

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121815\
 Data File : BF083890.D
 Acq On : 18 Dec 2015 2:00
 Operator : UM/SJ
 Sample : PB87315BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87315BL

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-BF121815.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	5.070	259	261	270	rBB	228990	299621	14.45%	2.163%
2	5.493	295	298	300	rBV	1301818	1385631	66.81%	10.003%
3	6.510	385	387	389	rBV	1221953	1197295	57.73%	8.643%
4	6.647	396	399	401	rBV	1374069	1475238	71.13%	10.650%
5	6.876	417	419	421	rBV	351562	320020	15.43%	2.310%
6	7.036	430	433	435	rBV	876169	1150135	55.45%	8.303%
7	7.436	466	468	470	rBV	932170	863060	41.61%	6.230%
8	8.156	529	531	533	rBV	536608	443093	21.36%	3.199%
9	9.230	623	625	628	rBV	1602600	1663752	80.22%	12.011%
10	9.905	682	684	687	rBB	453321	483601	23.32%	3.491%
11	10.694	751	753	756	rBV	1048550	1035524	49.93%	7.475%
12	11.391	812	814	817	rBV	636768	529480	25.53%	3.822%
13	12.979	950	953	955	rBV	2133218	2074096	100.00%	14.973%
14	14.031	1043	1045	1047	rBV	479286	520367	25.09%	3.757%
15	15.471	1168	1171	1180	rBV	264478	411332	19.83%	2.969%

