

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: zhng@umich.edu                        *
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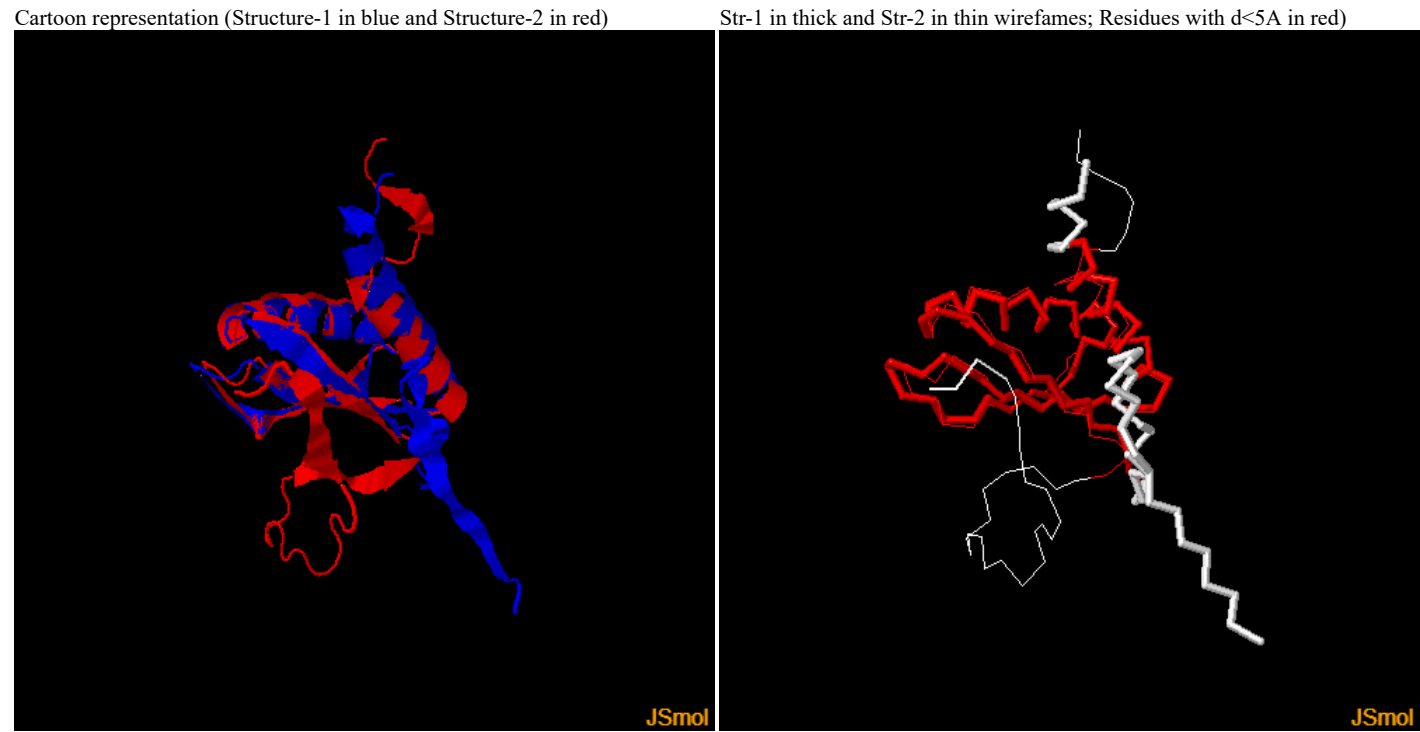
Structure1: A690420      Length= 108
Structure2: B690420      Length= 108 (by which all scores are normalized)
Number of residues in common= 108
RMSD of the common residues= 10.999

TM-score    = 0.6435 (d0= 3.82)
MaxSub-score= 0.6112 (d0= 3.50)
GDT-TS-score= 0.6296 %(d<1)=0.4167 %(d<2)=0.6389 %(d<4)=0.6852 %(d<8)=0.7778
GDT-HA-score= 0.4769 %(d<0.5)=0.1667 %(d<1)=0.4167 %(d<2)=0.6389 %(d<4)=0.6852

----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      -0.2073525651  0.2142509350  -0.1119444473  -0.9703427114
2      -3.8397426394  -0.6318781691  -0.7734344629  -0.0502902681
3      -0.1487872678  -0.7448667775  0.6239131128  -0.2364443092

Superposition in the TM-score: Length(d<5.0)= 74 RMSD= 1.29
(":" denotes the residue pairs of distance < 5.0 Angstrom)
MSNQTCVENEVCEACGCAGEIGFIIREGDDVAEVSFLFGSDKAHLEGKLAEYISLAKQVYANVEYEVAPVADNATELHARFKFEVSAEKLIFELKTRALARLEHHHHHH
:
MSNQTCVENEVCEACGCAGEIGFIIREGDDVAEVSFLFGSDKAHLEGKLAEYISLAKQVYANVEYEVAPVADNATELHARFKFEVSAEKLIFELKTRALARLEHHHHHH
12345678901234567890123456789012345678901234567890123456789012345678901234567890123456789012345678
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Visualization of TM-score superposition



Document downloads

- Click [A690420.pdb](#) to download the first structure that you submitted.
- Click [B690420.pdb](#) to download the second structure that you submitted.
- Click [C690420.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C690420.pdb' to view the structure).
- Click [D690420.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D690420.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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