

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043016\
 Data File : BF086518.D
 Acq On : 30 Apr 2016 12:37
 Operator : UM/SJ
 Sample : PB90117BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB90117BL

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-BF042716.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	4.955	248	251	257	rVB	56754	91577	1.93%	0.251%
2	5.367	284	287	289	rBV	3640194	4124719	87.11%	11.320%
3	6.407	375	378	380	rBV	3753098	4061832	85.78%	11.147%
4	6.533	387	389	392	rBV	3890609	3795535	80.16%	10.416%
5	6.761	407	409	412	rBV	734852	815389	17.22%	2.238%
6	6.921	420	423	425	rBV	3225329	2990665	63.16%	8.207%
7	7.333	456	459	461	rBV	2335853	2590719	54.71%	7.110%
8	8.053	519	522	524	rBV	1109216	1143251	24.14%	3.137%
9	9.127	613	616	619	rBV	3493514	4087905	86.33%	11.219%
10	9.802	672	675	678	rBV	1375530	1308213	27.63%	3.590%
11	10.590	741	744	761	rBV	2701844	2967138	62.66%	8.143%
12	11.288	802	805	807	rBV	964309	1288418	27.21%	3.536%
13	12.876	941	944	946	rBV	3532314	4735179	100.00%	12.995%
14	13.916	1032	1035	1048	rBV	901306	1385123	29.25%	3.801%
15	15.311	1154	1157	1167	rBV	664509	1053074	22.24%	2.890%

