

Troy Van Voorhis

List of Publications by Year in descending order

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122
papers

15,585
citations

38742

50
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18130

120
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125
all docs

125
docs citations

125
times ranked

15366
citing authors

#	ARTICLE	IF	CITATIONS
1	Heterogeneous Pair Approximation of Methanol Oxidation on TiO ₂ Reveals Two Reaction Pathways. <i>Journal of Physical Chemistry C</i> , 2022, 126, 1845-1856.	3.1	4
2	Diabatic Valence-Hole States in the C ₂ Molecule: "Putting Humpty Dumpty Together Again". <i>Journal of Physical Chemistry A</i> , 2022, 126, 3090-3100.	2.5	8
3	Understanding the Dipole Moment of Liquid Water from a Self-Attractive Hartree Decomposition. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 6-12.	4.6	14
4	Colloidal CdSe nanocrystals are inherently defective. <i>Nature Communications</i> , 2021, 12, 890.	12.8	22
5	Accurate Electronic Excitation Energies in Full-Valence Active Space via Bootstrap Embedding. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 3335-3347.	5.3	8
6	Software for the frontiers of quantum chemistry: An overview of developments in the Q-Chem 5 package. <i>Journal of Chemical Physics</i> , 2021, 155, 084801.	3.0	518
7	Maximizing TADF via Conformational Optimization. <i>Journal of Physical Chemistry A</i> , 2021, 125, 7644-7654.	2.5	11
8	Machine learning dynamic correlation in chemical kinetics. <i>Journal of Chemical Physics</i> , 2021, 155, 144107.	3.0	2
9	Resolving the Triexciton Recombination Pathway in CdSe/CdS Nanocrystals through State-Specific Correlation Measurements. <i>Nano Letters</i> , 2021, 21, 7457-7464.	9.1	13
10	Photoswitchable Sol-Gel Transitions and Catalysis Mediated by Polymer Networks with Coumarin-Decorated Cu ₂₄ L ₂₄ Metal-Organic Cages as Junctions. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 2784-2792.	13.8	58
11	Photoswitchable Sol-Gel Transitions and Catalysis Mediated by Polymer Networks with Coumarin-Decorated Cu ₂₄ L ₂₄ Metal-Organic Cages as Junctions. <i>Angewandte Chemie</i> , 2020, 132, 2806-2814.	2.0	12
12	An N-Heterocyclic Carbene-Derived Distonic Radical Cation. <i>Angewandte Chemie</i> , 2020, 132, 3980-3983.	2.0	4
13	An N-Heterocyclic Carbene-Derived Distonic Radical Cation. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 3952-3955.	13.8	16
14	A pyridinic Fe-N ₄ macrocycle models the active sites in Fe/N-doped carbon electrocatalysts. <i>Nature Communications</i> , 2020, 11, 5283.	12.8	286
15	Bootstrap embedding with an unrestricted mean-field bath. <i>Journal of Chemical Physics</i> , 2020, 153, 214101.	3.0	5
16	Understanding Disorder in 2D Materials: The Case of Carbon Doping of Silicene. <i>Nano Letters</i> , 2020, 20, 6336-6343.	9.1	8
17	Electrochemically mediated carbon dioxide separation with quinone chemistry in salt-concentrated aqueous media. <i>Nature Communications</i> , 2020, 11, 2278.	12.8	71
18	Investigation of External Quantum Efficiency Roll-Off in OLEDs Using the Mean-Field Steady-State Kinetic Model. <i>Journal of Physical Chemistry C</i> , 2020, 124, 14424-14431.	3.1	2

