

class11

2024-02-25

```
# Change this for YOUR results dir name
results_dir <- "HIVPRDIMER_23119/"

# File names for all PDB models
pdb_files <- list.files(path=results_dir,
                        pattern="*.pdb",
                        full.names = TRUE)

# Print our PDB file names
basename(pdb_files)
```

```
## [1] "HIVPRDIMER_23119_unrelaxed_rank_001_alphafold2_multimer_v3_model_5_seed_000.pdb"
## [2] "HIVPRDIMER_23119_unrelaxed_rank_002_alphafold2_multimer_v3_model_1_seed_000.pdb"
## [3] "HIVPRDIMER_23119_unrelaxed_rank_003_alphafold2_multimer_v3_model_4_seed_000.pdb"
## [4] "HIVPRDIMER_23119_unrelaxed_rank_004_alphafold2_multimer_v3_model_2_seed_000.pdb"
## [5] "HIVPRDIMER_23119_unrelaxed_rank_005_alphafold2_multimer_v3_model_3_seed_000.pdb"
```

```
library(bio3d)

# Read all data from Models
# and superpose/fit coords
pdbs <- pdbaln(pdb_files, fit=TRUE, exefile="msa")
```

```
## Reading PDB files:
## HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_001_alphafold2_multimer_v3_model_5_seed_000.pdb
## HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_002_alphafold2_multimer_v3_model_1_seed_000.pdb
## HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_003_alphafold2_multimer_v3_model_4_seed_000.pdb
## HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_004_alphafold2_multimer_v3_model_2_seed_000.pdb
## HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_005_alphafold2_multimer_v3_model_3_seed_000.pdb
## .....
##
## Extracting sequences
##
## pdb/seq: 1   name: HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_001_alphafold2_multimer_v3_model_5_seed_000.pdb
## pdb/seq: 2   name: HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_002_alphafold2_multimer_v3_model_1_seed_000.pdb
## pdb/seq: 3   name: HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_003_alphafold2_multimer_v3_model_4_seed_000.pdb
## pdb/seq: 4   name: HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_004_alphafold2_multimer_v3_model_2_seed_000.pdb
## pdb/seq: 5   name: HIVPRDIMER_23119//HIVPRDIMER_23119_unrelaxed_rank_005_alphafold2_multimer_v3_model_3_seed_000.pdb
```

