

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092515\
 Data File : BF081806.D
 Acq On : 25 Sep 2015 16:44
 Operator : UM/IZ
 Sample : PB85698BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85698BL

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-BF092415.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	5.143	260	263	266	rBV	328191	397940	9.26%	1.350%
2	5.543	295	298	300	rBV	2384183	2227077	51.83%	7.553%
3	6.560	384	387	389	rBV	2139461	2061217	47.97%	6.991%
4	6.709	397	400	402	rBV	2540448	2345020	54.58%	7.953%
5	6.949	418	421	423	rBV	835488	928579	21.61%	3.149%
6	7.097	432	434	437	rBV	2470446	2494134	58.05%	8.459%
7	7.509	467	470	472	rBV	2087712	1996292	46.46%	6.771%
8	8.229	531	533	536	rBV	1408354	1252660	29.15%	4.249%
9	9.303	625	627	630	rBV	2860255	4142912	96.42%	14.051%
10	9.989	684	687	689	rBV	1748543	1527926	35.56%	5.182%
11	10.778	753	756	758	rBV	1266694	1464761	34.09%	4.968%
12	11.475	814	817	819	rBV	1877943	1627109	37.87%	5.519%
13	13.064	953	956	958	rBV	4293792	4296793	100.00%	14.573%
14	14.115	1045	1048	1050	rBV	1681653	1577610	36.72%	5.351%
15	15.578	1173	1176	1179	rBV	749159	1144386	26.63%	3.881%

