

Wenlan Liu

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7352907/publications.pdf>

Version: 2024-02-01

18
papers

1,010
citations

623734

14
h-index

888059

17
g-index

19
all docs

19
docs citations

19
times ranked

1428
citing authors

#	ARTICLE	IF	CITATIONS
1	Molecular Origin of Balanced Bipolar Transport in Neat Layers of the Emitter CzDBA. <i>Advanced Materials Technologies</i> , 2021, 6, 2000120.	5.8	9
2	Intrinsic efficiency limits in low-bandgap non-fullerene acceptor organic solar cells. <i>Nature Materials</i> , 2021, 20, 378-384.	27.5	257
3	Impact of Acceptor Quadrupole Moment on Charge Generation and Recombination in Blends of IDTâ€Based Nonâ€Fullerene Acceptors with PCE10 as Donor Polymer. <i>Advanced Energy Materials</i> , 2021, 11, 2100839.	19.5	23
4	Reduced Intrinsic Nonâ€Radiative Losses Allow Roomâ€Temperature Triplet Emission from Purely Organic Emitters. <i>Advanced Materials</i> , 2021, 33, e2101844.	21.0	28
5	Chemical Design Rules for Nonâ€Fullerene Acceptors in Organic Solar Cells. <i>Advanced Energy Materials</i> , 2021, 11, 2102363.	19.5	38
6	Chemical Design Rules for Nonâ€Fullerene Acceptors in Organic Solar Cells (<i>Adv. Energy Mater.</i>) Tj ETQq0 0 0 rgBT /Qyerlock_210 Tf 50 5	19.5	210
7	Long-range exciton diffusion in molecular non-fullerene acceptors. <i>Nature Communications</i> , 2020, 11, 5220.	12.8	204
8	Origin of the Ĩâ€“Ĩ Spacing Change upon Doping of Semiconducting Polymers. <i>Journal of Physical Chemistry C</i> , 2018, 122, 27983-27990.	3.1	25
9	A model hamiltonian tuned toward high level <i>ab initio</i> calculations to describe the character of excitonic states in perylenebisimide aggregates. <i>Journal of Computational Chemistry</i> , 2018, 39, 1979-1989.	3.3	14
10	Functionalized Nickel Oxide Hole Contact Layers: Work Function versus Conductivity. <i>ACS Applied Materials & Interfaces</i> , 2017, 9, 39821-39829.	8.0	37
11	Influence of a polarizable surrounding on the electronically excited states of aggregated perylene materials. <i>Journal of Computational Chemistry</i> , 2016, 37, 1601-1610.	3.3	14
12	A general ansatz for constructing quasi-diabatic states in electronically excited aggregated systems. <i>Journal of Chemical Physics</i> , 2015, 143, 084106.	3.0	30
13	Identification of Ultrafast Relaxation Processes As a Major Reason for Inefficient Exciton Diffusion in Perylene-Based Organic Semiconductors. <i>Journal of the American Chemical Society</i> , 2014, 136, 9327-9337.	13.7	56
14	Explicitly correlated internally contracted multireference coupled-cluster singles and doubles theory: ic-MRCCSD(F12â€). <i>Chemical Physics Letters</i> , 2013, 565, 122-127.	2.6	25
15	Ultrafast Exciton Self-Trapping upon Geometry Deformation in Perylene-Based Molecular Aggregates. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 792-796.	4.6	123
16	Comparison of the electronic structure of different peryleneâ€based dyeâ€aggregates. <i>Journal of Computational Chemistry</i> , 2012, 33, 1544-1553.	3.3	55
17	Assessment of TDâ€DFTâ€and TDâ€HFâ€based approaches for the prediction of exciton coupling parameters, potential energy curves, and electronic characters of electronically excited aggregates. <i>Journal of Computational Chemistry</i> , 2011, 32, 1971-1981.	3.3	70
18	Ultrafast Energy Transfer Triggers Ionization Energy Offset Dependence of Quantum Efficiency in Low-bandgap Non-fullerene Acceptor Solar Cells. , 0, , ,		0