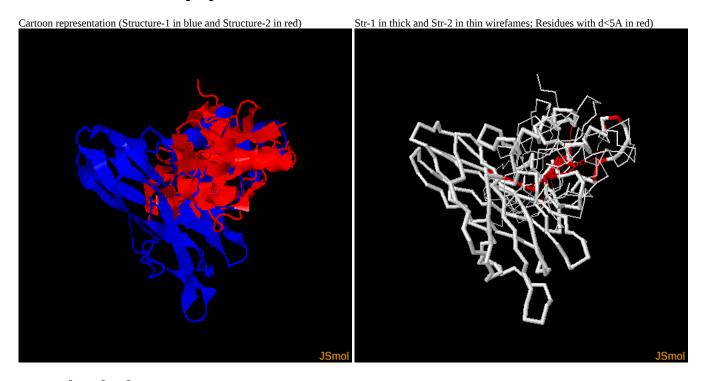
4/15/2018 TM-score Server

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
                        Length= 212
Length= 209 (by which all scores are normalized)
Structure1: A220475
Structure2: B220475
Number of residues in common=
                               74
RMSD of the common residues=
                                11.897
           = 0.1022
                      (d0 = 5.38)
TM-score
MaxSub-score= 0.0441 (d0= 3.50)
GDT-TS-score= 0.0658 %(d<1)=0.0239 %(d<2)=0.0335 %(d<4)=0.0766 %(d<8)=0.1292
GDT-HA-score= 0.0383 %(d<0.5)=0.0191 %(d<1)=0.0239 %(d<2)=0.0335 %(d<4)=0.0766
 ----- rotation matrix to rotate Chain-1 to Chain-2 -----
                                                        u(i.3)
                         u(i.1)
                                        u(i.2)
      -46.4519540137 -0.2758608536
                                      0.5023295515
                                                      0.8194911904
1
       21.6539349064
                     -0.8444908903
                                      0.2805190795
       39.9245972342
                      -0.4590598279
                                     -0.8179083492
                                                     0.3468284975
Superposition in the TM-score: Length(d<5.0)= 13 RMSD= 2.91 (":" denotes the residue pairs of distance < 5.0 Angstrom)
.
PVANADVVFDFONYTAKAGDEVTVDVLVDSKNKPISADVKFKVDSPLTIEEIDKESLAFNTTVTNAILGANFKSLDDKGEPLVPKDGAAVFTLYVNVPANTPDGTYYVGFNGKNEVHKSNDGSQFTVASKNGAITV
```

Visualization of TM-score superposition



Document downloads

- Click <u>A220475.pdb</u> to download the first structure that you submitted.
- Click <u>B220475.pdb</u> to download the second structure that you submitted.
- Click <u>C220475.pdb</u> to download the superposed structure in C-alpah trace (This file is in a Rasmol script format, you can run 'rasmol -script C220475.pdb' to view the structure).
- Click <u>D220475.pdb</u> to download the superposed structure in full-atom (This file is in a Rasmol script format, you can run 'rasmol -script D220475.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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