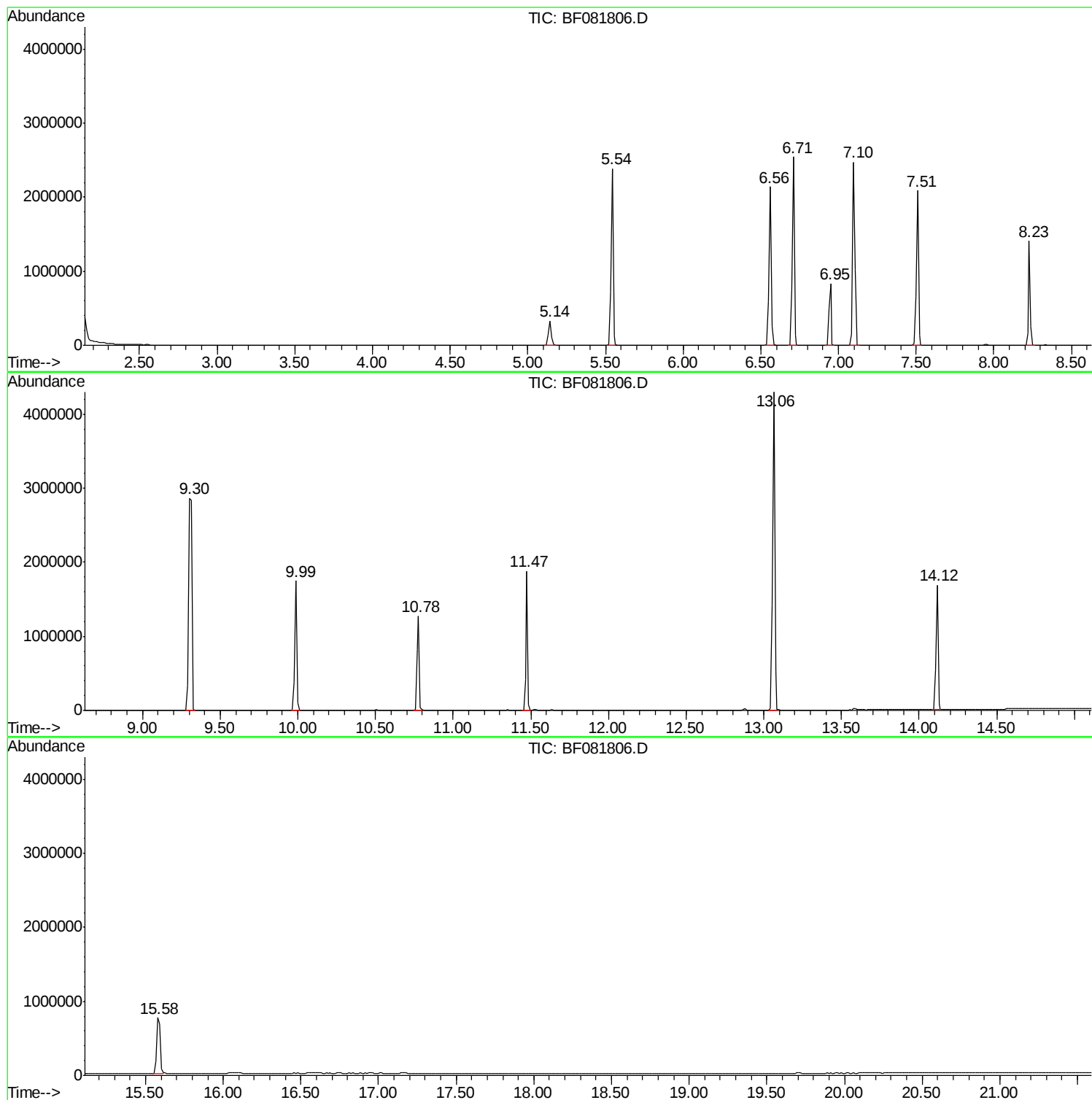
Sum of corrected areas: 29484416

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.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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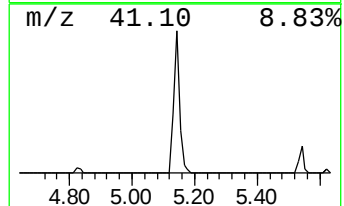
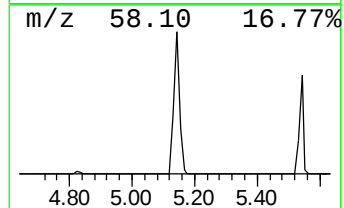
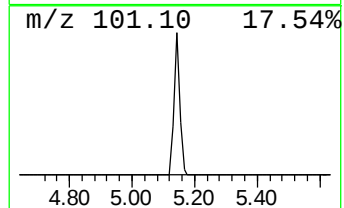
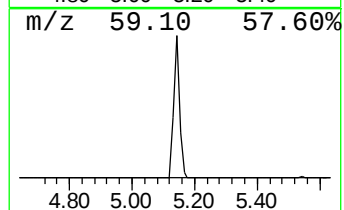
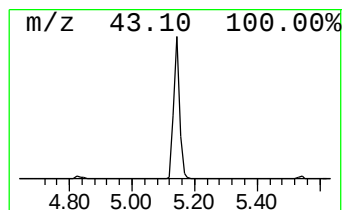
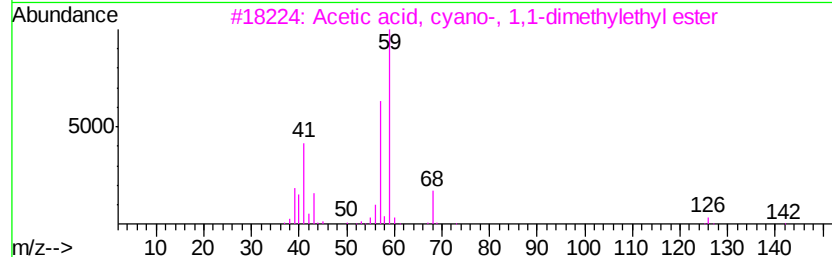
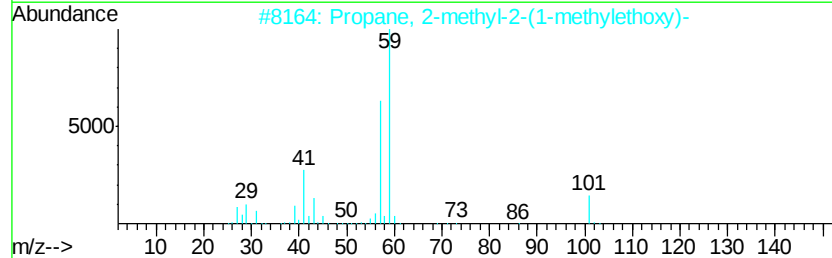
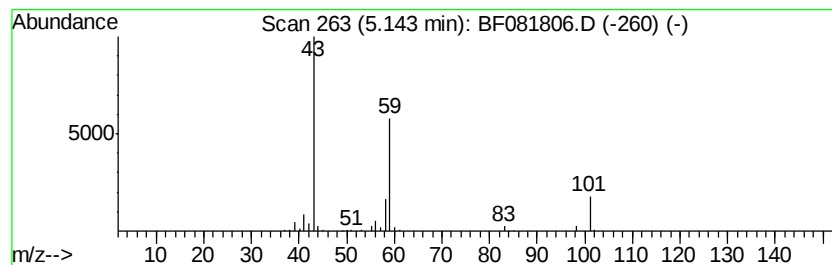
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	8.57 ng	397940	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
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			Acetone	58	C3H6O	000067-64-1	9



