

Data Path : Z:\HPCHEM1\BNA_F\Data\BF033117\
 Data File : BF094092.D
 Acq On : 31 Mar 2017 13:26
 Operator : SJ/MA
 Sample : PB97953BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB97953BL

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-BF032217.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	4.057	364	369	379	rVB	317171	425343	7.80%	1.046%
2	4.446	429	435	438	rBV	4130344	4396764	80.58%	10.815%
3	5.504	608	615	618	rBV	3491427	4497800	82.43%	11.064%
4	5.645	633	639	642	rBV	4096650	4604872	84.39%	11.327%
5	5.898	677	682	685	rBV	985563	978889	17.94%	2.408%
6	6.075	707	712	716	rBV	3336196	3667654	67.22%	9.022%
7	6.545	785	792	795	rBV	2374735	2959859	54.24%	7.281%
8	7.398	931	937	941	rBV	1238440	1332918	24.43%	3.279%
9	8.728	1155	1163	1166	rBV	4193168	5456511	100.00%	13.422%
10	9.539	1294	1301	1306	rBV2	1305909	1440965	26.41%	3.545%
11	10.516	1459	1467	1470	rBV	2382788	2778218	50.92%	6.834%
12	11.363	1604	1611	1615	rBV	1299992	1393047	25.53%	3.427%
13	13.351	1939	1949	1952	rBV	3973143	4739025	86.85%	11.657%
14	14.615	2158	2164	2178	rVB	1081857	1164562	21.34%	2.865%
15	16.245	2436	2441	2450	rVB	674072	816638	14.97%	2.009%

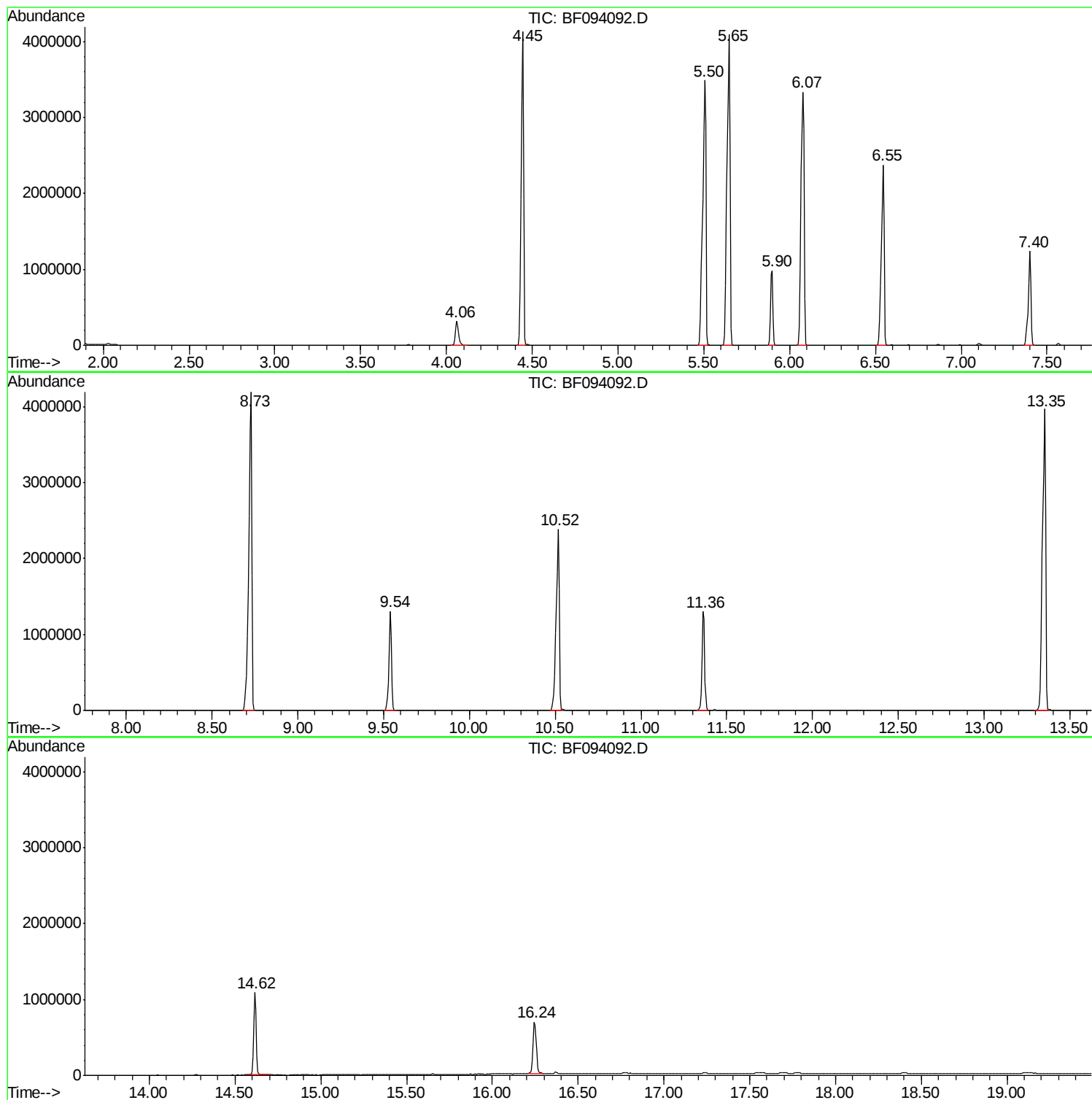
Sum of corrected areas: 40653065

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.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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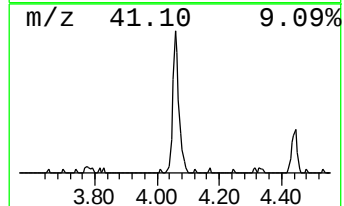
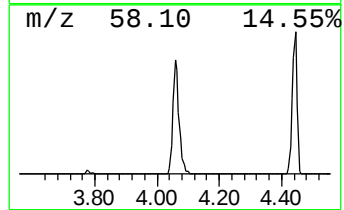
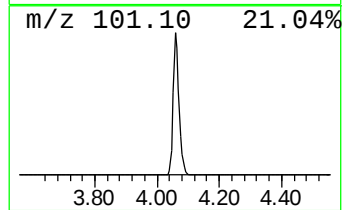
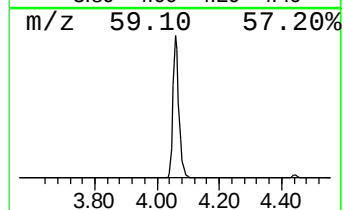
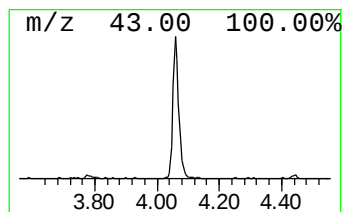
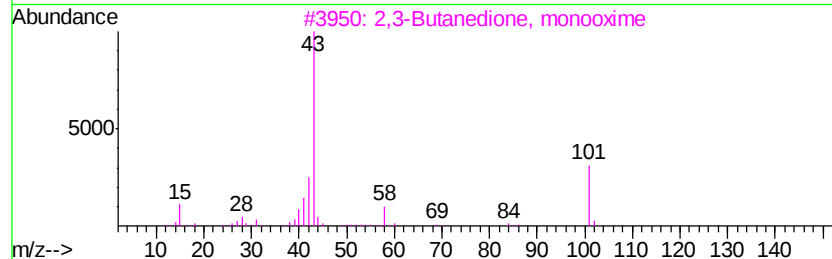
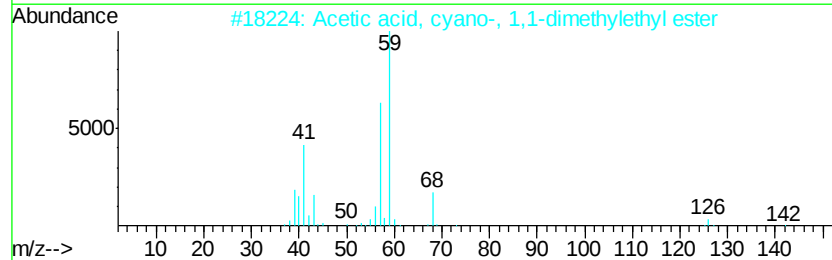
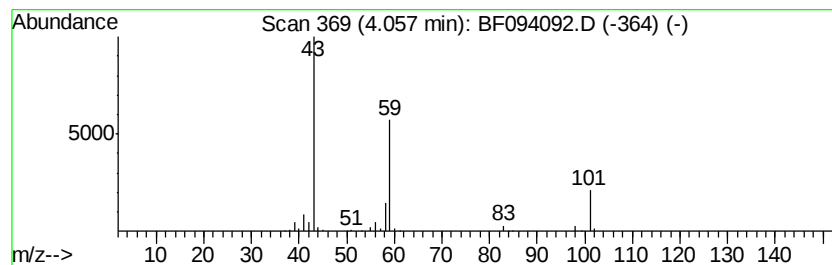
