

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111715\
 Data File : BF082987.D
 Acq On : 18 Nov 2015 1:27
 Operator : UM/IZ
 Sample : G4407-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-15

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-BF110715.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	4.899	244	246	249	rVB	776475	1029541	15.19%	2.478%
2	5.356	283	286	288	rBV	2597357	2476961	36.55%	5.962%
3	6.396	375	377	379	rBV	1851655	1526037	22.52%	3.673%
4	6.522	385	388	391	rBV	4215129	3876364	57.20%	9.330%
5	6.750	406	408	410	rBV	1307778	1059314	15.63%	2.550%
6	6.910	419	422	424	rBV	4460720	3890938	57.42%	9.365%
7	7.322	455	458	460	rBV	3425718	3814512	56.29%	9.181%
8	8.042	518	521	523	rBV	1335936	1373934	20.27%	3.307%
9	9.116	612	615	618	rBV	4575232	5799151	85.57%	13.958%
10	9.790	671	674	676	rBV	1835474	1601399	23.63%	3.854%
11	10.579	740	743	746	rBV	3690562	3821954	56.40%	9.199%
12	11.265	801	803	806	rBV2	1382289	1610221	23.76%	3.876%
13	12.865	940	943	945	rBV	6159422	6776851	100.00%	16.311%
14	13.905	1031	1034	1036	rBV	1312699	1655982	24.44%	3.986%
15	14.762	1107	1109	1111	rBV	48890	68944	1.02%	0.166%
16	15.300	1153	1156	1158	rBV	837638	1165935	17.20%	2.806%

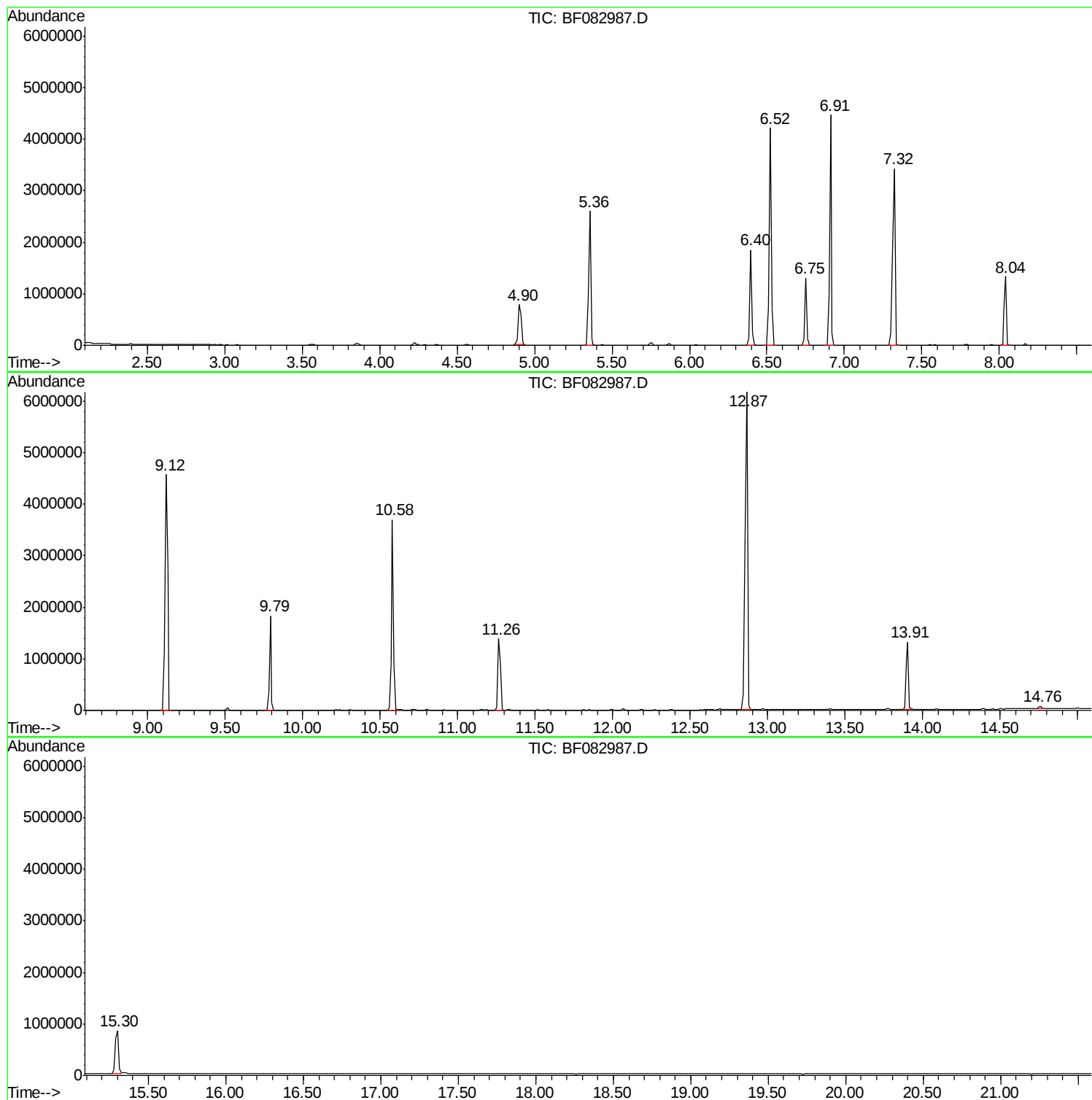
Sum of corrected areas: 41548038

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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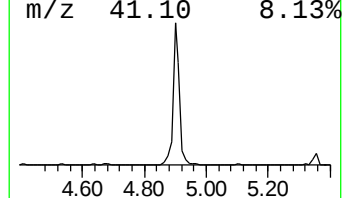
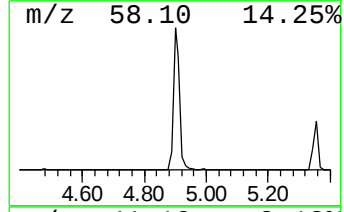
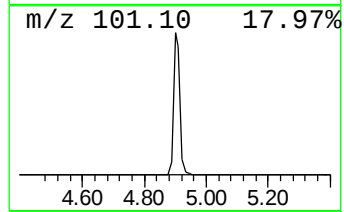
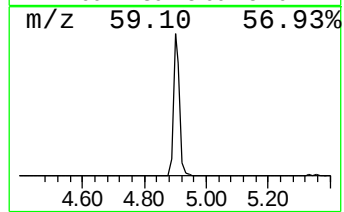
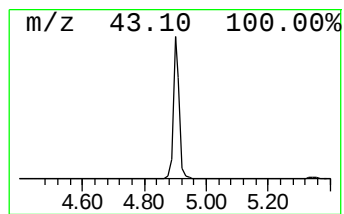
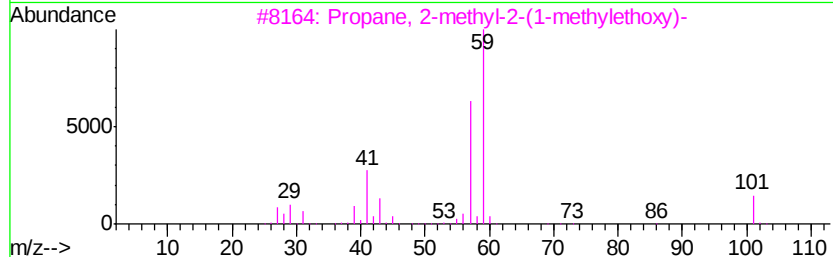
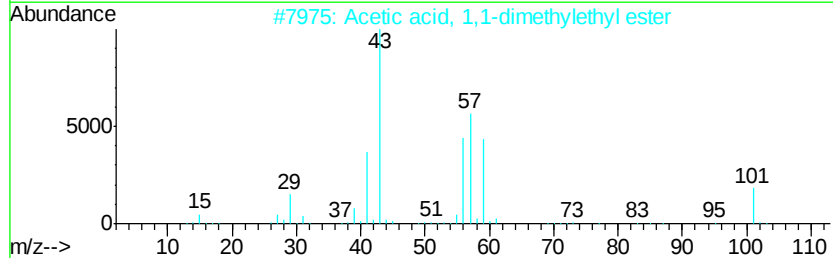
