

Veaceslav Coropceanu

List of Publications by Year in descending order

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132
docs citations

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times ranked

15798
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Charge Transport in Organic Semiconductors. <i>Chemical Reviews</i> , 2007, 107, 926-952. | 47.7 | 3,853 |
| 2 | Charge-Transfer and Energy-Transfer Processes in π -Conjugated Oligomers and Polymers: A Molecular Picture. <i>Chemical Reviews</i> , 2004, 104, 4971-5004. | 47.7 | 2,539 |
| 3 | Molecular Understanding of Organic Solar Cells: The Challenges. <i>Accounts of Chemical Research</i> , 2009, 42, 1691-1699. | 15.6 | 1,291 |
| 4 | Effect of Electronic Polarization on Charge-Transport Parameters in Molecular Organic Semiconductors. <i>Journal of the American Chemical Society</i> , 2006, 128, 9882-9886. | 13.7 | 756 |
| 5 | Design rules for minimizing voltage losses in high-efficiency organic solar cells. <i>Nature Materials</i> , 2018, 17, 703-709. | 27.5 | 701 |
| 6 | Up-Conversion Intersystem Crossing Rates in Organic Emitters for Thermally Activated Delayed Fluorescence: Impact of the Nature of Singlet vs Triplet Excited States. <i>Journal of the American Chemical Society</i> , 2017, 139, 4042-4051. | 13.7 | 698 |
| 7 | Charge Transport Properties in Discotic Liquid Crystals: A Quantum-Chemical Insight into Structure-Property Relationships. <i>Journal of the American Chemical Society</i> , 2004, 126, 3271-3279. | 13.7 | 464 |
| 8 | Delocalization of exciton and electron wavefunction in non-fullerene acceptor molecules enables efficient organic solar cells. <i>Nature Communications</i> , 2020, 11, 3943. | 12.8 | 458 |
| 9 | The Vibrational Reorganization Energy in Pentacene: Molecular Influences on Charge Transport. <i>Journal of the American Chemical Society</i> , 2002, 124, 7918-7919. | 13.7 | 425 |
| 10 | The Impact of Molecular Orientation on the Photovoltaic Properties of a Phthalocyanine/Fullerene Heterojunction. <i>Advanced Functional Materials</i> , 2012, 22, 2987-2995. | 14.9 | 298 |
| 11 | Exciton-Dissociation and Charge-Recombination Processes in Pentacene/C ₆₀ Solar Cells: Theoretical Insight into the Impact of Interface Geometry. <i>Journal of the American Chemical Society</i> , 2009, 131, 15777-15783. | 13.7 | 275 |
| 12 | Intersystem Crossing Processes in Nonplanar Aromatic Heterocyclic Molecules. <i>Journal of Physical Chemistry A</i> , 2007, 111, 10490-10499. | 2.5 | 261 |
| 13 | Prediction of Remarkable Ambipolar Charge-Transport Characteristics in Organic Mixed-Stack Charge-Transfer Crystals. <i>Journal of the American Chemical Society</i> , 2012, 134, 2340-2347. | 13.7 | 245 |
| 14 | A unified description of non-radiative voltage losses in organic solar cells. <i>Nature Energy</i> , 2021, 6, 799-806. | 39.5 | 235 |
| 15 | Ultralow Doping in Organic Semiconductors: Evidence of Trap Filling. <i>Physical Review Letters</i> , 2012, 109, 176601. | 7.8 | 231 |
| 16 | Charge-transfer electronic states in organic solar cells. <i>Nature Reviews Materials</i> , 2019, 4, 689-707. | 48.7 | 229 |
| 17 | Impact of interfacial molecular orientation on radiative recombination and charge generation efficiency. <i>Nature Communications</i> , 2017, 8, 79. | 12.8 | 198 |
| 18 | High stability and luminescence efficiency in donor-acceptor neutral radicals not following the Aufbau principle. <i>Nature Materials</i> , 2019, 18, 977-984. | 27.5 | 181 |

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 19 | Impact of Perfluorination on the Charge-Transport Parameters of Oligoacene Crystals. <i>Journal of the American Chemical Society</i> , 2009, 131, 1502-1512. | 13.7 | 174 |
