

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090215\
 Data File : BF081341.D
 Acq On : 2 Sep 2015 11:19
 Operator : UM/IZ
 Sample : PB85343BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85343BL

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-BF090115.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	4.456	248	251	266	rBV	619537	1076857	57.94%	6.974%
2	4.947	291	294	296	rBV	1245643	1682635	90.53%	10.898%
3	6.044	388	390	393	rBV	1189218	1684803	90.65%	10.912%
4	6.147	397	399	402	rBV	1276624	1552876	83.55%	10.057%
5	6.387	418	420	422	rBV	409580	348123	18.73%	2.255%
6	6.547	431	434	436	rBV	907454	1218681	65.57%	7.893%
7	6.959	468	470	473	rBV	898527	946016	50.90%	6.127%
8	7.679	530	533	535	rBV	408522	442659	23.82%	2.867%
9	8.753	625	627	630	rBV	1296949	1783000	95.93%	11.548%
10	9.416	682	685	687	rBV	557855	493251	26.54%	3.195%
11	10.193	751	753	756	rBV2	928784	1074708	57.82%	6.960%
12	10.879	811	813	815	rBV	655701	536020	28.84%	3.472%
13	12.468	949	952	954	rBV	1961649	1858585	100.00%	12.037%
14	13.485	1039	1041	1044	rBV	494343	431279	23.20%	2.793%
15	14.800	1153	1156	1158	rBV	285233	310744	16.72%	2.013%

