

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090315\
 Data File : BF081384.D
 Acq On : 3 Sep 2015 19:12
 Operator : UM/IZ
 Sample : PB85376BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85376BL

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-BF090115.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	4.422	244	248	259	rBV	602967	903813	50.83%	6.764%
2	4.924	289	292	303	rVB	1097157	1098463	61.78%	8.221%
3	6.033	387	389	392	rBV	1042199	1011889	56.91%	7.573%
4	6.136	396	398	400	rBV	1296279	1119122	62.94%	8.376%
5	6.365	416	418	421	rBV	438962	384835	21.64%	2.880%
6	6.525	429	432	434	rBV	1420665	1281966	72.10%	9.595%
7	6.947	466	469	471	rBV	962383	1109002	62.37%	8.300%
8	7.656	528	531	533	rBV	560311	500469	28.15%	3.746%
9	8.742	623	626	628	rBV	1844839	1778161	100.00%	13.308%
10	9.393	681	683	686	rVB	416631	513300	28.87%	3.842%
11	10.182	749	752	754	rBV	730245	692541	38.95%	5.183%
12	10.868	809	812	814	rBV	450244	553220	31.11%	4.140%
13	12.445	948	950	953	rBV	1371888	1642451	92.37%	12.293%
14	13.474	1037	1040	1042	rBV	401916	481040	27.05%	3.600%
15	14.777	1152	1154	1157	rBV	292124	291130	16.37%	2.179%

