

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
 Data File : BF085011.D
 Acq On : 11 Feb 2016 11:37
 Operator : SJ/IZ
 Sample : PB88297BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88297BL

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-BF020116.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	1.735	10	13	48	rVB	1837866	3669713	75.54%	9.216%
2	4.753	274	277	286	rBV	532369	912894	18.79%	2.293%
3	5.199	314	316	319	rBV	3090445	3524318	72.55%	8.851%
4	6.273	407	410	412	rBV	3289068	4120841	84.83%	10.349%
5	6.376	417	419	422	rBV	3436758	3491762	71.88%	8.769%
6	6.604	437	439	442	rBV	917480	834979	17.19%	2.097%
7	6.764	450	453	455	rBV	3301948	2931571	60.35%	7.362%
8	7.176	487	489	492	rBV	1838920	2414659	49.71%	6.064%
9	7.896	550	552	554	rBV	1309391	1155370	23.78%	2.902%
10	8.982	643	647	649	rBV	3710351	4671422	96.16%	11.732%
11	9.645	702	705	707	rBV	1487219	1324024	27.26%	3.325%
12	10.433	771	774	777	rBV	2779549	2777755	57.18%	6.976%
13	11.119	832	834	837	rBV	1212114	1273507	26.22%	3.198%
14	12.719	971	974	976	rVB	4043653	4857808	100.00%	12.200%
15	13.748	1062	1064	1067	rBV	1016617	1102242	22.69%	2.768%
16	15.108	1181	1183	1186	rBV	626876	755056	15.54%	1.896%

