

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020717\
 Data File : BF092833.D
 Acq On : 7 Feb 2017 17:34
 Operator : SJ/MA
 Sample : PB96777BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96777BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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	5.093	260	263	273	rVB	1012351	1606030	39.57%	4.962%
2	5.516	297	300	302	rBV	4066734	4058588	100.00%	12.539%
3	6.544	387	390	392	rBV	3748938	3738770	92.12%	11.551%
4	6.670	399	401	404	rBV	3315675	3766571	92.80%	11.637%
5	6.899	419	421	424	rBV	1009871	996895	24.56%	3.080%
6	7.059	432	435	438	rBV	3339596	2947036	72.61%	9.105%
7	7.470	468	471	473	rBV	2776378	2656221	65.45%	8.206%
8	8.190	531	534	536	rBV	1395276	1202920	29.64%	3.716%
9	9.265	625	628	630	rBV	3890942	3462123	85.30%	10.696%
10	9.939	685	687	690	rVB	986113	1150647	28.35%	3.555%
11	10.739	754	757	759	rBV2	1194286	1672372	41.21%	5.167%
12	11.425	815	817	820	rBV	682113	931220	22.94%	2.877%
13	13.014	954	956	959	rBV	2376028	2539978	62.58%	7.847%
14	14.065	1046	1048	1051	rBV	742055	843581	20.79%	2.606%
15	15.517	1172	1175	1184	rBV	478243	794902	19.59%	2.456%

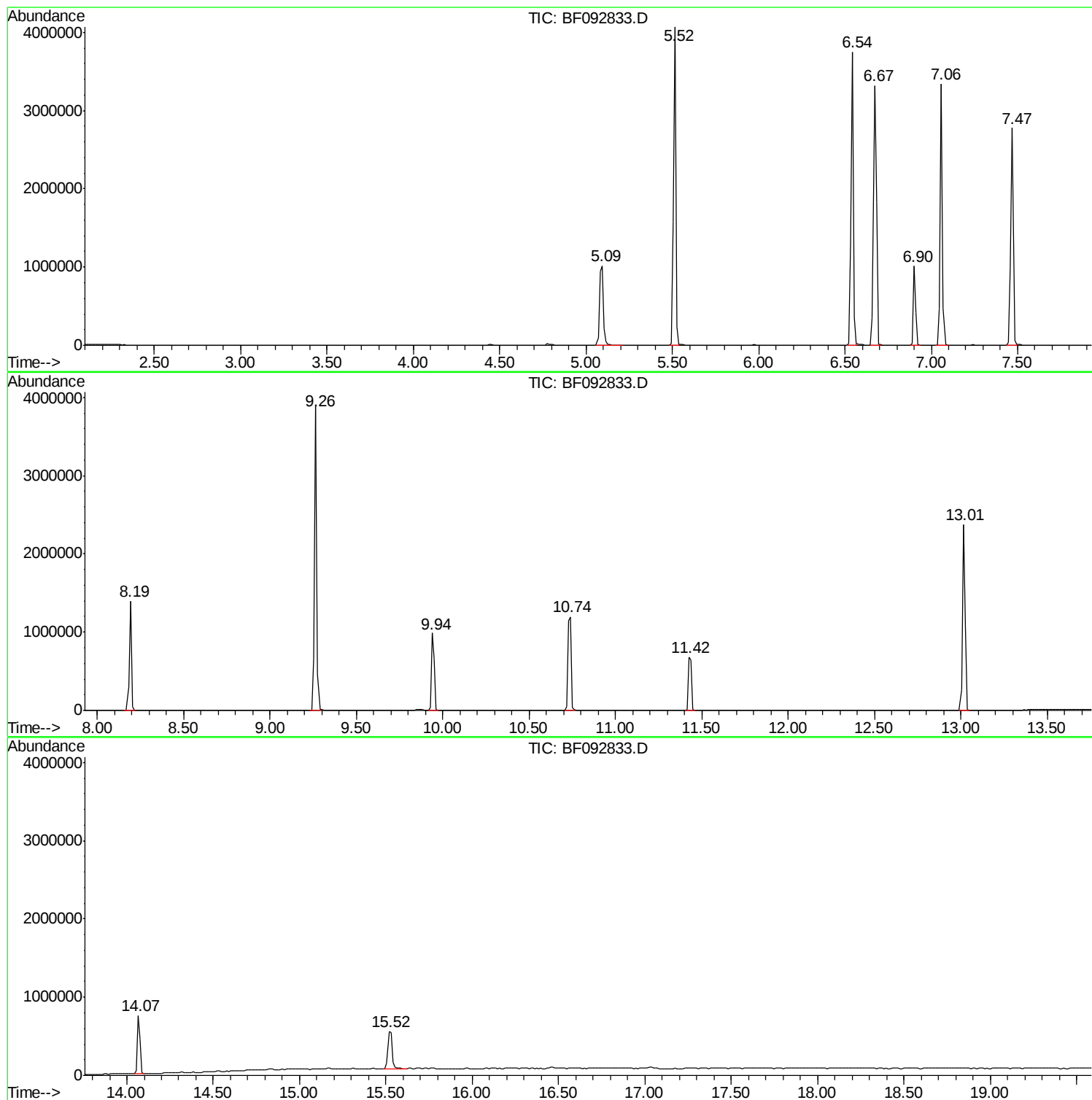
Sum of corrected areas: 32367854

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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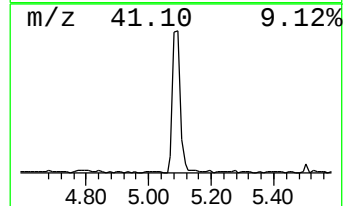
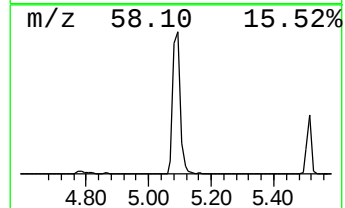
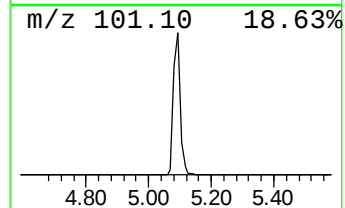
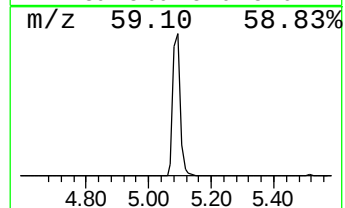
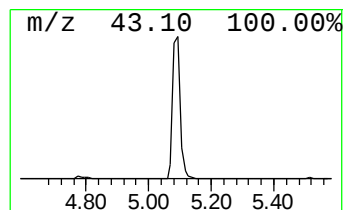
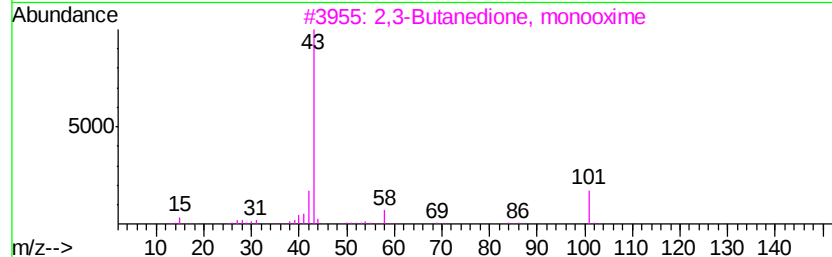
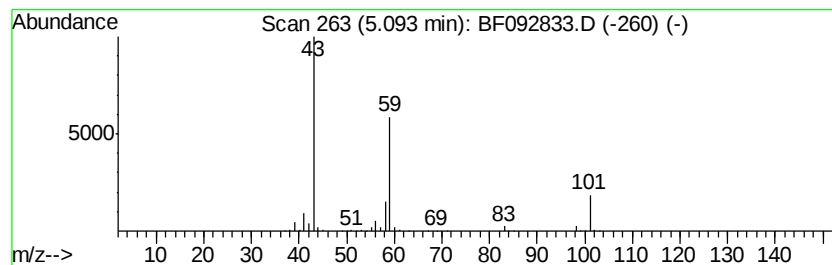
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	32.22 ng	1606030	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9
