## Robin Grotjahn

List of Publications by Year in descending order

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933447 1199594 12 853 10 12 citations h-index g-index papers 12 12 12 1063 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	A Look at Realâ€World Transitionâ€Metal Thermochemistry and Kinetics with Local Hybrid Functionals. Israel Journal of Chemistry, 2023, 63, .	2.3	8
2	Reliable TDDFT Protocol Based on a Local Hybrid Functional for the Prediction of Vibronic Phosphorescence Spectra Applied to Tris(2, $2\hat{a}\in^2$ -bipyridine)-Metal Complexes. Journal of Physical Chemistry A, 2021, 125, 7099-7110.	2.5	13
3	Assessment of hybrid functionals for singlet and triplet excitations: Why do some local hybrid functionals perform so well for triplet excitation energies?. Journal of Chemical Physics, 2021, 155, 124108.	3.0	14
4	Electronic States of 2,3-Diamino-1,4-naphthoquinone and Its N-Alkylated Derivatives. Journal of Physical Chemistry C, 2020, 124, 60-69.	3.1	12
5	Two Ï€â€Electrons Make the Difference: From BODIPY to BODIIM Switchable Fluorescent Dyes. Chemistry - A European Journal, 2020, 26, 1422-1428.	3.3	18
6	A Local Hybrid Functional with Wide Applicability and Good Balance between (De)Localization and Left–Right Correlation. Journal of Chemical Theory and Computation, 2020, 16, 5645-5657.	5.3	54
7	Validation of Local Hybrid Functionals for Excited States: Structures, Fluorescence, Phosphorescence, and Vibronic Spectra. Journal of Chemical Theory and Computation, 2020, 16, 5821-5834.	<b>5.</b> 3	21
8	Evaluation of Local Hybrid Functionals for Electric Properties: Dipole Moments and Static and Dynamic Polarizabilities. Journal of Physical Chemistry A, 2020, 124, 8346-8358.	2.5	15
9	TURBOMOLE: Modular program suite for <i>ab initio</i> quantum-chemical and condensed-matter simulations. Journal of Chemical Physics, 2020, 152, 184107.	3.0	616
10	Development and Implementation of Excited-State Gradients for Local Hybrid Functionals. Journal of Chemical Theory and Computation, 2019, 15, 5508-5522.	5 <b>.</b> 3	21
11	Development of a TDDFT-Based Protocol with Local Hybrid Functionals for the Screening of Potential Singlet Fission Chromophores. Journal of Chemical Theory and Computation, 2017, 13, 4984-4996.	5.3	57
12	A novel model for smectic liquid crystals: Elastic anisotropy and response to a steady-state flow. Journal of Chemical Physics, 2016, 145, 164903.	3.0	4