

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122215\
 Data File : BF084016.D
 Acq On : 23 Dec 2015 4:16
 Operator : UM/SJ
 Sample : G4916-08
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20151214MGP1209BRV46.5

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-BF122215.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	4.487	208	210	213	rBB	62070	79454	4.33%	0.751%
2	5.001	253	255	262	rBB	76604	82092	4.48%	0.775%
3	5.436	291	293	296	rBV	373274	507337	27.68%	4.792%
4	6.476	382	384	386	rBV	228754	292238	15.94%	2.760%
5	6.601	393	395	397	rBV	1225527	1086615	59.28%	10.264%
6	6.830	413	415	417	rBV	288960	250653	13.67%	2.368%
7	6.990	426	429	431	rBV	831600	1040807	56.78%	9.831%
8	7.390	462	464	467	rBV	752004	735393	40.12%	6.946%
9	8.110	525	527	530	rBV	400122	346379	18.90%	3.272%
10	9.196	619	622	624	rBV	1675415	1833144	100.00%	17.316%
11	9.870	679	681	683	rBV	385662	387966	21.16%	3.665%
12	10.659	747	750	752	rBV	1052462	1067234	58.22%	10.081%
13	11.356	809	811	813	rBV	299332	398197	21.72%	3.761%
14	12.934	947	949	952	rBV	1492576	1796129	97.98%	16.966%
15	13.985	1039	1041	1044	rBV	372531	388020	21.17%	3.665%
16	15.414	1164	1166	1172	rVB	199761	295021	16.09%	2.787%