pdbs

```
##                               1           .           .           .           .           50
```

```

## [Truncated_Name:1]HIVPRDIMER PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGI
## [Truncated_Name:2]HIVPRDIMER PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGI
## [Truncated_Name:3]HIVPRDIMER PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGI
## [Truncated_Name:4]HIVPRDIMER PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGI
## [Truncated_Name:5]HIVPRDIMER PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGI
##
## 1 . . . . 50
##
## 51 . . . . 100
## [Truncated_Name:1]HIVPRDIMER GGFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNFP
## [Truncated_Name:2]HIVPRDIMER GGFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNFP
## [Truncated_Name:3]HIVPRDIMER GGFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNFP
## [Truncated_Name:4]HIVPRDIMER GGFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNFP
## [Truncated_Name:5]HIVPRDIMER GGFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNFP
##
## 51 . . . . 100
##
## 101 . . . . 150
## [Truncated_Name:1]HIVPRDIMER QITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGIG
## [Truncated_Name:2]HIVPRDIMER QITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGIG
## [Truncated_Name:3]HIVPRDIMER QITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGIG
## [Truncated_Name:4]HIVPRDIMER QITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGIG
## [Truncated_Name:5]HIVPRDIMER QITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWPKMIGGIG
##
## 101 . . . . 150
##
## 151 . . . . 198
## [Truncated_Name:1]HIVPRDIMER GFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNF
## [Truncated_Name:2]HIVPRDIMER GFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNF
## [Truncated_Name:3]HIVPRDIMER GFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNF
## [Truncated_Name:4]HIVPRDIMER GFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNF
## [Truncated_Name:5]HIVPRDIMER GFIKVRQYDQILIEICGHKAIGTVLVGPTPVNIIGNLLTQIGCTLNF
##
## 151 . . . . 198

```

```

