

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020316\
 Data File : BF084794.D
 Acq On : 3 Feb 2016 19:42
 Operator : SJ/IZ
 Sample : PB88152BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88152BL

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-BF020116.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	4.841	239	241	254	rVB	706644	1019878	22.58%	2.735%
2	5.287	277	280	282	rBV	3373267	3897831	86.31%	10.454%
3	6.339	369	372	374	rBV	3559873	4009526	88.78%	10.754%
4	6.453	379	382	384	rBV	3993920	3574241	79.14%	9.586%
5	6.682	400	402	405	rVB	1151386	1029055	22.79%	2.760%
6	6.842	413	416	418	rBV	3461203	3050739	67.55%	8.182%
7	7.253	449	452	454	rBV	2408587	2296324	50.85%	6.159%
8	7.973	512	515	517	rBV	1488825	1425223	31.56%	3.823%
9	9.047	606	609	612	rBV	2940121	4252785	94.17%	11.406%
10	9.722	665	668	671	rVB	1698844	1545777	34.23%	4.146%
11	10.511	734	737	739	rBV	2861356	2734029	60.54%	7.333%
12	11.196	795	797	800	rBV	1291923	1512913	33.50%	4.058%
13	12.796	934	937	939	rBV	3263170	4516132	100.00%	12.113%
14	13.825	1025	1027	1030	rBV	1024357	1430891	31.68%	3.838%
15	15.208	1145	1148	1151	rBV	873057	989045	21.90%	2.653%

