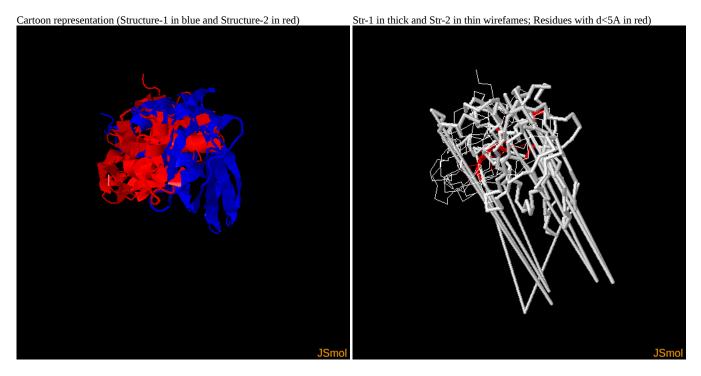
4/15/2018 TM-score Server

## **TM-score Results**

```
TM-SCORE
  A scoring function to assess the similarity of protein structures
  Based on statistics:
  0.0 < TM-score < 0.17, random structural similarity
0.5 < TM-score < 1.00, in about the same fold
Reference: Yang Zhang and Jeffrey Skolnick, Proteins 2004 57: 702-710
  For comments, please email to: zhng@umich.edu
                       Length= 215
Length= 209 (by which all scores are normalized)
Structure1: A735896
Structure2: B735896
Number of residues in common= 69
RMSD of the common residues=
                                11.878
           = 0.0961
                      (d0 = 5.38)
TM-score
MaxSub-score= 0.0430 (d0= 3.50)
GDT-TS-score= 0.0622 %(d<1)=0.0239 %(d<2)=0.0335 %(d<4)=0.0622 %(d<8)=0.1292
GDT-HA-score= 0.0359 %(d<0.5)=0.0239 %(d<1)=0.0239 %(d<2)=0.0335 %(d<4)=0.0622
 ----- rotation matrix to rotate Chain-1 to Chain-2
                                                        u(i.3)
i
                         u(i.1)
                                        u(i.2)
                       0.8105900163
        7.7107086384
                                      0.0230008372 -0.5851621886
1
      -52.7555597529
                       0.3666185917
                                      0.7592568600
       13.7562278410
                       0.4566559248
                                     -0.6503844878
                                                      0.6070134960
Superposition in the TM-score: Length(d<5.0)= 15 \, RMSD= 3.23 (":" denotes the residue pairs of distance < 5.0 Angstrom)
.
PVANADVVFDF0NYTAKAGDEVTVDVLVDSKNKPISAXDVKFKVDSPLTIEEIDKESLAFNTTVXTNXAILGANFKSLDDKGEPLVPKDGAAVFTLYVNVPANTPDGTYYVGFNGKNEVHKSNDGS0FTVASKNGA
```

## Visualization of TM-score superposition



## **Document downloads**

- Click <u>A735896.pdb</u> to download the first structure that you submitted.
- Click <u>B735896.pdb</u> to download the second structure that you submitted.
- Click <u>C735896.pdb</u> to download the superposed structure in C-alpah trace (This file is in a Rasmol script format, you can run 'rasmol -script C735896.pdb' to view the structure).
- Click <u>D735896.pdb</u> to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D735896.pdb' to view the structure).

## **Reference:**

- Y. Zhang, J. Skolnick, Scoring function for automated assessment of protein structure template quality, Proteins, 57: 702-710 (2004).
- J. Xu, Y. Zhang, How significant is a protein structure similarity with TM-score=0.5? Bioinformatics, 26, 889-895 (2010).

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