

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
 Data File : BF080162.D
 Acq On : 25 Jun 2015 13:22
 Operator : TP/IZ
 Sample : PB84184BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84184BL

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-BF061715.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	5.510	209	212	215	rBV	362355	345315	24.00%	3.513%
2	5.910	244	247	249	rBV	781685	948240	65.89%	9.647%
3	6.310	279	282	284	rBV	23819	29852	2.07%	0.304%
4	6.893	331	333	336	rBV	686111	941436	65.42%	9.578%
5	7.064	345	348	350	rBV	1120314	1023881	71.15%	10.417%
6	7.293	366	368	370	rVB	214489	197677	13.74%	2.011%
7	7.453	380	382	385	rVB	913341	778571	54.10%	7.921%
8	7.853	415	417	419	rBV	772575	623596	43.33%	6.345%
9	8.584	479	481	484	rBV	302369	267545	18.59%	2.722%
10	9.659	573	575	578	rBV	1196068	1156397	80.36%	11.765%
11	10.356	634	636	639	rBV	342881	302370	21.01%	3.076%
12	11.156	703	706	708	rBV2	588147	734689	51.05%	7.475%
13	11.865	766	768	770	rBV	397951	331422	23.03%	3.372%
14	13.556	914	916	919	rBV	1579790	1439053	100.00%	14.641%
15	14.837	1026	1028	1031	rBV	349777	356301	24.76%	3.625%
16	16.722	1190	1193	1197	rBV	257661	352562	24.50%	3.587%

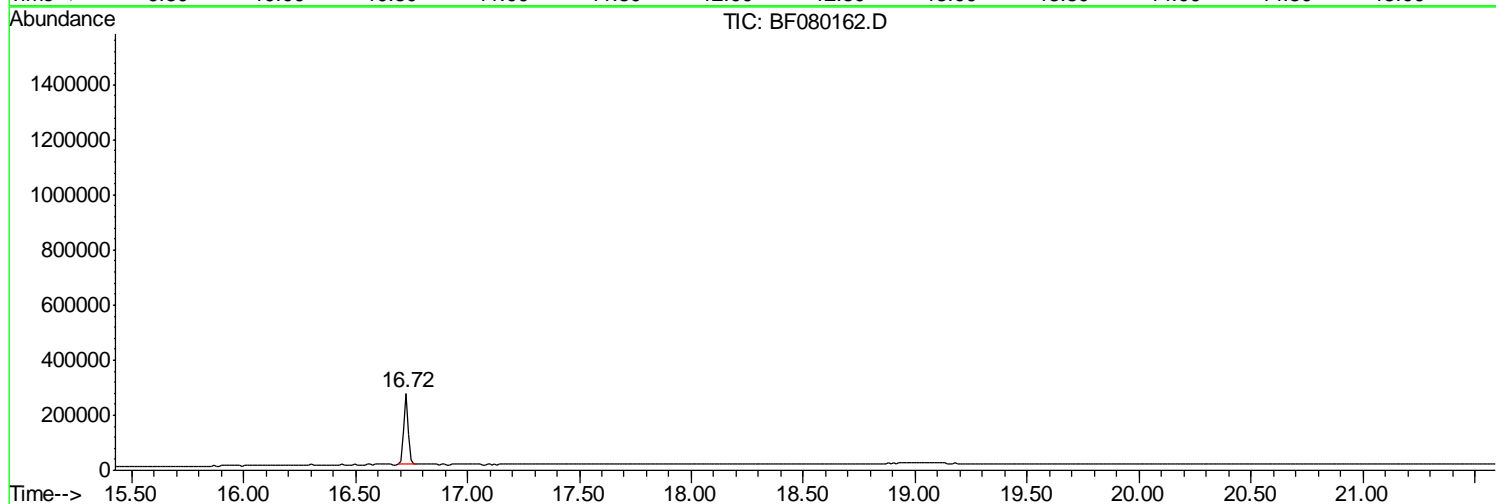
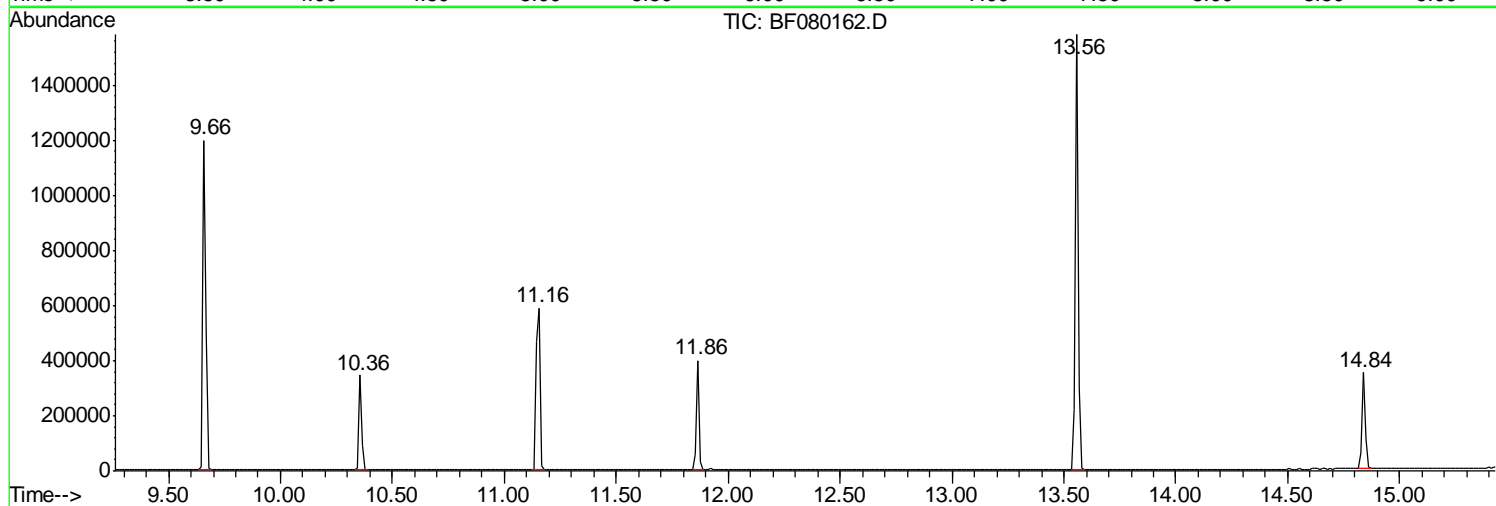
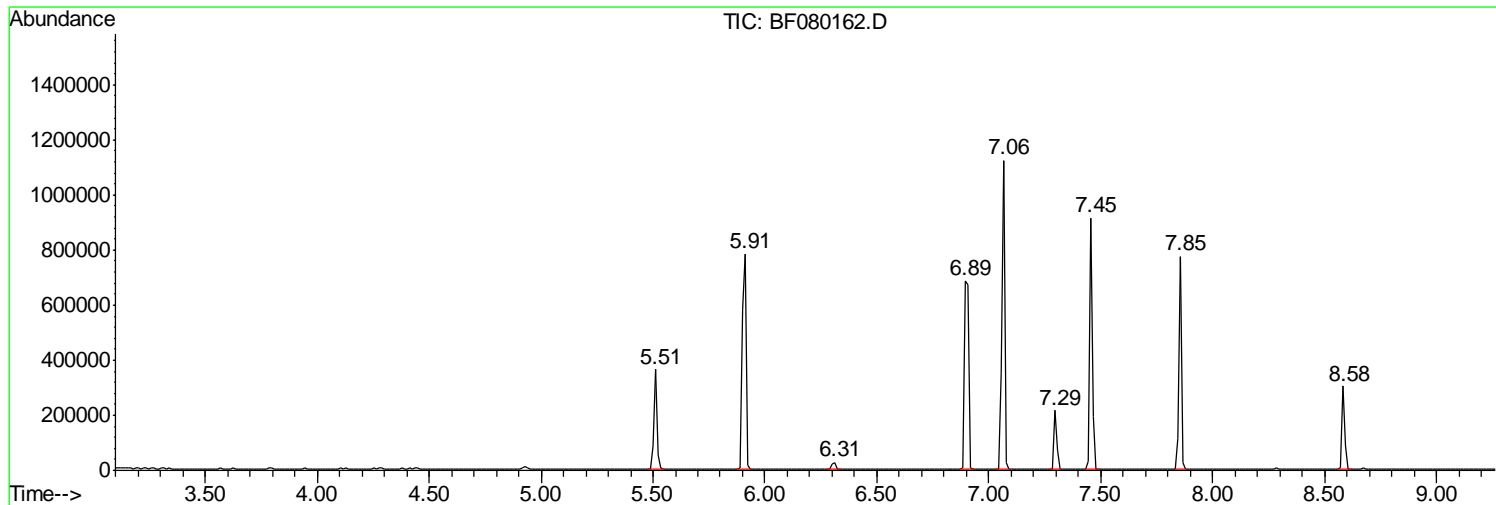
Sum of corrected areas: 9828907

