

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122215\
 Data File : BF084013.D
 Acq On : 23 Dec 2015 2:53
 Operator : UM/SJ
 Sample : G4916-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20151214MGP209BRV46.5

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-BF122215.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	4.487	208	210	213	rBB	56846	69954	3.55%	0.628%
2	5.001	253	255	263	rBB	93680	102142	5.18%	0.917%
3	5.436	291	293	296	rBV	431622	588531	29.84%	5.284%
4	6.476	382	384	386	rBV	252247	354363	17.97%	3.182%
5	6.601	393	395	397	rBV	1295576	1155999	58.62%	10.379%
6	6.830	413	415	417	rBV	262233	240045	12.17%	2.155%
7	6.990	426	429	431	rBV	767401	1039796	52.73%	9.336%
8	7.390	462	464	467	rBV	762270	734056	37.22%	6.591%
9	8.110	525	527	530	rBV	384861	336857	17.08%	3.024%
10	9.196	619	622	624	rBV	1616426	1816383	92.11%	16.308%
11	9.870	679	681	683	rBV	364290	382228	19.38%	3.432%
12	10.659	747	750	752	rBV	1109182	1118623	56.73%	10.044%
13	11.356	809	811	813	rBV	303995	406168	20.60%	3.647%
14	12.934	947	949	952	rBV	1585787	1972001	100.00%	17.706%
15	13.985	1039	1041	1044	rBV	417671	444352	22.53%	3.990%
16	15.414	1164	1166	1173	rBV	261374	376241	19.08%	3.378%

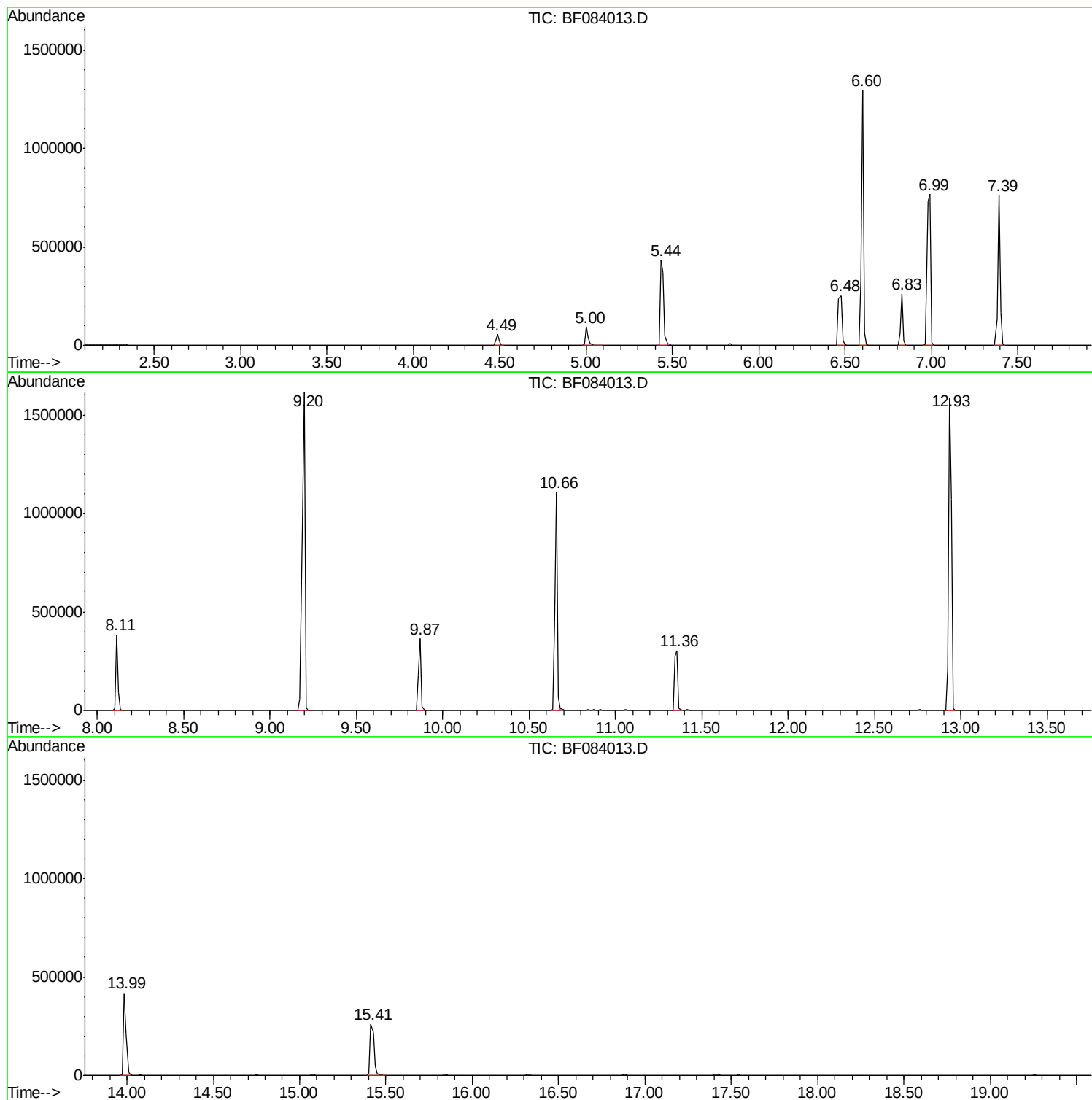
Sum of corrected areas: 11137739

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122215.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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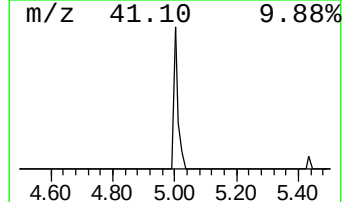
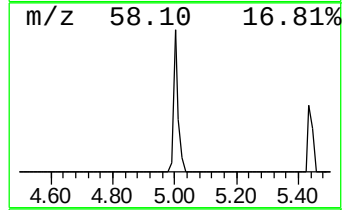
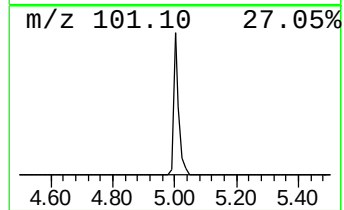
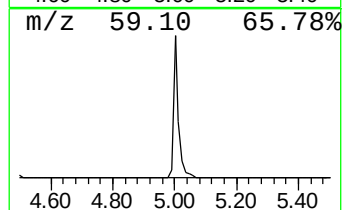
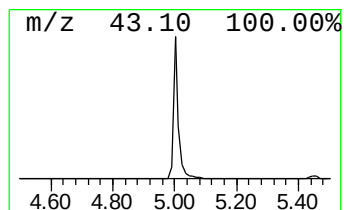
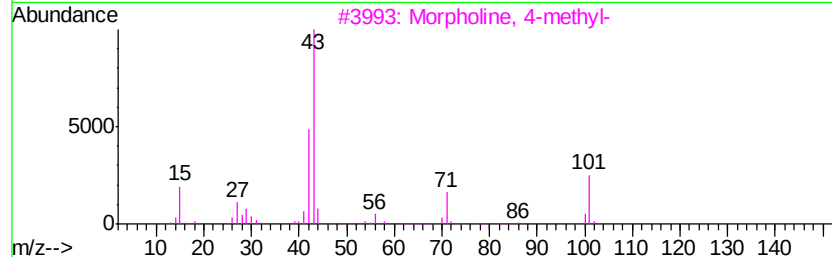
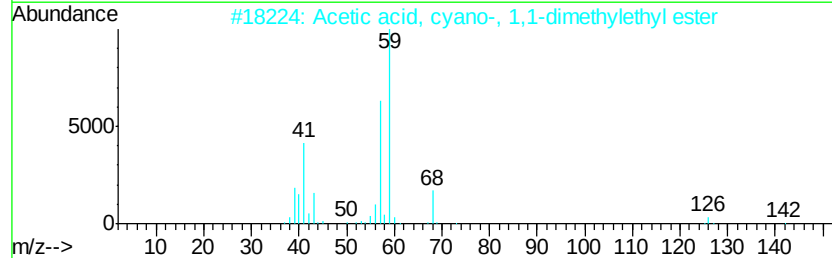
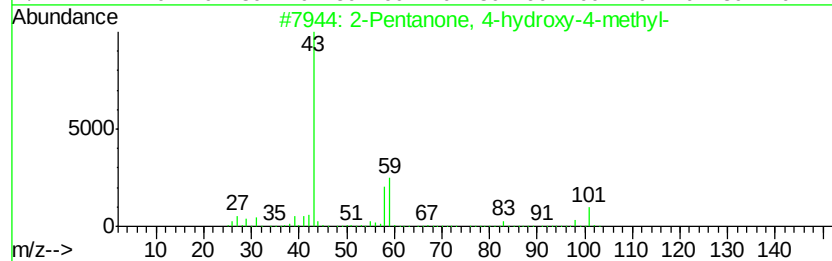
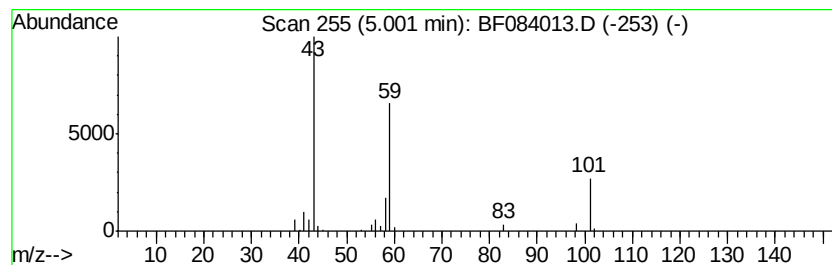