| 20 | Charge Transport Parameters of the Pentathienoacene Crystal. <i>Journal of the American Chemical Society</i> , 2007, 129, 13072-13081. | 13.7 | 153 |
| 21 | Bistetracene: An Air-Stable, High-Mobility Organic Semiconductor with Extended Conjugation. <i>Journal of the American Chemical Society</i> , 2014, 136, 9248-9251. | 13.7 | 150 |
| 22 | Vibronic Coupling in the Ground and Excited States of Oligoacene Cations. <i>Journal of Physical Chemistry B</i> , 2006, 110, 18904-18911. | 2.6 | 140 |
| 23 | Intervalence Transitions in the Mixed-Valence Monocations of Bis(triarylamines) Linked with Vinylene and Phenylene-Vinylene Bridges. <i>Journal of the American Chemical Society</i> , 2005, 127, 16900-16911. | 13.7 | 135 |
| 24 | Interaction of Charge Carriers with Lattice Vibrations in Oligoacene Crystals from Naphthalene to Pentacene. <i>Journal of the American Chemical Society</i> , 2010, 132, 14437-14446. | 13.7 | 128 |
| 25 | Assessing the nature of the charge-transfer electronic states in organic solar cells. <i>Nature Communications</i> , 2018, 9, 5295. | 12.8 | 126 |
| 26 | Delocalization in Platinum-Alkynyl Systems: A Metal-Bridged Organic Mixed-Valence Compound. <i>Journal of the American Chemical Society</i> , 2004, 126, 11782-11783. | 13.7 | 121 |
| 27 | Design of Efficient Ambipolar Host Materials for Organic Blue Electrophosphorescence: Theoretical Characterization of Hosts Based on Carbazole Derivatives. <i>Journal of the American Chemical Society</i> , 2011, 133, 17895-17900. | 13.7 | 116 |
| 28 | Asymmetric electron acceptor enables highly luminescent organic solar cells with certified efficiency over 18%. <i>Nature Communications</i> , 2022, 13, 2598. | 12.8 | 113 |
| 29 | A comparative theoretical study of exciton-dissociation and charge-recombination processes in oligothiophene/fullerene and oligothiophene/peryleneimide complexes for organic solar cells. <i>Journal of Materials Chemistry</i> , 2011, 21, 1479. | 6.7 | 112 |
| 30 | Electronic Coupling in Tetraanisylarylenediamine Mixed-Valence Systems: The Interplay between Bridge Energy and Geometric Factors. <i>Journal of the American Chemical Society</i> , 2005, 127, 8508-8516. | 13.7 | 107 |
| 31 | Interaction of Charge Carriers with Lattice Vibrations in Organic Molecular Semiconductors: Naphthalene as a Case Study. <i>Journal of Physical Chemistry C</i> , 2009, 113, 4679-4686. | 3.1 | 102 |
| 32 | Static and Dynamic Energetic Disorders in the C ₆₀ , PC ₆₁ BM, C ₇₀ , and PC ₇₁ BM Fullerenes. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 3657-3662. | 4.6 | 101 |
| 33 | Factors Governing Intercalation of Fullerenes and Other Small Molecules Between the Side Chains of Semiconducting Polymers Used in Solar Cells. <i>Advanced Energy Materials</i> , 2012, 2, 1208-1217. | 19.5 | 97 |
| 34 | Closely Stacked Oligo(phenylene ethynylene)s: Effect of π -Stacking on the Electronic Properties of Conjugated Chromophores. <i>Journal of the American Chemical Society</i> , 2012, 134, 7176-7185. | 13.7 | 96 |
| 35 | Phosphine Oxide Derivatives as Hosts for Blue Phosphors: A Joint Theoretical and Experimental Study of Their Electronic Structure. <i>Chemistry of Materials</i> , 2010, 22, 247-254. | 6.7 | 95 |
| 36 | Charge-Transfer States in Organic Solar Cells: Understanding the Impact of Polarization, Delocalization, and Disorder. <i>ACS Applied Materials & Interfaces</i> , 2017, 9, 18095-18102. | 8.0 | 90 |

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| 37 | Influence of Structural Dynamics on Polarization Energies in Anthracene Single Crystals. <i>Journal of Physical Chemistry C</i> , 2010, 114, 20678-20685. | 3.1 | 86 |