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19	Gas-Phase Ethylene Polymerization by Single-Site Cr Centers in a Metal-Organic Framework. <i>ACS Catalysis</i> , 2020, 10, 3864-3870.	11.2	17
20	Bootstrap Embedding For Large Molecular Systems. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 5035-5046.	5.3	11
21	Lock-and-Key Exciplexes for Thermally Activated Delayed Fluorescence. <i>Organic Materials</i> , 2020, 02, 001-010.	2.0	7
22	Toward Prediction of Nonradiative Decay Pathways in Organic Compounds II: Two Internal Conversion Channels in BODIPYs. <i>Journal of Physical Chemistry C</i> , 2020, 124, 3925-3938.	3.1	51
23	Innenteilbild: Photoswitchable Sol-Gel Transitions and Catalysis Mediated by Polymer Networks with Coumarin-Decorated Cu ₂ L ₂ Metal-Organic Cages as Junctions (<i>Angew. Tj E</i>) 2021 1 0.784314	10.1	14
24	Dominance of Exciton Lifetime in the Stability of Phosphorescent Dyes. <i>Advanced Optical Materials</i> , 2019, 7, 1901048.	7.3	23
25	Large Increase in External Quantum Efficiency by Dihedral Angle Tuning in a Sky-Blue Thermally Activated Delayed Fluorescence Emitter. <i>Advanced Optical Materials</i> , 2019, 7, 1900476.	7.3	25
26	Using SCF metadynamics to extend density matrix embedding theory to excited states. <i>Journal of Chemical Physics</i> , 2019, 151, 034112.	3.0	15
27	Bootstrap Embedding for Molecules. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 4497-4506.	5.3	22
28	Atom-Based Bootstrap Embedding For Molecules. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 6368-6374.	4.6	14
29	Implementation of the Many-Pair Expansion for Systematically Improving Density Functional Calculations of Molecules. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 1089-1101.	5.3	5
30	Toward Prediction of Nonradiative Decay Pathways in Organic Compounds I: The Case of Naphthalene Quantum Yields. <i>Journal of Physical Chemistry C</i> , 2019, 123, 15394-15402.	3.1	41
31	Extracting Design Principles for Efficient Thermally Activated Delayed Fluorescence (TADF) from a Simple Four-State Model. <i>Chemistry of Materials</i> , 2019, 31, 6995-7006.	6.7	84
32	A Heterogeneous Kinetics Model for Triplet Exciton Transfer in Solid-State Upconversion. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 3147-3152.	4.6	24
33	Half-Projected \ddot{f} Self-Consistent Field For Electronic Excited States. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 2954-2965.	5.3	25
34	Polymers with Side Chain Porosity for Ultraporous and Plasticization Resistant Materials for Gas Separations. <i>Advanced Materials</i> , 2019, 31, e1807871.	21.0	64
35	Unraveling the Fate of Host Excitons in Host-Guest Phosphorescent Organic Light-Emitting Diodes. <i>Journal of Physical Chemistry C</i> , 2019, 123, 10311-10318.	3.1	10
36	Robust gold nanorods stabilized by bidentate N-heterocyclic-carbene-thiolate ligands. <i>Nature Chemistry</i> , 2019, 11, 57-63.	13.6	109

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37	Triplet Tuning: A Novel Family of Non-Empirical Exchange-Correlation Functionals. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 1226-1241.	5.3	34
38	QM/MM Study of Static and Dynamic Energetic Disorder in the Emission Layer of an Organic Light-Emitting Diode. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1329-1334.	4.6	23
39	Non-radiative deactivation of cytosine derivatives at elevated temperature. <i>Molecular Physics</i> , 2018, 116, 2591-2598.	1.7	2
40	Molecular Design of Deep Blue Thermally Activated Delayed Fluorescence Materials Employing a Homoconjugative Triptycene Scaffold and Dihedral Angle Tuning. <i>Chemistry of Materials</i> , 2018, 30, 1462-1466.	6.7	71
41	Self-Attractive Hartree Decomposition: Partitioning Electron Density into Smooth Localized Fragments. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 92-103.	5.3	7
42	Polydiacetylene functionalized with charged termini for device-free colorimetric detection of malathion. <i>Journal of Colloid and Interface Science</i> , 2018, 528, 27-35.	9.4	16
43	Band-Like Charge Photogeneration at a Crystalline Organic Donor/Acceptor Interface. <i>Advanced Energy Materials</i> , 2018, 8, 1701494.	19.5	23
44	Incremental embedding: A density matrix embedding scheme for molecules. <i>Journal of Chemical Physics</i> , 2018, 149, 194108.	3.0	17
45	Morphology of Passivating Organic Ligands around a Nanocrystal. <i>Journal of Physical Chemistry C</i> , 2018, 122, 26267-26274.	3.1	34
46	Polyaniline Nanofiber Electrodes for Reversible Capture and Release of Mercury(II) from Water. <i>Journal of the American Chemical Society</i> , 2018, 140, 14413-14420.	13.7	87
47	Surface States Mediate Triplet Energy Transfer in Nanocrystal-Acene Composite Systems. <i>Journal of the American Chemical Society</i> , 2018, 140, 7543-7553.	13.7	88
48	Insights into Magneto-Optics of Helical Conjugated Polymers. <i>Journal of the American Chemical Society</i> , 2018, 140, 6501-6508.	13.7	76
49	Superhydrophobic, Surfactant-Doped, Conducting Polymers for Electrochemically Reversible Adsorption of Organic Contaminants. <i>Advanced Functional Materials</i> , 2018, 28, 1801466.	14.9	33
50	Charge Transfer in Molecular Materials. , 2018, , 1-31.		4
51	Exploring Low Internal Reorganization Energies for Silicene Nanoclusters. <i>Physical Review Applied</i> , 2018, 9, .	3.8	7
52	A Semiconducting Conjugated Radical Polymer: Ambipolar Redox Activity and Faraday Effect. <i>Journal of the American Chemical Society</i> , 2018, 140, 10881-10889.	13.7	41
53	Long-range interactions from the many-pair expansion: A different avenue to dispersion in DFT. <i>Journal of Chemical Physics</i> , 2017, 146, 024111.	3.0	6
54	Molecular dynamics and charge transport in organic semiconductors: a classical approach to modeling electron transfer. <i>Chemical Science</i> , 2017, 8, 2597-2609.	7.4	13