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.06	8.69 ng	425343	1,4-Dichlorobenzene-d4	5.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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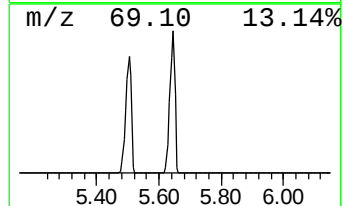
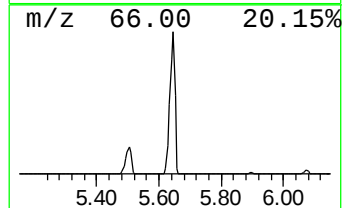
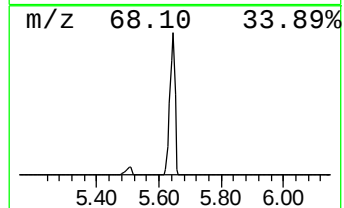
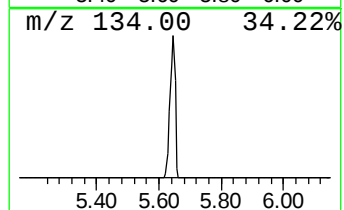
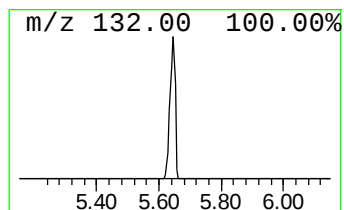
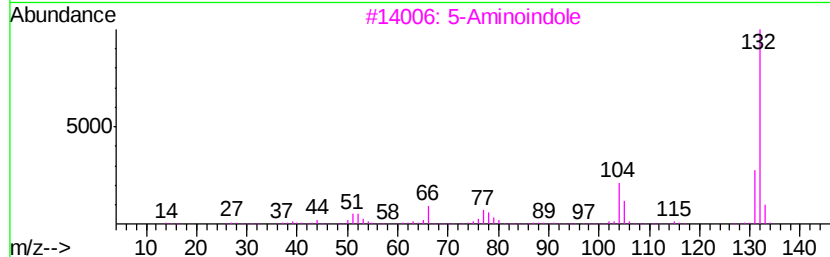
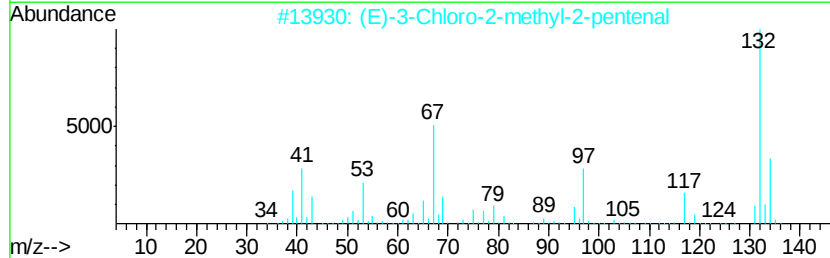
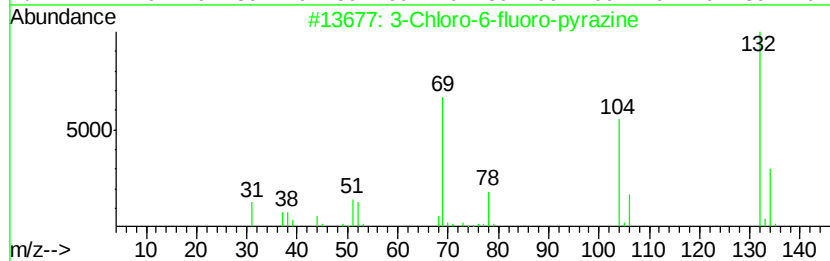
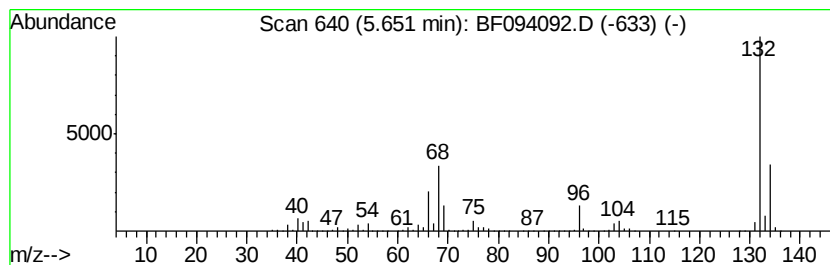
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Peak Number 2 unknown5.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.65	94.08 ng	4604870	1,4-Dichlorobenzene-d4	5.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-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...	4.06	8.7	ng	425343	1	5.90	978889	20.0
unknown5.65	5.65	94.1	ng	4604870	1	5.90	978889	20.0