

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
 Data File : BF083012.D
 Acq On : 18 Nov 2015 16:06
 Operator : UM/IZ
 Sample : PB86622BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86622BL

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	4.887	242	245	255	rVB	709868	917733	14.90%	2.467%
2	5.333	281	284	287	rBV	2882597	2664222	43.26%	7.161%
3	6.373	373	375	378	rBV	2363359	2606112	42.32%	7.005%
4	6.499	383	386	388	rBV	3274670	2887940	46.89%	7.763%
5	6.727	404	406	408	rVB	1148852	933948	15.16%	2.510%
6	6.887	417	420	422	rBV	3686975	3317509	53.87%	8.917%
7	7.299	452	456	458	rBV	2031212	2770313	44.98%	7.446%
8	8.019	516	519	521	rBV	1046418	1178973	19.14%	3.169%
9	9.093	610	613	616	rBV	4516734	4401477	71.47%	11.831%
10	9.768	669	672	674	rBV	1530342	1426282	23.16%	3.834%
11	10.556	738	741	743	rBV	1988951	2653942	43.09%	7.134%
12	11.242	799	801	804	rVB	2280050	1946086	31.60%	5.231%
13	12.831	937	940	943	rBV	5259662	6158645	100.00%	16.554%
14	13.871	1028	1031	1034	rBV	2140718	1896948	30.80%	5.099%
15	15.265	1150	1153	1156	rBV	1249459	1442873	23.43%	3.878%

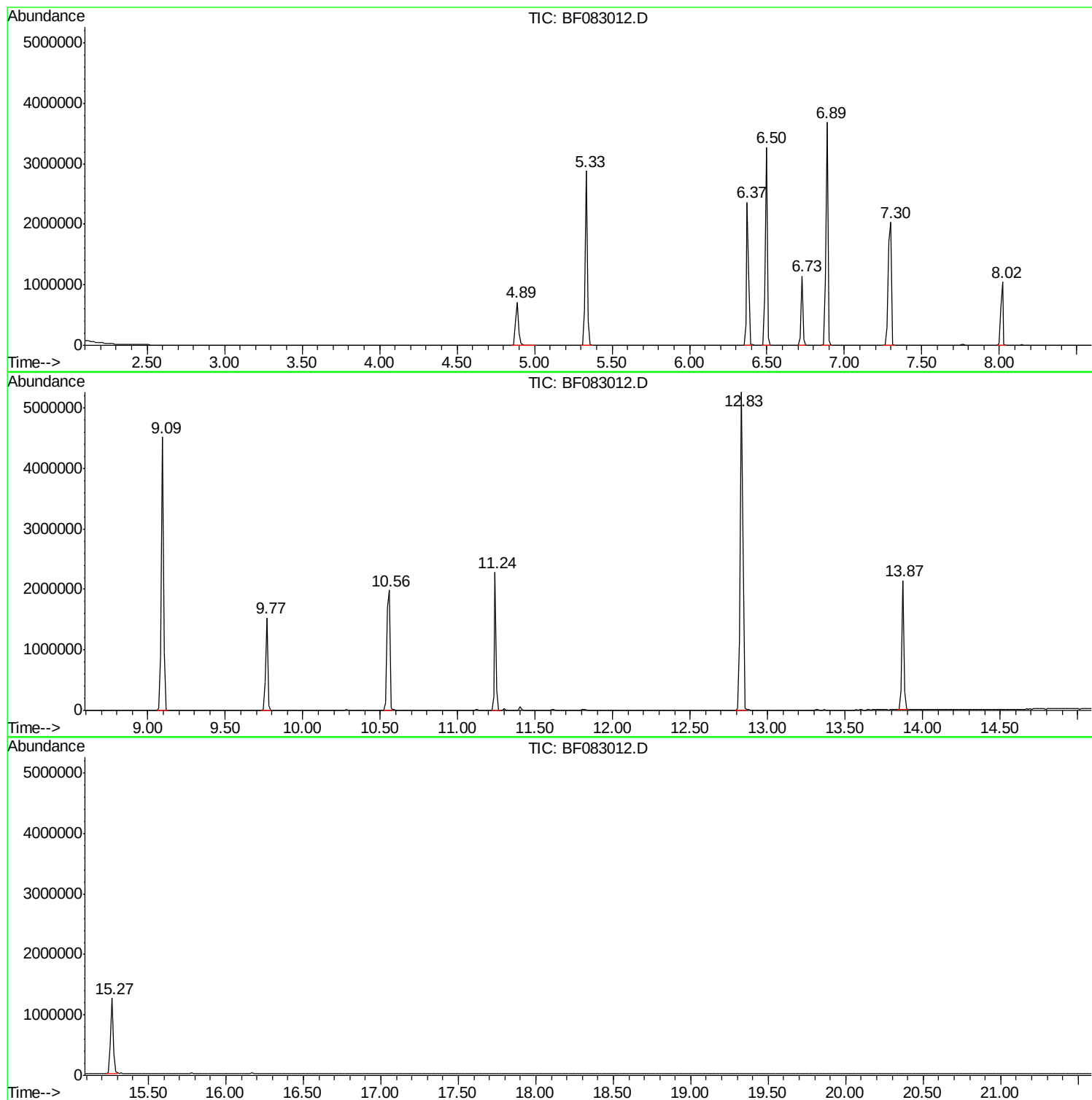
Sum of corrected areas: 37203003

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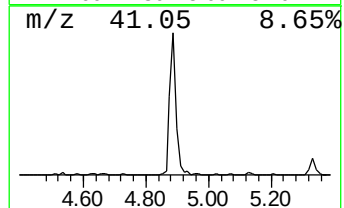
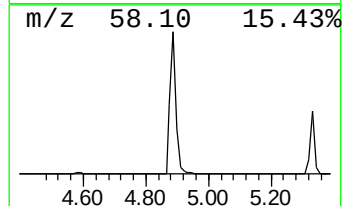
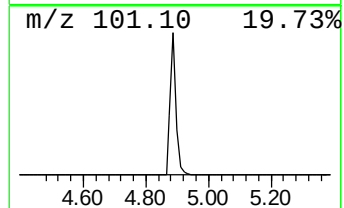
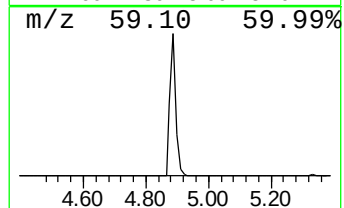
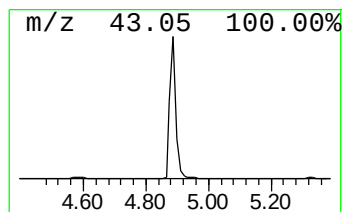
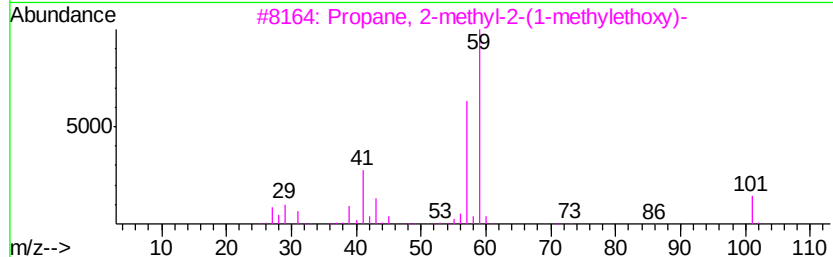
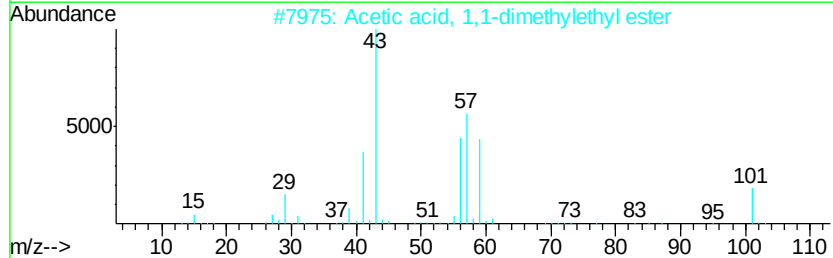
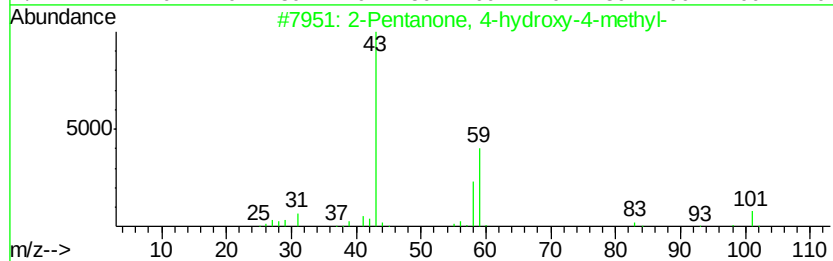
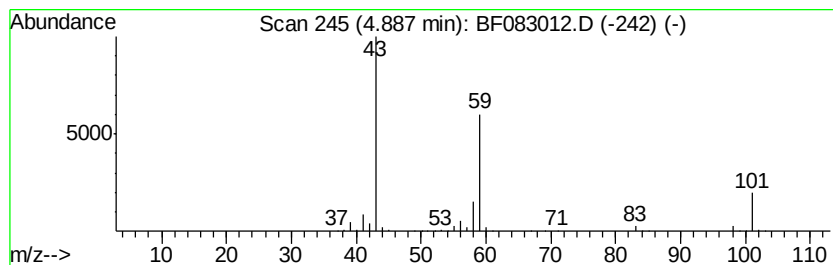
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	19.65 ng	917733	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	