

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102315\
 Data File : BF082479.D
 Acq On : 23 Oct 2015 21:50
 Operator : UM/IZ
 Sample : PB86215BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86215BL

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-BF101015.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	4.926	241	244	258	rBV	659418	1099924	38.33%	5.245%
2	5.360	279	282	284	rBV	1578027	1814636	63.23%	8.654%
3	6.400	370	373	381	rBV	1731497	1739287	60.61%	8.294%
4	6.526	381	384	386	rBV	1568819	1940066	67.61%	9.252%
5	6.754	402	404	406	rBV	464098	543728	18.95%	2.593%
6	6.903	415	417	420	rBV	1609573	1513371	52.74%	7.217%
7	7.326	451	454	456	rBV	1071123	1192107	41.54%	5.685%
8	8.035	514	516	519	rBV	578996	737263	25.69%	3.516%
9	9.120	608	611	614	rBV	2352916	2409747	83.97%	11.492%
10	9.795	667	670	673	rBV	931952	934179	32.55%	4.455%
11	10.595	737	740	742	rBV	1058205	1385720	48.29%	6.608%
12	11.281	798	800	803	rBV	889914	970115	33.81%	4.626%
13	12.881	937	940	942	rBV	2859236	2869703	100.00%	13.685%
14	13.967	1033	1035	1038	rBV	813860	990760	34.52%	4.725%
15	15.498	1166	1169	1179	rBV	495034	828806	28.88%	3.952%

