

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100615\
 Data File : BF082083.D
 Acq On : 6 Oct 2015 14:36
 Operator : UM/IZ
 Sample : PB85944BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85944BL

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	266	rBV	685721	907875	31.82%	4.559%
2	5.486	290	293	295	rBV	1833553	1838558	64.43%	9.232%
3	6.514	380	383	385	rBV	1730311	1811851	63.50%	9.098%
4	6.652	392	395	397	rBV	1744173	1967622	68.95%	9.880%
5	6.880	413	415	426	rVB	521111	469157	16.44%	2.356%
6	7.040	426	429	431	rBV	1488566	1577579	55.29%	7.921%
7	7.452	462	465	467	rBV	961037	1126813	39.49%	5.658%
8	8.172	525	528	530	rBV	508291	584118	20.47%	2.933%
9	9.246	619	622	624	rBV	2389955	2318910	81.26%	11.644%
10	9.920	679	681	684	rBV	637752	721267	25.28%	3.622%
11	10.720	748	751	753	rBV	1160282	1445081	50.64%	7.256%
12	11.418	809	812	814	rBV	564540	737892	25.86%	3.705%
13	13.006	948	951	953	rBV	2473626	2853520	100.00%	14.328%
14	14.058	1040	1043	1045	rBV	679180	814094	28.53%	4.088%
15	15.510	1167	1170	1179	rBV	476121	741113	25.97%	3.721%

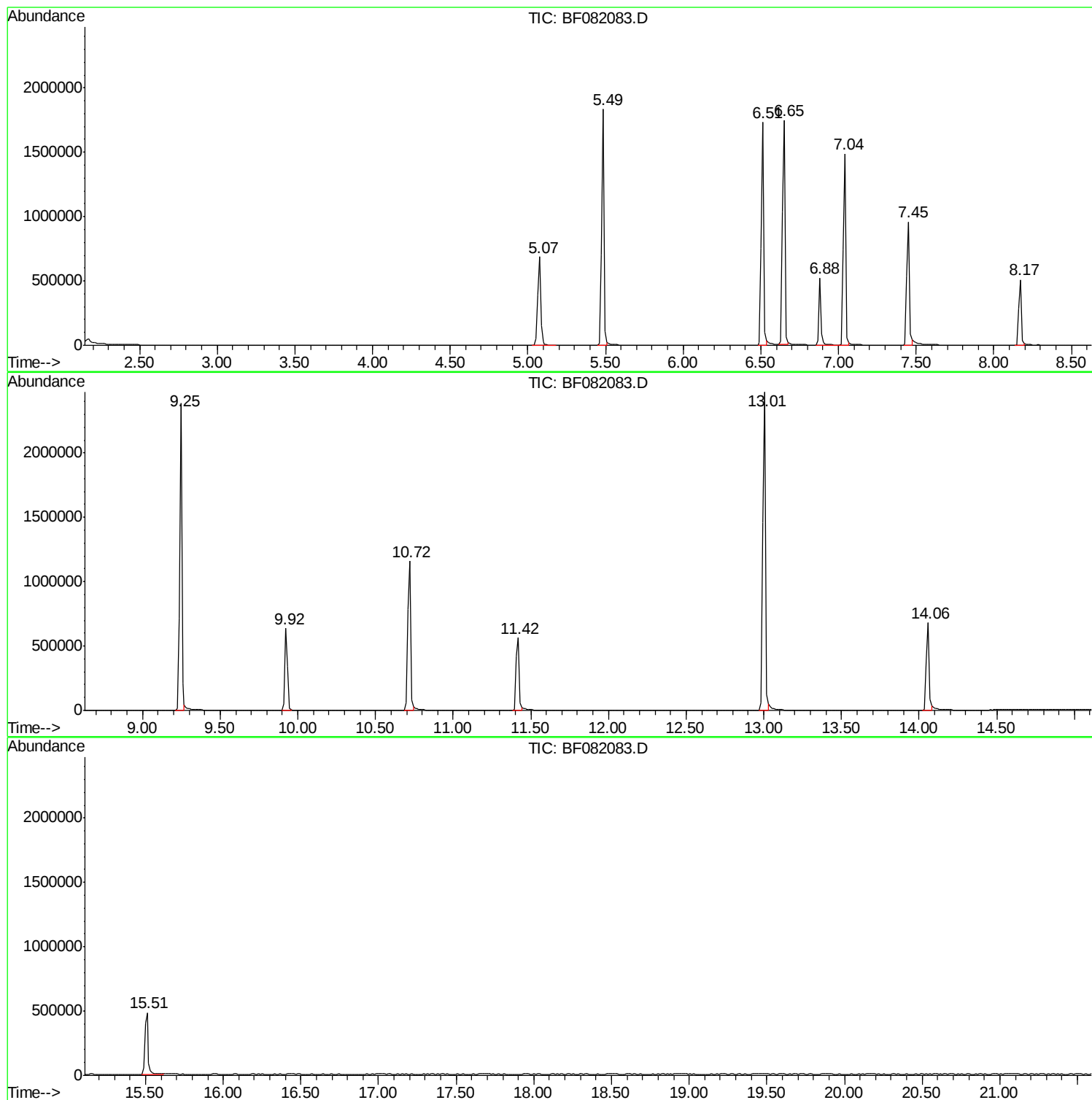
Sum of corrected areas: 19915450

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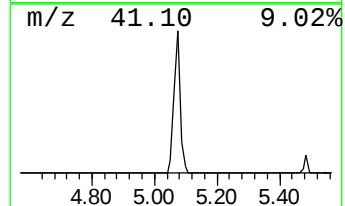
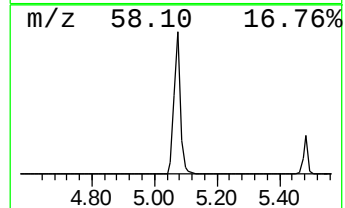
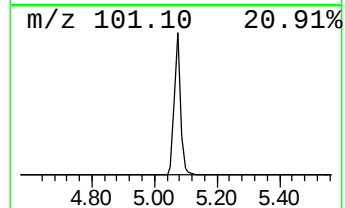
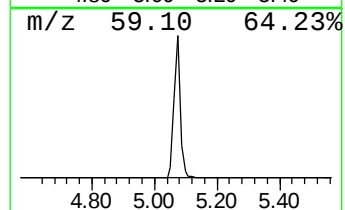
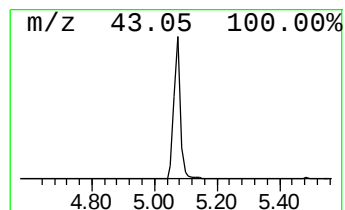
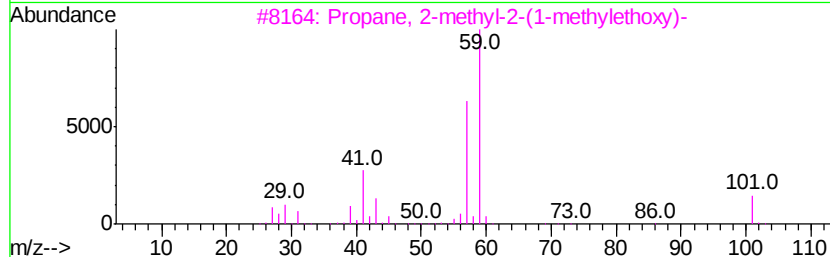
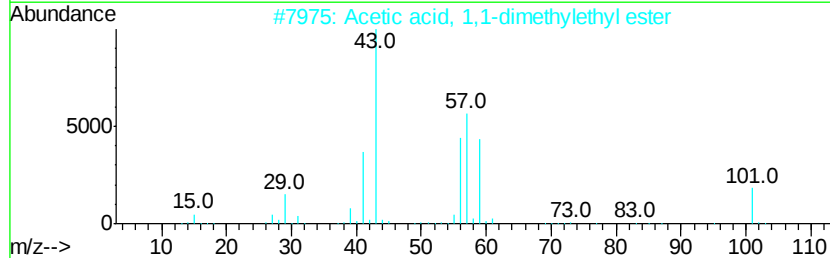
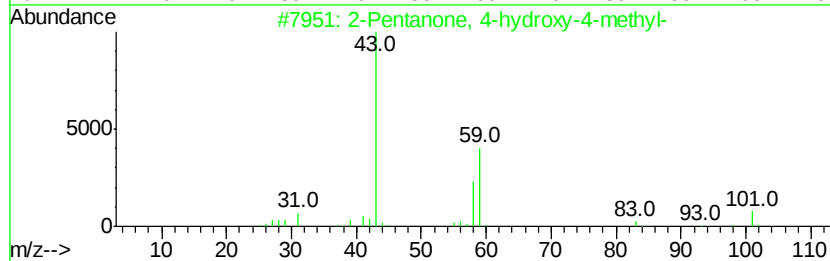
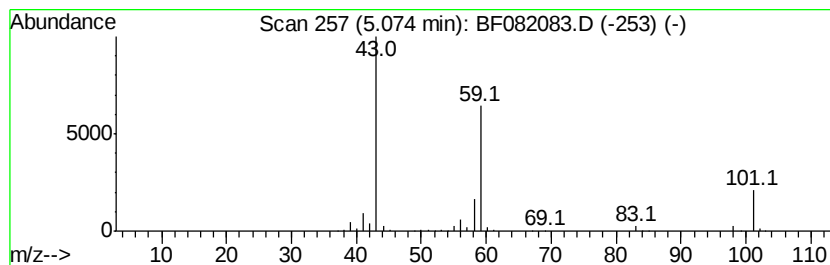
