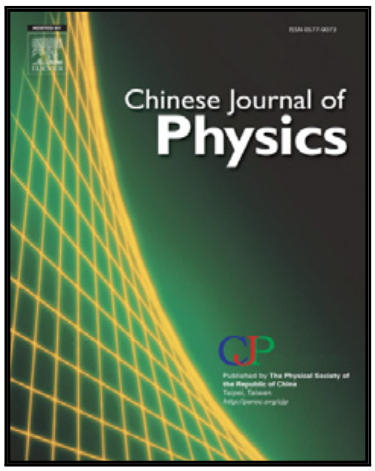

First-principle calculations of structural, electronic and magnetic investigations of $Mn_2RuGe_{1-x}Sn_x$ quaternary Heusler alloys

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Highlights

- DFT calculations for Mn_2RuGe, Mn_2RuSn and their quaternary $Mn_2RuGe_{1-x}Sn_x$ Heusler alloys.
- Structural and magnetic investigation for Mn_2RuGe and Mn_2RuSn.
- Half metallic behavior is explained via spin polarization results.
- Electronic properties for quaternary Heusler alloys were studied.