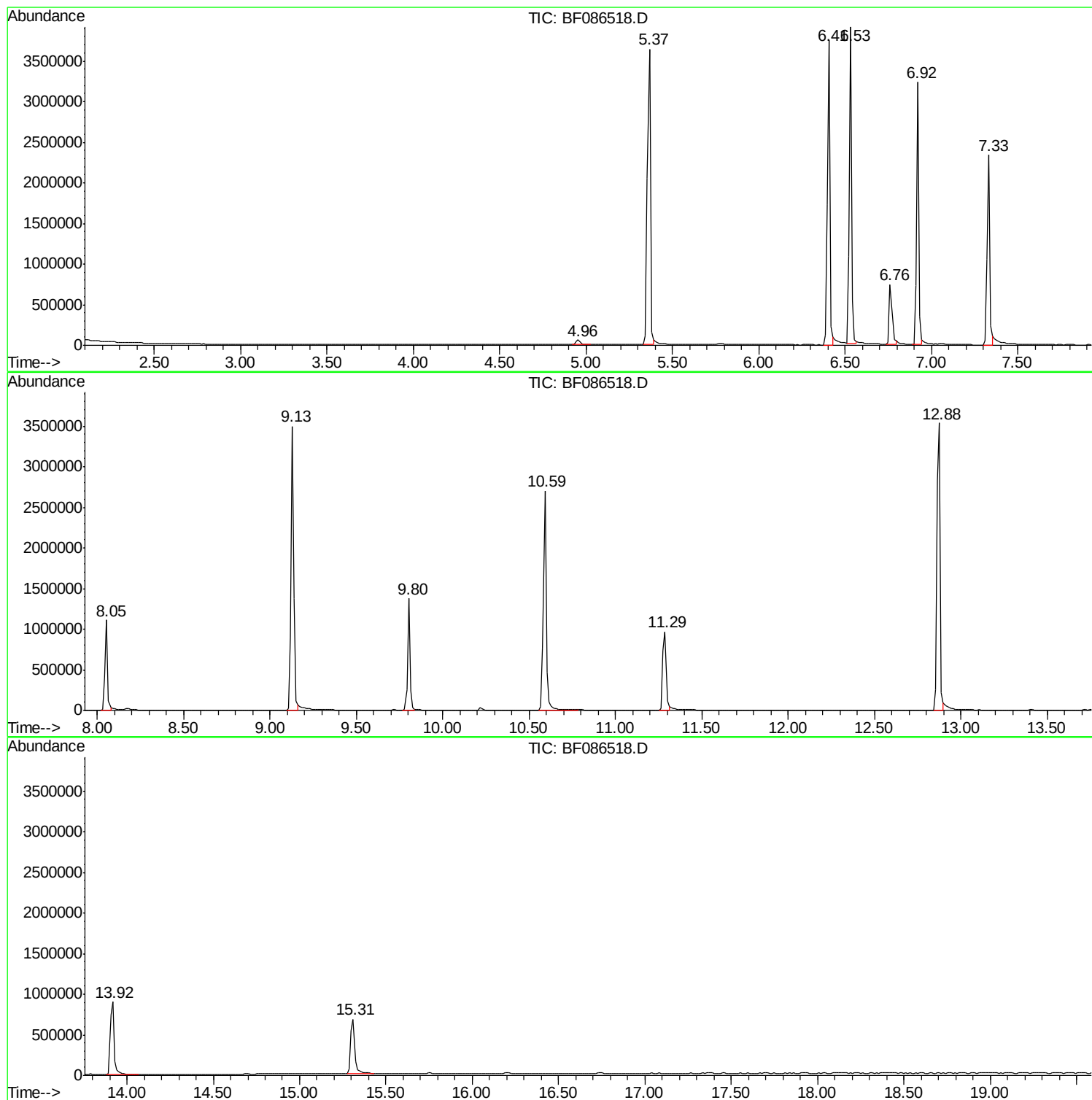
Sum of corrected areas: 36438737

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.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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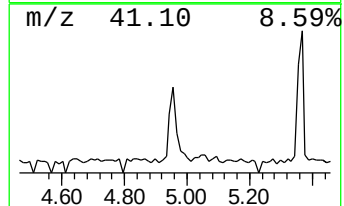
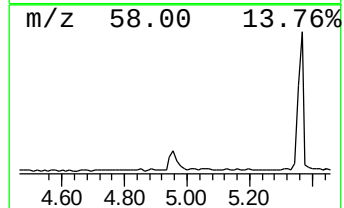
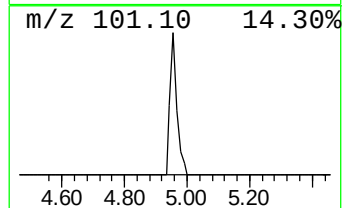
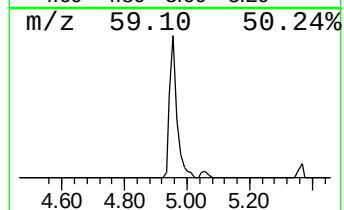
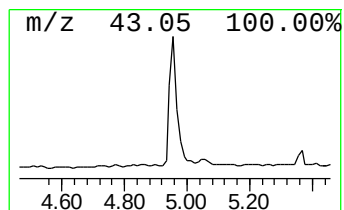
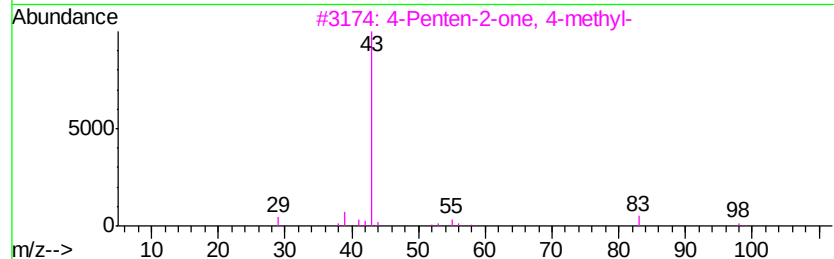
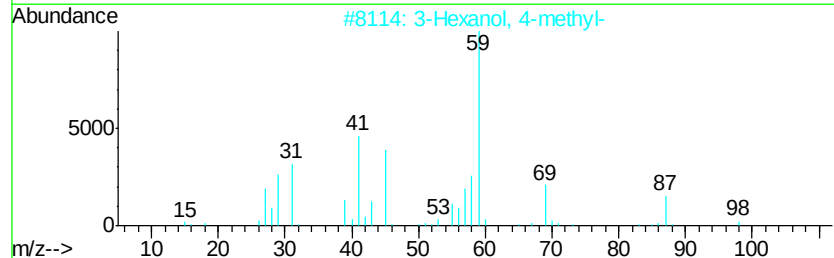
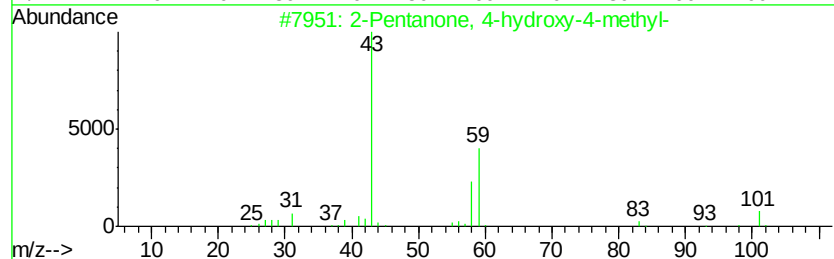
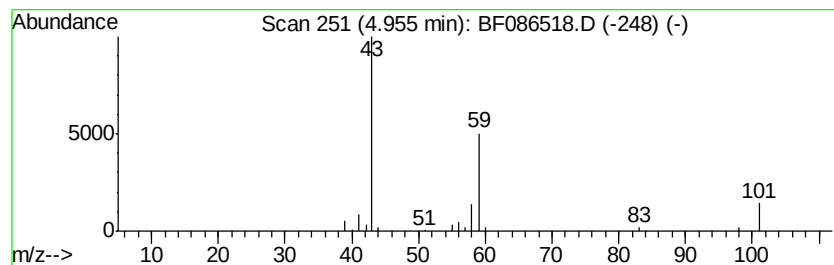
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	2.25 ng	91577	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5			Hexanamide	115	C6H13NO	000628-02-4	9



