

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062618\
 Data File : BF106853.D
 Acq On : 27 Jun 2018 5:33
 Operator : JU/SJ
 Sample : J3639-06
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 062018-DBRI-E-U-S

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-BF061918.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	4.090	294	298	303	rBB	17712	20858	1.63%	0.284%
2	5.184	482	484	492	rBB	15366	16571	1.30%	0.226%
3	5.595	551	554	570	rBV	353467	331062	25.92%	4.506%
4	6.595	721	724	737	rBV2	261392	252508	19.77%	3.437%
5	6.731	743	747	751	rBV	663059	587922	46.04%	8.001%
6	6.954	782	785	789	rBB	288815	252948	19.81%	3.443%
7	7.113	808	812	816	rBV	933401	858809	67.25%	11.688%
8	7.525	877	882	885	rBV	665595	648594	50.79%	8.827%
9	8.242	1000	1004	1020	rBB	300073	295206	23.12%	4.018%
10	9.313	1182	1186	1189	rBV	1268433	1159290	90.78%	15.778%
11	9.701	1250	1252	1259	rBB	23982	23414	1.83%	0.319%
12	9.995	1299	1302	1311	rVB	302465	298505	23.37%	4.063%
13	10.654	1411	1414	1417	rBV	29975	21797	1.71%	0.297%
14	10.789	1434	1437	1446	rBV	440245	413904	32.41%	5.633%
15	11.489	1552	1556	1564	rBV	363479	313696	24.56%	4.269%
16	12.871	1789	1791	1794	rBB	18692	13544	1.06%	0.184%
17	13.077	1821	1826	1829	rBV	1332663	1277046	100.00%	17.380%
18	13.954	1969	1975	1984	rBV	13877	19289	1.51%	0.263%
19	14.142	2003	2007	2013	rVB	330543	325364	25.48%	4.428%
20	15.671	2260	2267	2274	rBV2	151152	217378	17.02%	2.958%

