

Anna Miglio

List of Publications by Year in descending order

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19
papers

1,544
citations

623188

14
h-index

794141

19
g-index

19
all docs

19
docs citations

19
times ranked

2418
citing authors

#	ARTICLE	IF	CITATIONS
1	Identification and design principles of low hole effective mass p-type transparent conducting oxides. Nature Communications, 2013, 4, 2292.	5.8	507
2	How Does Chemistry Influence Electron Effective Mass in Oxides? A High-Throughput Computational Analysis. Chemistry of Materials, 2014, 26, 5447-5458.	3.2	127
3	Band widths and gaps from the Tran-Blaha functional: Comparison with many-body perturbation theory. Physical Review B, 2013, 87, .	1.1	125
4	High-Mobility Bismuth-based Transparent <i>p</i> -Type Oxide from High-Throughput Material Screening. Chemistry of Materials, 2016, 28, 30-34.	3.2	118
5	Phase Diagrams and Stability of Lead-Free Halide Double Perovskites Cs ₂ BB ₂ X ₆ : B = Sb and Bi, B ²⁺ = Cu, Ag, and Au, and X = Cl, Br, and I. Journal of Physical Chemistry C, 2018, 122, 158-170.	1.5	114
6	High-Throughput Design of Non-oxide p-Type Transparent Conducting Materials: Data Mining, Search Strategy, and Identification of Boron Phosphide. Chemistry of Materials, 2017, 29, 2568-2573.	3.2	109
7	Superionic Diffusion through Frustrated Energy Landscape. Chem, 2019, 5, 2450-2460.	5.8	92
8	Predominance of non-adiabatic effects in zero-point renormalization of the electronic band gap. Npj Computational Materials, 2020, 6, .	3.5	65
9	Quasiparticles and phonon satellites in spectral functions of semiconductors and insulators: Cumulants applied to the full first-principles theory and the Fröhlich polaron. Physical Review B, 2018, 97, .	1.1	60
10	First-principles study of Ce -doped silicate nitride phosphors: Neutral excitation, Stokes shift, and luminescent center identification. Physical Review B, 2016, 93, .	1.1	49
11	First-principles study of the luminescence of Eu -doped phosphors. Physical Review B, 2017, 96, .	1.1	49
12	Assessment of First-Principles and Semiempirical Methodologies for Absorption and Emission Energies of Ce ³⁺ -Doped Luminescent Materials. Advanced Optical Materials, 2017, 5, 1600997.	3.6	35
13	Local Bonding Influence on the Band Edge and Band Gap Formation in Quaternary Chalcopyrites. Advanced Science, 2017, 4, 1700080.	5.6	35
14	<i>Ab initio</i> study of luminescence in Ce-doped Lu_2O_3 : The role of oxygen vacancies on emission color and thermal quenching behavior. Physical Review Materials, 2018, 2, .	0.9	16
15	<i>Ab-initio</i> study of oxygen vacancy stability in bulk and Cerium-doped lutetium oxyorthosilicate. Journal of Luminescence, 2018, 204, 499-505.	1.5	13
16	Beyond the one-dimensional configuration coordinate model of photoluminescence. Physical Review B, 2019, 100, .	1.1	10
17	Importance of Long-Range Channel Sr Displacements for the Narrow Emission in Sr ₂ Al ₂ O ₂ N ₂ :Eu ²⁺ Phosphor. Advanced Optical Materials, 2021, 9, 2100649.	3.6	10
18	Fröhlich polaron effective mass and localization length in cubic materials: Degenerate and anisotropic electronic bands. Physical Review B, 2021, 104, .	1.1	8

#	ARTICLE	IF	CITATIONS
19	Design rule for the emission linewidth of Eu ²⁺ http://www.w3.org/1998/Math/MathML" display="inline" id="d1e1136" altimg="si27.svg"><math>2</math>+</math>-activated phosphors. <i>Journal of Luminescence</i>, 2020, 224, 117258.	1.5	7