

Class10

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Comparative structure analysis of ADK

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

msa

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

bio3d-view

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

True

Search and Retrieve ADK structures

```
library(bio3d)

aa <- get.seq("lake_A")
```

Warning in get.seq("lake_A"): Removing existing file: seqs.fasta

Fetching... Please wait. Done.

```
aa
```

```

      1      .      .      .      .      .      .      60
pdb|1AKE|A  MRIILLGAPGAGKGTQAQFIMEKYGIPQISTGDMLRAAVKSGSELGKQAKDIMDAGKLV
      1      .      .      .      .      .      .      60

      61      .      .      .      .      .      .      120
pdb|1AKE|A  DELVIALVKERIAQEDCRNGFLLDGFRTIPQADAMKEAGINVDYVLEFDVPDELIVDRI
      61      .      .      .      .      .      .      120

      121      .      .      .      .      .      .      180
pdb|1AKE|A  VGRRVHAPSGRVYHVKNPPKVEGKDDVTGEELTTRKDDQEETVRKRLVEYHQMTPALIG
      121      .      .      .      .      .      .      180

      181      .      .      .      214
pdb|1AKE|A  YYSKEAEAGNTKYAKVDGTPVAEVRADLEKILG
      181      .      .      .      214

```

Call:

```
read.fasta(file = outfile)
```

Class:

```
fasta
```

Alignment dimensions:

```
1 sequence rows; 214 position columns (214 non-gap, 0 gap)
```

```
+ attr: id, ali, call
```

Q13. How many amino acids are in this sequence, i.e. how long is this sequence?

214

We can run BLAST with this sequence:

```

# Blast or hmmer search
# b <- blast.pdb(aa)

# hits <- plot(b)

```

Let's see what is in our `hits` object.

```

# hits$ pdb.id

hits <- NULL

```

```
hits$pdb.id <- c('1AKE_A', '6S36_A', '6RZE_A', '3HPR_A', '1E4V_A', '5EJE_A', '1E4Y_A', '3X2S_A', '1E4Z_A')
```

Now we can download all of these PDB structures:

```
# Download related PDB files
files <- get.pdb(hits$pdb.id, path="pdbs", split=TRUE, gzip=TRUE)
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/1AKE.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/6S36.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/6RZE.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/3HPR.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/1E4V.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/5EJE.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/1E4Y.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/3X2S.pdb exists. Skipping download
```

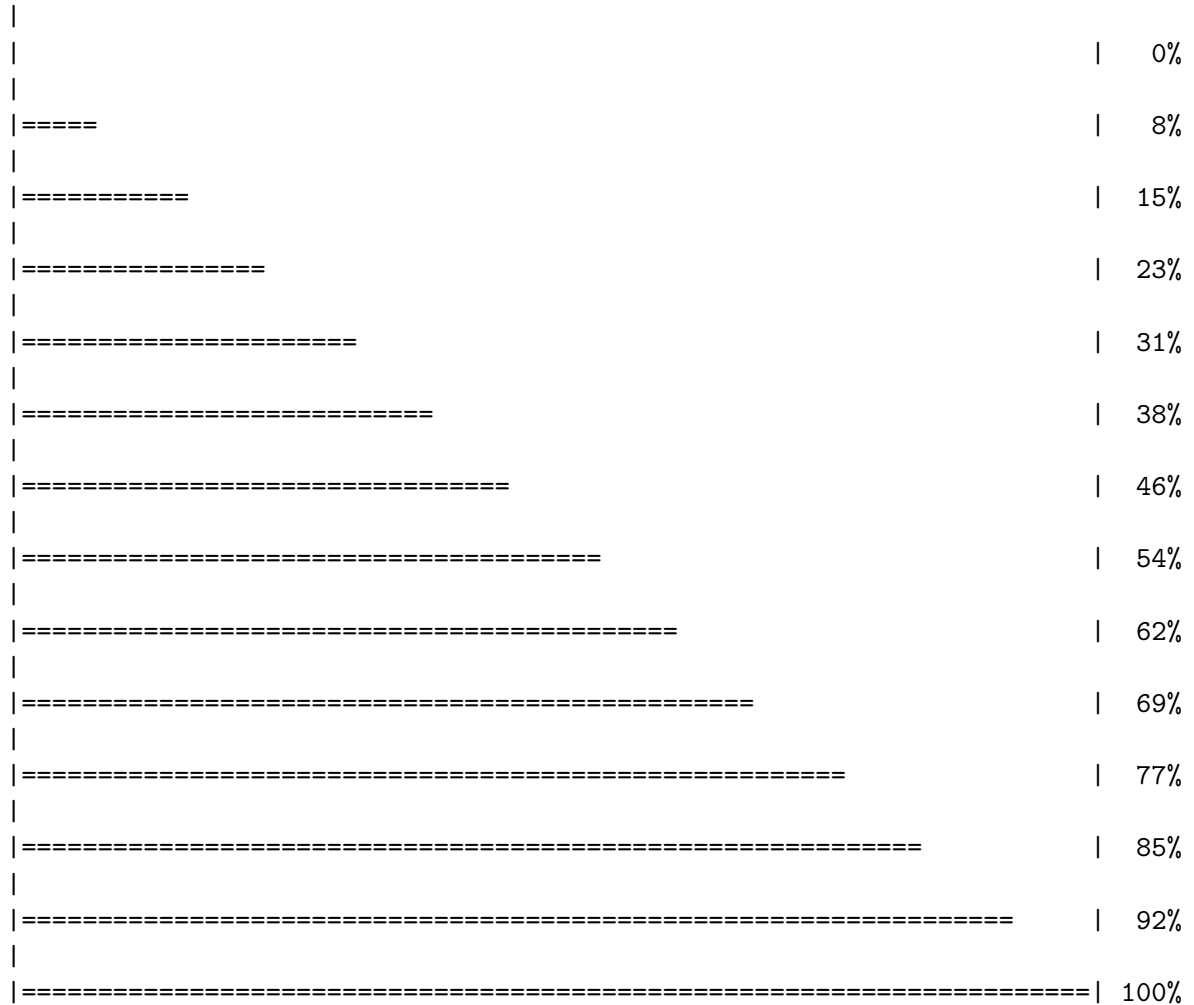
```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/6HAP.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/6HAM.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):
pdbs/4K46.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):  
pdbs/3GMT.pdb exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE):  
pdbs/4PZL.pdb exists. Skipping download
```