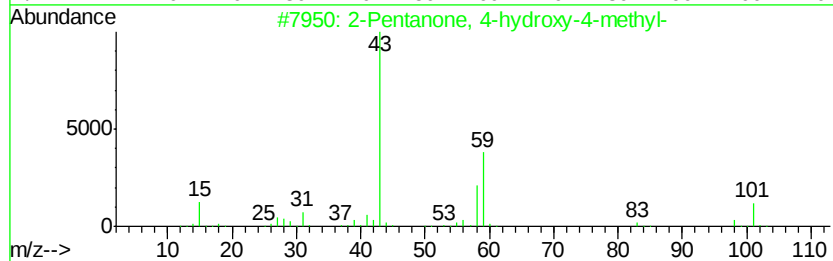
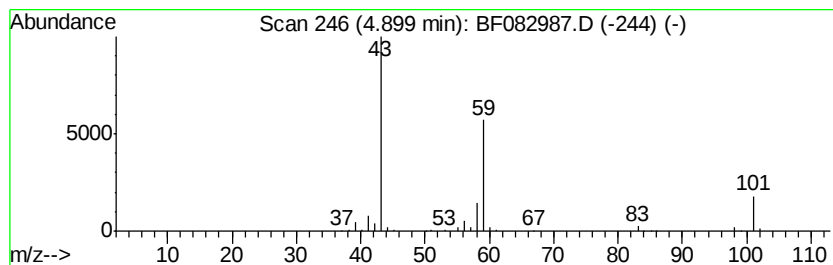
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	19.44 ng	1029540	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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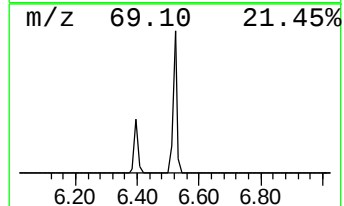
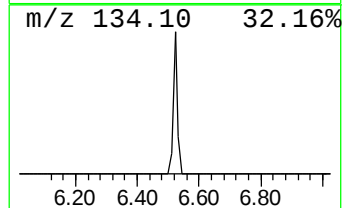
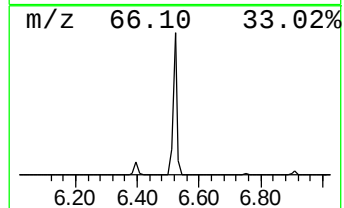
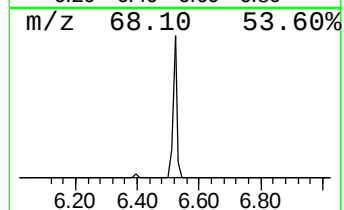
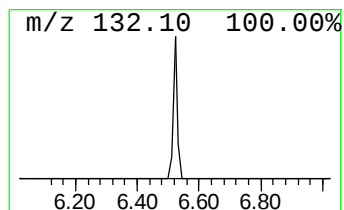
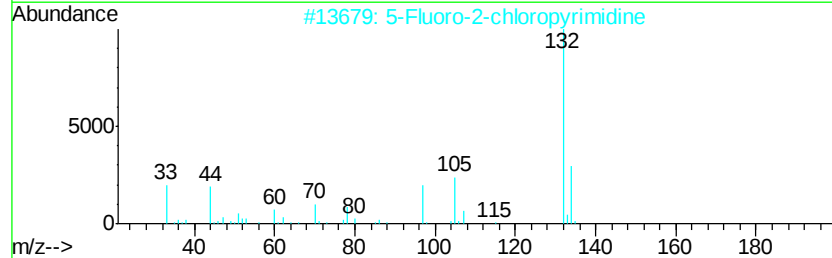
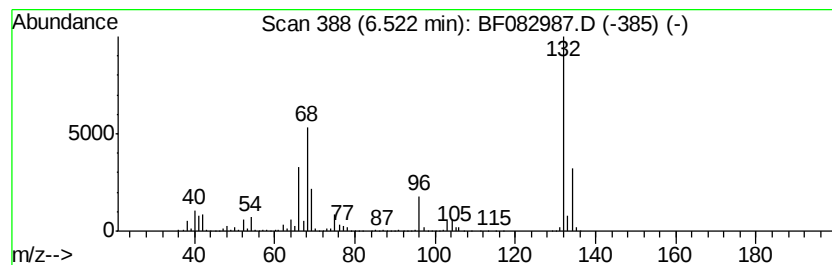
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Peak Number 2 unknown6.52 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	73.19 ng	3876360	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	19.4	ng	1029540	1	6.75	1059310	20.0
unknown6.52	6.52	73.2	ng	3876360	1	6.75	1059310	20.0