

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082615\
 Data File : BF081218.D
 Acq On : 26 Aug 2015 18:12
 Operator : UM/IZ
 Sample : PB85191BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85191BL

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-BF081715.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	4.617	249	252	255	rBV	511012	807069	56.82%	7.146%
2	5.086	290	293	295	rBV	984507	1030737	72.57%	9.127%
3	6.160	385	387	390	rBV	859852	899867	63.35%	7.968%
4	6.275	395	397	400	rBV	855616	889758	62.64%	7.879%
5	6.515	416	418	420	rBV	390144	327650	23.07%	2.901%
6	6.675	429	432	434	rBV	956146	1108312	78.03%	9.814%
7	7.086	465	468	470	rBV	944165	919783	64.75%	8.145%
8	7.795	528	530	533	rBV	296884	400498	28.20%	3.546%
9	8.880	622	625	628	rBV	1495528	1420427	100.00%	12.578%
10	9.543	680	683	685	rBV	519434	456068	32.11%	4.038%
11	10.321	749	751	754	rBV2	553621	562316	39.59%	4.979%
12	11.006	809	811	814	rBV2	376594	427788	30.12%	3.788%
13	12.595	947	950	953	rBV	1422670	1330455	93.67%	11.781%
14	13.624	1038	1040	1042	rBV	435856	405354	28.54%	3.589%
15	14.950	1153	1156	1159	rBV	294480	307199	21.63%	2.720%

