

# 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	Electronic structure and transport properties of coupled CdS/ZnSe quantum dots. <i>Journal of Physics Condensed Matter</i> , 2021, 33, 125002.	1.8	3
2	Magnetic Signatures of Hydroxyl- and Water-Terminated Neutral and Tetra-Anionic Mn <sub>12</sub> -Acetate. <i>Journal of Computational Chemistry</i> , 2019, 40, 2301-2308.	3.3	8
3	Interpretation and Automatic Generation of Fermi-Orbital Descriptors. <i>Journal of Computational Chemistry</i> , 2019, 40, 2843-2857.	3.3	21
4	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
5	Analytic atomic gradients in the fermi-Làwdin orbital self-interaction correction. <i>Journal of Computational Chemistry</i> , 2019, 40, 820-825.	3.3	16
6	Fermi-Làwdin 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
7	Self-consistent self-interaction corrected density functional theory calculations for atoms using Fermi-Làwdin orbitals: Optimized Fermi-orbital descriptors for Li-Kr. <i>Journal of Chemical Physics</i> , 2017, 147, 164107.	3.0	39
8	Symmetry Breaking within Fermi-Làwdin Orbital Self-Interaction Corrected Density Functional Theory. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 5823-5828.	5.3	6
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	Charge transfer from and to manganese phthalocyanine: bulk materials and interfaces. <i>Beilstein Journal of Nanotechnology</i> , 2017, 8, 1601-1615.	2.8	11
11	Electronic structure, transport, and collective effects in molecular layered systems. <i>Beilstein Journal of Nanotechnology</i> , 2017, 8, 2094-2105.	2.8	3
12	Electronic properties of the charge transfer material MnPc/F4TCNQ. <i>Journal of Chemical Physics</i> , 2016, 145, 114702.	3.0	11
13	A gate controlled molecular switch based on picene-F4TCNQ charge-transfer material. <i>Nanoscale</i> , 2014, 6, 14508-14513.	5.6	6
14	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
15	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
16	Phthalocyanine dimers in a blend: Spectroscopic and theoretical studies of MnPc <sup>+</sup> /F16CoPc <sup>-</sup> . <i>Journal of Chemical Physics</i> , 2013, 138, 024707.	3.0	17
17	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
18	Systematic theoretical investigation of the phthalocyanine based dimer: MnPc <sub>n</sub> $\text{xmlns:mml="http://www.w3.org/1998/Math/MathML" display="block">\text{mml:msup}\text{mml:mrow}$ $\text{mml:mrow}\text{mml:mi}^{\wedge}\text{mml:mi}\text{mml:mo}+\text{mml:mo}\text{mml:mrow}\text{mml:msup}\text{mml:math}/\text{F}\text{mml:math}$ $\text{xmlns:mml="http://www.w3.org/1998/Math/MathML" display="block">\text{mml:msub}\text{mml:mrow}$ $\text{mml:mn}16\text{mml:mn}\text{mml:msub}\text{mml:math}\text{CoPc}\text{mml:math}$ $\text{xmlns:mml="http://www.w3.org/1998/Math/MathML" display="block">\text{mml:msup}\text{mml:mrow}$ $\text{mml:math}$	3.2	8

#	ARTICLE	IF	CITATIONS
19	Hybrid States and Charge Transfer at a Phthalocyanine Heterojunction: Physical Review Letters, 2012, 109, 027601.		
20	Electronic properties of 1,2;8,9-dibenzopentacene thin films: A joint experimental and theoretical study. Physical Review B, 2012, 86, .	3.2	8
21	Versatile Simulation Tool and Novel Measurement Method for Electrical Characterization of Semiconductors. Solid State Phenomena, 2009, 156-158, 241-246.	0.3	4
22	Silicon nano particles: Surface characterization, defects and electronic properties. Physica Status Solidi (B): Basic Research, 2008, 245, 959-962.	1.5	7