

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081915\
 Data File : BF081085.D
 Acq On : 19 Aug 2015 17:14
 Operator : UM/IZ
 Sample : PB85071BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85071BL

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-BF081715.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	4.687	243	245	252	rBV	125431	138382	21.97%	2.802%
2	5.144	283	285	288	rBV	198260	208407	33.08%	4.220%
3	6.219	377	379	382	rBV	241000	216180	34.32%	4.378%
4	6.344	388	390	393	rBV	265095	217802	34.58%	4.411%
5	6.584	409	411	413	rBV	508602	414947	65.87%	8.403%
6	6.744	422	425	427	rBV	126160	152918	24.28%	3.097%
7	7.144	459	460	463	rBV	92011	123224	19.56%	2.495%
8	7.876	521	524	526	rBV	561787	545188	86.55%	11.040%
9	8.950	616	618	620	rBV	297875	240021	38.10%	4.860%
10	9.613	674	676	679	rBV	495849	558340	88.63%	11.306%
11	10.390	742	744	747	rBB2	99949	120531	19.13%	2.441%
12	11.088	802	805	807	rBV	735340	621460	98.65%	12.585%
13	12.665	941	943	946	rBB	273539	231923	36.82%	4.696%
14	13.705	1031	1034	1036	rBV	645253	629937	100.00%	12.756%
15	15.042	1148	1151	1154	rBV	431761	503798	79.98%	10.202%
16	15.877	1222	1224	1230	rBV4	3024	7534	1.20%	0.153%
17	16.620	1285	1289	1293	rBV4	1703	7659	1.22%	0.155%

