

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101218\
 Data File : BF109873.D
 Acq On : 13 Oct 2018 9:37
 Operator : JU/SJ
 Sample : PB113699BL
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB113699BL

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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	4.904	475	479	497	rBV	155289	233120	7.86%	1.291%
2	5.322	546	550	572	rBV	1307971	1504734	50.73%	8.336%
3	6.340	719	723	740	rBV	1252161	1624389	54.76%	8.999%
4	6.463	740	744	769	rVB	1641191	1796876	60.58%	9.955%
5	6.693	780	783	805	rVV	413263	549118	18.51%	3.042%
6	6.845	805	809	842	rVB	1773798	2045171	68.95%	11.330%
7	7.257	875	879	934	rBV	1182686	1687695	56.90%	9.350%
8	7.975	997	1001	1024	rBV	504051	691284	23.30%	3.830%
9	9.051	1178	1184	1215	rBV	2672992	2966253	100.00%	16.433%
10	9.722	1294	1298	1324	rBV	629593	709917	23.93%	3.933%
11	10.504	1428	1431	1457	rBV	586342	853279	28.77%	4.727%
12	11.198	1546	1549	1574	rBV	389211	590725	19.91%	3.273%
13	12.780	1814	1818	1841	rBV	1881616	1942335	65.48%	10.760%
14	13.827	1992	1996	2018	rBV	358677	485210	16.36%	2.688%
15	15.227	2229	2234	2254	rBV	195507	370573	12.49%	2.053%

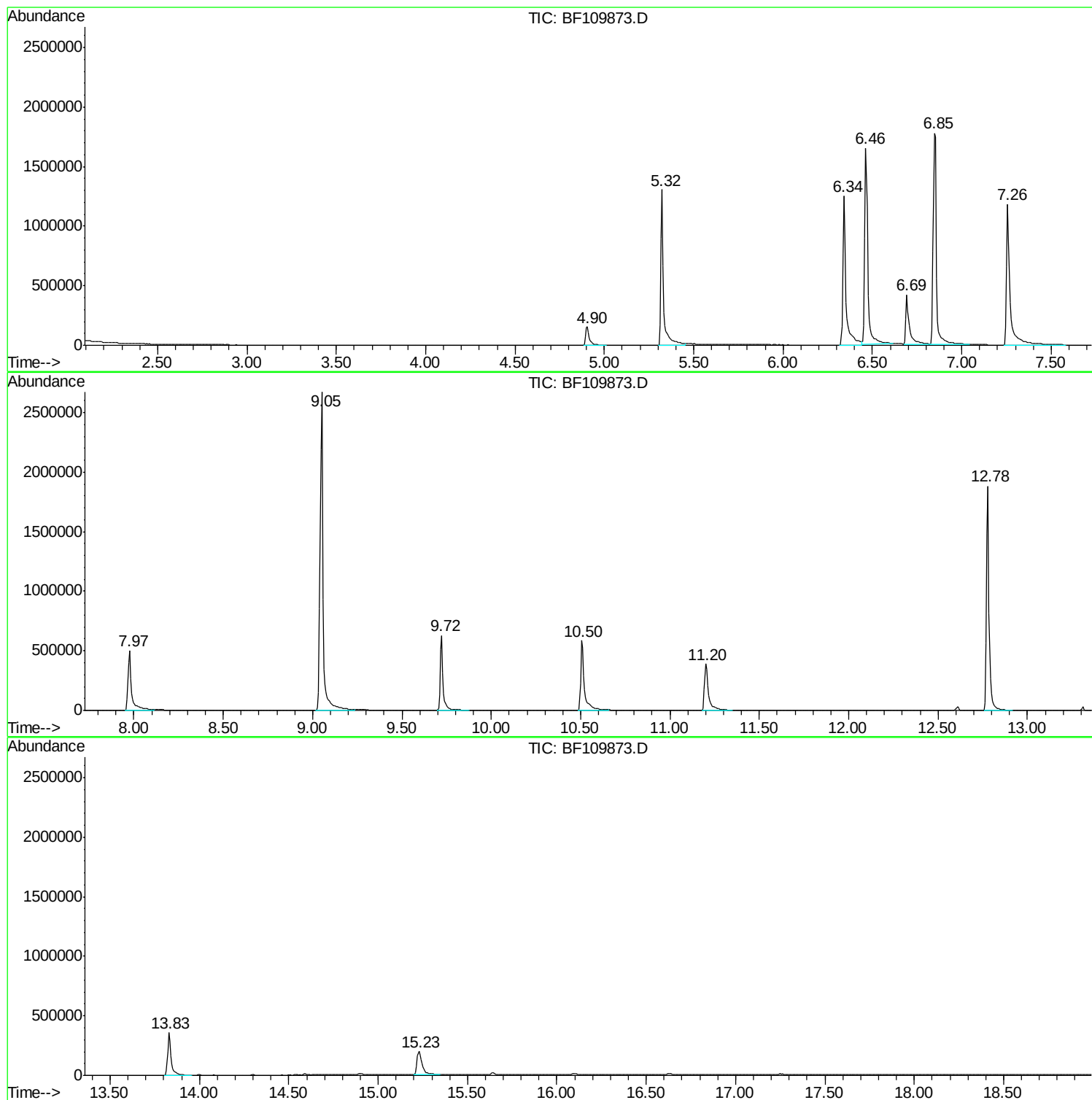
Sum of corrected areas: 18050679

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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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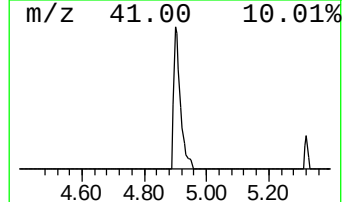
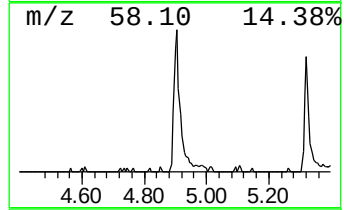
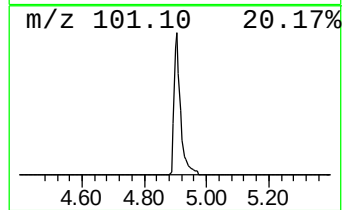
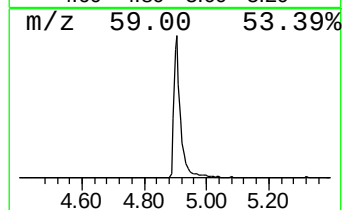
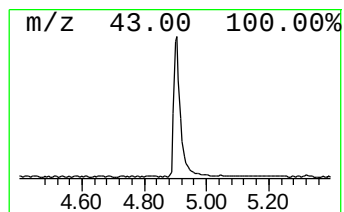
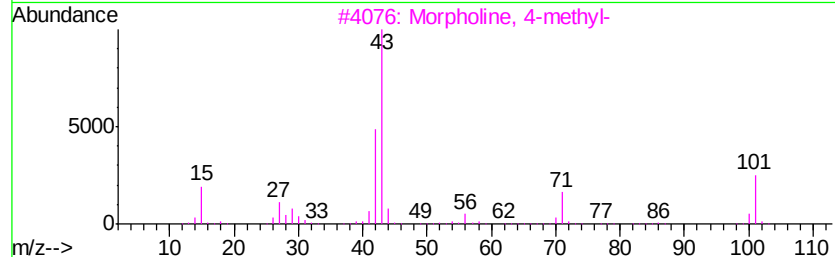
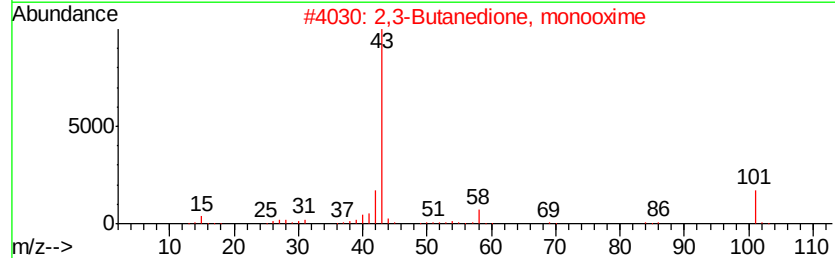
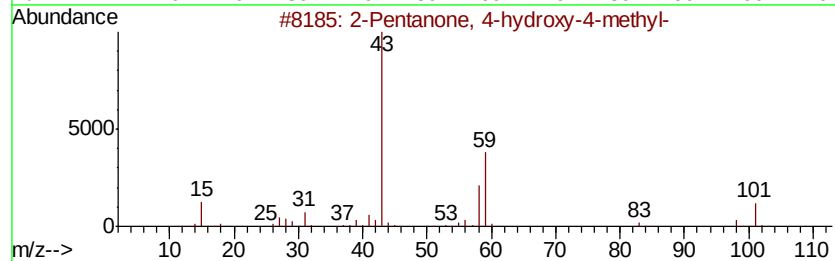
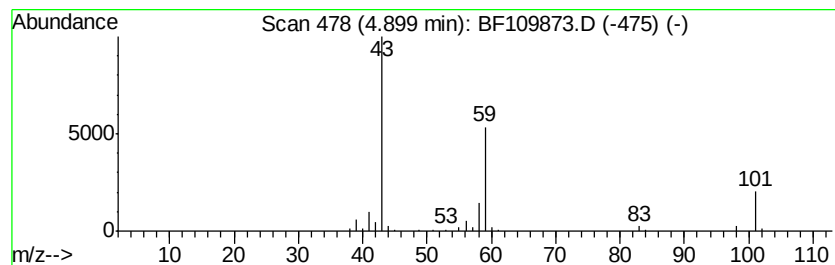