| 38 | Electronic Couplings in Organic Mixed-Valence Compounds: The Contribution of Photoelectron Spectroscopy. <i>Journal of the American Chemical Society</i> , 2004, 126, 2727-2731. | 13.7 | 85 |
| 39 | Effect of Solid-State Polarization on Charge-Transfer Excitations and Transport Levels at Organic Interfaces from a Screened Range-Separated Hybrid Functional. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 3277-3283. | 4.6 | 84 |
| 40 | A New Class of Mixed-Valence Systems with Orbitally Degenerate Organic Redox Centers. Examples Based on Hexa-Rhenium Molecular Prisms. <i>Journal of the American Chemical Society</i> , 2006, 128, 12592-12593. | 13.7 | 83 |
| 41 | Polymorphism in the 1:1 Charge-Transfer Complex DBTTF-TCNQ and Its Effects on Optical and Electronic Properties. <i>Advanced Electronic Materials</i> , 2016, 2, 1600203. | 5.1 | 83 |
| 42 | Impact of Electron Delocalization on the Nature of the Charge-Transfer States in Model Pentacene/C ₆₀ Interfaces: A Density Functional Theory Study. <i>Journal of Physical Chemistry C</i> , 2014, 118, 27648-27656. | 3.1 | 80 |
| 43 | Understanding the Density Functional Dependence of DFT-Calculated Electronic Couplings in Organic Semiconductors. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 919-924. | 4.6 | 79 |
| 44 | Electronic Properties of Mixed-Stack Organic Charge-Transfer Crystals. <i>Journal of Physical Chemistry C</i> , 2014, 118, 14150-14156. | 3.1 | 79 |
| 45 | Isolation and Crystal Structures of Two Singlet Bis(Triarylamine) Dications with Nonquinoidal Geometries. <i>Journal of the American Chemical Society</i> , 2006, 128, 1812-1817. | 13.7 | 78 |
| 46 | To bend or not to bend – are heteroatom interactions within conjugated molecules effective in dictating conformation and planarity?. <i>Materials Horizons</i> , 2016, 3, 333-339. | 12.2 | 78 |
| 47 | Theoretical Study of the Local and Charge-Transfer Excitations in Model Complexes of Pentacene-C ₆₀ Using Tuned Range-Separated Hybrid Functionals. <i>Journal of Chemical Theory and Computation</i> , 2014, 10, 2379-2388. | 5.3 | 77 |
| 48 | Intramolecular Electron-Transfer Rates in Mixed-Valence Triarylamines: Measurement by Variable-Temperature ESR Spectroscopy and Comparison with Optical Data. <i>Journal of the American Chemical Society</i> , 2009, 131, 1717-1723. | 13.7 | 75 |
| 49 | Vibronic Coupling in Organic Semiconductors: The Case of Fused Polycyclic Benzene-Thiophene Structures. <i>Chemistry - A European Journal</i> , 2006, 12, 2073-2080. | 3.3 | 74 |
| 50 | Bis[bis(4-alkoxyphenyl)amino] Derivatives of Dithienylethene, Bithiophene, Dithienothiophene and Dithienopyrrole: Palladium-Catalysed Synthesis and Highly Delocalised Radical Cations. <i>Chemistry - A European Journal</i> , 2007, 13, 9637-9646. | 3.3 | 72 |
| 51 | Tuning Delocalization in the Radical Cations of 1,4-Bis[4-(diarylamino)styryl]benzenes, 2,5-Bis[4-(diarylamino)styryl]thiophenes, and 2,5-Bis[4-(diarylamino)styryl]pyrroles through Substituent Effects. <i>Journal of the American Chemical Society</i> , 2012, 134, 10146-10155. | 13.7 | 72 |
| 52 | Temperature-Mediated Polymorphism in Molecular Crystals: The Impact on Crystal Packing and Charge Transport. <i>Chemistry of Materials</i> , 2015, 27, 112-118. | 6.7 | 72 |
| 53 | Mode-selective vibrational modulation of charge transport in organic electronic devices. <i>Nature Communications</i> , 2015, 6, 7880. | 12.8 | 72 |
| 54 | Quantum-Chemical Approach to Electronic Coupling: Application to Charge Separation and Charge Recombination Pathways in a Model Molecular Donor-Acceptor System for Organic Solar Cells. <i>Journal of Physical Chemistry C</i> , 2008, 112, 3429-3433. | 3.1 | 69 |

