.90%	7.780%
10	8.014	1923	1941	1965	rBV2	371562	868488	36.29%	7.449%
11	8.297	1984	2029	2082	rBV4	71120	738949	30.87%	6.338%
12	8.577	2100	2116	2139	rVB	575443	1214967	50.76%	10.421%
13	10.088	2571	2586	2615	rBV	1286092	2393417	100.00%	20.528%
14	11.406	2982	2996	3019	rBV	837209	1448535	60.52%	12.424%
15	12.403	3292	3306	3323	rBV	885665	1447362	60.47%	12.414%
16	13.348	3587	3600	3616	rBV	929404	1549423	64.74%	13.289%

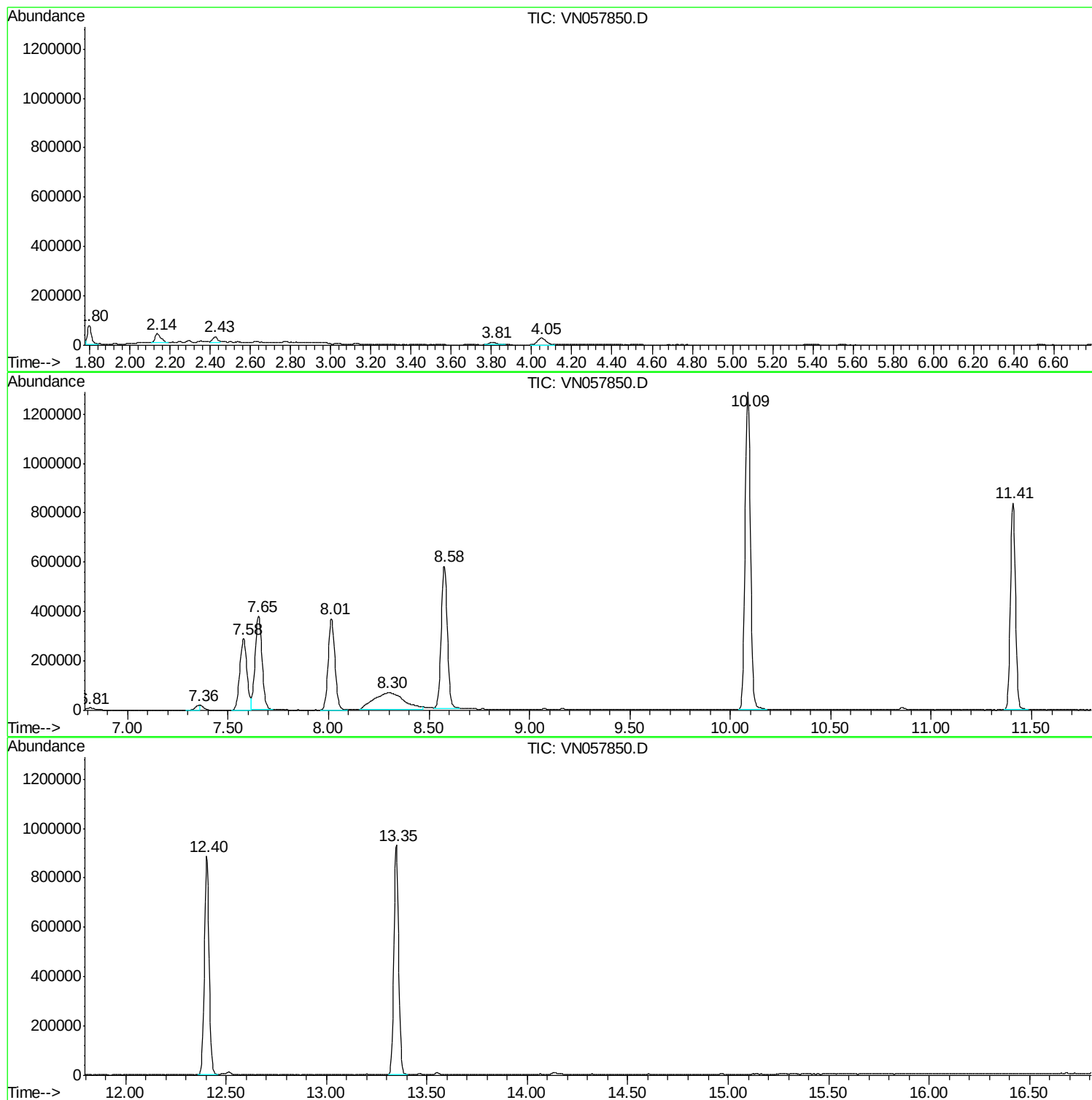
Sum of corrected areas: 11659152

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
Quant Title : SW846 8260

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



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Peak Number 1 Propene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.80	5.17 ug/l	93856	Pentafluorobenzene	7.65	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Cyclopropane	42	C3H6	000075-19-4	40
3	Methylacrylonitrile	67	C4H5N	000126-98-7	22
4	3-Butenenitrile	67	C4H5N	000109-75-1	12
5	2-Butenenitrile	67	C4H5N	004786-20-3	10

Peak Number 2 Acetic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.30	30.41 ug/l	738949	1,4-Difluorobenzene	8.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid	60	C2H4O2	000064-19-7	90
2	Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	45
3	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	9
4	Propanal, 3-methoxy-	88	C4H8O2	002806-84-0	9
5	Ammonium acetate	77	C2H7NO2	000631-61-8	9

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.80	5.2	ug/l	93856	1	7.65	907131	50.0
Acetic acid	8.30	30.4	ug/l	738949	2	8.58	1214970	50.0