

Hao Wu

List of Publications by Year in descending order

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Version: 2024-02-01

22
papers

2,738
citations

516710

16
h-index

713466

21
g-index

24
all docs

24
docs citations

24
times ranked

2261
citing authors

#	ARTICLE	IF	CITATIONS
1	Markov models of molecular kinetics: Generation and validation. <i>Journal of Chemical Physics</i> , 2011, 134, 174105.	3.0	968
2	Boltzmann generators: Sampling equilibrium states of many-body systems with deep learning. <i>Science</i> , 2019, 365, .	12.6	332
3	VAMPnets for deep learning of molecular kinetics. <i>Nature Communications</i> , 2018, 9, 5.	12.8	330
4	Data-Driven Model Reduction and Transfer Operator Approximation. <i>Journal of Nonlinear Science</i> , 2018, 28, 985-1010.	2.1	192
5	Multiensemble Markov models of molecular thermodynamics and kinetics. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, E3221-30.	7.1	173
6	Variational Approach for Learning Markov Processes from Time Series Data. <i>Journal of Nonlinear Science</i> , 2020, 30, 23-66.	2.1	156
7	Estimation and uncertainty of reversible Markov models. <i>Journal of Chemical Physics</i> , 2015, 143, 174101.	3.0	102
8	Variational Koopman models: Slow collective variables and molecular kinetics from short off-equilibrium simulations. <i>Journal of Chemical Physics</i> , 2017, 146, 154104.	3.0	100
9	Combining experimental and simulation data of molecular processes via augmented Markov models. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, 8265-8270.	7.1	93
10	Markov state models from short non-equilibrium simulations—Analysis and correction of estimation bias. <i>Journal of Chemical Physics</i> , 2017, 146, .	3.0	51
11	Variational selection of features for molecular kinetics. <i>Journal of Chemical Physics</i> , 2019, 150, 194108.	3.0	46
12	Deeptime: a Python library for machine learning dynamical models from time series data. <i>Machine Learning: Science and Technology</i> , 2022, 3, 015009.	5.0	37
13	Identification of kinetic order parameters for non-equilibrium dynamics. <i>Journal of Chemical Physics</i> , 2019, 150, 164120.	3.0	31
14	Rapid Calculation of Molecular Kinetics Using Compressed Sensing. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 2771-2783.	5.3	24
15	Optimal Data-Driven Estimation of Generalized Markov State Models for Non-Equilibrium Dynamics. <i>Computation</i> , 2018, 6, 22.	2.0	20
16	Normalizing field flows: Solving forward and inverse stochastic differential equations using physics-informed flow models. <i>Journal of Computational Physics</i> , 2022, 461, 111202.	3.8	15
17	Gaussian Markov transition models of molecular kinetics. <i>Journal of Chemical Physics</i> , 2015, 142, 084104.	3.0	14
18	Flooding and Overflow Mitigation Using Deep Reinforcement Learning Based on Koopman Operator of Urban Drainage Systems. <i>Water Resources Research</i> , 2022, 58, .	4.2	13

#	ARTICLE	IF	CITATIONS
19	Kernel Embedding Based Variational Approach for Low-Dimensional Approximation of Dynamical Systems. Computational Methods in Applied Mathematics, 2021, 21, 635-659.	0.8	7
20	Maximum margin clustering for state decomposition of metastable systems. Neurocomputing, 2015, 164, 5-22.	5.9	2
21	$\hat{\mu}$ -Dependent controllability for two time-scale systems. Tsinghua Science and Technology, 2009, 14, 271-280.	6.1	0
22	Adaptive hypersonic flight control via back-stepping and Kriging estimation. , 2011, , .		0