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| 55 | Polymers with Carbazole-Oxadiazole Side Chains as Ambipolar Hosts for Phosphorescent Light-Emitting Diodes. <i>Chemistry of Materials</i> , 2011, 23, 4002-4015. | 6.7 | 67 |
| 56 | Description of the Charge Transfer States at the Pentacene/C ₆₀ Interface: Combining Range-Separated Hybrid Functionals with the Polarizable Continuum Model. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2616-2621. | 4.6 | 66 |
| 57 | Vibration-Assisted Crystallization Improves Organic/Dielectric Interface in Organic Thin-Film Transistors. <i>Advanced Materials</i> , 2013, 25, 6956-6962. | 21.0 | 65 |
| 58 | Probing Charge Transport in π -Stacked Fluorene-Based Systems. <i>Journal of Physical Chemistry B</i> , 2006, 110, 9482-9487. | 2.6 | 64 |
| 59 | Electronic and Optical Properties of 4-H-Cyclopenta[2,1-b:3,4-b ²]bithiophene Derivatives and Their 4-Heteroatom-Substituted Analogues: A Joint Theoretical and Experimental Comparison. <i>Journal of Physical Chemistry B</i> , 2010, 114, 14397-14407. | 2.6 | 64 |
| 60 | Electronic-Structure Theory of Organic Semiconductors: Charge-Transport Parameters and Metal/Organic Interfaces. <i>Annual Review of Materials Research</i> , 2013, 43, 63-87. | 9.3 | 62 |
| 61 | Design and synthesis of two-dimensional covalent organic frameworks with four-arm cores: prediction of remarkable ambipolar charge-transport properties. <i>Materials Horizons</i> , 2019, 6, 1868-1876. | 12.2 | 62 |
| 62 | Role of band states and trap states in the electrical properties of organic semiconductors: Hopping versus mobility edge model. <i>Physical Review B</i> , 2013, 87, . | 3.2 | 57 |
| 63 | Thermal Narrowing of the Electronic Bandwidths in Organic Molecular Semiconductors: Impact of the Crystal Thermal Expansion. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 3325-3329. | 4.6 | 56 |
| 64 | Charge-Transfer States at Organic/Organic Interfaces: Impact of Static and Dynamic Disorders. <i>Advanced Energy Materials</i> , 2019, 9, 1803926. | 19.5 | 54 |
| 65 | Theoretical Investigation of Triscarbazole Derivatives As Host Materials for Blue Electrophosphorescence: Effects of Topology. <i>Chemistry of Materials</i> , 2011, 23, 5223-5230. | 6.7 | 53 |
| 66 | Intrinsic charge transport in single crystals of organic molecular semiconductors: A theoretical perspective. <i>MRS Bulletin</i> , 2013, 38, 57-64. | 3.5 | 53 |
| 67 | Experimental Reorganization Energies of Pentacene and Perfluoropentacene: Effects of Perfluorination. <i>Journal of Physical Chemistry C</i> , 2013, 117, 22428-22437. | 3.1 | 53 |
| 68 | Theoretical study of substitution effects on molecular reorganization energy in organic semiconductors. <i>Journal of Chemical Physics</i> , 2011, 135, 104703. | 3.0 | 52 |
| 69 | On the Molecular Origin of Charge Separation at the Donor-Acceptor Interface. <i>Advanced Energy Materials</i> , 2018, 8, 1702232. | 19.5 | 51 |
| 70 | Symmetry effects on nonlocal electron-phonon coupling in organic semiconductors. <i>Physical Review B</i> , 2012, 85, . | 3.2 | 48 |
| 71 | Nonlocal electron-phonon coupling in the pentacene crystal: Beyond the Γ -point approximation. <i>Journal of Chemical Physics</i> , 2012, 137, 164303. | 3.0 | 48 |
| 72 | Triisopropylsilylethynyl-Functionalized Graphene-Like Fragment Semiconductors: Synthesis, Crystal Packing, and Density Functional Theory Calculations. <i>Chemistry - A European Journal</i> , 2013, 19, 17907-17916. | 3.3 | 48 |