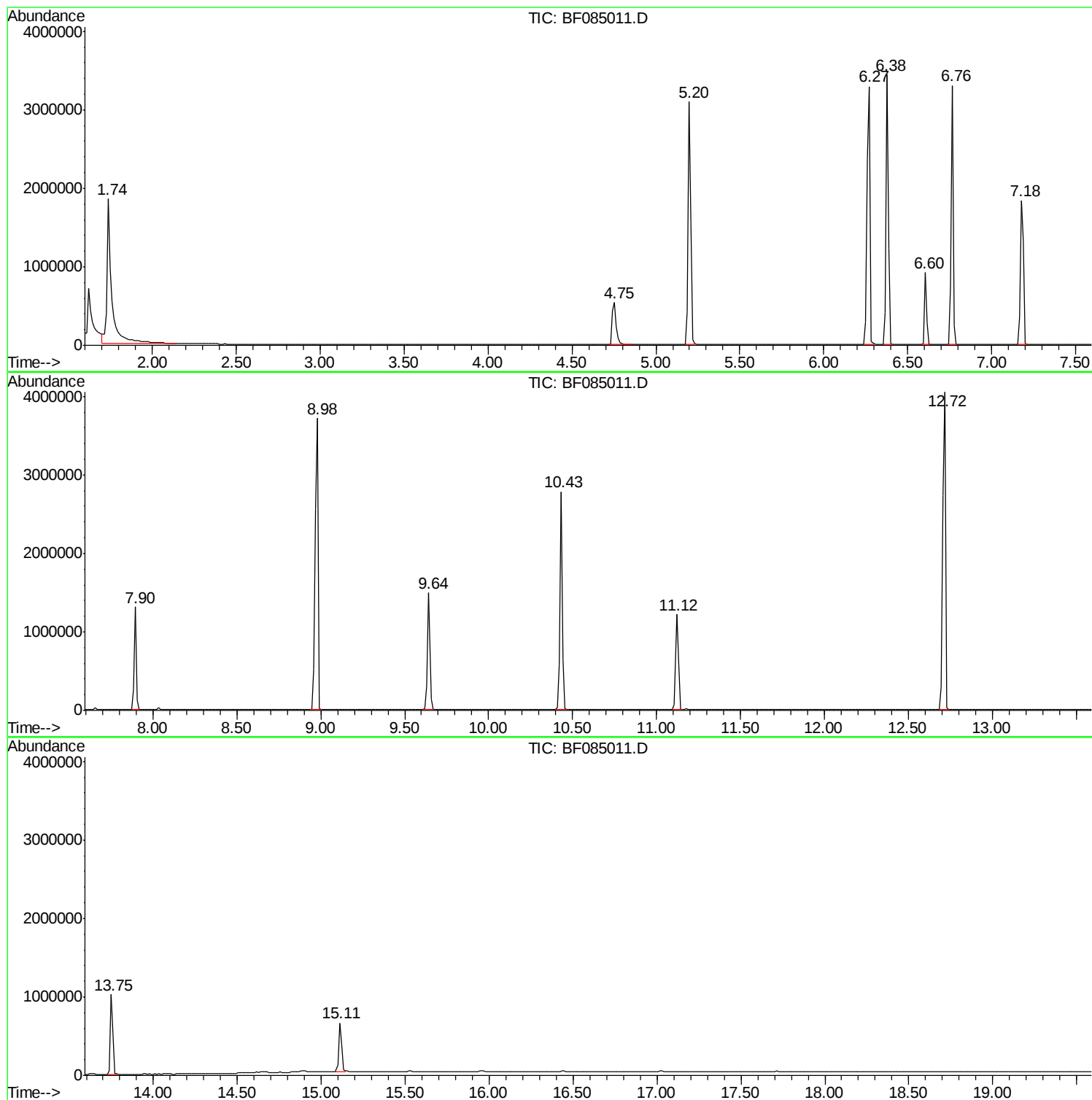
Sum of corrected areas: 39817921

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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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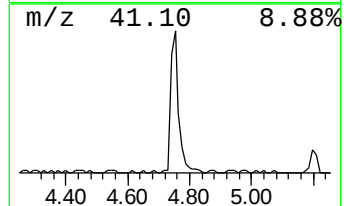
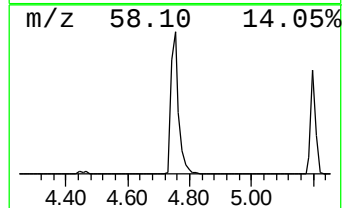
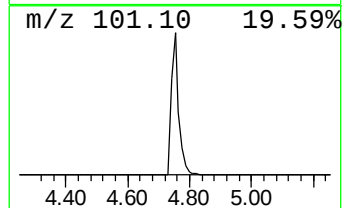
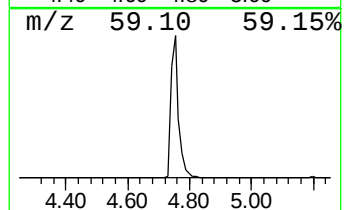
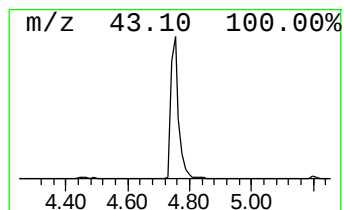
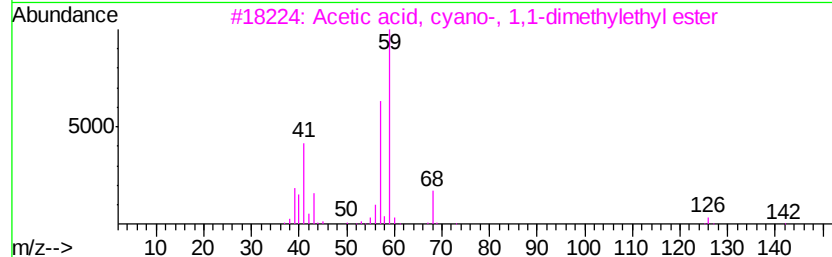
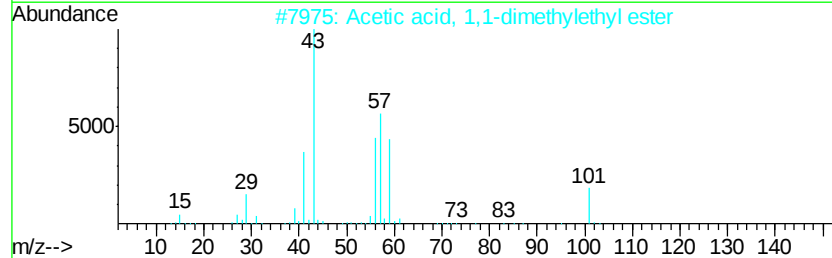
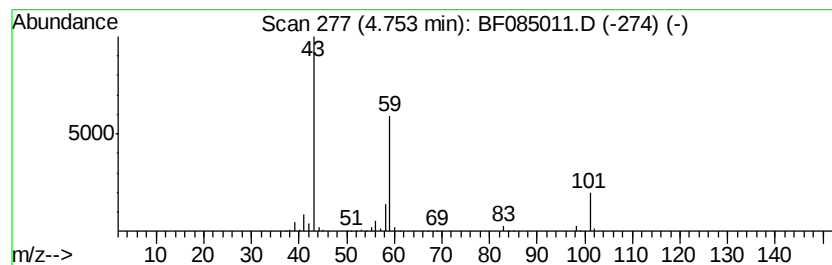
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.75	21.87 ng	912894	1,4-Dichlorobenzene-d4	6.60

