

## TM-score Results

```
*****
*                                     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                           *
*****
```

```
Structure1: A952383      Length= 338
Structure2: B952383      Length= 338 (by which all scores are normalized)
Number of residues in common= 338
RMSD of the common residues= 18.374
```

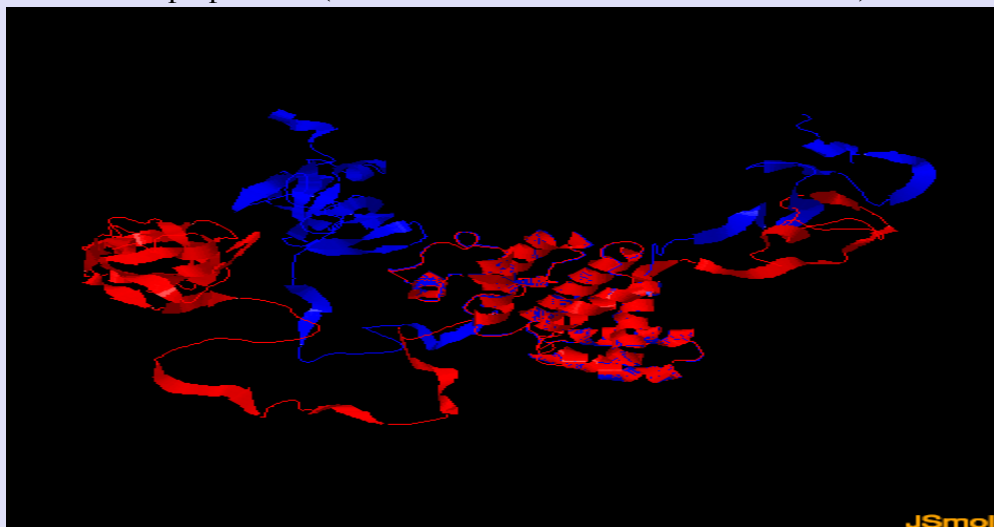
```
TM-score = 0.5600 (d0= 6.71)
MaxSub-score= 0.5402 (d0= 3.50)
GDT-TS-score= 0.5466 %(d<1)=0.5385 %(d<2)=0.5385 %(d<4)=0.5414 %(d<8)=0.5680
GDT-HA-score= 0.5392 %(d<0.5)=0.5385 %(d<1)=0.5385 %(d<2)=0.5385 %(d<4)=0.5414
```

```
----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      -23.1043923245 -0.9431572642 -0.0712425050 0.3246211336
2       91.6544324992 -0.1228424255 0.9823119352 -0.1413258666
3       18.9284987167 -0.3088108052 -0.1731697651 -0.9352262395
```

```
Superposition in the TM-score: Length(d<5.0)=183 RMSD= 0.21
(":" denotes the residue pairs of distance < 5.0 Angstrom)
GPLGMSAEGYQYRALYDYKKEREEDIDLHLGDILTVNKGSLVALGFSQGQEARPEEIGWLNQYNETTGERGDFPGTYVEYI
GPLGMSAEGYQYRALYDYKKEREEDIDLHLGDILTVNKGSLVALGFSQGQEARPEEIGWLNQYNETTGERGDFPGTYVEYI
1234567890123456789012345678901234567890123456789012345678901234567890123456789012
```

## Visualization

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



## Document downloads

- Click [A952383.pdb](#) to download the first structure that you submitted.
- Click [B952383.pdb](#) to download the second structure that you submitted.
- Click [C952383.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C952383.pdb' to view the structure).
- Click [D952383.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D952383.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).

[Back to TM-score Homepage](#)