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| 73 | Defect-Driven Interfacial Electronic Structures at an Organic/Metal-Oxide Semiconductor Heterojunction. <i>Advanced Materials</i> , 2014, 26, 4711-4716. | 21.0 | 46 |
| 74 | Organic Neutral Radical Emitters: Impact of Chemical Substitution and Electronic-State Hybridization on the Luminescence Properties. <i>Journal of the American Chemical Society</i> , 2020, 142, 17782-17786. | 13.7 | 46 |
| 75 | π-Stacked Oligo(phenylene vinylene)s Based on Pseudo-Geminal Substituted [2.2]Paracyclophanes: Impact of Interchain Geometry and Interactions on the Electronic Properties. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 11629-11632. | 13.8 | 44 |
| 76 | Dipolar Second-Order Nonlinear Optical Chromophores Containing Ferrocene, Octamethylferrocene, and Ruthenocene Donors and Strong π-Acceptors: Crystal Structures and Comparison of π-Donor Strengths. <i>Organometallics</i> , 2009, 28, 1350-1357. | 2.3 | 43 |
| 77 | Borderline Class II/III Ligand-Centered Mixed Valency in a Porphyrinic Molecular Rectangle. <i>Inorganic Chemistry</i> , 2005, 44, 5789-5797. | 4.0 | 42 |
| 78 | Donor Conjugated Polymers with Polar Side Chain Groups: The Role of Dielectric Constant and Energetic Disorder on Photovoltaic Performance. <i>Advanced Functional Materials</i> , 2018, 28, 1803418. | 14.9 | 42 |
| 79 | Photophysical Properties of an Alkyne-Bridged Bis(zinc porphyrin)-Perylene Bis(dicarboximide) Derivative. <i>Journal of Physical Chemistry A</i> , 2009, 113, 10826-10832. | 2.5 | 41 |
| 80 | Electronic Structure and Charge-Transport Parameters of Functionalized Tetracene Crystals: Impact of Partial Fluorination and Alkyl or Alkoxy Derivatization. <i>Chemistry of Materials</i> , 2009, 21, 3593-3601. | 6.7 | 41 |
| 81 | Hexaazatriphenylene (HAT) versus tri-HAT: The Bigger the Better?. <i>Chemistry - A European Journal</i> , 2011, 17, 10312-10322. | 3.3 | 40 |
| 82 | Every Atom Counts: Elucidating the Fundamental Impact of Structural Change in Conjugated Polymers for Organic Photovoltaics. <i>Chemistry of Materials</i> , 2018, 30, 2995-3009. | 6.7 | 39 |
| 83 | Hyperfluorescence-Based Emission in Purely Organic Materials: Suppression of Energy-Loss Mechanisms via Alignment of Triplet Excited States. , 2020, 2, 1412-1418. | | 39 |
| 84 | Dipolar Ferrocene and Ruthenocene Second-Order Nonlinear Optical Chromophores: A Time-Dependent Density Functional Theory Investigation of Their Absorption Spectra. <i>Organometallics</i> , 2013, 32, 6061-6068. | 2.3 | 38 |
| 85 | Charge-Transport Properties of F ₆ TNAP-Based Charge-Transfer Cocrystals. <i>Advanced Functional Materials</i> , 2019, 29, 1904858. | 14.9 | 36 |
| 86 | Resolving Atomic-Scale Interactions in Nonfullerene Acceptor Organic Solar Cells with Solid-State NMR Spectroscopy, Crystallographic Modelling, and Molecular Dynamics Simulations. <i>Advanced Materials</i> , 2022, 34, e2105943. | 21.0 | 36 |
| 87 | Nonlocal electron-phonon coupling in organic semiconductor crystals: The role of acoustic lattice vibrations. <i>Journal of Chemical Physics</i> , 2013, 138, 204713. | 3.0 | 34 |
| 88 | Purely Organic Emitters for Multiresonant Thermally Activated Delay Fluorescence: Design of Highly Efficient Sulfur and Selenium Derivatives. , 2022, 4, 440-447. | | 33 |
| 89 | Packing and Disorder in Substituted Fullerenes. <i>Journal of Physical Chemistry C</i> , 2016, 120, 17242-17250. | 3.1 | 28 |
| 90 | A polarized response. <i>Nature Materials</i> , 2006, 5, 929-930. | 27.5 | 27 |

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| 91 | Impact of Active Layer Morphology on Bimolecular Recombination Dynamics in Organic Solar Cells. <i>Journal of Physical Chemistry C</i> , 2017, 121, 24954-24961. | 3.1 | 26 |