5		1,2-Butanediol, (.+/-.)-	90	C4H10O2	026171-83-5	9



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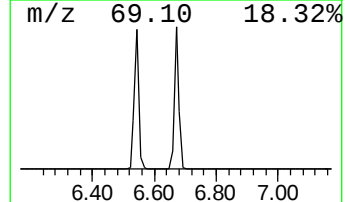
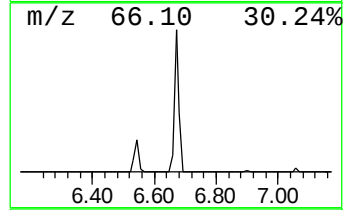
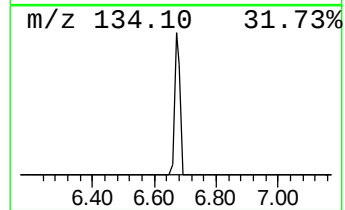
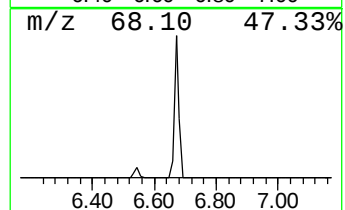
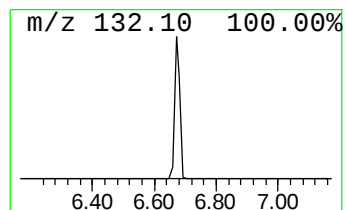
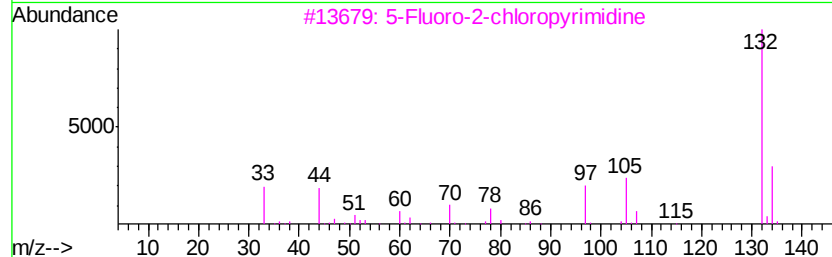
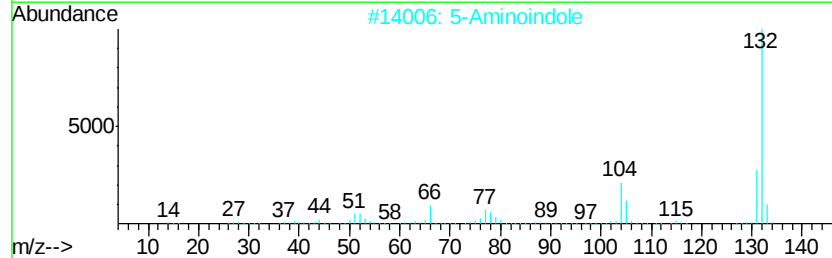
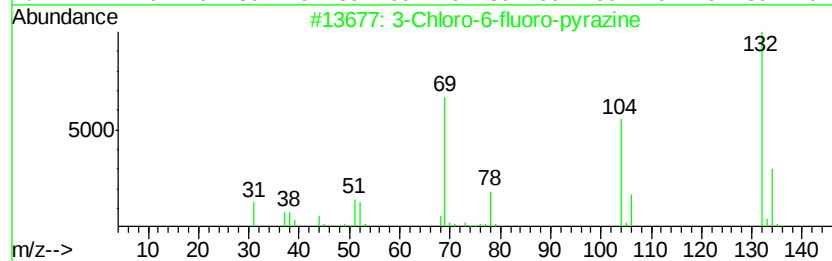
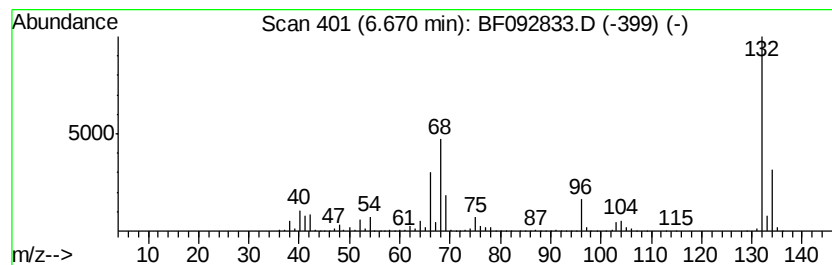
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Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	75.57 ng	3766570	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5			Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	32.2	ng	1606030	1	6.90	996895	20.0
unknown6.67	6.67	75.6	ng	3766570	1	6.90	996895	20.0