

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090115\
 Data File : BF081312.D
 Acq On : 1 Sep 2015 19:33
 Operator : UM/IZ
 Sample : PB85326BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85326BL

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	263	rBV	609199	1102057	56.91%	6.793%
2	4.947	291	294	296	rBV	1368173	1777280	91.77%	10.956%
3	6.056	388	391	393	rBV	1614353	1931447	99.73%	11.906%
4	6.159	397	400	402	rBV	1245981	1686701	87.09%	10.397%
5	6.387	418	420	422	rBV	426678	361106	18.65%	2.226%
6	6.547	431	434	436	rBV	1064770	1295380	66.89%	7.985%
7	6.959	468	470	473	rBV	841734	1026833	53.02%	6.330%
8	7.679	530	533	535	rBV	458726	462658	23.89%	2.852%
9	8.765	624	628	630	rBV	1432324	1936643	100.00%	11.938%
10	9.416	682	685	687	rBV	571787	494817	25.55%	3.050%
11	10.193	751	753	756	rBV2	811980	1122224	57.95%	6.918%
12	10.879	811	813	816	rBV	609621	512588	26.47%	3.160%
13	12.468	949	952	955	rBV	1749267	1727039	89.18%	10.646%
14	13.485	1039	1041	1044	rBV	451388	438489	22.64%	2.703%
15	14.799	1154	1156	1159	rBV	364001	347331	17.93%	2.141%

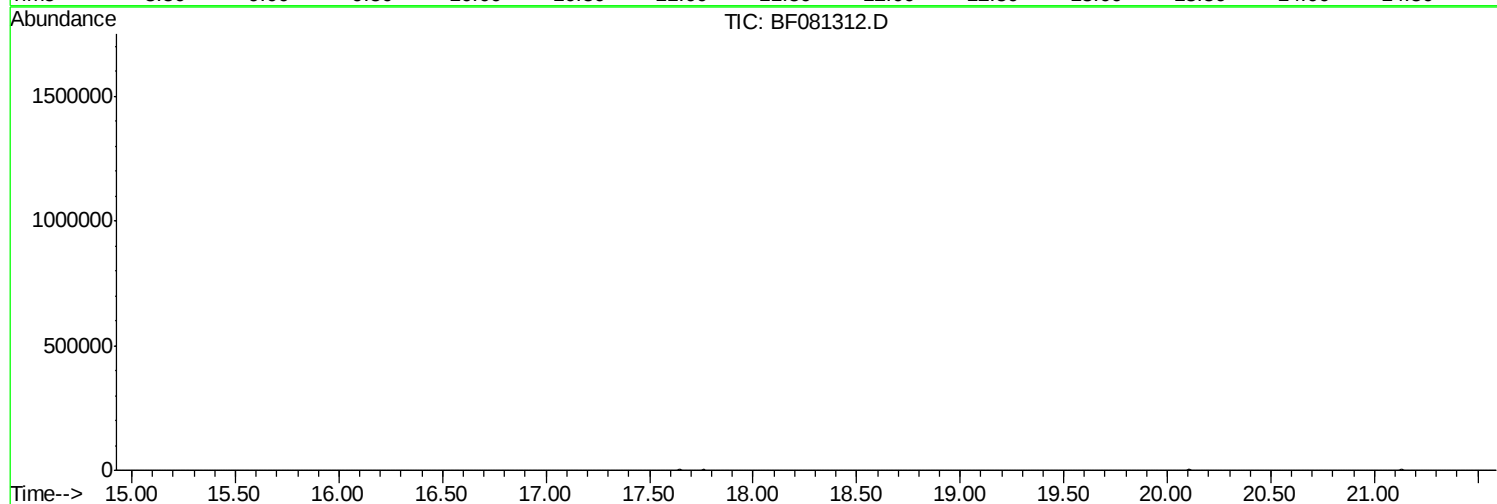
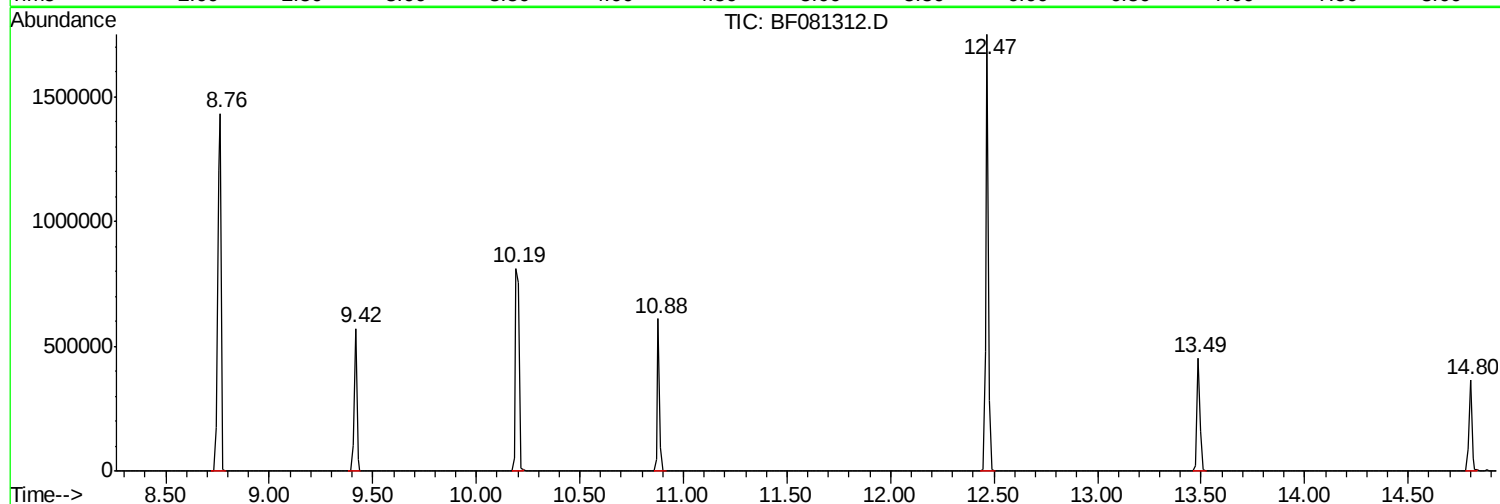
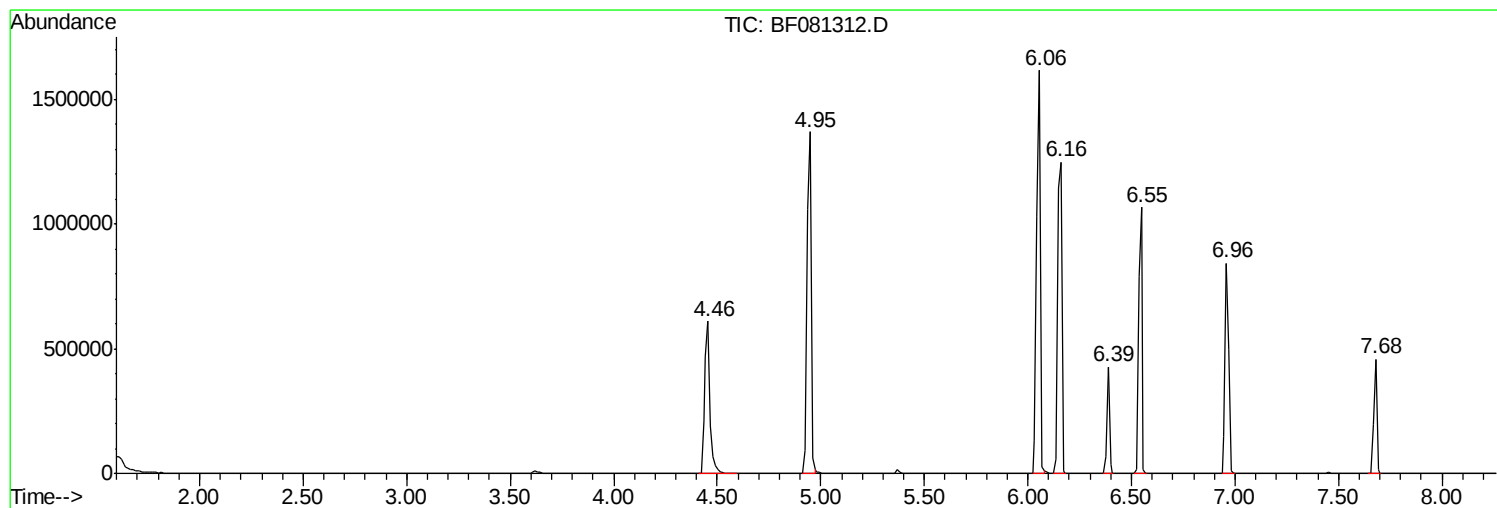
Sum of corrected areas: 16222593

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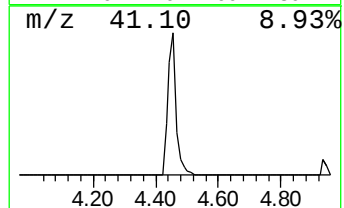