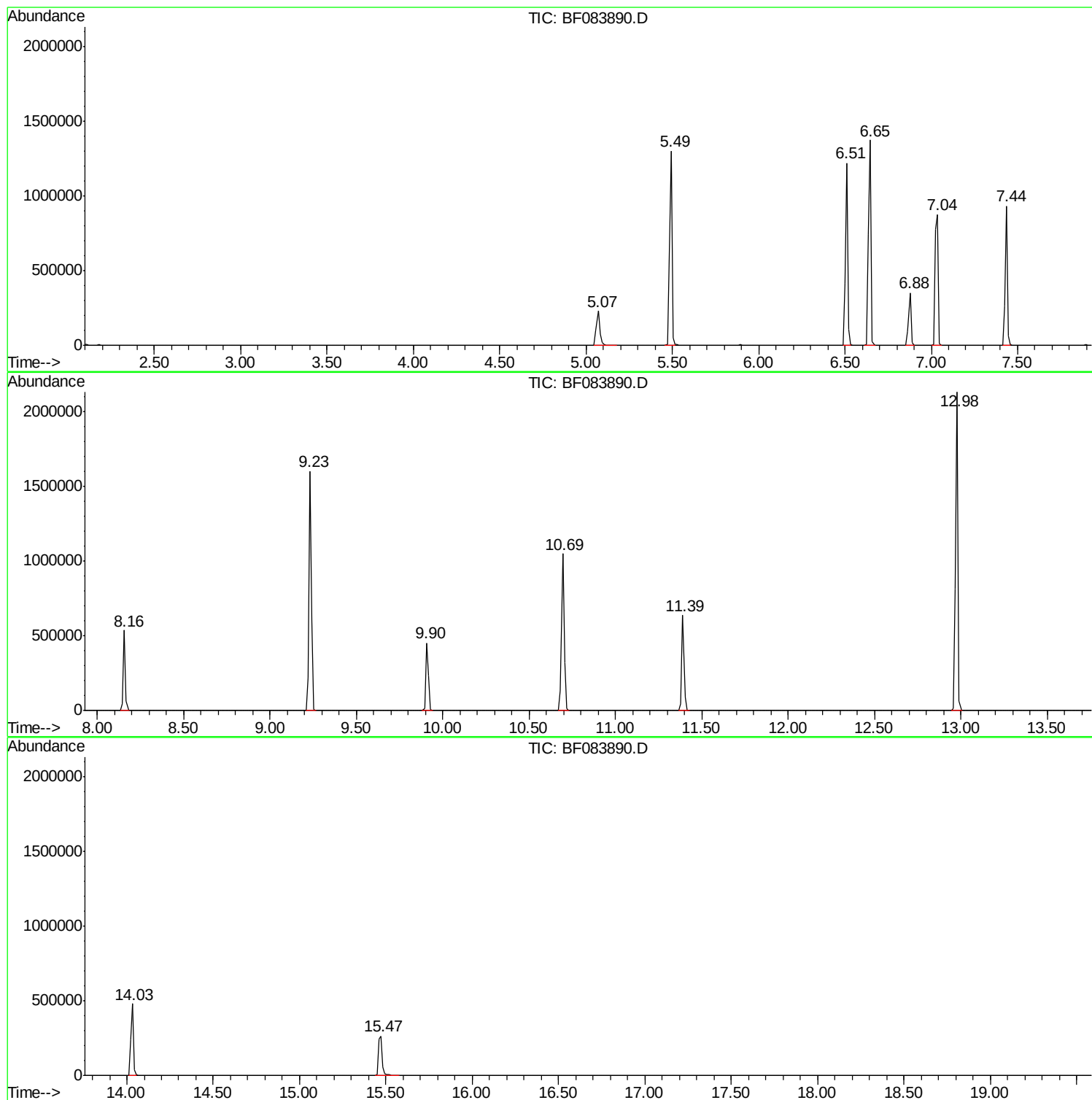
Sum of corrected areas: 13852245

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.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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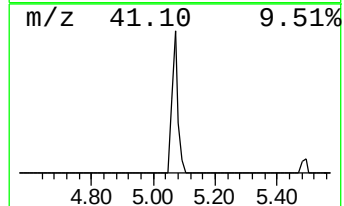
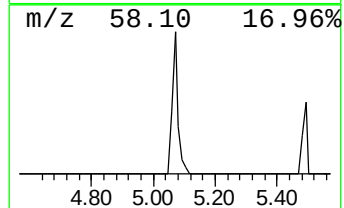
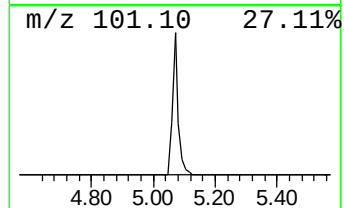
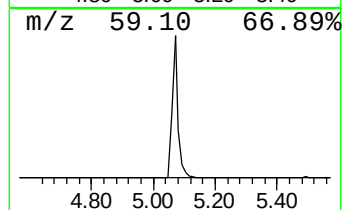
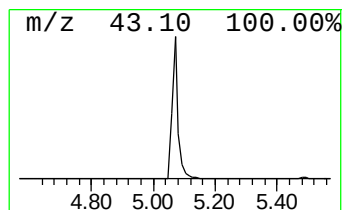
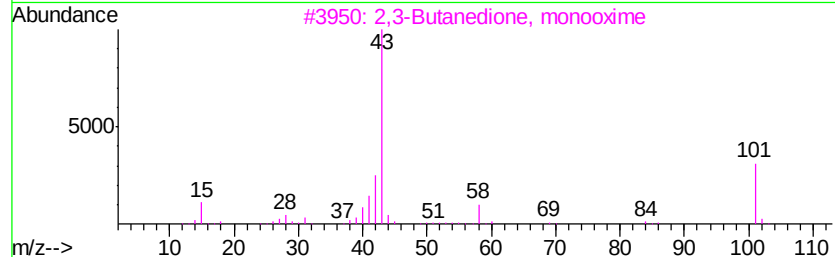
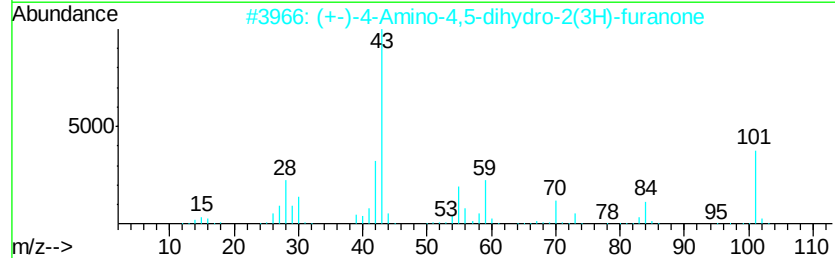
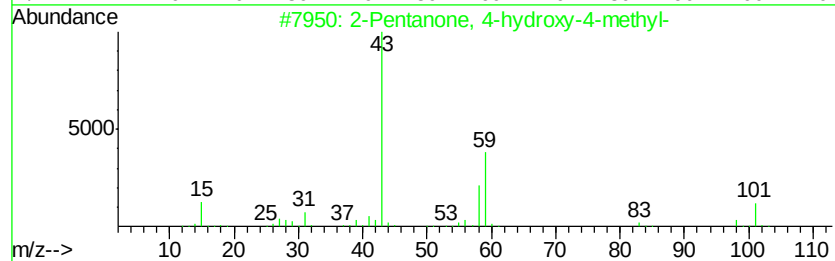
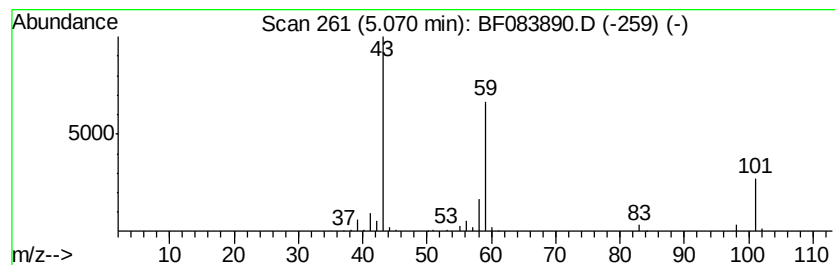
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	18.73 ng	299621	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5		Acetone	58	C3H6O	000067-64-1	10



