

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030816\
 Data File : BF085260.D
 Acq On : 8 Mar 2016 17:21
 Operator : UM/SJ
 Sample : PB88766BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88766BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.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	5.167	249	252	260	rBV	671016	847269	21.55%	2.598%
2	5.567	284	287	290	rBV	3043008	3132059	79.67%	9.603%
3	6.584	373	376	378	rBV	3263622	3469363	88.25%	10.637%
4	6.722	386	388	391	rBV	2757349	3470503	88.28%	10.641%
5	6.962	406	409	411	rVB	926106	862344	21.94%	2.644%
6	7.122	420	423	425	rBV	2242982	2883926	73.36%	8.842%
7	7.522	455	458	460	rBV	2271467	2311256	58.79%	7.087%
8	8.242	518	521	524	rBV	1309057	1147054	29.18%	3.517%
9	9.316	612	615	618	rBV	3027129	3931210	100.00%	12.054%
10	10.002	672	675	677	rBV	1379228	1276749	32.48%	3.915%
11	10.791	741	744	746	rBV	2339795	2549169	64.84%	7.816%
12	11.488	802	805	807	rBV	1262503	1304635	33.19%	4.000%
13	13.077	941	944	946	rBV	2636413	3747652	95.33%	11.491%
14	14.117	1033	1035	1038	rBV	1043948	956458	24.33%	2.933%
15	15.580	1160	1163	1166	rBV	499656	724829	18.44%	2.222%

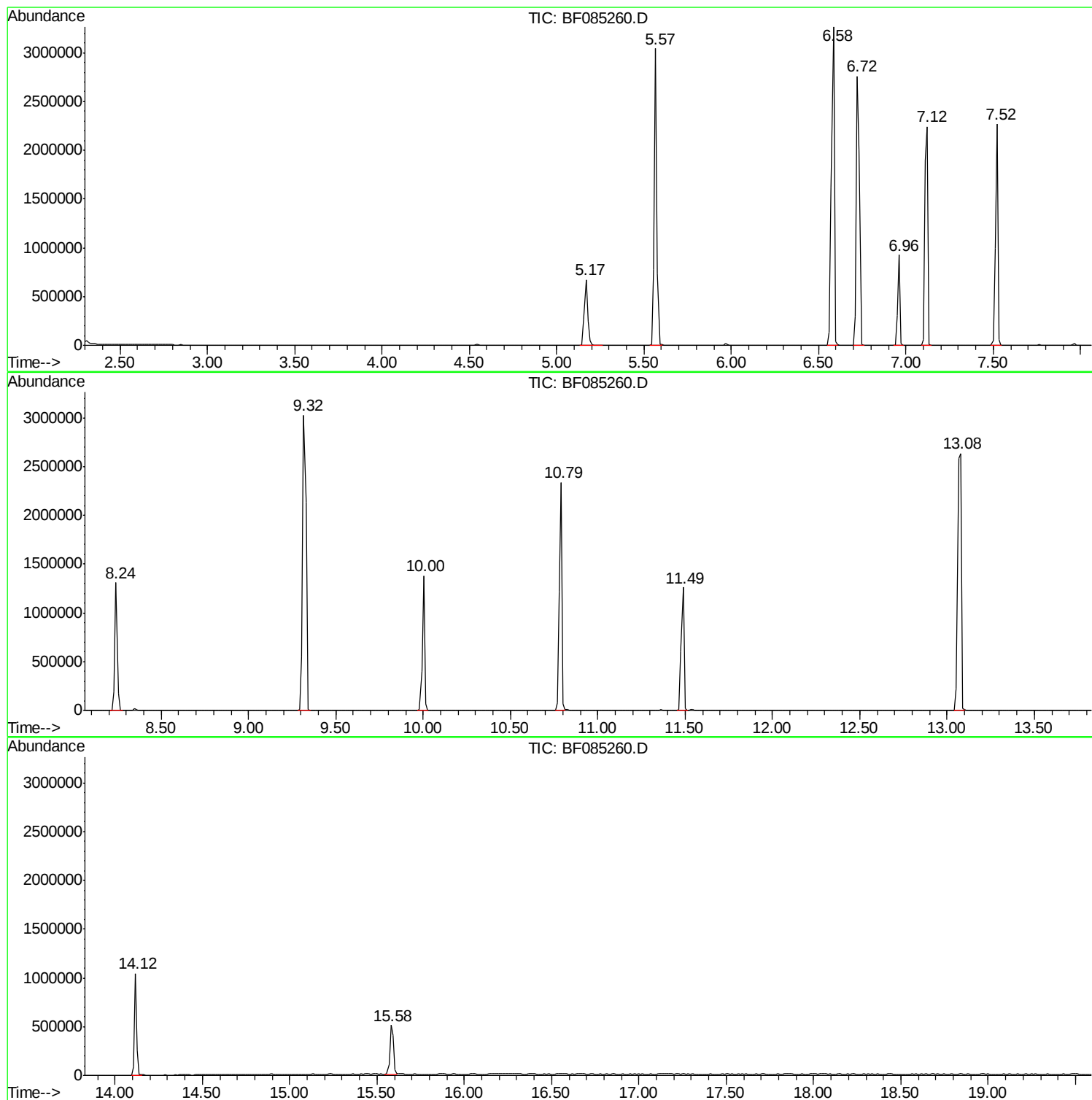
Sum of corrected areas: 32614476

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.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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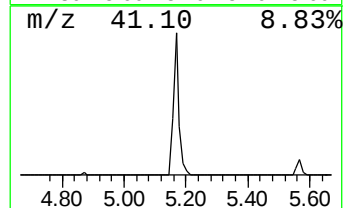
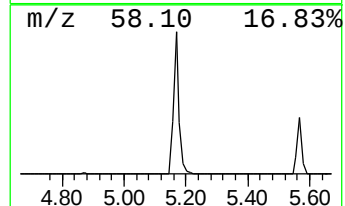
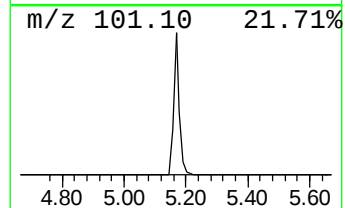
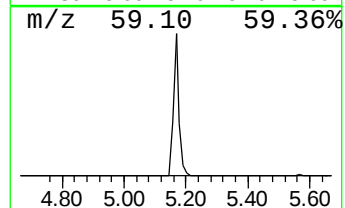
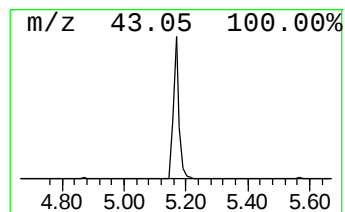
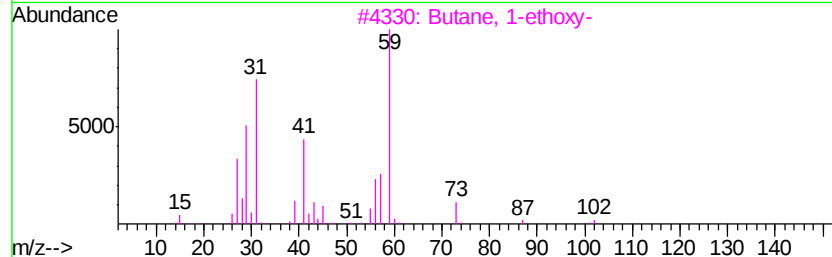
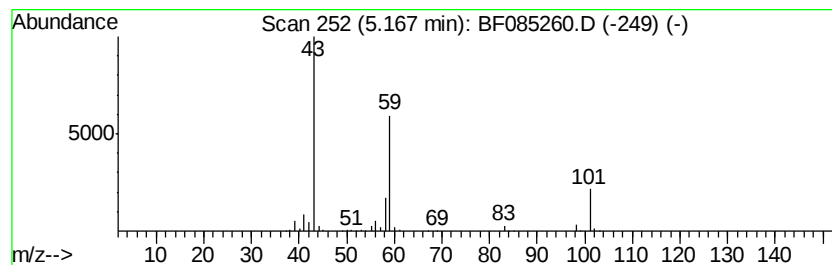