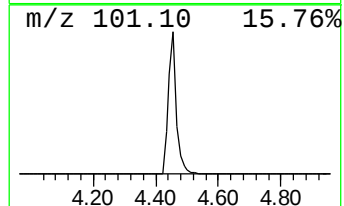
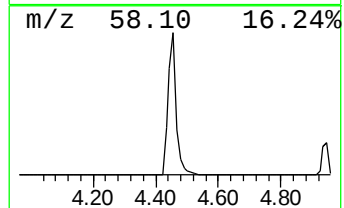
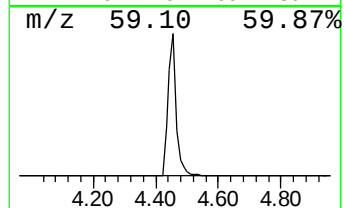
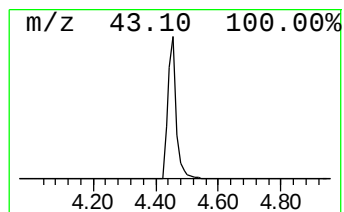
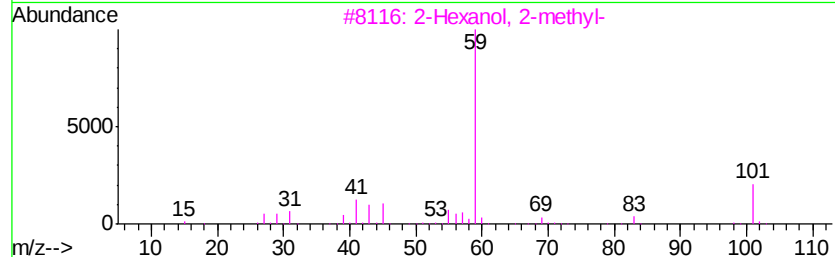
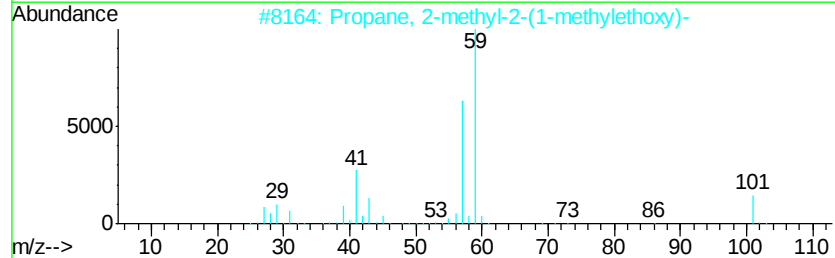
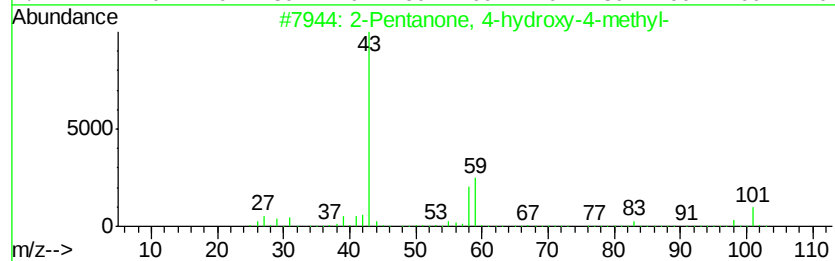
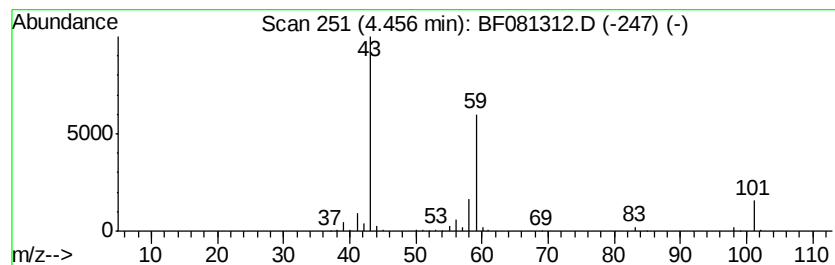
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	61.04 ng	1102060	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	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-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23	



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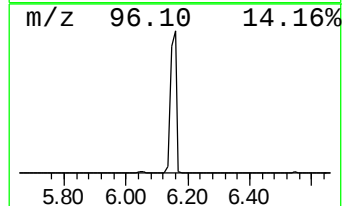
