

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122415\
 Data File : BF084088.D
 Acq On : 24 Dec 2015 19:31
 Operator : UM/SJ
 Sample : PB87482BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87482BL

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-BF122415.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	4.979	251	253	261	rBV	124941	158458	12.75%	1.701%
2	5.413	290	291	294	rBV	587942	849698	68.37%	9.121%
3	6.453	380	382	384	rBV	682092	875830	70.48%	9.401%
4	6.567	390	392	395	rBV	1051385	898256	72.28%	9.642%
5	6.796	410	412	415	rBV	336920	276924	22.28%	2.973%
6	6.956	424	426	428	rBV	722811	779555	62.73%	8.368%
7	7.356	459	461	464	rBV	438272	529076	42.57%	5.679%
8	8.076	522	524	527	rBV	266280	351772	28.31%	3.776%
9	9.162	616	619	621	rBV	1162318	1242710	100.00%	13.340%
10	9.836	675	678	680	rBV	433189	422259	33.98%	4.533%
11	10.625	745	747	749	rBV	703026	707485	56.93%	7.594%
12	11.322	806	808	810	rBV	325848	417632	33.61%	4.483%
13	12.534	912	914	917	rBB	11079	16158	1.30%	0.173%
14	12.899	944	946	948	rBV	1363812	1186629	95.49%	12.738%
15	13.814	1024	1026	1030	rVB	12779	14454	1.16%	0.155%
16	13.951	1036	1038	1041	rBV	333457	351698	28.30%	3.775%
17	15.368	1160	1162	1172	rVB	144421	237347	19.10%	2.548%

