# Class06 Homework

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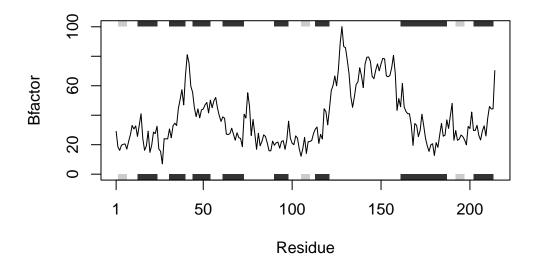
# Can you improve this analysis code?

#### **Explanation for this function:**

This function that I created take the input string, which is the protein code, and it reads the PDB file of the input string. Then, we use the trim.pdb function to create a smaller PDB object. After that, it will take the output of s.chainA and accessing the atom column and b element. Finally, the function creates a plot of the output from s.b and label x as Bfactor and y as residue.

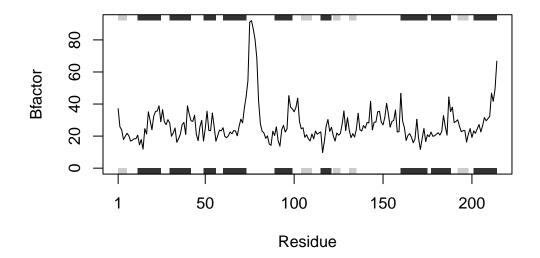
```
library(bio3d)
analysis_fun <- function(string){
   s.name <- read.pdb(string)
   s.chainA <- trim.pdb(s.name, chain="A", elety="CA")
   s.b <- s.chainA$atom$b
   plotb3(s.b, sse=s.chainA, typ="l", ylab="Bfactor")
}
analysis_fun("4AKE")</pre>
```

Note: Accessing on-line PDB file



# analysis\_fun("1AKE")

Note: Accessing on-line PDB file PDB has ALT records, taking A only, rm.alt=TRUE



# analysis\_fun("1E4Y")

Note: Accessing on-line PDB file

