

Class10 –Comparative structure analysis of Adenylate Kinase

Eric Garcia Ordunez (A16316409)

Comparative analysis of ADK

ADK (adenylate Kinase) is an important drug target and we would love to know how it works - i.e. molecular structure

There has been lots of work done on this protein due to its importance including lots of crystal structures.

We will begin with getting an example ADK sequence

Q 10: 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?

True

```
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  VGRRVHAPSGRVYHVKNPPKVEGKDDVTGEELTTRKDDQEETVRKRLVEYHQMTAPLIG
      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 now run BLAST with this sequence

```
#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',
```

Now we can download all these PDB structure files:

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

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

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

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

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

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

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

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

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

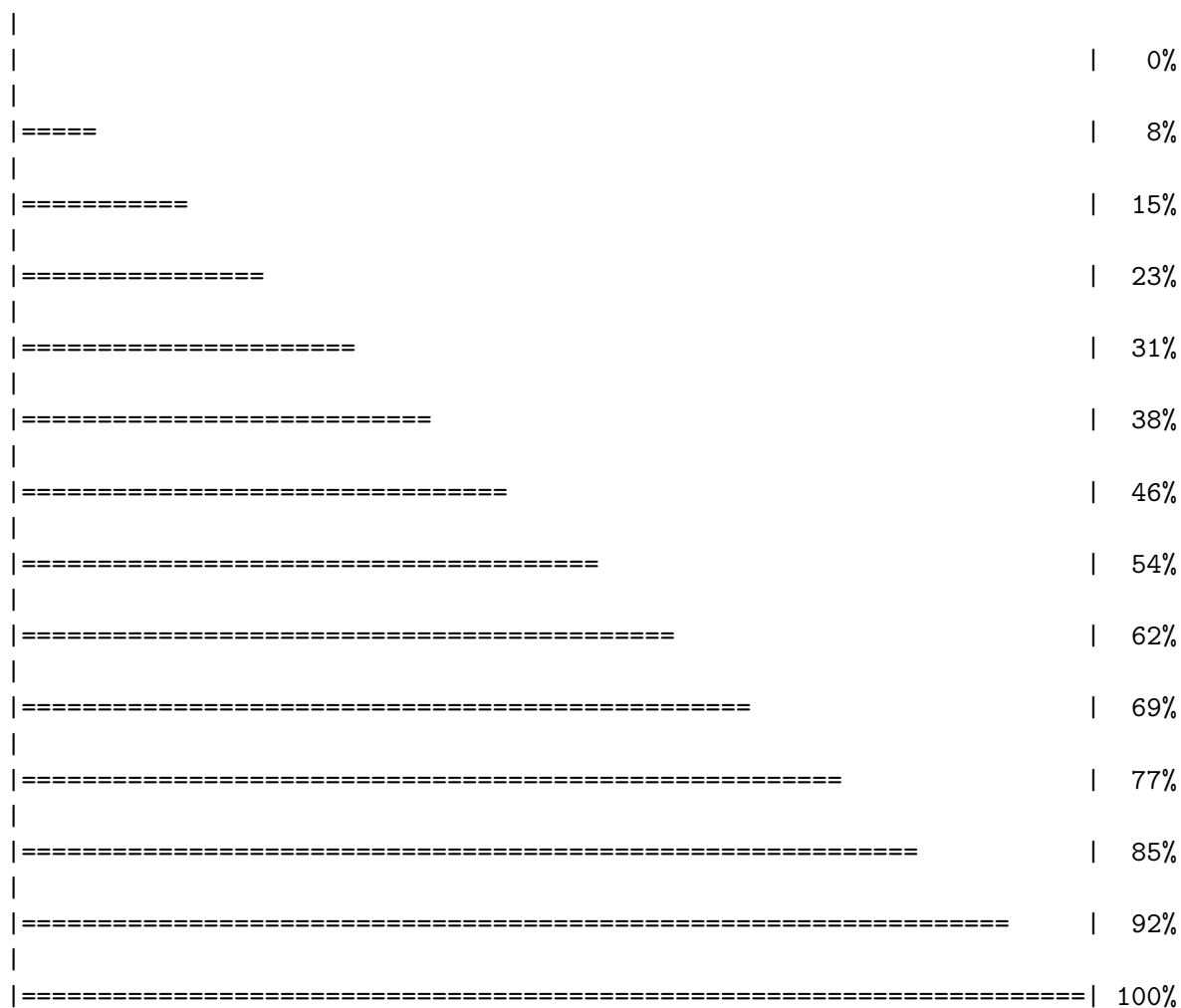
```
Warning in get.pdb(hits$pdb.id, path = "pdb", split = TRUE, gzip = TRUE):
pdb/6HAP.pdb.gz exists. Skipping download
```

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

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

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

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



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

```
# Align related PDBs
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
```