## Call:
##   pdbaln(files = pdb_files, fit = TRUE, exefile = "msa")
##
## Class:
##   pdba, fasta
##
## Alignment dimensions:
##   5 sequence rows; 198 position columns (198 non-gap, 0 gap)
##
## + attr: xyz, resno, b, chain, id, ali, resid, sse, call

```

```
rd <- rmsd(pdba, fit=T)
```

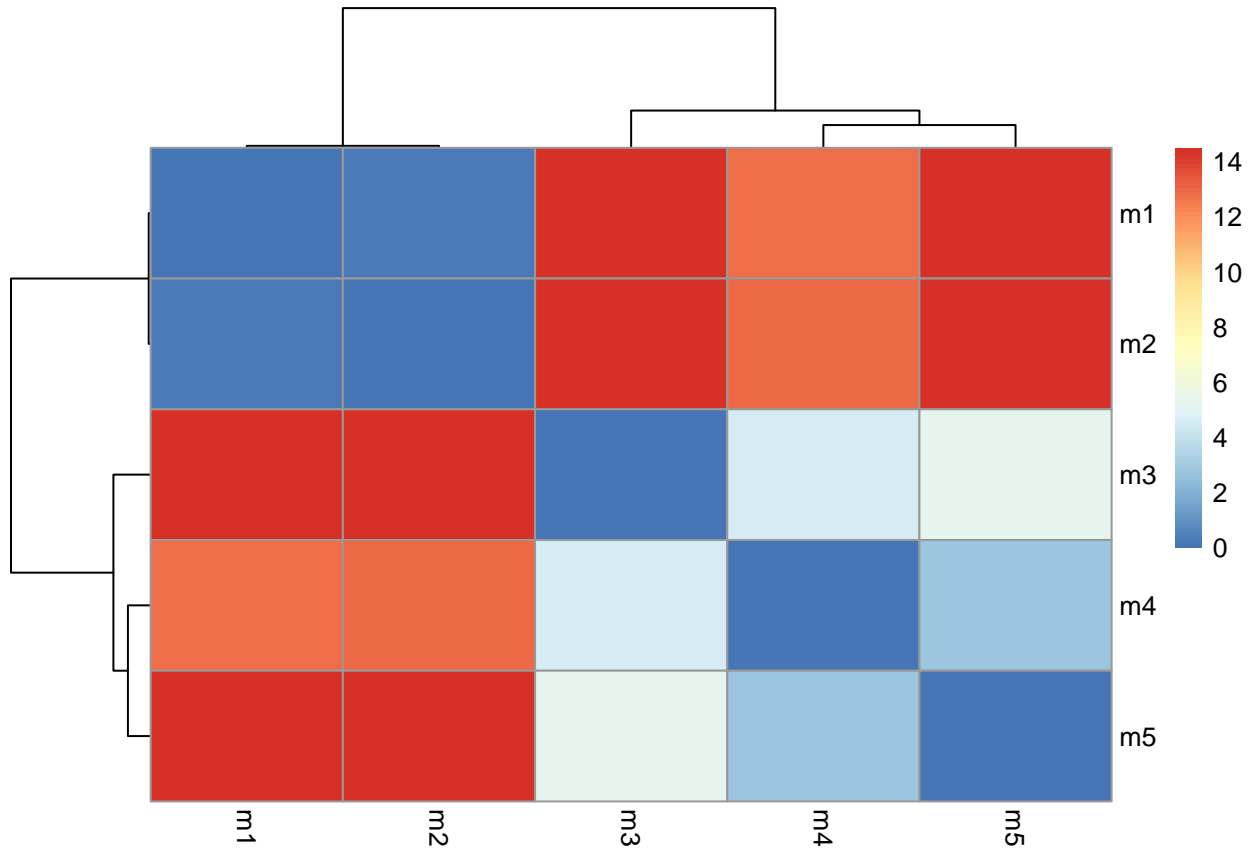
```
## Warning in rmsd(pdba, fit = T): No indices provided, using the 198 non NA positions
```

```
range(rd)
```

```
## [1] 0.000 14.507
```

```
library(pheatmap)

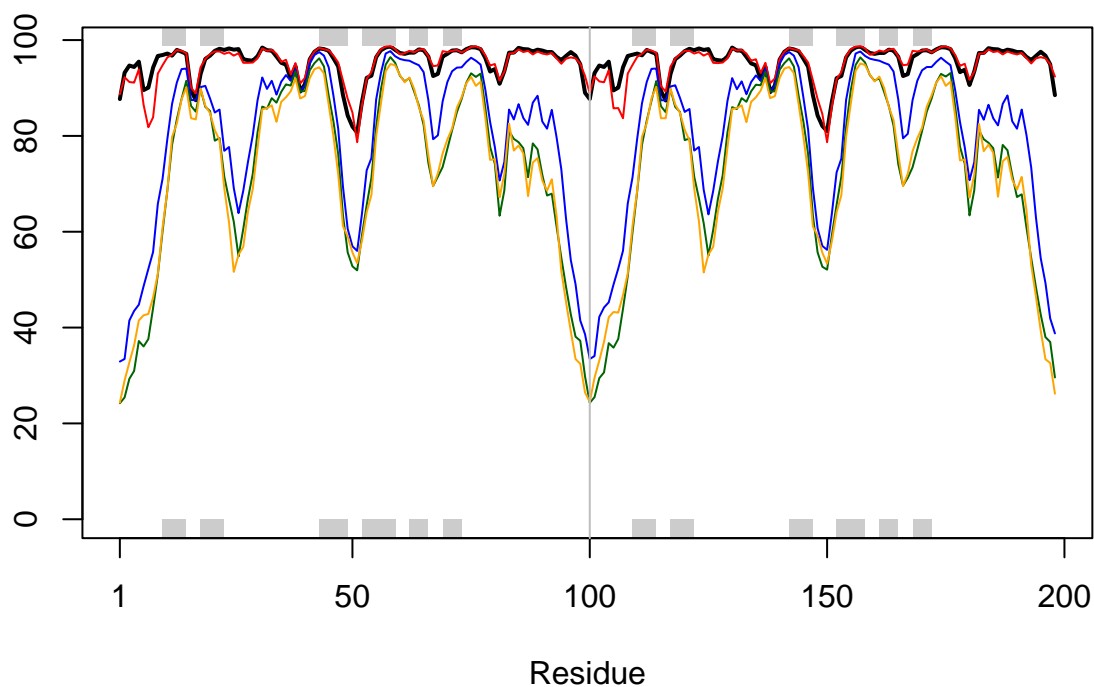
colnames(rd) <- paste0("m",1:5)
rownames(rd) <- paste0("m",1:5)
pheatmap(rd)
```



```
# Read a reference PDB structure
pdb <- read.pdb("1hsg")
```

```
## Note: Accessing on-line PDB file
```

```
plotb3(pdb$b[1,], typ="l", lwd=2, sse=pdb)
points(pdb$b[2,], typ="l", col="red")
points(pdb$b[3,], typ="l", col="blue")
points(pdb$b[4,], typ="l", col="darkgreen")
points(pdb$b[5,], typ="l", col="orange")
abline(v=100, col="gray")
```



```
core <- core.find(pdb)
```

```
## core size 197 of 198 vol = 5017.583
## core size 196 of 198 vol = 4299.462
## core size 195 of 198 vol = 4030.786
## core size 194 of 198 vol = 3797.241
## core size 193 of 198 vol = 3567.126
## core size 192 of 198 vol = 3378.469
## core size 191 of 198 vol = 3249.342
## core size 190 of 198 vol = 3149.254
## core size 189 of 198 vol = 3070.29
## core size 188 of 198 vol = 2993.999
## core size 187 of 198 vol = 2917.618
## core size 186 of 198 vol = 2865.321
## core size 185 of 198 vol = 2835.031
## core size 184 of 198 vol = 2825.584
## core size 183 of 198 vol = 2833.979
## core size 182 of 198 vol = 2894.691
## core size 181 of 198 vol = 2975.843
## core size 180 of 198 vol = 3026.495
## core size 179 of 198 vol = 3070.895
## core size 178 of 198 vol = 3121.204
## core size 177 of 198 vol = 3127.656
## core size 176 of 198 vol = 3102.311
## core size 175 of 198 vol = 3060.45
```

```