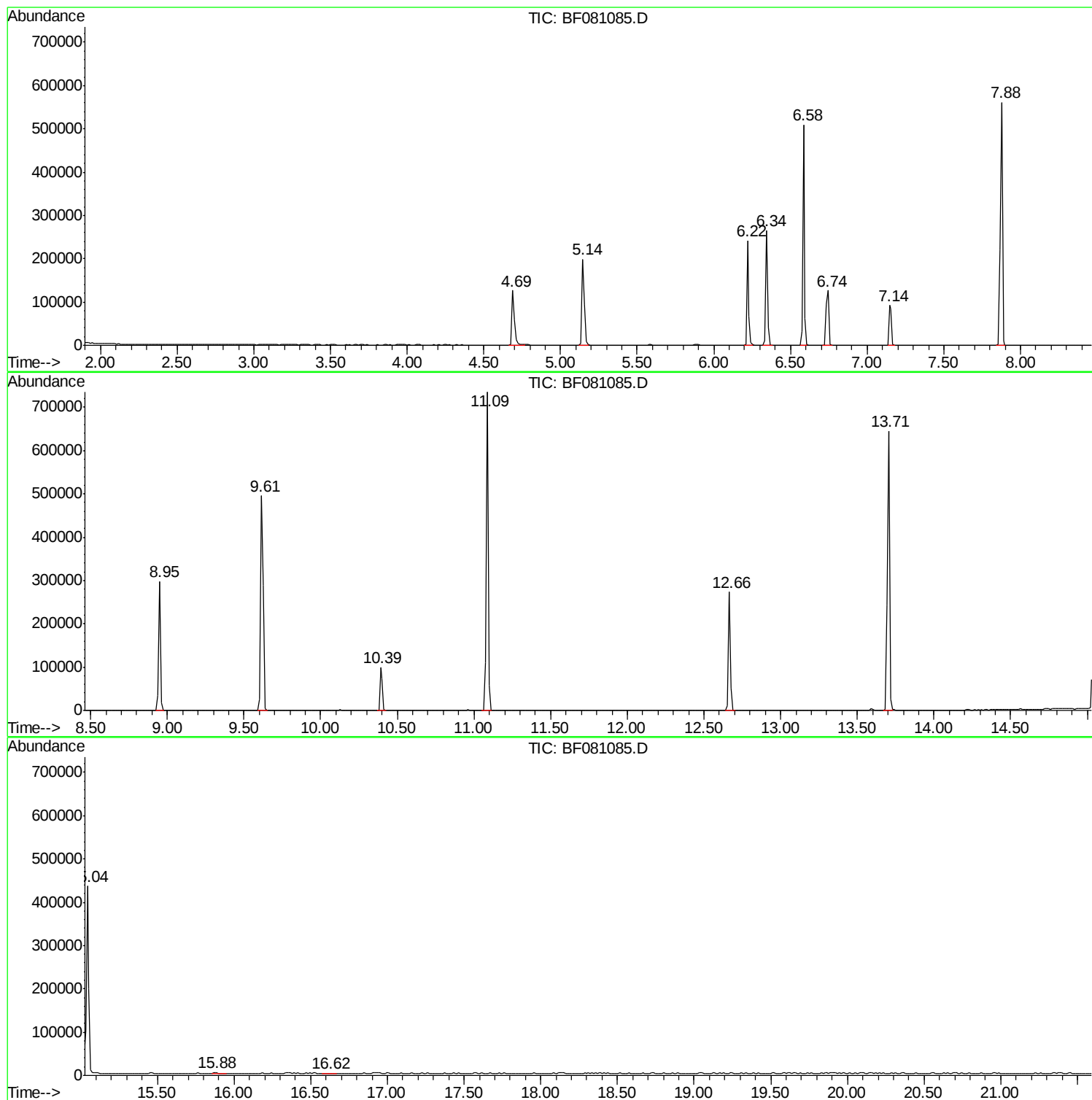
Sum of corrected areas: 4938251

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.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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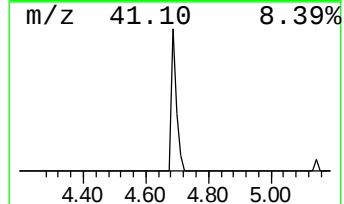
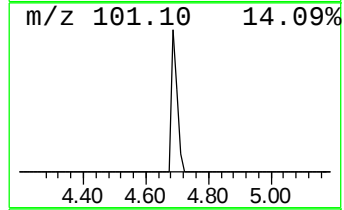
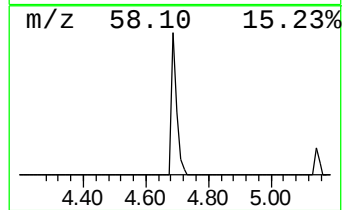
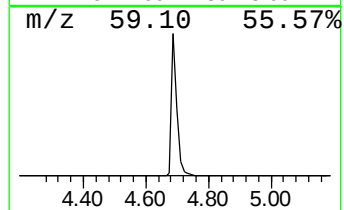
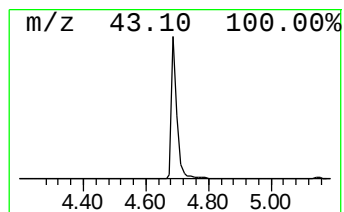
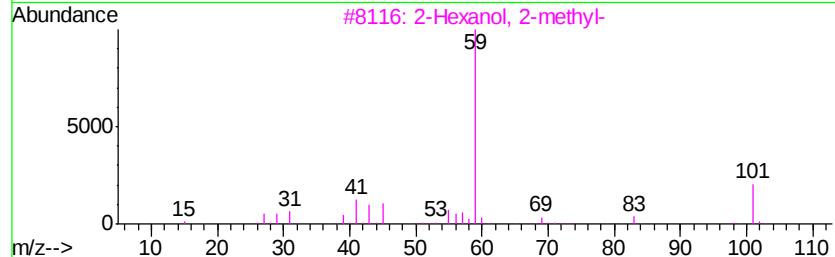
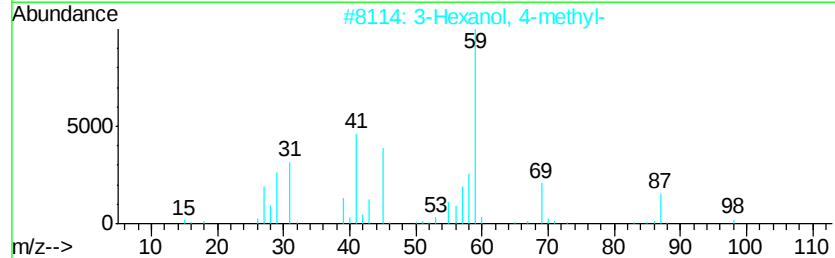
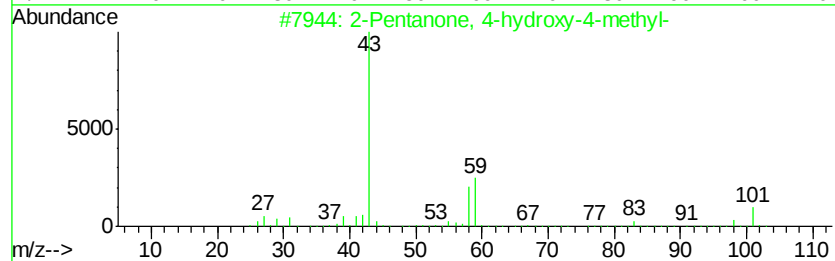
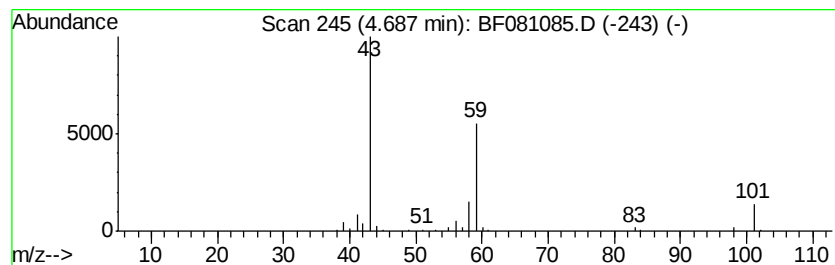
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.69	6.67 ng	138382	1,4-Dichlorobenzene-d4	6.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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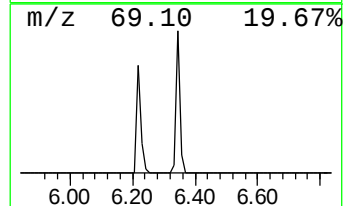