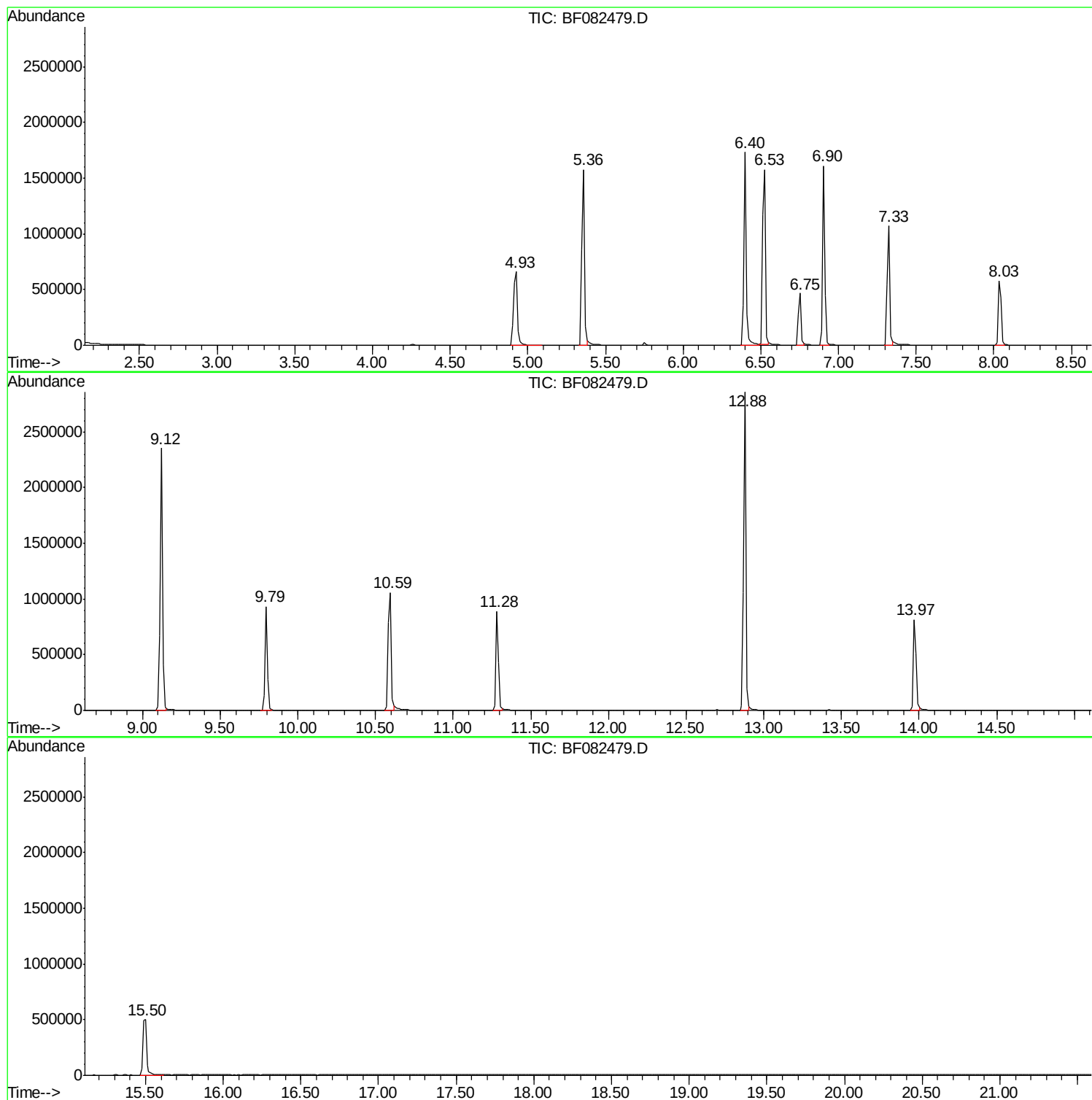
Sum of corrected areas: 20969412

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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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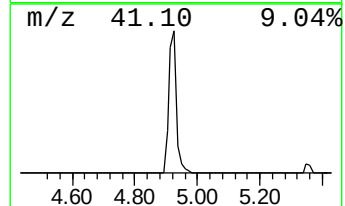
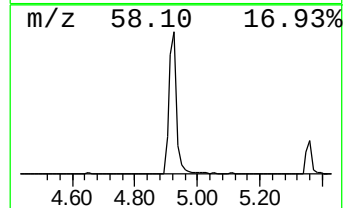
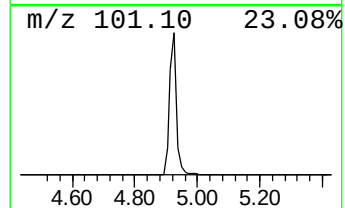
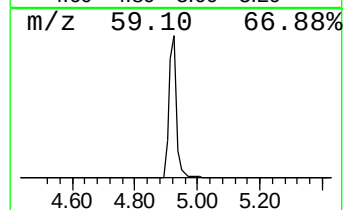
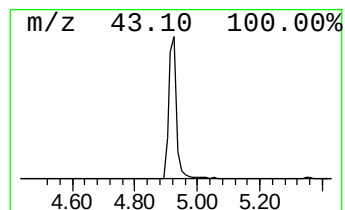
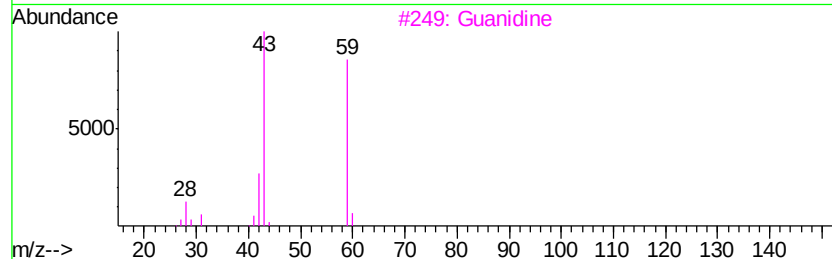
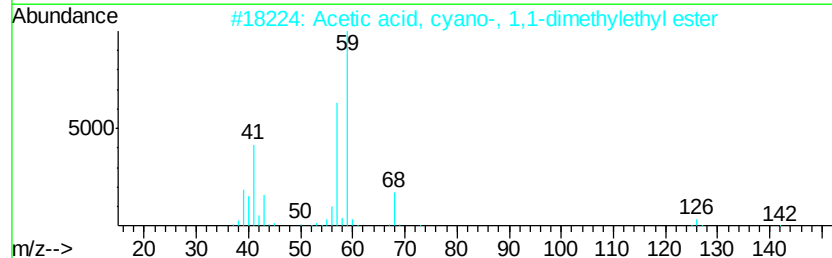
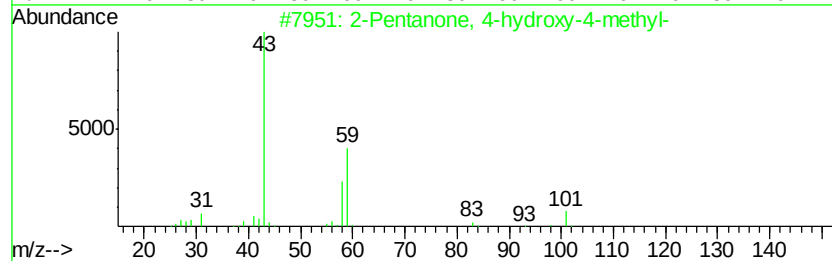
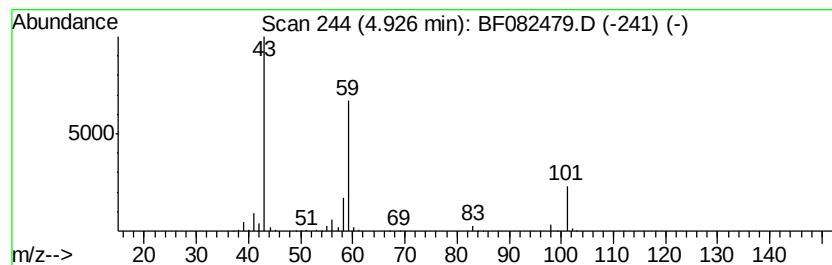
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	40.46 ng	1099920	1,4-Dichlorobenzene-d4	6.75

