

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: A567195      Length=  729
Structure2: B567195      Length=  729 (by which all scores are normalized)
Number of residues in common=  729
RMSD of the common residues=  38.003

TM-score    = 0.2655 (d0= 9.28)
MaxSub-score= 0.2505 (d0= 3.50)
GDT-TS-score= 0.2510 %(d<1)=0.2497 %(d<2)=0.2497 %(d<4)=0.2510 %(d<8)=0.2538
GDT-HA-score= 0.2500 %(d<0.5)=0.2497 %(d<1)=0.2497 %(d<2)=0.2497 %(d<4)=0.2510

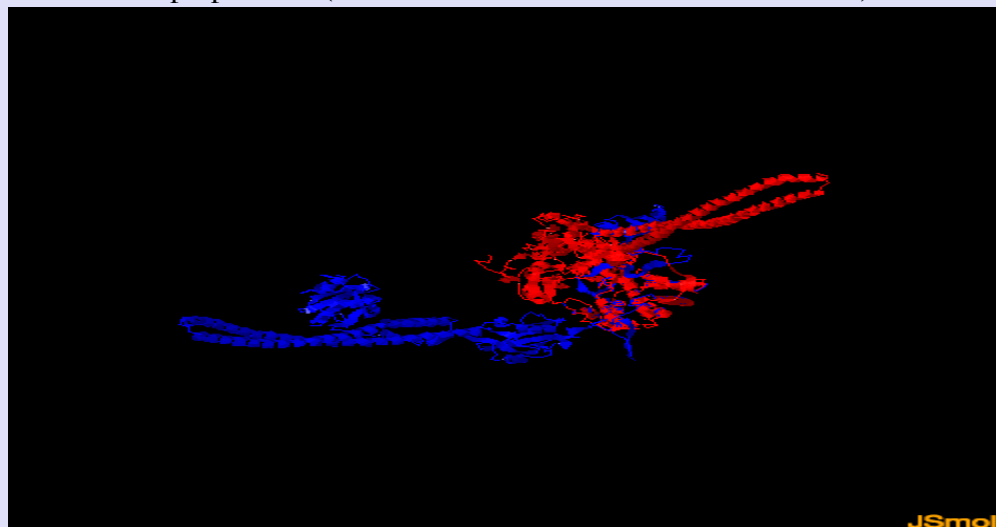
----- rotation matrix to rotate Chain-1 to Chain-2 -----
i          t(i)          u(i,1)          u(i,2)          u(i,3)
1    -49.0547345692    -0.3498685106    -0.6762565244    -0.6482816814
2     -9.0499848549     0.8927991216    -0.4502922737    -0.0121077167
3    269.7965204014    -0.2837283099    -0.5830214245     0.7613043181

Superposition in the TM-score: Length(d<5.0)=183  RMSD=  0.19
(":" denotes the residue pairs of distance < 5.0 Angstrom)
GPLGMSAEGYQYRALYDYKKEREEDIDLHLGDILTVNKGSLVALGFSQGQEARPEEIGWLNQYNETTGERGDFPGTYVEYI
GPLGMSAEGYQYRALYDYKKEREEDIDLHLGDILTVNKGSLVALGFSQGQEARPEEIGWLNQYNETTGERGDFPGTYVEYI
123456789012345678901234567890123456789012345678901234567890123456789012

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Visualization

TM-score superposition (Structure-1 in blue and Structure-2 in red)



Document downloads

- Click [A567195.pdb](#) to download the first structure that you submitted.
- Click [B567195.pdb](#) to download the second structure that you submitted.
- Click [C567195.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C567195.pdb' to view the structure).
- Click [D567195.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D567195.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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