Ab initio study of the pressure dependence of mechanic GeB2O4 (B = Mg, Zn and Cd) spinel crystals

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Citation Report

#	Article	IF	Citations
1	First-principles calculations to investigate elastic properties, ideal tensile strength and electronic properties of TiSi, Ti5Si3 and Ti5Si4. Chemical Physics Letters, 2022, 806, 139992.	2.6	2
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