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	8.51 ng	102142	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	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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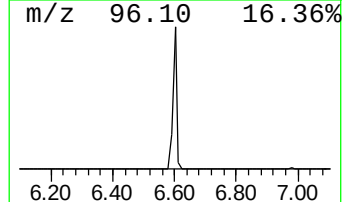
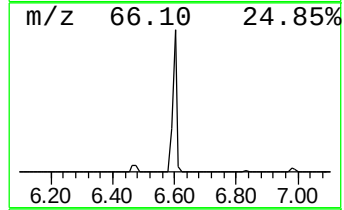
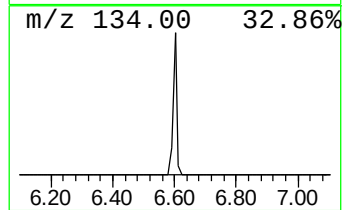
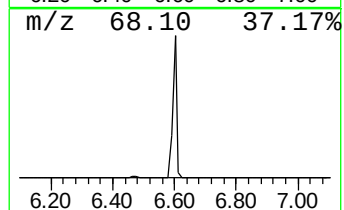
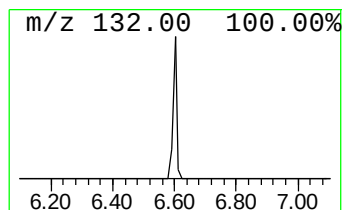
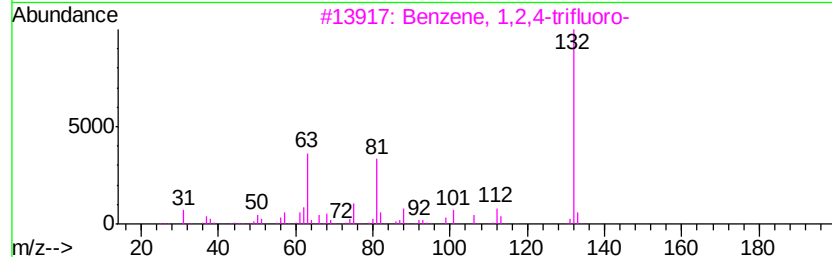
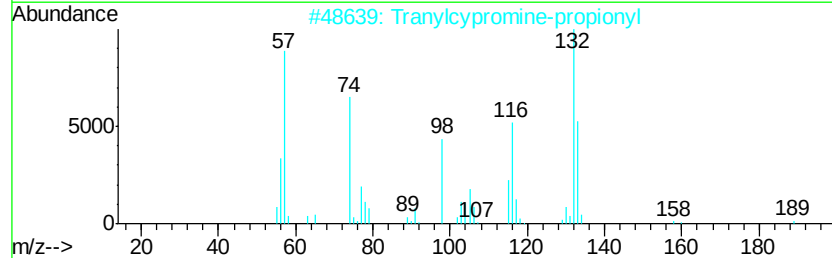
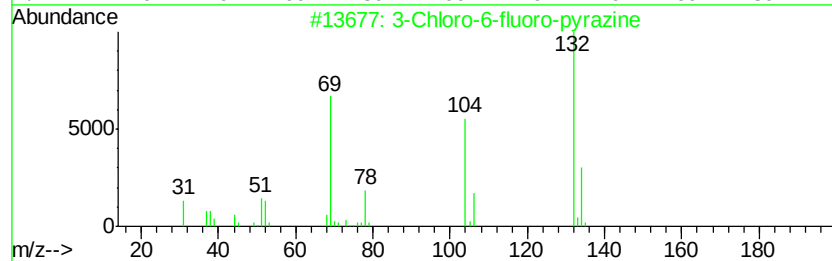
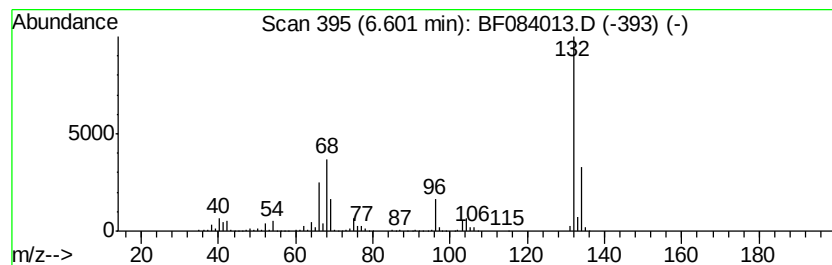
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Peak Number 3 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	96.32 ng	1156000	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	8.5	ng	102142	1	6.83	240045	20.0
unknown6.60	6.60	96.3	ng	1156000	1	6.83	240045	20.0