

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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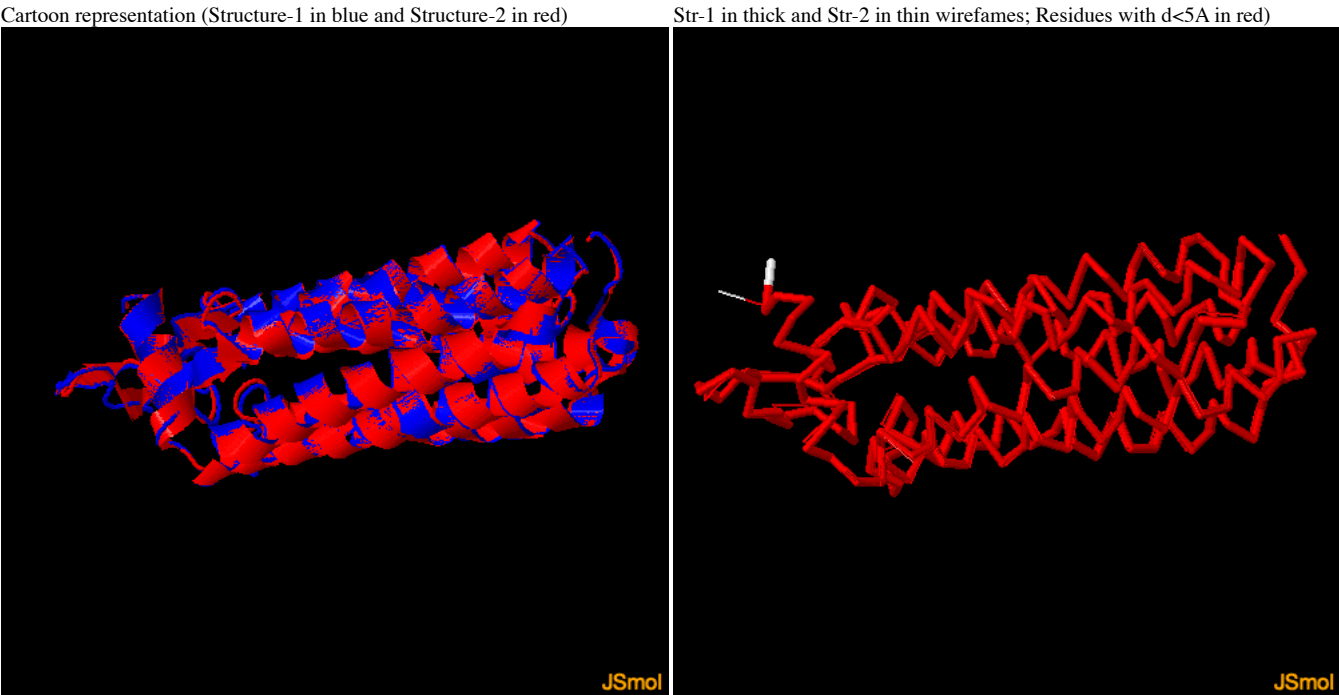
Structure1: A659661      Length= 260
Structure2: B659661      Length= 260 (by which all scores are normalized)
Number of residues in common= 260
RMSD of the common residues= 0.642

TM-score   = 0.9905  (d0= 5.96)
MaxSub-score= 0.9751  (d0= 3.50)
GDT-TS-score= 0.9904  %(d<1)=0.9692  %(d<2)=0.9962  %(d<4)=0.9962  %(d<8)=1.0000
GDT-HA-score= 0.9115  %(d<0.5)=0.6846  %(d<1)=0.9692  %(d<2)=0.9962  %(d<4)=0.9962

----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      -6.6407886887    0.2346782280    0.9709904092    0.0458667045
2      44.6807806514   -0.7820265845    0.2166125669   -0.5843914929
3      -1.0573249440   -0.5773738394    0.1012749777    0.8101745667

Superposition in the TM-score: Length(d<5.0)=259  RMSD= 0.53
(":" denotes the residue pairs of distance < 5.0 Angstrom)
AVTQRELFEFVLNDPLLASSLYINIALAGLSILLFVFMTRGLDDPRAKLIADVSTILVPVVSIASTGLASGLTISVLEMPAGHFAEGSSVMLGGEEVDGVVTMWGRYLTWALSTP
::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::::
AVTQRELFEFVLNDPLLASSLYINIALAGLSILLFVFMTRGLDDPRAKLIADVSTILVPVVSIASTGLASGLTISVLEMPAGHFAEGSSVMLGGEEVDGVVTMWGRYLTWALSTP
12345678901234567890123456789012345678901234567890123456789012345678901234567890123456789012345678901234567
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Visualization of TM-score superposition



Document downloads

- Click [A659661.pdb](#) to download the first structure that you submitted.
- Click [B659661.pdb](#) to download the second structure that you submitted.
- Click [C659661.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C659661.pdb' to view the structure).
- Click [D659661.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D659661.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).