

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121615\
 Data File : BF083849.D
 Acq On : 16 Dec 2015 15:41
 Operator : UM/SJ
 Sample : PB87334BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87334BL

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-BF121315.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	5.081	259	262	270	rVB	309070	496157	17.44%	2.448%
2	5.504	296	299	302	rBV	1827049	1836310	64.54%	9.059%
3	6.522	385	388	391	rBV	1617105	1767077	62.11%	8.718%
4	6.659	397	400	403	rBV	2095615	1969904	69.24%	9.718%
5	6.887	418	420	423	rBV	603365	515333	18.11%	2.542%
6	7.047	431	434	436	rBV	1974903	1693689	59.53%	8.356%
7	7.447	466	469	472	rBV	1153034	1222161	42.96%	6.029%
8	8.167	530	532	535	rBV	524464	667035	23.44%	3.291%
9	9.253	624	627	629	rBV	2628073	2845209	100.00%	14.036%
10	9.928	683	686	688	rBV	878222	824805	28.99%	4.069%
11	10.716	752	755	757	rBV	1536958	1626488	57.17%	8.024%
12	11.413	813	816	818	rBV	606664	819701	28.81%	4.044%
13	12.991	951	954	957	rBV	2370257	2671003	93.88%	13.177%
14	14.042	1043	1046	1048	rBV	834607	784924	27.59%	3.872%
15	15.482	1169	1172	1175	rBV	388870	530380	18.64%	2.617%

