

Valley splitting of monolayer $\text{Hf}_3\text{C}_2\text{O}_2$ by spin-orbit coupling effect:

First principle calculation using the *HSE06* methods

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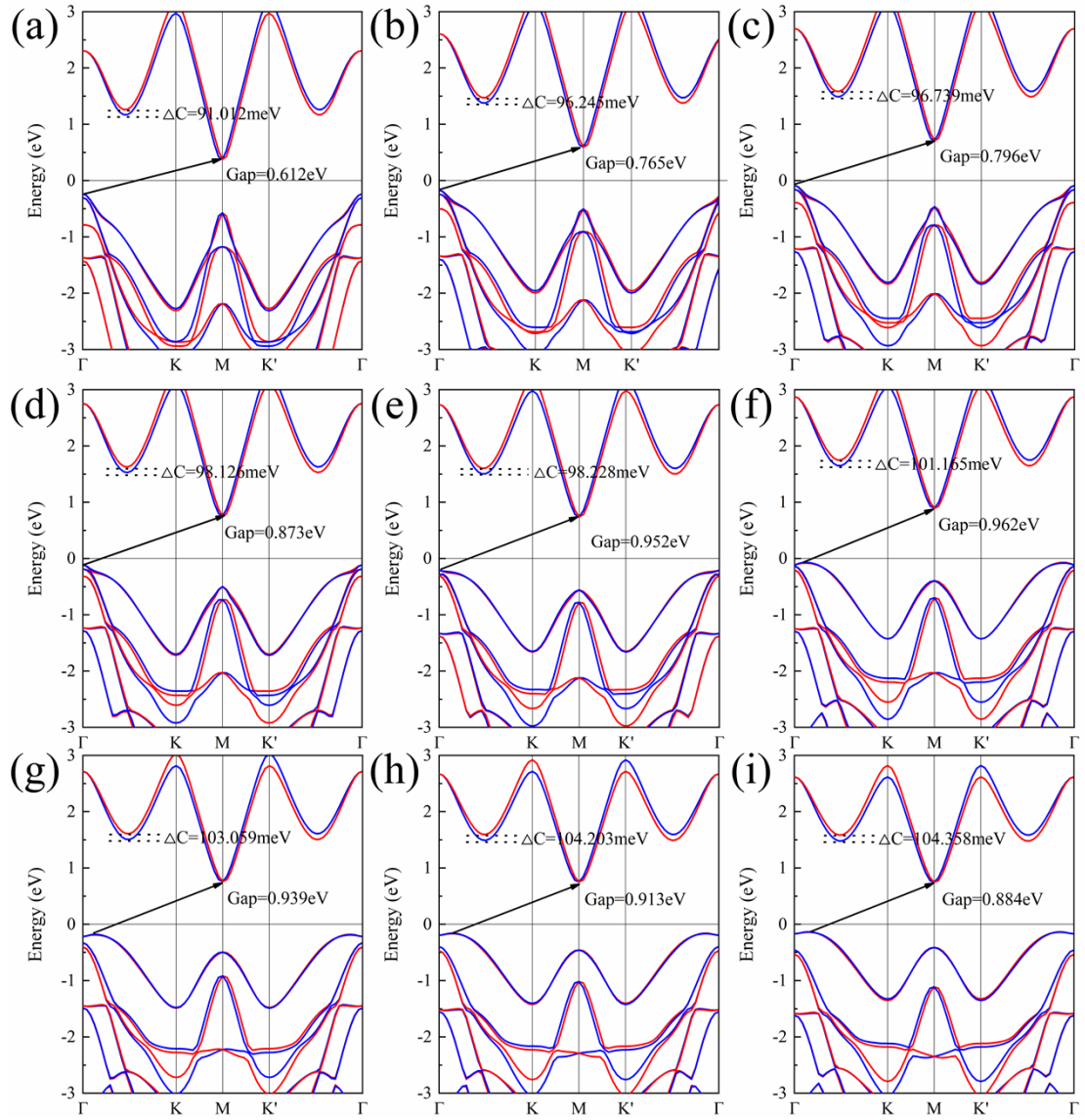


Figure S1. (a)~(i) The band structure of $1L$ $\text{Hf}_3\text{C}_2\text{O}_2$ under biaxial strain from -4% to 4%.

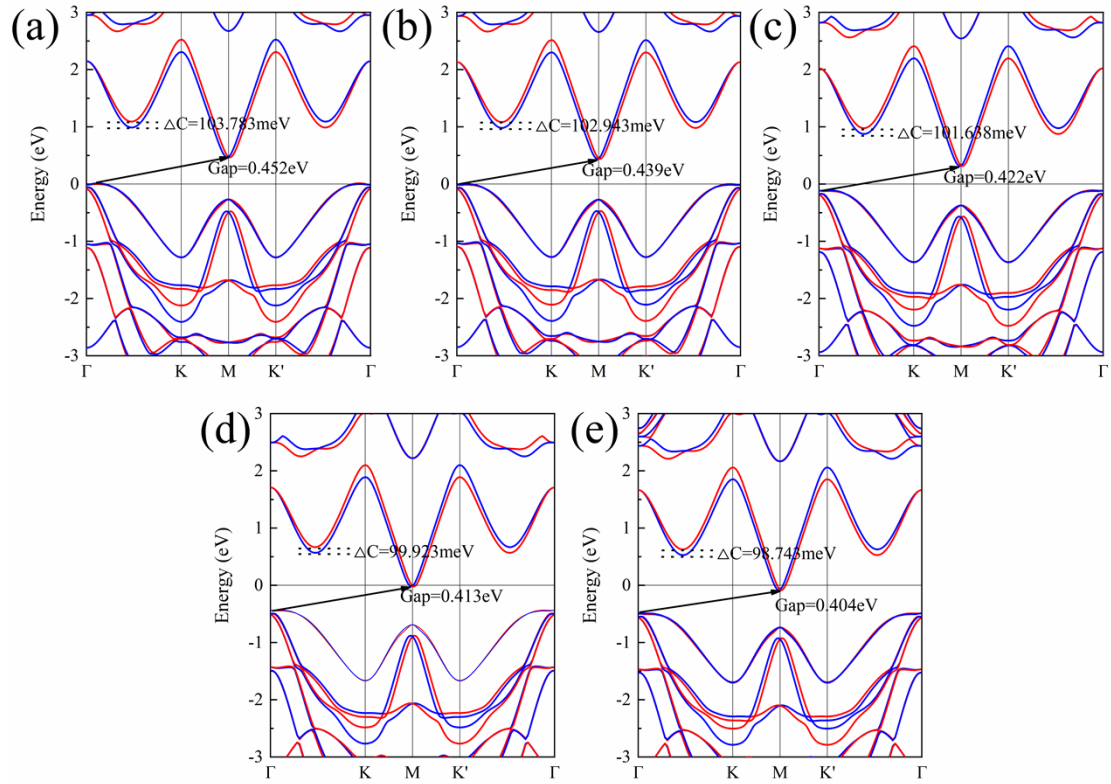


Figure S2. (a)~(e) The band structure of $IL\text{Hf}_3\text{C}_2\text{O}_2$ at doping concentrations ranging from -0.1 electron/unit cell to 0.1 electron/unit cell. (The interval is 0.05 electron/unit cell)