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55	Performance of Bootstrap Embedding for long-range interactions and 2D systems. <i>Molecular Physics</i> , 2017, 115, 2242-2253.	1.7	21
56	A structural and mechanistic study of I^- -clamp-mediated cysteine perfluoroarylation. <i>Scientific Reports</i> , 2017, 7, 7954.	3.3	20
57	Molecular-Level Insights into Oxygen Reduction Catalysis by Graphite-Conjugated Active Sites. <i>ACS Catalysis</i> , 2017, 7, 7680-7687.	11.2	33
58	Shorter Exciton Lifetimes via an External Heavy-Atom Effect: Alleviating the Effects of Bimolecular Processes in Organic Light-Emitting Diodes. <i>Advanced Materials</i> , 2017, 29, 1701987.	21.0	90
59	Mean field treatment of heterogeneous steady state kinetics. <i>Chemical Physics Letters</i> , 2017, 685, 185-190.	2.6	3
60	Speed Limit for Triplet-Exciton Transfer in Solid-State PbS Nanocrystal-Sensitized Photon Upconversion. <i>ACS Nano</i> , 2017, 11, 7848-7857.	14.6	130
61	The Impact of Carrier Delocalization and Interfacial Electric Field Fluctuations on Organic Photovoltaics. <i>Journal of Physical Chemistry C</i> , 2017, 121, 26629-26636.	3.1	7
62	$\langle i \rangle \text{f} \langle /i \rangle$ -SCF: A direct energy-targeting method to mean-field excited states. <i>Journal of Chemical Physics</i> , 2017, 147, 214104.	3.0	50
63	A hybrid memory kernel approach for condensed phase non-adiabatic dynamics. <i>Journal of Chemical Physics</i> , 2017, 147, 014108.	3.0	5
64	Quantum chemical approaches to [NiFe] hydrogenase. <i>Essays in Biochemistry</i> , 2017, 61, 293-303.	4.7	5
65	Bootstrap embedding: An internally consistent fragment-based method. <i>Journal of Chemical Physics</i> , 2016, 145, 074102.	3.0	51
66	Condensed phase electron transfer beyond the Condon approximation. <i>Journal of Chemical Physics</i> , 2016, 145, 214105.	3.0	12
67	Charge Recombination in Phosphorescent Organic Light-Emitting Diode Host-Guest Systems through QM/MM Simulations. <i>Journal of Physical Chemistry C</i> , 2016, 120, 19987-19994.	3.1	19
68	Salt Effect Accelerates Site-Selective Cysteine Bioconjugation. <i>ACS Central Science</i> , 2016, 2, 637-646.	11.3	36
69	Adiabatic Approximation in Explicit Solvent Models of RedOx Chemistry. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 5111-5116.	5.3	10
70	Many-electron expansion: A density functional hierarchy for strongly correlated systems. <i>Physical Review B</i> , 2016, 93, .	3.2	13
71	First-Principles Monte Carlo Simulations of Reaction Equilibria in Compressed Vapors. <i>ACS Central Science</i> , 2016, 2, 409-415.	11.3	16
72	Prediction of Excited-State Energies and Singlet-Triplet Gaps of Charge-Transfer States Using a Restricted Open-Shell Kohn-Sham Approach. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 3353-3359.	5.3	74