## core size 174 of 198 vol = 2993.84
## core size 173 of 198 vol = 2902.747
## core size 172 of 198 vol = 2841.824
## core size 171 of 198 vol = 2771.39
## core size 170 of 198 vol = 2708.164
## core size 169 of 198 vol = 2616.115
## core size 168 of 198 vol = 2540.663
## core size 167 of 198 vol = 2471.823
## core size 166 of 198 vol = 2396.567
## core size 165 of 198 vol = 2324.756
## core size 164 of 198 vol = 2258.532
## core size 163 of 198 vol = 2189.811
## core size 162 of 198 vol = 2118.531
## core size 161 of 198 vol = 2048.541
## core size 160 of 198 vol = 1964.22
## core size 159 of 198 vol = 1878.019
## core size 158 of 198 vol = 1802.026
## core size 157 of 198 vol = 1719.543
## core size 156 of 198 vol = 1640.479
## core size 155 of 198 vol = 1561.746
## core size 154 of 198 vol = 1490.107
## core size 153 of 198 vol = 1416.211
## core size 152 of 198 vol = 1345.494
## core size 151 of 198 vol = 1287.606
## core size 150 of 198 vol = 1225.523
## core size 149 of 198 vol = 1168.6
## core size 148 of 198 vol = 1123.809
## core size 147 of 198 vol = 1069.607
## core size 146 of 198 vol = 1028.33
## core size 145 of 198 vol = 986.295
## core size 144 of 198 vol = 947.191
## core size 143 of 198 vol = 910.624
## core size 142 of 198 vol = 868.922
## core size 141 of 198 vol = 829.982
## core size 140 of 198 vol = 788.548
## core size 139 of 198 vol = 749.234
## core size 138 of 198 vol = 713.554
## core size 137 of 198 vol = 679.035
## core size 136 of 198 vol = 639.012
## core size 135 of 198 vol = 599.236
## core size 134 of 198 vol = 556.226
## core size 133 of 198 vol = 521.307
## core size 132 of 198 vol = 484.526
## core size 131 of 198 vol = 453.614
## core size 130 of 198 vol = 422.947
## core size 129 of 198 vol = 404.641
## core size 128 of 198 vol = 397.064
## core size 127 of 198 vol = 371.629
## core size 126 of 198 vol = 355.609
## core size 125 of 198 vol = 334.859
## core size 124 of 198 vol = 313.691
## core size 123 of 198 vol = 291.489
## core size 122 of 198 vol = 268.734
## core size 121 of 198 vol = 245.865

```

```

