

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090115\
 Data File : BF081314.D
 Acq On : 1 Sep 2015 20:35
 Operator : UM/IZ
 Sample : PB85287BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85287BL

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	247	251	267	rBB	483394	910627	57.13%	6.729%
2	4.936	291	293	296	rBV	1040417	1383429	86.80%	10.223%
3	6.045	388	390	393	rBV	1071597	1382775	86.75%	10.218%
4	6.147	397	399	402	rBV	1098724	1321810	82.93%	9.767%
5	6.387	418	420	422	rBV	417745	358233	22.48%	2.647%
6	6.547	431	434	436	rBV	775723	1060076	66.51%	7.833%
7	6.959	468	470	473	rBV	785647	815664	51.17%	6.027%
8	7.679	530	533	535	rBV	457897	460827	28.91%	3.405%
9	8.753	625	627	630	rBV	1060138	1545403	96.96%	11.420%
10	9.416	682	685	687	rBV	570743	495969	31.12%	3.665%
11	10.194	751	753	756	rBV2	769348	896352	56.24%	6.623%
12	10.879	811	813	815	rBV	638173	527848	33.12%	3.900%
13	12.468	949	952	954	rBV	1749014	1593903	100.00%	11.778%
14	13.485	1039	1041	1044	rBV	473627	438766	27.53%	3.242%
15	14.800	1153	1156	1158	rBV	340537	341252	21.41%	2.522%

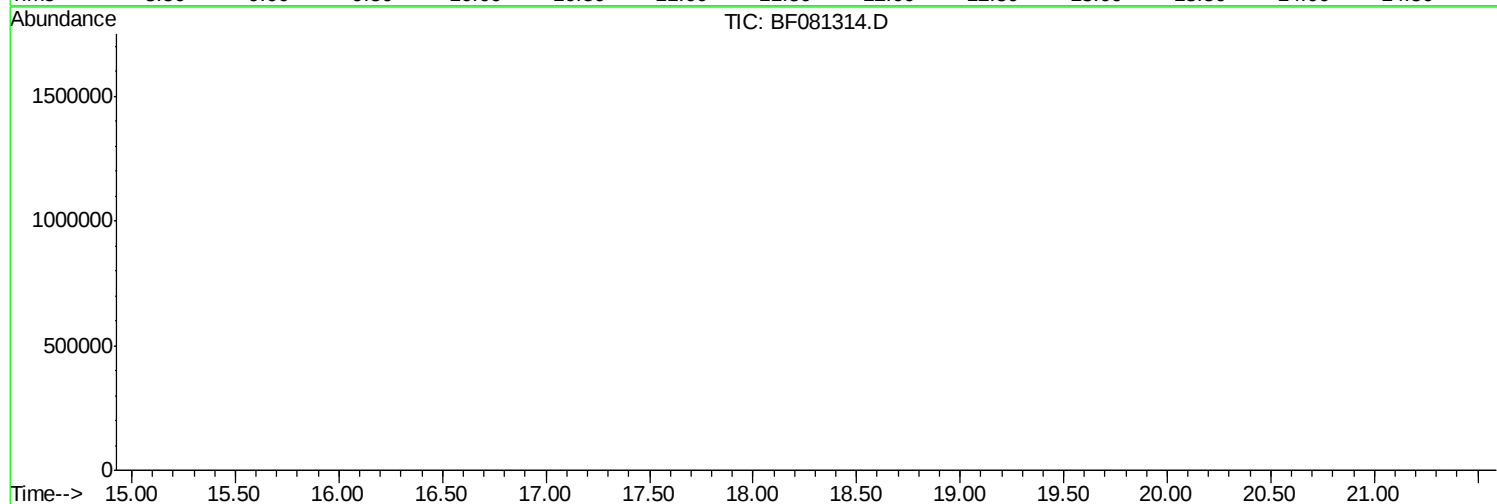
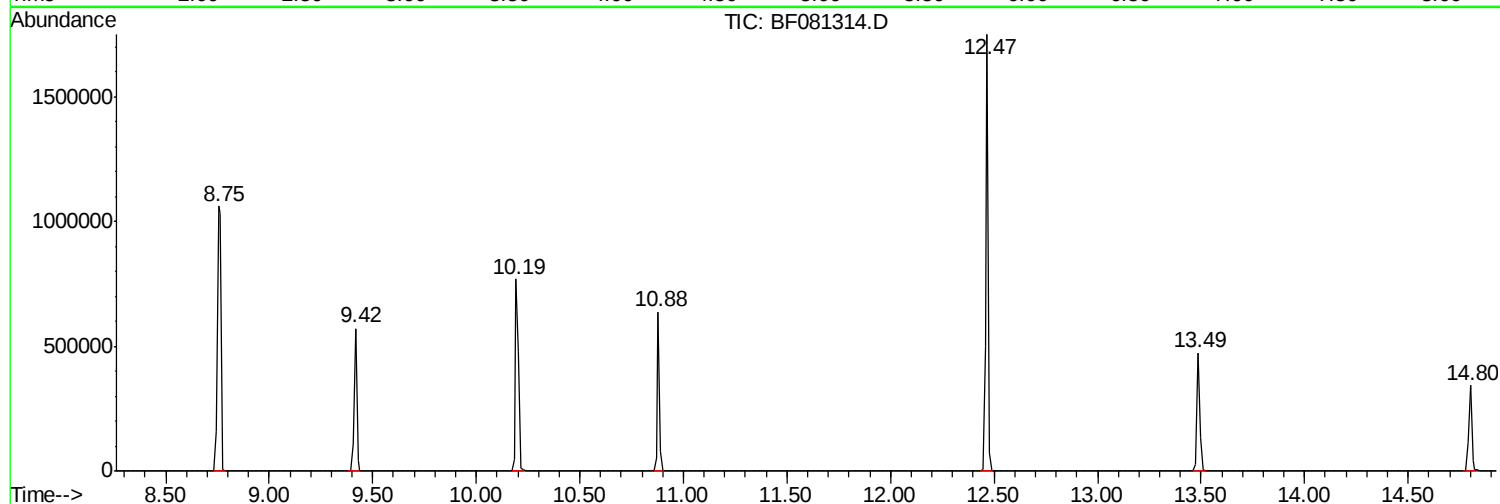
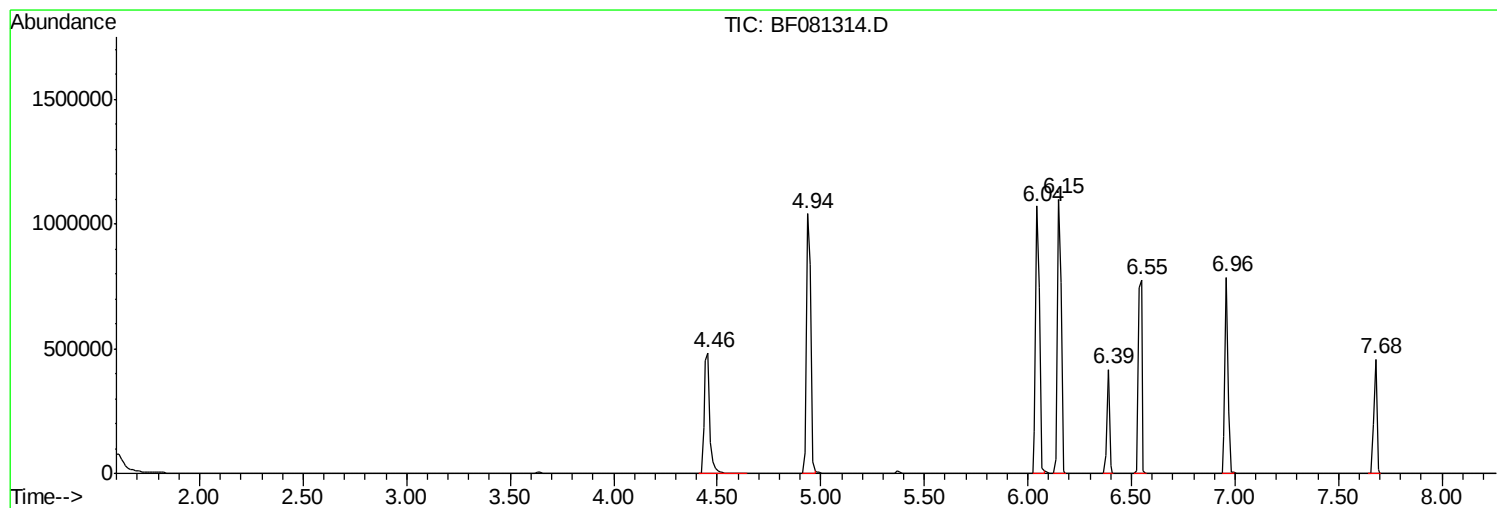
Sum of corrected areas: 13532934

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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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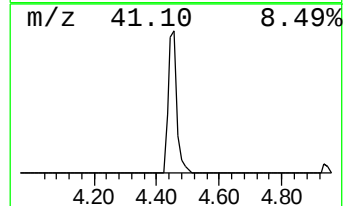
