

# Class 10: Structural BioInformatics II

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## Comparative Analysis of ADK

ADK (adenelate Kinase) is an important drug target and we would love to know how it worked - i.e. molecular mechanism

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

We will begin getting an example ADK sequence from the database. We will then use this to find all ADK structures in the PDB.

```
library(bio3d)
aa <- get.seq("1ake_A")
```

Warning in get.seq("1ake\_A"): Removing existing file: seqs.fasta

Fetching... Please wait. Done.

```
aa
```

```

      1      .      .      .      .      .      60
pdb|1AKE|A  MRIILLGAPGAGKGTQAQFIMEKYGIPQISTGMDLRAAVKSGSELGKQAKDIMDAGKLV
      1      .      .      .      .      .      60
      61      .      .      .      .      .      120
pdb|1AKE|A  DELVIALVKERIAQEDCRNGFLLDGFPRTIPQADAMKEAGINVDYVLEFDVPDELIVDRI
      61      .      .      .      .      .      120
      121      .      .      .      .      .      180
pdb|1AKE|A  VGRRVHAPSGRVYHVKFNPPKVEGKDDVTGEELTTRKDDQEETVRKRLVEYHQMTAPLIG
```



Let's see what is in our hits objects

```
# List out some 'top hits'
#head(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="pdbc", split=TRUE, gzip=TRUE)
```

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

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

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

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

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

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

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

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

Warning in get.pdb(hits\$ pdb.id, path = "pdbc", split = TRUE, gzip = TRUE):  
pdbc/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

	0%
=====	8%
=====	15%
=====	23%
=====	31%
=====	38%
=====	46%
=====	54%
=====	62%
=====	69%
=====	77%
=====	85%
=====	92%
=====	100%

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
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
[Truncated_Name:1] 1AKE_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
[Truncated_Name:2] 6S36_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
[Truncated_Name:3] 6RZE_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
[Truncated_Name:4] 3HPR_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
[Truncated_Name:5] 1E4V_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
[Truncated_Name:6] 5EJE_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDACKLVDELVIALVKE
[Truncated_Name:7] 1E4Y_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
[Truncated_Name:8] 3X2S_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDCGLVDELVIALVKE
[Truncated_Name:9] 6HAP_A.pdb      TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVRE
[Truncated_Name:10] 6HAM_A.pdb     TGDMLRAAIAKSGSELGKQAKDIMDAGKLVDEIIIALVKE
[Truncated_Name:11] 4K46_A.pdb     TGDMLRAAIAKAGTELGKQAKSVIDAGQLVSDDIILGLVKE

```

[Truncated_Name:12] 3GMT_A.pdb	TGDM	LRAAVKAGTPLGVEAKTYMDEGKLVPDSLIIGLVKE
[Truncated_Name:13] 4PZL_A.pdb	TGDM	IRETIKSGSALGQELKKVLDAGELVSDEFIIKIVKD
	****~*    ^* *^ **    *    ^*    ** *    ^^ ~*^^	
	41	80
	81	120
[Truncated_Name:1] 1AKE_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:2] 6S36_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:3] 6RZE_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:4] 3HPR_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:5] 1E4V_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:6] 5EJE_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:7] 1E4Y_A.pdb	RIAQEDCRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:8] 3X2S_A.pdb	RIAQEDSRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:9] 6HAP_A.pdb	RICQEDSRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:10] 6HAM_A.pdb	RICQEDSRNGFLLDGF	PRTIPQADAMKEAGINVDYVLEFD
[Truncated_Name:11] 4K46_A.pdb	RIAQDDCAKGFLLDGF	PRTIPQADGLKEVGVVVDYVIEFD
[Truncated_Name:12] 3GMT_A.pdb	RLKEADCANGYLF	DGFRTIAQADAMKEAGVAIDYVLEID
[Truncated_Name:13] 4PZL_A.pdb	RISKNCNNGFLLDGF	PRTIPQAQELDKLGVNIDYIVEVD
	*^    *    *~* ** ***** **    ^    *^ ~***~* *    *	
	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	VADSVIVERMAGRRAHLASGR	TYHNVYNPPKVEGKDDVTG
[Truncated_Name:12] 3GMT_A.pdb	VPFSEIIERMSGRRTHPASGR	TYHVKNPPKVEGKDDVTG
[Truncated_Name:13] 4PZL_A.pdb	VADNLLIERITGRRIH	PASGRTYHTKFNPPKVADKDDVTG
	*    ^^^ ^    *** *    *** **    ^*****    *** **	
	121	160
	161	200
[Truncated_Name:1] 1AKE_A.pdb	EELTTRKDDQEETVRKRLVEYH	QMTAPLIGYYSKEAEAGN
[Truncated_Name:2] 6S36_A.pdb	EELTTRKDDQEETVRKRLVEYH	QMTAPLIGYYSKEAEAGN
[Truncated_Name:3] 6RZE_A.pdb	EELTTRKDDQEETVRKRLVEYH	QMTAPLIGYYSKEAEAGN