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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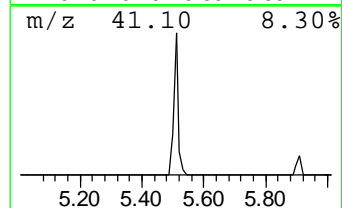
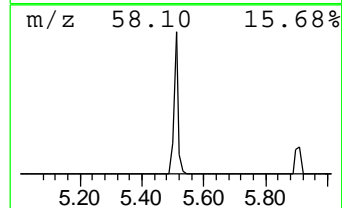
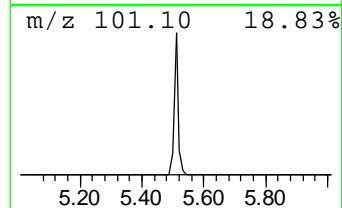
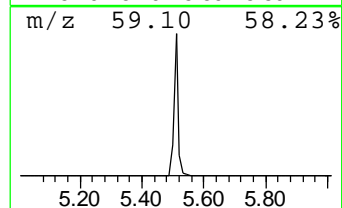
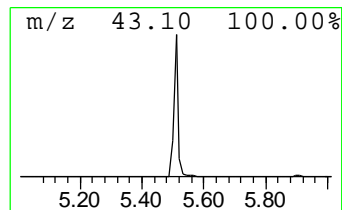
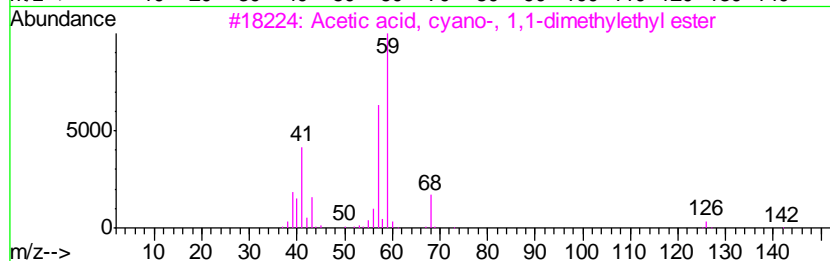
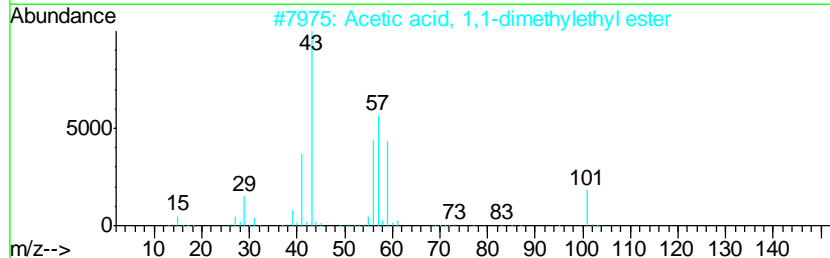
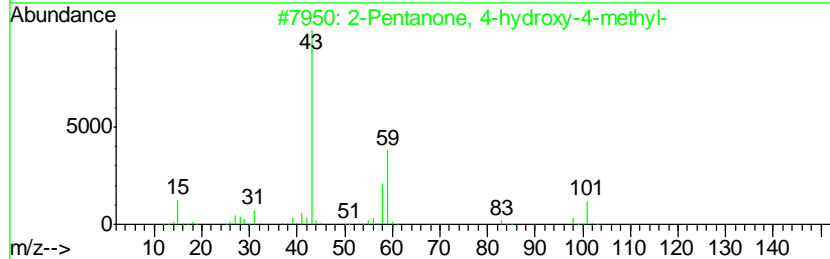
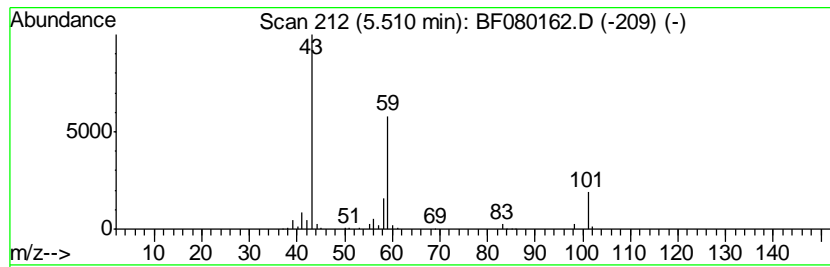
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 Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.51	34.94 ng	345315	1,4-Dichlorobenzene-d4	7.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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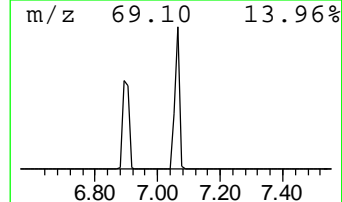