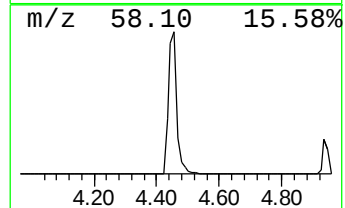
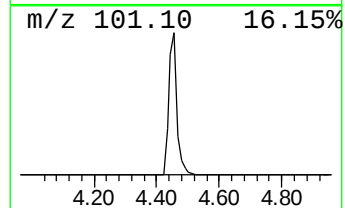
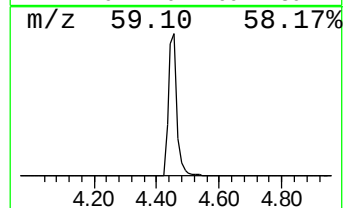
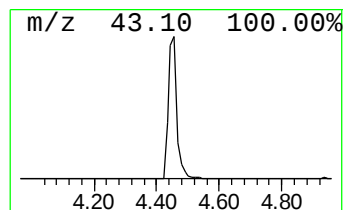
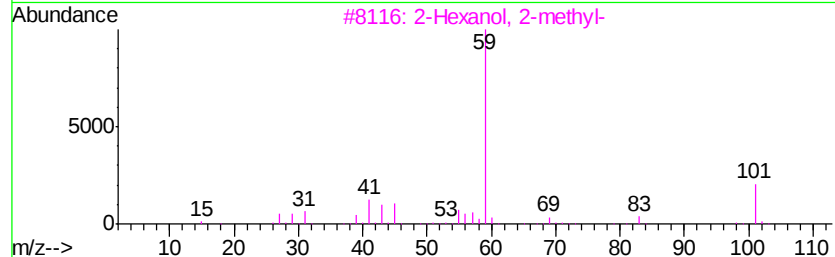
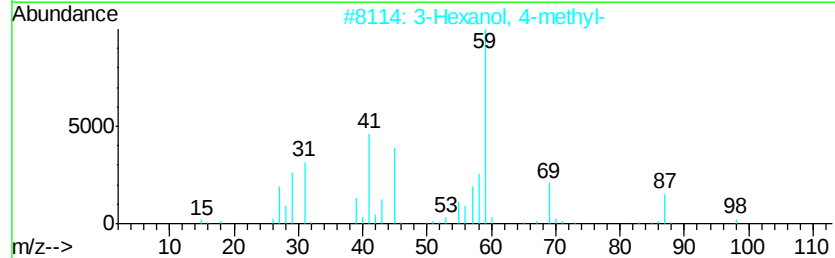
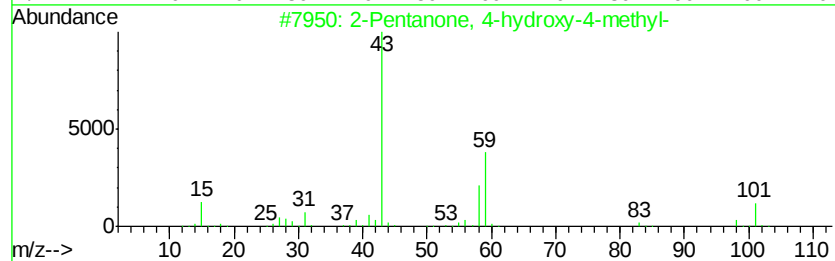
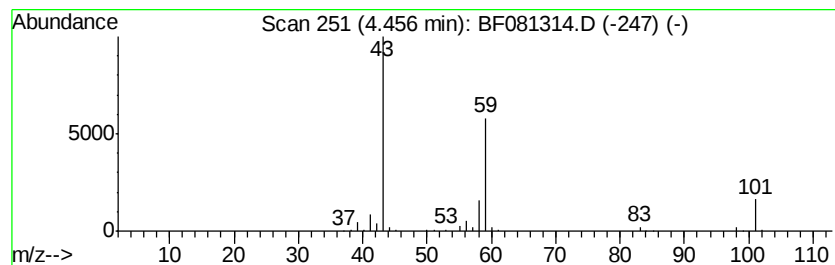
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	50.84 ng	910627	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	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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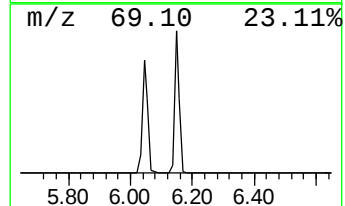
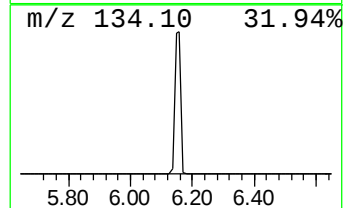
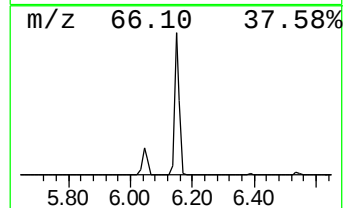
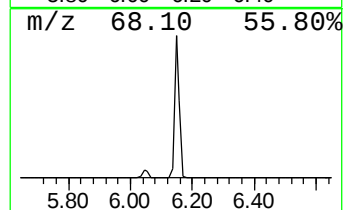
