

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062616\
 Data File : BF088451.D
 Acq On : 27 Jun 2016 12:51
 Operator : UM/SJ
 Sample : PB91593BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB91593BL

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-BF060716.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	1.678	7	8	52	rVB	915980	2209844	100.00%	13.790%
2	4.684	269	271	285	rBV	37943	80718	3.65%	0.504%
3	5.130	307	310	332	rBV	1217884	1516113	68.61%	9.461%
4	6.204	402	404	411	rBV	1165325	1328936	60.14%	8.293%
5	6.307	411	413	431	rVB	1594022	1582058	71.59%	9.872%
6	6.536	431	433	444	rBV	291197	317678	14.38%	1.982%
7	6.684	444	446	449	rBV	1066585	1226780	55.51%	7.655%
8	7.107	481	483	506	rBV	635131	957979	43.35%	5.978%
9	7.827	543	546	558	rBV	297500	458469	20.75%	2.861%
10	8.902	637	640	643	rBV	1672271	1854565	83.92%	11.573%
11	9.565	696	698	701	rBV	474077	535500	24.23%	3.342%
12	10.353	765	767	782	rBB2	379433	516847	23.39%	3.225%
13	11.051	825	828	830	rBV	395230	534761	24.20%	3.337%
14	12.628	963	966	969	rBV	1473282	1967401	89.03%	12.277%
15	13.668	1055	1057	1072	rBV2	373695	535357	24.23%	3.341%
16	15.028	1173	1176	1188	rBV	253059	401998	18.19%	2.509%

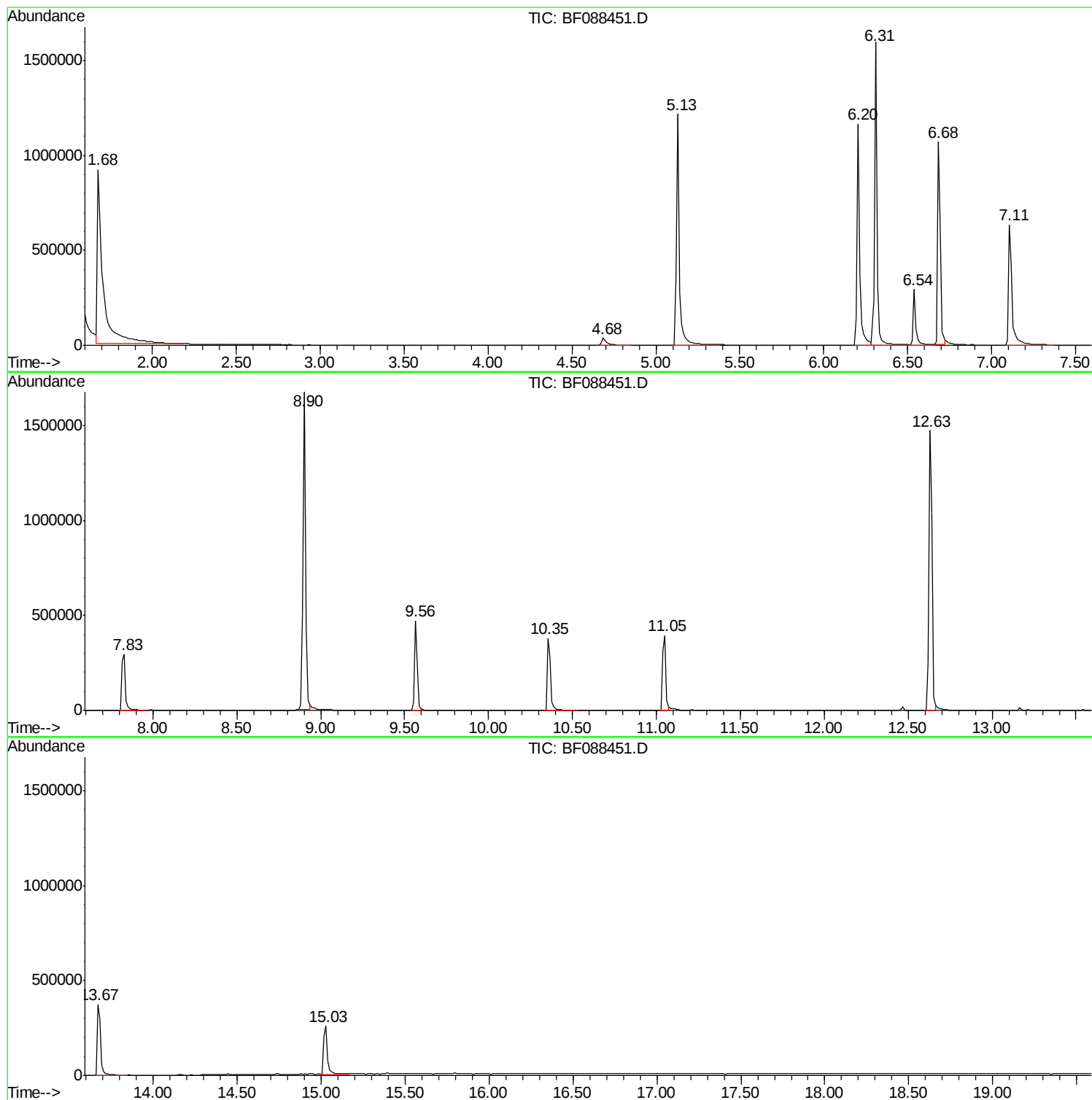
Sum of corrected areas: 16025004

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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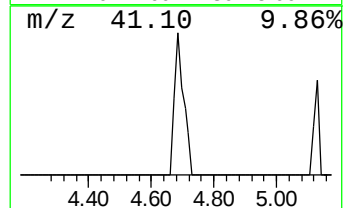
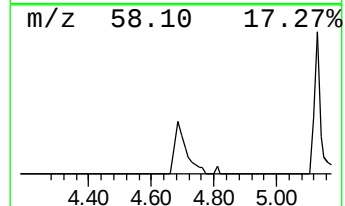
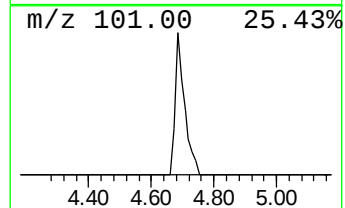
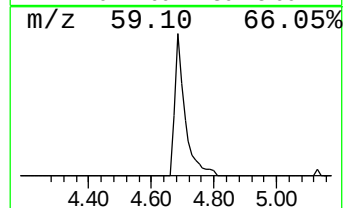
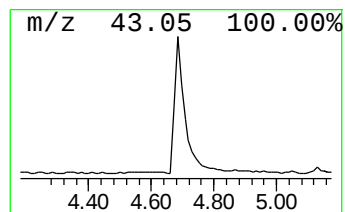
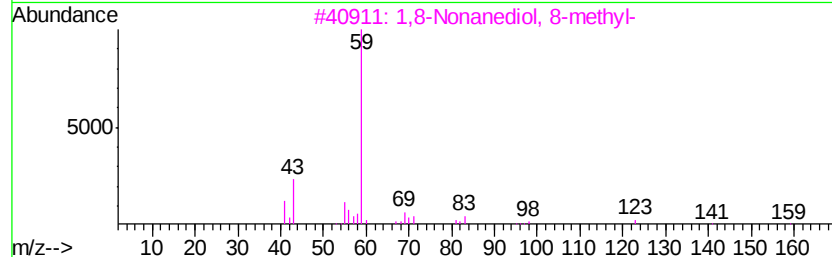
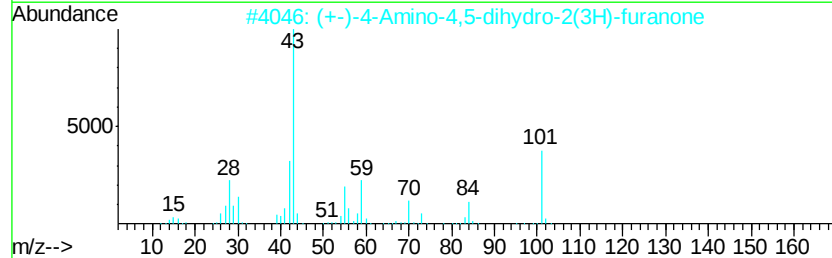
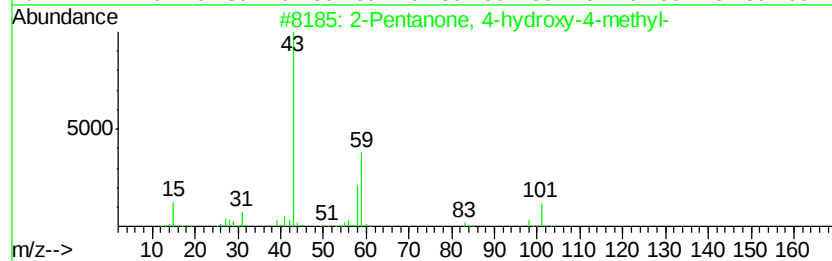
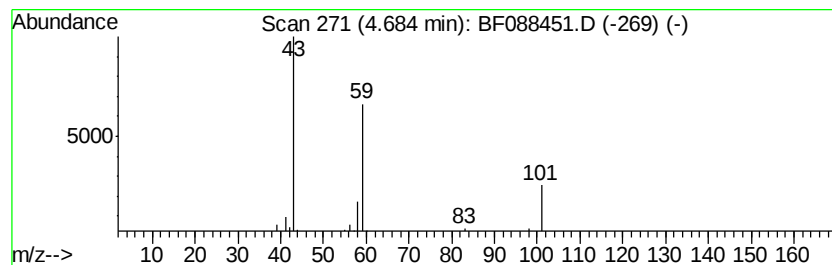
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.68	5.08 ng	80718	1,4-Dichlorobenzene-d4	6.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	28
