

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
 Data File : BF083013.D
 Acq On : 18 Nov 2015 16:35
 Operator : UM/IZ
 Sample : PB86689BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86689BL

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-BF110715.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	4.899	242	246	255	rBV	1611080	2363158	44.26%	5.391%
2	5.344	282	285	287	rBV	3723041	4524484	84.75%	10.323%
3	6.384	373	376	378	rBV	4122870	4680134	87.66%	10.678%
4	6.499	383	386	389	rBV	4141397	3910084	73.24%	8.921%
5	6.727	404	406	408	rBV	1211805	981490	18.38%	2.239%
6	6.887	417	420	422	rBV	3633900	3411954	63.91%	7.784%
7	7.299	452	456	458	rBV	2116780	2996874	56.13%	6.837%
8	8.019	516	519	521	rBV	1088206	1255750	23.52%	2.865%
9	9.093	610	613	616	rBV	4211856	4713103	88.28%	10.753%
10	9.768	669	672	674	rBV	1653673	1541237	28.87%	3.516%
11	10.556	738	741	743	rBV	3234064	3446524	64.56%	7.863%
12	11.242	799	801	804	rVB	1702704	1508923	28.26%	3.443%
13	12.831	937	940	943	rBV	4118616	5338835	100.00%	12.180%
14	13.871	1029	1031	1034	rBV	1793339	1589813	29.78%	3.627%
15	15.265	1150	1153	1156	rBV	1296480	1568904	29.39%	3.579%

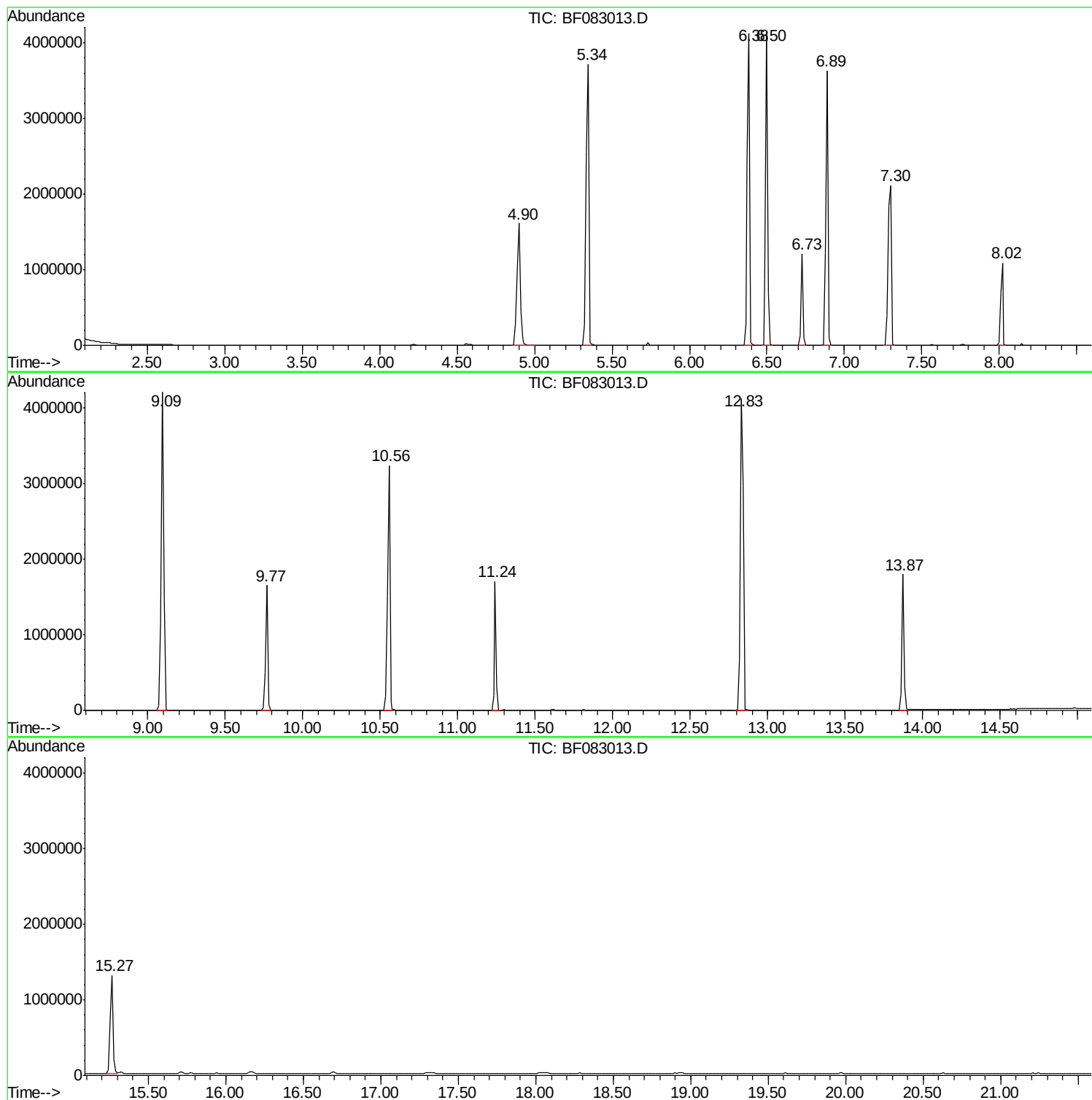
Sum of corrected areas: 43831267

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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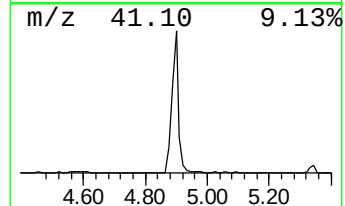
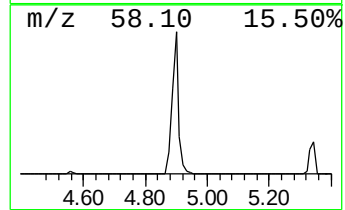
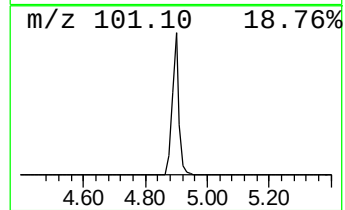
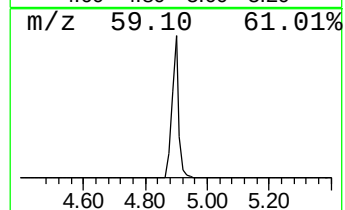
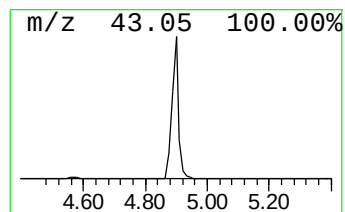
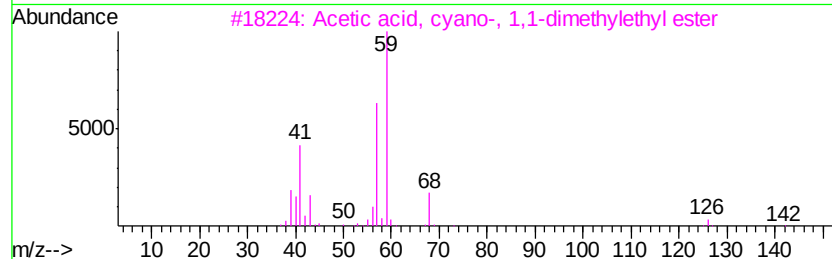
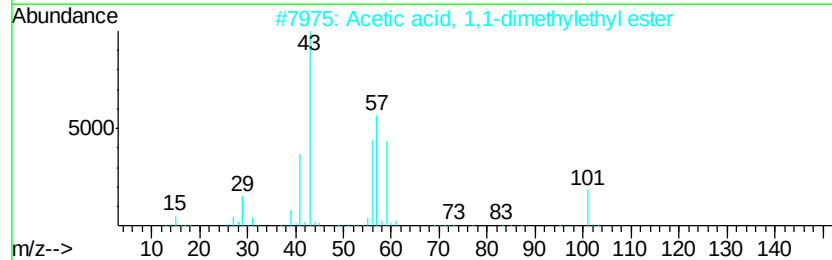
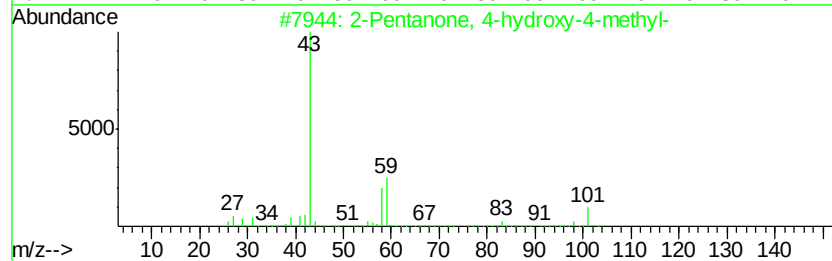