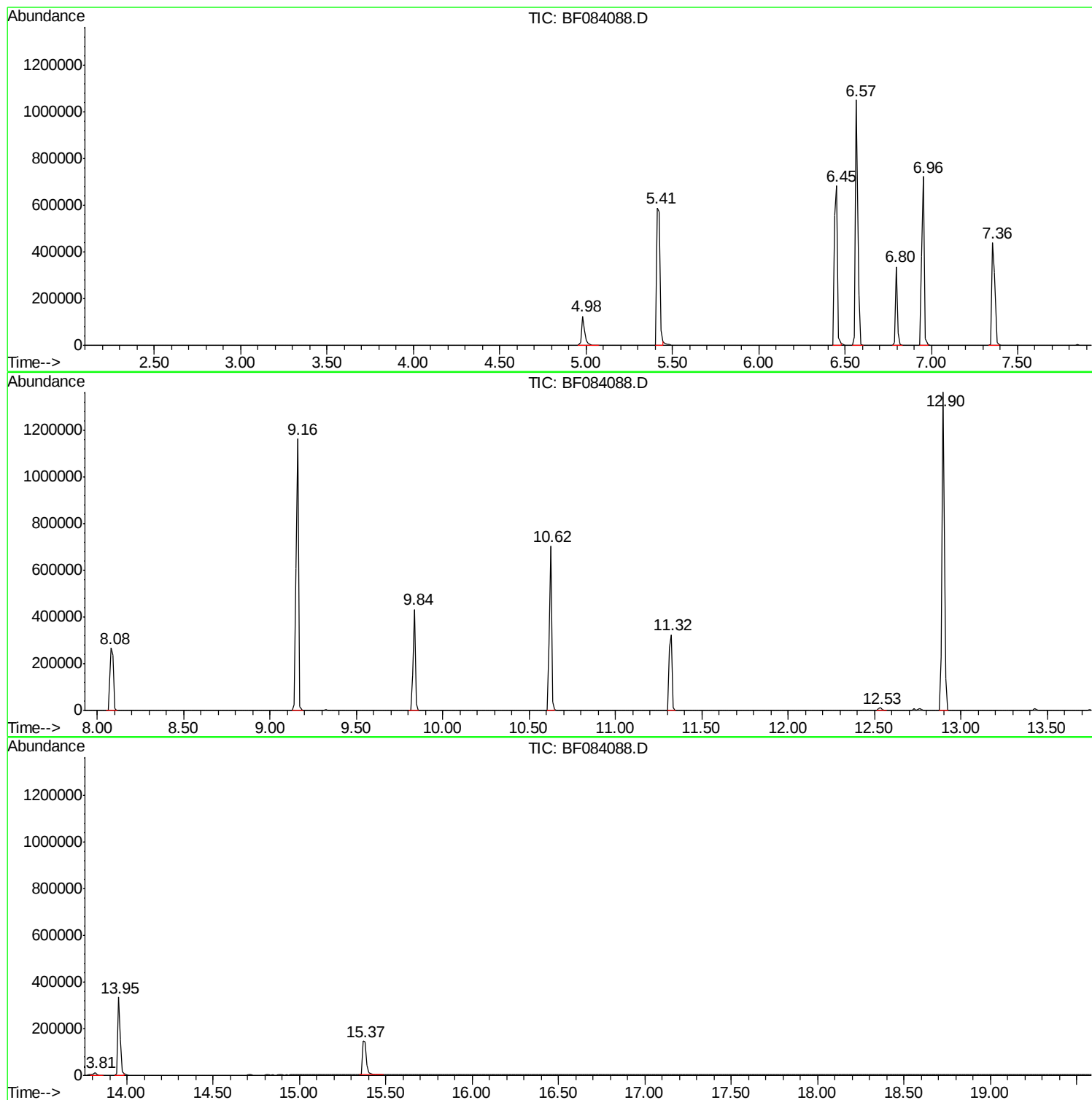
Sum of corrected areas: 9315941

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122415.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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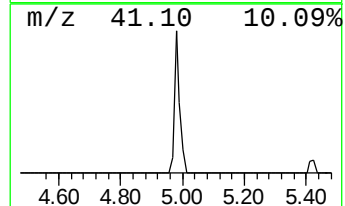
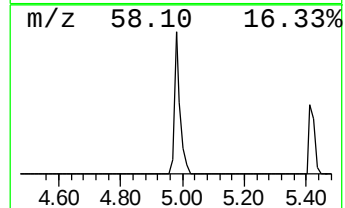
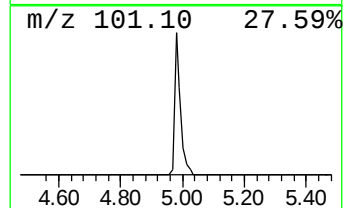
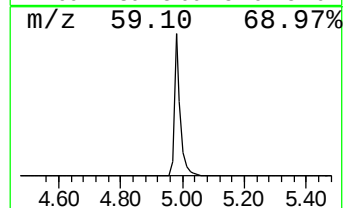
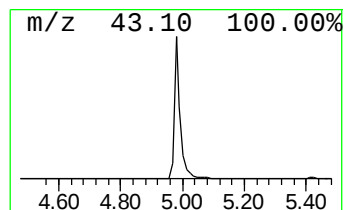
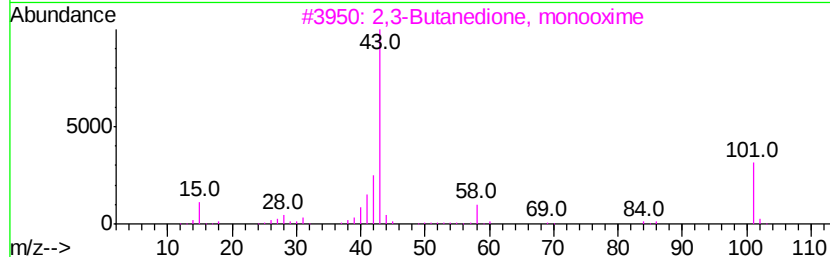
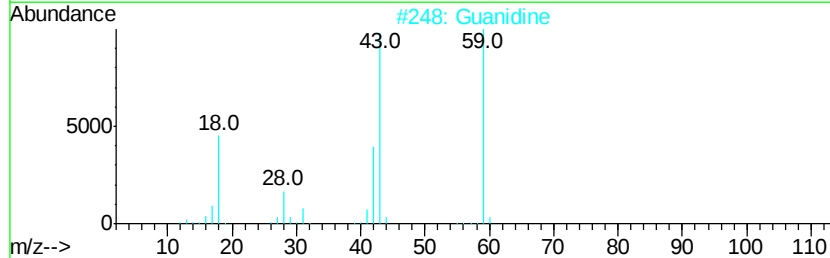
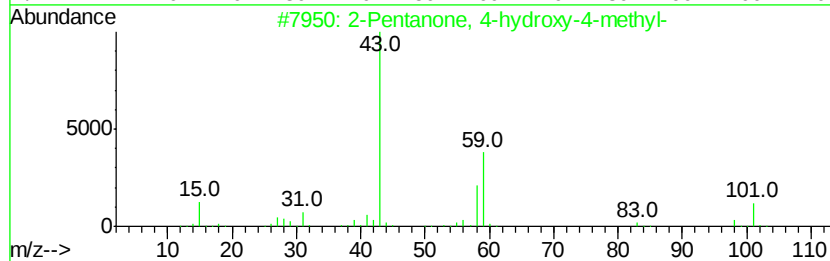
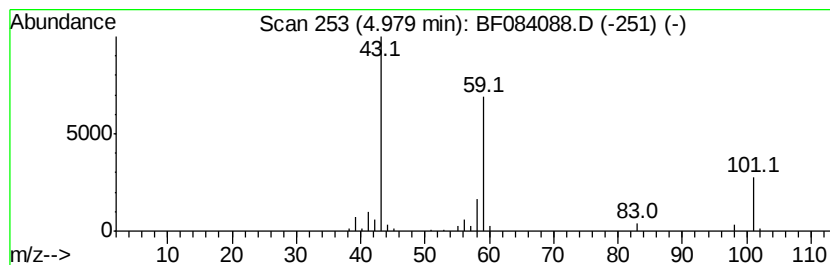
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	11.44 ng	158458	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Guanidine	59	CH5N3	000113-00-8	43
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12



