

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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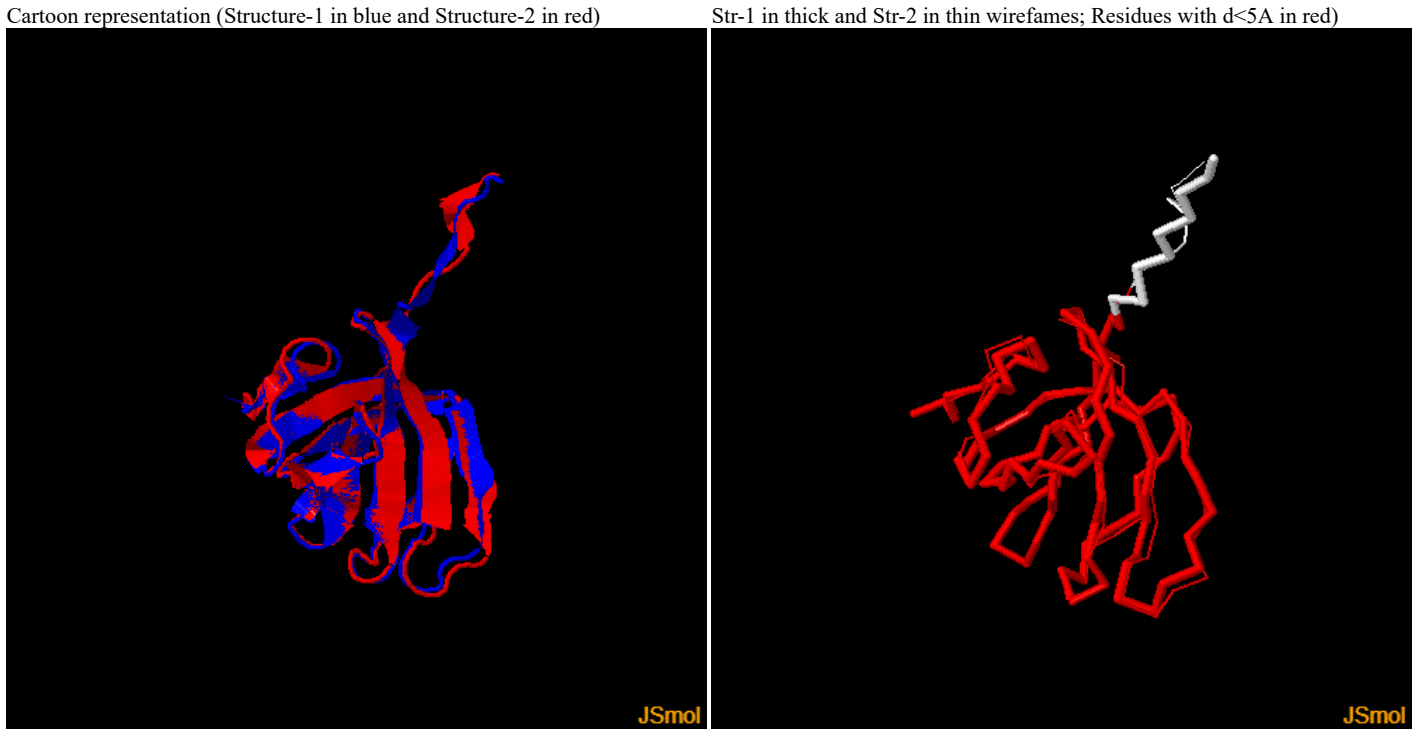
Structure1: A732053      Length= 109
Structure2: B732053      Length= 109 (by which all scores are normalized)
Number of residues in common= 109
RMSD of the common residues= 4.807

TM-score      = 0.8751 (d0= 3.84)
MaxSub-score= 0.8554 (d0= 3.50)
GDT-TS-score= 0.8761 %(d<1)=0.7064 %(d<2)=0.8991 %(d<4)=0.9266 %(d<8)=0.9725
GDT-HA-score= 0.7225 %(d<0.5)=0.3578 %(d<1)=0.7064 %(d<2)=0.8991 %(d<4)=0.9266

----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      0.6227171204 -0.8009108155 -0.2241043709 -0.5552648886
2     -0.5737271324 -0.5755268755  0.0321780659  0.8171495503
3      0.8916082558 -0.1652594357  0.9740337792 -0.1547498494

Superposition in the TM-score: Length(d<5.0)=100 RMSD= 1.00
(":" denotes the residue pairs of distance < 5.0 Angstrom)
MEPQLTKIVDIVENGQWANLKAKVIQLWENTHESISQVGLLGDETGIKFTIWKNALPLLEQGESYLLRSVVVGEYNDRFQVQVKNKSSIEKLSEPIEVGLEHHHHHH
::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::
MEPQLTKIVDIVENGQWANLKAKVIQLWENTHESISQVGLLGDETGIKFTIWKNALPLLEQGESYLLRSVVVGEYNDRFQVQVKNKSSIEKLSEPIEVGLEHHHHHH
123456789012345678901234567890123456789012345678901234567890123456789012345678901234567890123456789
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Visualization of TM-score superposition



Document downloads

- Click [A732053.pdb](#) to download the first structure that you submitted.
- Click [B732053.pdb](#) to download the second structure that you submitted.
- Click [C732053.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C732053.pdb' to view the structure).
- Click [D732053.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D732053.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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