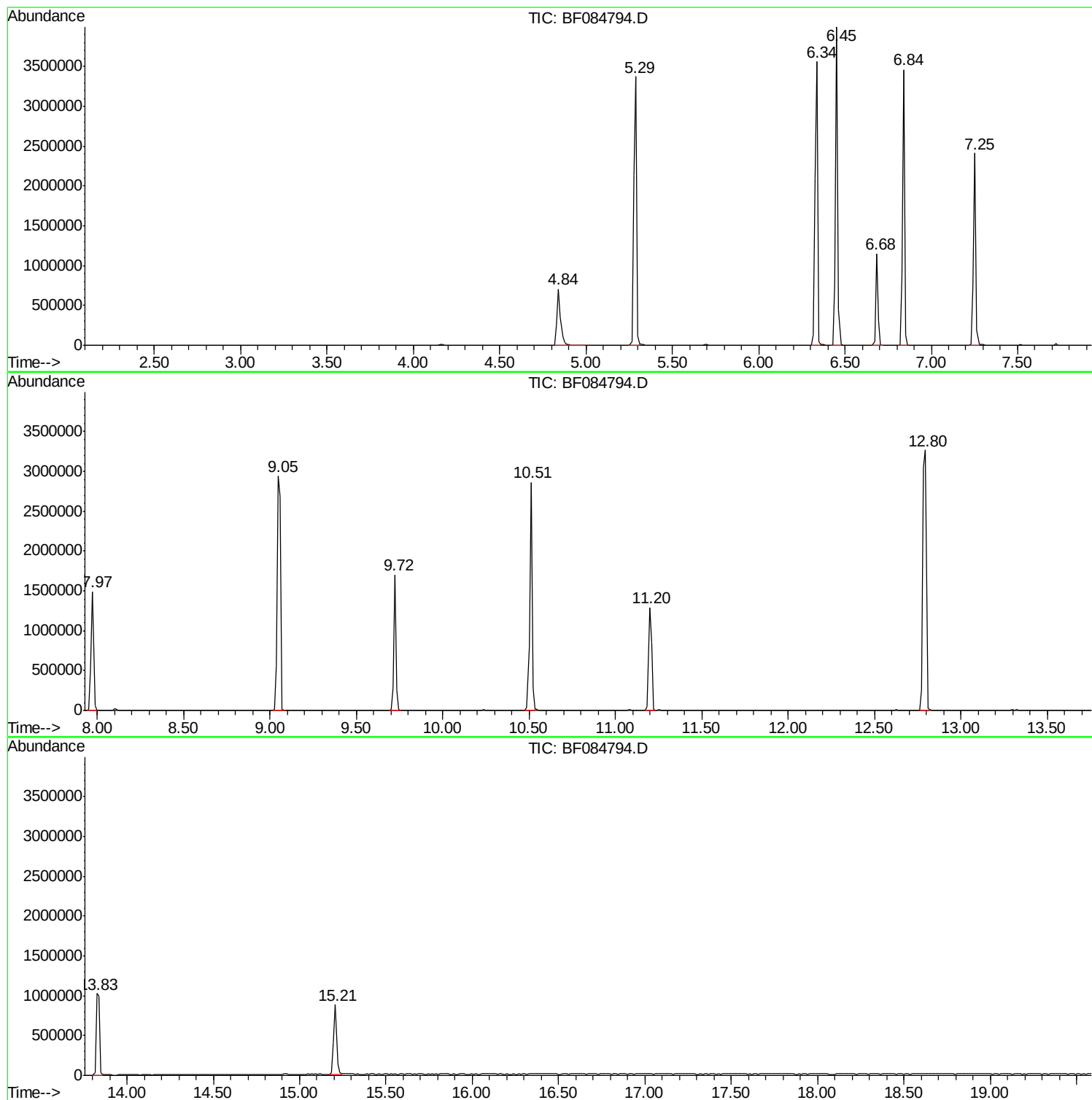
Sum of corrected areas: 37284389

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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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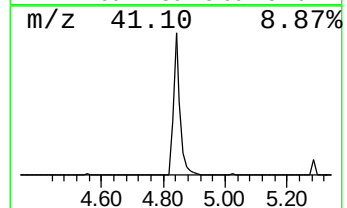
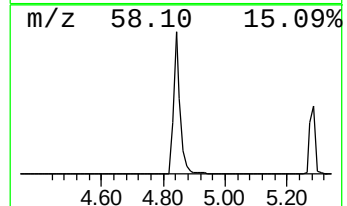
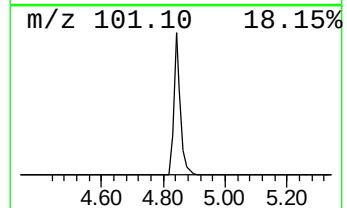
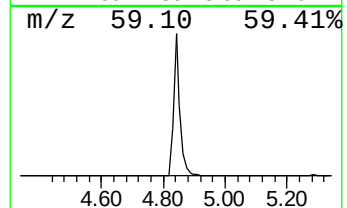
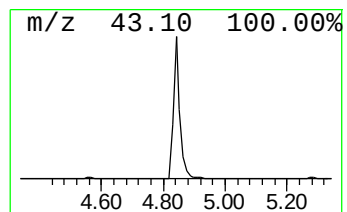
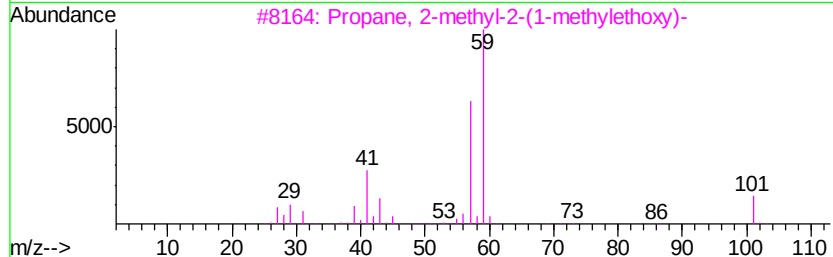
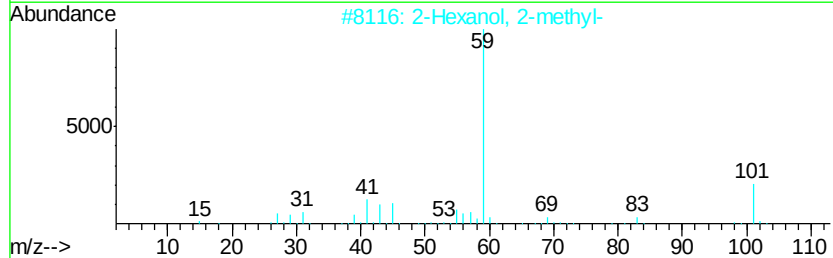
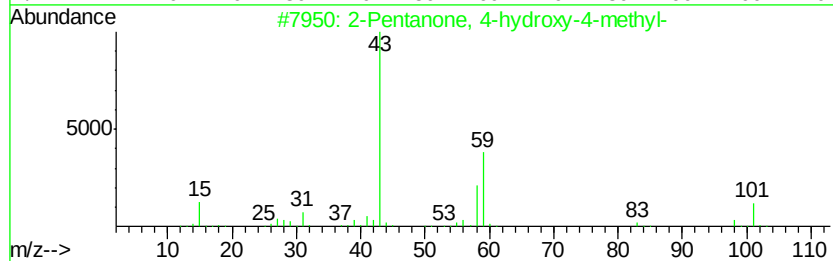
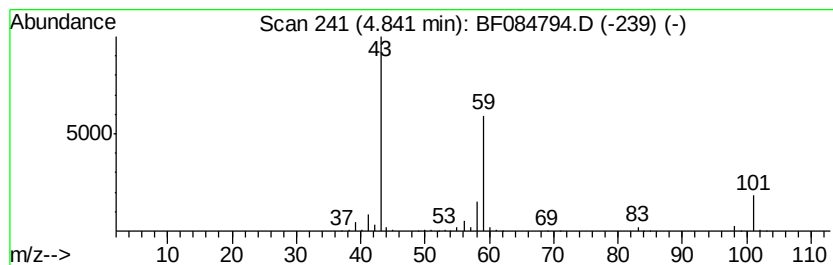
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	19.82 ng	1019880	1,4-Dichlorobenzene-d4	6.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25



