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Titolo	Pressure-Induced Phase Transitions in AB ₂ X ₄ Chalcogenide Compounds // edited by Francisco Javier Manjon, Ion Tiginyanu, Veaceslav Ursaki
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Note generali	Description based upon print version of record.
Nota di bibliografia	Includes bibliographical references at the end of each chapters and index.
Nota di contenuto	Introduction -- Part I Spinel-Structured AB Chalcogenide Compounds -- Part II Ordered-Vacancy AB Chalcogenide Compounds -- Part III Chalcogenide Compounds with other Types of Structures.
Sommario/riassunto	This book on pressure-induced phase transitions in AB ₂ X ₄ chalcogenide compounds deals with one important AmBnXp material. The interest in these materials is caused by their properties. The results are discussed for three main groups of structural families: cubic-spinel structures, defective tetragonal structures, and other structures like layered and wurtzite-type modifications. A systematic analysis of the behavior of cubic (spinel), tetragonal (defect chalcopyrites and stannites) and other crystal modifications of AB ₂ X ₄ compounds under hydrostatic pressure is performed. The behavior of Al ₂ Al ₂ S ₄ , Al ₂ Ga ₂ S ₄ , Al ₂ Al ₂ Se ₄ and Al ₂ Ga ₂ Se ₄ compounds with defective tetragonal structures, compounds with layered and wurtzite structures under

hydrostatic pressure and the pressure dependence of the band gap, lattice parameters, interatomic distances, vibrational modes and pressure-induced phase transitions is discussed. Many of these compounds, except oxide spinels, undergo a pressure-induced phase transition towards the rocksalt-type structure. The phase transition is preceded by disorder in the cation sublattice. The dependence of the transition pressure to the rocksalt-type structure as a function of the compound ionicity and the size criterion is analyzed. At high pressures, all ordered-vacancy compounds are found to exhibit a band anticrossing between several conduction bands that leads to a strong decrease of its pressure coefficient and consequently to a strong non-linear pressure dependence of the direct bandgap energy. Theoretical studies of phase transitions in several ordered-vacancy compounds reveal that the existence of ordered vacancies alter the cation-anion bond distances and their compressibilities. The book is written for students, Ph D. students and specialists in materials science, phase transitions and new materials.
