

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090616\
 Data File : BF089992.D
 Acq On : 6 Sep 2016 17:04
 Operator : UM/SJ
 Sample : PB93146BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93146BL

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-BF083116.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	11	13	62	rVB	3169398	8146112	100.00%	13.377%
2	4.764	275	278	292	rBV	678769	1313187	16.12%	2.157%
3	5.210	313	317	319	rBV	3670519	5094248	62.54%	8.366%
4	6.273	406	410	412	rBV	3937436	5517857	67.74%	9.061%
5	6.387	417	420	422	rBV	4569948	5339451	65.55%	8.768%
6	6.616	438	440	451	rVB	1243343	1320138	16.21%	2.168%
7	6.776	451	454	456	rBV	3229001	4169620	51.19%	6.847%
8	7.187	487	490	492	rBV	2916836	3578996	43.94%	5.877%
9	7.896	550	552	555	rBV	1325346	1659928	20.38%	2.726%
10	8.982	643	647	649	rBV	5541499	6038829	74.13%	9.917%
11	9.645	702	705	708	rBV	2126573	2067373	25.38%	3.395%
12	10.433	771	774	776	rBV	3554746	3947513	48.46%	6.483%
13	11.119	831	834	836	rBV	2241774	2296858	28.20%	3.772%
14	12.708	969	973	975	rBV	5122299	6689749	82.12%	10.986%
15	13.725	1060	1062	1065	rBV	1780846	2056543	25.25%	3.377%
16	15.074	1177	1180	1182	rBV	1443214	1657756	20.35%	2.722%

