

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050916\
 Data File : BF086833.D
 Acq On : 9 May 2016 21:07
 Operator : UM/SJ
 Sample : PB90374BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB90374BL

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-BF050916.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	4.864	241	243	250	rBV	226217	380318	7.06%	0.887%
2	5.299	278	281	283	rBV	3726542	4528211	84.03%	10.558%
3	6.350	370	373	375	rBV	4337256	4560740	84.63%	10.634%
4	6.464	381	383	385	rBV	4547887	4636786	86.05%	10.811%
5	6.693	400	403	405	rBV	774400	1022810	18.98%	2.385%
6	6.842	414	416	419	rBV	3404151	3276080	60.79%	7.639%
7	7.264	450	453	455	rBV	2589833	3151553	58.48%	7.348%
8	7.973	513	515	518	rBV	1341675	1337377	24.82%	3.118%
9	9.059	606	610	612	rBV	4101003	5202051	96.54%	12.129%
10	9.722	666	668	671	rBV	1584221	1551929	28.80%	3.618%
11	10.510	734	737	740	rBV2	2682191	3360511	62.36%	7.835%
12	11.208	795	798	800	rBV	1253337	1656228	30.73%	3.862%
13	12.785	933	936	939	rBV	4126144	5388752	100.00%	12.564%
14	13.825	1025	1027	1030	rBV	1437917	1556470	28.88%	3.629%
15	15.208	1145	1148	1150	rBV	834840	1278983	23.73%	2.982%