```

[Truncated_Name:4] 3HPR_A.pdb      EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:5] 1E4V_A.pdb      EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[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      EDLVIREDDKEETVRLARLGVYHNQTAPLIAYYGKEAEAGN
[Truncated_Name:12] 3GMT_A.pdb      EPLVQRDDKEETVKKRLDVYEAQTKPLITYYGDWARRGA
[Truncated_Name:13] 4PZL_A.pdb      EPLITRTDDNEDTVKQRLSVYHAQTAKLIDFYRNFSSNT
                                     * * * * * ^ * * * * * ^ *
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:

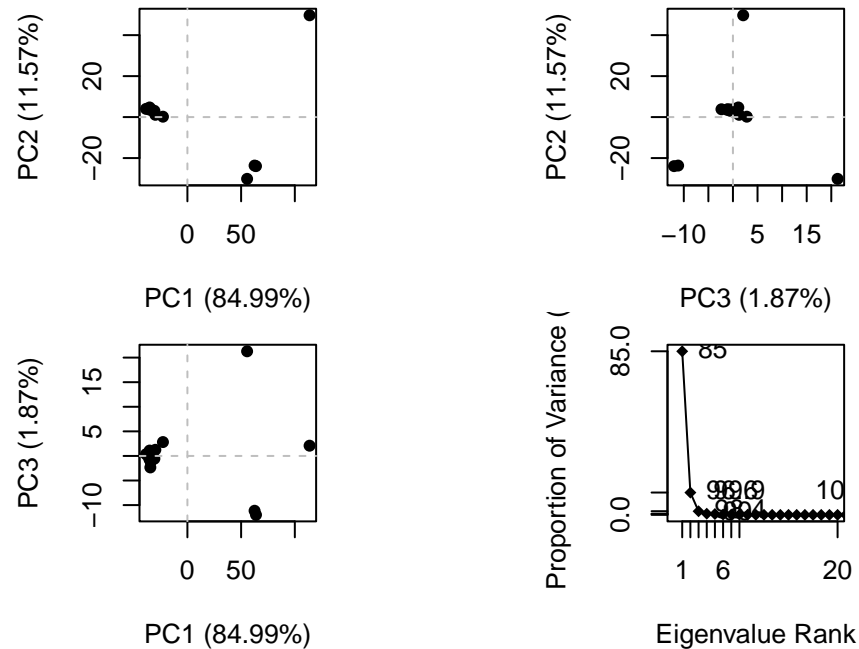
```
13 sequence rows; 227 position columns (204 non-gap, 23 gap)
```

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

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

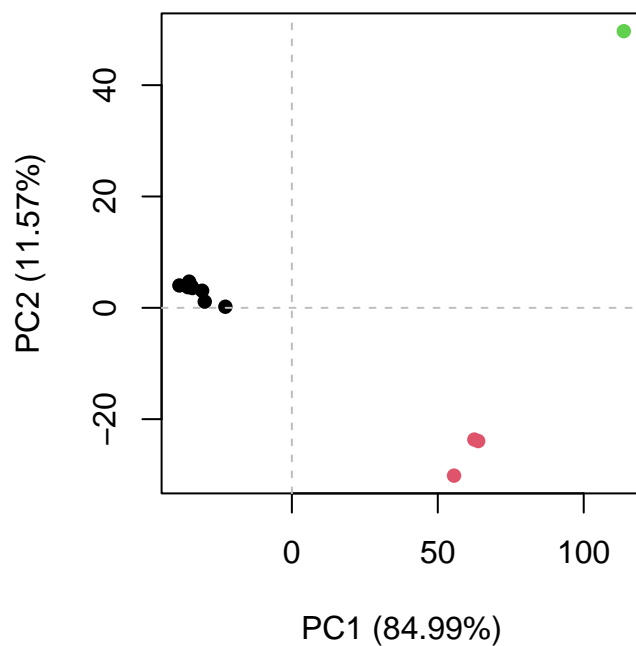
We can cluster the structures by RMSD (or any other method). Function `rmsd()` will calculate all pairwise RMSD values of the structural ensemble. This facilitates clustering analysis based on the pairwise structural deviation:

```
# 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 <- 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.

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

```
Total Frames#: 34
Total XYZs#: 612, (Atoms#: 204)
```

```
[1] 26.787 52.261 40.414 <...> 15.653 53.622 42.018 [20808]
```

```
+ attr: Matrix DIM = 34 x 612
```