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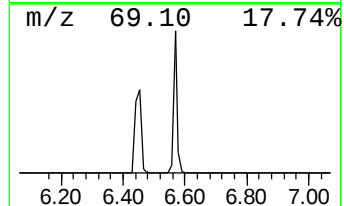
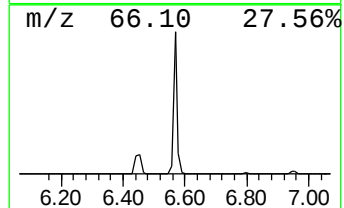
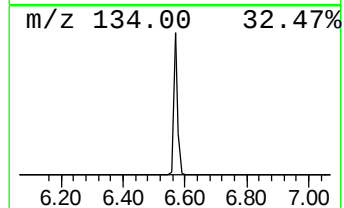
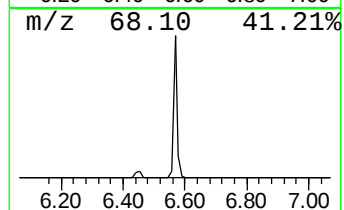
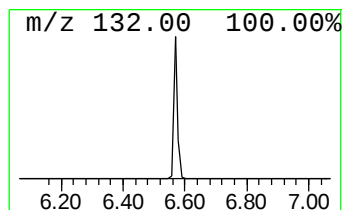
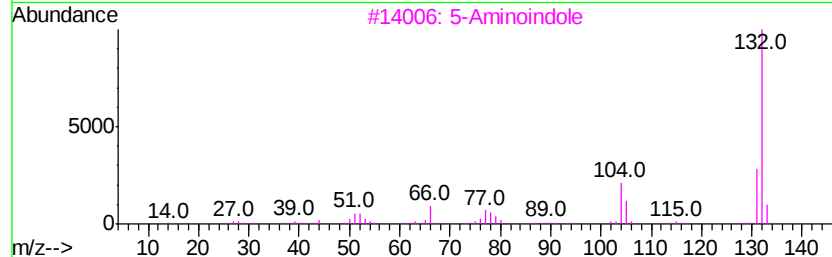
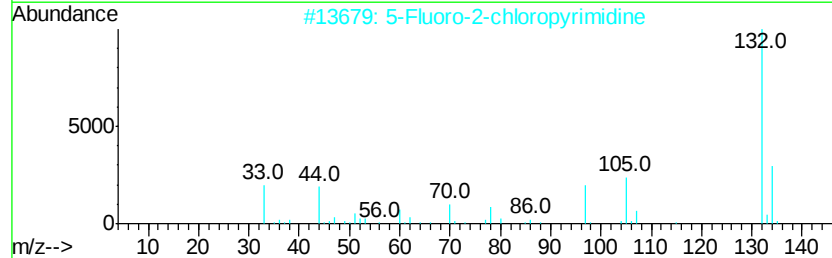
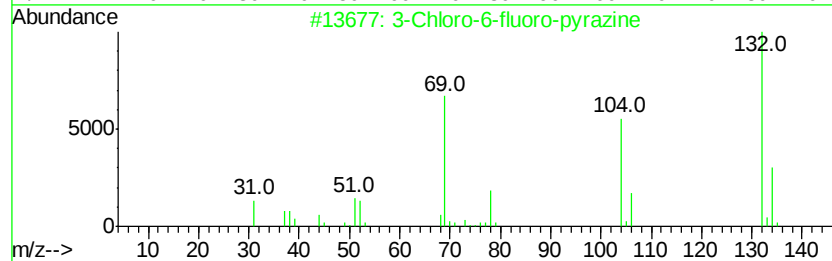
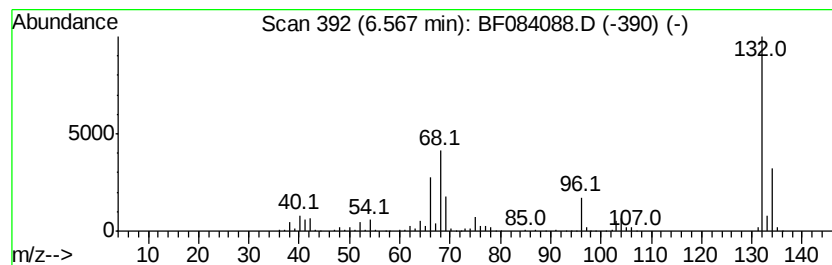
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Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	64.87 ng	898256	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.98	11.4	ng	158458	1	6.80	276924	20.0
unknown6.57	6.57	64.9	ng	898256	1	6.80	276924	20.0