Reassigning the Pressure-Induced Phase Transitions of Perovskite

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Citation Report

#	Article	IF	CITATIONS
1	Hybrid perovskites under pressure: Present and future directions. Journal of Applied Physics, 2022, 132, .	2.5	4
2	Residual Strain Evolution Induced by Crystallization Kinetics During Antiâ€Solvent Spin Coating in Organic–Inorganic Hybrid Perovskite. Advanced Science, 2023, 10, .	11.2	5
3	Novel high-temperature phase and crystal structure evolution of CsCuBr ₃ halide identified by neutron powder diffraction. CrystEngComm, 2023, 25, 4417-4426.	2.6	2
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5	Pressure-Induced Phase Transition versus Amorphization in Hybrid Methylammonium Lead Bromide Perovskite. Journal of Physical Chemistry C, 2023, 127, 12821-12826.	3.1	5
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8	Building Blockâ€Inspired Hybrid Perovskite Derivatives for Ferroelectric Channel Layers with Gate‶unable Memory Behavior. Angewandte Chemie, 0, , .	2.0	0
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10	Lattice and Electronic Structural Evolutions in Compressed Multilayer MnPS ₃ . Journal of Physical Chemistry C, 2023, 127, 17186-17193.	3.1	0
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13	Halogen Substitution Regulates High Temperature Dielectric Switch in Leadâ€Free Chiral Hybrid Perovskites. Chemistry - A European Journal, 2024, 30, .	3.3	0
14	Boosting Photoluminescence in MAPbBr ₃ Single Crystals through Laser-Based Surface Modification. ACS Photonics, 0, , .	6.6	0
15	Cation Dynamics as Structure Explorer in Hybrid Perovskites─The Case of MAPbI ₃ . Crystal Growth and Design, 0, , .	3.0	0
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17	Probing Local Structural Phase Transition at the Surface of (BA) ₂ PbI ₄ via Interlayer Exciton Emission. Advanced Functional Materials, 0, , .	14.9	0
18	Mechanical Properties and Temperature and Pressure Effects on the Crystal Structure of (2S,3S,4R)-2,3,4-(Trinitratomethyl)-1-nitroazetidine (TMNA) by Single-Crystal X-ray Diffractometry and Density Functional Theory, Journal of Physical Chemistry C, 2024, 128, 927-940.	3.1	0

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19	Pressure-induced band gap engineering and enhanced optoelectronic properties of non-toxic Ca-based perovskite CsCaCl3: Insights from density functional theory. Computational Condensed Matter, 2024, 38, e00879.	2.1	2
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21	Computational highâ€pressure chemistry: Ab initio simulations of atoms, molecules, and extended materials in the gigapascal regime. Wiley Interdisciplinary Reviews: Computational Molecular Science, 2024, 14, .	14.6	0
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