

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
 Data File : BF085350.D
 Acq On : 10 Mar 2016 23:21
 Operator : UM/SJ
 Sample : H1744-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-3

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-BF030316.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	5.133	246	249	254	rVB	61189	94846	2.56%	0.397%
2	5.533	282	284	287	rBV	1439591	1522945	41.12%	6.369%
3	6.562	371	374	376	rBV	745219	930133	25.11%	3.890%
4	6.699	384	386	389	rBV	2078305	2904905	78.43%	12.148%
5	6.939	404	407	409	rBV	1056889	869616	23.48%	3.637%
6	7.099	418	421	423	rVB	3086647	2939252	79.36%	12.291%
7	7.499	453	456	459	rBV	1993655	2103812	56.80%	8.798%
8	8.219	517	519	522	rBV	947654	996856	26.91%	4.169%
9	9.305	611	614	616	rBV	3092960	3703847	100.00%	15.489%
10	9.979	670	673	675	rBV	1049244	907124	24.49%	3.793%
11	10.768	739	742	744	rBV	2020167	1922802	51.91%	8.041%
12	11.465	801	803	805	rBV	890533	801008	21.63%	3.350%
13	12.014	849	851	860	rVB2	22731	43651	1.18%	0.183%
14	13.054	939	942	944	rBV	2246105	2821780	76.19%	11.800%
15	14.105	1031	1034	1036	rBV	462926	650264	17.56%	2.719%
16	15.568	1158	1162	1164	rBV	443504	700568	18.91%	2.930%