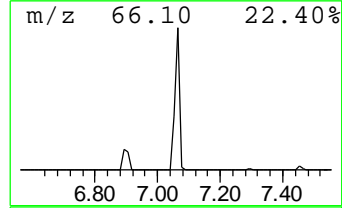
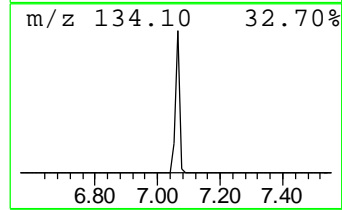
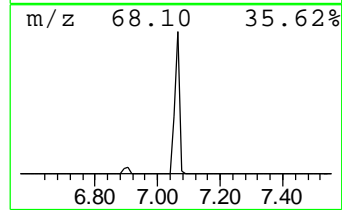
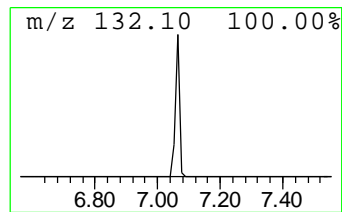
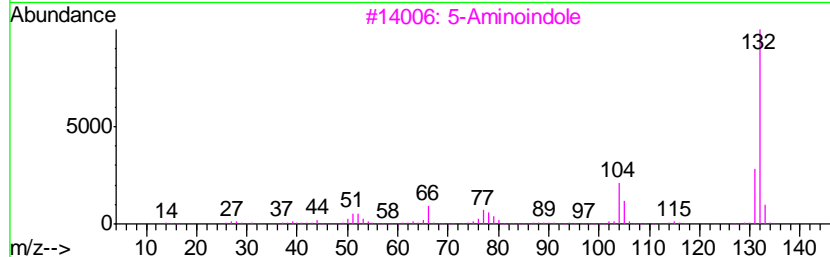
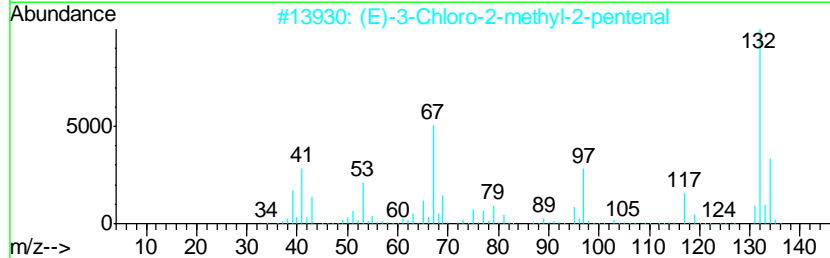
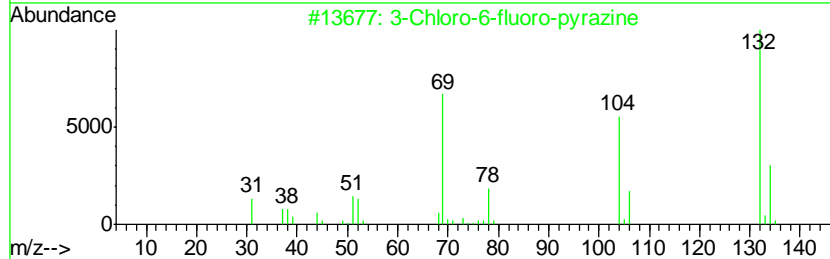
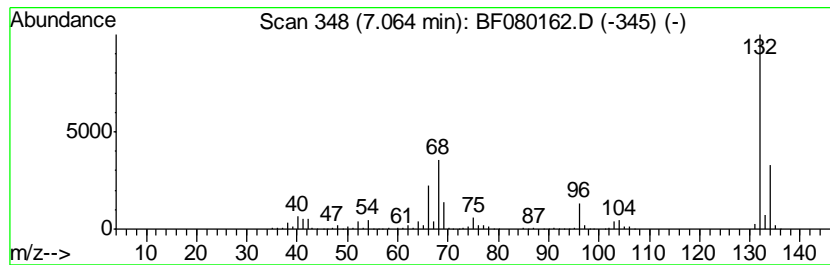
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 Peak Number 3 unknown7.06 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	103.59 ng	1023880	1,4-Dichlorobenzene-d4	7.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.51	34.9	ng	345315	1	7.29	197677	20.0
unknown7.06	7.06	103.6	ng	1023880	1	7.29	197677	20.0