```

****~*  ~* *~ **  *  ~*  ** *  ^^ ~*~^
41      .      .      .      80

81      .      .      .      120
[Truncated_Name:1] 1AKE_A.pdb  RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[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  RISKNCNNGFLLDGVPR TIPQAQELDKLG VNI DYIVEVD
*~  *  *~* ** ***** **  ^  *~ ~**~*~* *
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 TYHNVYNPPKVEGKDDVTG
[Truncated_Name:12] 3GMT_A.pdb  VPFSEIIERMSGRR THPASGR TYHV KFNPPKVEGKDDVTG
[Truncated_Name:13] 4PZL_A.pdb  VADNLLIERITGRRIHPASGR TYHTKFNPPKVADKDDVTG
*  ^^^ ^ *** *  *** **  ^***** *** **
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      EELTTRKDDQEECVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:7] 1E4Y_A.pdb      EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:8] 3X2S_A.pdb      EELTTRKDDQEETVRKRLCEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:9] 6HAP_A.pdb      EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:10] 6HAM_A.pdb      EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:11] 4K46_A.pdb      EDLVIREDDKEETVLARLGVYHNQ TAPLIAYYGKEAEAGN
[Truncated_Name:12] 3GMT_A.pdb      EPLVQRDDDKKEETVKKRLDVYEA QTKPLITYYGDWARRGA
[Truncated_Name:13] 4PZL_A.pdb      EPLITRTDDNEDTVKQRLSVYHAQ TAKLIDFYRNFSSNTNT
                                     * * * * * ^ * * * * * ^ *
                                     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
```

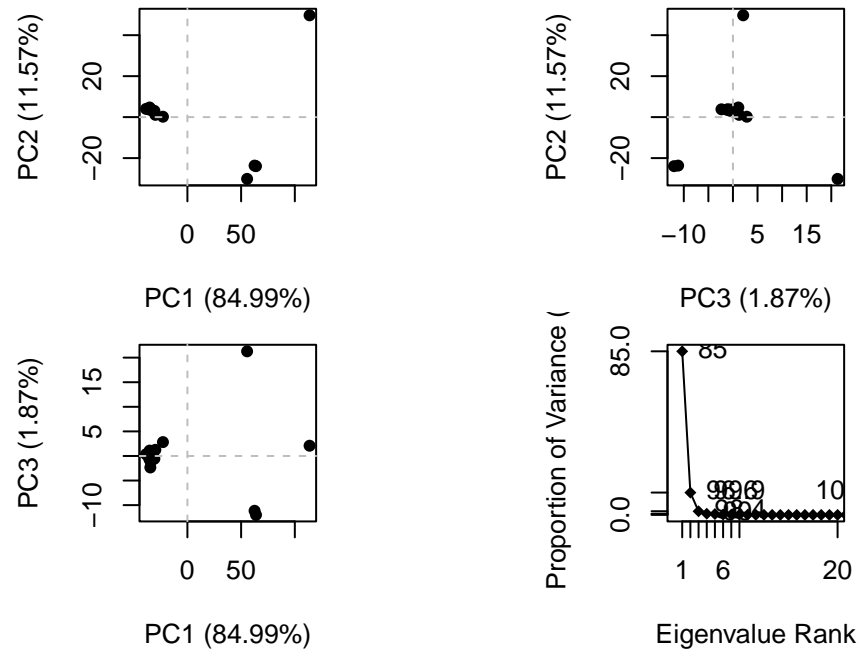
Alignment dimensions:

```
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 superposed we can perform all sorts of analysis on them. Lets do PCA...


```
# Perform PCA
pc.xray <- pca(pdbbs)
plot(pc.xray)
```



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

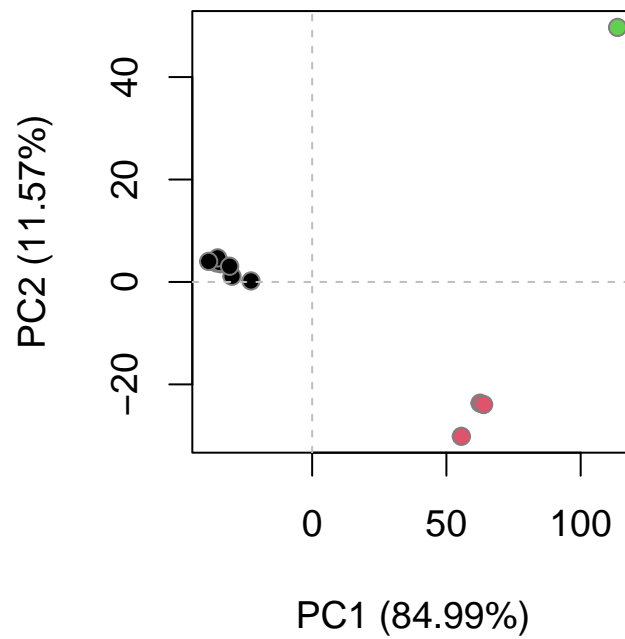
We can cluster the structures by RMSD (or any other method).

```
# Calculate RMSD
rd <- rmsd(pdbbs)
```

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

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

plot(pc.xray, 1:2, col="grey50", bg=grps.rd, pch=21, cex=1)
```



This is going to take the PCA and make a movie- also called a trajectory of the major difference (i.e. structural displacements of ADK)

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