

Hongping Li

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Study on the pyrolysis mechanism of unsaturated fatty acid: A combined density functional theory and experimental study. <i>International Journal of Energy Research</i> , 2022, 46, 2029-2040.	4.5	10
2	Facile synthesized fluorine substituted benzothiadiazole based dopant-free hole transport material for high efficiency perovskite solar cell. <i>Dyes and Pigments</i> , 2021, 184, 108786.	3.7	15
3	The impact of B-site antisite defects on the magnetic and electronic properties in double perovskite Pb ₂ FeOsO ₆ . <i>Ceramics International</i> , 2021, 47, 992-1001.	4.8	2
4	Insight into the Mechanism of Glycerol Dehydration and Subsequent Pyridine Synthesis. <i>ACS Sustainable Chemistry and Engineering</i> , 2021, 9, 3095-3103.	6.7	23
5	Passivation functionalized phenothiazine-based hole transport material for highly efficient perovskite solar cell with efficiency exceeding 22%. <i>Chemical Engineering Journal</i> , 2021, 410, 128328.	12.7	83
6	Spin Polarization-Assisted Dopant Segregation at a Coherent Phase Boundary. <i>ACS Nano</i> , 2021, 15, 19938-19944.	14.6	6
7	Highly efficient perovskite solar cells based on symmetric hole transport material constructed with indaceno[1,2-b:5,6-b']dithiophene core building block. <i>Journal of Energy Chemistry</i> , 2020, 43, 98-103.	12.9	31
8	Fluorine-Substituted Benzotriazole Core Building Block-Based Highly Efficient Hole-Transporting Materials for Mesoporous Perovskite Solar Cells. <i>Solar Rrl</i> , 2020, 4, 1900362.	5.8	16
9	Benzo[1,2- <i>c</i> : <i>i</i> :4,5- <i>c</i> : <i>i</i>]-dithiophene-4,8-dione (BDD) Core Building Block Based Dopant-Free Hole-Transport Materials for Efficient and Stable Perovskite Solar Cell. <i>ACS Applied Energy Materials</i> , 2020, 3, 10333-10339.	5.1	3
10	A half-metallic ferrimagnet of CeCu ₃ Cr ₄ O ₁₂ with 4f itinerant electron. <i>Applied Physics Letters</i> , 2020, 117, 132404.	3.3	1
11	Cyclic Compound Formation Mechanisms during Pyrolysis of Typical Aliphatic Acidic Amino Acids. <i>ACS Sustainable Chemistry and Engineering</i> , 2020, 8, 16968-16978.	6.7	32
12	Ferrimagnetic semiconductor of CaCu ₃ Fe ₂ V ₂ O ₁₂ with direct bandgap. <i>Chemical Physics Letters</i> , 2020, 759, 137910.	2.6	0
13	Ferrimagnetic semiconductor with a direct bandgap. <i>Applied Physics Letters</i> , 2020, 116, .	3.3	7
14	Unraveling the effect of B-site antisite defects on the electronic and magnetic properties of the quadruple perovskite CaCu ₃ Fe ₂ Nb ₂ O ₁₂ . <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 3059-3065.	2.8	5
15	Highly efficient phenothiazine 5,5-dioxide-based hole transport materials for planar perovskite solar cells with a PCE exceeding 20%. <i>Journal of Materials Chemistry A</i> , 2019, 7, 9510-9516.	10.3	60
16	Direct Determination of Atomic Structure and Magnetic Coupling of Magnetite Twin Boundaries. <i>ACS Nano</i> , 2018, 12, 2662-2668.	14.6	30
17	Highly Efficient Phenoxazine Core Unit Based Hole Transport Materials for Hysteresis-Free Perovskite Solar Cells. <i>ACS Applied Materials & Interfaces</i> , 2018, 10, 36608-36614.	8.0	41
18	Molecular Engineering of Triphenylamine-Based Non-Fullerene Electron-Transport Materials for Efficient Rigid and Flexible Perovskite Solar Cells. <i>ACS Applied Materials & Interfaces</i> , 2018, 10, 38970-38977.	8.0	34

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19	First-principles study of negative thermal expansion mechanism in A-site-ordered perovskite $\text{SrCu}_3\text{Fe}_4\text{O}_{12}$. RSC Advances, 2015, 5, 1801-1807.	3.6	10
20	A-site ordering induced suppression of magnetic cluster glass and dielectric anomaly in $\text{La}_2\text{BiCoMnO}_6$. Applied Physics Letters, 2012, 100, .	3.3	11
21	Mechanism of A-B intersite charge transfer and negative thermal expansion in A-site-ordered perovskite $\text{LaCu}_3\text{Fe}_4\text{O}_{12}$. Journal of Applied Physics, 2012, 111, 103718.	2.5	17
22	Electrical conductivity optimization in electrolyte-free fuel cells by single-component $\text{Ce}_{0.8}\text{Sm}_{0.2}\text{O}_{2-\delta}$ - $\text{Li}_{0.15}\text{Ni}_{0.45}\text{Zn}_{0.4}$ layer. RSC Advances, 2012, 2, 3828.	3.6	57
23	A-Site-Doping Enhanced B-Site Ordering and Correlated Magnetic Property in $\text{La}_2\text{BiCoMnO}_6$. Journal of Physical Chemistry C, 2012, 116, 16841-16847.	3.1	46
24	First-principle investigation of magnetic coupling mechanism in hypothesized A-site-ordered perovskite $\text{YMn}_3\text{Sc}_4\text{O}_{12}$. Journal of Computational Chemistry, 2012, 33, 82-87.	3.3	4
25	Covalent State and the Electronic and Transport Properties of $\text{CaCu}_3\text{Ni}_4\text{O}_{12}$: A First-Principles Study. Journal of Physical Chemistry C, 2011, 115, 2366-2370.	3.1	10
26	First-principles investigation of A-B intersite charge transfer and correlated electrical and magnetic properties in $\text{BiCu}_3\text{Fe}_4\text{O}_{12}$. Journal of Computational Chemistry, 2011, 32, 1235-1240.	3.3	10
27	Strain control of orbital polarization and correlated metal-insulator transition in $\text{La}_2\text{CoMnO}_6$ from first principles. Applied Physics Letters, 2011, 99, .	3.3	21
28	The peculiar magnetic property evolution along $\text{RCu}_3\text{Mn}_4\text{O}_{12}$ (R=Y, La, Pr, Nd, Sm, Eu, Gd, Tb, Dy, Ho, Er, Tm, Yb, Lu). Journal of Applied Physics, 2011, 110, 074107.	3.3	2