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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: A182408      Length= 484
Structure2: B182408      Length= 484 (by which all scores are normalized)
Number of residues in common= 242
RMSD of the common residues= 0.000

TM-score = 0.5000 (d0= 7.83)
MaxSub-score= 0.5000 (d0= 3.50)
GDT-TS-score= 0.5000 %(d<1)=0.5000 %(d<2)=0.5000 %(d<4)=0.5000 %(d<8)=0.5000
GDT-HA-score= 0.5000 %(d<0.5)=0.5000 %(d<1)=0.5000 %(d<2)=0.5000 %(d<4)=0.5000

```

```

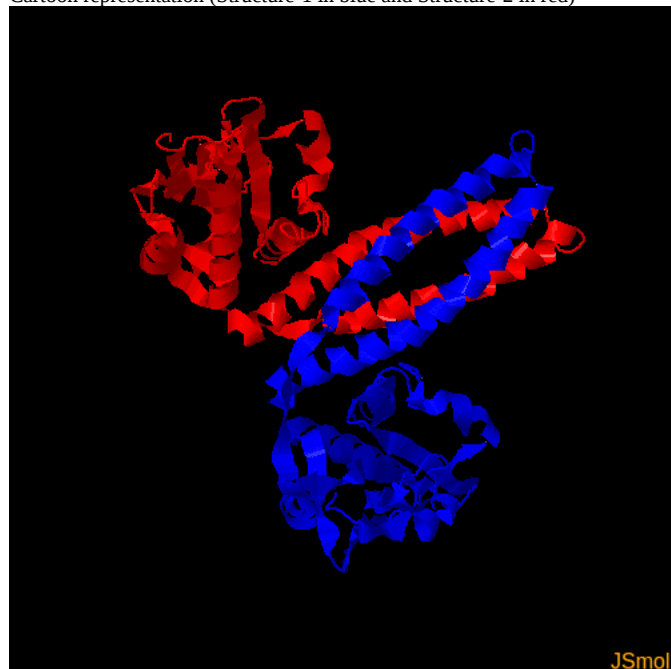
----- rotation matrix to rotate Chain-1 to Chain-2 -----
i          t(i)          u(i,1)          u(i,2)          u(i,3)
1      0.0000000000      1.0000000000      0.0000000000      0.0000000000
2     -0.0000000000      0.0000000000      1.0000000000      0.0000000000
3      0.0000000000      0.0000000000      0.0000000000      1.0000000000

```

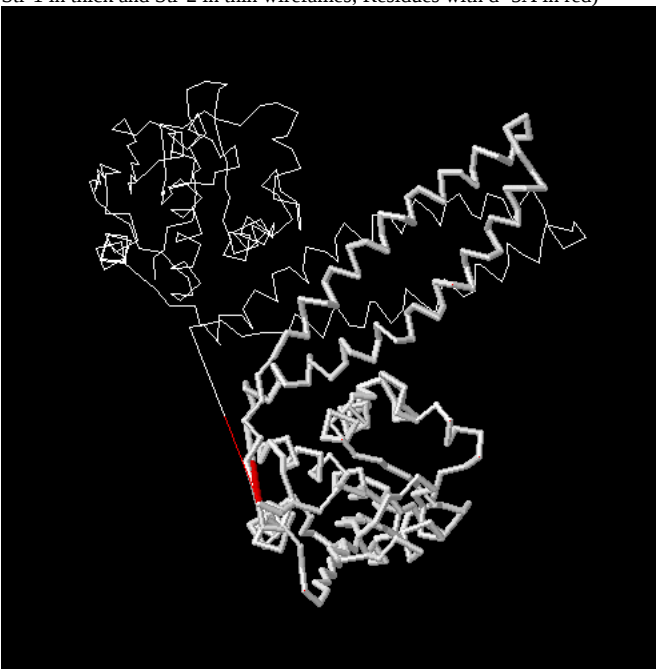
superposition in the m score: Length(a=5.0)=242 RMSD= 0  
(":" denotes the residue pairs of distance < 5.0 Angstrom)

[illegible]

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



Str-1 in thick and Str-2 in thin wireframes; Residues with  $d < 5\text{\AA}$  in red)



- Click [A182408.pdb](#) to download the first structure that you submitted.
- Click [B182408.pdb](#) to download the second structure that you submitted.
- Click [C182408.pdb](#) to download the superposed structure in C-alpha trace (This file is in a Rasmol script format, you can run 'rasmol -script C182408.pdb' to view the structure).
- Click [D182408.pdb](#) to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D182408.pdb' to view the structure).

- 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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