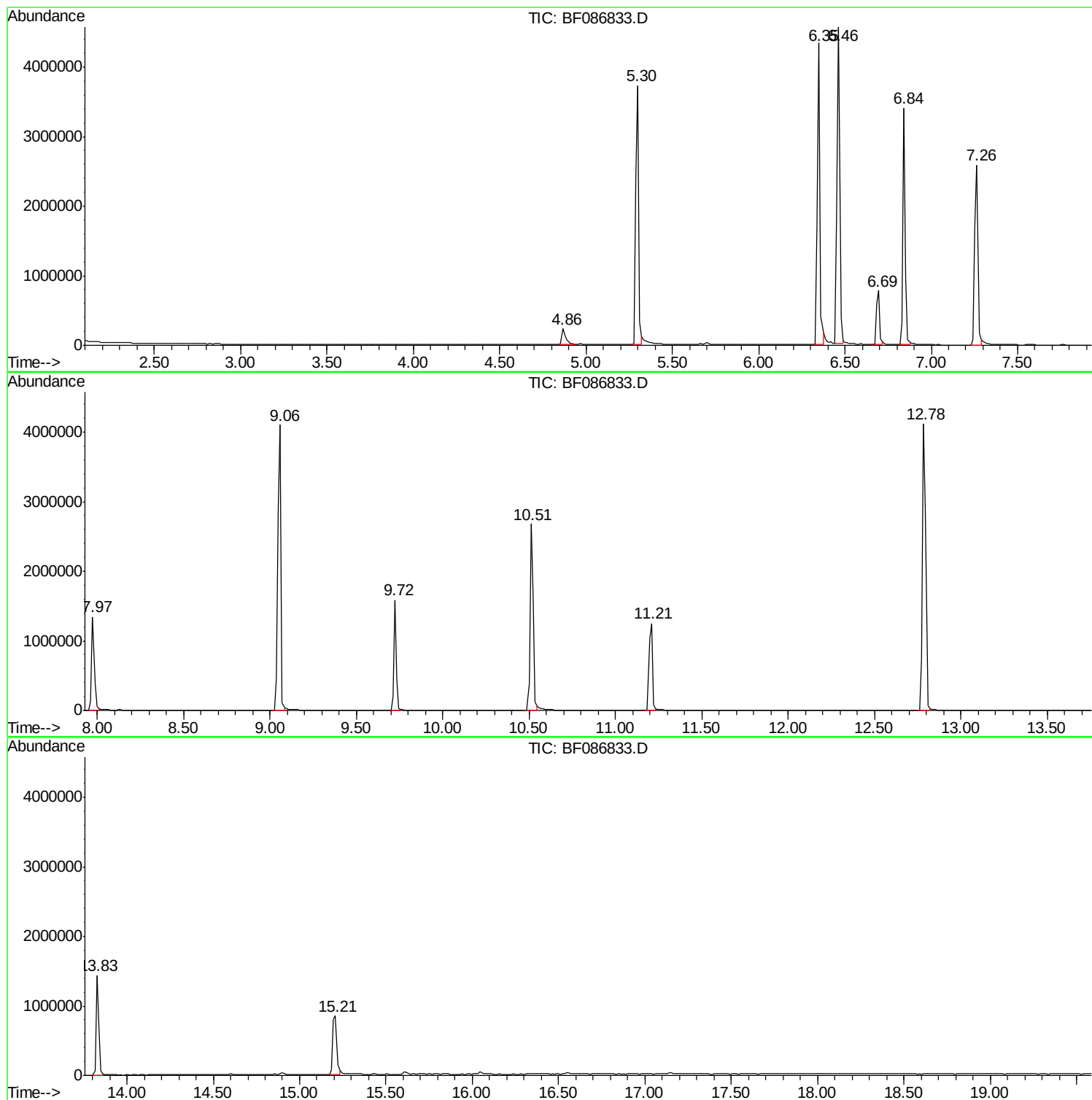
Sum of corrected areas: 42888799

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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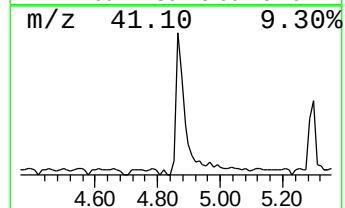
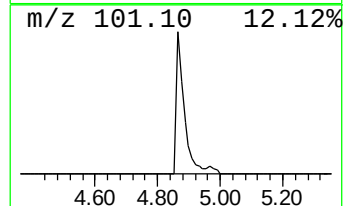
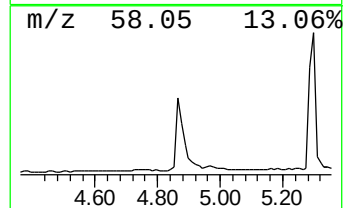
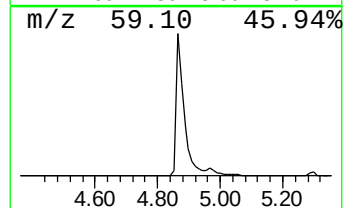
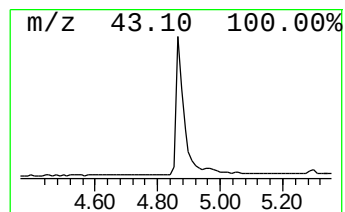
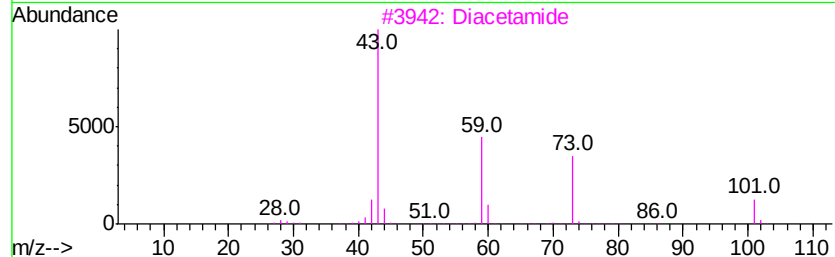
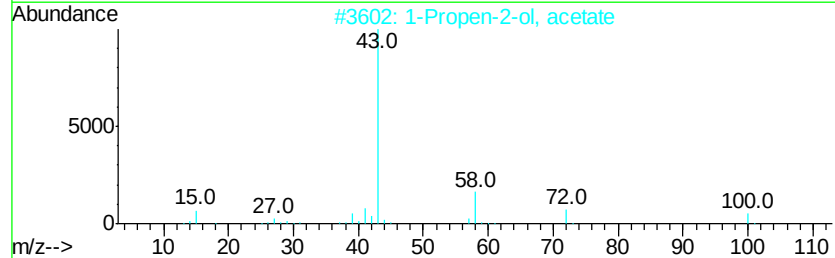
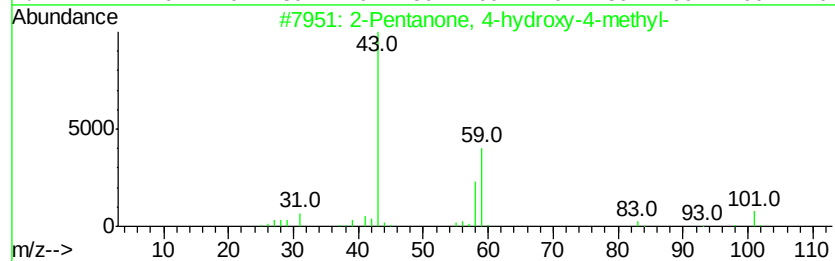
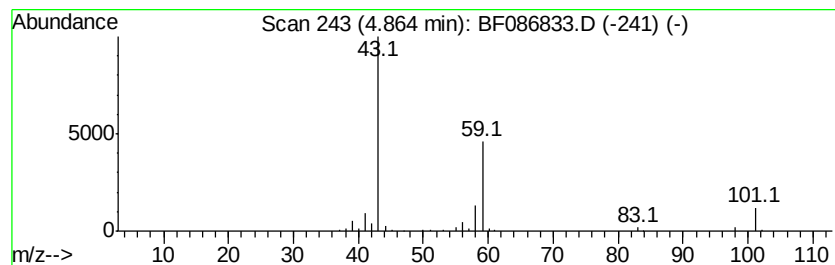
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	7.44 ng	380318	1,4-Dichlorobenzene-d4	6.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
3	Diacetamide	101	C4H7NO2	000625-77-4	9	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	
5	2-Pentanone, 5-(acetyloxy)-	144	C7H12O3	005185-97-7	9	