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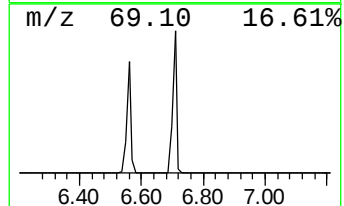
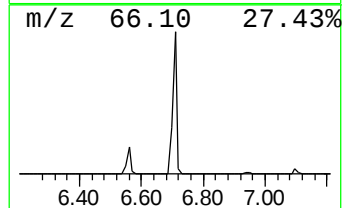
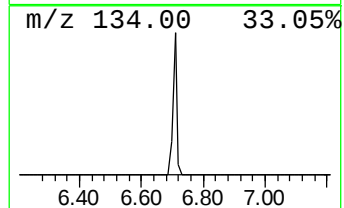
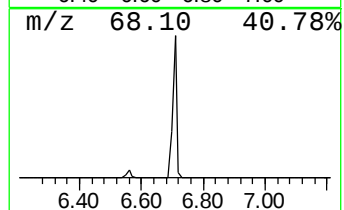
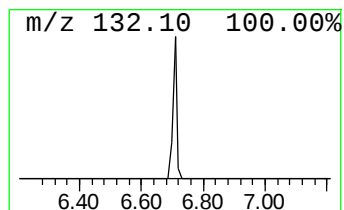
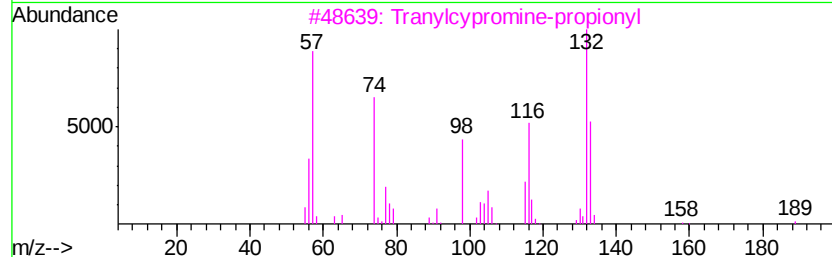
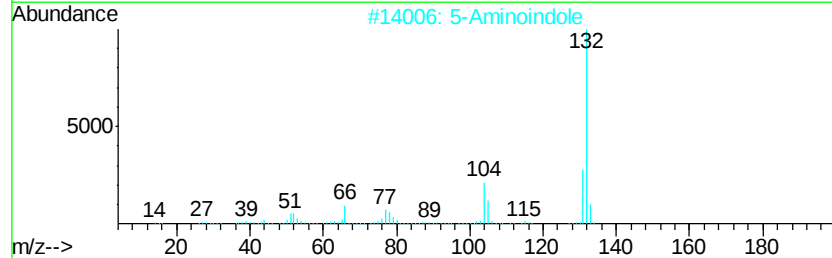
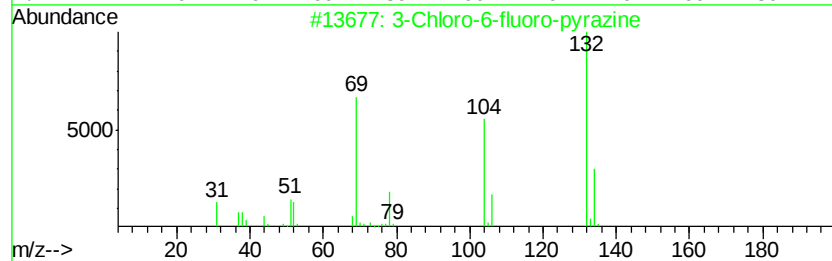
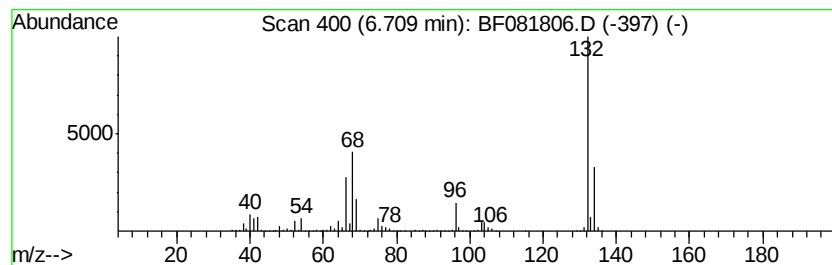
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Peak Number 2 unknown6.71 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	50.51 ng	2345020	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	8.6	ng	397940	1	6.95	928579	20.0
unknown6.71	6.71	50.5	ng	2345020	1	6.95	928579	20.0