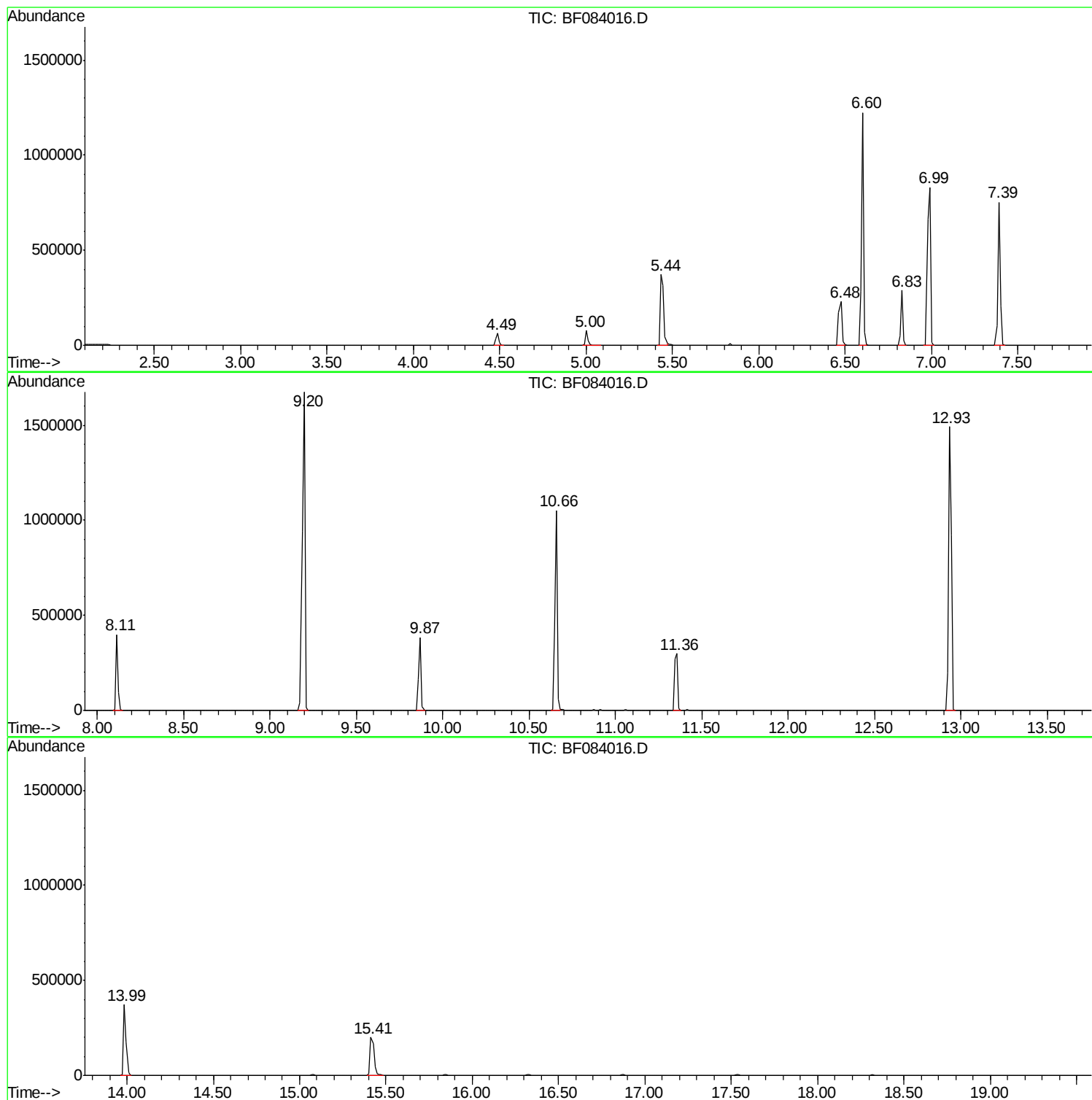
Sum of corrected areas: 10586679

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122215.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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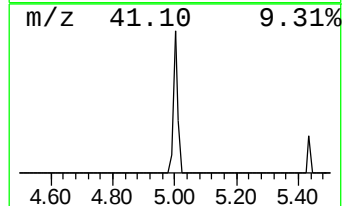
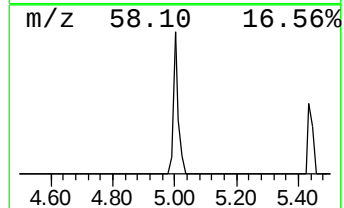
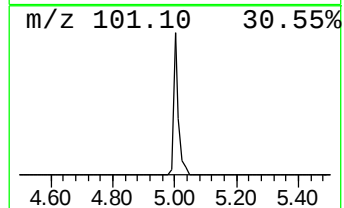
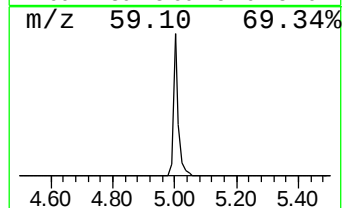
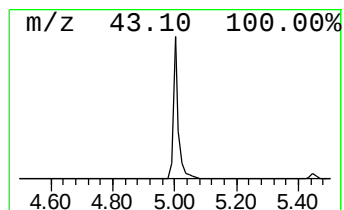
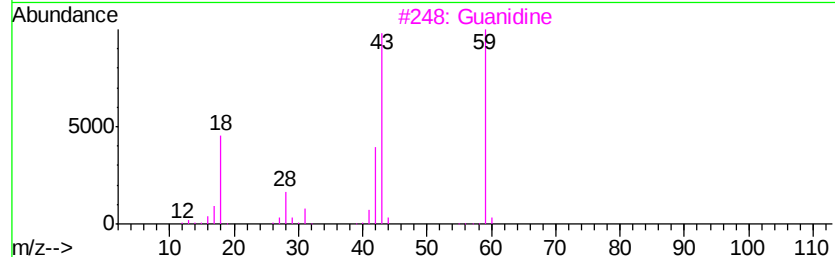
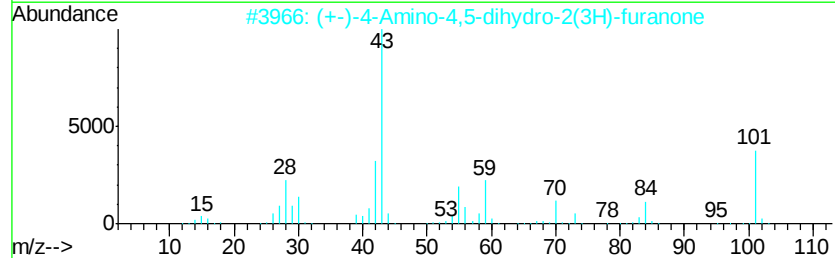
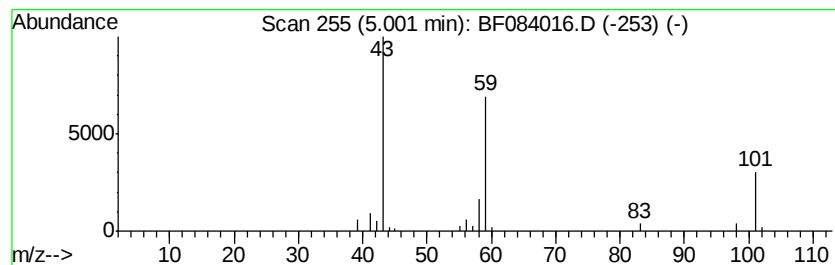
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	6.55 ng	82092	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	47
3		Guanidine	59	CH5N3	000113-00-8	43
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
5		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17



