

Yingzhuo Lun

List of Publications by Year in descending order

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Version: 2024-02-01

12
papers

1,299
citations

1307594

7
h-index

1199594

12
g-index

12
all docs

12
docs citations

12
times ranked

2015
citing authors

#	ARTICLE	IF	CITATIONS
1	Cation and anion immobilization through chemical bonding enhancement with fluorides for stable halide perovskite solar cells. <i>Nature Energy</i> , 2019, 4, 408-415.	39.5	831
2	Liquid medium annealing for fabricating durable perovskite solar cells with improved reproducibility. <i>Science</i> , 2021, 373, 561-567.	12.6	227
3	The Spacer Cations Interplay for Efficient and Stable Layered 2D Perovskite Solar Cells. <i>Advanced Energy Materials</i> , 2020, 10, 1901566.	19.5	89
4	Manipulation of current rectification in van der Waals ferroionic CuInP ₂ S ₆ . <i>Nature Communications</i> , 2022, 13, 574.	12.8	60
5	Thickness-Dependent In-Plane Polarization and Structural Phase Transition in van der Waals Ferroelectric CuInP ₂ S ₆ . <i>Small</i> , 2020, 16, e1904529.	10.0	50
6	Atomically Asymmetric Inversion Scales up to Mesoscopic Single-Crystal Monolayer Flakes. <i>ACS Nano</i> , 2020, 14, 13834-13840.	14.6	11
7	Size-dependent strain-engineered nanostructures in MoS ₂ monolayer investigated by atomic force microscopy. <i>Nanotechnology</i> , 2021, 32, 465703.	2.6	8
8	Van der Waals direction transformation induced by shear strain in layered PdSe ₂ . <i>Extreme Mechanics Letters</i> , 2021, 44, 101231.	4.1	7
9	Screening piezoelectricity in determination of flexoelectric coefficient at nanoscale. <i>Mechanics of Materials</i> , 2020, 150, 103591.	3.2	6
10	Asymmetric mechanical properties in ferroelectrics driven by flexo-deformation effect. <i>Journal of the Mechanics and Physics of Solids</i> , 2022, 164, 104891.	4.8	6
11	Visualization of Strain-Engineered Nanopattern in Center-Confined Mesoscopic WS ₂ Monolayer Flakes. <i>Journal of Physical Chemistry C</i> , 2022, 126, 7184-7192.	3.1	3
12	Near-zero Poisson's ratio and suppressed mechanical anisotropy in strained black phosphorene/SnSe van der Waals heterostructure: a first-principles study. <i>Applied Mathematics and Mechanics (English)</i> Tj ETQq0 0 0 88BT /Overlock 10 Tf		