

Torsten Hahn

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6697295/publications.pdf>

Version: 2024-02-01

22
papers

420
citations

840776

11
h-index

713466

21
g-index

22
all docs

22
docs citations

22
times ranked

478
citing authors

#	ARTICLE	IF	CITATIONS
1	Crystal Growth, Structure, and Transport Properties of the Charge-Transfer Salt Picene/2,3,5,6-Tetrafluoro-7,7,8,8-tetracyanoquinodimethane. <i>Crystal Growth and Design</i> , 2014, 14, 1338-1346.	3.0	66
2	Stretched or noded orbital densities and self-interaction correction in density functional theory. <i>Journal of Chemical Physics</i> , 2019, 150, 174102.	3.0	46
3	Self-consistent self-interaction corrected density functional theory calculations for atoms using Fermi-Löwdin orbitals: Optimized Fermi-orbital descriptors for Li-Kr. <i>Journal of Chemical Physics</i> , 2017, 147, 164107.	3.0	39
4	Fermi-Löwdin orbital self-interaction corrected density functional theory: Ionization potentials and enthalpies of formation. <i>Journal of Computational Chemistry</i> , 2018, 39, 2463-2471.	3.3	35
5	Hybrid States and Charge Transfer at a Phthalocyanine Heterojunction: $F + \text{MnPc} + \text{CoPc}$ <i>Physical Review Letters</i> , 2012, 109, 027601.	3.2	32
6	Synthesis and properties of new 9,10-anthraquinone derived compounds for molecular electronics. <i>New Journal of Chemistry</i> , 2013, 37, 601-610.	2.8	28
7	Optical and magneto-optical properties of metal phthalocyanine and metal porphyrin thin films. <i>Journal of Physics Condensed Matter</i> , 2014, 26, 104201.	1.8	27
8	Interpretation and Automatic Generation of Fermi-Orbital Descriptors. <i>Journal of Computational Chemistry</i> , 2019, 40, 2843-2857.	3.3	21
9	The Role of Self-Interaction Corrections, Vibrations, and Spin-Orbit in Determining the Ground Spin State in a Simple Heme. <i>Magnetochemistry</i> , 2017, 3, 31.	2.4	18
10	Phthalocyanine dimers in a blend: Spectroscopic and theoretical studies of $\text{MnPc} + \text{F16CoPc}$. <i>Journal of Chemical Physics</i> , 2013, 138, 024707.	3.0	17
11	Analytic atomic gradients in the fermi-Löwdin orbital self-interaction correction. <i>Journal of Computational Chemistry</i> , 2019, 40, 820-825.	3.3	16
12	Electronic properties of the charge transfer material MnPc/F4TCNQ. <i>Journal of Chemical Physics</i> , 2016, 145, 114702.	3.0	11
13	Charge transfer from and to manganese phthalocyanine: bulk materials and interfaces. <i>Beilstein Journal of Nanotechnology</i> , 2017, 8, 1601-1615.	2.8	11
14	Electronic properties of 1,2;8,9-dibenzopentacene thin films: A joint experimental and theoretical study. <i>Physical Review B</i> , 2012, 86, .	3.2	8
15	Electronic properties of the phthalocyanine based dimer: $\text{MnPc} + \text{F16CoPc}$. <i>Journal of Chemical Physics</i> , 2013, 138, 024707.	3.2	8
16	Magnetic Signatures of Hydroxyl- and Water-Terminated Neutral and Tetra-Anionic Mn ₁₂ -Acetate. <i>Journal of Computational Chemistry</i> , 2019, 40, 2301-2308.	3.3	8
17	Silicon nano particles: Surface characterization, defects and electronic properties. <i>Physica Status Solidi (B): Basic Research</i> , 2008, 245, 959-962.	1.5	7
18	A gate controlled molecular switch based on picene-F4TCNQ charge-transfer material. <i>Nanoscale</i> , 2014, 6, 14508-14513.	5.6	6

#	ARTICLE	IF	CITATIONS
19	Symmetry Breaking within Fermiâ€“LÃ¼wdin Orbital Self-Interaction Corrected Density Functional Theory. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 5823-5828.	5.3	6
20	Versatile Simulation Tool and Novel Measurement Method for Electrical Characterization of Semiconductors. <i>Solid State Phenomena</i> , 2009, 156-158, 241-246.	0.3	4
21	Electronic structure, transport, and collective effects in molecular layered systems. <i>Beilstein Journal of Nanotechnology</i> , 2017, 8, 2094-2105.	2.8	3
22	Electronic structure and transport properties of coupled CdS/ZnSe quantum dots. <i>Journal of Physics Condensed Matter</i> , 2021, 33, 125002.	1.8	3