

class09

Phoebe LI

2/13/2022

```
# import the csv file
db <- read.csv("Data Export Summary.csv", row.names = 1)
db
```

	X.ray	NMR	EM	Multiple.methods	Neutron	Other	Total
## Protein (only)	144433	11881	6732	182	70	32	163330
## Protein/Oligosaccharide	8543	31	1125	5	0	0	9704
## Protein/NA	7621	274	2165	3	0	0	10063
## Nucleic acid (only)	2396	1399	61	8	2	1	3867
## Other	150	31	3	0	0	0	184
## Oligosaccharide (only)	11	6	0	1	0	4	22

```
# Xray structure percent
xray.percent <- sum(db$X.ray)/sum(db$Total)*100
round(xray.percent, 2)
```

```
## [1] 87.17
```

```
# EM structure percent
EM.percent <- sum(db$EM)/sum(db$Total)*100
round(EM.percent, 2)
```

```
## [1] 5.39
```

Q1: What percentage of structures in the PDB are solved by X-Ray and Electron Microscopy.

x-ray:87.2% EM:5.39%

Q2: What proportion of structures in the PDB are protein?

```
percent.protein <- (db$Total[1])/sum(db$Total)
round(percent.protein*100, 2)
```

```
## [1] 87.26
```

Q3: Type HIV in the PDB website search box on the home page and determine how many HIV-1 protease structures are in the current PDB?

4486 Structures

Q4: Water molecules normally have 3 atoms. Why do we see just one atom per water molecule in this structure?

Because It does not show the H atom in this structure.

```
install.packages("bio3d", repos="http://cran.us.r-project.org")
```

Introduction to Bio3D in R

```
##  
## The downloaded binary packages are in  
## /var/folders/9d/xssg21015fq5rb8769f22wfw0000gn/T//RtmptUT7iY/downloaded_packages  
  
library(bio3d)
```

Reading PDB file data into R

```
pdb <- read.pdb("1hsg")  
  
## Note: Accessing on-line PDB file  
  
pdb  
  
##  
## Call: read.pdb(file = "1hsg")  
##  
## Total Models#: 1  
## Total Atoms#: 1686, XYZs#: 5058 Chains#: 2 (values: A B)  
##  
## Protein Atoms#: 1514 (residues/Calpha atoms#: 198)  
## Nucleic acid Atoms#: 0 (residues/phosphate atoms#: 0)  
##  
## Non-protein/nucleic Atoms#: 172 (residues: 128)  
## Non-protein/nucleic resid values: [ HOH (127), MK1 (1) ]  
##  
## Protein sequence:  
## PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYD  
## QILIEICGHKAIGTVLVGPTPVNIIGRNLLTQIGCTLNFPQITLWQRPLVTIKIGGQLKE  
## ALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYDQILIEICGHKAIGTVLVGPTP  
## VNIIGRNLLTQIGCTLNF  
##  
## + attr: atom, xyz, seqres, helix, sheet,  
## calpha, remark, call
```

```
attributes(pdb)
```

```
## $names
## [1] "atom" "xyz" "seqres" "helix" "sheet" "calpha" "remark" "call"
##
## $class
## [1] "pdb" "sse"
```

```
head(pdb$atom)
```

```
##   type eleno elety alt resid chain resno insert      x      y      z o      b
## 1 ATOM     1     N <NA>  PRO     A      1 <NA> 29.361 39.686 5.862 1 38.10
## 2 ATOM     2    CA <NA>  PRO     A      1 <NA> 30.307 38.663 5.319 1 40.62
## 3 ATOM     3     C <NA>  PRO     A      1 <NA> 29.760 38.071 4.022 1 42.64
## 4 ATOM     4     O <NA>  PRO     A      1 <NA> 28.600 38.302 3.676 1 43.40
## 5 ATOM     5    CB <NA>  PRO     A      1 <NA> 30.508 37.541 6.342 1 37.87
## 6 ATOM     6    CG <NA>  PRO     A      1 <NA> 29.296 37.591 7.162 1 38.40
##   segid elesy charge
## 1 <NA>     N  <NA>
## 2 <NA>     C  <NA>
## 3 <NA>     C  <NA>
## 4 <NA>     O  <NA>
## 5 <NA>     C  <NA>
## 6 <NA>     C  <NA>
```

Q7: How many amino acid residues are there in this pdb object?

```
PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYD QILIE-
ICGHKAIGTVLVGPTPVNIIGRNLLTQIGCTLNFPQITLWQRPLVTIKIGGQLKE ALLDTGAD-
DTVLEEMSLPGRWKPKMIGGIGGFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGRNLLTQIGCTLNF
```

198 amino acid residues

Q8: Name one of the two non-protein residues?

HOH

Q9: How many protein chains are in this structure?

2 chains

Q10. Which of the packages above is found only on BioConductor and not CRAN?

2 of them the biocManager and msa

Q11. Which of the above packages is not found on BioConductor or CRAN?:

bio3D

Q12. True or False? Functions from the devtools package can be used to install packages from GitHub and BitBucket?

True