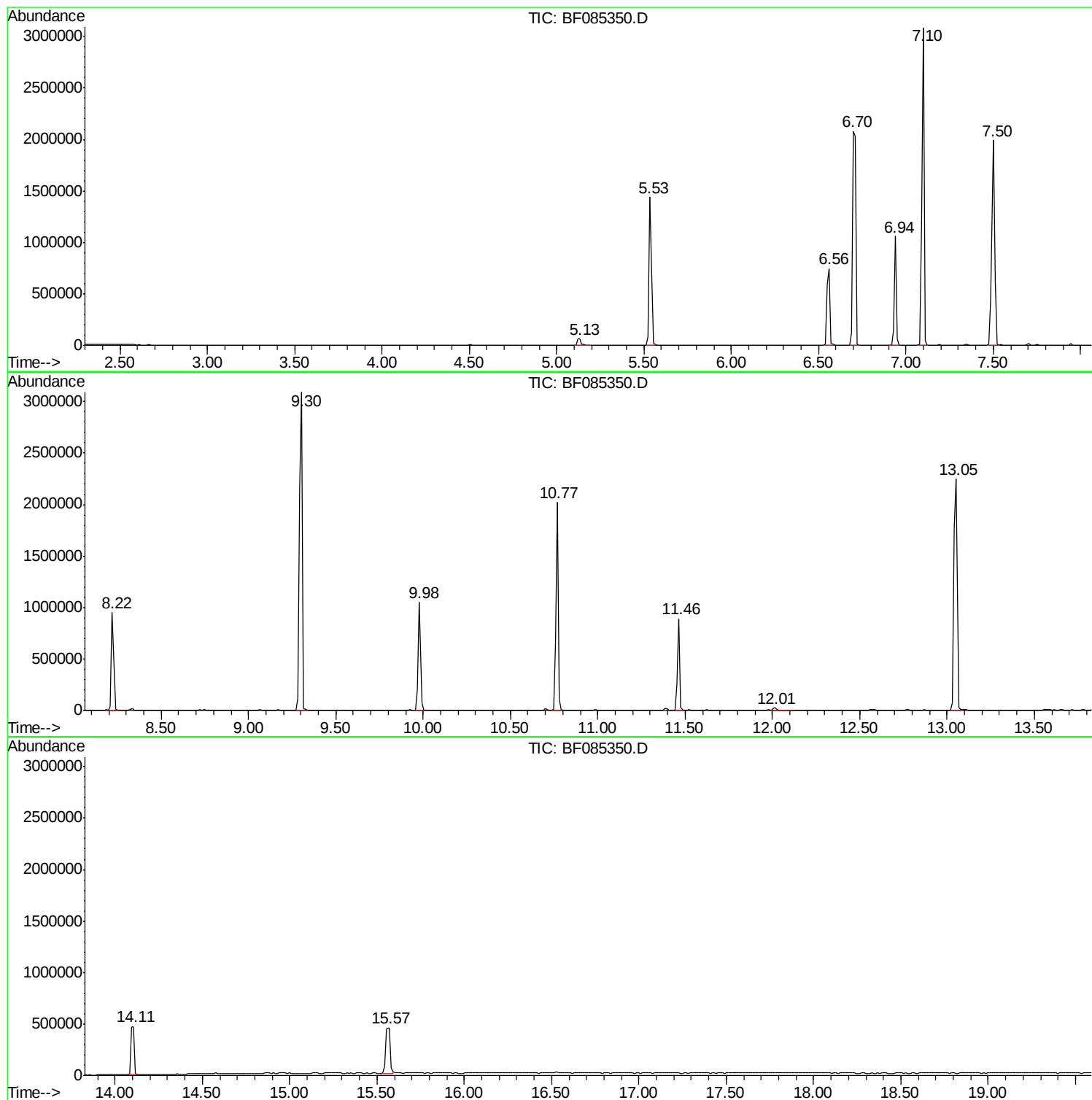
Sum of corrected areas: 23913409

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.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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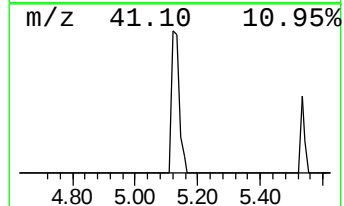
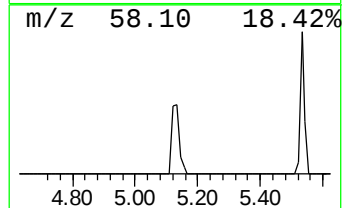
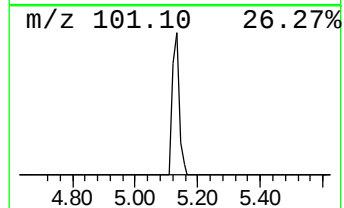
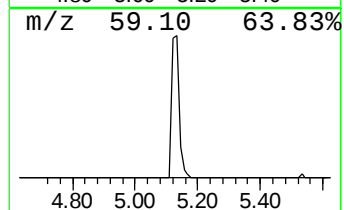
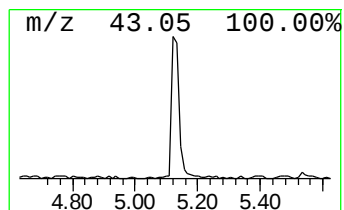
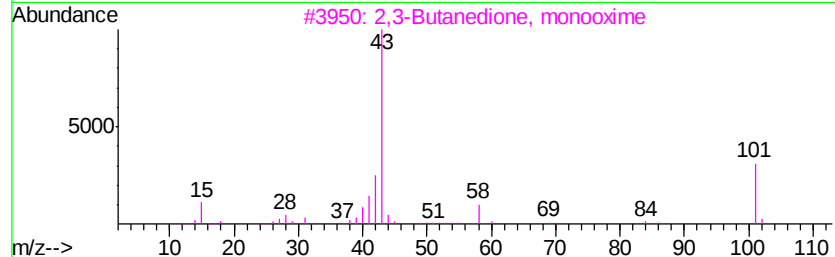
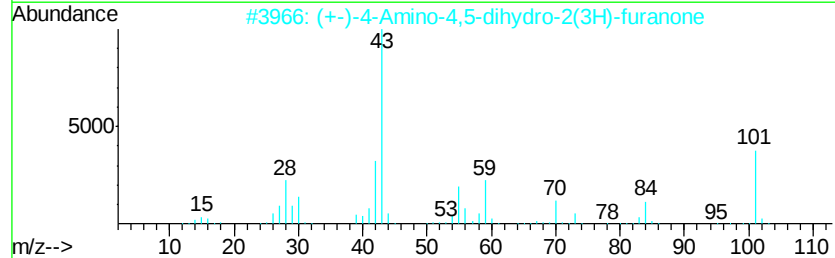
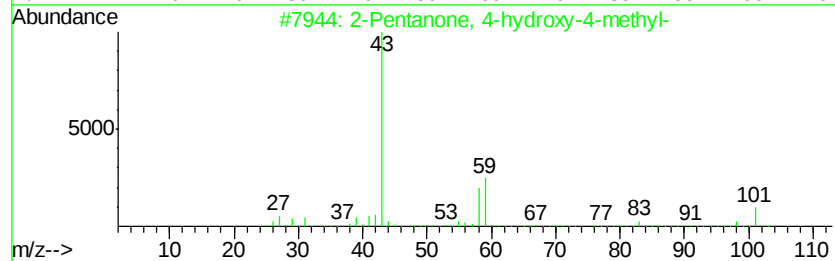
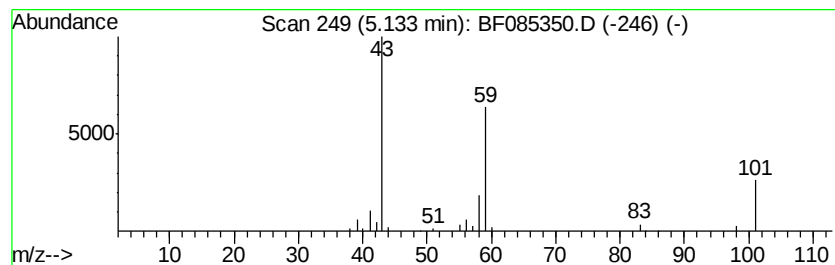
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	2.18 ng	94846	1,4-Dichlorobenzene-d4	6.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12