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	38.70 ng	907875	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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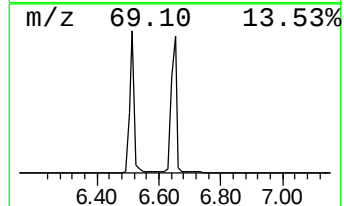
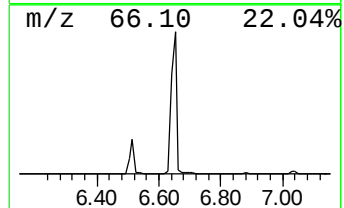
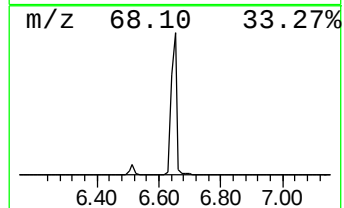
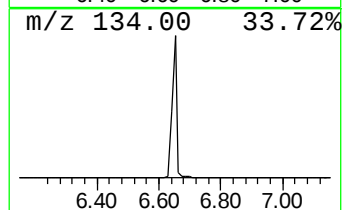
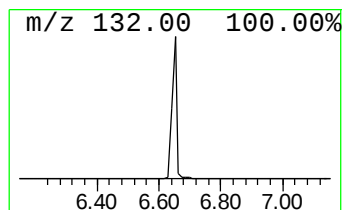
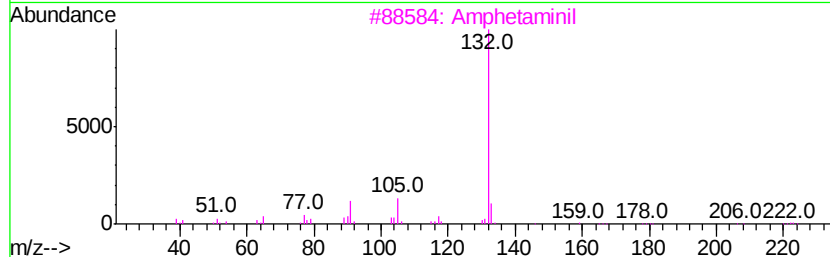
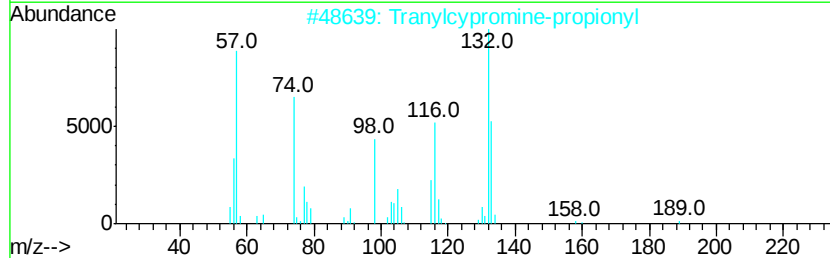
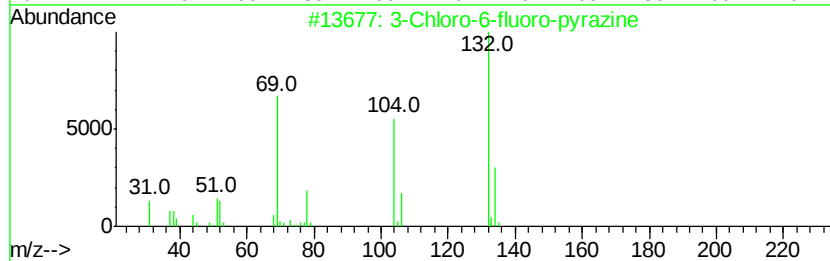
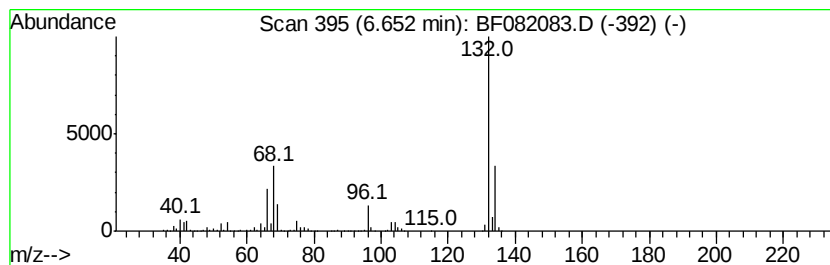
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Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	83.88 ng	1967620	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	38.7	ng	907875	1	6.88	469157	20.0
unknown6.65	6.65	83.9	ng	1967620	1	6.88	469157	20.0