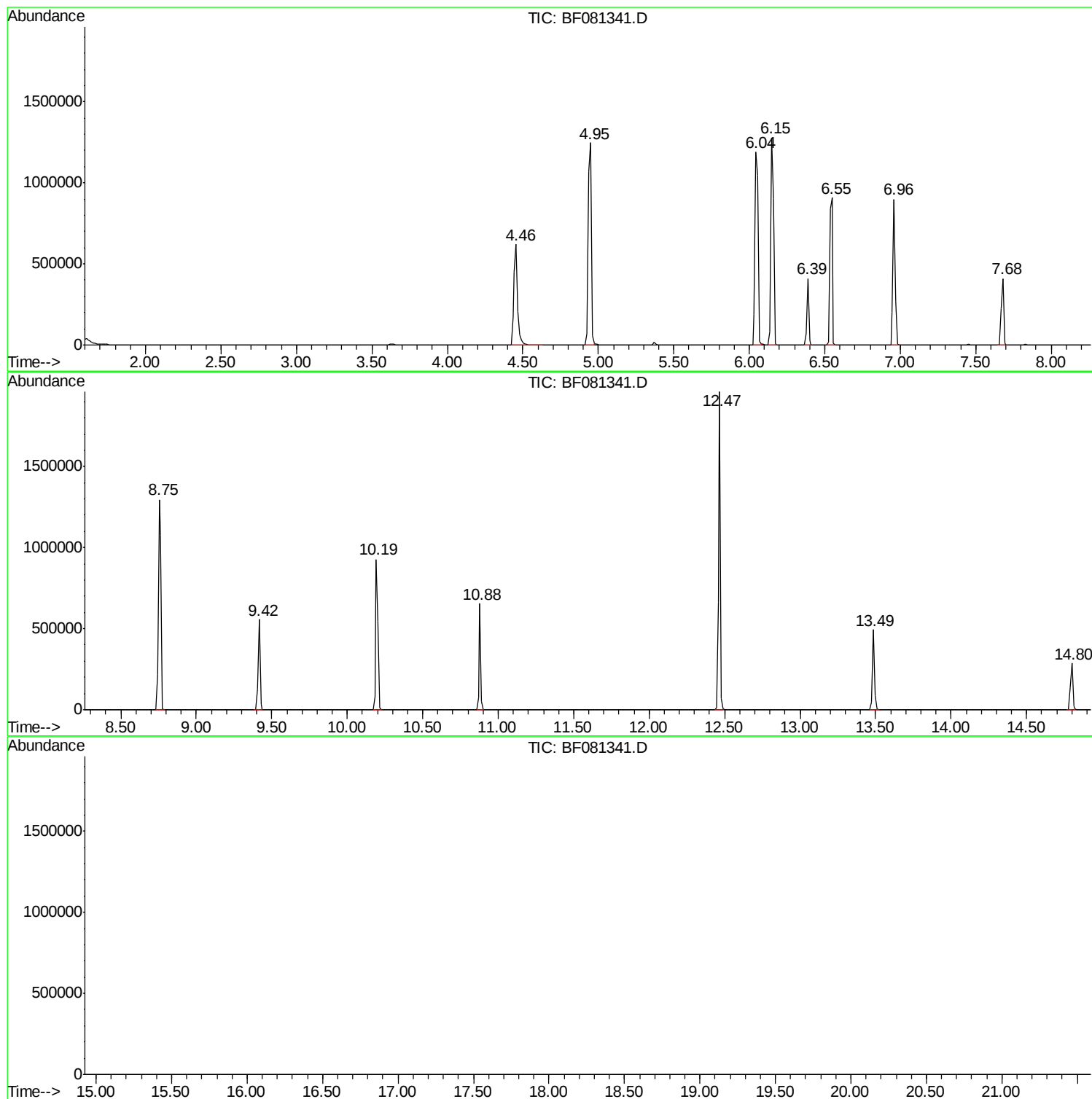
Sum of corrected areas: 15440237

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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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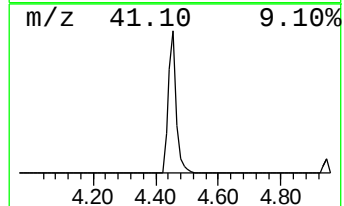
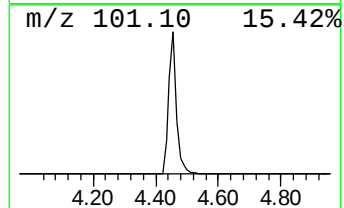
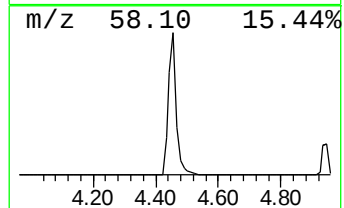
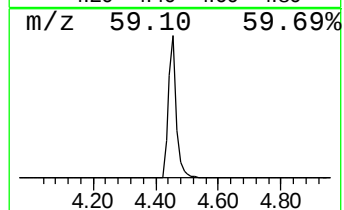
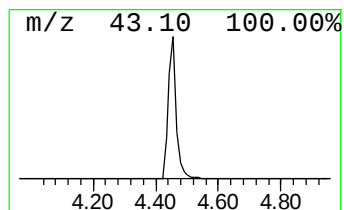
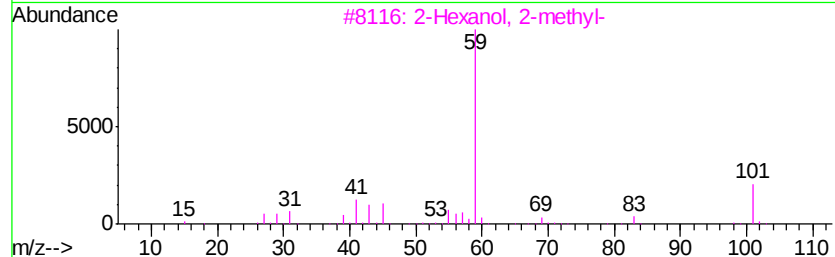
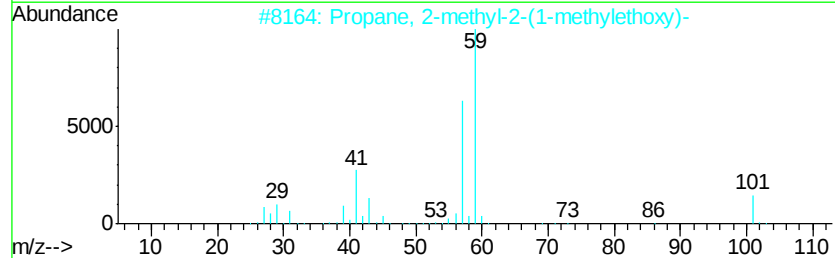
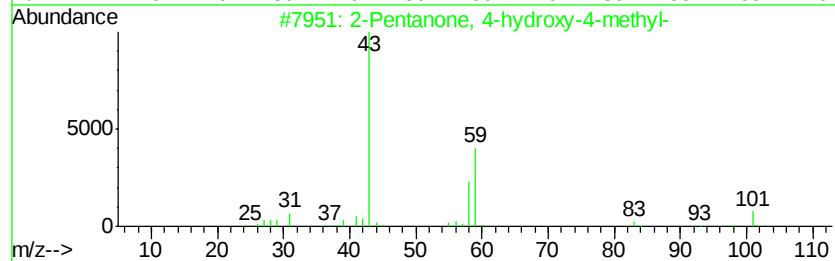
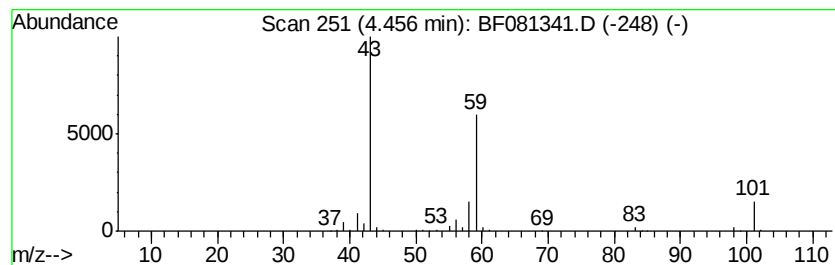
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	61.87 ng	1076860	1,4-Dichlorobenzene-d4	6.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O		017348-59-3	36
3	2-Hexanol, 2-methyl-	116	C7H16O		000625-23-0	33
4	3-Octanol	130	C8H18O		000589-98-0	33
5	3-Hexanol, 4-methyl-	116	C7H16O		000615-29-2	33