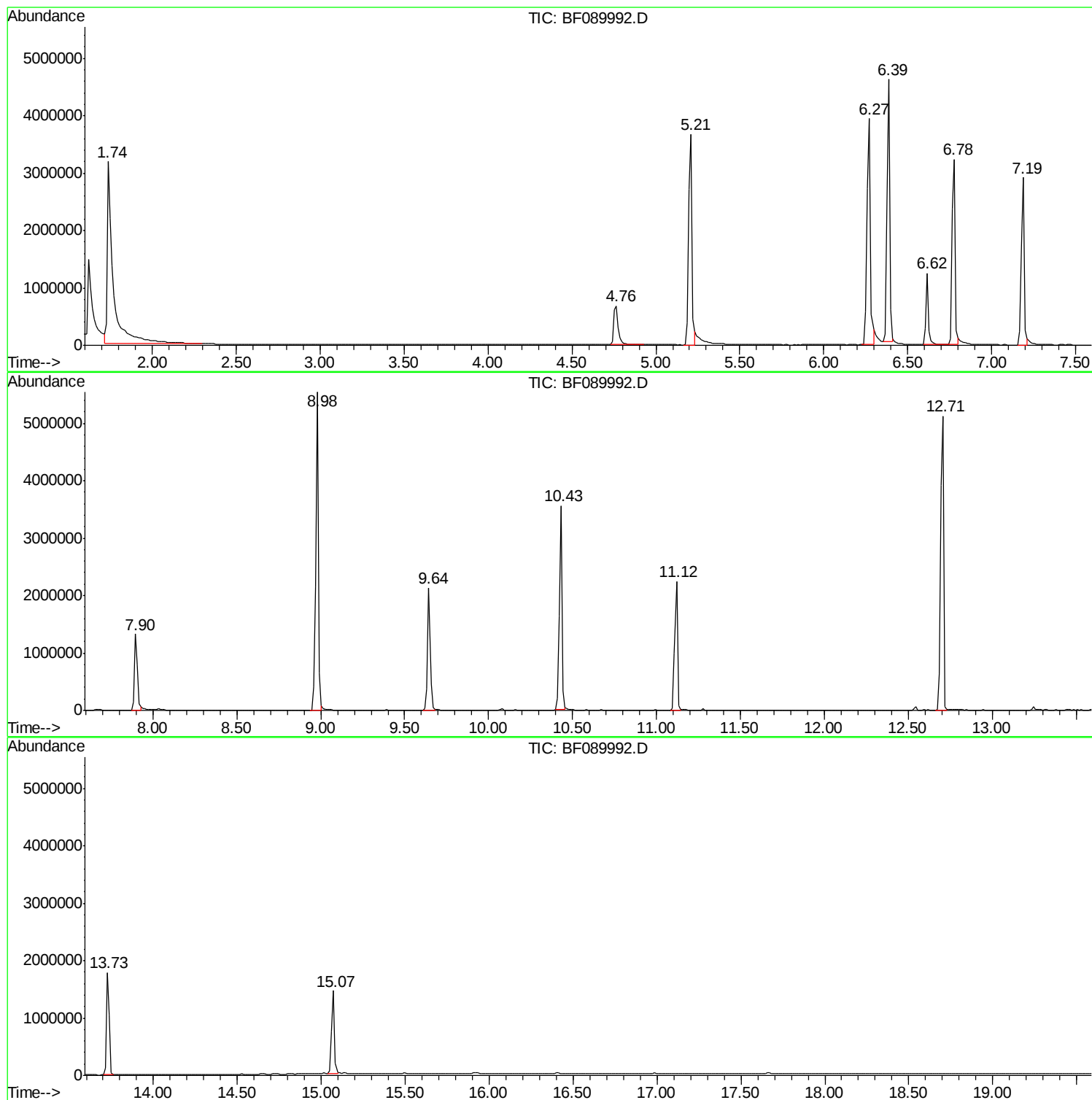
Sum of corrected areas: 60894158

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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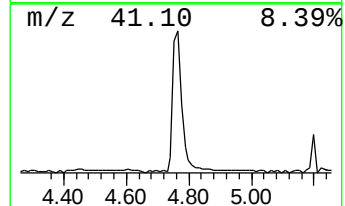
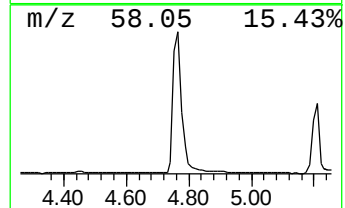
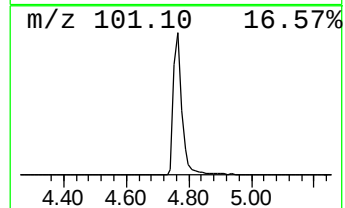
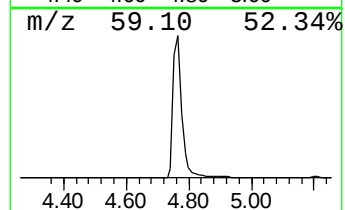
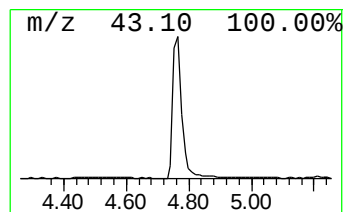
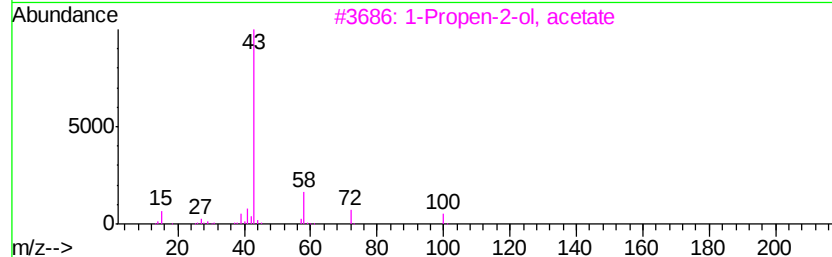
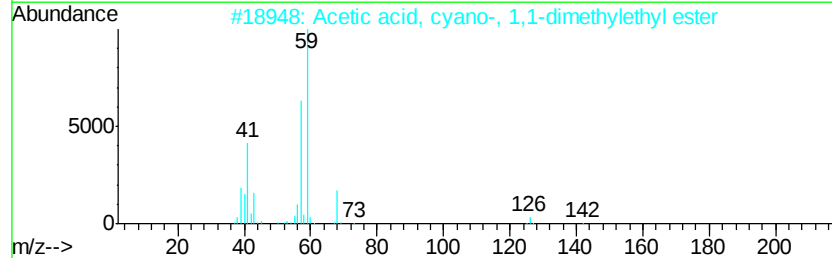
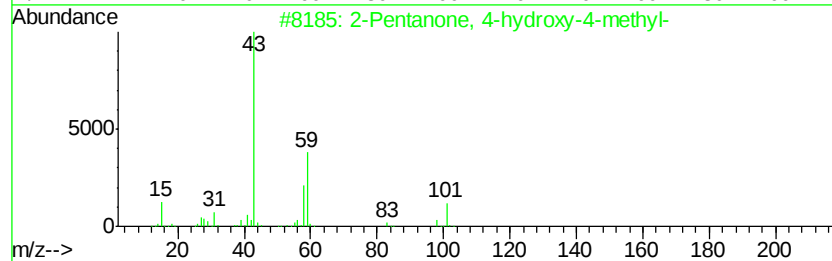
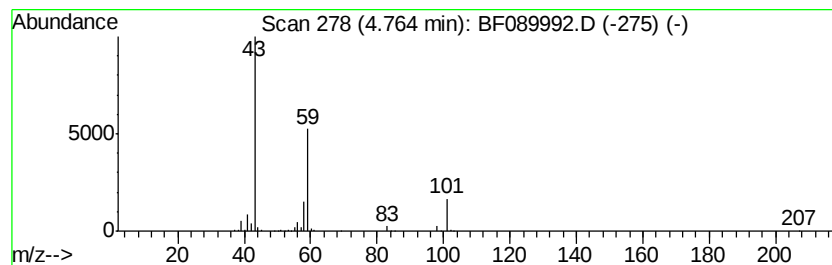
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.76	19.89 ng	1313190	1,4-Dichlorobenzene-d4	6.62

