

Andrea Cepellotti

List of Publications by Year in descending order

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Version: 2024-02-01

17
papers

2,943
citations

687220

13
h-index

940416

16
g-index

18
all docs

18
docs citations

18
times ranked

5266
citing authors

#	ARTICLE	IF	CITATIONS
1	Two-dimensional materials from high-throughput computational exfoliation of experimentally known compounds. <i>Nature Nanotechnology</i> , 2018, 13, 246-252.	15.6	1,317
2	Thermal Conductivity of Graphene and Graphite: Collective Excitations and Mean Free Paths. <i>Nano Letters</i> , 2014, 14, 6109-6114.	4.5	449
3	AiiDA: automated interactive infrastructure and database for computational science. <i>Computational Materials Science</i> , 2016, 111, 218-230.	1.4	399
4	Phonon hydrodynamics in two-dimensional materials. <i>Nature Communications</i> , 2015, 6, 6400.	5.8	385
5	Perspective on <i>ab initio</i> phonon thermal transport. <i>Journal of Applied Physics</i> , 2019, 126, .	1.1	76
6	Thermal Transport in Crystals as a Kinetic Theory of Relaxons. <i>Physical Review X</i> , 2016, 6, .	2.8	72
7	Optomechanical Measurement of Thermal Transport in Two-Dimensional MoSe ₂ Lattices. <i>Nano Letters</i> , 2019, 19, 3143-3150.	4.5	43
8	Rippling ultrafast dynamics of suspended 2D monolayers, graphene. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, E6555-E6561.	3.3	41
9	Boltzmann Transport in Nanostructures as a Friction Effect. <i>Nano Letters</i> , 2017, 17, 4675-4682.	4.5	38
10	Generalization of Fourier's Law into Viscous Heat Equations. <i>Physical Review X</i> , 2020, 10, .	2.8	36
11	Bayesian force fields from active learning for simulation of inter-dimensional transformation of stanene. <i>Npj Computational Materials</i> , 2021, 7, .	3.5	33
12	Transport waves as crystal excitations. <i>Physical Review Materials</i> , 2017, 1, .	0.9	20
13	NH ₃ -NO Coadsorption System on Pt(111). I. Structure of the Mixed Layer. <i>Journal of Physical Chemistry C</i> , 2013, 117, 21186-21195.	1.5	14
14	NH ₃ -NO Coadsorption System on Pt(111). II. Intermolecular Interaction. <i>Journal of Physical Chemistry C</i> , 2013, 117, 21196-21202.	1.5	10
15	Interband tunneling effects on materials transport properties using the first principles Wigner distribution. <i>Materials Today Physics</i> , 2021, 19, 100412.	2.9	9
16	On the Kinetic Theory of Thermal Transport in Crystals. , 2018, , 1-42.		0
17	On the Kinetic Theory of Thermal Transport in Crystals. , 2020, , 767-808.		0