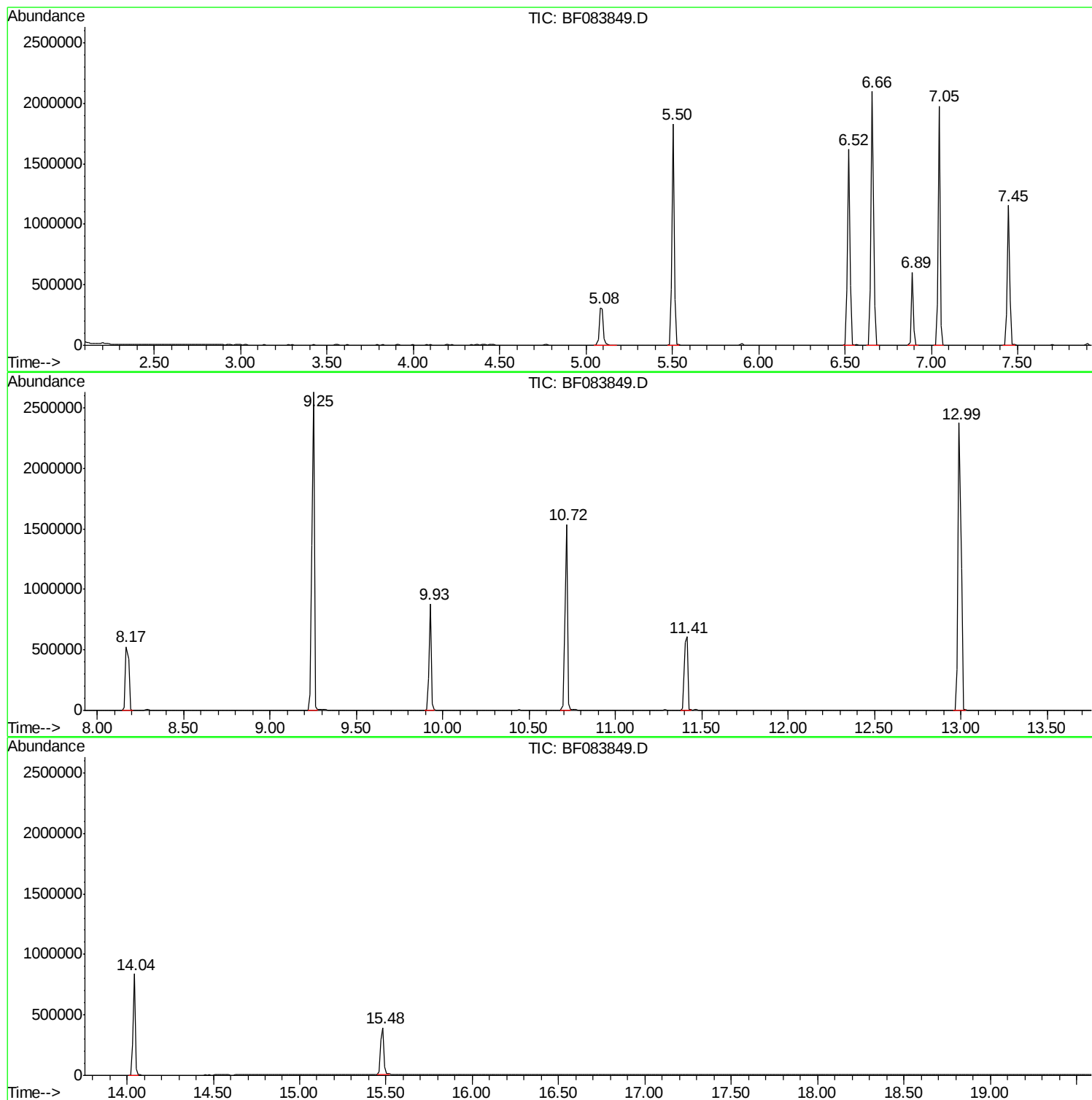
Sum of corrected areas: 20270176

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.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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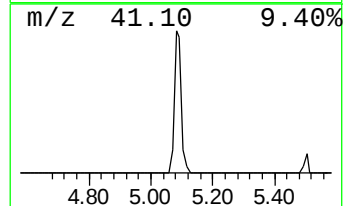
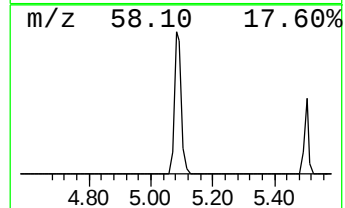
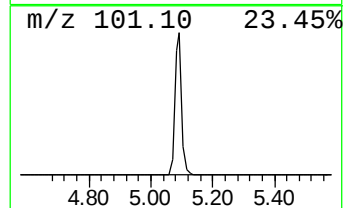
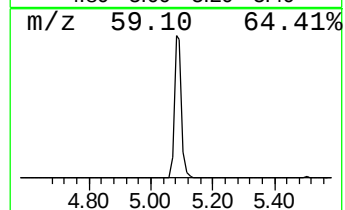
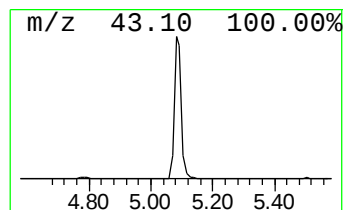
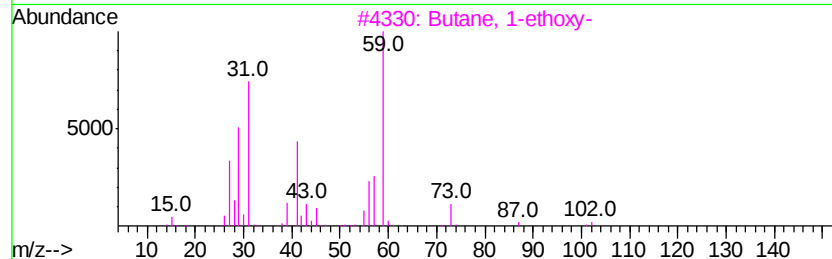
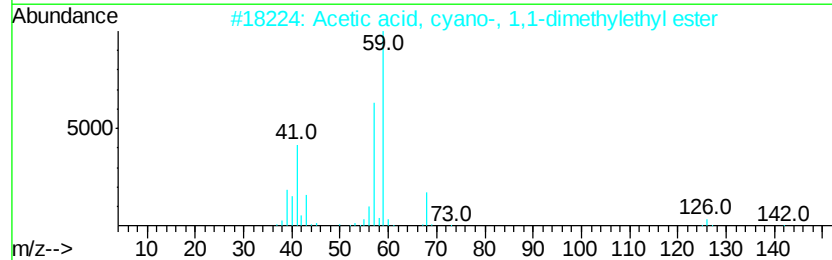
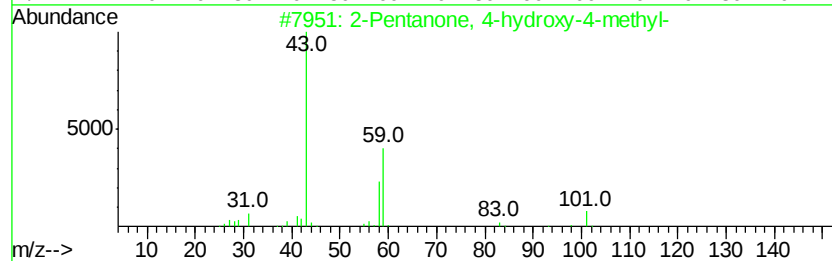
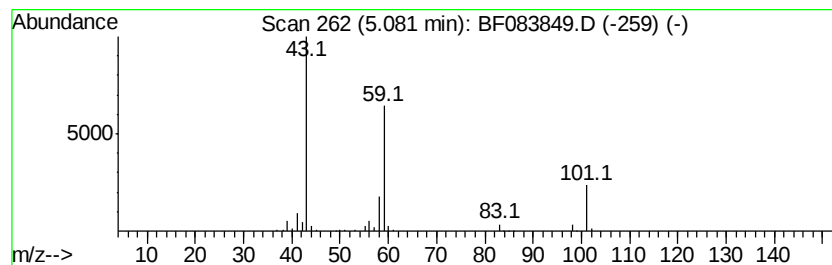
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	19.26 ng	496157	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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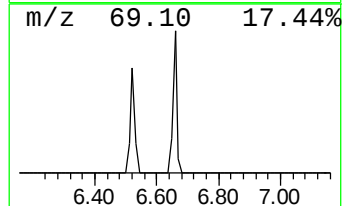