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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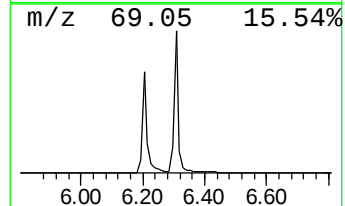
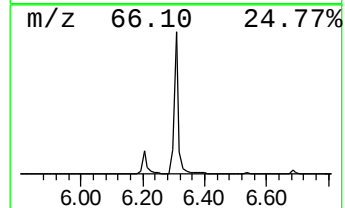
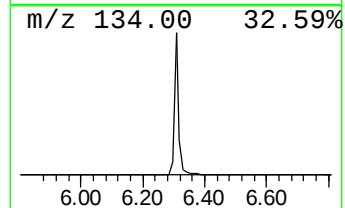
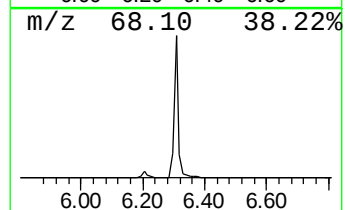
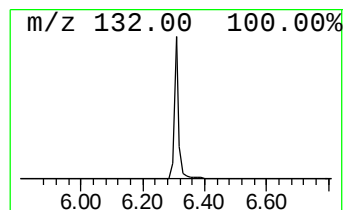
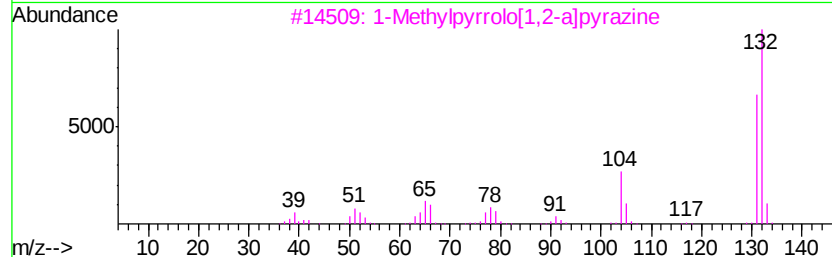
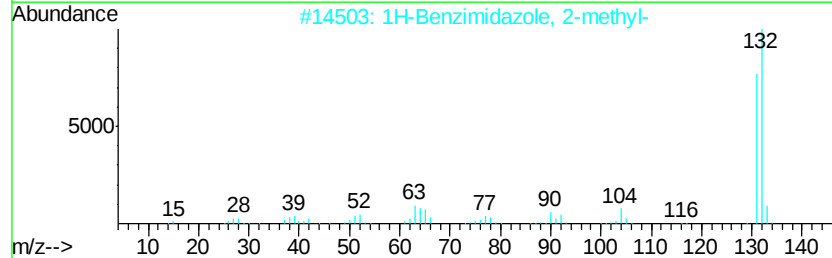
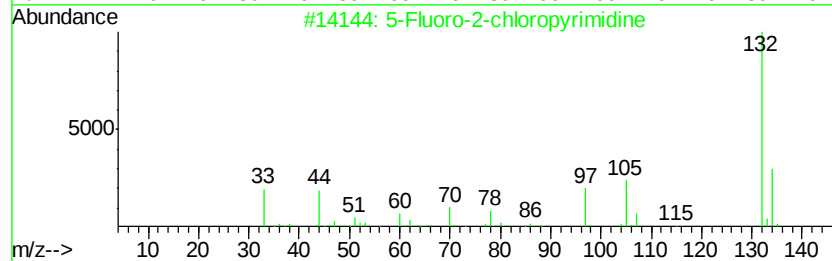
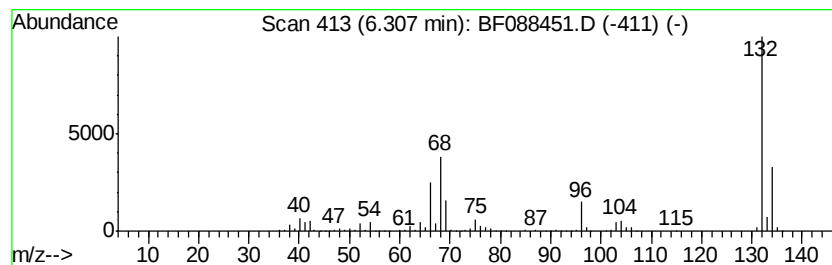
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Peak Number 3 unknown6.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	99.60 ng	1582060	1,4-Dichlorobenzene-d4	6.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	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...	4.68	5.1	ng	80718	1	6.54	317678	20.0
unknown6.31	6.31	99.6	ng	1582060	1	6.54	317678	20.0