

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100215\
 Data File : BF082003.D
 Acq On : 3 Oct 2015 4:34
 Operator : UM/IZ
 Sample : PB85910BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85910BL

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-BF092815.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	5.074	253	257	270	rBV	772159	1048706	36.20%	4.954%
2	5.486	290	293	295	rBV	1477598	1867874	64.47%	8.823%
3	6.514	380	383	385	rBV	1663090	1932785	66.71%	9.130%
4	6.651	392	395	397	rBV	2086971	2015288	69.56%	9.520%
5	6.891	413	416	427	rBV	327323	534510	18.45%	2.525%
6	7.040	427	429	432	rBV	1758399	1574244	54.34%	7.436%
7	7.452	463	465	488	rBV	1121739	1265089	43.67%	5.976%
8	8.172	526	528	537	rBV	735383	746453	25.77%	3.526%
9	9.257	620	623	625	rBV	1730379	2417330	83.44%	11.419%
10	9.932	679	682	685	rBV	935291	874944	30.20%	4.133%
11	10.720	749	751	754	rBV	1241148	1405894	48.53%	6.641%
12	11.418	810	812	826	rBV	569640	916407	31.63%	4.329%
13	13.018	949	952	973	rBV	2770987	2897095	100.00%	13.685%
14	14.069	1042	1044	1061	rBV	807488	960607	33.16%	4.538%
15	15.532	1169	1172	1186	rBV	442535	712747	24.60%	3.367%

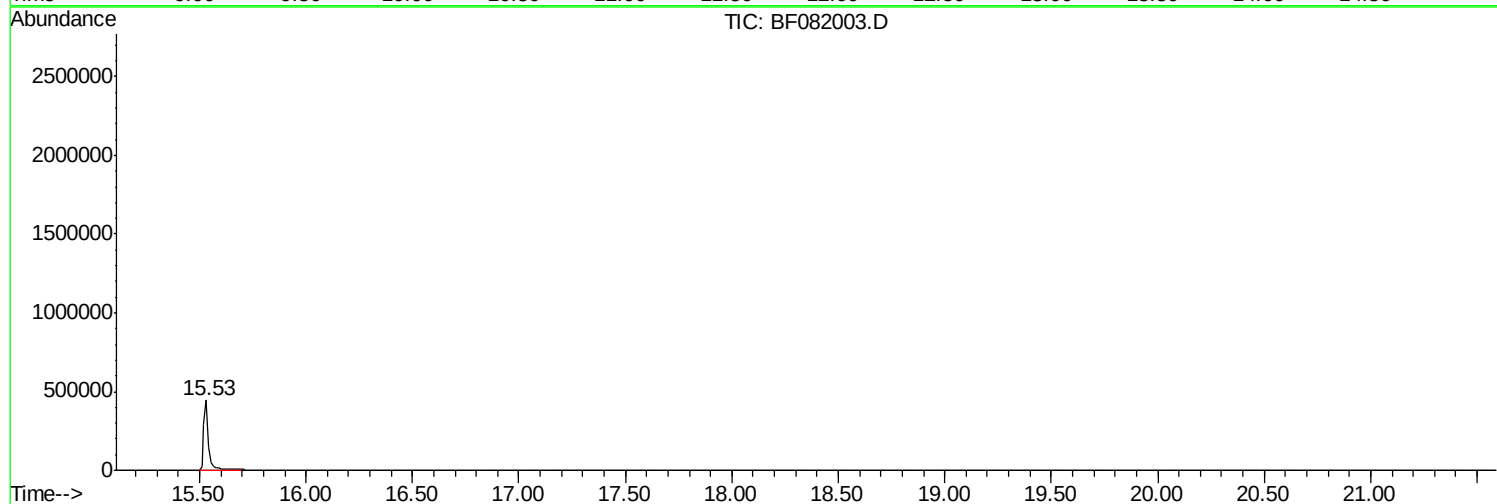
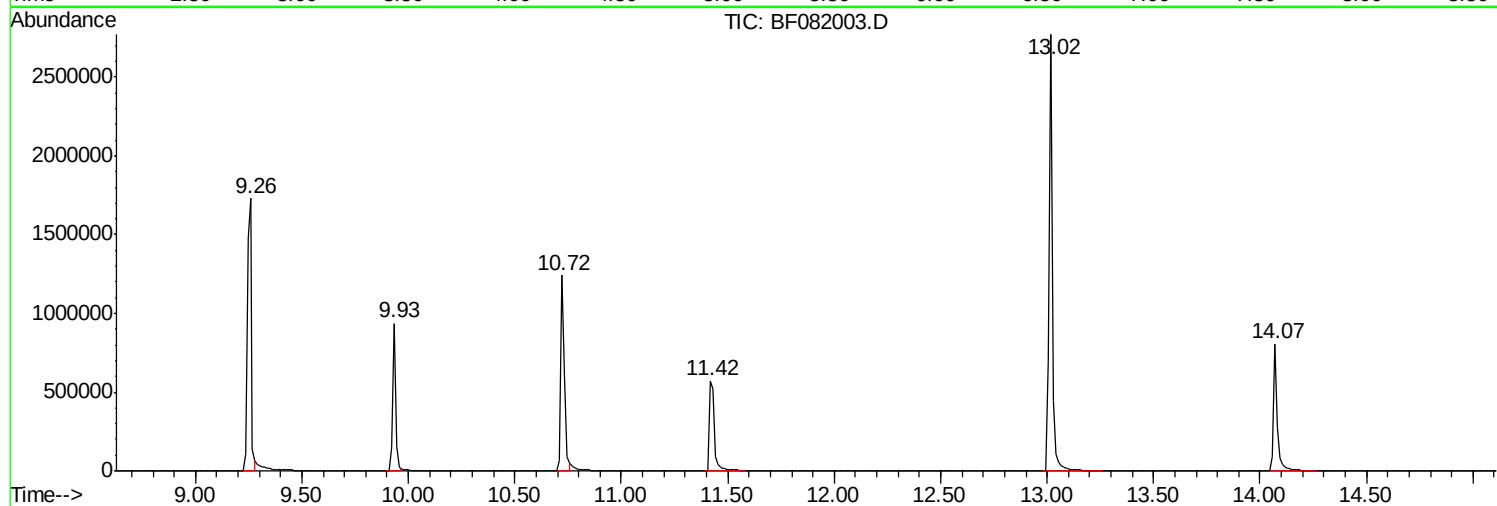
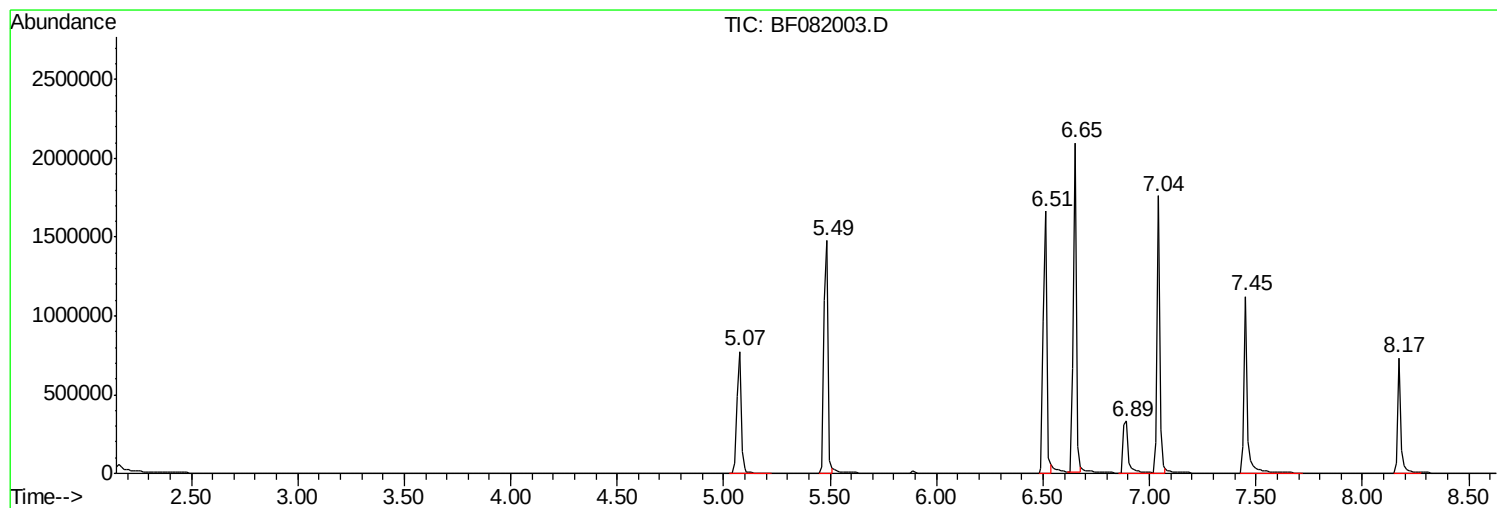
Sum of corrected areas: 21169973