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73	Ī-Clamp-mediated cysteine conjugation. <i>Nature Chemistry</i> , 2016, 8, 120-128.	13.6	236
74	Solid-state infrared-to-visible upconversion sensitized by colloidal nanocrystals. <i>Nature Photonics</i> , 2016, 10, 31-34.	31.4	418
75	Communication: CDFT-CI couplings can be unreliable when there is fractional charge transfer. <i>Journal of Chemical Physics</i> , 2015, 143, 231102.	3.0	26
76	Density matrix embedding in an antisymmetrized geminal power bath. <i>Journal of Chemical Physics</i> , 2015, 143, 024107.	3.0	46
77	Time-dependent projected Hartree-Fock. <i>Journal of Chemical Physics</i> , 2015, 142, 124103.	3.0	20
78	Spin-dependent charge transfer state design rules in organic photovoltaics. <i>Nature Communications</i> , 2015, 6, 6415.	12.8	83
79	Why many semiempirical molecular orbital theories fail for liquid water and how to fix them. <i>Journal of Computational Chemistry</i> , 2015, 36, 934-939.	3.3	16
80	The Role of Electron-Hole Separation in Thermally Activated Delayed Fluorescence in Donor-Acceptor Blends. <i>Journal of Physical Chemistry C</i> , 2015, 119, 25591-25597.	3.1	45
81	Thermally Activated Delayed Fluorescence Materials Based on Homoconjugation Effect of Donor-Acceptor Triptycenes. <i>Journal of the American Chemical Society</i> , 2015, 137, 11908-11911.	13.7	331
82	Advances in molecular quantum chemistry contained in the Q-Chem 4 program package. <i>Molecular Physics</i> , 2015, 113, 184-215.	1.7	2,561
83	Extended MÅller-Plesset perturbation theory for dynamical and static correlations. <i>Journal of Chemical Physics</i> , 2014, 141, 164117.	3.0	31
84	Resummed memory kernels in generalized system-bath master equations. <i>Journal of Chemical Physics</i> , 2014, 141, 054112.	3.0	24
85	Singlet fission efficiency in tetracene-based organic solar cells. <i>Applied Physics Letters</i> , 2014, 104, .	3.3	79
86	A transferable model for singlet-fission kinetics. <i>Nature Chemistry</i> , 2014, 6, 492-497.	13.6	402
87	Energy harvesting of non-emissive triplet excitons in tetracene by emissive PbS nanocrystals. <i>Nature Materials</i> , 2014, 13, 1039-1043.	27.5	235
88	A pathway to diphosphorus from the dissociation of photoexcited tetraphosphorus. <i>RSC Advances</i> , 2013, 3, 23166.	3.6	14
89	Charge Transfer or J-Coupling? Assignment of an Unexpected Red-Shifted Absorption Band in a Naphthalenediimide-Based Metal-Organic Framework. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 453-458.	4.6	65
90	Systematic Parametrization of Polarizable Force Fields from Quantum Chemistry Data. <i>Journal of Chemical Theory and Computation</i> , 2013, 9, 452-460.	5.3	164

