

# Jens JÃ¸rgen Mortensen

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/43398/publications.pdf>

Version: 2024-02-01

10  
papers

2,536  
citations

1163065

8  
h-index

1372553

10  
g-index

10  
all docs

10  
docs citations

10  
times ranked

3954  
citing authors

#	ARTICLE	IF	CITATIONS
1	The atomic simulation environment—a Python library for working with atoms. <i>Journal of Physics Condensed Matter</i> , 2017, 29, 273002.	1.8	1,933
2	Recent progress of the Computational 2D Materials Database (C2DB). <i>2D Materials</i> , 2021, 8, 044002.	4.4	218
3	Time-dependent density-functional theory in the projector augmented-wave method. <i>Journal of Chemical Physics</i> , 2008, 128, 244101.	3.0	187
4	RPBE-vdW Description of Benzene Adsorption on Au(111). <i>Topics in Catalysis</i> , 2010, 53, 378-383.	2.8	57
5	Local Bayesian optimizer for atomic structures. <i>Physical Review B</i> , 2019, 100, .	3.2	57
6	MyQueue: Task and workflow scheduling system. <i>Journal of Open Source Software</i> , 2020, 5, 1844.	4.6	36
7	Efficient first principles simulation of electron scattering factors for transmission electron microscopy. <i>Ultramicroscopy</i> , 2019, 197, 16-22.	1.9	29
8	Interfacing CRYSTAL/AMBER to Optimize QM/MM Lennard-Jones Parameters for Water and to Study Solvation of TiO <sub>2</sub> Nanoparticles. <i>Molecules</i> , 2018, 23, 2958.	3.8	9
9	Numerical quality control for DFT-based materials databases. <i>Npj Computational Materials</i> , 2022, 8, .	8.7	6
10	Computational exfoliation of atomically thin one-dimensional materials with application to Majorana bound states. <i>Physical Review Materials</i> , 2022, 6, .	2.4	4