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Quant Method : H:\HPCHEM1\BNA_F\Method\8270-BF092815.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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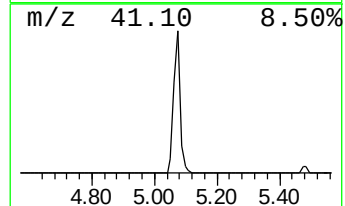
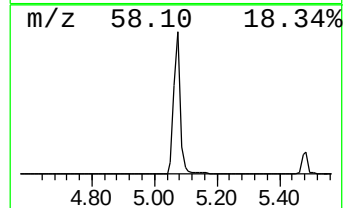
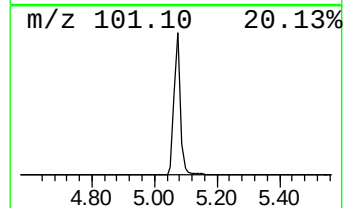
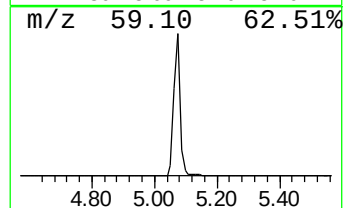
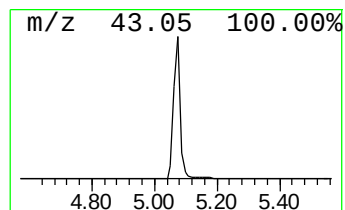
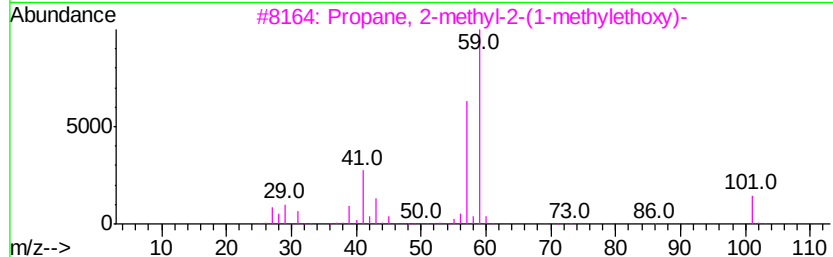
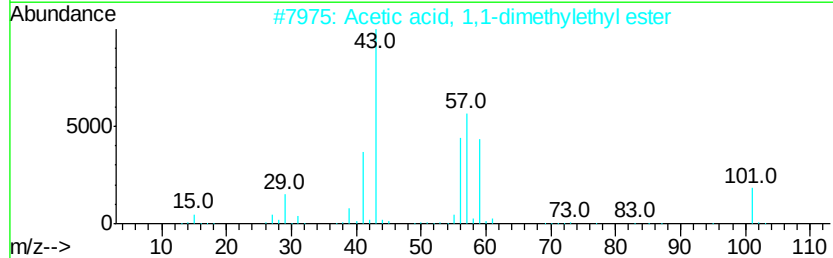
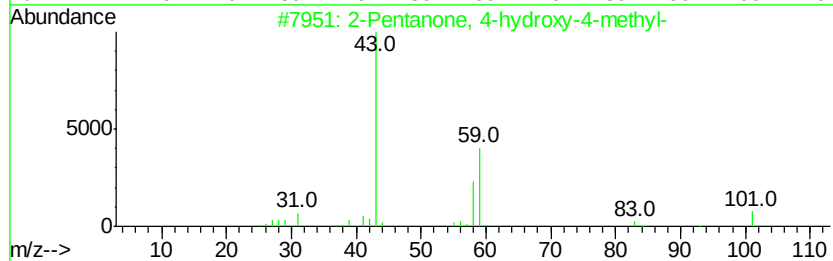
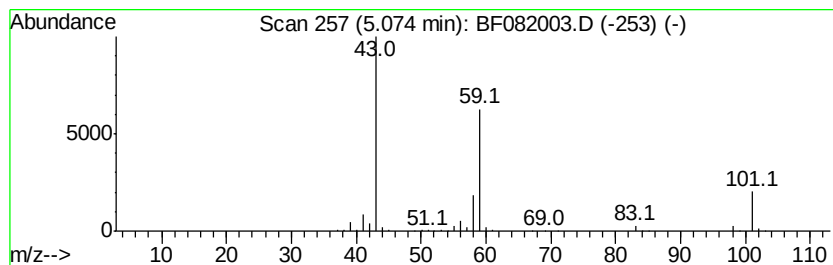
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	39.24 ng	1048710	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		Acetone	58	C3H6O	000067-64-1	9



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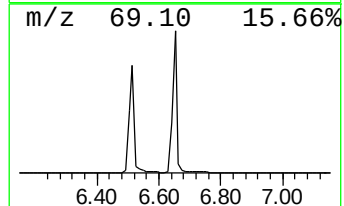