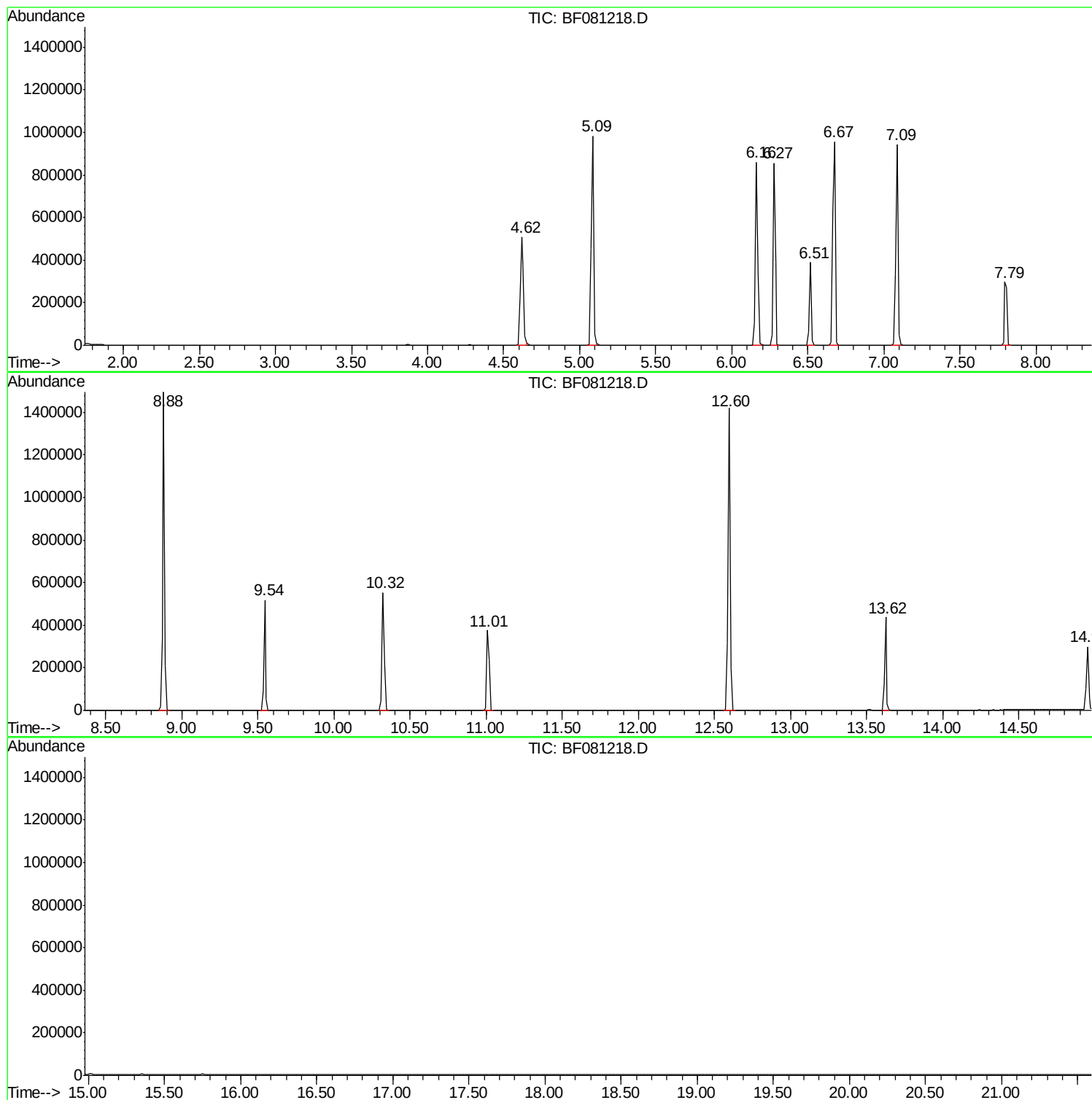
Sum of corrected areas: 11293281

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.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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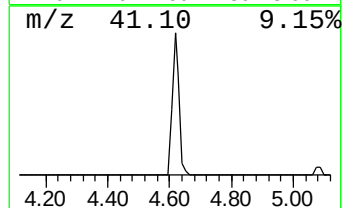
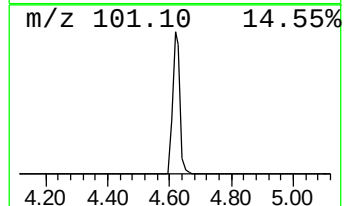
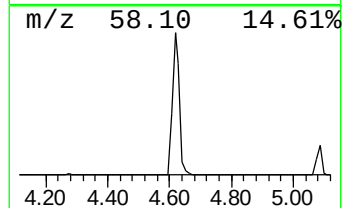
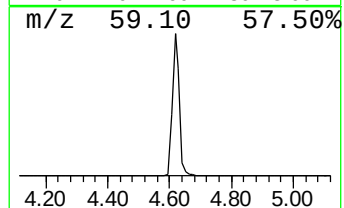
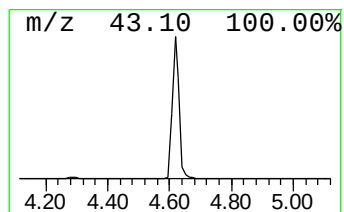
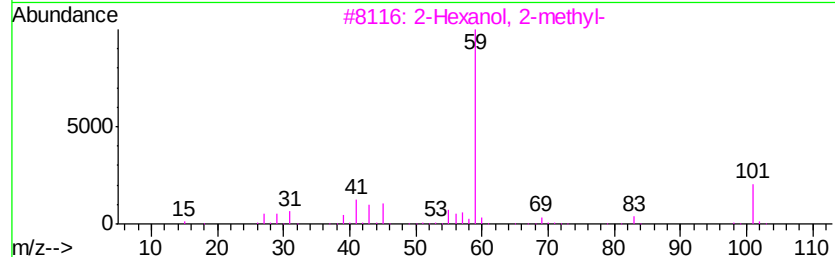
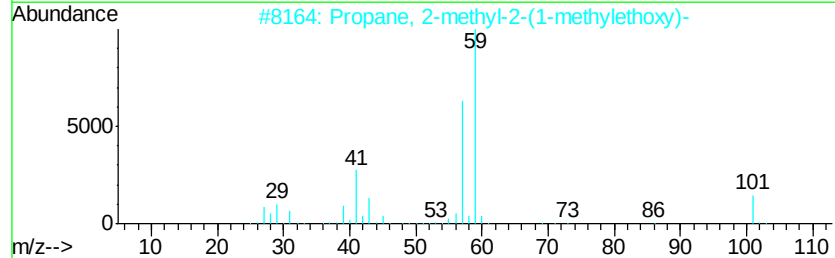
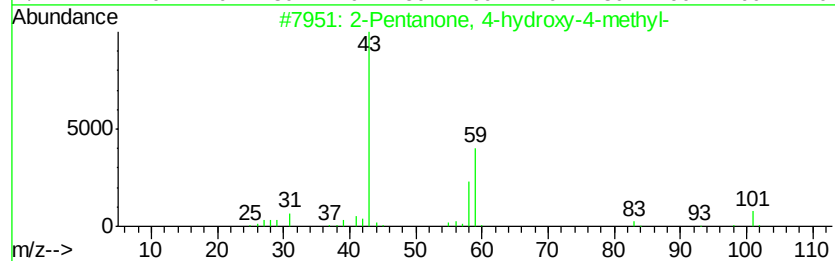
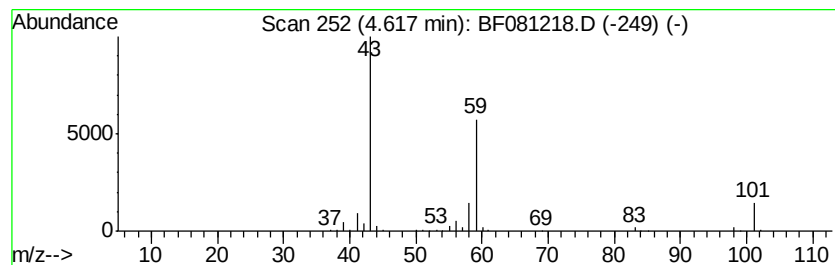
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.62	49.26 ng	807069	1,4-Dichlorobenzene-d4	6.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42		
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33		
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32		