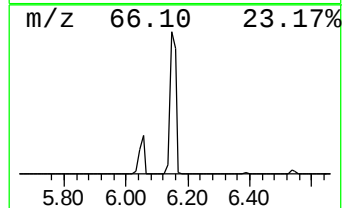
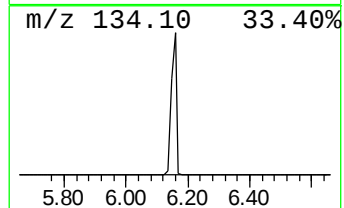
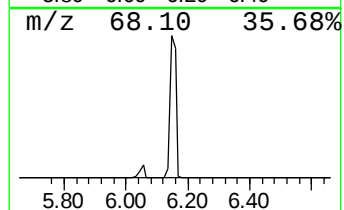
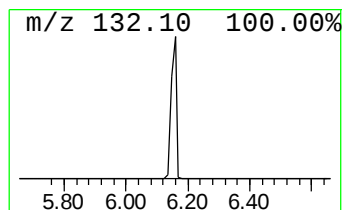
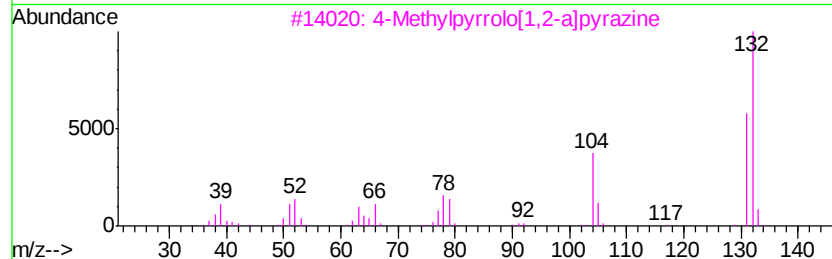
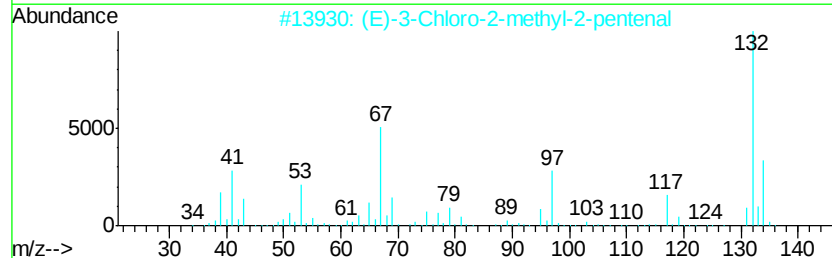
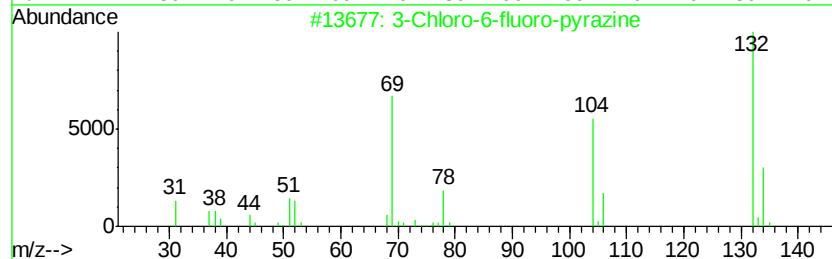
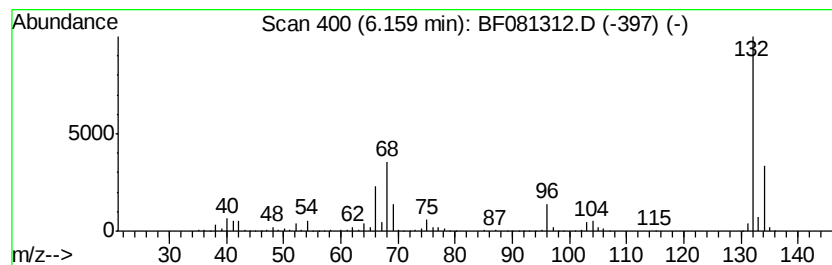
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Peak Number 2 unknown6.16 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.16	93.42 ng	1686700	1,4-Dichlorobenzene-d4	6.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.46	61.0	ng	1102060	1	6.39	361106	20.0
unknown6.16	6.16	93.4	ng	1686700	1	6.39	361106	20.0