

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: A807922      Length= 61
Structure2: B807922      Length= 61 (by which all scores are normalized)
Number of residues in common= 61
RMSD of the common residues= 4.105
```

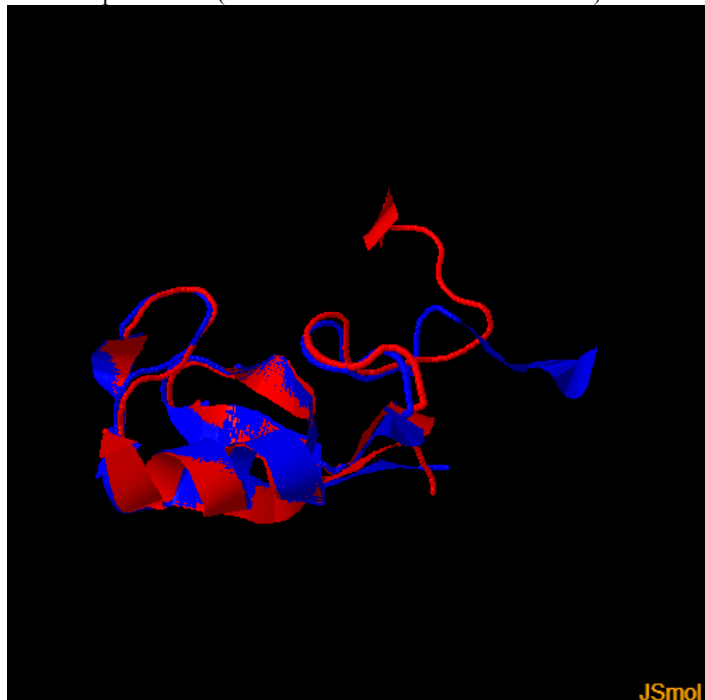
```
TM-score    = 0.8293 (d0= 2.64)
MaxSub-score= 0.8448 (d0= 3.50)
GDT-TS-score= 0.8689 %(d<1)=0.7705 %(d<2)=0.8689 %(d<4)=0.9016 %(d<8)=0.9344
GDT-HA-score= 0.7418 %(d<0.5)=0.4262 %(d<1)=0.7705 %(d<2)=0.8689 %(d<4)=0.9016
```

```
----- rotation matrix to rotate Chain-1 to Chain-2 -----
i      t(i)      u(i,1)      u(i,2)      u(i,3)
1      -6.9984097960  0.7877803716 -0.0136924767 -0.6158040291
2      -11.2968651468 -0.4210719006  0.7176987321 -0.5546232816
3      -8.2451381614  0.4495559373  0.6962191078  0.5596234567
```

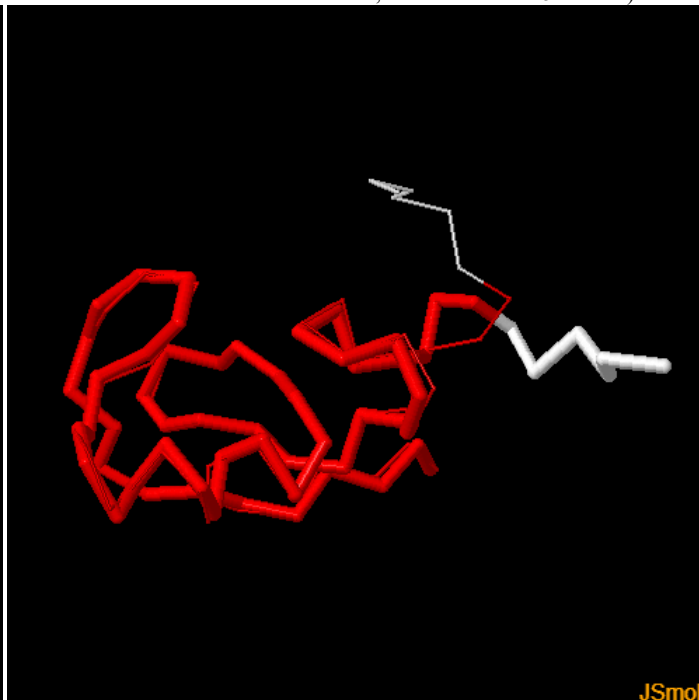
```
Superposition in the TM-score: Length(d<5.0)= 55 RMSD= 0.96
(":" denotes the residue pairs of distance < 5.0 Angstrom)
SHMPTSEEDLCPICYAHPISAVFQPCGHKSKACINQHLMNKNKDCFFCKTTIVSVEDWEKG
:
SHMPTSEEDLCPICYAHPISAVFQPCGHKSKACINQHLMNKNKDCFFCKTTIVSVEDWEKG
1234567890123456789012345678901234567890123456789012345678901
```

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<5Å in red



Document downloads

- Click [A807922.pdb](#) to download the first structure that you submitted.
- Click [B807922.pdb](#) to download the second structure that you submitted.
- Click [C807922.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C807922.pdb' to view the structure).
- Click [D807922.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D807922.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](#)