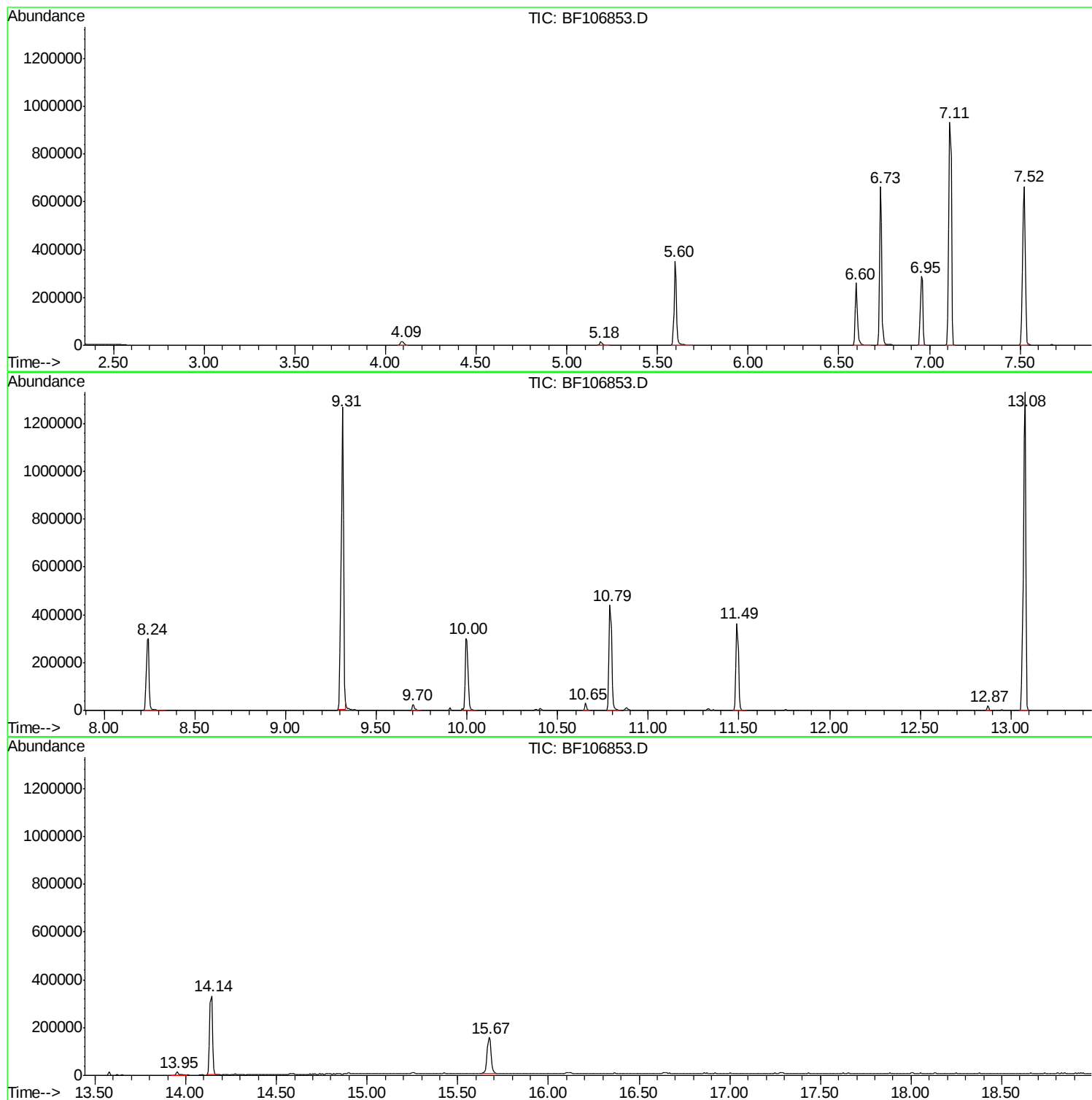
Sum of corrected areas: 7347705

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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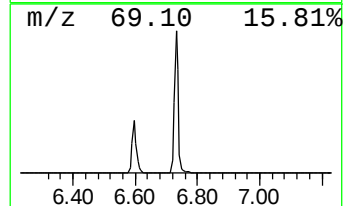
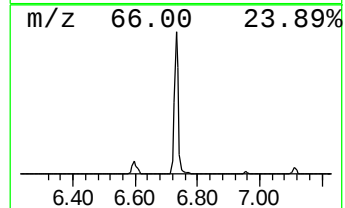
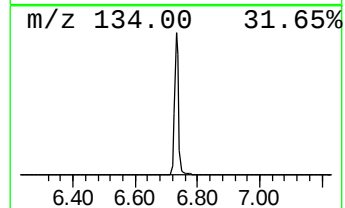
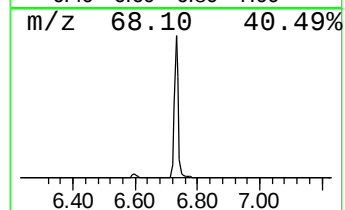
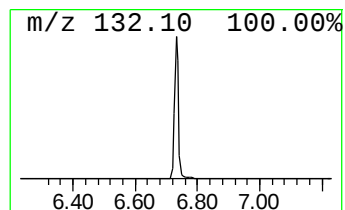
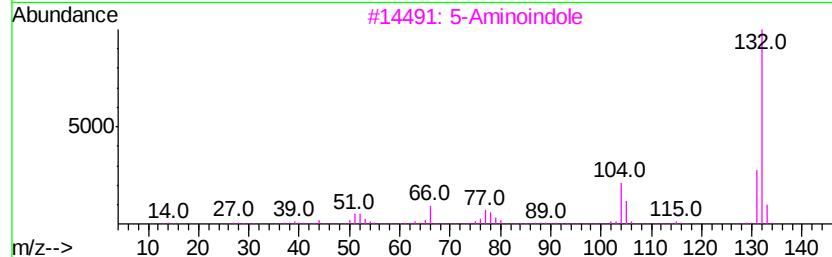
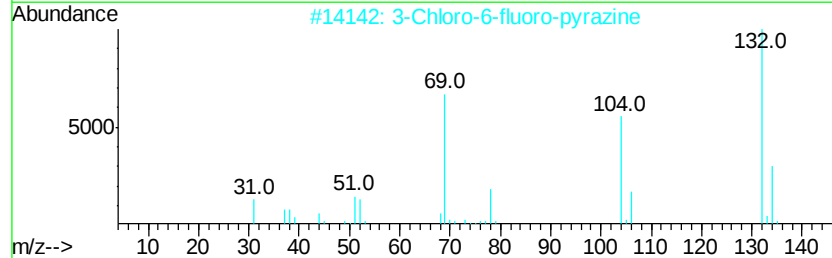
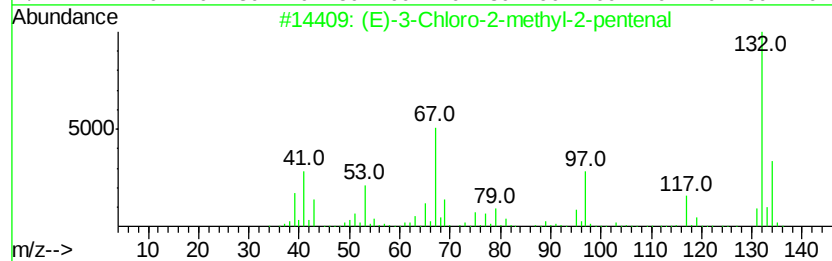
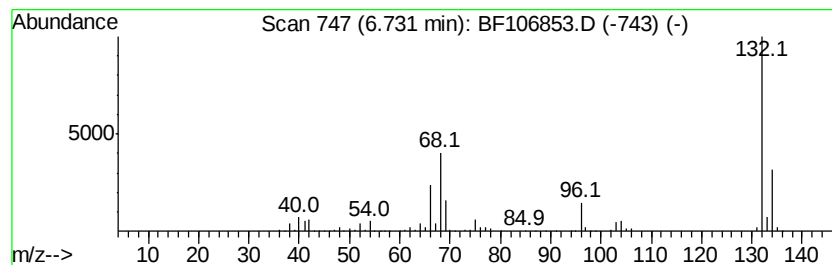
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.73 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	46.49 ng	587922	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
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
unknown6.73	6.73	46.5	ng	587922	1	6.95	252948	20.0