Theoretical insights into the structural, optoelectronic, thermoelectric, and thermodynamic behavior of novel quaternary LiZrCoX (X = Ge, Sn) compounds based on first-principles study

MOSTAFA MOHAMED SALAHELDIN ABDELKHALEK ELEWA ,Meena Kumari;Jisha Annie Abraham;Ramesh Sharma;Debidatta Behera;S. K. Mukherjee;Murefah mana Al-Anazyf ;Mohammed S. Alqahtanig

Abstract

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