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, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		Guanidine	59	CH5N3	000113-00-8	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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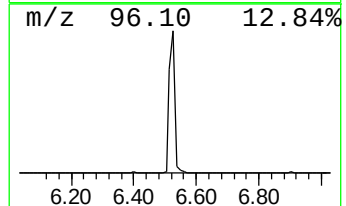
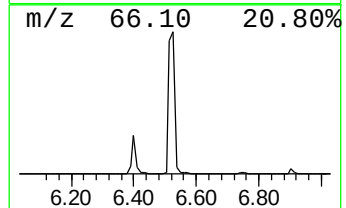
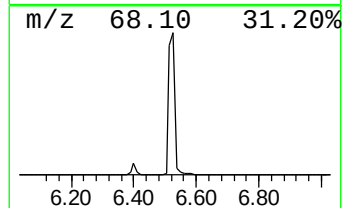
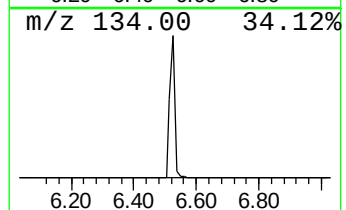
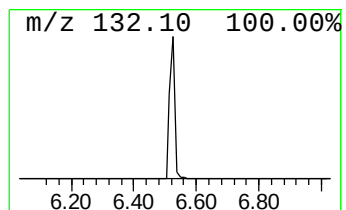
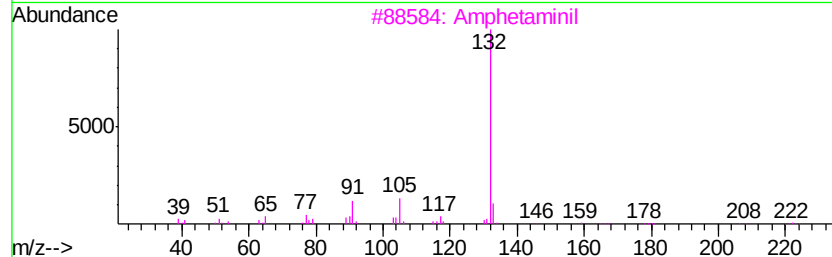
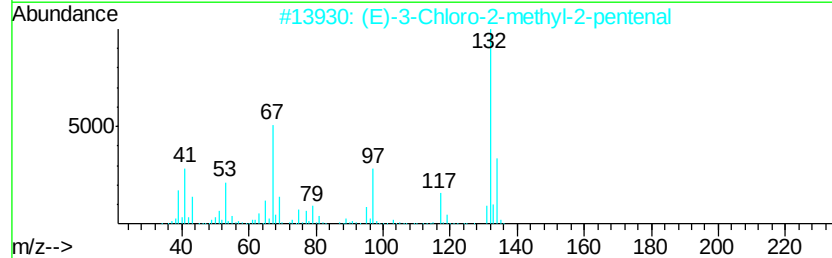
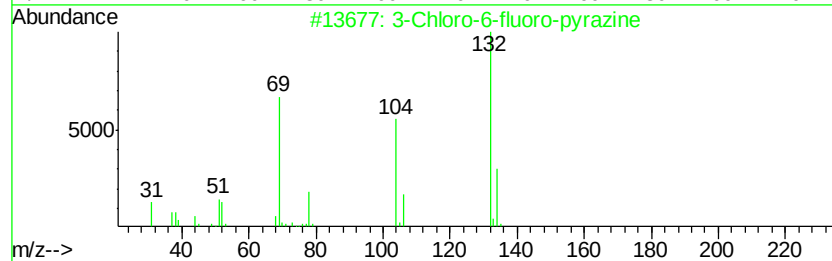
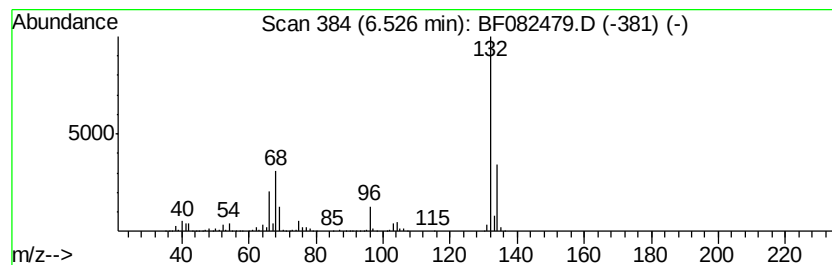
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Peak Number 2 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	71.36 ng	1940070	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.93	40.5	ng	1099920	1	6.75	543728	20.0
unknown6.53	6.53	71.4	ng	1940070	1	6.75	543728	20.0