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Computational investigations of the band-structure, thermodynamic and optical features of thorium-based oxide ThGeO_4 using the FP-APW+LO approach

F. Chiker, H. Khachai, C. Mathieu, S. Bin-Omran, Belkacem Kada, Xiao-Wei Sun, Sandeep, D.P. Rai, R. Khenata

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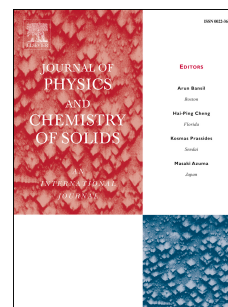
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F. Chiker ^a, H. Khachai ^a, C. Mathieu ^b, S. Bin-Omran ^c, Belkacem Kada ^d, Xiao-Wei Sun ^{e, *}, Sandeep ^f, D. P. Rai ^g, R. Khenata ^{h,*}

^a *Laboratory of Study of Materials and Instruments of Optics, Faculty of Sciences Exacts, University of Djilali Liabes, Sidi Bel Abbes, Algeria*

^b *Faculty of Jean Perrin, University of Artois, SP 18-62307 Lens Cedex, France*

^c *Department of Physics and Astronomy, College of Science, King Saud University, P.O. Box 2455, Riyadh 11451, Saudi Arabia*

^d *Department of Aeronautical Engineering, Faculty of Engineering, King Abdulaziz University, PO. Box 80204, 21589, Jeddah, Saudi Arabia*

^e *School of Mathematics and Physics, Lanzhou Jiaotong University, Lanzhou 730070, China*

^f *Department of Physics, Mizoram University, Aizawl, Mizoram 796004, India*

^g *Department of Physics, Pachhunga University College, Aizawl-796001, India*

^h *Laboratoire de Physique Quantique et de Modélisation Mathématique de la Matière (LPQ3M), Université de Mascara, 29000 Mascara, Algeria*

* Authors to whom correspondence should be addressed

E-mail addresses: sunxw_lzjtu@yeah.net (X.-W. Sun), khenata_rabah@yahoo.fr (R. Khenata)

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