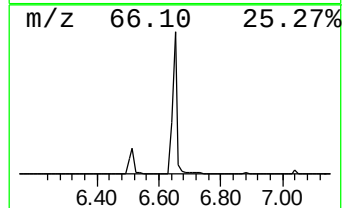
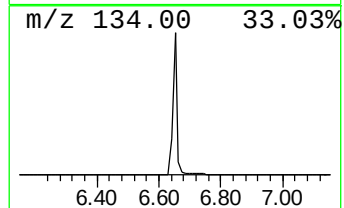
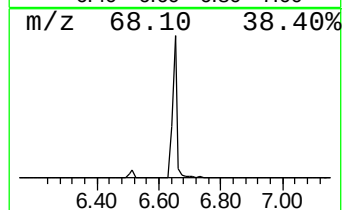
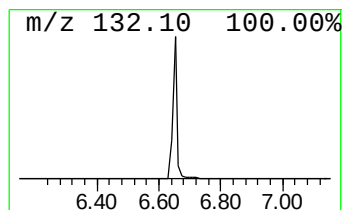
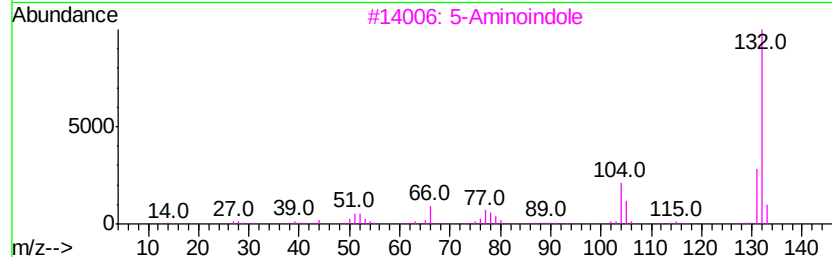
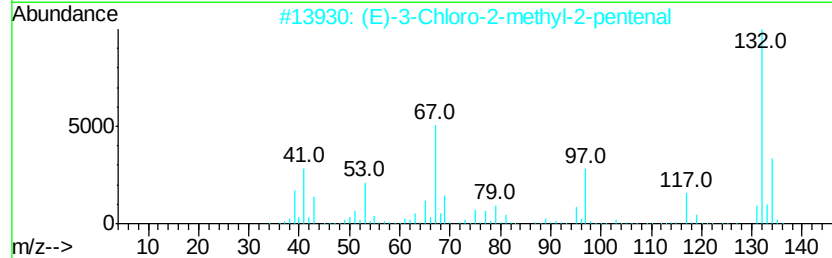
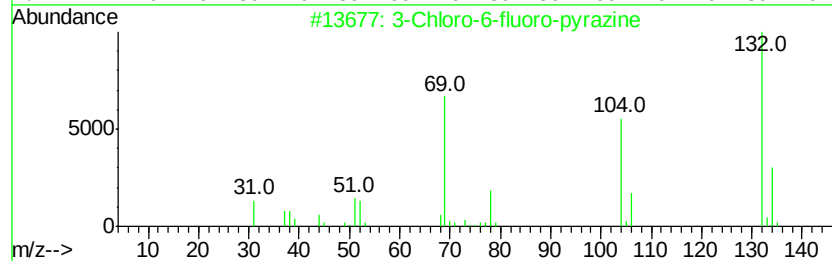
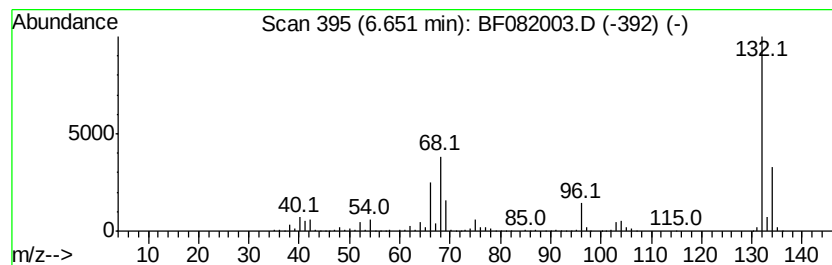
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Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	75.41 ng	2015290	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
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...	5.07	39.2	ng	1048710	1	6.89	534510	20.0
unknown6.65	6.65	75.4	ng	2015290	1	6.89	534510	20.0