

Nils Edvin Richard Zimmermann

List of Publications by Year in descending order

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Version: 2024-02-01

21
papers

1,501
citations

516710

16
h-index

752698

20
g-index

29
all docs

29
docs citations

29
times ranked

2015
citing authors

#	ARTICLE	IF	CITATIONS
1	Benchmarking Coordination Number Prediction Algorithms on Inorganic Crystal Structures. <i>Inorganic Chemistry</i> , 2021, 60, 1590-1603.	4.0	31
2	Local structure order parameters and site fingerprints for quantification of coordination environment and crystal structure similarity. <i>RSC Advances</i> , 2020, 10, 6063-6081.	3.6	57
3	The Materials Project: Accelerating Materials Design Through Theory-Driven Data and Tools. , 2020, , 1751-1784.		14
4	High-throughput assessment of hypothetical zeolite materials for their synthesizability and industrial deployability. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2019, 234, 437-450.	0.8	12
5	PyCDT: A Python toolkit for modeling point defects in semiconductors and insulators. <i>Computer Physics Communications</i> , 2018, 226, 165-179.	7.5	142
6	Electrostatic Estimation of Intercalant Jump-Diffusion Barriers Using Finite-Size Ion Models. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 628-634.	4.6	16
7	The Materials Project: Accelerating Materials Design Through Theory-Driven Data and Tools. , 2018, , 1-34.		11
8	NaCl nucleation from brine in seeded simulations: Sources of uncertainty in rate estimates. <i>Journal of Chemical Physics</i> , 2018, 148, 222838.	3.0	62
9	Matminer: An open source toolkit for materials data mining. <i>Computational Materials Science</i> , 2018, 152, 60-69.	3.0	446
10	Quantifying local coordination environment and structural similarity through order parameter-based site fingerprints and their application to machine learning. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2018, 74, a209-a209.	0.1	0
11	Assessing Local Structure Motifs Using Order Parameters for Motif Recognition, Interstitial Identification, and Diffusion Path Characterization. <i>Frontiers in Materials</i> , 2017, 4, .	2.4	54
12	History and Utility of Zeolite Framework-Type Discovery from a Data-Science Perspective. <i>Crystal Growth and Design</i> , 2016, 16, 3043-3048.	3.0	41
13	Transport in Nanoporous Materials Including MOFs: The Applicability of Fick's Laws. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 14580-14583.	13.8	90
14	Nucleation of NaCl from Aqueous Solution: Critical Sizes, Ion-Attachment Kinetics, and Rates. <i>Journal of the American Chemical Society</i> , 2015, 137, 13352-13361.	13.7	151
15	Transport into Nanosheets: Diffusion Equations Put to Test. <i>Journal of Physical Chemistry C</i> , 2013, 117, 7384-7390.	3.1	43
16	Predicting Local Transport Coefficients at Solid-Gas Interfaces. <i>Journal of Physical Chemistry C</i> , 2012, 116, 18878-18883.	3.1	37
17	Surface Barriers of Hydrocarbon Transport Triggered by Ideal Zeolite Structures. <i>Journal of Physical Chemistry C</i> , 2012, 116, 3677-3683.	3.1	46
18	Adsorption and diffusion in zeolites: the pitfall of isotypic crystal structures. <i>Molecular Simulation</i> , 2011, 37, 986-989.	2.0	13

#	ARTICLE	IF	CITATIONS
19	On the Effects of the External Surface on the Equilibrium Transport in Zeolite Crystals. Journal of Physical Chemistry C, 2010, 114, 300-310.	3.1	42
20	Path Sampling Calculation of Methane Diffusivity in Natural Gas Hydrates from a Water-Vacancy Assisted Mechanism. Journal of the American Chemical Society, 2008, 130, 17342-17350.	13.7	124
21	In-Depth Study of the Influence of Host Framework Flexibility on the Diffusion of Small Gas Molecules in One-Dimensional Zeolitic Pore Systems. Journal of Physical Chemistry C, 2007, 111, 17370-17381.	3.1	59