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	8.49 ng	233120	1,4-Dichlorobenzene-d4	6.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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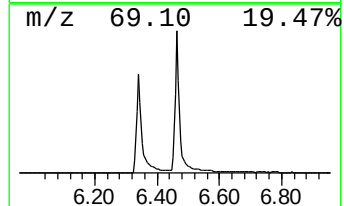
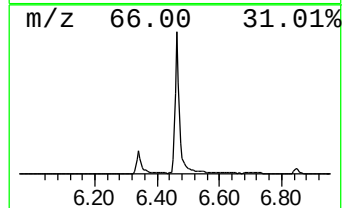
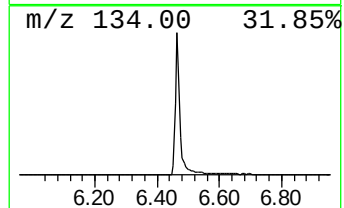
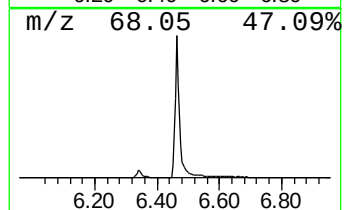
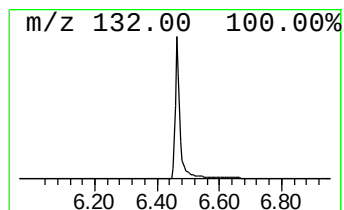
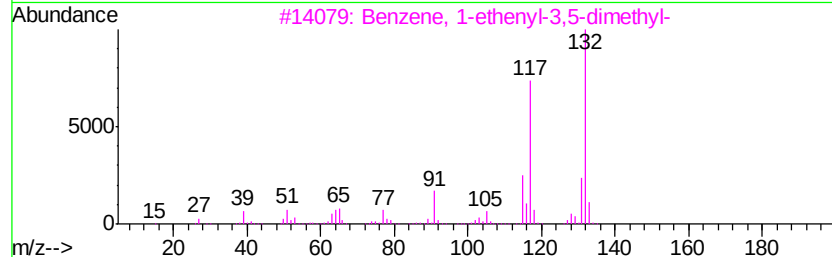
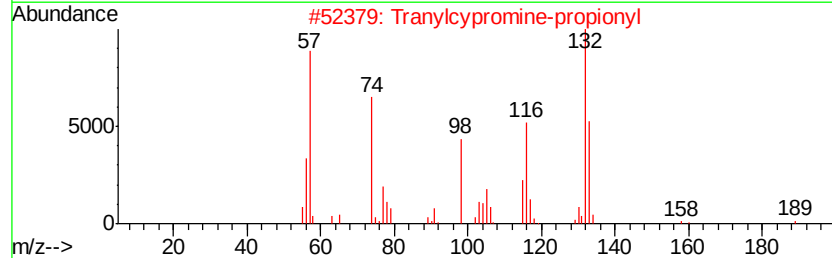
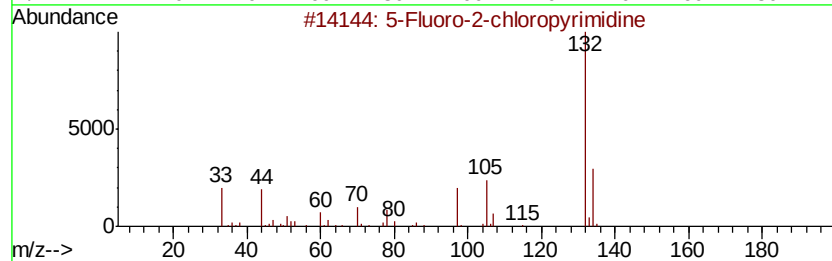
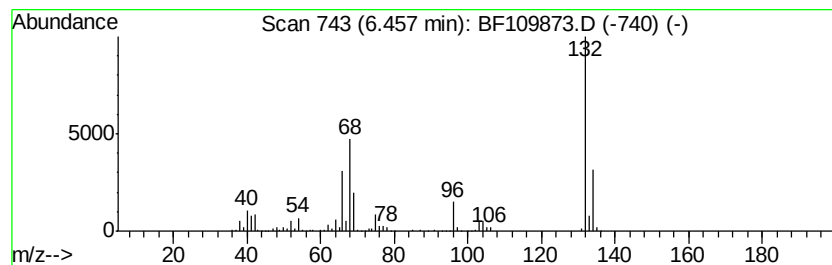
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Peak Number 2 unknown6.46 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	65.45 ng	1796880	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	8.5	ng	233120	1	6.69	549118	20.0
unknown6.46	6.46	65.5	ng	1796880	1	6.69	549118	20.0