

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110416\
 Data File : BF091014.D
 Acq On : 4 Nov 2016 9:26
 Operator : UM/SJ
 Sample : PB94512BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94512BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF110316.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.778	7	8	51	rVB	2484060	6829816	100.00%	12.840%
2	4.864	276	278	290	rBV	653233	1150562	16.85%	2.163%
3	5.298	313	316	319	rBV	3614619	4881421	71.47%	9.177%
4	6.350	405	408	410	rBV	3854337	4341119	63.56%	8.161%
5	6.476	416	419	421	rBV	4110792	5197067	76.09%	9.770%
6	6.704	437	439	441	rBV	863978	975333	14.28%	1.834%
7	6.853	450	452	455	rBV	3207881	3736291	54.71%	7.024%
8	7.264	486	488	491	rBV	1964614	3085414	45.18%	5.800%
9	7.984	549	551	554	rBV	1054745	1252844	18.34%	2.355%
10	9.070	642	646	648	rBV	4824376	5491421	80.40%	10.324%
11	9.745	701	705	707	rBV	1278563	1657433	24.27%	3.116%
12	10.533	770	774	776	rBV	3333629	3661947	53.62%	6.884%
13	11.219	832	834	837	rBV	1797602	1702027	24.92%	3.200%
14	12.808	970	973	976	rBV	4943902	6298660	92.22%	11.841%
15	13.848	1062	1064	1067	rBV	1141644	1501811	21.99%	2.823%
16	15.242	1183	1186	1188	rBV	1100175	1336329	19.57%	2.512%
17	15.368	1194	1197	1208	rVB2	26000	92737	1.36%	0.174%

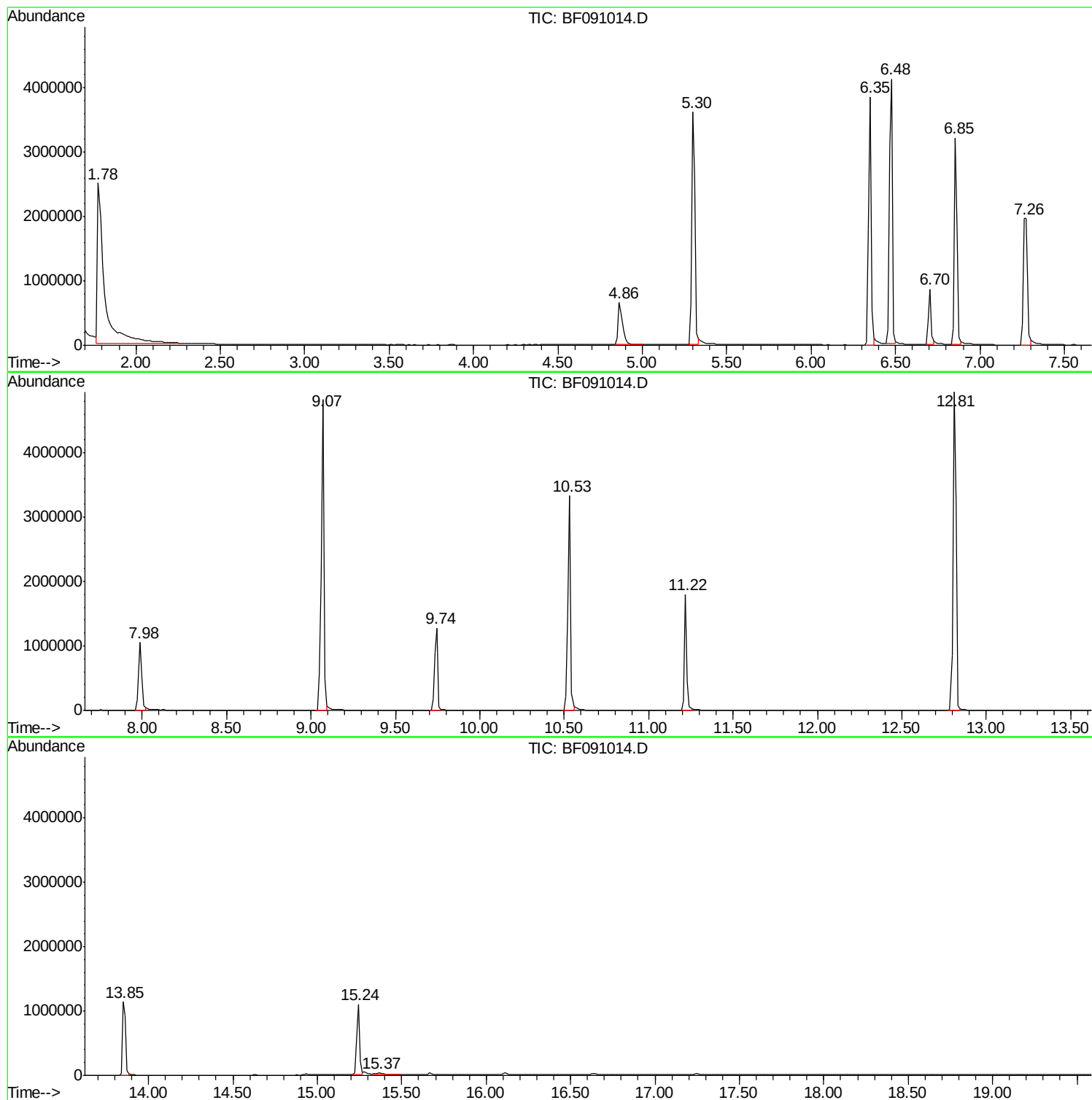
Sum of corrected areas: 53192232

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.86	23.59 ng	1150560	1,4-Dichlorobenzene-d4	6.70	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25

Peak Number 2 unknown6.48 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.48	106.57 ng	5197070	1,4-Dichlorobenzene-d4	6.70	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3	Amphetaminil	250	C17H18N2	017590-01-1	10
4	Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.86	23.6	ng	1150560	1	6.70	975333	20.0
unknown6.48	6.48	106.6	ng	5197070	1	6.70	975333	20.0