Now I want to align and superpose these structures which are all over the place.

```
pdbs <- pdbaln(files, fit=TRUE, exefile="msa")
```

Reading PDB files:

```

pdbs/split_chain/1AKE_A.pdb
pdbs/split_chain/6S36_A.pdb
pdbs/split_chain/6RZE_A.pdb
pdbs/split_chain/3HPR_A.pdb
pdbs/split_chain/1E4V_A.pdb
pdbs/split_chain/5EJE_A.pdb
pdbs/split_chain/1E4Y_A.pdb
pdbs/split_chain/3X2S_A.pdb
pdbs/split_chain/6HAP_A.pdb
pdbs/split_chain/6HAM_A.pdb
pdbs/split_chain/4K46_A.pdb
pdbs/split_chain/3GMT_A.pdb
pdbs/split_chain/4PZL_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
..  PDB has ALT records, taking A only, rm.alt=TRUE
.... PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
...

```

Extracting sequences

```

pdb/seq: 1   name: pdbs/split_chain/1AKE_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 2   name: pdbs/split_chain/6S36_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 3   name: pdbs/split_chain/6RZE_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 4   name: pdbs/split_chain/3HPR_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 5   name: pdbs/split_chain/1E4V_A.pdb
pdb/seq: 6   name: pdbs/split_chain/5EJE_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 7   name: pdbs/split_chain/1E4Y_A.pdb
pdb/seq: 8   name: pdbs/split_chain/3X2S_A.pdb
pdb/seq: 9   name: pdbs/split_chain/6HAP_A.pdb
pdb/seq: 10  name: pdbs/split_chain/6HAM_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 11  name: pdbs/split_chain/4K46_A.pdb
    PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 12  name: pdbs/split_chain/3GMT_A.pdb

```

pdb/seq: 13 name: pdbs/split_chain/4PZL_A.pdb

Let's have a look at our pdbs object.