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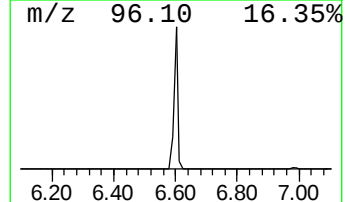
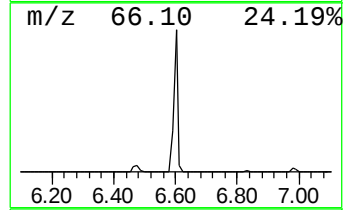
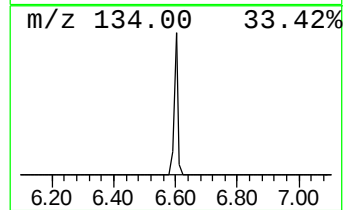
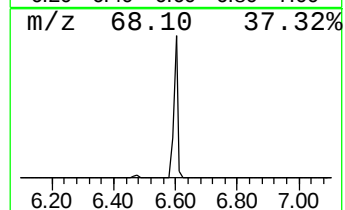
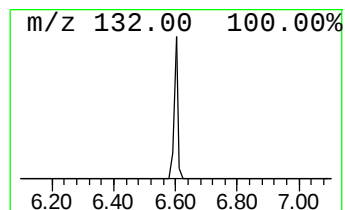
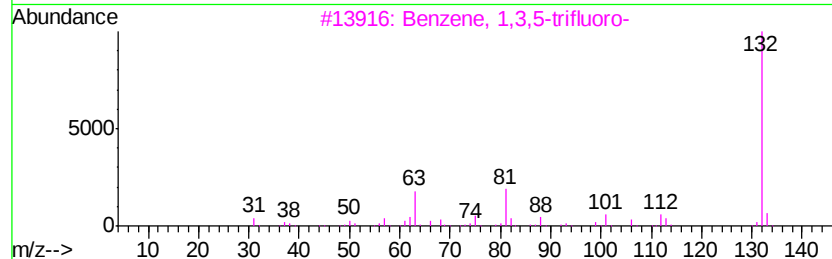
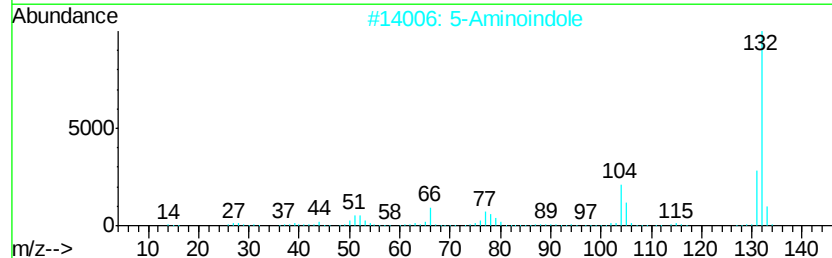
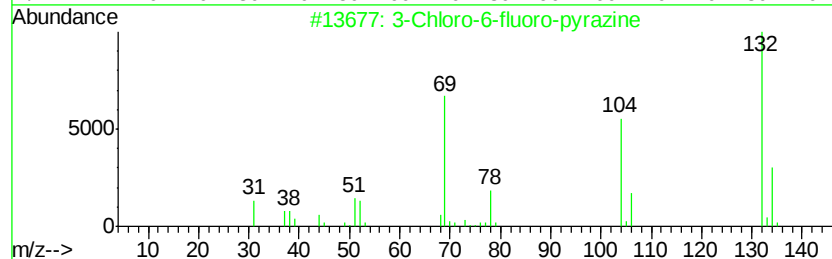
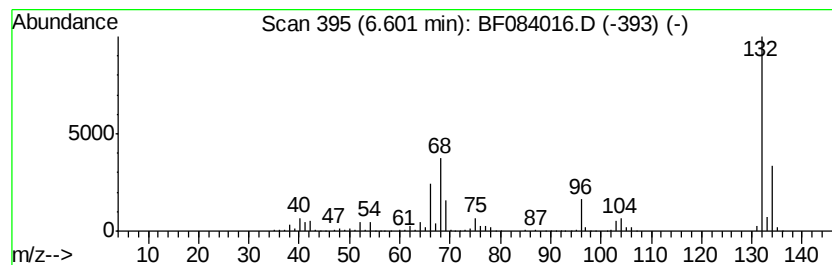
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Peak Number 3 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	86.70 ng	1086620	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	5	Aminoindole	132	C8H8N2	005192-03-0	12
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	6.5	ng	82092	1	6.83	250653	20.0
unknown6.60	6.60	86.7	ng	1086620	1	6.83	250653	20.0