

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120516\
 Data File : BF091433.D
 Acq On : 5 Dec 2016 22:18
 Operator : UM/SJ
 Sample : H5838-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB120216

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-BF112916.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	5.030	237	240	242	rBV	371967	535648	8.23%	1.480%
2	5.441	274	276	279	rBV	1784858	1895221	29.12%	5.238%
3	6.470	364	366	369	rBV	1376236	1168036	17.94%	3.228%
4	6.619	376	379	381	rBV	3760131	4137842	63.57%	11.435%
5	6.847	397	399	401	rBV	998504	815445	12.53%	2.254%
6	7.007	410	413	415	rBV	3681290	3387139	52.04%	9.361%
7	7.419	445	449	451	rBV	1856477	2915480	44.79%	8.057%
8	8.127	509	511	514	rVB	833084	1118650	17.19%	3.091%
9	9.213	603	606	608	rBV	5571335	5376894	82.61%	14.859%
10	9.888	662	665	667	rBV	1536995	1402301	21.54%	3.875%
11	10.676	731	734	736	rBV	3627166	3719593	57.14%	10.279%
12	11.362	792	794	797	rVB	971544	1220824	18.76%	3.374%
13	12.962	931	934	936	rBV	5829253	6509062	100.00%	17.988%
14	14.002	1022	1025	1027	rBV	998651	1240590	19.06%	3.428%
15	15.419	1146	1149	1152	rBV	493933	742706	11.41%	2.053%

