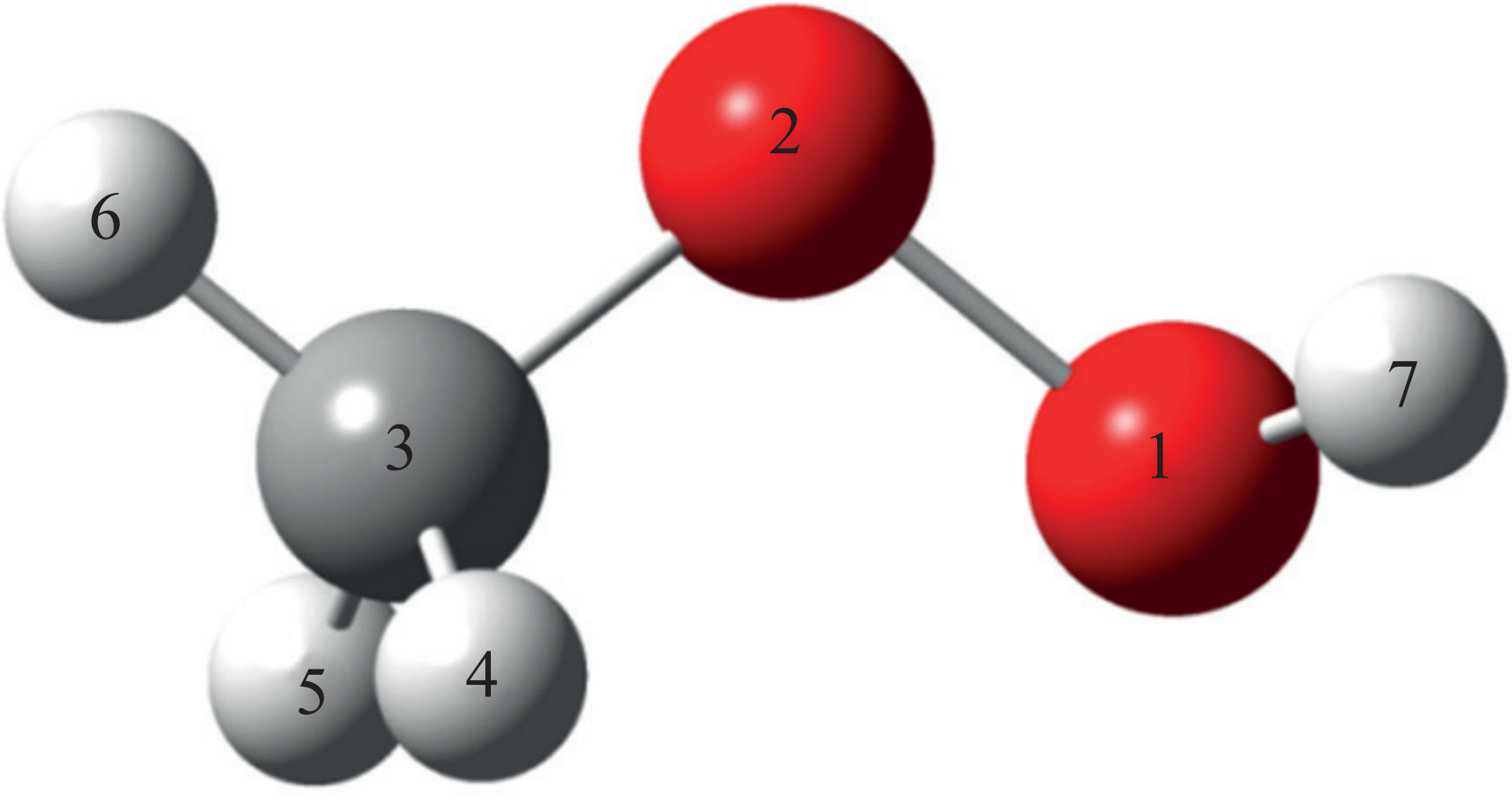

Fig. 2. Equilibrium configuration of the MHP molecule, calculated at the CCSD(T)/Aug-cc-pVTZ level of theory