

Dirk Lebiedz

List of Publications by Year in descending order

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Version: 2024-02-01

25
papers

555
citations

687363

13
h-index

642732

23
g-index

25
all docs

25
docs citations

25
times ranked

549
citing authors

#	ARTICLE	IF	CITATIONS
1	Flow curvature manifold and energy of generalized Li ⁺ Arnold systems. <i>Chaos, Solitons and Fractals</i> , 2022, 161, 112354.	5.1	3
2	Towards Differential Geometric Characterization of Slow Invariant Manifolds in Extended Phase Space: Sectional Curvature and Flow Invariance. <i>SIAM Journal on Applied Dynamical Systems</i> , 2018, 17, 732-753.	1.6	3
3	An Optimization Approach to Kinetic Model Reduction for Combustion Chemistry. <i>Flow, Turbulence and Combustion</i> , 2014, 92, 885-902.	2.6	3
4	A robust optimization approach to experimental design for model discrimination of dynamical systems. <i>Mathematical Programming</i> , 2013, 141, 405-433.	2.4	7
5	Simplified Reaction Models for Combustion in Gas Turbine Combustion Chambers. <i>Fluid Mechanics and Its Applications</i> , 2013, , 161-182.	0.2	0
6	A Continuation Method for the Efficient Solution of Parametric Optimization Problems in Kinetic Model Reduction. <i>SIAM Journal of Scientific Computing</i> , 2013, 35, A1584-A1603.	2.8	6
7	Robust Optimal Design of Synthetic Biological Networks. <i>Methods in Molecular Biology</i> , 2012, 813, 45-55.	0.9	3
8	A Variational Principle for Computing Slow Invariant Manifolds in Dissipative Dynamical Systems. <i>SIAM Journal of Scientific Computing</i> , 2011, 33, 703-720.	2.8	16
9	Geometric Criteria for Model Reduction in Chemical Kinetics via Optimization of Trajectories. <i>Lecture Notes in Computational Science and Engineering</i> , 2011, , 241-252.	0.3	4
10	Minimal curvature trajectories: Riemannian geometry concepts for slow manifold computation in chemical kinetics. <i>Journal of Computational Physics</i> , 2010, 229, 6512-6533.	3.8	22
11	Entropy-Related Extremum Principles for Model Reduction of Dissipative Dynamical Systems. <i>Entropy</i> , 2010, 12, 706-719.	2.2	22
12	Predicted Auxiliary Navigation Mechanism of Peritrichously Flagellated Chemotactic Bacteria. <i>PLoS Computational Biology</i> , 2010, 6, e1000717.	3.2	42
13	Role of Translational Coupling in Robustness of Bacterial Chemotaxis Pathway. <i>PLoS Biology</i> , 2009, 7, e1000171.	5.6	54
14	Oscillatory NAD(P)H Waves and Calcium Oscillations in Neutrophils? A Modeling Study of Feasibility. <i>Biophysical Journal</i> , 2009, 96, 417-428.	0.5	15
15	Protein exchange dynamics at chemoreceptor clusters in <i>Escherichia coli</i> . <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 6403-6408.	7.1	75
16	Dependence of Bacterial Chemotaxis on Gradient Shape and Adaptation Rate. <i>PLoS Computational Biology</i> , 2008, 4, e1000242.	3.2	97
17	Targeting characteristic wave properties in reaction-diffusion systems by optimization of external forcing. <i>Physical Review E</i> , 2007, 76, 056211.	2.1	2
18	Influence of Coadsorbates on the NO Dissociation on a Rhodium(311) Surface. <i>ChemPhysChem</i> , 2005, 6, 2513-2521.	2.1	15

#	ARTICLE	IF	CITATIONS
19	Automatic network coupling analysis for dynamical systems based on detailed kinetic models. <i>Physical Review E</i> , 2005, 72, 041911.	2.1	9
20	Coverage dependence of oxygen decomposition and surface diffusion on rhodium (111): A DFT study. <i>Journal of Chemical Physics</i> , 2005, 122, 034710.	3.0	56
21	Influence of initial oxygen coverage and magnetic moment on the NO decomposition on rhodium (111). <i>Journal of Chemical Physics</i> , 2005, 122, 154702.	3.0	29
22	Specific external forcing of spatiotemporal dynamics in reaction-diffusion systems. <i>Chaos</i> , 2005, 15, 023901.	2.5	5
23	Linear relationship between activation energies and reaction energies for coverage-dependent dissociation reactions on rhodium surfaces. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 2552.	2.8	21
24	Dynamic control and information processing in chemical reaction systems by tuning self-organization behavior. <i>Chaos</i> , 2004, 14, 611-616.	2.5	7
25	Manipulation of Self-Aggregation Patterns and Waves in a Reaction-Diffusion System by Optimal Boundary Control Strategies. <i>Physical Review Letters</i> , 2003, 91, 208301.	7.8	39