## core size 120 of 198 vol = 236.559
## core size 119 of 198 vol = 218.641
## core size 118 of 198 vol = 201.313
## core size 117 of 198 vol = 183.861
## core size 116 of 198 vol = 167.249
## core size 115 of 198 vol = 151.276
## core size 114 of 198 vol = 137.843
## core size 113 of 198 vol = 124.983
## core size 112 of 198 vol = 112.07
## core size 111 of 198 vol = 101.394
## core size 110 of 198 vol = 91.994
## core size 109 of 198 vol = 82.201
## core size 108 of 198 vol = 74.644
## core size 107 of 198 vol = 70.256
## core size 106 of 198 vol = 64.859
## core size 105 of 198 vol = 58.745
## core size 104 of 198 vol = 54.966
## core size 103 of 198 vol = 49.885
## core size 102 of 198 vol = 45.389
## core size 101 of 198 vol = 41.648
## core size 100 of 198 vol = 38.714
## core size 99 of 198 vol = 36.289
## core size 98 of 198 vol = 33.698
## core size 97 of 198 vol = 28.156
## core size 96 of 198 vol = 23.583
## core size 95 of 198 vol = 19.899
## core size 94 of 198 vol = 16.637
## core size 93 of 198 vol = 12.448
## core size 92 of 198 vol = 9.42
## core size 91 of 198 vol = 8.296
## core size 90 of 198 vol = 5.783
## core size 89 of 198 vol = 4.006
## core size 88 of 198 vol = 2.903
## core size 87 of 198 vol = 2.24
## core size 86 of 198 vol = 1.765
## core size 85 of 198 vol = 1.408
## core size 84 of 198 vol = 1.164
## core size 83 of 198 vol = 0.969
## core size 82 of 198 vol = 0.833
## core size 81 of 198 vol = 0.675
## core size 80 of 198 vol = 0.579
## core size 79 of 198 vol = 0.529
## core size 78 of 198 vol = 0.456
## FINISHED: Min vol ( 0.5 ) reached

```

```
core.inds <- print(core, vol=0.5)
```

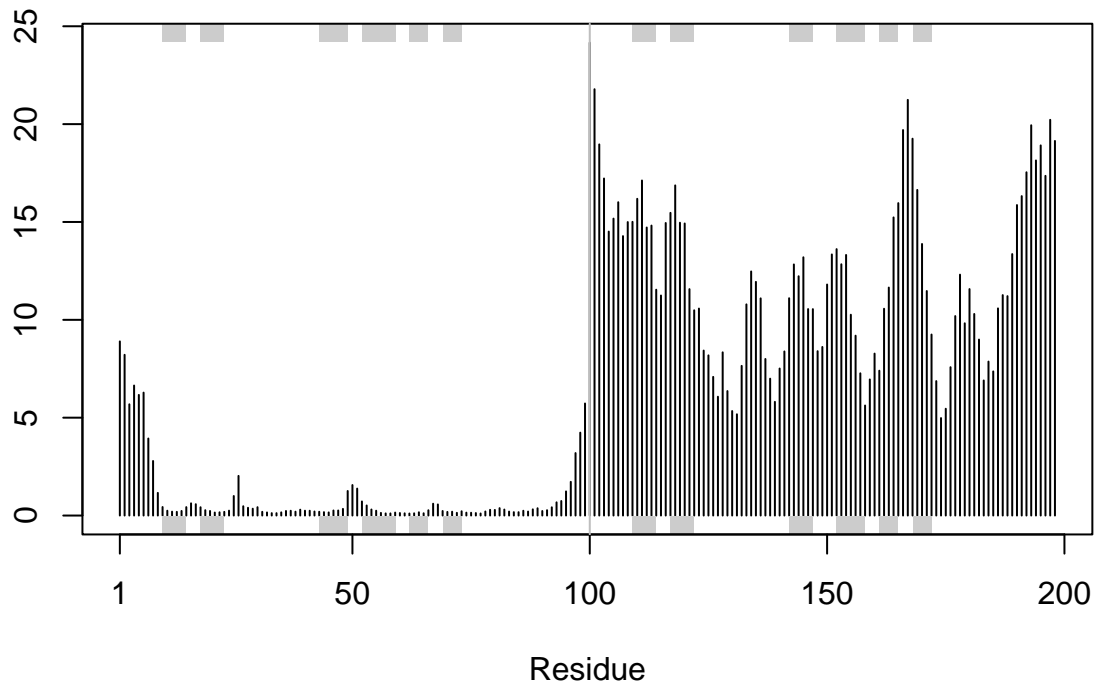
```

## # 79 positions (cumulative volume <= 0.5 Angstrom^3)
##   start end length
## 1    10  24     15
## 2    27  48     22
## 3    53  94     42

```

```
xyz <- pdbfit(pdb, core.inds, outpath="corefit_structures")
rf <- rmsf(xyz)

plotb3(rf, sse=pdb)
abline(v=100, col="gray", ylab="RMSF")
```



```
library(jsonlite)

# Listing of all PAE JSON files
pae_files <- list.files(path=results_dir,
                        pattern=".*model.*\\.json",
                        full.names = TRUE)

pae1 <- read_json(pae_files[1],simplifyVector = TRUE)
pae5 <- read_json(pae_files[5],simplifyVector = TRUE)
attributes(pae1)

## $names
## [1] "plddt" "max_pae" "pae" "ptm" "iptm"

