

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092915\
 Data File : BF081889.D
 Acq On : 29 Sep 2015 20:03
 Operator : UM/IZ
 Sample : PB85717BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85717BL

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-BF092815.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	5.120	257	261	273	rBV	546613	829241	27.29%	3.986%
2	5.520	294	296	312	rBV	1531800	1649599	54.29%	7.930%
3	6.549	383	386	388	rBV	1370743	1635942	53.84%	7.864%
4	6.686	396	398	400	rBV	1702922	1664828	54.79%	8.003%
5	6.926	417	419	430	rVB	251653	409640	13.48%	1.969%
6	7.074	430	432	449	rBV	1391776	1401264	46.11%	6.736%
7	7.486	466	468	500	rBV	1000644	1221312	40.19%	5.871%
8	8.206	529	531	552	rBV	481626	613422	20.19%	2.949%
9	9.280	623	625	628	rBV	1923998	2704120	88.99%	13.000%
10	9.966	682	685	687	rBV	915022	919576	30.26%	4.421%
11	10.755	751	754	774	rBV	1493716	1748148	57.53%	8.404%
12	11.452	813	815	828	rBV	850540	978290	32.19%	4.703%
13	13.041	952	954	957	rBV	2816811	3038755	100.00%	14.608%
14	14.092	1044	1046	1061	rBV	894336	1096860	36.10%	5.273%
15	15.567	1171	1175	1190	rBV	474367	890681	29.31%	4.282%

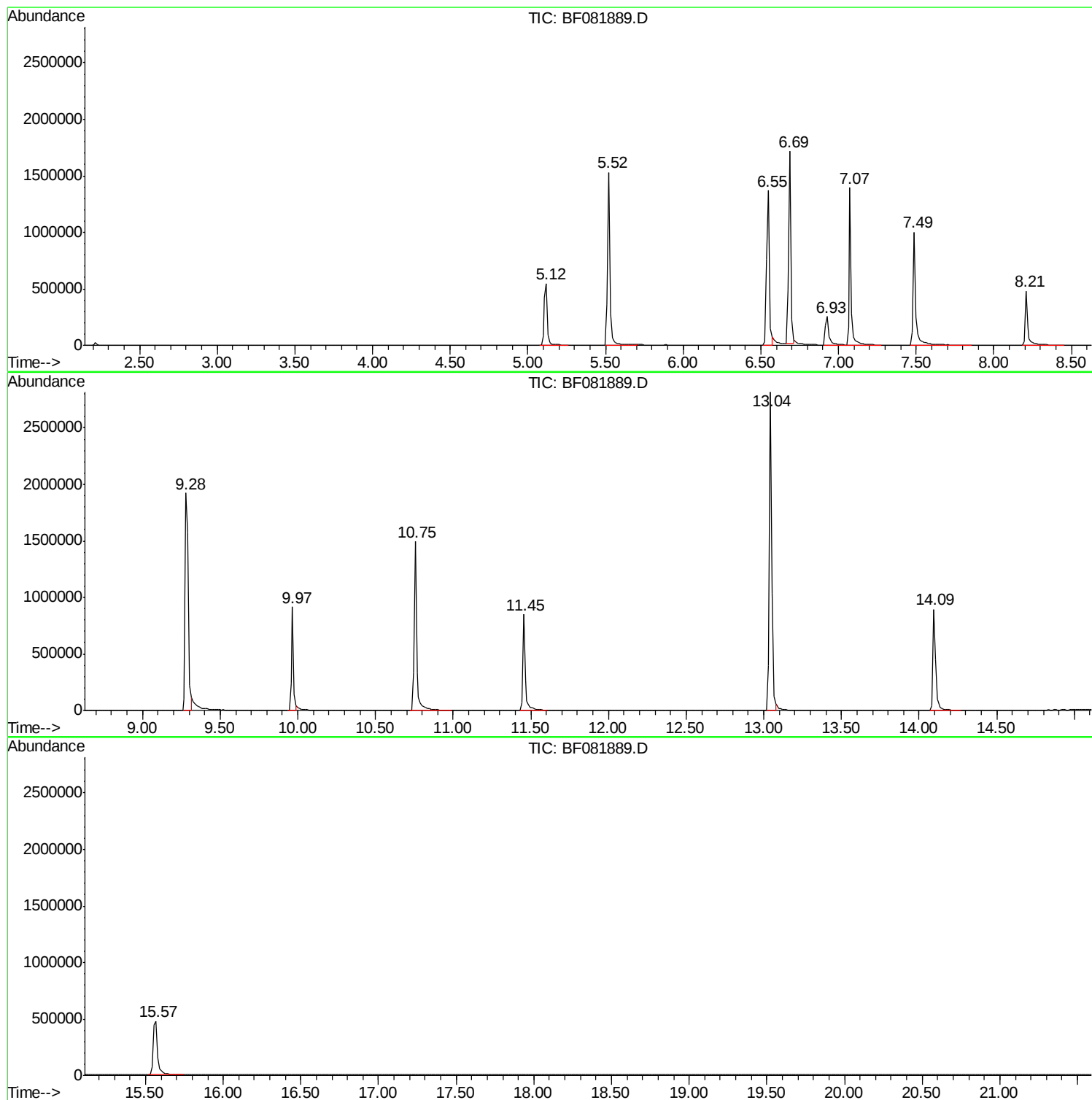
Sum of corrected areas: 20801678

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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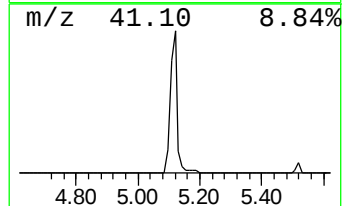
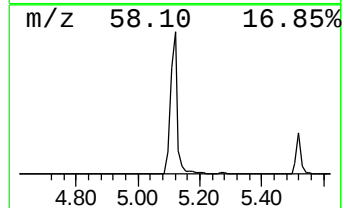
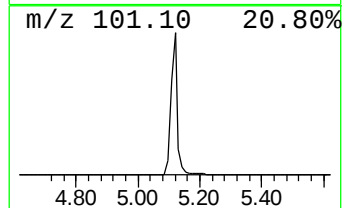
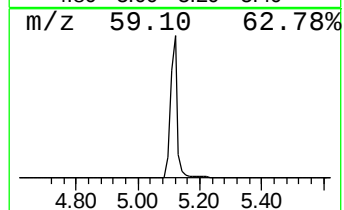
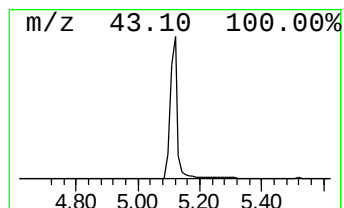
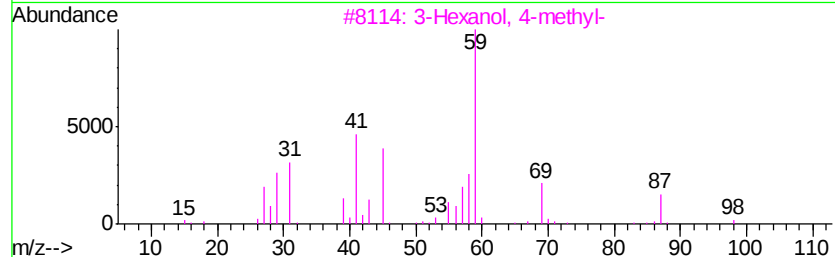
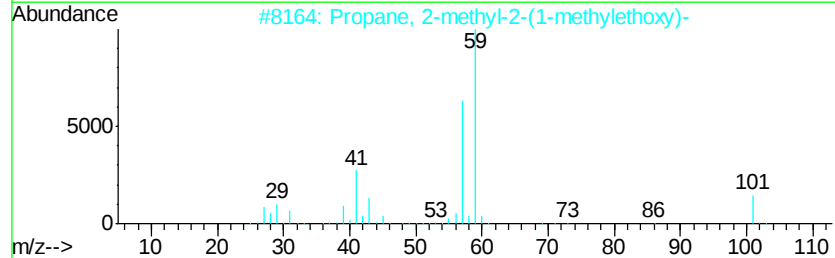
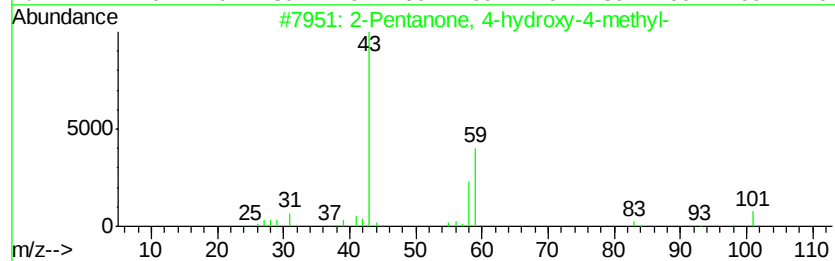
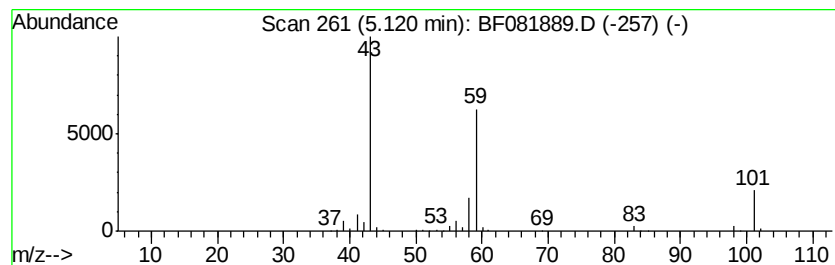