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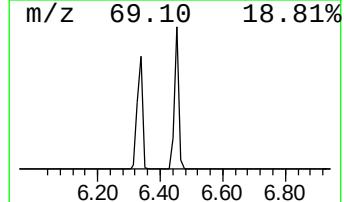
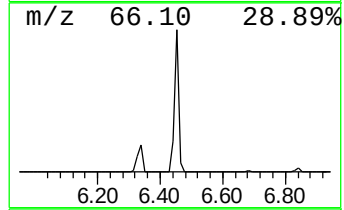
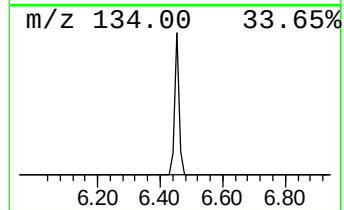
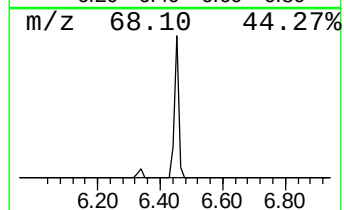
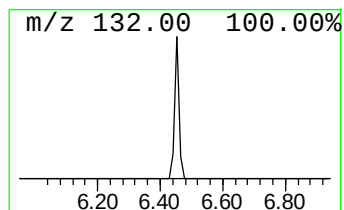
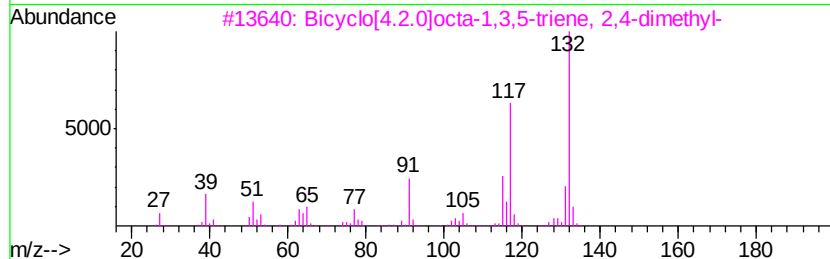
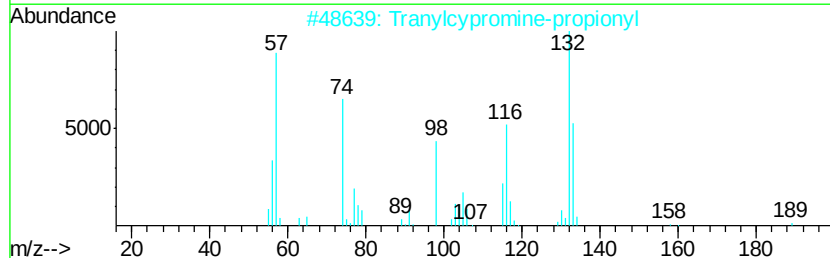
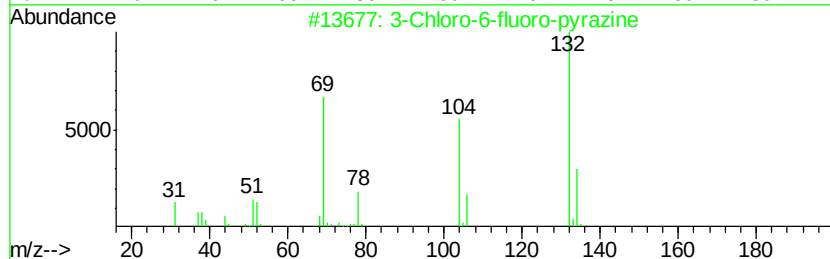
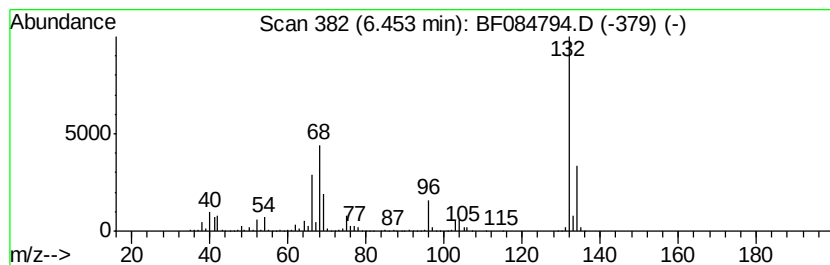
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Peak Number 2 unknown6.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	69.47 ng	3574240	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene, ...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.84	19.8	ng	1019880	1	6.68	1029060	20.0
unknown6.45	6.45	69.5	ng	3574240	1	6.68	1029060	20.0