

Wolfram Hergert

List of Publications by Year in descending order

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29

papers

703

citations

567281

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docs citations

30

times ranked

1230

citing authors

#	ARTICLE	IF	CITATIONS
1	Exchange coupling in transition metal monoxides: Electronic structure calculations. <i>Physical Review B</i> , 2009, 80, .	3.2	107
2	Defect-Induced Magnetism in Solids. <i>IEEE Transactions on Magnetics</i> , 2013, 49, 4668-4674.	2.1	87
3	Oxygen-vacancy-induced local ferromagnetism as a driving mechanism in enhancing the magnetic response of ferrites. <i>Physical Review B</i> , 2014, 89, .	3.2	80
4	Numerical solution of the relativistic single-site scattering problem for the Coulomb and the Mathieu potential. <i>Journal of Physics Condensed Matter</i> , 2015, 27, 435202.	1.8	40
5	Magnetic properties of defect-free and oxygen-deficient cubic $\text{Sr}_{2-\delta}\text{FeMoO}_6$. <i>Physical Review B</i> , 2015, 92, .	3.2	32
6	Magnetic and Electronic Properties of Complex Oxides from First-principles. <i>Physica Status Solidi (B): Basic Research</i> , 2020, 257, 1900671.	1.5	31
7	Toward Versatile $\text{Sr}_2\text{FeMoO}_6$ -Based Spintronics by Exploiting Nanoscale Defects. <i>ACS Applied Materials & Interfaces</i> , 2016, 8, 20440-20447.	8.0	30
8	Defect-induced Magnetism in Nonmagnetic Oxides: Basic Principles, Experimental Evidence, and Possible Devices with ZnO and TiO_2 . <i>Physica Status Solidi (B): Basic Research</i> , 2020, 257, 1900623.	1.5	26
9	Room-temperature pressure-induced surface ferromagnetism: First-principles study. <i>Physical Review B</i> , 2011, 84, .	3.2	22
10	First-principles study of uniaxial strained and bent ZnO wires. <i>Physical Review B</i> , 2014, 89, .	3.2	22
11	Variation of magnetic properties of $\text{Sr}_2\text{FeMoO}_6$ due to oxygen vacancies. <i>Journal of Physics Condensed Matter</i> , 2018, 30, 305801.	1.8	21
12	Effect of hydrostatic pressure and uniaxial strain on the electronic structure of $\text{Pb}_1\text{Mo}_{18}$. <i>Physical Review B</i> , 2015, 92, .	3.2	18
13	Chromium point defects in hexagonal BaTiO_3 . A comparative study of first-principles calculations and experiments. <i>Physical Review B</i> , 2015, 91, .	3.2	17
14	GTPack: A Mathematica Group Theory Package for Application in Solid-State Physics and Photonics. <i>Frontiers in Physics</i> , 2018, 6, .	2.1	17
15	Study of charged defects for substitutionally doped chromium in hexagonal barium titanate from first-principles theory. <i>Physica Status Solidi - Rapid Research Letters</i> , 2014, 8, 527-531.	2.4	15
16	Orientation-dependent electron-energy-loss spectroscopy of TiO_2 : A comparison of theory and experiment. <i>Physical Review B</i> , 2006, 73, .	3.2	14
17	Quasicrystals and their Approximants in 2D Ternary Oxides. <i>Physica Status Solidi (B): Basic Research</i> , 2020, 257, 1900624.	1.5	13
18	Analysis of the scattering properties of 3D non-spherical plasmonic nanoparticles accounting for non-local effects. <i>Journal of Modern Optics</i> , 2018, 65, 1778-1786.	1.3	11

#	ARTICLE	IF	CITATIONS
19	Producing ZnFe ₂ O ₄ thin films from ZnO/FeO multilayers. <i>Applied Surface Science</i> , 2017, 393, 256-261.	6.1	10
20	Titanium 3d ferromagnetism with perpendicular anisotropy in defective anatase. <i>Physical Review B</i> , 2020, 101, .	3.2	10
21	First-principles investigations of the magnetic phase diagram of $Gd_{x}Sr_{y}O_{z}$. <i>Physical Review B</i> , 2019, 99, .		
22	Elastic anomalies and long/short-range ordering effects: A first-principles investigation of the Ag _x Pd _{1-x} solid solution. <i>Physical Review B</i> , 2012, 86, .	3.2	8
23	Application of Generalized Mie Theory to EELS Calculations as a Tool for Optimization of Plasmonic Structures. <i>Plasmonics</i> , 2016, 11, 865-874.	3.4	7
24	Tuning the probability of defect formation via substrate strains in $Sr_xO_{2.4}$ films. <i>Physical Review Materials</i> , 2018, 2, .		
25	Theoretical investigation of iron incorporation in hexagonal barium titanate. <i>Physical Review B</i> , 2019, 100, .	3.2	6
26	Influence of surface band bending on a narrow band gap semiconductor: Tunneling atomic force studies of graphite with Bernal and rhombohedral stacking orders. <i>Physical Review Materials</i> , 2021, 5, .	2.4	5
27	<i>Ab initio</i> study of the p-hole magnetism at polar surfaces of ZnO: the role of correlations. <i>Journal of Physics Condensed Matter</i> , 2016, 28, 016003.	1.8	4
28	Electronic and Magnetic Properties of BaFeO ₃ on the Pt(111) Surface in a Quasicrystalline Approximant Structure. <i>Physica Status Solidi (B): Basic Research</i> , 2020, 257, 1900649.	1.5	4
29	Magnetic Structure of Bulk GdMnO ₃ : Influence of Strain. <i>Physica Status Solidi (B): Basic Research</i> , 2020, 257, 1900632.	1.5	3