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	40.49 ng	829241	1,4-Dichlorobenzene-d4	6.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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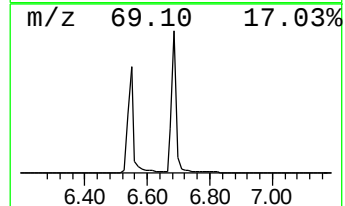
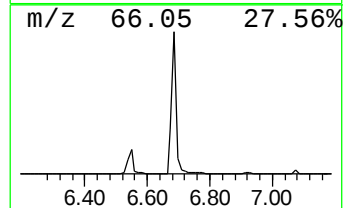
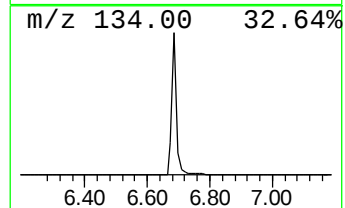
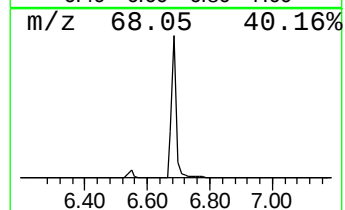
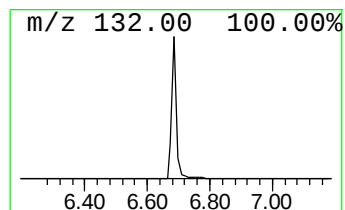
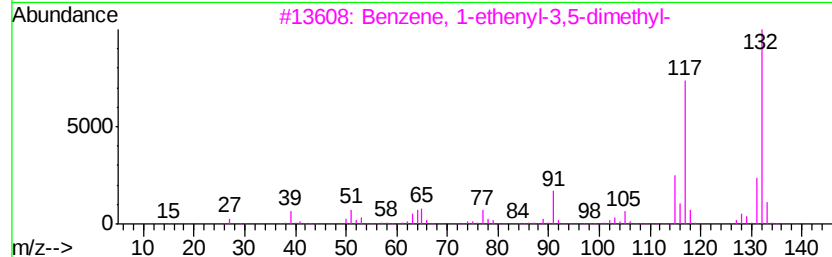
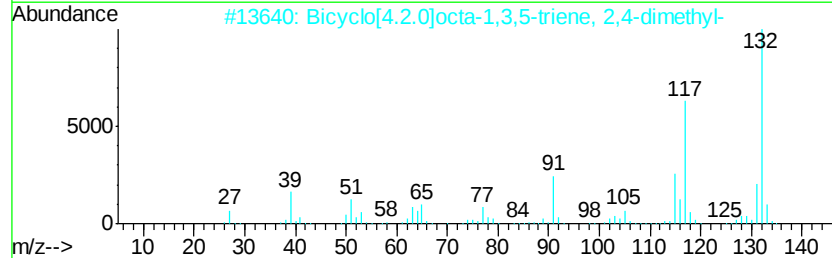
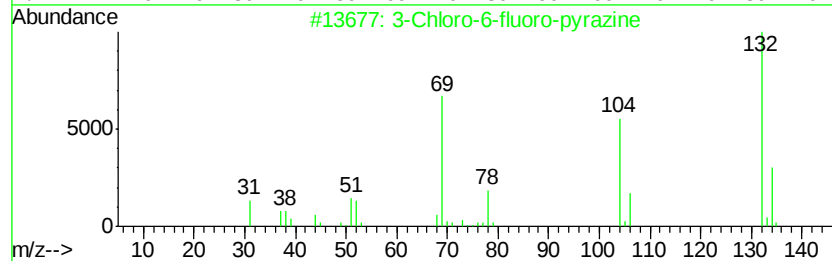
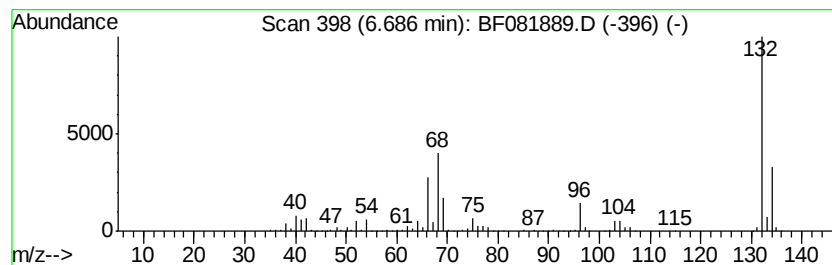
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Peak Number 2 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	81.28 ng	1664830	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	40.5	ng	829241	1	6.93	409640	20.0
unknown6.69	6.69	81.3	ng	1664830	1	6.93	409640	20.0