

## 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                       *
*****
```

```
Structure1: A528290      Length=   94
Structure2: B528290      Length=   94 (by which all scores are normalized)
Number of residues in common=   34
RMSD of the common residues=   5.729
```

```
TM-score    = 0.1870 (d0= 3.52)
MaxSub-score= 0.1784 (d0= 3.50)
GDT-TS-score= 0.2154 %(d<1)=0.1809 %(d<2)=0.1809 %(d<4)=0.1809 %(d<8)=0.3191
GDT-HA-score= 0.1755 %(d<0.5)=0.1596 %(d<1)=0.1809 %(d<2)=0.1809 %(d<4)=0.1809
```

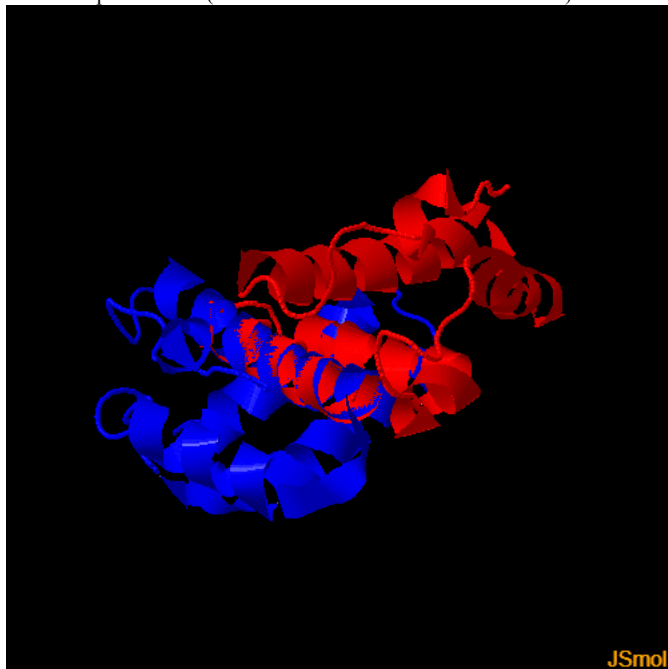
```
----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      -9.5895507140  0.9368562759 -0.1678872644  0.3067803526
2      -5.9157216824 -0.1855743430 -0.9821962501  0.0292008477
3      -6.2278703019  0.2964160614 -0.0842875598 -0.9513322899
```

```
Superposition in the TM-score: Length(d<5.0)= 17 RMSD= 0.41
(":" denotes the residue pairs of distance < 5.0 Angstrom)
```

```
SFGDWAEKFLKSKEADGVSVSQLNSYKNYCRNHLSPLYMKSLSEILPADIQSIINETKLAKNTLKAIRNTASQIFRLAIENRAIDFNPADYVRI-----
:
-----SFGDWAEKFLKSKEADGVSVSQLNSYKNYCRNHLSPLYMKSLSEILPADIQSIINETKLAKNTLKAIRNTASQIFRLAIENRAIDFNPADYVRI
123456789012345678901234567890123456789012345678901234567890123456789012345678901234567890123456789
```

## Visualization of TM-score superposition

Cartoon representation (Structure-1 in blue and Structure-2 in red)



Str-1 in thick and Str-2 in thin wireframes; Residues with d<5A in red



## Document downloads

- Click [A528290.pdb](#) to download the first structure that you submitted.
- Click [B528290.pdb](#) to download the second structure that you submitted.
- Click [C528290.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C528290.pdb' to view the structure).
- Click [D528290.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D528290.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).

