

Xiaokun Gu

List of Publications by Year in descending order

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48
papers

3,597
citations

218677

26
h-index

223800

46
g-index

48
all docs

48
docs citations

48
times ranked

4856
citing authors

#	ARTICLE	IF	CITATIONS
1	Flexible n-type thermoelectric materials by organic intercalation of layered transition metal dichalcogenide TiS ₂ . Nature Materials, 2015, 14, 622-627.	27.5	612
2	Measurement Techniques for Thermal Conductivity and Interfacial Thermal Conductance of Bulk and Thin Film Materials. Journal of Electronic Packaging, Transactions of the ASME, 2016, 138, .	1.8	328
3	Phonon transport in single-layer transition metal dichalcogenides: A first-principles study. Applied Physics Letters, 2014, 105, .	3.3	309
4	<i>Colloquium</i>: Phononic thermal properties of two-dimensional materials. Reviews of Modern Physics, 2018, 90, .	45.6	238
5	First-principles prediction of phononic thermal conductivity of silicene: A comparison with graphene. Journal of Applied Physics, 2015, 117, .	2.5	204
6	A new regime of nanoscale thermal transport: Collective diffusion increases dissipation efficiency. Proceedings of the National Academy of Sciences of the United States of America, 2015, 112, 4846-4851.	7.1	170
7	Probing Anisotropic Thermal Conductivity of Transition Metal Dichalcogenides MX ₂ (M = Tj ETQq1 1 0.784314 rgBT / Over 21.0 163	21.0	163
8	Layer thickness-dependent phonon properties and thermal conductivity of MoS ₂ . Journal of Applied Physics, 2016, 119, .	2.5	136
9	Simultaneous measurement of thermal conductivity and heat capacity of bulk and thin film materials using frequency-dependent transient thermoreflectance method. Review of Scientific Instruments, 2013, 84, 034902.	1.3	120
10	Lattice thermal conductivity of organic-inorganic hybrid perovskite CH ₃ NH ₃ PbI ₃ . Applied Physics Letters, 2016, 108, .	3.3	97
11	Bottom-up Design of Three-Dimensional Carbon-Honeycomb with Superb Specific Strength and High Thermal Conductivity. Nano Letters, 2017, 17, 179-185.	9.1	95
12	Temperature Dependence of Anisotropic Thermal Conductivity Tensor of Bulk Black Phosphorus. Advanced Materials, 2017, 29, 1603297.	21.0	89
13	Revealing the Origins of 3D Anisotropic Thermal Conductivities of Black Phosphorus. Advanced Electronic Materials, 2016, 2, 1600040.	5.1	85
14	A Review of Simulation Methods in Micro/Nanoscale Heat Conduction. ES Energy & Environments, 2018, , .	1.1	78
15	Revisiting phonon-phonon scattering in single-layer graphene. Physical Review B, 2019, 100, .	3.2	71
16	Anisotropic Tuning of Graphite Thermal Conductivity by Lithium Intercalation. Journal of Physical Chemistry Letters, 2016, 7, 4744-4750.	4.6	69
17	Dielectric Mismatch Mediates Carrier Mobility in Organic-Intercalated Layered TiS ₂ . Nano Letters, 2015, 15, 6302-6308.	9.1	62
18	The energy efficiency of interfacial solar desalination. Applied Energy, 2021, 302, 117581.	10.1	60

