

Lab10_inclass

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ADK (Adenelate Kinase) important drug target, know its molecular mechanism

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

```
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  MRIILLGAPGAGKGTQAQFIMEKYGIPQISTGDMRLRAAVKSGSELGKQAKDIMDAGKLV
      1      .      .      .      .      .      .      60

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

     121      .      .      .      .      .      .      180
pdb|1AKE|A  VGRRVHAPSGRVYHVKNPPKVEGKDDVTGEELTTRKDDQEETVRKRLVEYHQM
     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
```

We can now run BLAST with this sequence

```
#b <- blast.pdb(aa)
```

```
#same as plot.blast(b)
```

```
#hits <- plot(b)
```

Let's see what is in our hits object

```
#hits$pdb.id
```

Now we can download all these top hits PDB structure files:

```
# Download related PDB files
```

```
#?get.pdb
```

```
hits <- NULL
```

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

```
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 = "pdbs", split = TRUE, gzip = TRUE):
pdbs/3HPR.pdb.gz exists. Skipping download

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

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

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

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

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

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

	0%
=====	8%
=====	15%
=====	23%



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

```
# Align related PDBs; fit on top of each other, exefile run alignment on the same package
pdbbs <- pdbaln(files, fit = TRUE, exefile="msa")
```

Reading PDB files:

```
pdbbs/split_chain/1AKE_A.pdb
pdbbs/split_chain/6S36_A.pdb
pdbbs/split_chain/6RZE_A.pdb
pdbbs/split_chain/3HPR_A.pdb
pdbbs/split_chain/1E4V_A.pdb
pdbbs/split_chain/5EJE_A.pdb
pdbbs/split_chain/1E4Y_A.pdb
pdbbs/split_chain/3X2S_A.pdb
pdbbs/split_chain/6HAP_A.pdb
pdbbs/split_chain/6HAM_A.pdb
pdbbs/split_chain/4K46_A.pdb
pdbbs/split_chain/3GMT_A.pdb
pdbbs/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
```

	1	.	.	.	40
[Truncated_Name:1] 1AKE_A.pdb	-----	MRIILLGAPGAGKGTQAQFIMEKYGIPQIS			
[Truncated_Name:2] 6S36_A.pdb	-----	MRIILLGAPGAGKGTQAQFIMEKYGIPQIS			
[Truncated_Name:3] 6RZE_A.pdb	-----	MRIILLGAPGAGKGTQAQFIMEKYGIPQIS			
[Truncated_Name:4] 3HPR_A.pdb	-----	MRIILLGAPGAGKGTQAQFIMEKYGIPQIS			
[Truncated_Name:5] 1E4V_A.pdb	-----	MRIILLGAPVAGKGTQAQFIMEKYGIPQIS			
[Truncated_Name:6] 5EJE_A.pdb	-----	MRIILLGAPGAGKGTQAQFIMEKYGIPQIS			
[Truncated_Name:7] 1E4Y_A.pdb	-----	MRIILLGALVAGKGTQAQFIMEKYGIPQIS			

```

[Truncated_Name:8] 3X2S_A.pdb -----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:9] 6HAP_A.pdb -----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:10] 6HAM_A.pdb -----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:11] 4K46_A.pdb -----MRIILLGAPGAGKGTQAQFIMAKFGIPQIS
[Truncated_Name:12] 3GMT_A.pdb -----MRLILLGAPGAGKGTQANFIKEKFGIPQIS
[Truncated_Name:13] 4PZL_A.pdb TENLYFQSNAMRIILLGAPGAGKGTQAKIIEQKYNIAHIS
                                **~*****  *  *~ *  **
1                                .  .  . 40

