

Galen T Craven

List of Publications by Year in descending order

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Version: 2024-02-01

24
papers

530
citations

567281

15
h-index

642732

23
g-index

25
all docs

25
docs citations

25
times ranked

271
citing authors

#	ARTICLE	IF	CITATIONS
1	Lagrangian Descriptors of Thermalized Transition States on Time-Varying Energy Surfaces. <i>Physical Review Letters</i> , 2015, 115, 148301.	7.8	68
2	Electron transfer across a thermal gradient. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, 9421-9429.	7.1	50
3	Deconstructing field-induced ketene isomerization through Lagrangian descriptors. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 4008-4018.	2.8	49
4	Transition state geometry of driven chemical reactions on time-dependent double-well potentials. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 30270-30281.	2.8	37
5	Lagrangian descriptors of driven chemical reaction manifolds. <i>Physical Review E</i> , 2017, 96, 022222.	2.1	35
6	Electron transfer at thermally heterogeneous molecule-metal interfaces. <i>Journal of Chemical Physics</i> , 2017, 146, .	3.0	31
7	Electron-Transfer-Induced Thermal and Thermoelectric Rectification. <i>Physical Review Letters</i> , 2018, 121, 247704.	7.8	31
8	Chemical reactions induced by oscillating external fields in weak thermal environments. <i>Journal of Chemical Physics</i> , 2015, 142, 074108.	3.0	28
9	Communication: Transition state trajectory stability determines barrier crossing rates in chemical reactions induced by time-dependent oscillating fields. <i>Journal of Chemical Physics</i> , 2014, 141, 041106.	3.0	27
10	Electrothermal Transistor Effect and Cyclic Electronic Currents in Multithermal Charge Transfer Networks. <i>Physical Review Letters</i> , 2017, 118, 207201.	7.8	24
11	Persistence of transition-state structure in chemical reactions driven by fields oscillating in time. <i>Physical Review E</i> , 2014, 89, 040801.	2.1	23
12	Machine learning approaches for structural and thermodynamic properties of a Lennard-Jones fluid. <i>Journal of Chemical Physics</i> , 2020, 153, 104502.	3.0	22
13	Wiedemann-Franz Law for Molecular Hopping Transport. <i>Nano Letters</i> , 2020, 20, 989-993.	9.1	21
14	Transition state theory for activated systems with driven anharmonic barriers. <i>Journal of Chemical Physics</i> , 2017, 147, 074104.	3.0	19
15	Electron-transfer-induced and phononic heat transport in molecular environments. <i>Journal of Chemical Physics</i> , 2017, 147, 124101.	3.0	18
16	Structure of a tractable stochastic mimic of soft particles. <i>Soft Matter</i> , 2014, 10, 5350-5361.	2.7	11
17	Stochastic dynamics of penetrable rods in one dimension: Occupied volume and spatial order. <i>Journal of Chemical Physics</i> , 2013, 138, 244901.	3.0	9
18	Ex Machina Determination of Structural Correlation Functions. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 4372-4378.	4.6	7

#	ARTICLE	IF	CITATIONS
19	Effective Surface Coverage of Coarse-Grained Soft Matter. <i>Journal of Physical Chemistry B</i> , 2014, 118, 14092-14102.	2.6	5
20	Stochastic dynamics of penetrable rods in one dimension: Entangled dynamics and transport properties. <i>Journal of Chemical Physics</i> , 2015, 142, 154906.	3.0	5
21	Upside/Downside statistical mechanics of nonequilibrium Brownian motion. I. Distributions, moments, and correlation functions of a free particle. <i>Journal of Chemical Physics</i> , 2018, 148, 044101.	3.0	4
22	Upside/Downside statistical mechanics of nonequilibrium Brownian motion. II. Heat transfer and energy partitioning of a free particle. <i>Journal of Chemical Physics</i> , 2018, 149, 104103.	3.0	3
23	Nonequilibrium structure in sequential assembly. <i>Physical Review E</i> , 2015, 92, 052108.	2.1	2
24	Heat transport induced by electron transfer: A general temperature quantum calculation. <i>Journal of Chemical Physics</i> , 2021, 155, 194104.	3.0	1