#	ARTICLE	IF	CITATIONS
19	PHONON TRANSPORT AND THERMAL CONDUCTIVITY IN TWO-DIMENSIONAL MATERIALS. Annual Review of Heat Transfer, 2016, 19, 1-65.	1.0	57
20	On the influence of junction structures on the mechanical and thermal properties of carbon honeycombs. Carbon, 2017, 119, 278-286.	10.3	56
21	Stable planar single-layer hexagonal silicene under tensile strain and its anomalous Poisson's ratio. Applied Physics Letters, 2014, 104, 081902.	3.3	49
22	Unification of nonequilibrium molecular dynamics and the mode-resolved phonon Boltzmann equation for thermal transport simulations. Physical Review B, 2020, 101, .	3.2	49
23	Thermal conductivity of single-layer $\text{MoS}_2(1-x)\text{Se}_2x$ alloys from molecular dynamics simulations with a machine-learning-based interatomic potential. Computational Materials Science, 2019, 165, 74-81.	3.0	46
24	Thermal conductivity prediction by atomistic simulation methods: Recent advances and detailed comparison. Journal of Applied Physics, 2021, 130, .	2.5	36
25	Thermal conductivity modeling of hybrid organic-inorganic crystals and superlattices. Nano Energy, 2017, 41, 394-407.	16.0	32
26	Anisotropic thermal transport in van der Waals layered alloys $\text{WSe}_2(1-x)\text{Te}_2x$. Applied Physics Letters, 2018, 112, .	3.3	32
27	High-temperature phonon transport properties of SnSe from machine-learning interatomic potential. Journal of Physics Condensed Matter, 2021, 33, 405401.	1.8	24
28	Anomalous thermal transport in metallic transition-metal nitrides originated from strong electron-phonon interactions. Materials Today Physics, 2020, 15, 100256.	6.0	22
29	Thermal conductivity of silicon at elevated temperature: Role of four-phonon scattering and electronic heat conduction. International Journal of Heat and Mass Transfer, 2020, 160, 120165.	4.8	21
30	Phonon transport in single-layer $\text{Mo}_x\text{W}_x\text{S}_2$ alloy embedded with WS_2 nanodomains. Physical Review B, 2016, 94, .	3.2	18
31	Thermal conductivity of $\text{MoS}_2/\text{MoSe}_2$ heterostructures: The role of lattice mismatch, interlayer rotation and species intermixing. International Journal of Heat and Mass Transfer, 2019, 143, 118583.	4.8	17
32	Phonon transmission across Mg/MgO interface: A first-principles-based atomistic Green's function study. Physical Review B, 2015, 91, .	3.2	16
33	Anisotropic Thermal Transport in Organic-Inorganic Hybrid Crystal $\text{ZnTe}(\text{en})_{0.5}$. Journal of Physical Chemistry C, 2015, 119, 28300-28308.	3.1	16
34	Thermal conductivity of dielectric nanowires with different cross-section shapes. Chinese Physics B, 2007, 16, 3777-3782.	1.3	15
35	Effect of the accuracy of interatomic force constants on the prediction of lattice thermal conductivity. Computational Materials Science, 2017, 138, 368-376.	3.0	15
36	Seeking for Low Thermal Conductivity Atomic Configurations in SiGe Alloys with Bayesian Optimization. ES Energy & Environments, 2020, , .	1.1	14

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37	Shape dependence of slip length on patterned hydrophobic surfaces. Applied Physics Letters, 2011, 99, .	3.3	13
38	Thermal conductivity of hexagonal Si, Ge, and Si _{1-x} Ge _x alloys from first-principles. Journal of Applied Physics, 2018, 123, .	2.5	12
39	Impact of thermostat on interfacial thermal conductance prediction from non-equilibrium molecular dynamics simulations. Chinese Physics B, 2022, 31, 056301.	1.4	10
40	A minimal Tersoff potential for diamond silicon with improved descriptions of elastic and phonon transport properties. Journal of Physics Condensed Matter, 2020, 32, 135901.	1.8	9
41	Monitoring anharmonic phonon transport across interfaces in one-dimensional lattice chains. Physical Review E, 2020, 101, 022133.	2.1	8
42	A scattering rate model for accelerated evaluation of lattice thermal conductivity bypassing anharmonic force constants. Journal of Applied Physics, 2019, 125, .	2.5	6
43	GPU_PBTE: an efficient solver for three and four phonon scattering rates on graphics processing units. Journal of Physics Condensed Matter, 2021, 33, 495901.	1.8	6
44	Electronic band structure of carbon honeycombs. Materials Today Physics, 2018, 5, 72-77.	6.0	5
45	Black Phosphorus: Revealing the Origins of 3D Anisotropic Thermal Conductivities of Black Phosphorus (Adv. Electron. Mater. 5/2016). Advanced Electronic Materials, 2016, 2, .	5.1	4
46	A New Regime of Nanoscale Thermal Transport: Collective Diffusion Counteracts Dissipation Inefficiency. Springer Proceedings in Physics, 2015, , 341-344.	0.2	3
47	Mechanical and thermal properties of nanomaterials at sub-50nm dimensions characterized using coherent EUV beams. , 2015, , .		0
48	HIGH TEMPERATURE THERMAL CONDUCTIVITY OF SILICON FROM MACHINE-LEARNING-BASED INTERATOMIC POTENTIAL. , 2018, , .		0