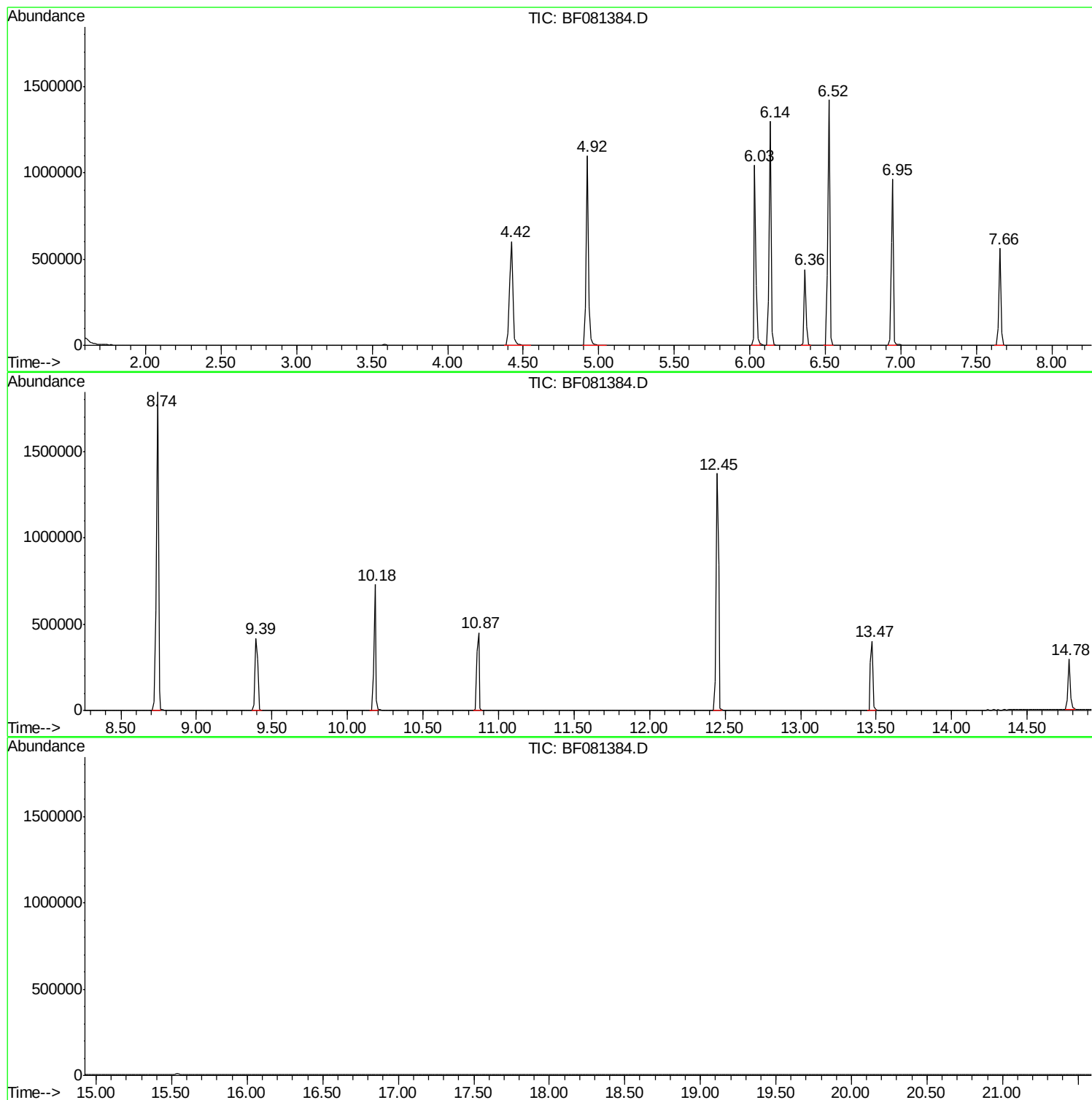
Sum of corrected areas: 13361402

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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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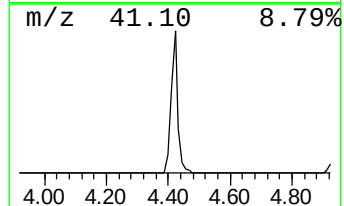
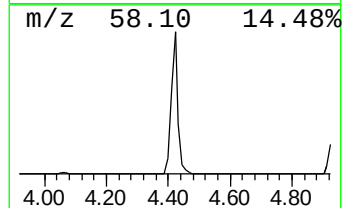
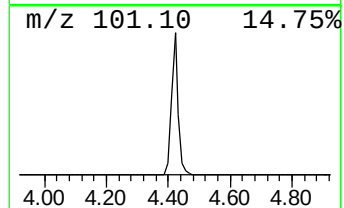
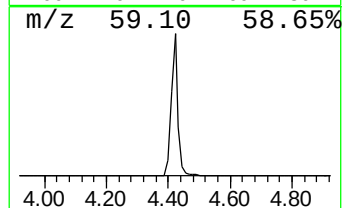
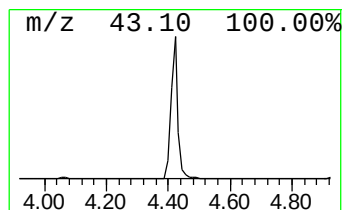
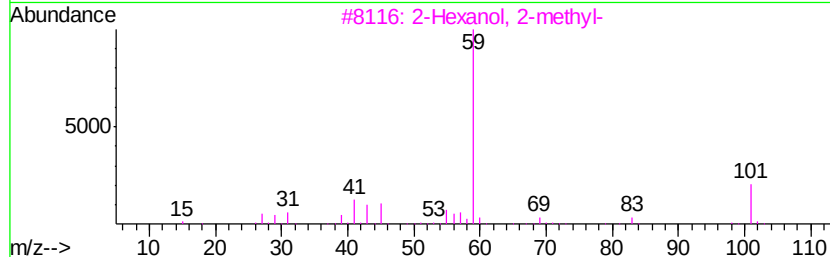
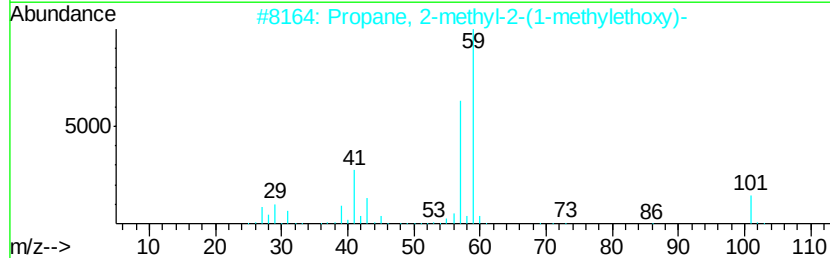
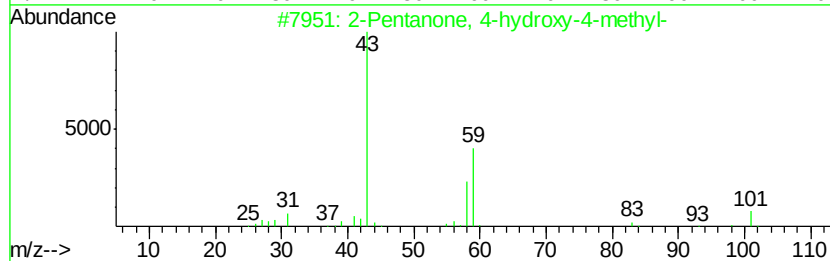
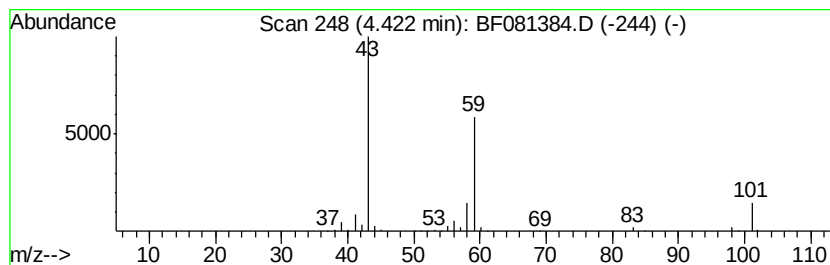
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.42	46.97 ng	903813	1,4-Dichlorobenzene-d4	6.36

