

Tahereh Nemati Aram

List of Publications by Year in descending order

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Version: 2024-02-01

21
papers

384
citations

759233

12
h-index

794594

19
g-index

21
all docs

21
docs citations

21
times ranked

336
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 1 | Singlet fission molecules among known compounds: finding a few needles in a haystack. <i>Energy and Environmental Science</i> , 2019, 12, 2412-2416. | 30.8 | 74 |
| 2 | On the Largest Possible Mobility of Molecular Semiconductors and How to Achieve It. <i>Advanced Functional Materials</i> , 2020, 30, 2001906. | 14.9 | 45 |
| 3 | Practical Computation of the Charge Mobility in Molecular Semiconductors Using Transient Localization Theory. <i>Journal of Physical Chemistry C</i> , 2019, 123, 6989-6997. | 3.1 | 40 |
| 4 | Modeling charge transport in high-mobility molecular semiconductors: Balancing electronic structure and quantum dynamics methods with the help of experiments. <i>Journal of Chemical Physics</i> , 2020, 152, 190902. | 3.0 | 33 |
| 5 | Novel thermally activated delayed fluorescence materials by high-throughput virtual screening: going beyond donor-acceptor design. <i>Journal of Materials Chemistry C</i> , 2021, 9, 3324-3333. | 5.5 | 27 |
| 6 | Bright Frenkel Excitons in Molecular Crystals: A Survey. <i>Chemistry of Materials</i> , 2021, 33, 3368-3378. | 6.7 | 22 |
| 7 | High-throughput virtual screening for organic electronics: a comparative study of alternative strategies. <i>Journal of Materials Chemistry C</i> , 2021, 9, 13557-13583. | 5.5 | 20 |
| 8 | Organic materials repurposing, a data set for theoretical predictions of new applications for existing compounds. <i>Scientific Data</i> , 2022, 9, 54. | 5.3 | 16 |
| 9 | Strategies to reduce the dynamic disorder in molecular semiconductors. <i>Materials Horizons</i> , 2020, 7, 2922-2928. | 12.2 | 14 |
| 10 | Modeling of molecular photocells: Application to two-level photovoltaic system with electron-hole interaction. <i>Journal of Chemical Physics</i> , 2016, 145, 124116. | 3.0 | 13 |
| 11 | Simple model of a coherent molecular photocell. <i>Journal of Chemical Physics</i> , 2016, 144, 134102. | 3.0 | 12 |
| 12 | Charge separation in organic solar cells: Effects of Coulomb interaction, recombination and hole propagation. <i>Europhysics Letters</i> , 2016, 115, 18003. | 2.0 | 12 |
| 13 | Predictive Model of Charge Mobilities in Organic Semiconductor Small Molecules with Force-Matched Potentials. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 3494-3503. | 5.3 | 12 |
| 14 | Quantum modeling of two-level photovoltaic systems. <i>EPJ Photovoltaics</i> , 2017, 8, 85503. | 1.6 | 11 |
| 15 | The impact of long-range electron-hole interaction on the charge separation yield of molecular photocells. <i>Journal of Chemical Physics</i> , 2017, 146, 034103. | 3.0 | 10 |
| 16 | Impact of electron-phonon coupling on the quantum yield of photovoltaic devices. <i>Journal of Chemical Physics</i> , 2020, 152, 044109. | 3.0 | 9 |
| 17 | Quantitative Hole Mobility Simulation and Validation in Substituted Acenes. <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 5530-5537. | 4.6 | 7 |
| 18 | Impact of offset energies on the yield of interfacial charge separation in molecular photocells. <i>Journal of Chemical Physics</i> , 2018, 149, 064102. | 3.0 | 4 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 19 | Influence of Fermi velocity engineering on electronic and optical properties of graphene superlattices. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2015, 379, 974-978. | 2.1 | 3 |
| 20 | Quantum Two-Level Model for Excitonic Solar Cells. , 0, , . | | 0 |
| 21 | Feasibility of p-Doped Molecular Crystals as Transparent Conductive Electrodes via Virtual Screening. <i>Chemistry of Materials</i> , 2022, 34, 4050-4061. | 6.7 | 0 |