# Per-residue pLDDT scores
# same as B-factor of PDB..
head(pae1$plddt)
```

```
## [1] 87.69 93.19 94.69 94.38 95.50 89.56
```

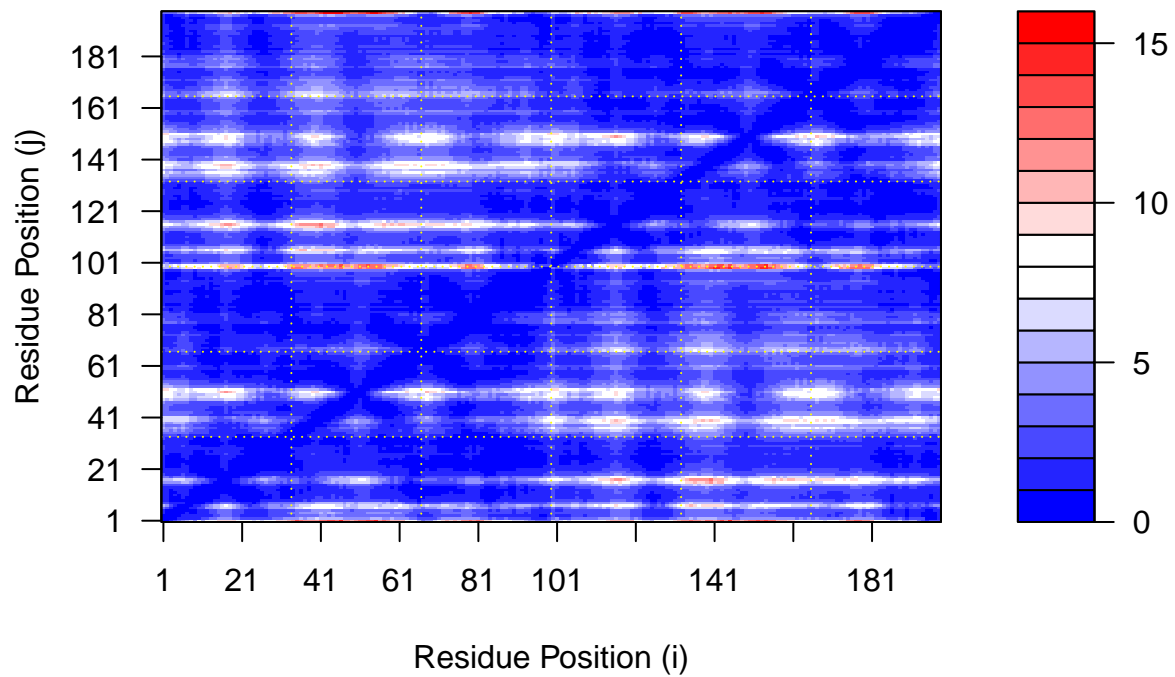
```
pae1$max_pae
```

```
## [1] 15.89844
```

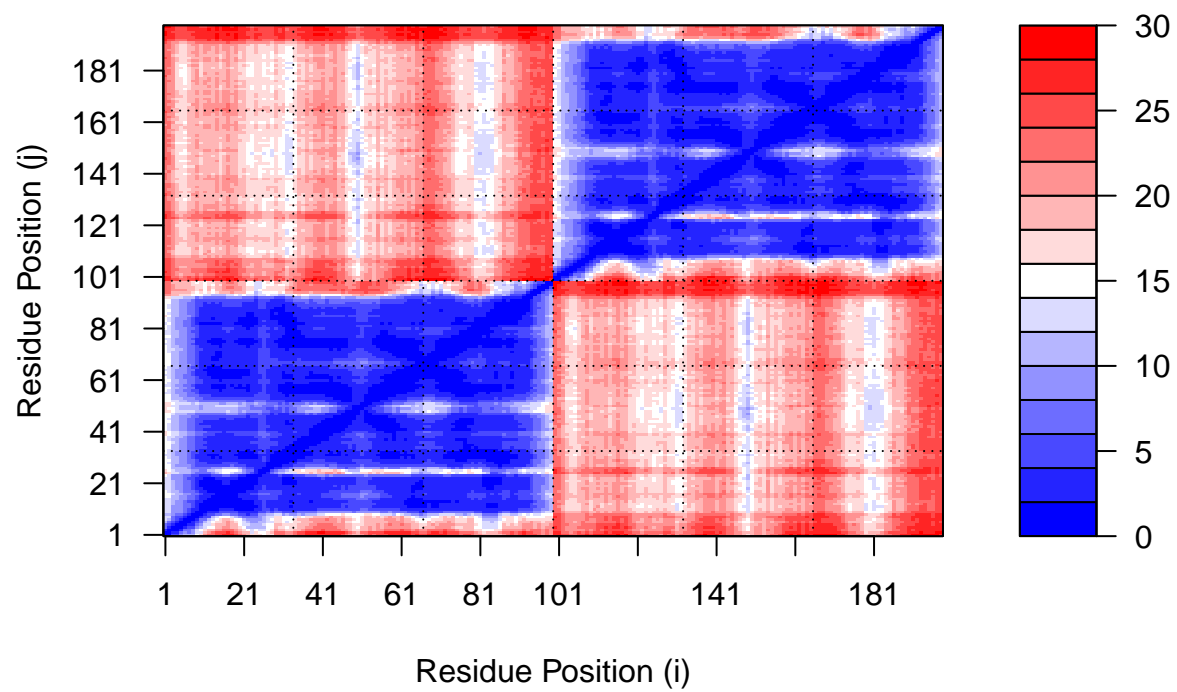
```
pae5$max_pae
```

```
## [1] 29.25
```

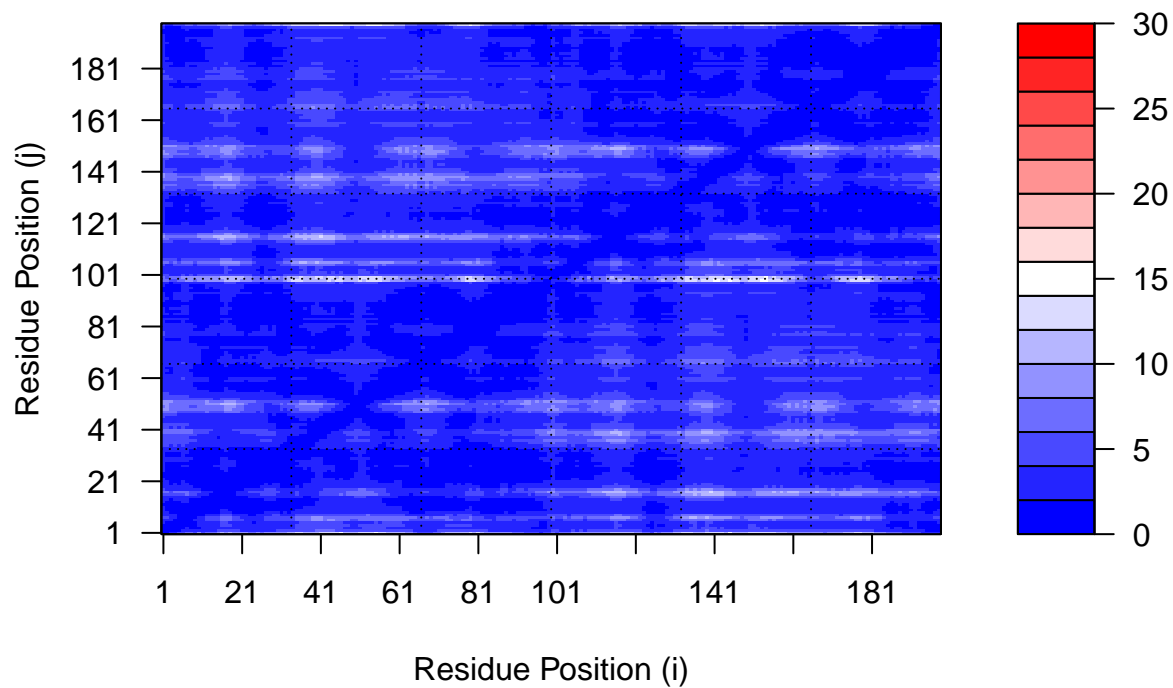
```
plot.dmat(pae1$pae,  
          xlab="Residue Position (i)",  
          ylab="Residue Position (j)")
```



```
plot.dmat(pae5$pae,  
          xlab="Residue Position (i)",  
          ylab="Residue Position (j)",  
          grid.col = "black",  
          zlim=c(0,30))
```

```
plot.dmat(pae1$paes,