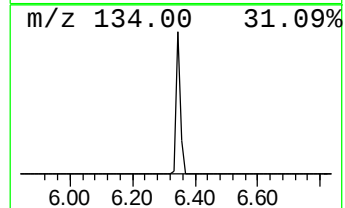
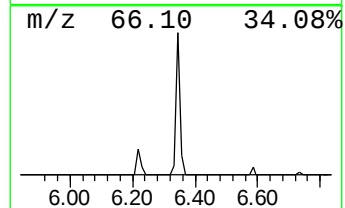
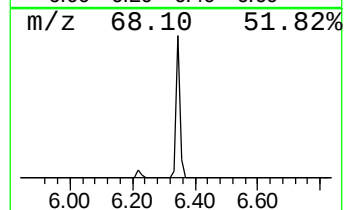
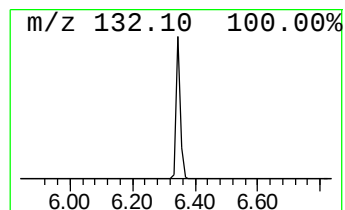
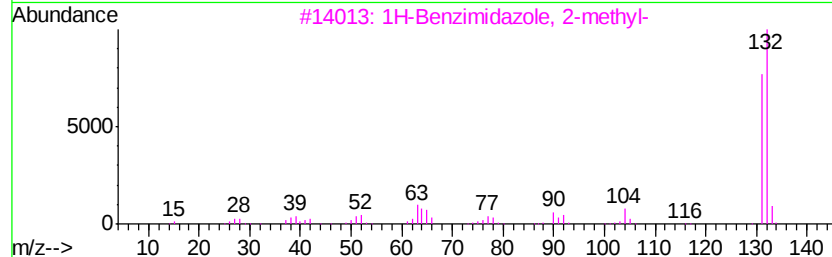
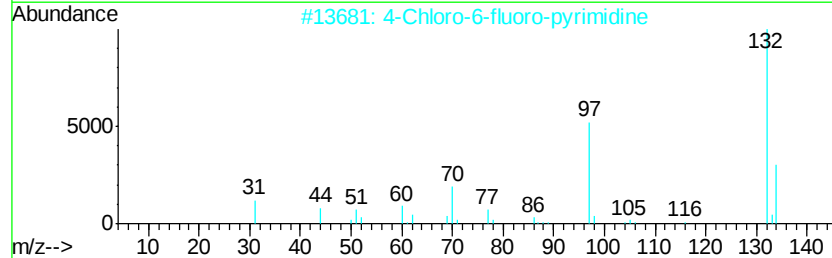
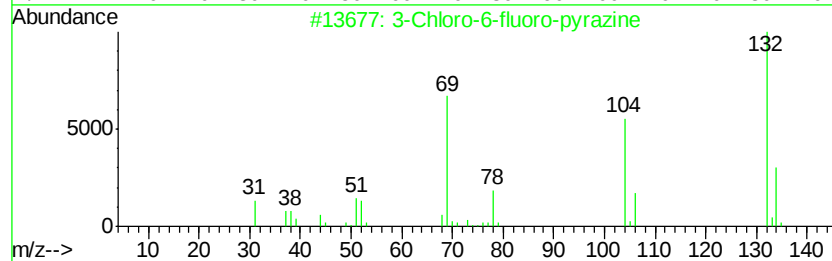
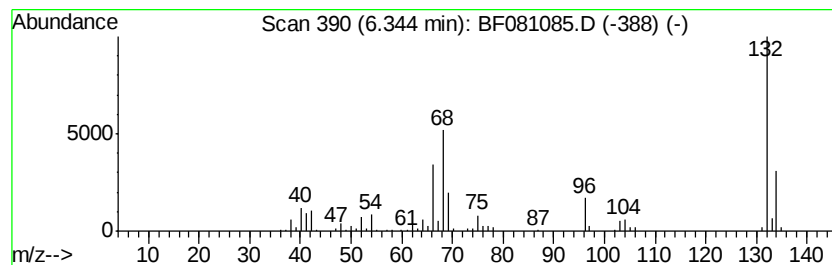
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Peak Number 2 unknown6.34 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	10.50 ng	217802	1,4-Dichlorobenzene-d4	6.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	11



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.69	6.7	ng	138382	1	6.58	414947	20.0
unknown6.34	6.34	10.5	ng	217802	1	6.58	414947	20.0