

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100515\
 Data File : BF082045.D
 Acq On : 5 Oct 2015 14:52
 Operator : UM/IZ
 Sample : PB85905BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85905BL

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	5.086	255	258	266	rBV	237145	337368	10.19%	1.614%
2	5.486	290	293	295	rBV	1643324	1571447	47.46%	7.517%
3	6.503	380	382	385	rBV	1191766	1455674	43.96%	6.963%
4	6.651	392	395	397	rBV	1342276	1721758	52.00%	8.236%
5	6.880	413	415	424	rBV	570539	582251	17.58%	2.785%
6	7.040	426	429	431	rBV	2083804	2087003	63.03%	9.983%
7	7.452	462	465	467	rBV	1329864	1486373	44.89%	7.110%
8	8.172	525	528	530	rBV	635947	714671	21.58%	3.419%
9	9.246	619	622	625	rBV	2819069	2971964	89.75%	14.216%
10	9.920	678	681	684	rBV	727212	865299	26.13%	4.139%
11	10.709	748	750	753	rBV2	922144	1224378	36.98%	5.857%
12	11.418	809	812	814	rBV	602333	838798	25.33%	4.012%
13	13.006	948	951	953	rBV	2815554	3311304	100.00%	15.840%
14	14.058	1040	1043	1054	rBV	579179	920951	27.81%	4.405%
15	15.498	1166	1169	1180	rBV	538768	816008	24.64%	3.903%