  xlab="Residue Position (i)",
  ylab="Residue Position (j)",
  grid.col = "black",
  zlim=c(0,30))
```



```
aln_file <- list.files(path=results_dir,
                      pattern=".a3m$",
                      full.names = TRUE)
aln_file

## [1] "HIVPRDIMER_23119//HIVPRDIMER_23119.a3m"

aln <- read.fasta(aln_file[1], to.upper = TRUE)

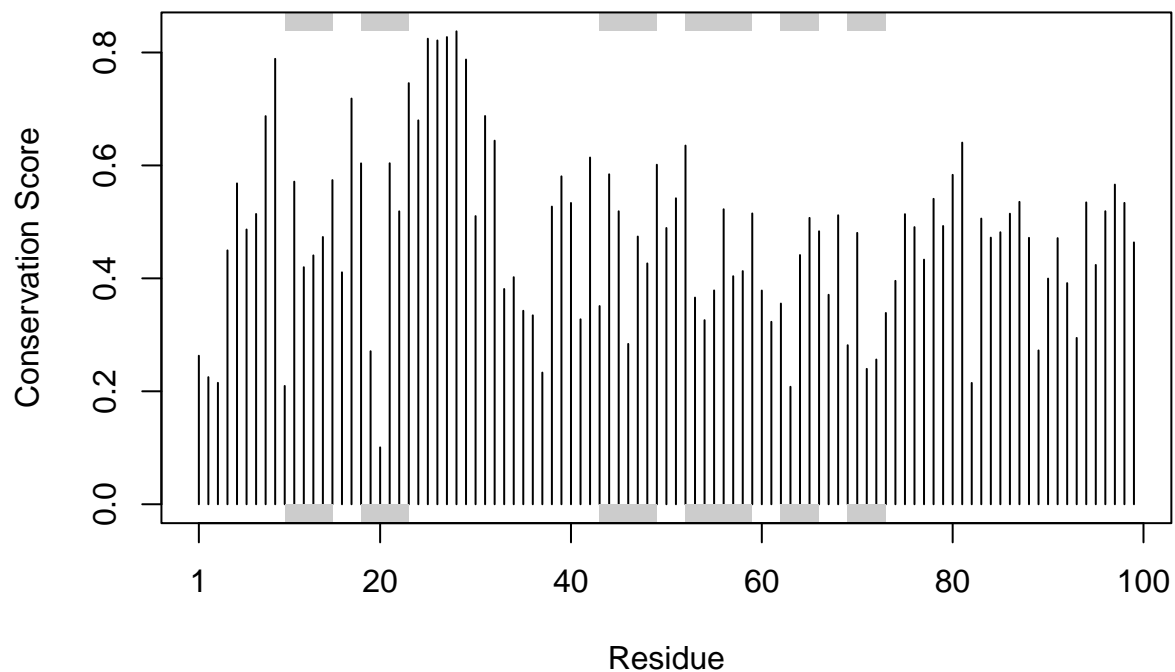
## [1] " ** Duplicated sequence id's: 101 **"
## [2] " ** Duplicated sequence id's: 101 **"

dim(aln$ali)

## [1] 5378 132

sim <- conserv(aln)

plotb3(sim[1:99], sse=trim.pdb(pdb, chain="A"),
      ylab="Conservation Score")
```



```
con <- consensus(aln, cutoff = 0.9)
con$seq
```

```
## [1] "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-"
## [19] "-" "-" "-" "-" "-" "-" "D" "T" "G" "A" "-" "-" "-" "-" "-" "-" "-"
## [37] "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-"
## [55] "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-"
## [73] "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-"
## [91] "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-"
## [109] "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-" "-"
## [127] "-" "-" "-" "-" "-" "-"
```

```
m1.pdb <- read.pdb(pdb_files[1])
occ <- vec2resno(c(sim[1:99], sim[1:99]), m1.pdb$atom$resno)
write.pdb(m1.pdb, o=occ, file="m1_conserv.pdb")
```