| 92 | All-Polymer Solar Cells: Impact of the Length of the Branched Alkyl Side Chains on the Polymer Acceptors on the Interchain Packing and Electronic Properties in Amorphous Blends. <i>Chemistry of Materials</i> , 2019, 31, 6239-6248. | 6.7 | 26 |
| 93 | Electronic structure of the pentacene-gold interface: A density-functional theory study. <i>Organic Electronics</i> , 2009, 10, 1571-1578. | 2.6 | 25 |
| 94 | Electronic Properties of 1,5-Diaminonaphthalene:Tetrahalo-1,4-benzoquinone Donor-Acceptor Cocrystals. <i>Journal of Physical Chemistry C</i> , 2017, 121, 23633-23641. | 3.1 | 25 |
| 95 | Discovery of Non-linear Optical Materials by Function-Based Screening of Multi-component Solids. <i>CheM</i> , 2018, 4, 150-161. | 11.7 | 25 |
| 96 | Thermally Activated Delayed Fluorescence Sensitization for Highly Efficient Blue Fluorescent Emitters. <i>Advanced Functional Materials</i> , 2020, 30, 2005898. | 14.9 | 25 |
| 97 | Impact of exact exchange in the description of the electronic structure of organic charge-transfer molecular crystals. <i>Physical Review B</i> , 2014, 90, . | 3.2 | 24 |
| 98 | Toward a Robust Quantum-Chemical Description of Organic Mixed-Valence Systems. <i>Journal of Physical Chemistry C</i> , 2014, 118, 3925-3934. | 3.1 | 23 |
| 99 | Radiative and Nonradiative Recombinations in Organic Radical Emitters: The Effect of Guest-Host Interactions. <i>Advanced Functional Materials</i> , 2020, 30, 2002916. | 14.9 | 23 |
| 100 | The Role of Intermolecular Interactions on the Performance of Organic Thermally Activated Delayed Fluorescence (TADF) Materials. <i>Advanced Optical Materials</i> , 2021, 9, 2002135. | 7.3 | 22 |
| 101 | Energetic fluctuations in amorphous semiconducting polymers: Impact on charge-carrier mobility. <i>Journal of Chemical Physics</i> , 2017, 147, 134904. | 3.0 | 21 |
| 102 | Characterizing the Polymer:Fullerene Intermolecular Interactions. <i>Chemistry of Materials</i> , 2016, 28, 1446-1452. | 6.7 | 20 |
| 103 | Electronic Structure of Multicomponent Organic Molecular Materials: Evaluation of Range-Separated Hybrid Functionals. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 3712-3719. | 5.3 | 20 |
| 104 | Impact of Phonon Dispersion on Nonlocal Electron-Phonon Couplings in Organic Semiconductors: The Naphthalene Crystal as a Case Study. <i>Journal of Physical Chemistry C</i> , 2018, 122, 44-49. | 3.1 | 18 |
| 105 | Charge-Transport Properties of the 1,4-Diiodobenzene Crystal: A Quantum-Mechanical Study. <i>Chemistry of Materials</i> , 2008, 20, 5832-5838. | 6.7 | 17 |
| 106 | Langmuir-Blodgett Thin Films of Diketopyrrolopyrrole-Based Amphiphiles. <i>ACS Applied Materials & Interfaces</i> , 2018, 10, 11995-12004. | 8.0 | 17 |
| 107 | Quaternary Charge-Transfer Solid Solutions: Electronic Tunability through Stoichiometry. <i>Chemistry of Materials</i> , 2019, 31, 6598-6604. | 6.7 | 17 |
| 108 | Suppression of Concentration Quenching in Ortho-Substituted Thermally Activated Delayed Fluorescence Emitters. <i>Advanced Theory and Simulations</i> , 2020, 3, 1900185. | 2.8 | 17 |

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|-----|---|------|-----------|
| 109 | Understanding the effects of electronic polarization and delocalization on charge-transport levels in oligoacene systems. <i>Journal of Chemical Physics</i> , 2017, 146, 224705. | 3.0 | 16 |
| 110 | Optical conductivity and optical effective mass in a high-mobility organic semiconductor: Implications for the nature of charge transport. <i>Physical Review B</i> , 2014, 90, . | 3.2 | 15 |