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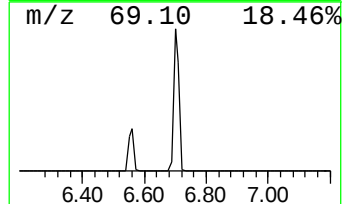
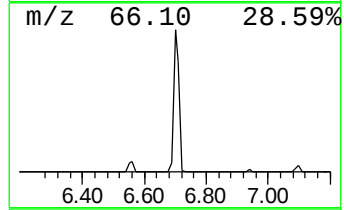
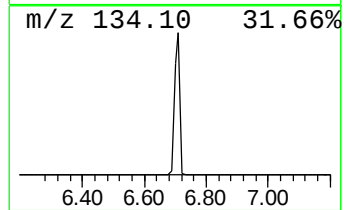
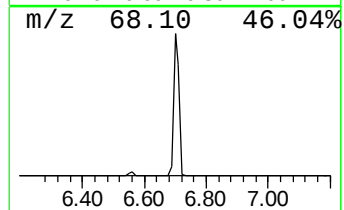
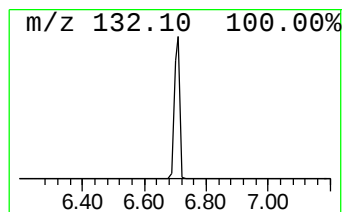
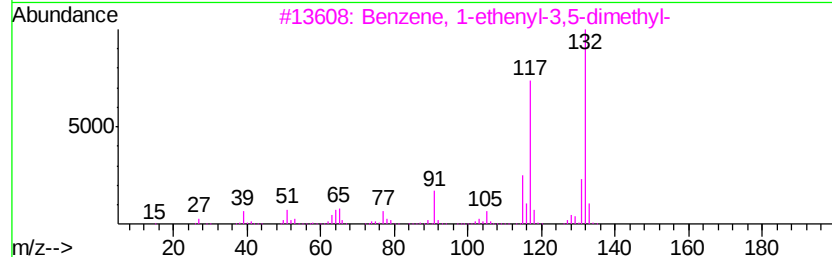
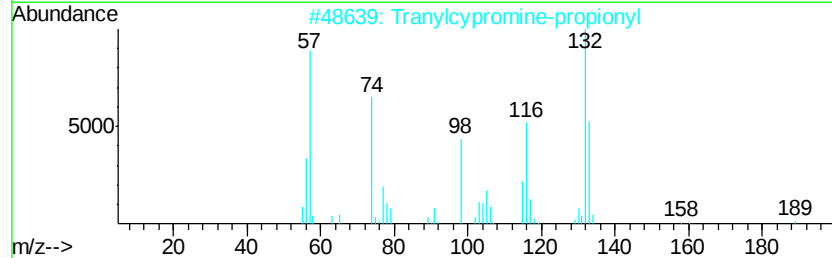
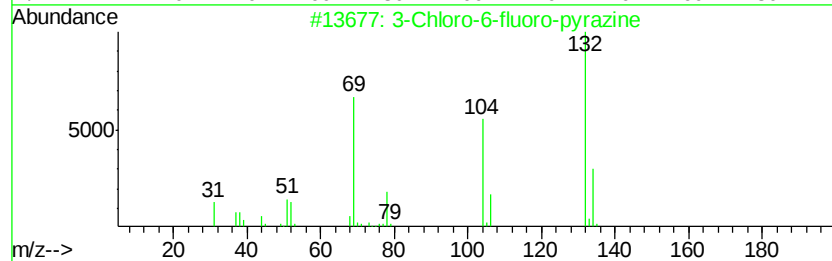
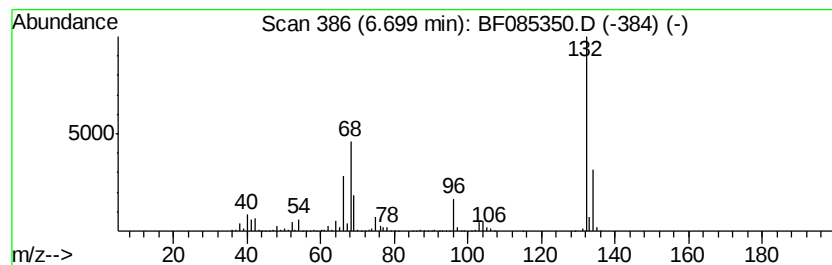
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Peak Number 2 unknown6.70 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	66.81 ng	2904910	1,4-Dichlorobenzene-d4	6.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	2.2	ng	94846	1	6.94	869616	20.0
unknown6.70	6.70	66.8	ng	2904910	1	6.94	869616	20.0