

Khuong P Ong

List of Publications by Year in descending order

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28
papers

2,027
citations

304743

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docs citations

29
times ranked

3964
citing authors

#	ARTICLE	IF	CITATIONS
1	Origin of giant electric-field-induced strain in faulted alkali niobate films. <i>Nature Communications</i> , 2022, 13, .	12.8	11
2	Giant piezoelectricity in oxide thin films with nanopillar structure. <i>Science</i> , 2020, 369, 292-297.	12.6	86
3	Multi Band Gap Electronic Structure in CH ₃ NH ₃ PbI ₃ . <i>Scientific Reports</i> , 2019, 9, 2144.	3.3	26
4	Growth of centimeter-sized [(CH ₃) ₃ NH ₂] ₂ [Mn(HCOO) ₃] hybrid formate perovskite single crystals and Raman evidence of pressure-induced phase transitions. <i>New Journal of Chemistry</i> , 2017, 41, 151-159.	2.8	31
5	Perspective: <i>n</i> -type oxide thermoelectrics via visual search strategies. <i>APL Materials</i> , 2016, 4, .	5.1	42
6	Transparent conducting properties of SrSnO ₃ and ZnSnO ₃ . <i>APL Materials</i> , 2015, 3, 062505.	5.1	65
7	Enhanced photovoltaic effects and switchable conduction behavior in BiFe0.6Sc0.4O ₃ thin films. <i>Acta Materialia</i> , 2015, 88, 83-90.	7.9	37
8	Stable Ferroelectric Perovskite Structure with Giant Axial Ratio and Polarization in Epitaxial BiFe _{0.6} Ga _{0.4} O ₃ Thin Films. <i>ACS Applied Materials & Interfaces</i> , 2015, 7, 2648-2653.	8.0	38
9	Mechanical Origin of the Structural Phase Transition in Methylammonium Lead Iodide CH ₃ NH ₃ PbI ₃ . <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 681-685.	4.6	63
10	Stabilization of 4H hexagonal phase in gold nanoribbons. <i>Nature Communications</i> , 2015, 6, 7684.	12.8	215
11	Ferroelectricity of CH ₃ NH ₃ PbI ₃ Perovskite. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 1155-1161.	4.6	295
12	Structural Evolution in Methylammonium Lead Iodide CH ₃ NH ₃ PbI ₃ . <i>Journal of Physical Chemistry A</i> , 2015, 119, 11033-11038.	2.5	66
13	Strain effects on the band gap and optical properties of perovskite SrSnO ₃ and BaSnO ₃ . <i>Applied Physics Letters</i> , 2014, 104, .	3.3	108
14	Simultaneous reduction in leakage current and enhancement in magnetic moment in BiFeO ₃ nanofibers via optimized Sn doping. <i>Physica Status Solidi - Rapid Research Letters</i> , 2014, 8, 653-657.	2.4	13
15	Structural Instability of Epitaxial (001) BiFeO ₃ Thin Films under Tensile Strain. <i>Scientific Reports</i> , 2014, 4, 4631.	3.3	27
16	Three-dimensional magnetism and coupling to the conduction electrons in PdCrO ₃ . <i>Physical Review B</i> , 2012, 85, .	3.2	14
17	Origin of a Tetragonal BiFeO ₃ Phase with a Giant <i>c/a</i> Ratio on SrTiO ₃ Substrates. <i>Advanced Functional Materials</i> , 2012, 22, 937-942.	14.9	61
18	Analysis of the thermoelectric properties of ZnO. <i>Physical Review B</i> , 2011, 83, .	3.2	265

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19	Origin of anisotropy and metallic behavior in delafossite PdCoO_2 . Physical Review B, 2010, 81, .	3.2	49
20	The Influence of Out-of-Plane Deformation on the Band Gap of Graphene Nanoribbons. Journal of Physical Chemistry C, 2010, 114, 12749-12753.	3.1	21
21	Unusual Transport and Strongly Anisotropic Thermopower in PtCoO_2 and PdCoO_2 . Physical Review Letters, 2010, 104, 176601.	7.8	82
22	Relative effects of all chemical elements on the electrical conductivity of metal and alloys: An alternative to Norbury-Linde rule. Journal of Alloys and Compounds, 2009, 478, 345-354.	5.5	15
23	Electrical conductivity and performance of doped LaCrO_3 perovskite oxides for solid oxide fuel cells. Journal of Power Sources, 2008, 176, 82-89.	7.8	167
24	Realizing the semiconducting state of delafossite AgFeO_2 by GGA+U calculations. Journal of Alloys and Compounds, 2008, 449, 366-370.	5.5	20
25	Origin of the light green color and electronic ground state of LaCrO_3 . Physical Review B, 2008, 77, .	3.2	63
26	Optimization of electrical conductivity of LaCrO_3 through doping: A combined study of molecular modeling and experiment. Applied Physics Letters, 2007, 90, 044109.	3.3	26
27	Electronic Structure and Optical Properties of AFeO_2 ($\text{A} = \text{Ag}, \text{Cu}$) within GGA Calculations. Chemistry of Materials, 2007, 19, 634-640.	6.7	95
28	Electronic properties of $\text{A}_{1-x}\text{Fe}_x\text{O}_2$ -site substituted lead zirconate titanate: Density functional calculations. Physical Review B, 2007, 76, .	3.2	25