| 111 | Impact of secondary donor units on the excited-state properties and thermally activated delayed fluorescence (TADF) efficiency of pentacarbazole-benzonitrile emitters. <i>Journal of Chemical Physics</i> , 2020, 153, 144708. | 3.0 | 14 |
| 112 | Impact of chemical modifications on the luminescence properties of organic neutral radical emitters. <i>Journal of Materials Chemistry C</i> , 2021, 9, 10794-10801. | 5.5 | 13 |
| 113 | Strong Suppression of Thermal Conductivity in the Presence of Long Terminal Alkyl Chains in Low-Disorder Molecular Semiconductors. <i>Advanced Materials</i> , 2021, 33, e2008708. | 21.0 | 12 |
| 114 | Lower limits for non-radiative recombination loss in organic donor/acceptor complexes. <i>Materials Horizons</i> , 2022, 9, 325-333. | 12.2 | 12 |
| 115 | Electronic, vibrational, and charge-transport properties of benzo[h]thienobenzothioophene-TCNQ co-crystals. <i>Materials Chemistry Frontiers</i> , 2020, 4, 3623-3631. | 5.9 | 11 |
| 116 | On the Physical Origins of Charge Separation at Donor-Acceptor Interfaces in Organic Solar Cells: Energy Bending versus Energy Disorder. <i>Advanced Theory and Simulations</i> , 2020, 3, 1900230. | 2.8 | 11 |
| 117 | Correlating Non-Geminate Recombination with Film Structure: A Comparison of Polythiophene: Fullerene Bilayer and Blend Films. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 3669-3676. | 4.6 | 9 |
| 118 | Electron-phonon coupling in anthracene-pyromellitic dianhydride. <i>Journal of Chemical Physics</i> , 2017, 146, 214705. | 3.0 | 9 |
| 119 | Organic thin films with charge-carrier mobility exceeding that of single crystals. <i>Journal of Materials Chemistry C</i> , 2017, 5, 10313-10319. | 5.5 | 9 |
| 120 | Characterization of the structural, mechanical, and electronic properties of fullerene mixtures: a molecular simulations description. <i>Journal of Materials Chemistry C</i> , 2018, 6, 3642-3650. | 5.5 | 8 |
| 121 | Electronic structure of confined carbyne from joint wavelength-dependent resonant Raman spectroscopy and density functional theory investigations. <i>Carbon</i> , 2022, 189, 276-283. | 10.3 | 8 |
| 122 | Theoretical description of the geometric and electronic structures of organic-organic interfaces in organic solar cells: a brief review. <i>Science China Chemistry</i> , 2014, 57, 1330-1339. | 8.2 | 6 |
| 123 | Charge Transport in Crystalline Organic Semiconductors. <i>Materials and Energy</i> , 2016, , 193-230. | 0.1 | 6 |
| 124 | Energy transfer processes in hyperfluorescent organic light-emitting diodes. <i>Journal of Materials Chemistry C</i> , 2022, 10, 4629-4636. | 5.5 | 6 |
| 125 | Bulk Heterojunction Solar Cells: Insight into Ternary Blends from a Characterization of the Intermolecular Packing and Electronic Properties in the Corresponding Binary Blends. <i>Advanced Theory and Simulations</i> , 2020, 3, 2000049. | 2.8 | 3 |
| 126 | Reply to Comment on Polymorphism in the 1:1 Charge-Transfer Complex DBTTF-TCNQ and Its Effects on Optical and Electronic Properties. <i>Advanced Electronic Materials</i> , 2017, 3, 1600521. | 5.1 | 2 |

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| 127 | Organic Thin-Film Transistors: Vibration-Assisted Crystallization Improves Organic/Dielectric Interface in Organic Thin-Film Transistors (Adv. Mater. 48/2013). Advanced Materials, 2013, 25, 7054-7054. | 21.0 | 0 |
| 128 | Charge-Transfer States at Organic-Organic Interfaces: Impact on Charge Recombination Processes. , 0, , . | | 0 |
| 129 | Organic charge-transfer compounds: complex interactions at the nanoscale. , 2019, , . | | 0 |
| 130 | Resolving atomic-scale interactions in non-fullerene acceptor organic solar cells by high-field NMR crystallography. , 0, , . | | 0 |