

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091718\
 Data File : VN051275.D
 Acq On : 17 Sep 2018 17:24
 Operator : MD\SY
 Sample : J4773-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CL-01-091418-A

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_N\METHODS\82N091718W.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	4.555	838	865	916	rBV	3743643	11768918	100.00%	52.103%
2	7.593	1791	1810	1821	rBV	326723	806214	6.85%	3.569%
3	7.667	1821	1833	1859	rVB	486216	1161093	9.87%	5.140%
4	8.030	1928	1946	1972	rBV	273194	653866	5.56%	2.895%
5	8.165	1972	1988	2004	rBV	143237	321044	2.73%	1.421%
6	8.587	2103	2119	2137	rBV	660820	1389490	11.81%	6.151%
7	9.272	2318	2332	2358	rBV	128271	257842	2.19%	1.142%
8	10.091	2571	2587	2622	rBV	1148027	2148430	18.26%	9.511%
9	10.767	2786	2797	2815	rVB2	69055	121145	1.03%	0.536%
10	11.169	2910	2922	2937	rBV	105982	176769	1.50%	0.783%
11	11.407	2983	2996	3020	rBV	817262	1487553	12.64%	6.586%
12	12.403	3292	3306	3325	rBV	630505	1094005	9.30%	4.843%
13	13.345	3586	3599	3616	rVB	561936	949561	8.07%	4.204%
14	15.133	4141	4155	4168	rBV	138311	251914	2.14%	1.115%

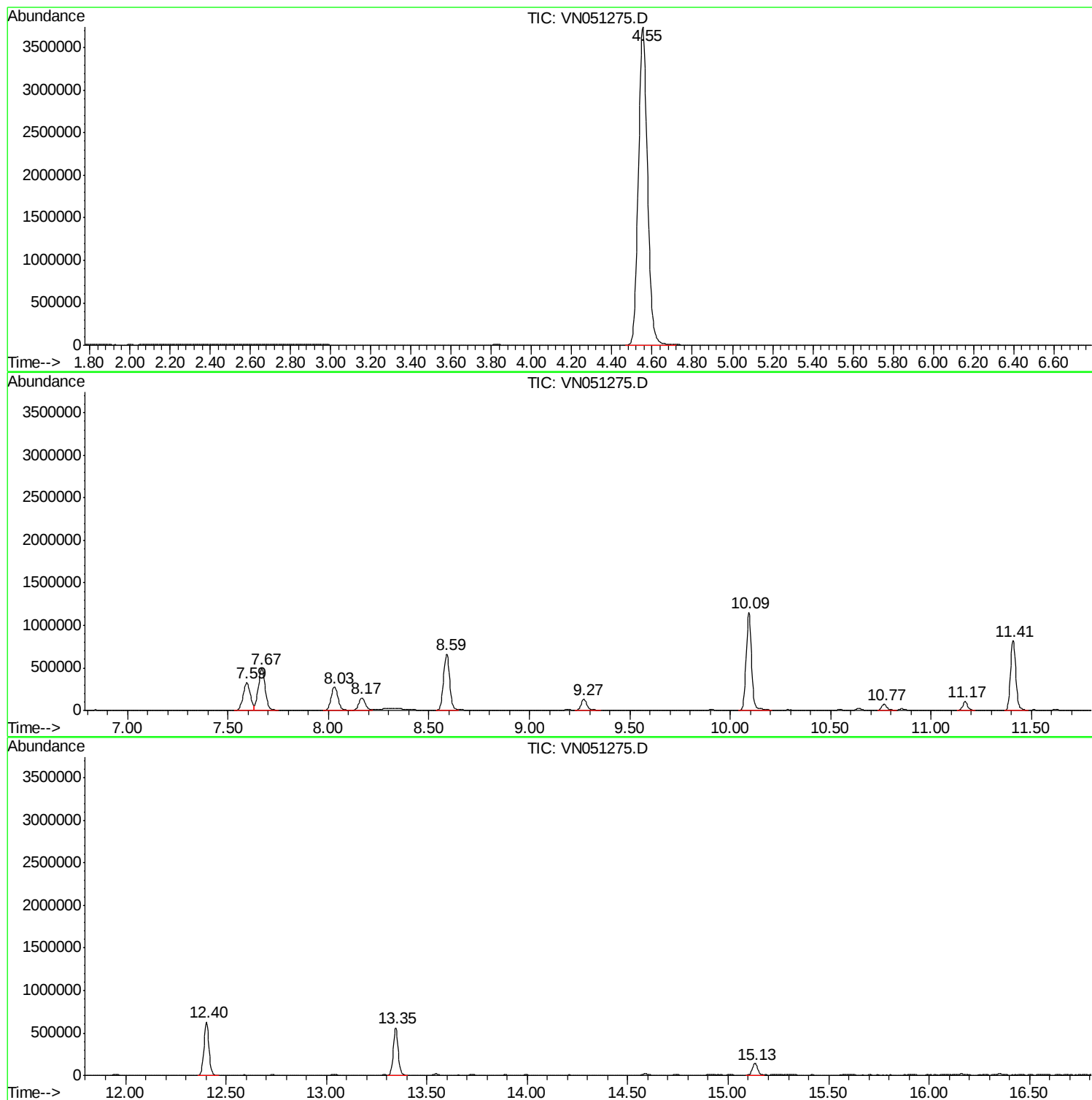
Sum of corrected areas: 22587844

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Quant Title : SW846 8260

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



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Peak Number 1 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.27	9.28 ug/l	257842	1,4-Difluorobenzene	8.59	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2	Isopropyl acetate	102	C5H10O2	000108-21-4	36
3	1-Butanamine	73	C4H11N	000109-73-9	9
4	Propanal, 2,3-dihydroxy-, (S)-	90	C3H6O3	000497-09-6	9
5	Isobutane	58	C4H10	000075-28-5	7

Peak Number 2 Butanoic acid, 1-methylethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.17	5.94 ug/l	176769	Chlorobenzene-d5	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
2	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	72
3	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
4	Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	50
5	Propanoic acid, 2-methyl-, penty...	158	C9H18O2	002445-72-9	50

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Propyl acetate	9.27	9.3	ug/l	257842	2	8.59	1389490	50.0
Butanoic acid, 1-...	11.17	5.9	ug/l	176769	3	11.41	1487550	50.0