

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101218\
 Data File : BF109876.D
 Acq On : 13 Oct 2018 10:59
 Operator : JU/SJ
 Sample : PB113650BL
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB113650BL

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	4.899	474	478	498	rBV	139618	212658	8.03%	1.296%
2	5.322	546	550	567	rBV	1076942	1310733	49.47%	7.986%
3	6.340	720	723	741	rBV	1163706	1470732	55.51%	8.961%
4	6.463	741	744	780	rVV	1490596	1744807	65.85%	10.631%
5	6.693	780	783	805	rVV	381009	543932	20.53%	3.314%
6	6.846	805	809	836	rVV	1625053	1849024	69.78%	11.266%
7	7.257	875	879	917	rBV	1069077	1498881	56.57%	9.133%
8	7.975	997	1001	1025	rBV	466089	654365	24.70%	3.987%
9	9.045	1179	1183	1206	rBV	2159455	2649692	100.00%	16.145%
10	9.722	1294	1298	1312	rBV	575783	655078	24.72%	3.991%
11	10.504	1428	1431	1452	rBV	465896	726544	27.42%	4.427%
12	11.198	1546	1549	1575	rBV	340343	537668	20.29%	3.276%
13	12.775	1813	1817	1842	rBV	1518444	1745508	65.88%	10.635%
14	13.827	1992	1996	2014	rBV	333635	464409	17.53%	2.830%
15	15.227	2229	2234	2253	rBV	181585	348221	13.14%	2.122%

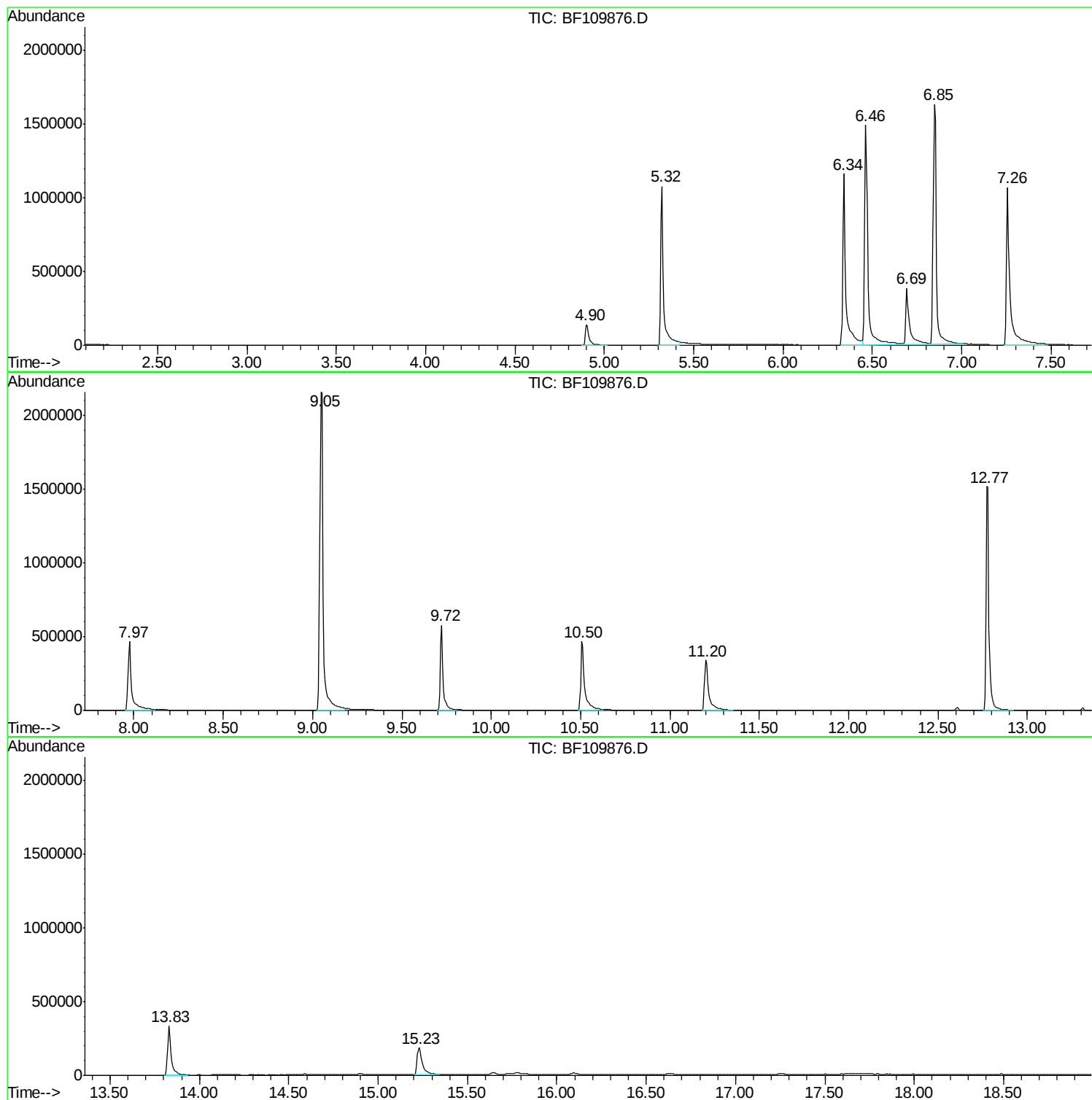
Sum of corrected areas: 16412252

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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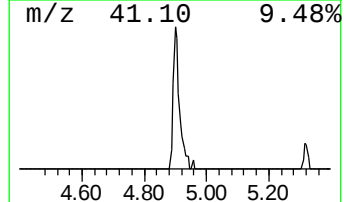
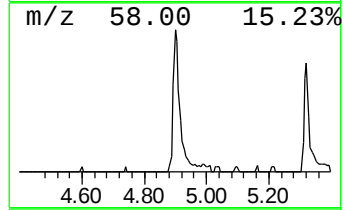
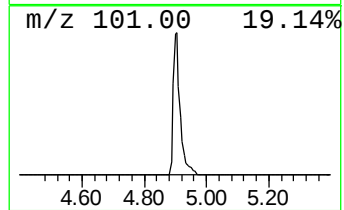
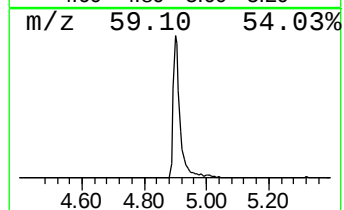
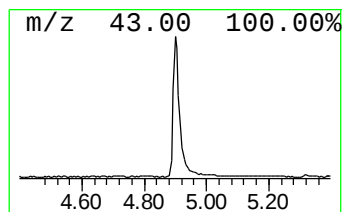
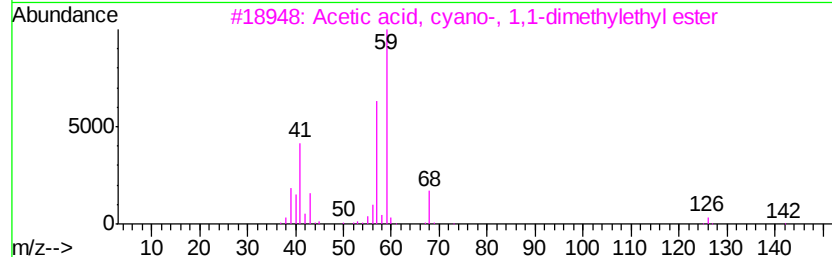
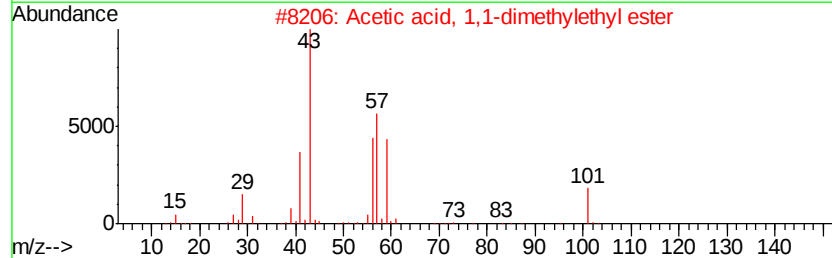
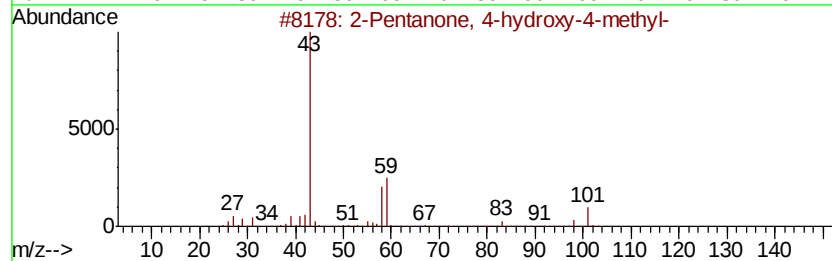
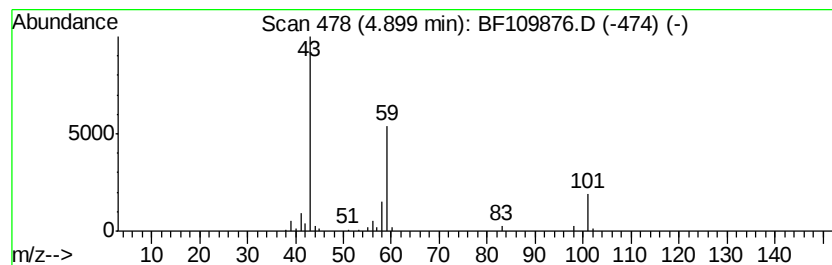