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	36
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	23



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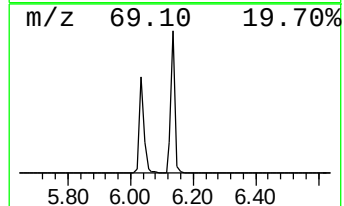
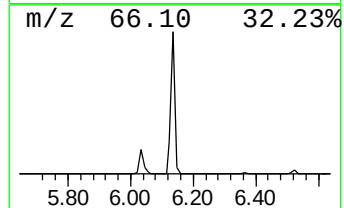
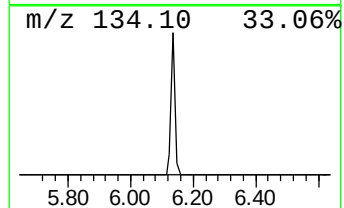
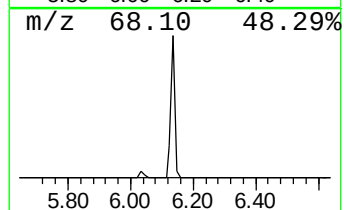
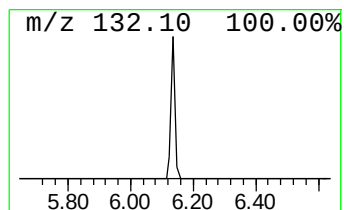
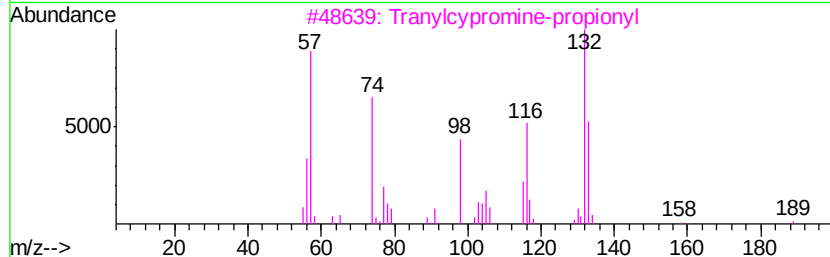
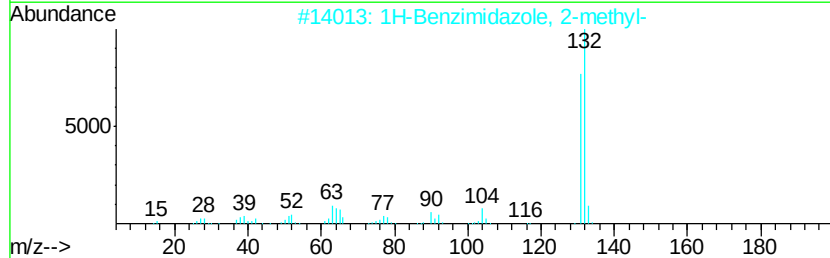
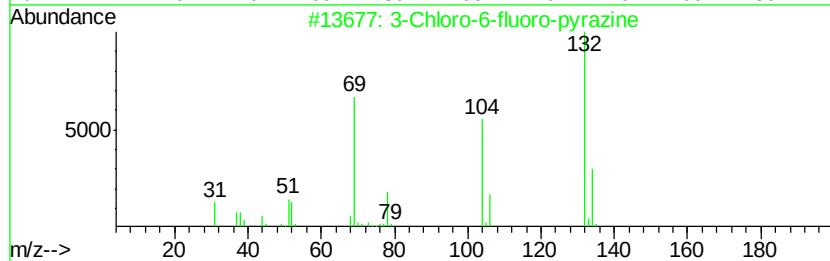
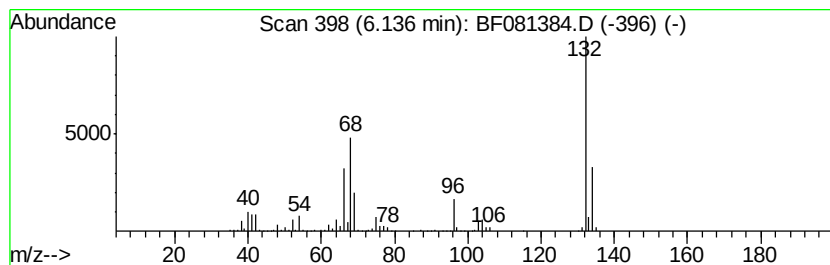
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Peak Number 2 unknown6.14 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	58.16 ng	1119120	1,4-Dichlorobenzene-d4	6.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	25
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.42	47.0	ng	903813	1	6.36	384835	20.0
unknown6.14	6.14	58.2	ng	1119120	1	6.36	384835	20.0