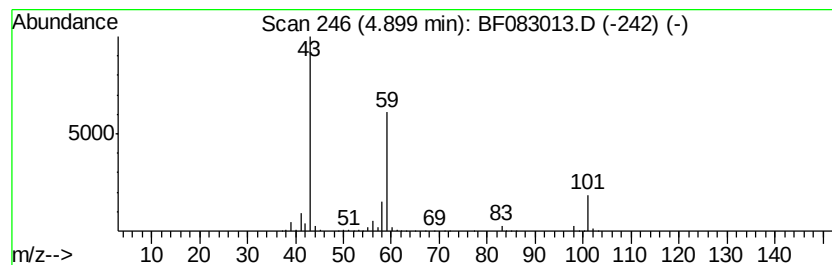
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	48.15 ng	2363160	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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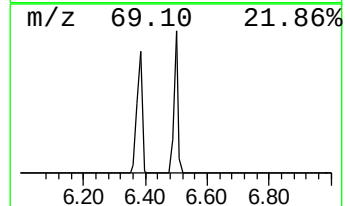
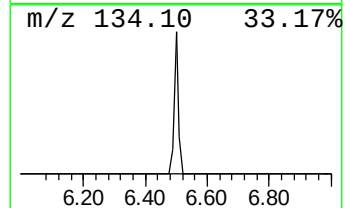
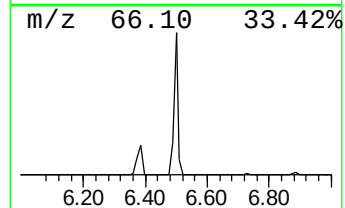
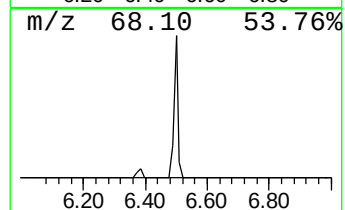
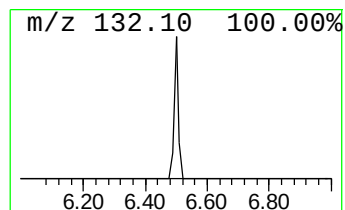
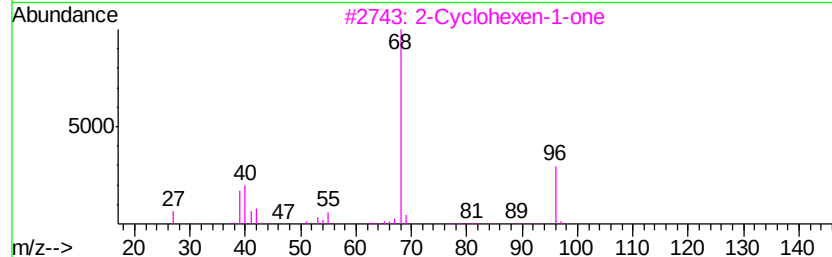
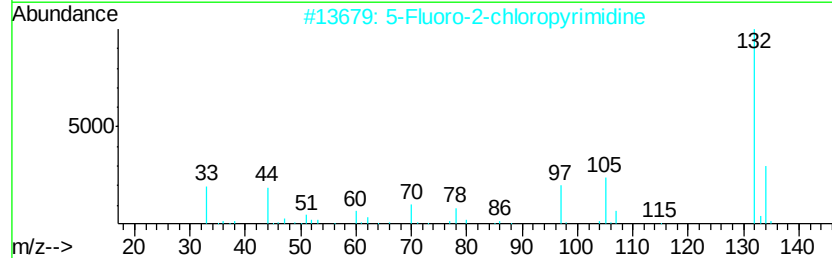
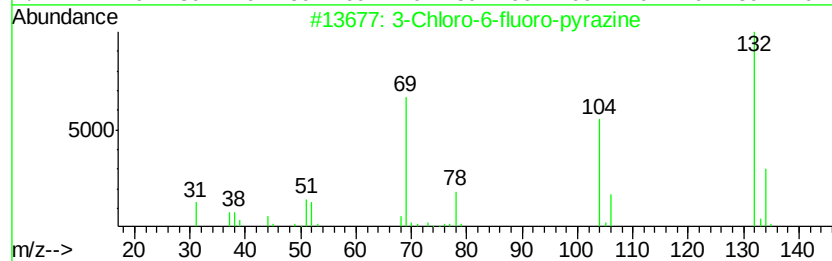
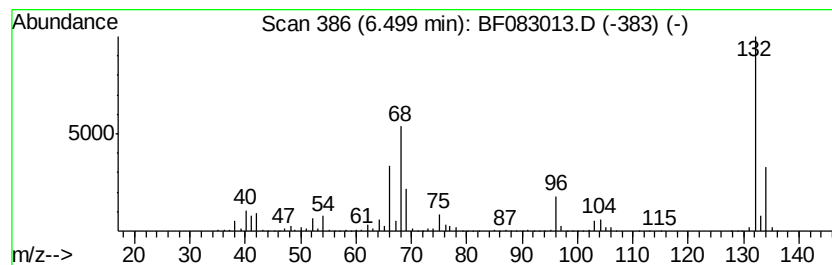
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Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	79.68 ng	3910080	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	48.1	ng	2363160	1	6.73	981490	20.0
unknown6.50	6.50	79.7	ng	3910080	1	6.73	981490	20.0