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	7.82 ng	212658	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	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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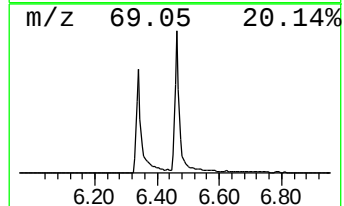
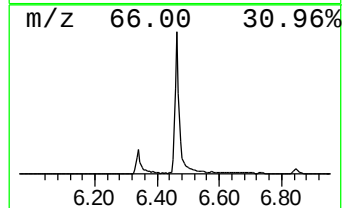
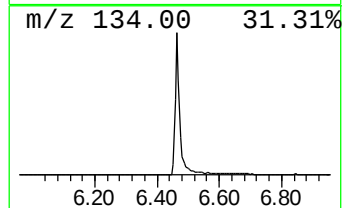
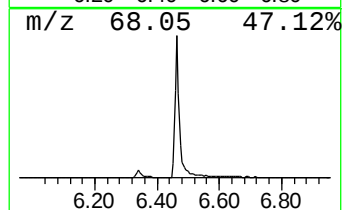
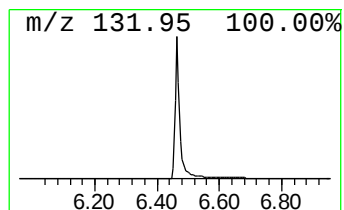
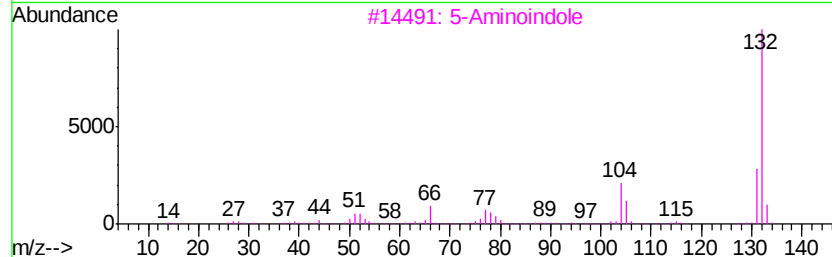
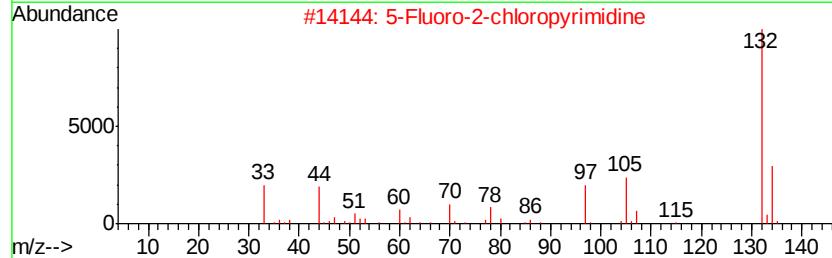
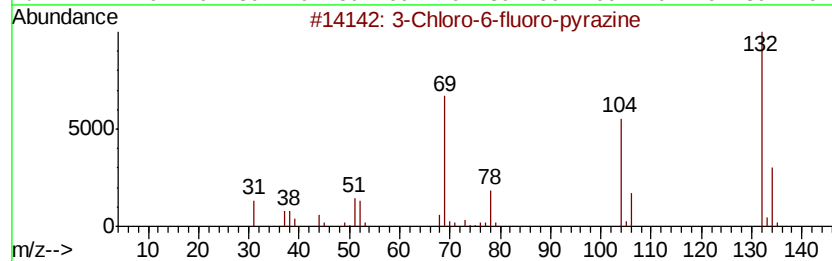
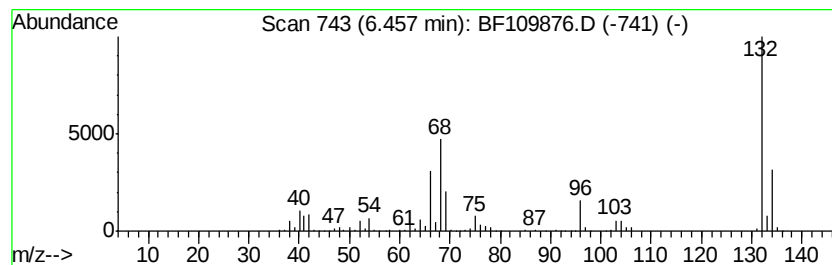
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Peak Number 2 unknown6.46 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	64.16 ng	1744810	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	7.8	ng	212658	1	6.69	543932	20.0
unknown6.46	6.46	64.2	ng	1744810	1	6.69	543932	20.0