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			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
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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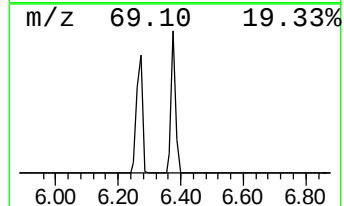
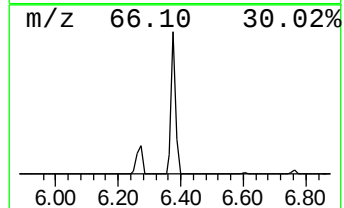
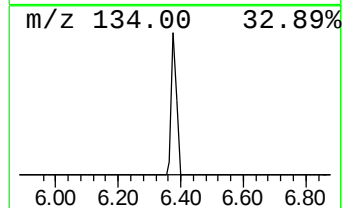
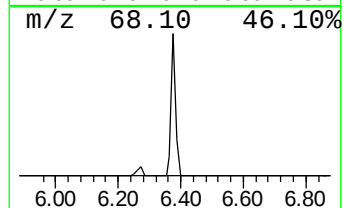
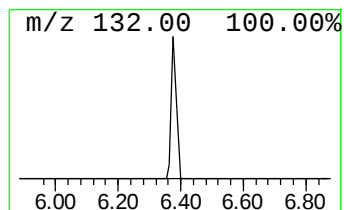
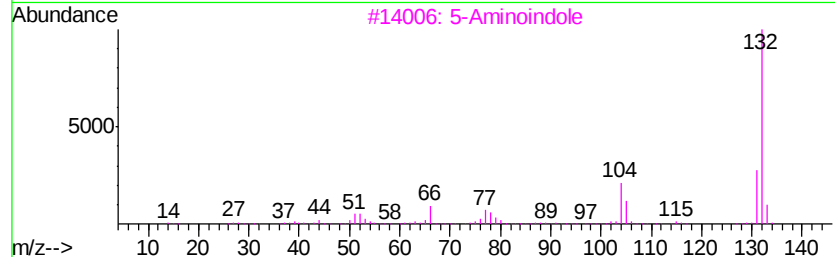
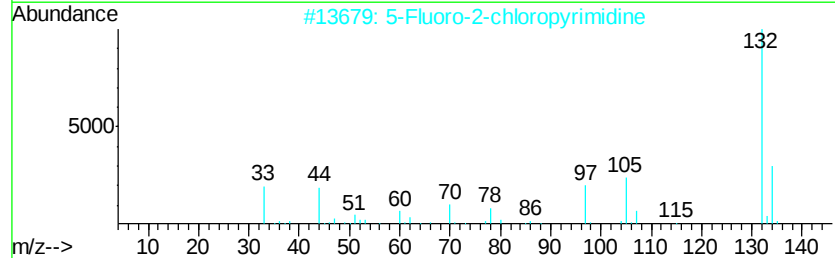
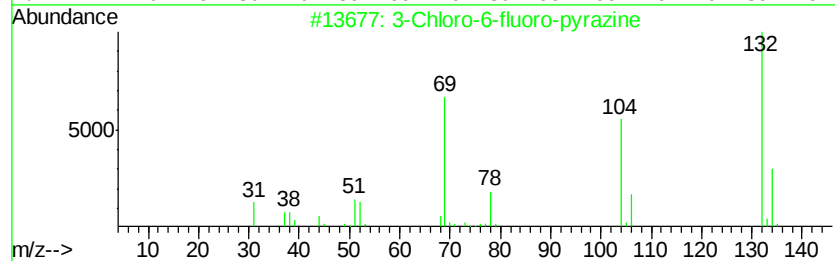
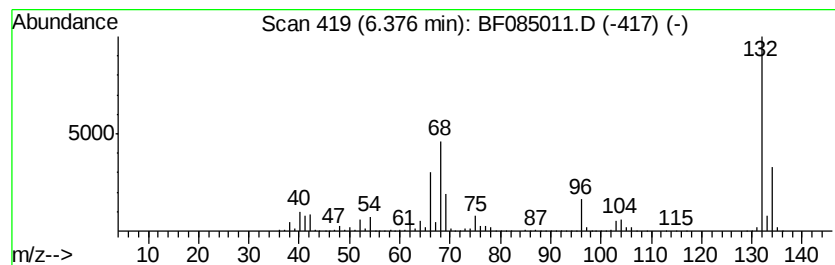
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Peak Number 3 unknown6.38 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	83.64 ng	3491760	1,4-Dichlorobenzene-d4	6.60

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	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.75	21.9	ng	912894	1	6.60	834979	20.0
unknown6.38	6.38	83.6	ng	3491760	1	6.60	834979	20.0