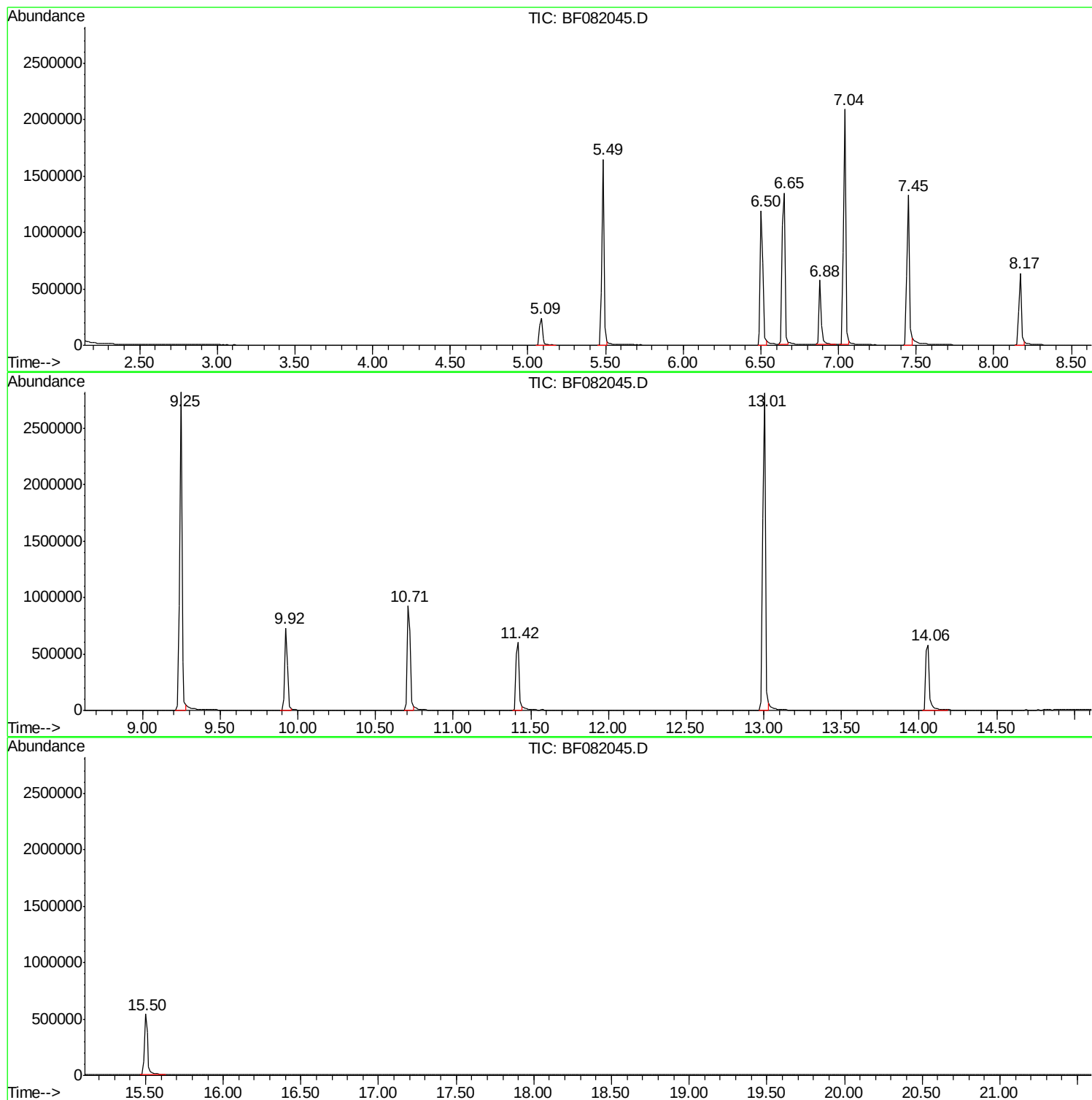
Sum of corrected areas: 20905247

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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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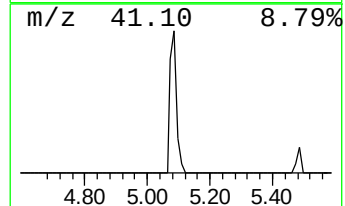
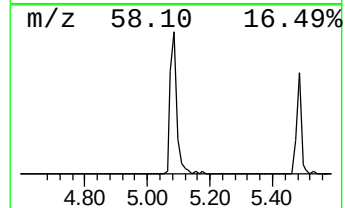
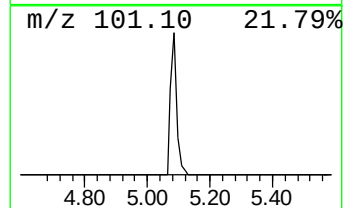
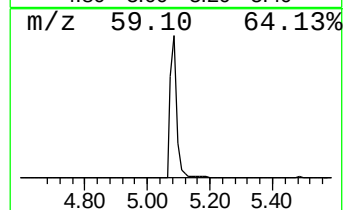
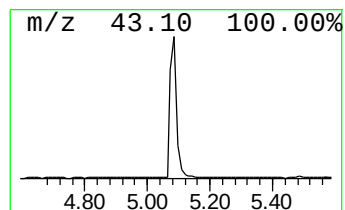
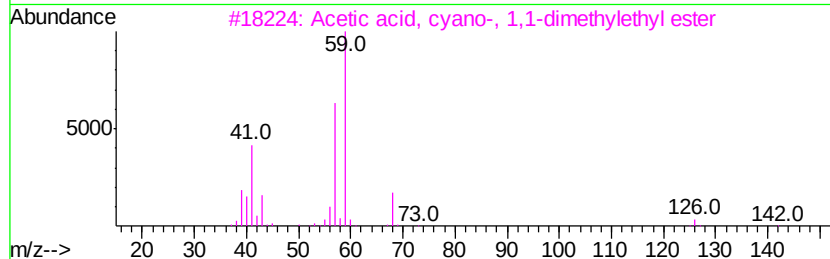
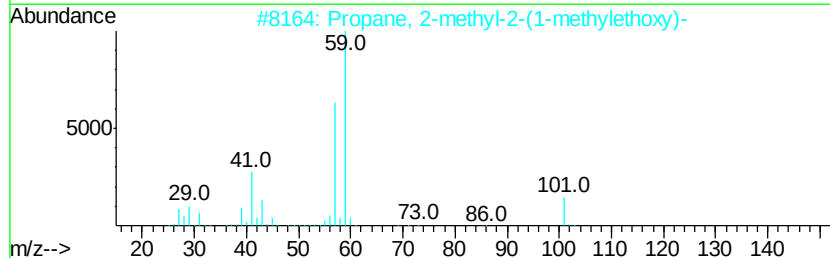
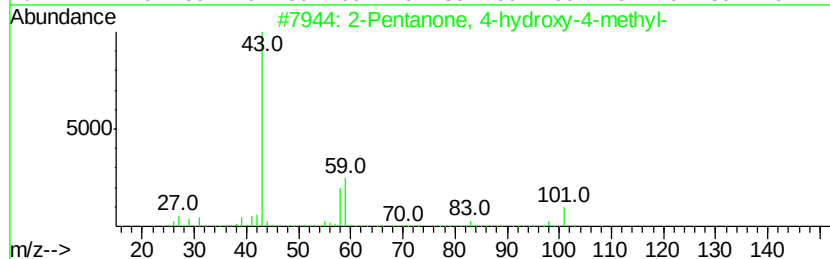
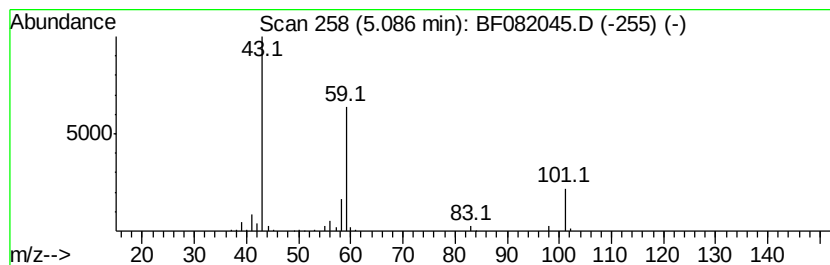
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	11.59 ng	337368	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	53
3		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Guanidine	59	CH5N3	000113-00-8	9



