

Lattice dynamic stability and electronic structures of ternary hydrides  $\text{La}_{1-x}\text{Y}_x\text{H}_3$  via first-principles cluster expansion

by

Prutthipong Tsuppayakorn-ae

Wiwittawin Sukmas

Prayoosak Pluengphon

Burapat Inceesungvorn

Piya Phansuke

Pungtip Kaewtubtim

Rajeev Ahuja

Thiti Bovornratanaraks

Wei Luo

RSC Advances 41, 2022

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