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CaTiO<sub>3</sub> And ...Molecular Dynamics Simulation In Perovskites CaTiO<sub>3</sub> And SrTiO<sub>3</sub> J. A. Souza<sup>1</sup> And J. P. Rino<sup>1</sup>, <sup>1</sup>Universidade Federal De São Carlos, DF, São Carlos, SP, Brazil We Propose An Effective Interatomic Potential In Order To Describe The Perovskite-type ATiO<sub>3</sub> (A = Ca, Apr 1th, 2024.

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