

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081518\
Data File : BN002359.D
Acq On : 16 Aug 2018 14:56
Operator : JU/SJ
Sample : PB111989BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK89

Integration Parameters: LSCINT.P

Integrator: RTE
Smoothing : OFF
Sampling : 1
Start Thrs: 0.2
Stop Thrs : 0
Filtering: 5
Min Area: 1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	26	30	41	rVB	292119	360958	11.74%	2.962%
2	7.675	791	797	808	rBV	468698	780966	25.40%	6.410%
3	10.446	1261	1268	1289	rBV	619182	1112501	36.18%	9.131%
4	14.310	1918	1925	1937	rBV2	1005682	1528151	49.70%	12.542%
5	17.057	2386	2392	2409	rBV	1567715	2263899	73.63%	18.580%
6	21.257	3100	3106	3118	rBV	2266975	3019324	98.20%	24.780%
7	23.363	3462	3464	3472	rVB6	26099	43856	1.43%	0.360%
8	23.474	3476	3483	3494	rVB2	1613670	3074727	100.00%	25.235%

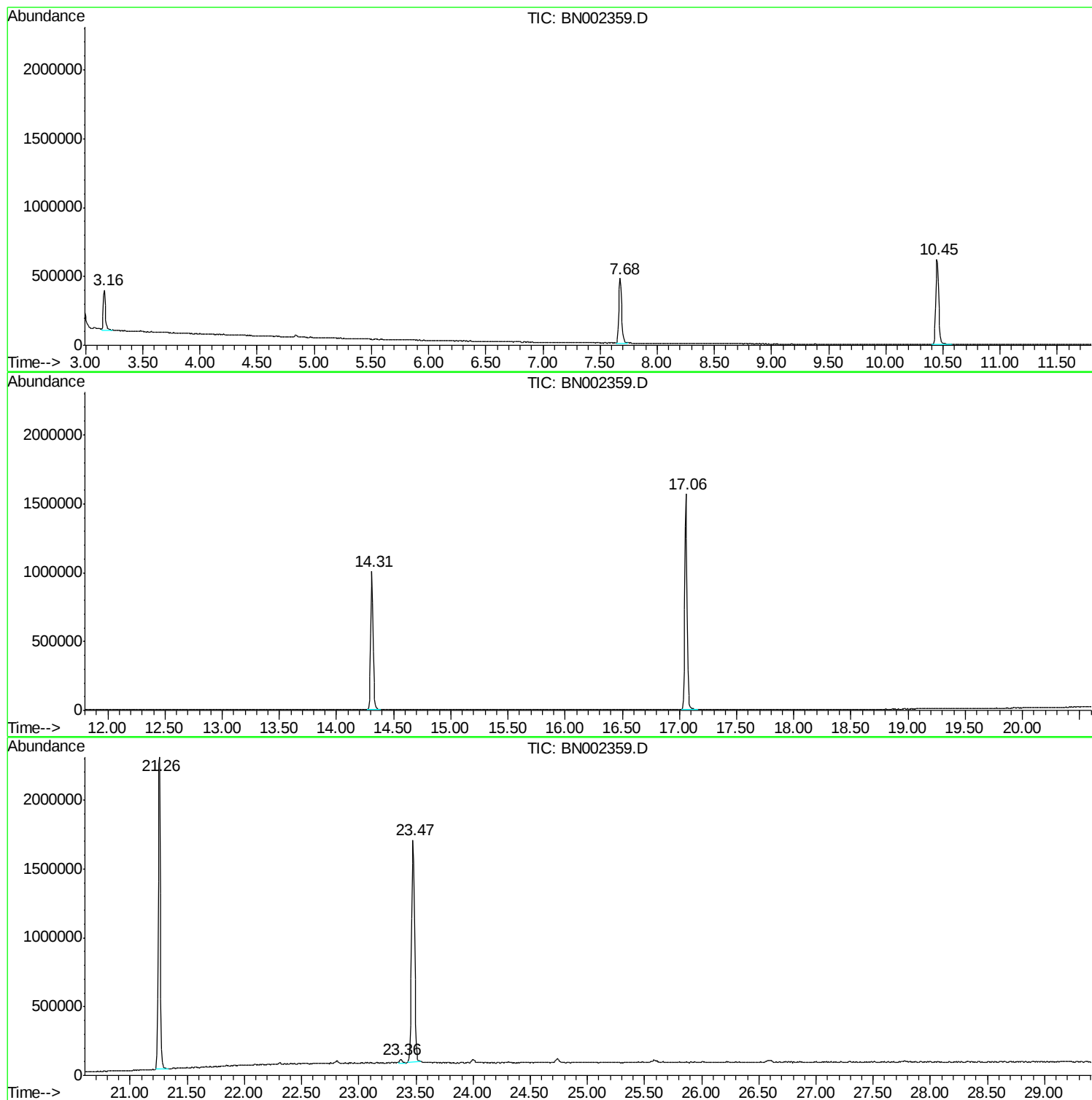
Sum of corrected areas: 12184382

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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



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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.16	9.24 ng/ul	360958	1,4-Dichlorobenzene-d4	7.68	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	91
2	Heptane	100	C7H16	000142-82-5	91
3	Heptane	100	C7H16	000142-82-5	90
4	Heptane	100	C7H16	000142-82-5	64
5	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	64

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.16	9.2	ng/ul	360958	1	7.68	780966	20.0