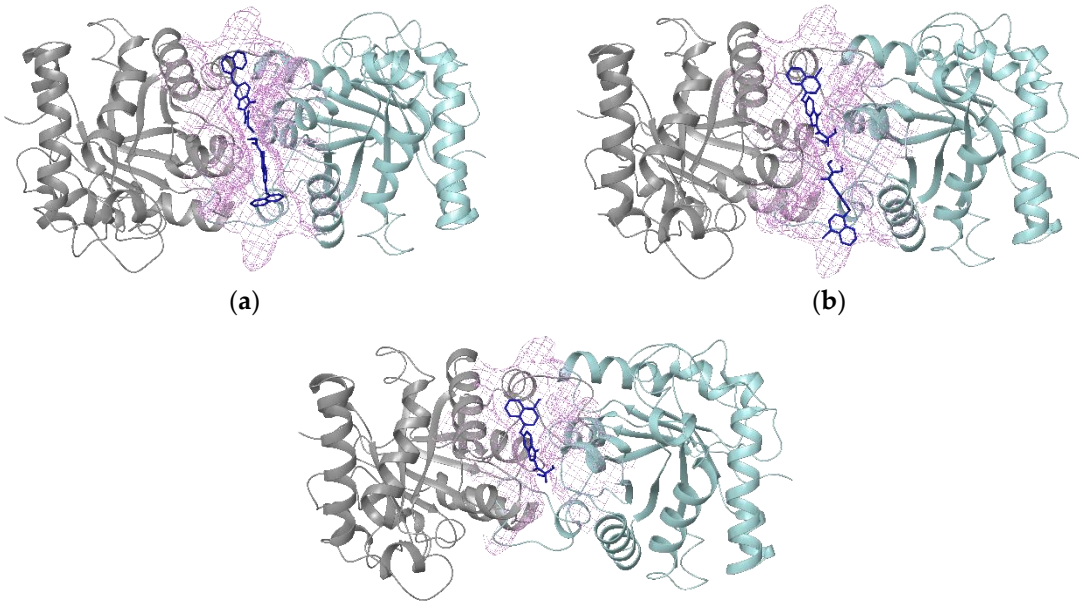
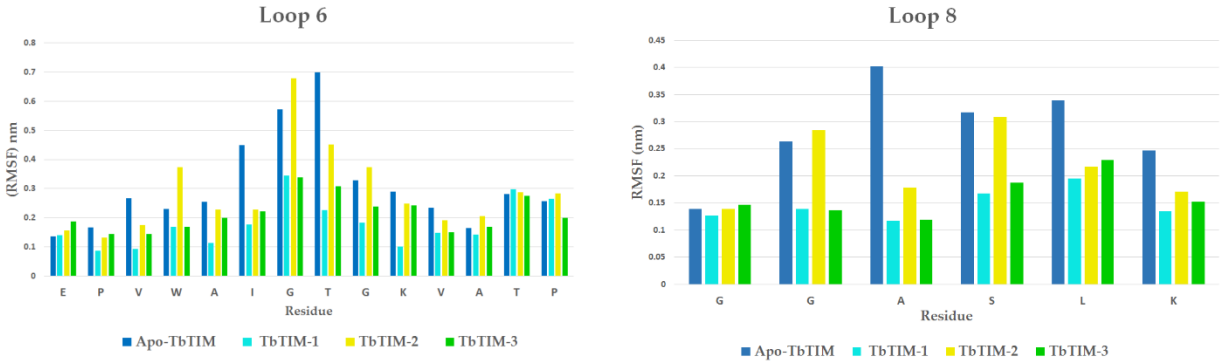


(**a**) (**b**)

(**c**)

Figure S1. Complete view of the binding mode of compounds (**a**) **1**, (**b**) **2**, and (**c**) **3**, at the dimer interface of TbTIM predicted by docking. For compounds **1** and **2** the two molecules are displayed. Molecular surface is drawn as a purple grid.



(a) **(b)**

Figure S2. RMSF comparison between Apo-TbTIM and the complexes with compounds **1**, **2** and **3**. (**a**) Loop 6 residues. (**b**) Loop 8 residues.