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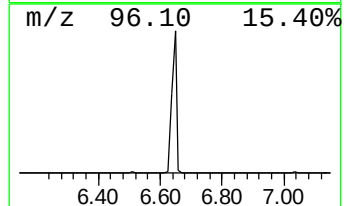
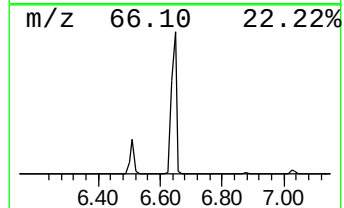
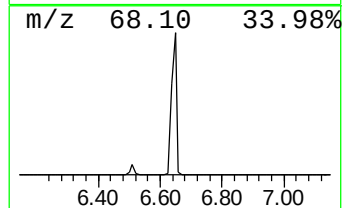
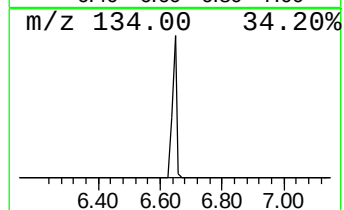
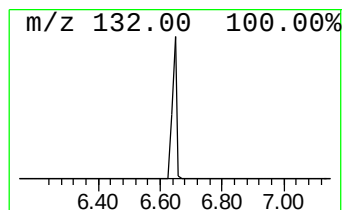
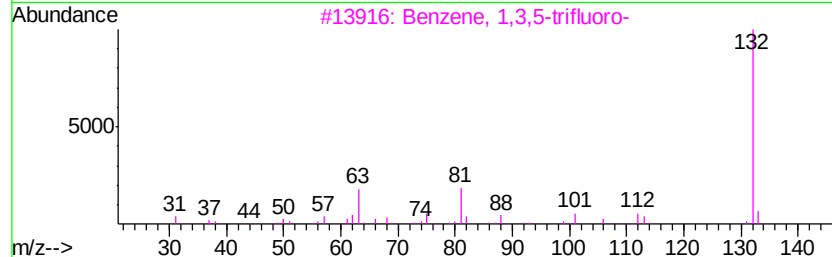
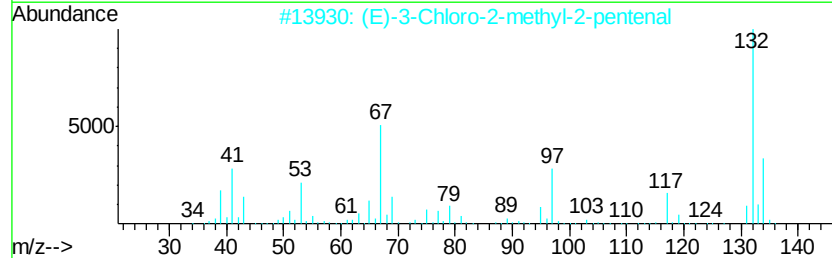
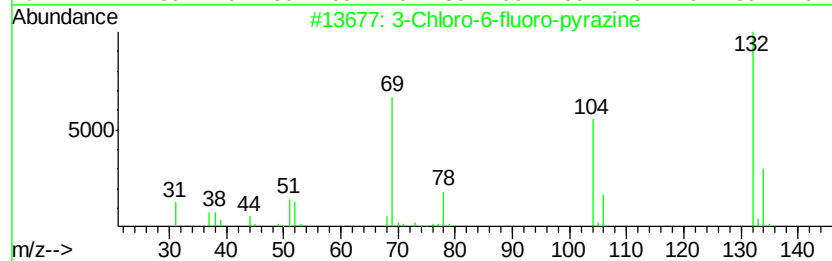
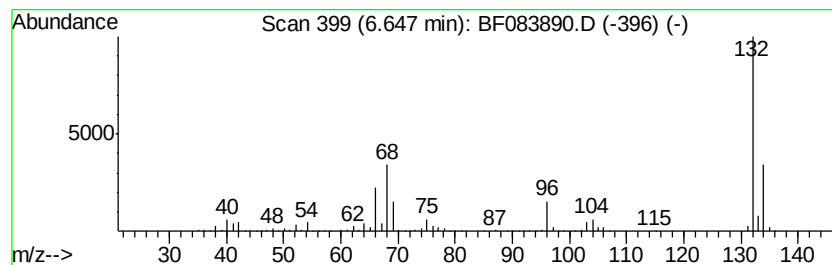
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Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	92.20 ng	1475240	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	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...	5.07	18.7	ng	299621	1	6.88	320020	20.0
unknown6.65	6.65	92.2	ng	1475240	1	6.88	320020	20.0