41                                .  .  . 80
[Truncated_Name:1] 1AKE_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
[Truncated_Name:2] 6S36_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
[Truncated_Name:3] 6RZE_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
[Truncated_Name:4] 3HPR_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
[Truncated_Name:5] 1E4V_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
[Truncated_Name:6] 5EJE_A.pdb TGDMLRAAVKSGSELGKQAKDIMDACKLVTDLVIALVKE
[Truncated_Name:7] 1E4Y_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVKE
[Truncated_Name:8] 3X2S_A.pdb TGDMLRAAVKSGSELGKQAKDIMDCGKLVTDLVIALVKE
[Truncated_Name:9] 6HAP_A.pdb TGDMLRAAVKSGSELGKQAKDIMDAGKLVTDLVIALVRE
[Truncated_Name:10] 6HAM_A.pdb TGDMLRAAIIKSGSELGKQAKDIMDAGKLVTDLIIIALVKE
[Truncated_Name:11] 4K46_A.pdb TGDMLRAAIIKAGTELGKQAKSVIDAGQLVSDDIILGLVKE
[Truncated_Name:12] 3GMT_A.pdb TGDMLRAAVKAGTPLGVEAKTYMDEGKLPVDSLIIIGLVKE
[Truncated_Name:13] 4PZL_A.pdb TGDMIRETIKSGSALGQELKKVLDAGELVSDEFIIKIVKD
****~*  ~* *~ **  *  ~*  ** *  ~ ~ ~*~
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 TIPQADGLKEVGVVVDYVIEFD
[Truncated_Name:12] 3GMT_A.pdb RLKEADCANGYLFDFPRTIAQADAMKEAGVAIDYVLEID
[Truncated_Name:13] 4PZL_A.pdb RISKNDCNNGFLLDGVPR TIPQAQELDKLGVNIDYIVEVD
*~  *  *~* ** ***** **  ^  *~ ~*~*~*
81                                .  .  . 120

```

	121	.	.	.	160
[Truncated_Name:1] 1AKE_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:2] 6S36_A.pdb	VPDELIVDKIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:3] 6RZE_A.pdb	VPDELIVDAIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:4] 3HPR_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDGTG				
[Truncated_Name:5] 1E4V_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:6] 5EJE_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:7] 1E4Y_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:8] 3X2S_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:9] 6HAP_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:10] 6HAM_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHVKNPPKVEGKDDVTG				
[Truncated_Name:11] 4K46_A.pdb	VADSVIVERMAGRRAHLASGRTYHNVNPPKVEGKDDVTG				
[Truncated_Name:12] 3GMT_A.pdb	VPFSEIIERMSGRRTHPASGRTYHVKNPPKVEGKDDVTG				
[Truncated_Name:13] 4PZL_A.pdb	VADNLLIERITGRIHPASGRTYHTKFNPPKVADKDDVTG				
	* ^^^ ^ *** * *** ** ^***** *** **				
	121	.	.	.	160
	161	.	.	.	200
[Truncated_Name:1] 1AKE_A.pdb	EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN				
[Truncated_Name:2] 6S36_A.pdb	EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN				
[Truncated_Name:3] 6RZE_A.pdb	EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN				
[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	EDLVIREDDKEETV LARLG VYHNQTAPLIAYYGKEAEAGN				
[Truncated_Name:12] 3GMT_A.pdb	EPLVQRDDDK EETVKKRLDVYEAQTKPLITYYGDWARRGA				
[Truncated_Name:13] 4PZL_A.pdb	EPLITRTDDNEDTVKQRLSVYHAQTAKLIDFYRNFSSTNT				
	* * * * * ^ * * * * * * * ^ *				
	161	.	.	.	200
	201	.	.	227	
[Truncated_Name:1] 1AKE_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:2] 6S36_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:3] 6RZE_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:4] 3HPR_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:5] 1E4V_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:6] 5EJE_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:7] 1E4Y_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				
[Truncated_Name:8] 3X2S_A.pdb	T--KYAKVDG TKPVAEVRADLEKILG-				