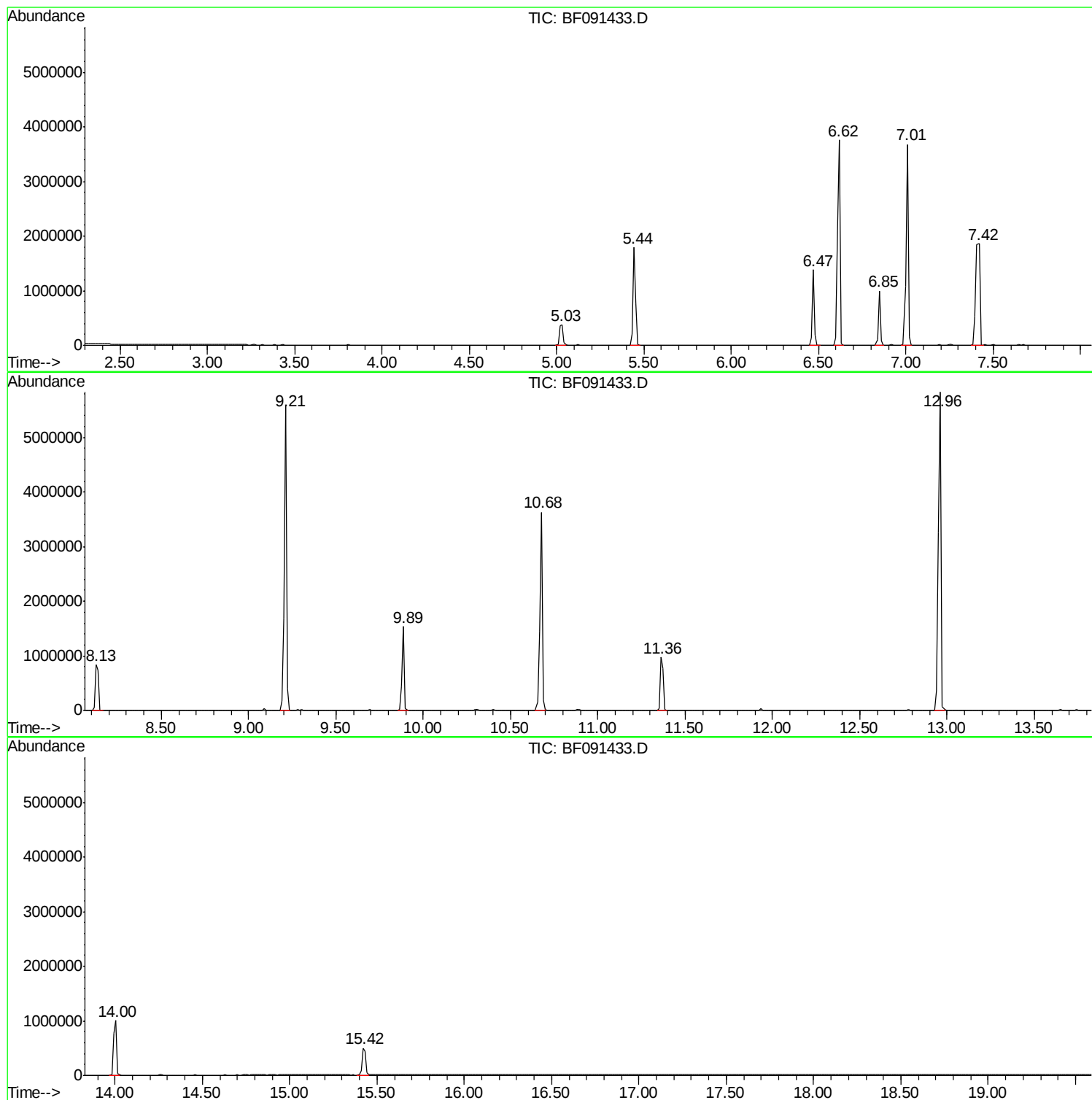
Sum of corrected areas: 36185431

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.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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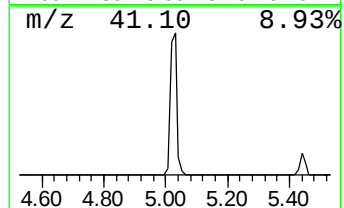
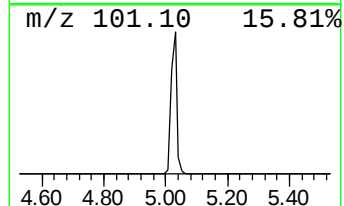
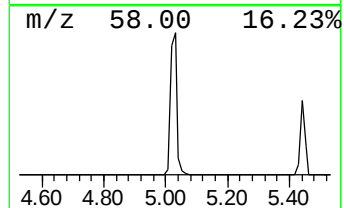
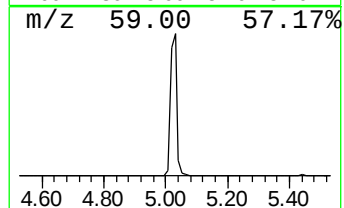
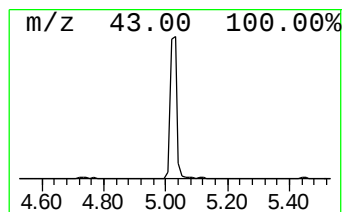
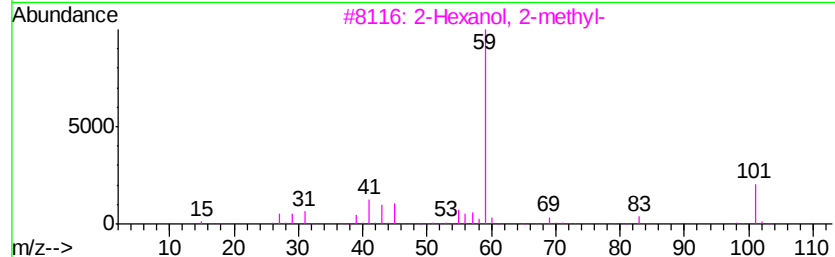
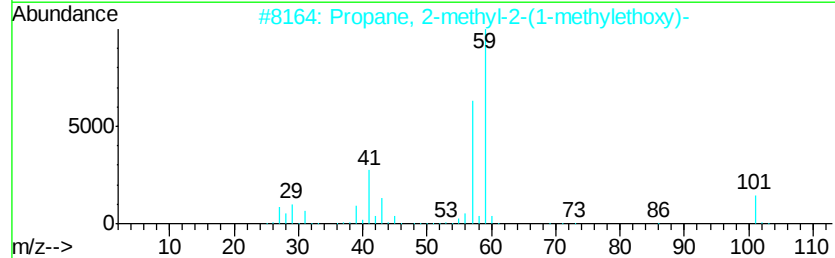
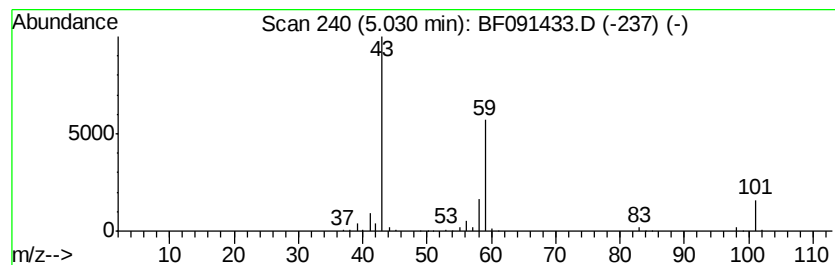
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	13.14 ng	535648	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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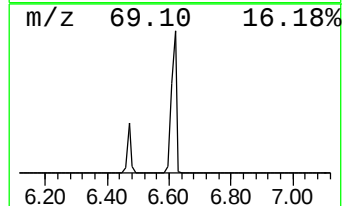
