

Zhaochuan Fan

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2151553/publications.pdf>

Version: 2024-02-01

17
papers

450
citations

759233

12
h-index

888059

17
g-index

19
all docs

19
docs citations

19
times ranked

873
citing authors

#	ARTICLE	IF	CITATIONS
1	A multiscale study of misfit dislocations in PbTe/PbSe(001) heteroepitaxy. Journal of Materials Research, 2019, 34, 2306-2314.	2.6	10
2	Controlling Nanoparticle Orientations in the Self-Assembly of Patchy Quantum Dot-Gold Heterostructural Nanocrystals. Journal of the American Chemical Society, 2019, 141, 6013-6021.	13.7	49
3	Thermally Induced Wurtzite to hBN Structural Transition. Physica Status Solidi - Rapid Research Letters, 2019, 13, 1800584.	2.4	3
4	Orientational Order in Self-Assembled Nanocrystal Superlattices. Journal of the American Chemical Society, 2019, 141, 1980-1988.	13.7	52
5	Ab initio phase diagram of PbSe crystals calculated with the random phase approximation. Physical Review B, 2018, 98, .	3.2	2
6	Self-Assembly of Quantum Dot-Gold Heterodimer Nanocrystals with Orientational Order. Nano Letters, 2018, 18, 5049-5056.	9.1	25
7	Fracture behaviors of brittle and ductile 2D carbon structures under uniaxial tensile stress. Carbon, 2017, 111, 486-492.	10.3	59
8	Atomistic understanding of cation exchange in PbS nanocrystals using simulations with pseudoligands. Nature Communications, 2016, 7, 11503.	12.8	48
9	Heat-induced transformation of CdSe-CdS-ZnS core-multishell quantum dots by Zn diffusion into inner layers. Chemical Communications, 2015, 51, 3320-3323.	4.1	20
10	Crystal structure, stability, and electronic properties of hydrated metal sulfates $MSO_4(H_2O)_n$ ($M=Ni$). Tj ETQq0 0 0 rgBT /Overlock 10 Tf 77-86.	3.8	13
11	Core-shell reconfiguration through thermal annealing in $Fe_xO/CoFe_2O_4$ ordered 2D nanocrystal arrays. Nanotechnology, 2014, 25, 055601.	2.6	9
12	A transferable force field for CdS-CdSe-PbS-PbSe solid systems. Journal of Chemical Physics, 2014, 141, 244503.	3.0	19
13	New Ab Initio Based Pair Potential for Accurate Simulation of Phase Transitions in ZnO. Journal of Physical Chemistry C, 2014, 118, 11050-11061.	3.1	45
14	Atomic Resolution Monitoring of Cation Exchange in CdSe-PbSe Heteronanocrystals during Epitaxial Solid-Vapor Growth. Nano Letters, 2014, 14, 3661-3667.	9.1	48
15	From Sphere to Multipod: Thermally Induced Transitions of CdSe Nanocrystals Studied by Molecular Dynamics Simulations. Journal of the American Chemical Society, 2013, 135, 5869-5876.	13.7	19
16	First-principles calculation of vibrational properties of $B_{12}As_2$ crystal. Physica Status Solidi (B): Basic Research, 2011, 248, 1242-1247.	1.5	7
17	First principles calculations of the vibrational properties of icosahedral solid boron oxygen $B_{12}O_2$. Physica B: Condensed Matter, 2011, 406, 297-303.	2.7	18