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91	A multireference perturbation method using non-orthogonal Hartree-Fock determinants for ground and excited states. <i>Journal of Chemical Physics</i> , 2013, 139, 174104.	3.0	29
92	Excitation energies and Stokes shifts from a restricted open-shell Kohn-Sham approach. <i>Journal of Chemical Physics</i> , 2013, 138, 164101.	3.0	78
93	Densities of states for disordered systems from free probability. <i>Physical Review B</i> , 2013, 88, .	3.2	4
94	Constrained Density Functional Theory. <i>Chemical Reviews</i> , 2012, 112, 321-370.	47.7	454
95	Benchmark Assessment of the Accuracy of Several van der Waals Density Functionals. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 1929-1934.	5.3	93
96	A Polarizable QM/MM Explicit Solvent Model for Computational Electrochemistry in Water. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 610-617.	5.3	71
97	Assessment of the \hat{T}^2 SCF density functional theory approach for electronic excitations in organic dyes. <i>Journal of Chemical Physics</i> , 2011, 134, 054128.	3.0	152
98	Molecular Insight Into the Energy Levels at the Organic Donor/Acceptor Interface: A Quantum Mechanics/Molecular Mechanics Study. <i>Journal of Physical Chemistry C</i> , 2011, 115, 14431-14436.	3.1	83
99	Direct-Coupling O ₂ Bond Forming a Pathway in Cobalt Oxide Water Oxidation Catalysts. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2200-2204.	4.6	177
100	Exciton/Charge-Transfer Electronic Couplings in Organic Semiconductors. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 594-601.	5.3	59
101	The Diabatic Picture of Electron Transfer, Reaction Barriers, and Molecular Dynamics. <i>Annual Review of Physical Chemistry</i> , 2010, 61, 149-170.	10.8	280
102	Nonlocal van der Waals density functional: The simpler the better. <i>Journal of Chemical Physics</i> , 2010, 133, 244103.	3.0	974
103	Electronic Properties of Disordered Organic Semiconductors via QM/MM Simulations. <i>Accounts of Chemical Research</i> , 2010, 43, 995-1004.	15.6	90
104	Implementation and assessment of a simple nonlocal van der Waals density functional. <i>Journal of Chemical Physics</i> , 2010, 132, 164113.	3.0	74
105	Nonlocal van der Waals Density Functional Made Simple. <i>Physical Review Letters</i> , 2009, 103, 063004.	7.8	329
106	On the Singlet-Triplet Splitting of Geminate Electron-Hole Pairs in Organic Semiconductors. <i>Journal of the American Chemical Society</i> , 2008, 130, 3420-3427.	13.7	100
107	Self-consistent implementation of a nonlocal van der Waals density functional with a Gaussian basis set. <i>Journal of Chemical Physics</i> , 2008, 129, 014106.	3.0	74
108	Rydberg energies using excited state density functional theory. <i>Journal of Chemical Physics</i> , 2008, 129, 124112.	3.0	49

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109	Extracting electron transfer coupling elements from constrained density functional theory. Journal of Chemical Physics, 2006, 125, 164105.	3.0	281
110	Advances in methods and algorithms in a modern quantum chemistry program package. Physical Chemistry Chemical Physics, 2006, 8, 3172-3191.	2.8	2,597
111	Constrained Density Functional Theory and Its Application in Long-Range Electron Transfer. Journal of Chemical Theory and Computation, 2006, 2, 765-774.	5.3	250
112	Direct Calculation of Electron Transfer Parameters through Constrained Density Functional Theory. Journal of Physical Chemistry A, 2006, 110, 9212-9218.	2.5	212
113	Direct optimization method to study constrained systems within density-functional theory. Physical Review A, 2005, 72, .	2.5	447
114	Semiclassical representations of electronic structure and dynamics. Journal of Chemical Physics, 2004, 120, 579-589.	3.0	10
115	GEOMETRIC DIRECT MINIMIZATION OF HARTREE-FOCK CALCULATIONS INVOLVING OPEN SHELL WAVEFUNCTIONS WITH SPIN RESTRICTED ORBITALS. Journal of Theoretical and Computational Chemistry, 2002, 01, 255-261.	1.8	10
116	A geometric approach to direct minimization. Molecular Physics, 2002, 100, 1713-1721.	1.7	137
117	Connections between coupled cluster and generalized valence bond theories. Journal of Chemical Physics, 2001, 115, 7814-7821.	3.0	54
118	Two-body coupled cluster expansions. Journal of Chemical Physics, 2001, 115, 5033-5040.	3.0	51
119	Q-Chem 2.0: a high-performance ab initio electronic structure program package. Journal of Computational Chemistry, 2000, 21, 1532-1548.	3.3	617
120	Benchmark variational coupled cluster doubles results. Journal of Chemical Physics, 2000, 113, 8873-8879.	3.0	114
121	A nonorthogonal approach to perfect pairing. Journal of Chemical Physics, 2000, 112, 5633-5638.	3.0	40
122	Q-Chem 2.0: a high-performance ab initio electronic structure program package. , 2000, 21, 1532.		2