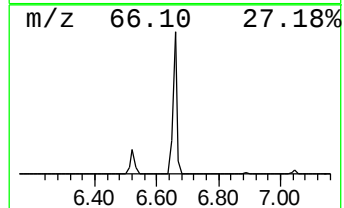
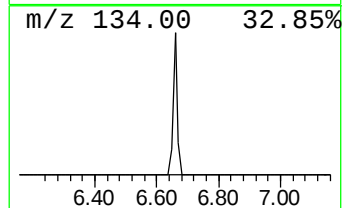
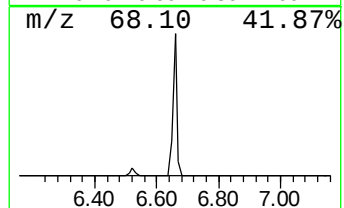
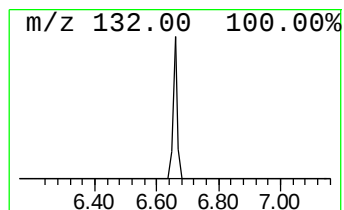
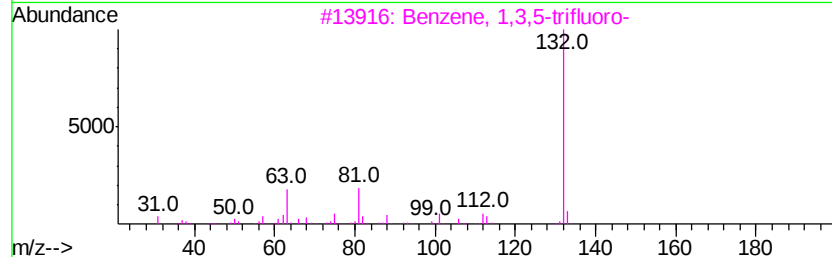
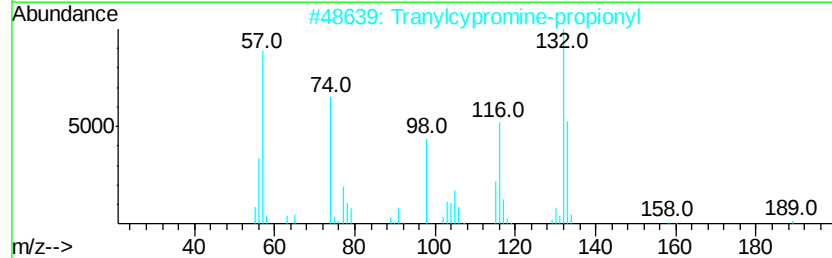
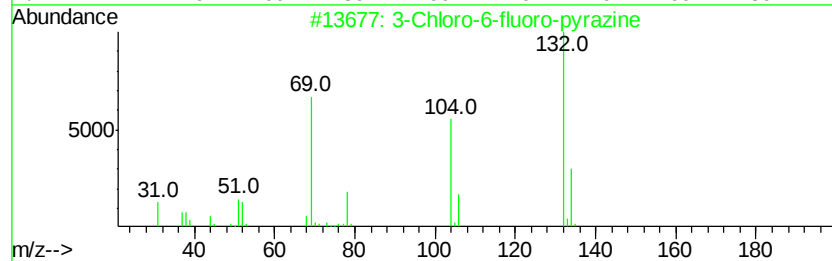
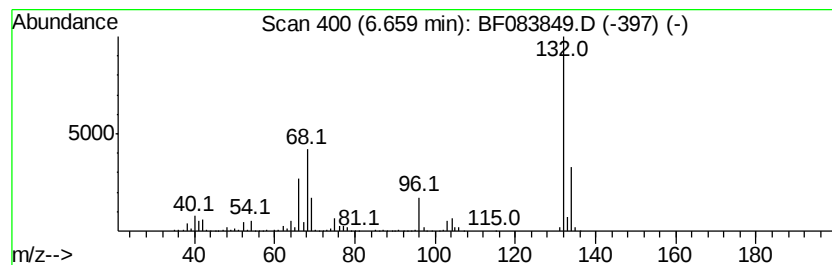
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	76.45 ng	1969900	1,4-Dichlorobenzene-d4	6.89

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		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.08	19.3	ng	496157	1	6.89	515333	20.0
unknown6.66	6.66	76.5	ng	1969900	1	6.89	515333	20.0