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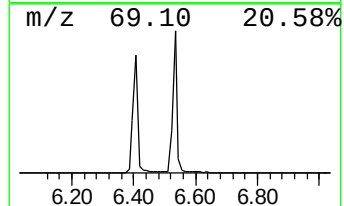
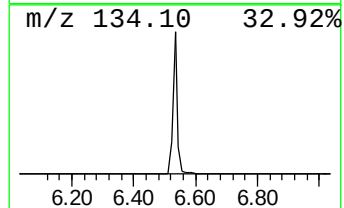
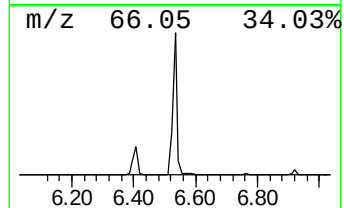
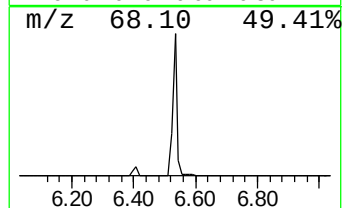
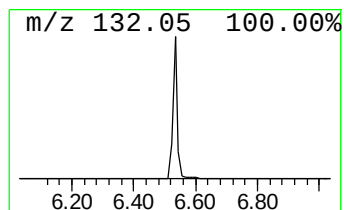
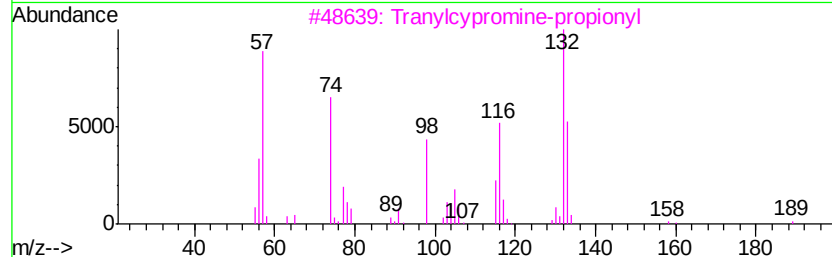
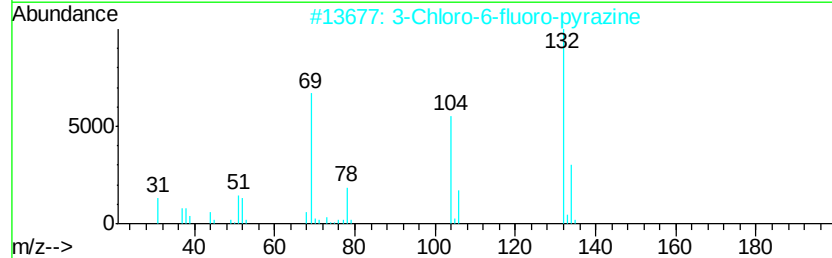
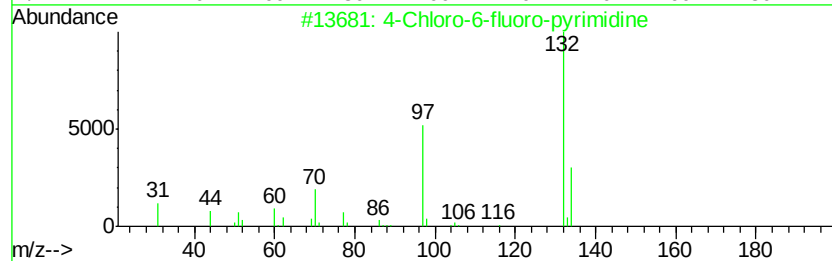
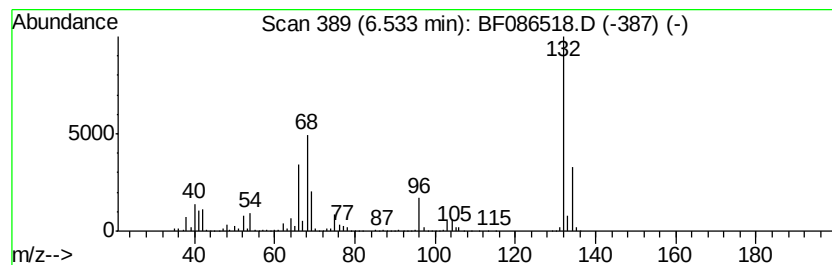
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Peak Number 2 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	93.10 ng	3795540	1,4-Dichlorobenzene-d4	6.76

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	16	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	2.3	ng	91577	1	6.76	815389	20.0
unknown6.53	6.53	93.1	ng	3795540	1	6.76	815389	20.0