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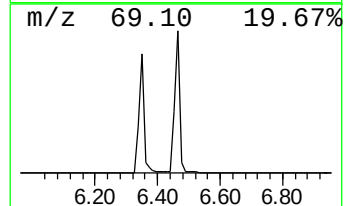
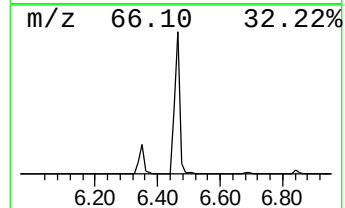
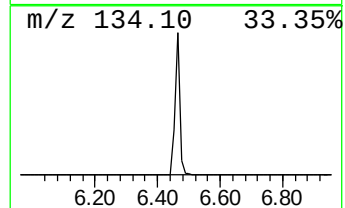
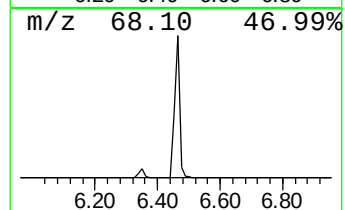
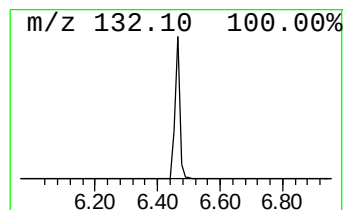
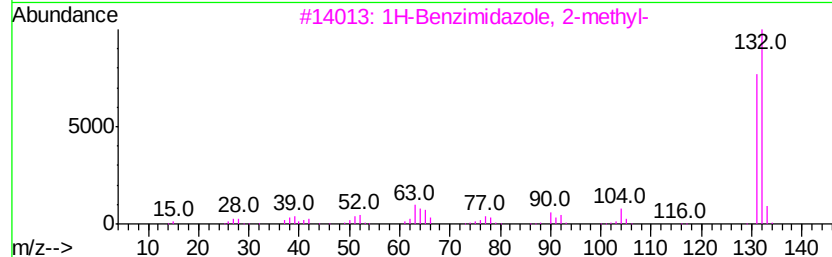
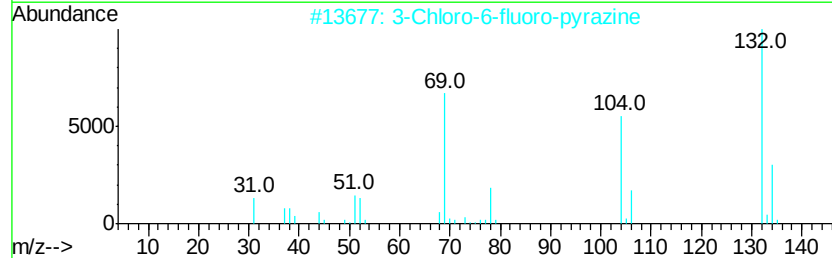
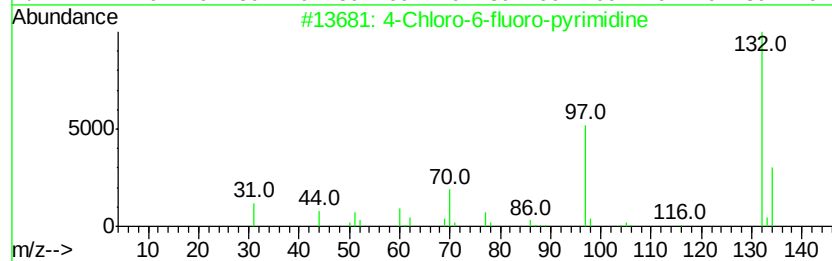
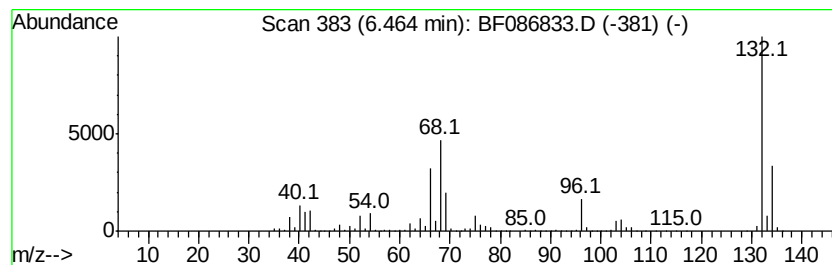
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Peak Number 2 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	90.67 ng	4636790	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.86	7.4	ng	380318	1	6.69	1022810	20.0
unknown6.46	6.46	90.7	ng	4636790	1	6.69	1022810	20.0