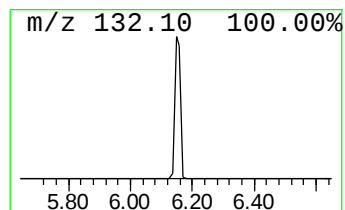
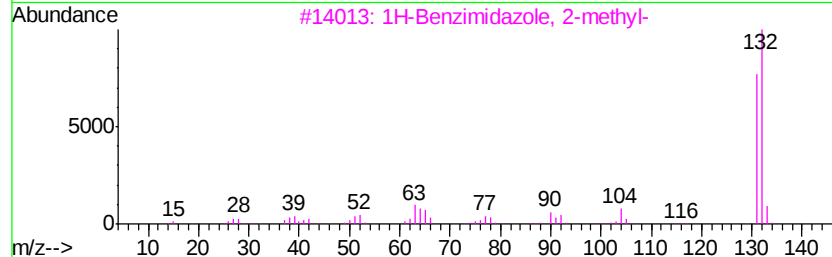
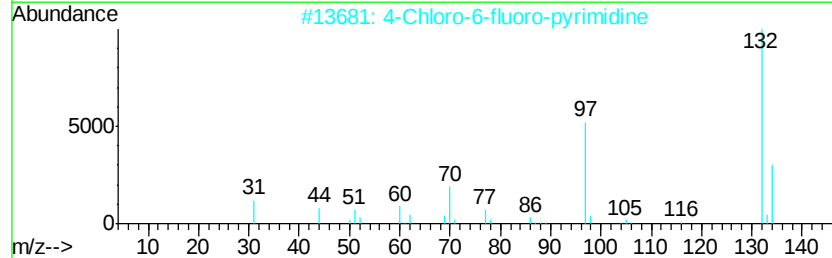
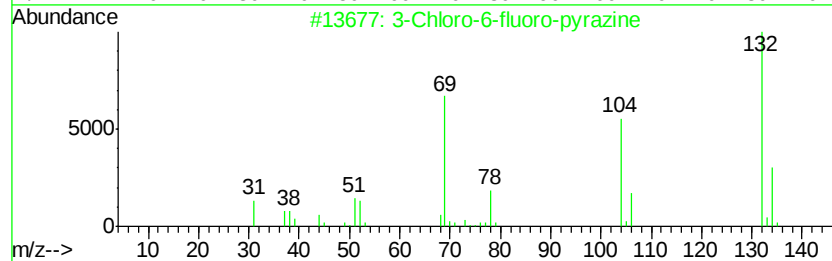
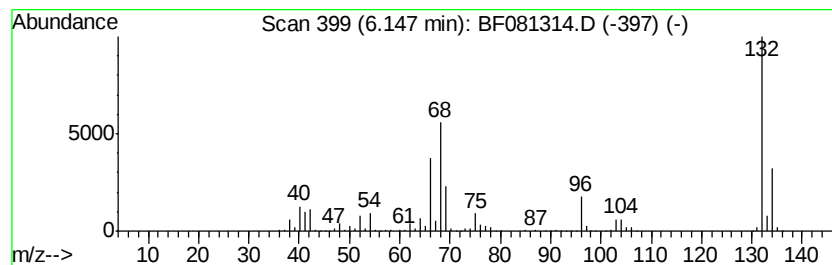
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Peak Number 2 unknown6.15 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.15	73.80 ng	1321810	1,4-Dichlorobenzene-d4	6.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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	22
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5			Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.46	50.8	ng	910627	1	6.39	358233	20.0
unknown6.15	6.15	73.8	ng	1321810	1	6.39	358233	20.0