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	25
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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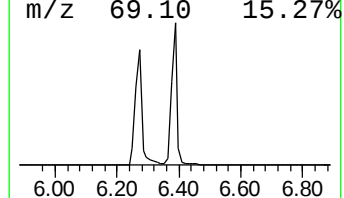
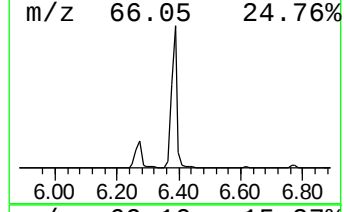
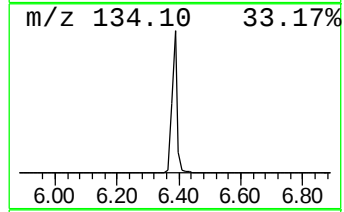
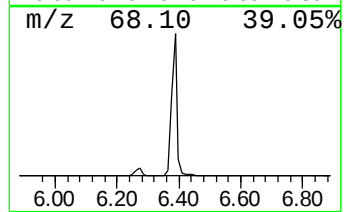
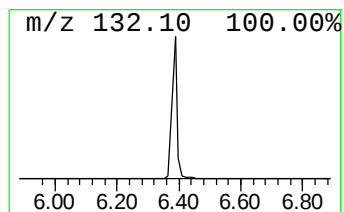
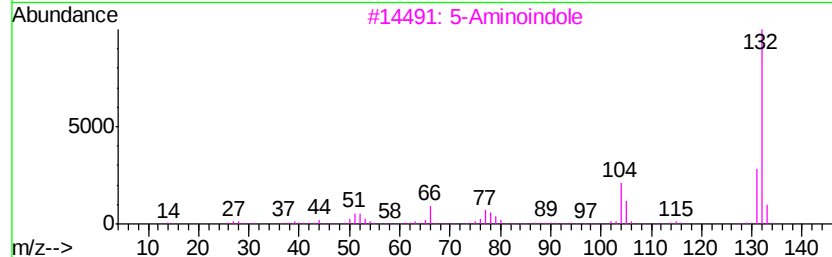
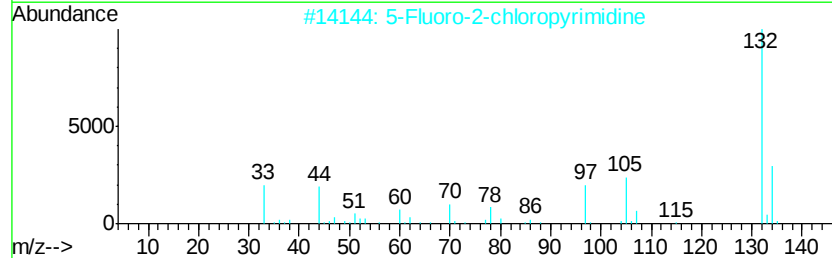
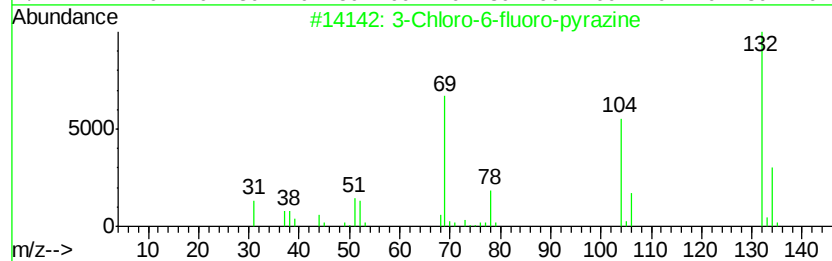
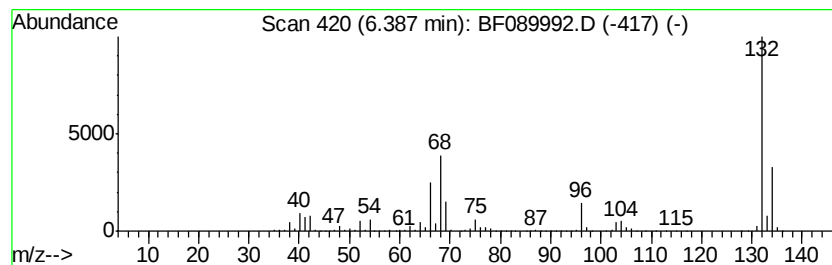
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Peak Number 3 unknown6.39 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	80.89 ng	5339450	1,4-Dichlorobenzene-d4	6.62

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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.76	19.9	ng	1313190	1	6.62	1320140	20.0
unknown6.39	6.39	80.9	ng	5339450	1	6.62	1320140	20.0