

## TM-score Results

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*                               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: zhngumich.edu                       *
*****
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Structure1: A298834      Length= 215
Structure2: B298834      Length= 209 (by which all scores are normalized)
Number of residues in common= 69
RMSD of the common residues= 11.796
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TM-score = 0.0980 (d0= 5.38)  
 MaxSub-score= 0.0423 (d0= 3.50)  
 GDT-TS-score= 0.0646 %(d<1)=0.0335 %(d<2)=0.0335 %(d<4)=0.0622 %(d<8)=0.1292  
 GDT-HA-score= 0.0383 %(d<0.5)=0.0239 %(d<1)=0.0335 %(d<2)=0.0335 %(d<4)=0.0622

```

----- rotation matrix to rotate Chain-1 to Chain-2 -----
i          t(i)          u(i,1)          u(i,2)          u(i,3)
1   -29.1188975710     0.6299042053     -0.3437253043     0.6964722589
2    -6.0024167148     0.7766700445     -0.2811739146     0.5636708896
3    72.8795169205     0.0020818834     0.8959878040     -0.4440737786

```

Superposition in the TM-score: Length(d<5.0)= 14    RMSD= 3.06

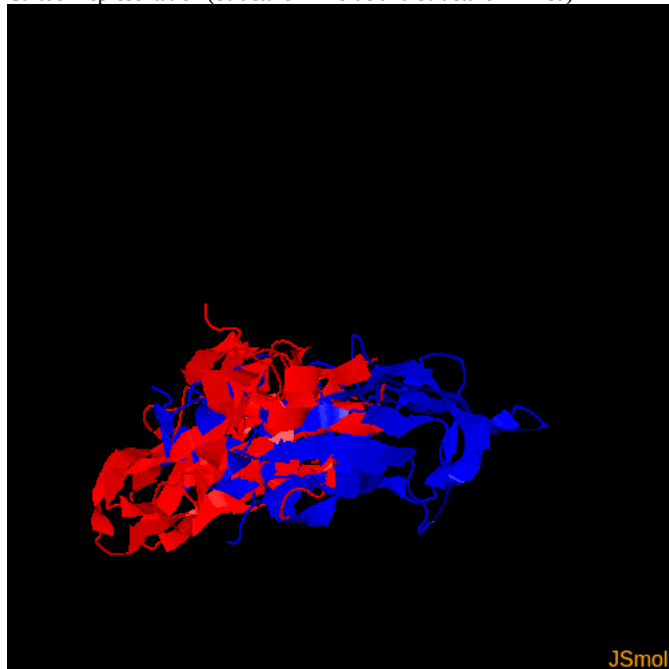
superposition in the m score: Length(a=5.0)= 14 RMSD= 5  
 (":" denotes the residue pairs of distance < 5.0 Angstrom)

PVANADVVFDFONYTAKAGDEVTVDLVDSKNKPI SAXDVKFKVDSPLTIEEDKESLAFNTTVXTNXAILGANFKSLDDKGEPLVPKDGA AVFTLYVNPANTPDGTY YVGFGKNEVHKSNDGSOFTVASKNGA

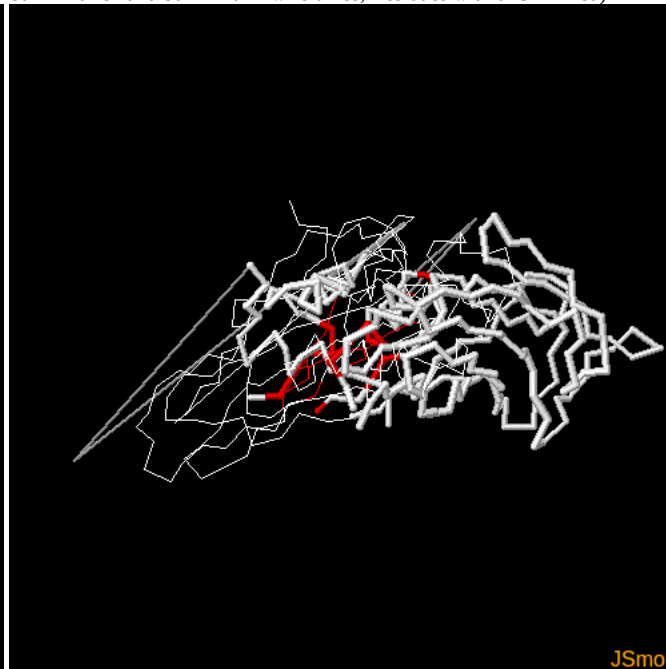
[illegible]

## Visualization of TM-score superposition

Cartoon representation (Structure-1 in blue and Structure-2 in red)



Str-1 in thick and Str-2 in thin wireframes; Residues with  $d < 5\text{\AA}$  in red)



## Document downloads

- Click [A298834.pdb](#) to download the first structure that you submitted.
- Click [B298834.pdb](#) to download the second structure that you submitted.
- Click [C298834.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C298834.pdb' to view the structure).
- Click [D298834.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D298834.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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