

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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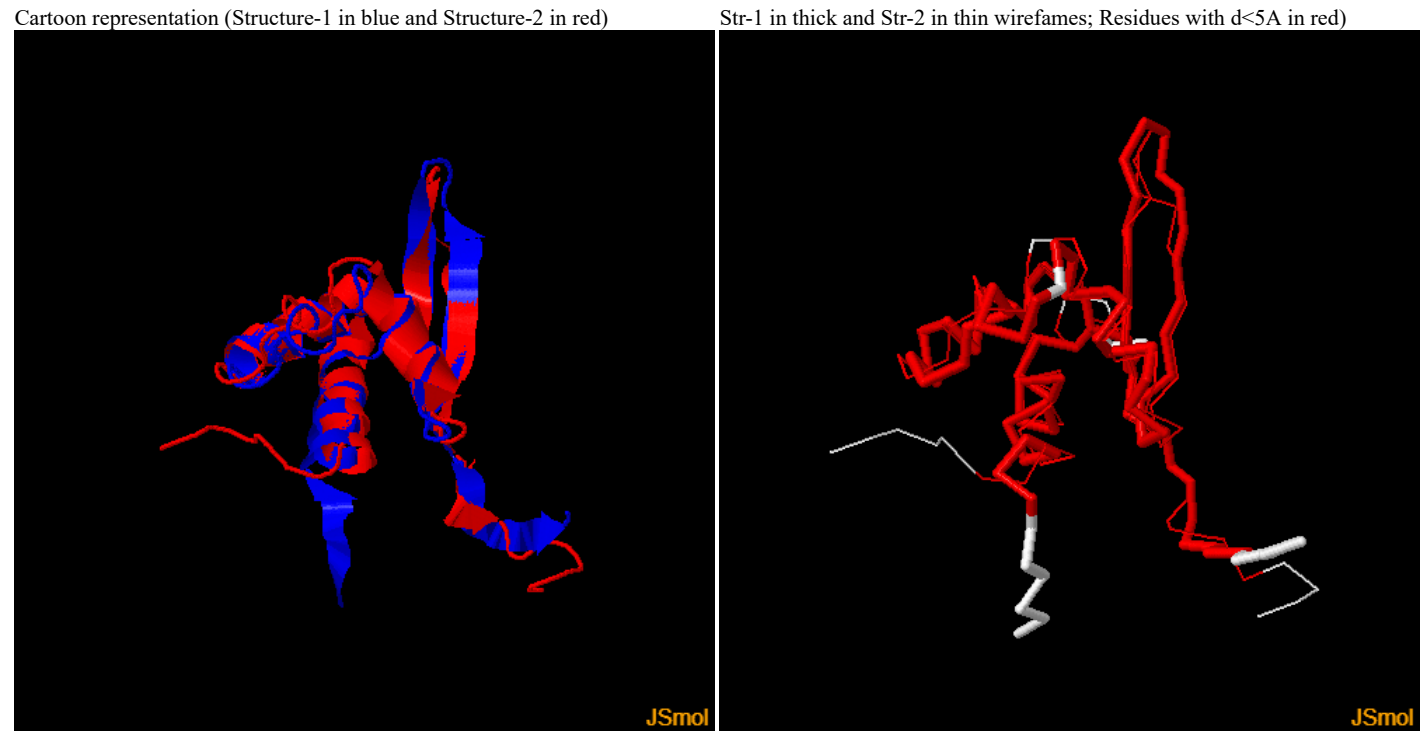
Structure1: A444159      Length= 92
Structure2: B444159      Length= 92 (by which all scores are normalized)
Number of residues in common= 92
RMSD of the common residues= 4.801

TM-score    = 0.7333 (d0= 3.48)
MaxSub-score= 0.6999 (d0= 3.50)
GDT-TS-score= 0.7255 %(d<1)=0.3913 %(d<2)=0.7065 %(d<4)=0.8587 %(d<8)=0.9457
GDT-HA-score= 0.5353 %(d<0.5)=0.1848 %(d<1)=0.3913 %(d<2)=0.7065 %(d<4)=0.8587

----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      3.0358118098 -0.2319123359  0.9429160684 -0.2390103687
2      -2.9946983979 0.8362843925  0.0677696901 -0.5440916136
3      1.6774951143 -0.4968350665 -0.3260621980 -0.8042626186

Superposition in the TM-score: Length(d<5.0)= 79 RMSD= 1.79
(":" denotes the residue pairs of distance < 5.0 Angstrom)
MPHKEKHPLQDMFTSAIEAVARDSGWAELSAVGSYLAKNDPSFDPRNWGHGRLSQMVKKLDLFTVQESRNGSKLHSEIRLRHDGLEHHHHHH
:
MPHKEKHPLQDMFTSAIEAVARDSGWAELSAVGSYLAKNDPSFDPRNWGHGRLSQMVKKLDLFTVQESRNGSKLHSEIRLRHDGLEHHHHHH
1234567890123456789012345678901234567890123456789012345678901234567890123456789012
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Visualization of TM-score superposition



Document downloads

- Click [A444159.pdb](#) to download the first structure that you submitted.
- Click [B444159.pdb](#) to download the second structure that you submitted.
- Click [C444159.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C444159.pdb' to view the structure).
- Click [D444159.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D444159.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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