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	19.65 ng	847269	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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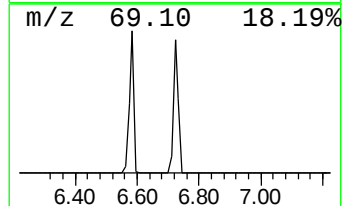
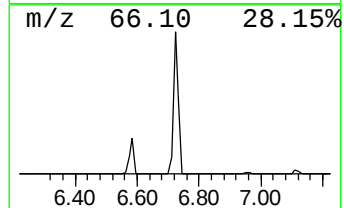
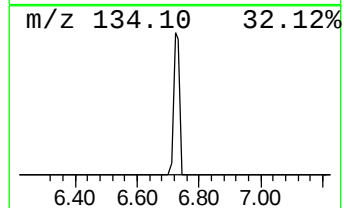
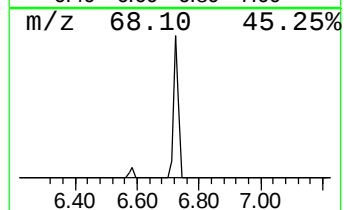
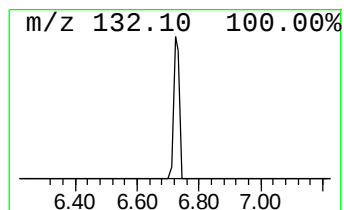
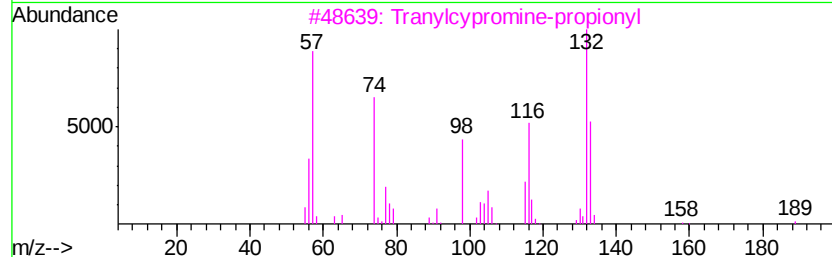
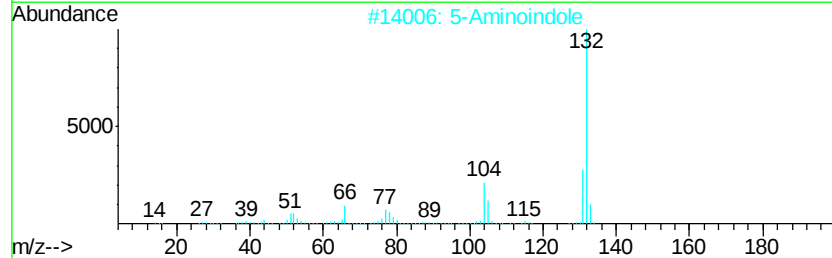
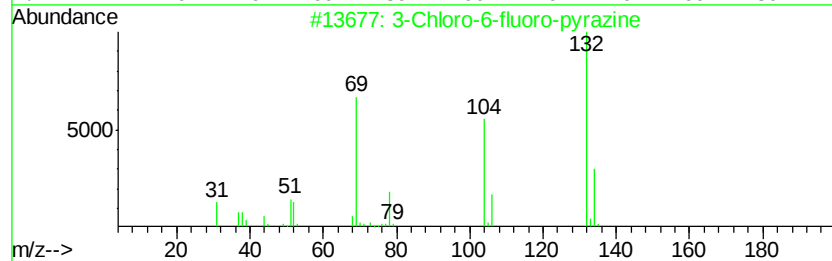
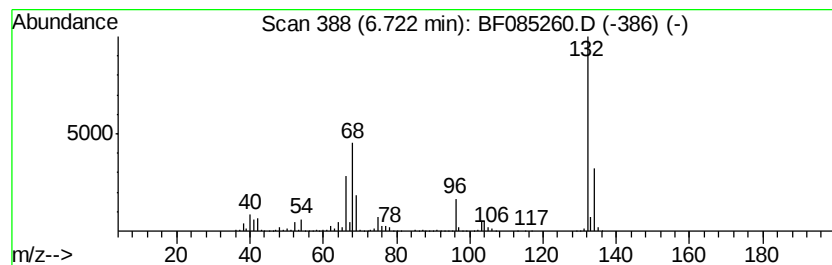
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Peak Number 2 unknown6.72 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	80.49 ng	3470500	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	19.6	ng	847269	1	6.96	862344	20.0
unknown6.72	6.72	80.5	ng	3470500	1	6.96	862344	20.0