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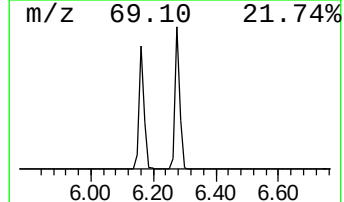
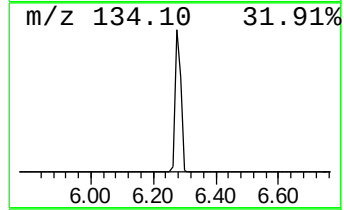
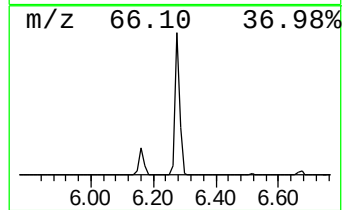
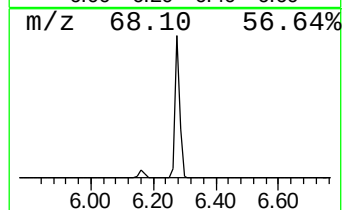
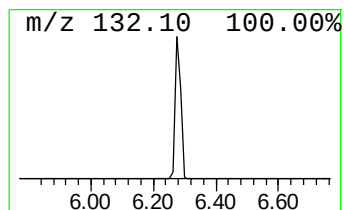
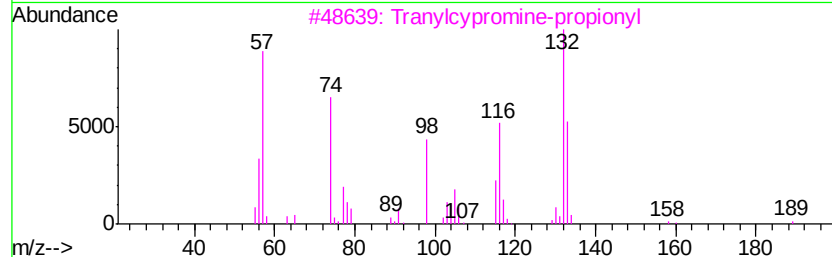
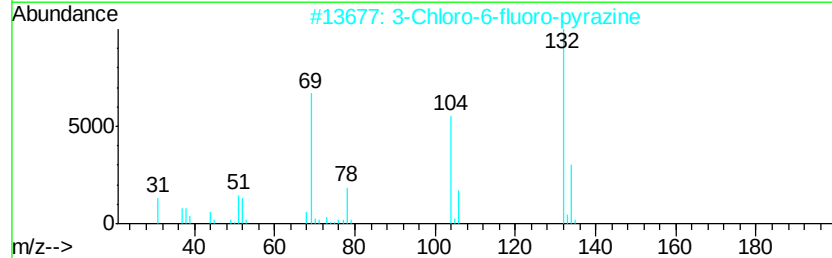
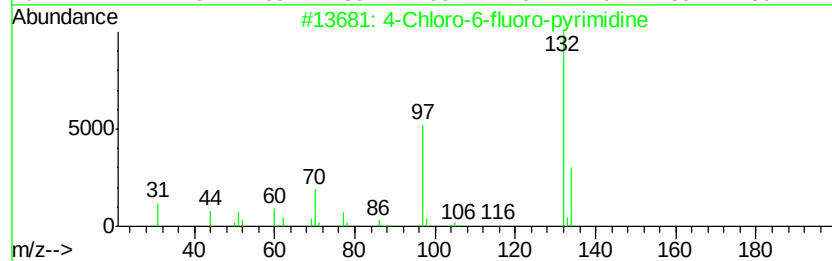
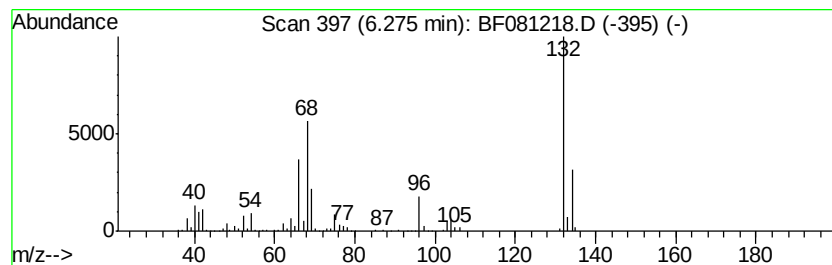
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Peak Number 2 unknown6.27 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.27	54.31 ng	889758	1,4-Dichlorobenzene-d4	6.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27	
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22	
4	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22	
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.62	49.3	ng	807069	1	6.51	327650	20.0
unknown6.27	6.27	54.3	ng	889758	1	6.51	327650	20.0