pdbs

```
[Truncated_Name:1] 1AKE_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:2] 6S36_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:3] 6RZE_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:4] 3HPR_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:5] 1E4V_A.pdb      1      .      .      .      40
-----MRIILLGAPVAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:6] 5EJE_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:7] 1E4Y_A.pdb      1      .      .      .      40
-----MRIILLGALVAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:8] 3X2S_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:9] 6HAP_A.pdb      1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:10] 6HAM_A.pdb     1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:11] 4K46_A.pdb     1      .      .      .      40
-----MRIILLGAPGAGKGTQAQFIMAKFGIPQIS
[Truncated_Name:12] 3GMT_A.pdb     1      .      .      .      40
-----MRLILLGAPGAGKGTQANFIKEKFGIPQIS
[Truncated_Name:13] 4PZL_A.pdb     1      .      .      .      40
TENLYFQSNAMRIILLGAPGAGKGTQAKIIEQKYNIAHIS
                **~*****  *****  *  *~ *  **
1      .      .      .      40

41      .      .      .      80
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDACKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDCGKLVTDLVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVRE
TGDMLRAAIAKSGSELGKQAKDIMDAGKLVTDLIIIALVKE
TGDMLRAAIAKAGTELGKQAKSVIDAGQLVSDDIILGLVKE
TGDMLRAAVKAGTPLGVEAKTYMDEGKLPVDSLIIGLVKE
TGDMIRETIKSGSALGQELKKVLDAGELVSDEFIIVKIVKD
****~*  ~* *~ **  *  ~*  ** *  ^^ ~*^^
41      .      .      .      80

81      .      .      .      120
RIAQEDCRNGFLLDGFPRTPQADAMKEAGINVDYVLEFD
```

[Truncated_Name:2] 6S36_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:3] 6RZE_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:4] 3HPR_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:5] 1E4V_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:6] 5EJE_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:7] 1E4Y_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:8] 3X2S_A.pdb	RIAQEDSRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:9] 6HAP_A.pdb	RICQEDSRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:10] 6HAM_A.pdb	RICQEDSRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:11] 4K46_A.pdb	RIAQDDCAKGFLLDGFPR TIPQADGLKEVG VVVVDYVIEFD
[Truncated_Name:12] 3GMT_A.pdb	RLKEADCANGYLF DGFPR TIAQADAMKEAGVAIDYVLEID
[Truncated_Name:13] 4PZL_A.pdb	RISKNDCNNGFLLDGVPR TIPQAQELDKLGVNIDYIVEVD
	*^ * *~* ** ***** ** ^ *^ ~**~* *
	81 . . . 120
	121 . . . 160
[Truncated_Name:1] 1AKE_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:2] 6S36_A.pdb	VPDELIVDKIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:3] 6RZE_A.pdb	VPDELIVDAIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:4] 3HPR_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDGTG
[Truncated_Name:5] 1E4V_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:6] 5EJE_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:7] 1E4Y_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:8] 3X2S_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:9] 6HAP_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:10] 6HAM_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:11] 4K46_A.pdb	VADSVIVERMAGRR AHLASGR TYHN VYNPPKVEGKDDVTG
[Truncated_Name:12] 3GMT_A.pdb	VPFSEIIERMSGRR THPASGR TYHV KFNPPKVEGKDDVTG
[Truncated_Name:13] 4PZL_A.pdb	VADNLLIERITGRRIHPASGR TYHTKF NPPKVADKDDVTG
	* ^^^ ^ *** * *** ** ^***** *** **
	121 . . . 160
	161 . . . 200
[Truncated_Name:1] 1AKE_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:2] 6S36_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:3] 6RZE_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:4] 3HPR_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:5] 1E4V_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:6] 5EJE_A.pdb	EELTTRKDDQEECVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:7] 1E4Y_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:8] 3X2S_A.pdb	EELTTRKDDQEETVRKRLCEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:9] 6HAP_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN
[Truncated_Name:10] 6HAM_A.pdb	EELTTRKDDQEETVRKRLVEYHQMTAP LIGYYSKEAEAGN

```

[Truncated_Name:11]4K46_A.pdb    EDLVIREDDKEETVRLARLGVYHNQTAPLIAYYGKEAEAGN
[Truncated_Name:12]3GMT_A.pdb    EPLVQRDDDDKEETVKKRLDVYEAQTKPLITYYGDWARRGA
[Truncated_Name:13]4PZL_A.pdb    EPLITRTDDNEDTVKQRLSVYHAQTAKLIDFYRNFSSNTNT
                                * * * ** *~ * ** * * ** ^*
                                161 . . . 200

                                201 . . 227
[Truncated_Name:1]1AKE_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:2]6S36_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:3]6RZE_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:4]3HPR_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:5]1E4V_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:6]5EJE_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:7]1E4Y_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:8]3X2S_A.pdb     T--KYAKVDGTPVAEVRADLEKILG-
[Truncated_Name:9]6HAP_A.pdb     T--KYAKVDGTPVCEVRADLEKILG-
[Truncated_Name:10]6HAM_A.pdb     T--KYAKVDGTPVCEVRADLEKILG-
[Truncated_Name:11]4K46_A.pdb     T--QYLKFDGTPKAVAEVSAELEKALA-
[Truncated_Name:12]3GMT_A.pdb     E-----NGLKAPA-----YRKISG-
[Truncated_Name:13]4PZL_A.pdb     KIPKYIKINGDQAVEKVSQDIFDQLNK
                                *
                                201 . . 227

