

Taku Watanabe

List of Publications by Year in descending order

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13
papers

1,942
citations

687363

13
h-index

1058476

14
g-index

14
all docs

14
docs citations

14
times ranked

2618
citing authors

#	ARTICLE	IF	CITATIONS
1	High-energy long-cycling all-solid-state lithium metal batteries enabled by silver-carbon composite anodes. <i>Nature Energy</i> , 2020, 5, 299-308.	39.5	932
2	Accelerating Applications of Metal-Organic Frameworks for Gas Adsorption and Separation by Computational Screening of Materials. <i>Langmuir</i> , 2012, 28, 14114-14128.	3.5	202
3	Thermal transport and grain boundary conductance in ultrananocrystalline diamond thin films. <i>Journal of Applied Physics</i> , 2006, 99, 114301.	2.5	139
4	Energetics of intrinsic point defects in uranium dioxide from electronic-structure calculations. <i>Journal of Nuclear Materials</i> , 2009, 384, 61-69.	2.7	127
5	Accurate Treatment of Electrostatics during Molecular Adsorption in Nanoporous Crystals without Assigning Point Charges to Framework Atoms. <i>Journal of Physical Chemistry C</i> , 2011, 115, 4824-4836.	3.1	106
6	Thermal transport properties of uranium dioxide by molecular dynamics simulations. <i>Journal of Nuclear Materials</i> , 2008, 375, 388-396.	2.7	87
7	Molecular chemisorption on open metal sites in $\text{Cu}_3(\text{benzenetricarboxylate})_2$: A spatially periodic density functional theory study. <i>Journal of Chemical Physics</i> , 2010, 133, 094509.	3.0	87
8	Computational identification of a metal organic framework for high selectivity membrane-based CO_2/CH_4 separations: $\text{Cu}(\text{hfipbb})(\text{H}_2\text{hfipbb})_0.5$. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 11389.	2.8	83
9	Thermal conductance across grain boundaries in diamond from molecular dynamics simulation. <i>Journal of Applied Physics</i> , 2007, 102, 063503.	2.5	47
10	Thermal Transport in Off-Stoichiometric Uranium Dioxide by Atomic Level Simulation. <i>Journal of the American Ceramic Society</i> , 2009, 92, 850-856.	3.8	47
11	Thermal transport in polyethylene and at polyethylene-diamond interfaces investigated using molecular dynamics simulation. <i>Journal of Physics Condensed Matter</i> , 2009, 21, 084219.	1.8	40
12	Highly Cyclable All-Solid-State Battery with Deposition-Type Lithium Metal Anode Based on Thin Carbon Black Layer. <i>Advanced Energy and Sustainability Research</i> , 2021, 2, 2100066.	5.8	23
13	Toward an Atomistically Informed Fuel Performance Code: Thermal Properties Using FRAPCON and Molecular Dynamics Simulation. <i>Nuclear Technology</i> , 2009, 165, 308-312.	1.2	6