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From Normal Insulator to Topological Insulator in Plumbene XIANG-LONG YU, JIANSHENG WU, South University of Science and Technology of China, Shenzhen — Plumbene, similar to silicene, has a buckled honeycomb structure with a larger band gap ($\sim 400 \text{ meV}$). However, all previous studies have shown that it is a normal insulator. Here, we perform first-principles calculations and employ a sixteen-band tight-binding model with nearest-neighbor and next-nearest-neighbor hopping terms to investigate electronic structures and topological properties of plumbene monolayer. We find that it can become a topological insulator with a large bulk gap ($\sim 200 \text{ meV}$) through electron doping, and the nontrivial state is very robust with respect to external strain. Plumbene can be an ideal candidate for realizing quantum spin Hall effect at room temperature. By investigating effects of external electric and magnetic fields on electronic structures and transport properties of plumbene, we present two rich phase diagrams with and without electron doping, and propose a theoretical design for a four-state spin-valley filter.

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