```

Call:

```
pdbaln(files = files, fit = TRUE, exefile = "msa")
```

Class:

```
pdbs, fasta
```

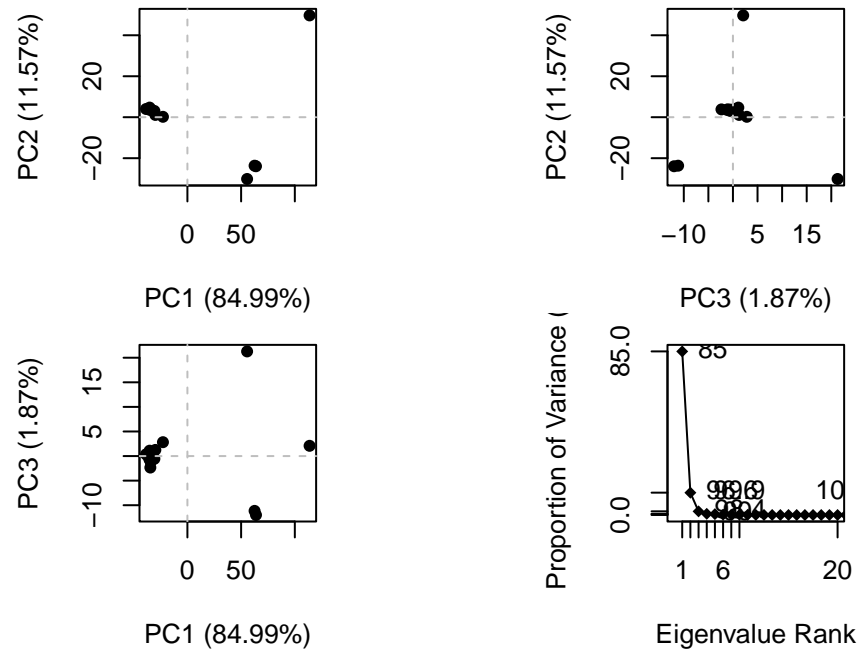
Alignment dimensions:

```
13 sequence rows; 227 position columns (204 non-gap, 23 gap)
```

```
+ attr: xyz, resno, b, chain, id, ali, resid, sse, call
```

Now we have our aligned and supersposed structures we can perform all sorts of analysis on them. Let's do PCA...

```
pc.xray <- pca(pdbs)
plot(pc.xray)
```

Results of PCA on Adenylate kinase X-ray structures. Each dot represents one PDB structure.

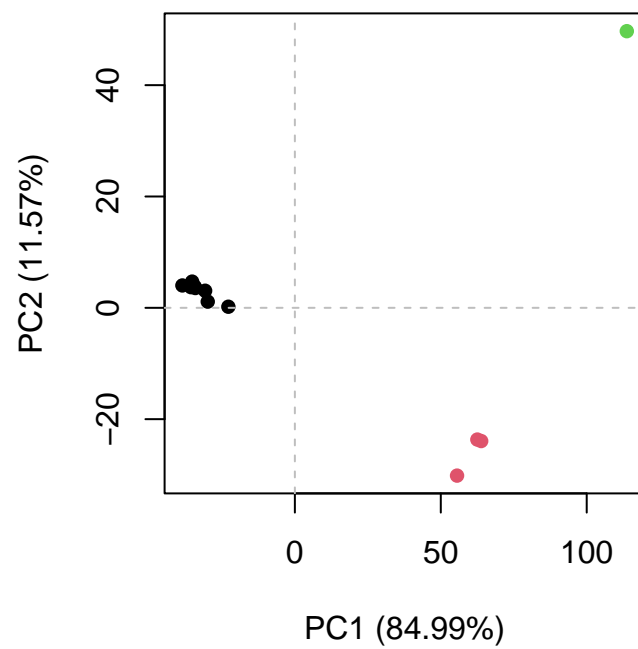
We can cluster the structures by RMSD (or any other method).

```
rd <- rmsd(pdbbs)
```

Warning in rmsd(pdbbs): No indices provided, using the 204 non NA positions

```
hc.rd <- hclust(dist(rd))
grps <- cutree(hc.rd, k=3)

plot(pc.xray, 1:2, col=grps)
```



Optional further visualization

We can make a wee movie - also called a trajectory of the major differences (i.e. structural displacements) of ADK.

```
# Visualizae first principal component  
pc1 <- mktrj(pc.xray, pc=1, file="pc_1.pdb")
```