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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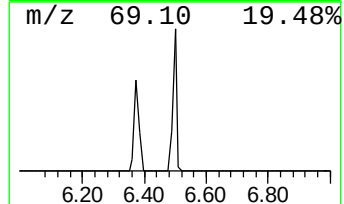
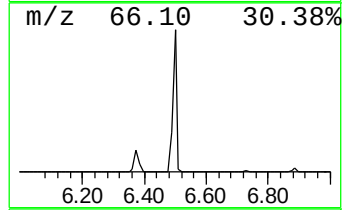
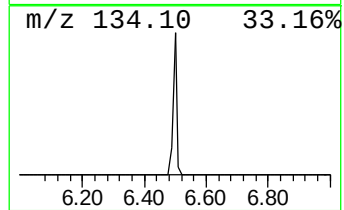
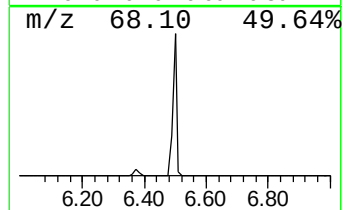
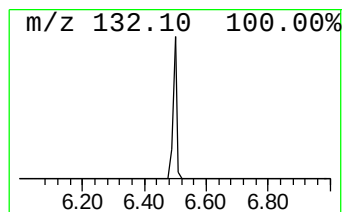
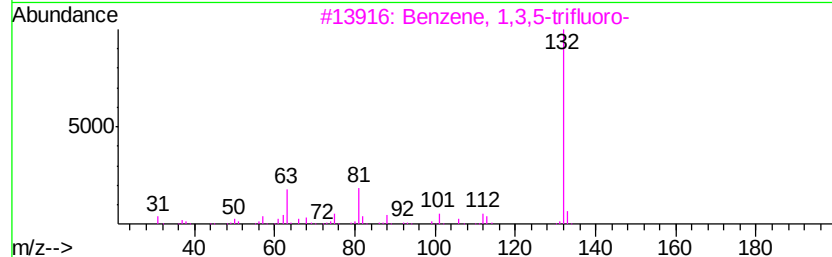
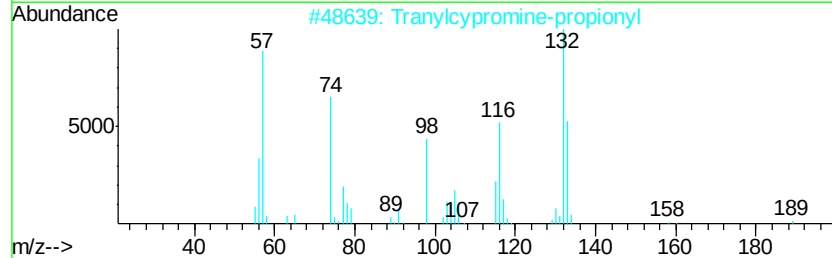
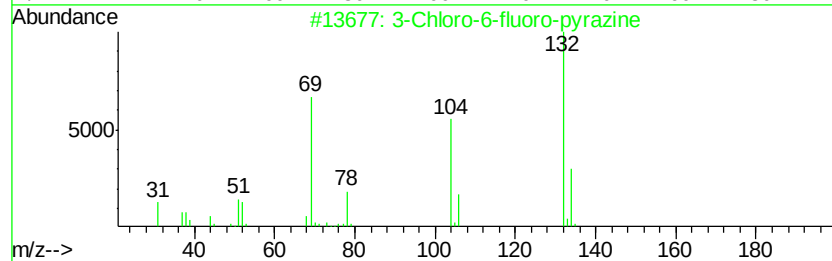
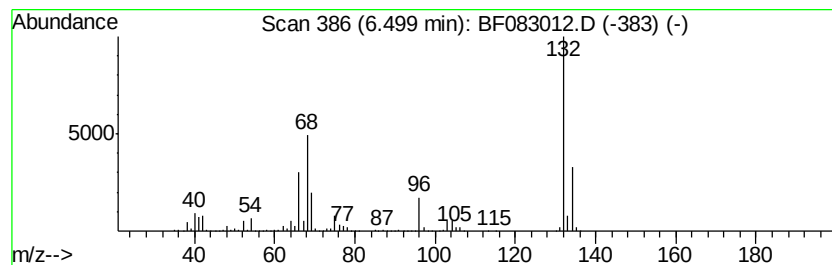
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Peak Number 2 unknown6.50 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	61.84 ng	2887940	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
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			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.89	19.6	ng	917733	1	6.73	933948	20.0
unknown6.50	6.50	61.8	ng	2887940	1	6.73	933948	20.0