

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011216\
 Data File : BF084432.D
 Acq On : 13 Jan 2016 1:56
 Operator : SJ/UM
 Sample : H1078-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TE-PMW-11S

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-BF123115.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	4.350	195	198	203	rVB2	54050	88955	1.10%	0.162%
2	5.013	253	256	264	rVB	970891	1088458	13.42%	1.985%
3	5.447	291	294	296	rBV	4154915	3783984	46.65%	6.902%
4	5.847	327	329	331	rBV	123629	114502	1.41%	0.209%
5	6.476	382	384	387	rBV	2326140	2212293	27.28%	4.035%
6	6.613	393	396	398	rBV	6664845	6255887	77.13%	11.410%
7	6.842	414	416	418	rBV	2008320	1677797	20.69%	3.060%
8	7.002	427	430	432	rBV	5488151	5854391	72.18%	10.678%
9	7.413	463	466	468	rBV	3699110	5135163	63.31%	9.366%
10	8.122	526	528	531	rBV	1818829	2052759	25.31%	3.744%
11	9.208	620	623	625	rBV	8644516	8110627	100.00%	14.793%
12	9.596	655	657	662	rVB3	66415	94974	1.17%	0.173%
13	9.813	674	676	679	rBV	88171	94830	1.17%	0.173%
14	9.882	679	682	684	rBV	2644517	2314984	28.54%	4.222%
15	10.316	717	720	722	rBV	144707	138583	1.71%	0.253%
16	10.671	748	751	754	rBV	4729126	4571516	56.36%	8.338%
17	10.785	760	761	766	rVB	81632	116042	1.43%	0.212%
18	11.368	809	812	814	rBV	2037465	2054331	25.33%	3.747%
19	12.957	948	951	953	rBV	6100745	6403176	78.95%	11.679%
20	13.997	1040	1042	1045	rBV	1387684	1362779	16.80%	2.486%
21	15.425	1164	1167	1170	rBV	1094602	1301526	16.05%	2.374%

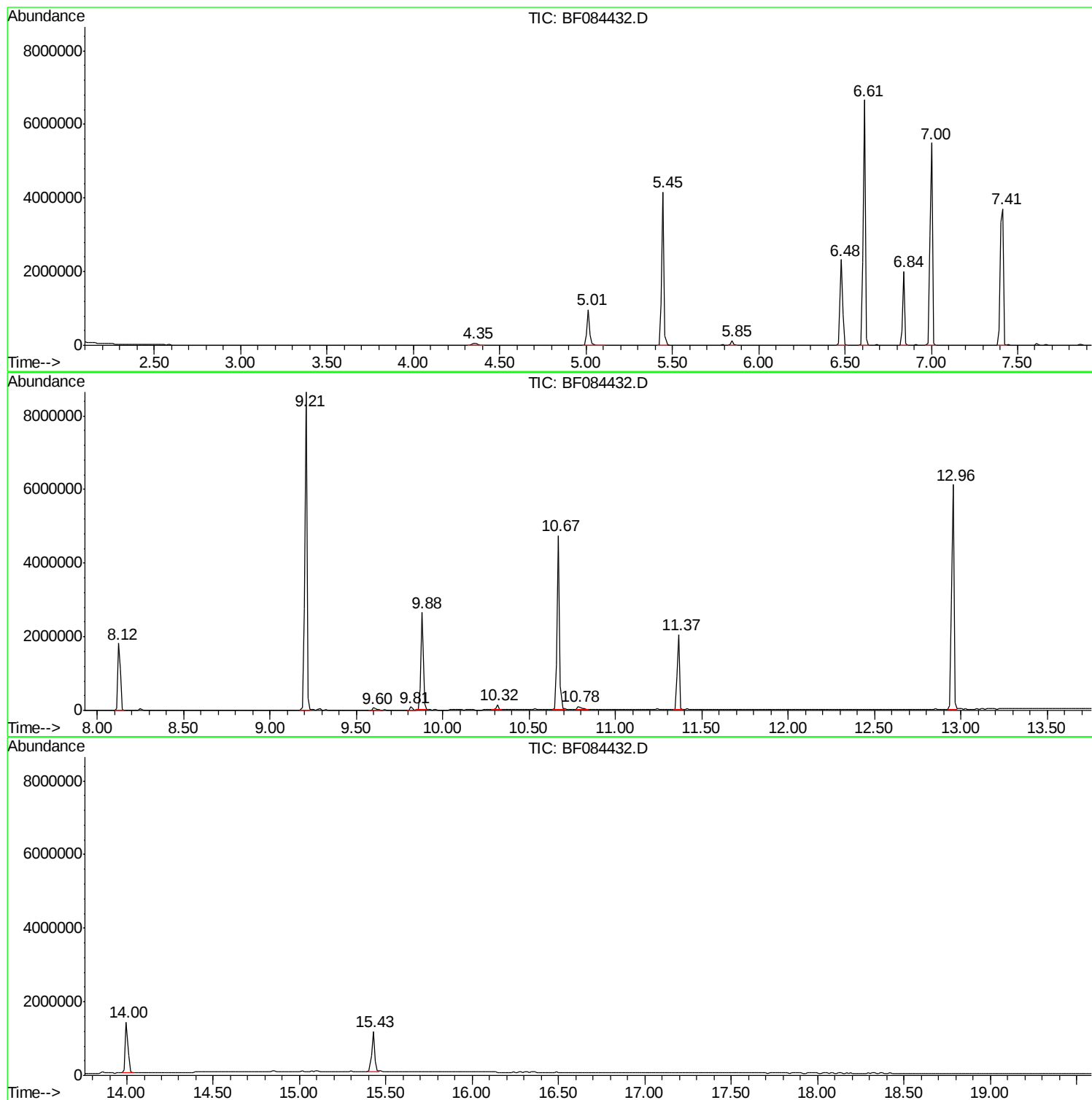
Sum of corrected areas: 54827557

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.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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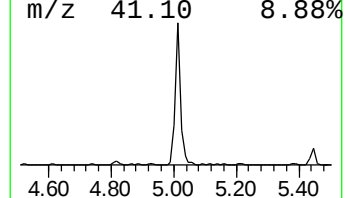
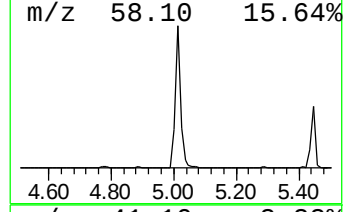
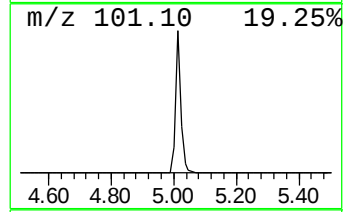
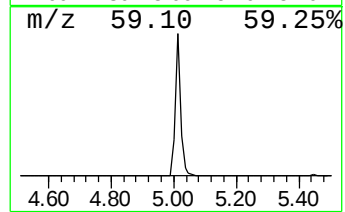
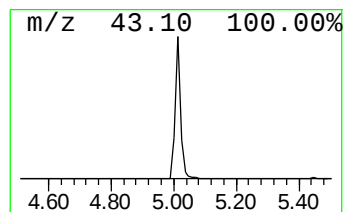
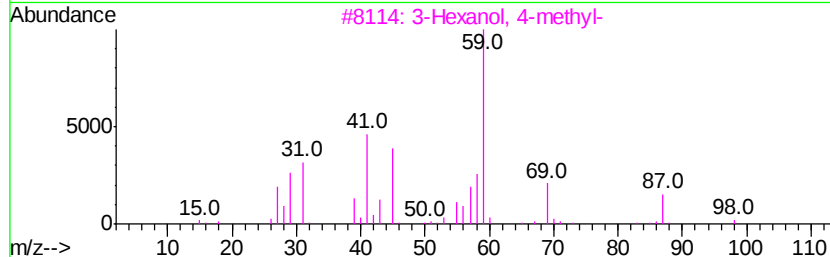
