

# Chiara Panosetti

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1114374/publications.pdf>

Version: 2024-02-01

10  
papers

229  
citations

1684188

5  
h-index

1588992

8  
g-index

10  
all docs

10  
docs citations

10  
times ranked

318  
citing authors

| #  | ARTICLE  | IF  | CITATIONS |
|----|--|-----|-----------|
| 1  | Systematic Comparison of Genetic Algorithm and Basin Hopping Approaches to the Global Optimization of Si(111) Surface Reconstructions. <i>Journal of Physical Chemistry A</i> , 2022, 126, 3043-3056.  | 2.5 | 5         |
| 2  | DFTB Modeling of Lithium-Intercalated Graphite with Machine-Learned Repulsive Potential. <i>Journal of Physical Chemistry A</i> , 2021, 125, 691-699.  | 2.5 | 11        |
| 3  | Accessing Structural, Electronic, Transport and Mesoscale Properties of Li-GICs via a Complete DFTB Model with Machine-Learned Repulsion Potential. <i>Materials</i> , 2021, 14, 6633.                 | 2.9 | 4         |
| 4  | Learning to Use the Force: Fitting Repulsive Potentials in Density-Functional Tight-Binding with Gaussian Process Regression. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 2181-2191. | 5.3 | 27        |
| 5  | Density Functional Tight Binding Modelling of Lithium Intercalated Graphite with Machine-Learned Repulsive Potential. <i>ECS Meeting Abstracts</i> , 2020, MA2020-02, 198-198.                         | 0.0 | 0         |
| 6  | Investigating Diffusion of Lithium Intercalated in Graphite By a Combination of Multiscale Modeling and NMR. <i>ECS Meeting Abstracts</i> , 2020, MA2020-02, 96-96.                                    | 0.0 | 0         |
| 7  | A Practical Guide to Surface Kinetic Monte Carlo Simulations. <i>Frontiers in Chemistry</i> , 2019, 7, 202.  | 3.6 | 154       |
| 8  | Atomic scale switches based on self-assembled surface magic clusters. <i>Applied Physics Letters</i> , 2018, 112, 253103.  | 3.3 | 1         |
| 9  | Global structure search for molecules on surfaces: Efficient sampling with curvilinear coordinates. <i>Journal of Chemical Physics</i> , 2016, 145, 084117.  | 3.0 | 11        |
| 10 | Global Materials Structure Search with Chemically Motivated Coordinates. <i>Nano Letters</i> , 2015, 15, 8044-8048.  | 9.1 | 16        |