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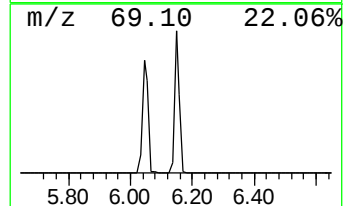
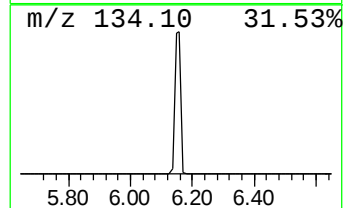
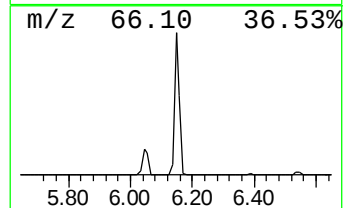
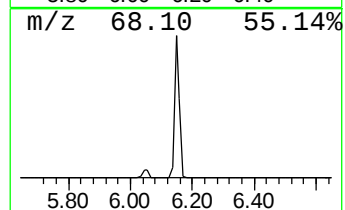
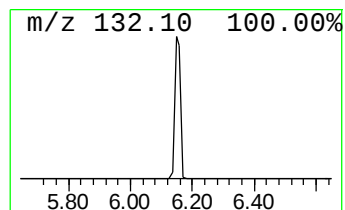
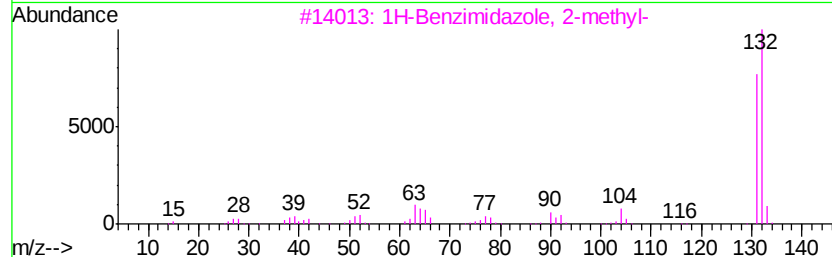
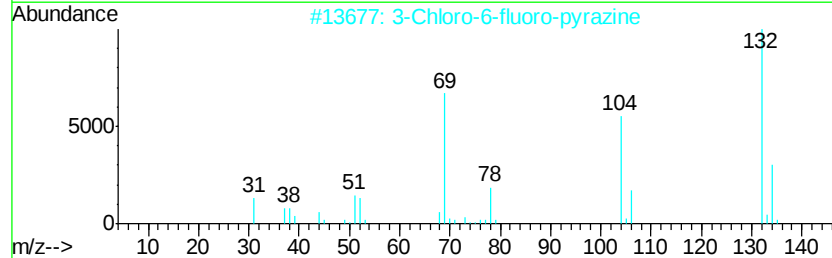
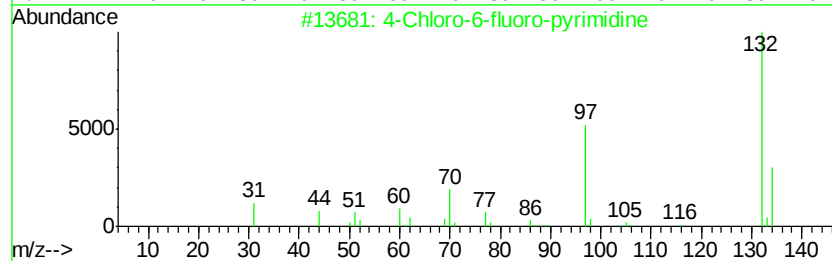
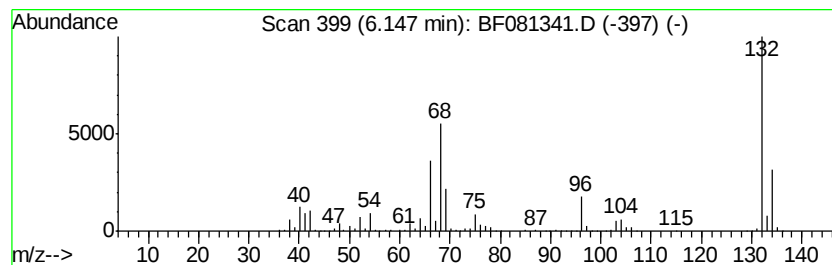
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Peak Number 2 unknown6.15 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.15	89.21 ng	1552880	1,4-Dichlorobenzene-d4	6.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.46	61.9	ng	1076860	1	6.39	348123	20.0
unknown6.15	6.15	89.2	ng	1552880	1	6.39	348123	20.0