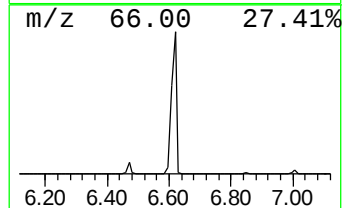
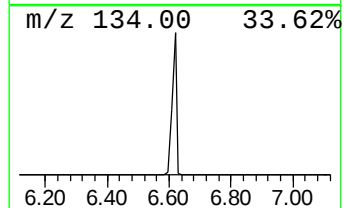
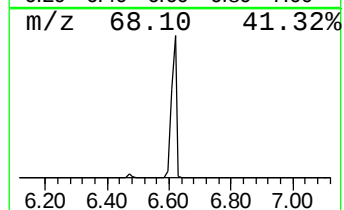
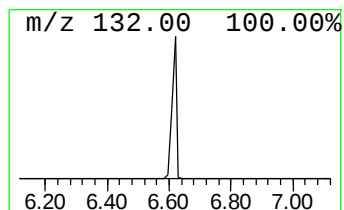
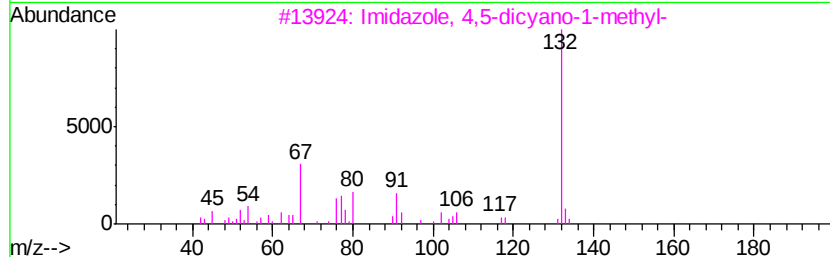
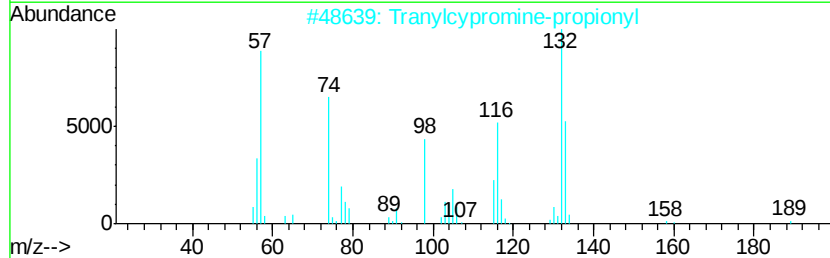
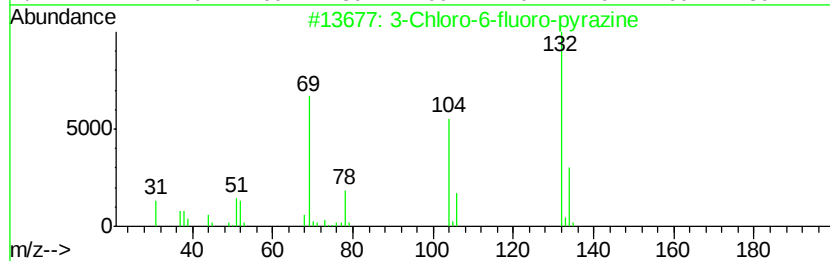
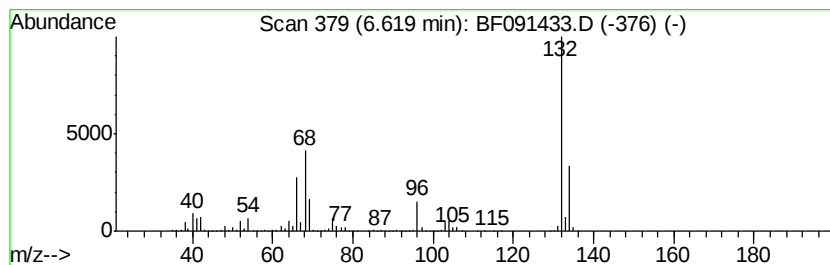
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	101.49 ng	4137840	1,4-Dichlorobenzene-d4	6.85

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		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	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...	5.03	13.1	ng	535648	1	6.85	815445	20.0
unknown6.62	6.62	101.5	ng	4137840	1	6.85	815445	20.0