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title: "Homework 6 Week 5"
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> Q6. How would you generalize the original code above to work with any set of input protein structures?

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```{r}
installing packages
install.packages("bio3d")

```

```{r}
The function below takes the B data and plots their subsequent factors for each
individual protein and reads the trends from the previous questions
protein_analysis <- function(files) {
 # load library and reading data files of the protein
 library(bio3d)
 proteins_data <- lapply(files, read.pdb)
 # B factor information being extracted
 b_factors_list <- lapply(proteins_data, function(protein) trim.pdb(protein, chain="A",
eley="CA")$atom$b)
 # Plot B-factors by using a for loop to iterate the protein structure
 for (index in seq_along(b_factors_list)) {
 plotb3(b_factors_list[[index]], sse=proteins_data[[index]]$atom, typ="l",
ylab="Bfactor", main=paste("Protein ", index))
 }
 # the plot above is the output as it shows the B factors for protein similarity. The for
loop allows the b factors and protein data to loop and make it more "readable"
 # dendrogram
 hierarchical_clustering <- hclust(dist(do.call(rbind, b_factors_list)))
 plot(hierarchical_clustering, main="Cluster Dendrogram")
}# the plot above is the output as it shows the B factors for protein similarity.
protein_files <- c("4AKE", "1AKE", "1E4Y")
protein files is the input
protein_analysis(protein_files)
```

}

```