```

[Truncated_Name:9]6HAP_A.pdb      T--KYAKVDGTPVCEVRADLEKILG-
[Truncated_Name:10]6HAM_A.pdb     T--KYAKVDGTPVCEVRADLEKILG-
[Truncated_Name:11]4K46_A.pdb     T--QYLKFDGTKAVAEVSAELEKALA-
[Truncated_Name:12]3GMT_A.pdb     E-----NGLKAPA-----YRKISG-
[Truncated_Name:13]4PZL_A.pdb     KIPKYIKINGDQAVEKVSQDIFDQLNK
                                   *
                                   .           .           227
                                   201

```

Call:

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

Class:

```
pdbs, fasta
```

Alignment dimensions:

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

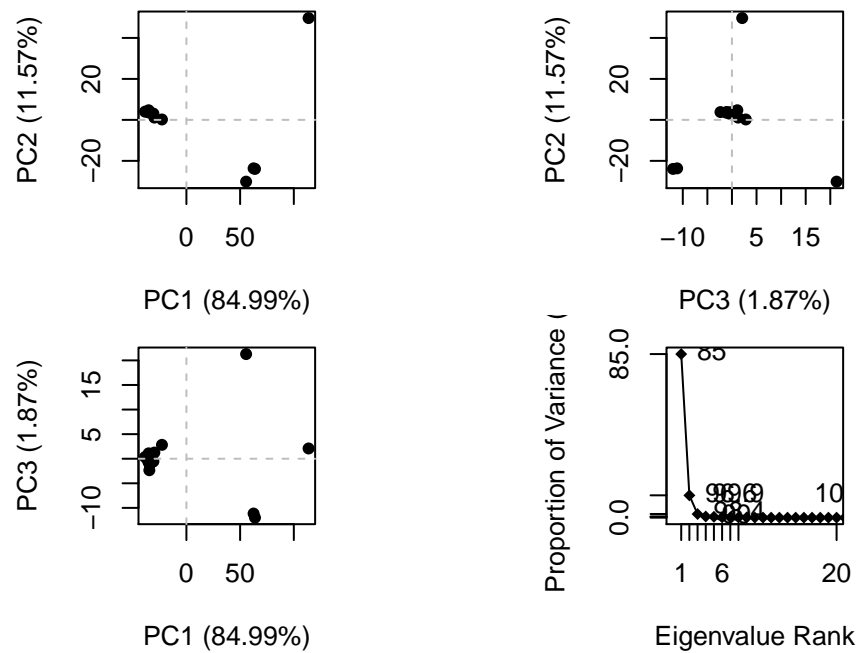
```
+ attr: xyz, resno, b, chain, id, ali, resid, sse, call
```

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(pdbs)
plot(pc.xray)

```

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

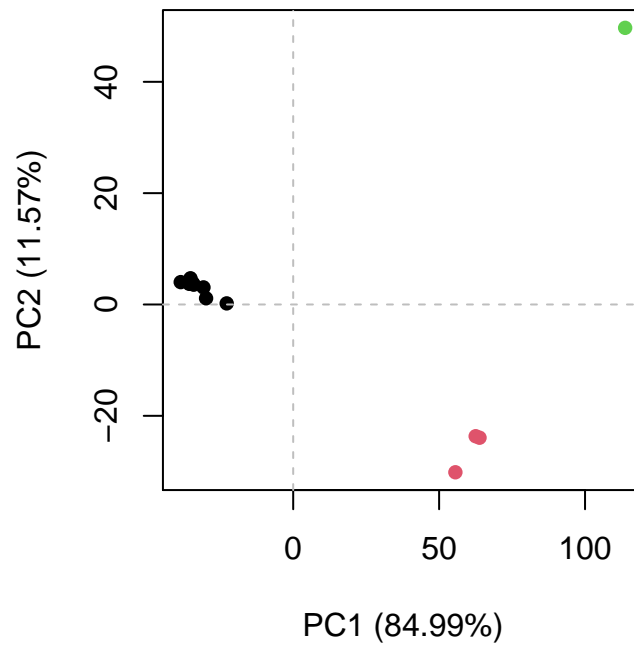
We can cluster the structures by RMSD (or any other method).

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

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

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

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



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

We can make a wee movie - a trajectory of the major differences of

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