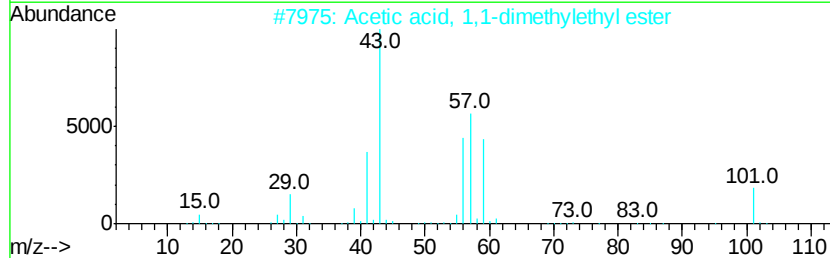
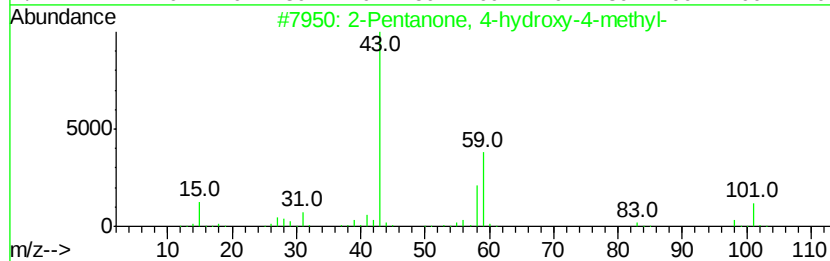
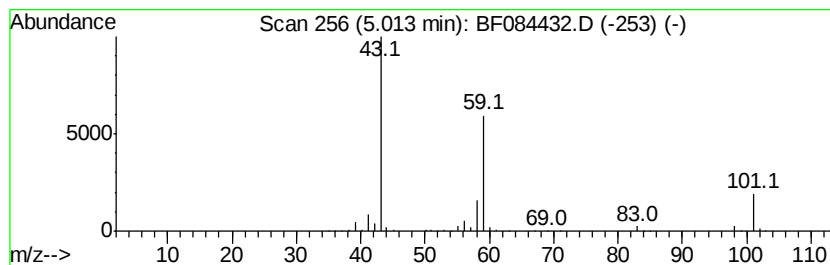
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.01	12.97 ng	1088460	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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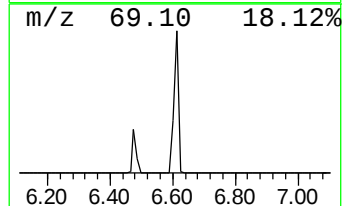
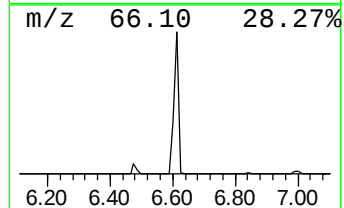
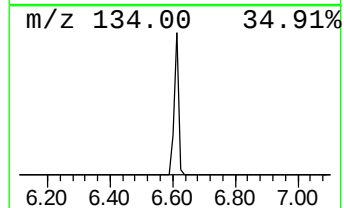
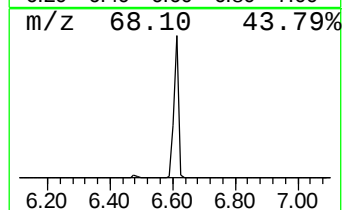
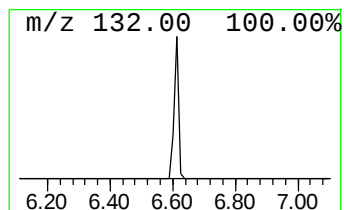
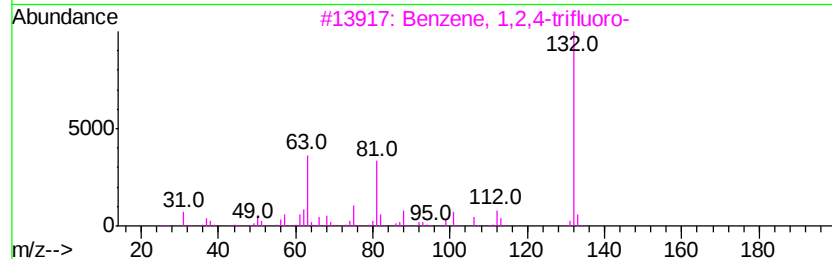
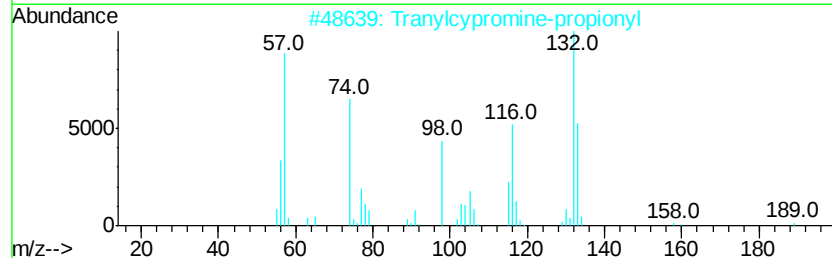
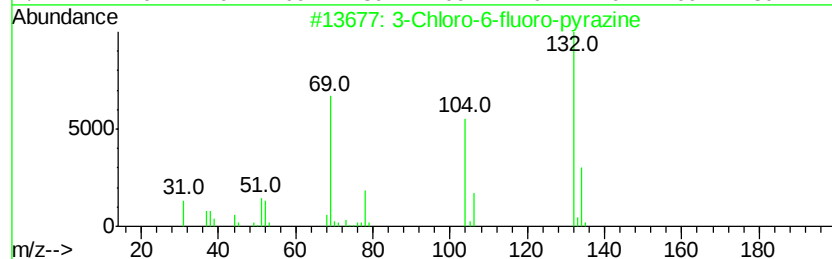
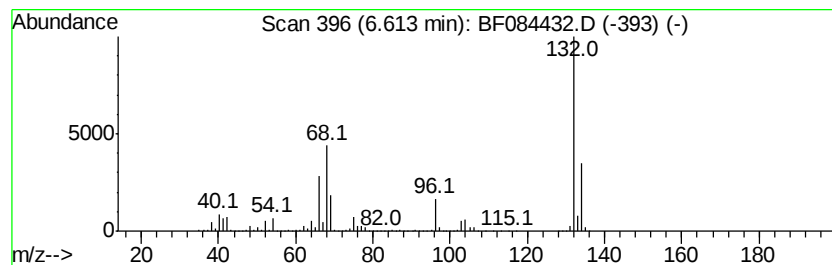
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Peak Number 2 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	74.57 ng	6255890	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	13.0	ng	1088460	1	6.84	1677800	20.0
unknown6.61	6.61	74.6	ng	6255890	1	6.84	1677800	20.0