

Alexander I Poteryaev

List of Publications by Year in descending order

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17
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840728

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times ranked

1026
citing authors

#	ARTICLE	IF	CITATIONS
1	Dynamical mean-field theory using Wannier functions: A flexible route to electronic structure calculations of strongly correlated materials. <i>Physical Review B</i> , 2006, 74, . Enhanced crystal-field splitting and orbital-selective coherence induced by strong correlations in V_2O_3 . <i>Physical Review B</i> , 2007, 76, .	3.2	177
2	Calculated phonon spectra of paramagnetic iron at the orbital-selective transition. <i>Physical Review B</i> , 2012, 85, .	3.2	129
3	Orbital-selective formation of local moments in Fe_2O_3 -iron: First-principles route to an effective model. <i>Physical Review B</i> , 2010, 81, .	3.2	97
4	Monoclinic M_1 phase of VO M_2 Mott-Hubbard versus band insulator. <i>Physical Review B</i> , 2012, 85, .	3.2	73
5	Calculation of photoemission spectra of the doped Mott insulator using LDA+DMFT(QMC). <i>European Physical Journal B</i> , 2000, 18, 55-61.	1.5	63
6	Effect of crystal-field splitting and interband hybridization on the metal-insulator transitions of strongly correlated systems. <i>Physical Review B</i> , 2008, 78, .	3.2	52
7	Unexpected 3+ valence of iron in FeO ₂ , a geologically important material lying between oxides and peroxides. <i>Scientific Reports</i> , 2017, 7, 13005.	3.3	47
8	Valence-band satellite in ferromagnetic nickel: LDA+DMFT study with exact diagonalization. <i>Physical Review B</i> , 2012, 85, .	3.2	28
9	Rotationally invariant exchange interaction: The case of paramagnetic iron. <i>Physical Review B</i> , 2012, 86, .	3.2	25
10	Magnetostriction and ferroelectric state in $AgCrS_2$. <i>Journal of Physics Condensed Matter</i> , 2015, 27, 165601.	1.8	11
11	Magnetic properties of Fe_2O_3 from CPA+DMFT perspectives. <i>Physical Review B</i> , 2016, 93, .	3.2	11
12	Hydrogenation-driven formation of local magnetic moments in FeO_2 . <i>Physical Review B</i> , 2018, 98, .	3.2	11
13	Electronic correlations and competing orders in multiorbital dimers: A cluster DMFT study. <i>Physical Review B</i> , 2019, 99, .	3.2	8
14	Specific heat of a binary alloy within the CPA+DMFT method. <i>Physical Review B</i> , 2015, 91, .	3.2	7
15	Momentum-resolved spectroscopy of correlated metals: A view from dynamical mean field theory. <i>Comptes Rendus Physique</i> , 2009, 10, 537-547.	0.9	5
16	Influence of Molecular Orbitals on Magnetic Properties of FeO ₂ Hx. <i>Molecules</i> , 2020, 25, 2211.	3.8	4