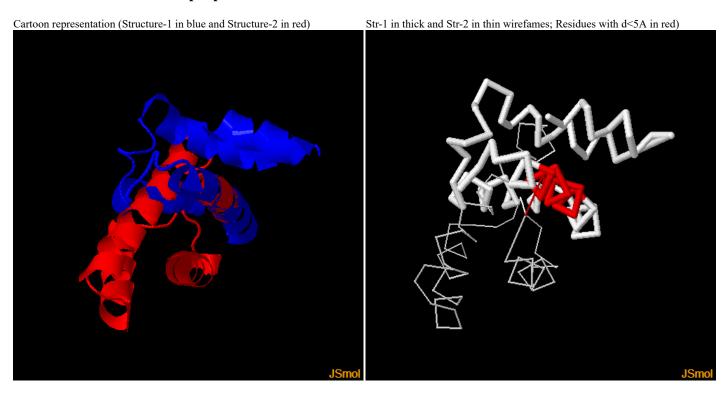
## **TM-score Results**

```
TM-SCORE
  A scoring function to assess the similarity of protein structures
  Based on statistics:
       0.0 < \text{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
Structure1: A853690
                   Length=
                           73
                          73 (by which all scores are normalized)
Structure2: B853690
                   Length=
Number of residues in common= 11
RMSD of the common residues=
TM-score
         = 0.1481 (d0= 3.00)
MaxSub-score= 0.1488 (d0= 3.50)
GDT-TS-score= 0.1473 %(d<1)=0.1370 %(d<2)=0.1507 %(d<4)=0.1507 %(d<8)=0.1507
GDT-HA-score= 0.1438 %(d<0.5)=0.1370 %(d<1)=0.1370 %(d<2)=0.1507 %(d<4)=0.1507
   ---- rotation matrix to rotate Chain-1 to Chain-2 -----
i
         t(i)
                    u(i,1)
                                u(i,2)
                                             u(i,3)
      4.2541864975
                  0.7122720658
                             -0.6971421019
                                           0.0816173631
      8.9488459818
                  0.4195807867 0.5161080436
                                           0.7467157764
      2.0418661158 -0.5626903835 -0.4976197113
                                           0.6601167740
Superposition in the TM-score: Length(d<5.0)= 11 RMSD= 0.41
(":" denotes the residue pairs of distance < 5.0 Angstrom)
VADDWLKQYANREKAIQHAIERFNTKPIQTIKKHDYQRFVDDISAQYSKNYVDSIVASTNMIFKYAYDTRLIK-------
                                            :::::::::
  ·-----REKAIQHAIERFNTKPIQTIKKHDYQRFVDDISAQYSKNYVDSIVASTNMIFKYAYDTRLIK
```

## Visualization of TM-score superposition



## **Document downloads**

- Click <u>A853690.pdb</u> to download the first structure that you submitted.
- Click <u>B853690.pdb</u> to download the second structure that you submitted.
- Click <u>C853690.pdb</u> to download the superposed structure in C-alpah trace (This file is in a Rasmol script format, you can run 'rasmol -script C853690.pdb' to view the structure).
- Click <u>D853690.pdb</u> to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D853690.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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