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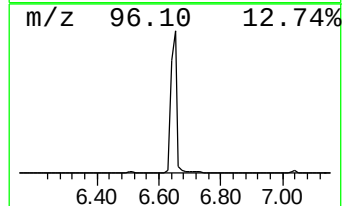
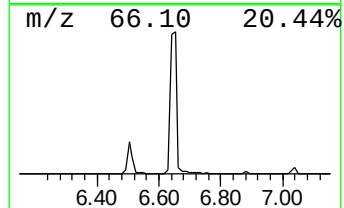
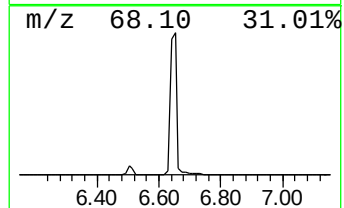
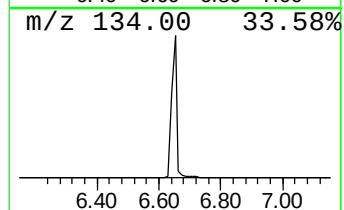
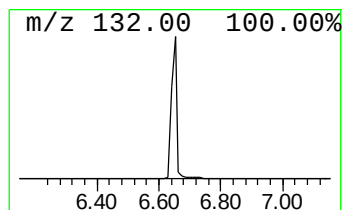
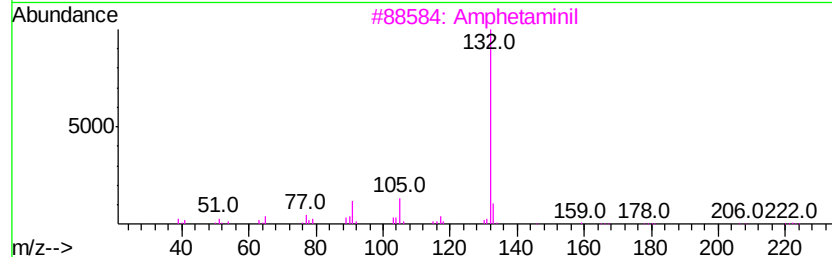
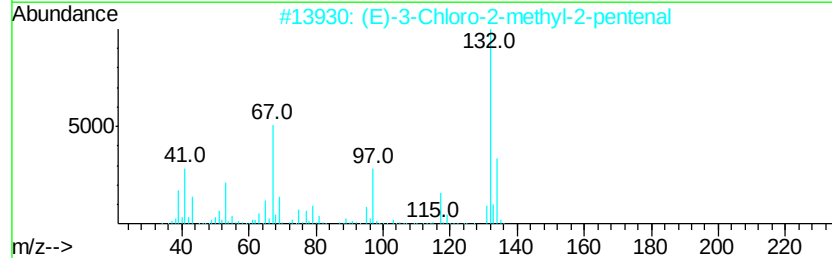
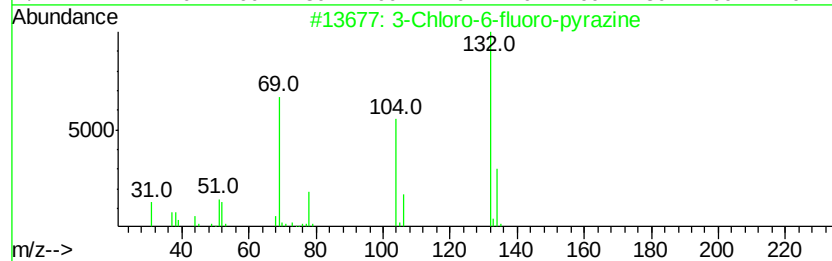
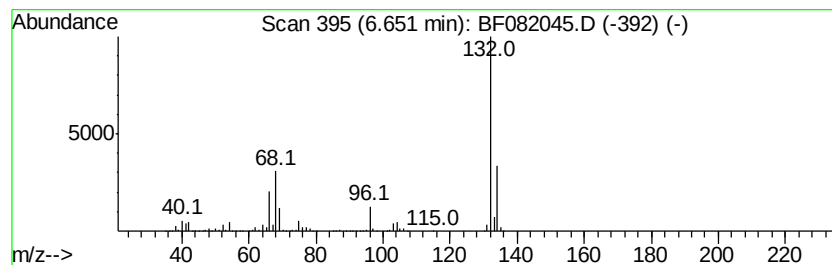
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Peak Number 2 unknown6.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	59.14 ng	1721760	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	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.09	11.6	ng	337368	1	6.88	582251	20.0
unknown6.65	6.65	59.1	ng	1721760	1	6.88	582251	20.0