

Konstantin Z Rushchanskii

List of Publications by Year in descending order

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15
papers

423
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840776

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times ranked

684
citing authors

#	ARTICLE	IF	CITATIONS
1	Ordering of Oxygen Vacancies and Related Ferroelectric Properties in HfO_2 . Physical Review Letters, 2021, 127, 087602.	7.8	31
2	Unconventional Ferroelectric Switching via Local Domain Wall Motion in Multiferroic Fe_2O_3 Films. Advanced Electronic Materials, 2020, 6, 1901134.	5.1	11
3	<i>Ab initio</i> phase diagrams of HfO , ZrO and Y_2O_3 : a comparative study. Faraday Discussions, 2019, 213, 321-337.	3.2	27
4	Phase-change memories (PCM) – Experiments and modelling: general discussion. Faraday Discussions, 2019, 213, 393-420.	3.2	7
5	Routes for increasing endurance and retention in HfO_2 -based resistive switching memories. Physical Review Materials, 2018, 2, .	2.4	15
6	Melting of Pb Charge Glass and Simultaneous Pb–Cr Charge Transfer in PbCrO_3 as the Origin of Volume Collapse. Journal of the American Chemical Society, 2015, 137, 12719-12728.	13.7	45
7	First-principles prediction of oxygen octahedral rotations in perovskite-structure EuTiO_3 . Physical Review B, 2012, 85, .	3.2	55
8	<i>Ab Initio</i> Phonon Structure of h-YMnO_3 in Low-Symmetry Ferroelectric Phase. Ferroelectrics, 2012, 426, 90-96.	0.6	4
9	Vacancy-mediated diffusion in biaxially strained Si. Applied Physics Letters, 2011, 98, 031908.	3.3	11
10	A multiferroic material to search for the permanent electric dipole moment of the electron. Nature Materials, 2010, 9, 649-654.	27.5	88
11	Stress enhanced self-diffusion in Si: Entropy effect in anisotropic elastic environment. Applied Physics Letters, 2008, 92, 152110.	3.3	6
12	Ferroelectricity, Nonlinear Dynamics, and Relaxation Effects in Monoclinic SnP_2S_6 . Physical Review Letters, 2007, 99, 207601.	7.8	54
13	Influence of pressure on the structural, dynamical, and electronic properties of the SnP_2S_6 layered crystal. Physical Review B, 2006, 73, .	3.2	23
14	Similarities of the band structure of In_4Se_3 and InSe under pressure and peculiarities of the creation of the band gap. Physica Status Solidi (B): Basic Research, 2006, 243, 592-609.	1.5	35
15	Band structure calculations on the layered compounds FeGa_2S_4 and NiGa_2S_4 . Journal of Physics and Chemistry of Solids, 2002, 63, 2019-2028.	4.0	13