S1 Table

Stiffness values in MWCENM

| Connection type | Stiffness ratio | Cutoff condition |
|----------------------------------|-----------------|--|
| Backbone (covalent) | 100 | Residue number (between ith and [i+1]th) |
| Backbone (non-bonded) | 1 | Residue number (between ith and [i+2/i+3]th) |
| Disulfide bond | 100 | PDB information |
| Hydrogen bond | 10 | HBPLUS * |
| Salt-bridge | 10 | Distance between charged residues < 4Å |
| Van der Waals force [†] | 1 | Non-bonded distance < 4Å |
| | $(4/d_{i,j})^8$ | $4\text{\AA} \leq \text{Nonbonded distance, } d_{i,j} < 8\text{\AA}$ |

*HBPLUS is a program used to calculate all possible hydrogen bonds within a protein.

[†]The stiffness ratio for the second range of Van der Waals force interaction is only fitted by the attractive term of the Lennard-Jones potential.