

# Thomas Wuest

## List of Publications by Year in descending order

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29  
papers

653  
citations

687363

13  
h-index

552781

26  
g-index

29  
all docs

29  
docs citations

29  
times ranked

397  
citing authors

#	ARTICLE	IF	CITATIONS
1	Generic, Hierarchical Framework for Massively Parallel Wang-Landau Sampling. <i>Physical Review Letters</i> , 2013, 110, 210603.	7.8	100
2	Versatile Approach to Access the Low Temperature Thermodynamics of Lattice Polymers and Proteins. <i>Physical Review Letters</i> , 2009, 102, 178101.	7.8	88
3	Collapse transitions in a flexible homopolymer chain: Application of the Wang-Landau algorithm. <i>Physical Review E</i> , 2010, 81, 011802.	2.1	79
4	The HP model of protein folding: A challenging testing ground for Wang-Landau sampling. <i>Computer Physics Communications</i> , 2008, 179, 124-127.	7.5	48
5	Generic folding and transition hierarchies for surface adsorption of hydrophobic-polar lattice model proteins. <i>Physical Review E</i> , 2013, 87, 012706.	2.1	44
6	Unraveling the Beautiful Complexity of Simple Lattice Model Polymers and Proteins Using Wang-Landau Sampling. <i>Journal of Statistical Physics</i> , 2011, 144, 638-651.	1.2	35
7	Alignment of radicals into chains by a Markov mechanism for polarity formation. <i>CrystEngComm</i> , 2002, 4, 432-439.	2.6	29
8	Monte Carlo simulations of the HP model (the $\alpha$ -helix model of protein folding). <i>Computer Physics Communications</i> , 2011, 182, 1896-1899.	7.5	29
9	Application of the Wang-Landau algorithm to the dimerization of glycophorin A. <i>Journal of Chemical Physics</i> , 2009, 130, 215106.	3.0	25
10	A Wang-Landau study of the phase transitions in a flexible homopolymer. <i>Computer Physics Communications</i> , 2009, 180, 587-589.	7.5	25
11	Prediction of Growth-Induced Polarity in Centrosymmetric Molecular Crystals Using Force Field Methods. <i>Chemistry of Materials</i> , 2005, 17, 85-94.	6.7	19
12	Numerical integration using Wang-Landau sampling. <i>Computer Physics Communications</i> , 2007, 177, 524-529.	7.5	18
13	Can Mono Domain Polar Molecular Crystals Exist?. <i>Crystal Growth and Design</i> , 2012, 12, 5211-5218.	3.0	15
14	Conformational transitions of a confined lattice protein: A Wang-Landau study. <i>Journal of Physics: Conference Series</i> , 2012, 402, 012048.	0.4	13
15	Influence of Solid Solution Formation on Polarity: Molecular Modeling Investigation of the System 4-Chloro-4-nitrostilbene/4,4-Dinitrostilbene. <i>Journal of Physical Chemistry B</i> , 2005, 109, 12582-12589.	2.6	12
16	A stochastic principle behind polar properties of condensed molecular matter. <i>New Journal of Chemistry</i> , 2013, 37, 2229.	2.8	12
17	How Symmetrical Molecules Can Induce Polarity: On the Paradox of Dilution. <i>Crystal Growth and Design</i> , 2005, 5, 93-97.	3.0	11
18	The role of chain-stiffness in lattice protein models: A replica-exchange Wang-Landau study. <i>Journal of Chemical Physics</i> , 2018, 149, 125101.	3.0	11

#	ARTICLE	IF	CITATIONS
19	Growth-induced polarity formation in solid solutions of organic molecules: Markov mean-field model and Monte Carlo simulations. <i>Journal of Chemical Physics</i> , 2005, 122, 084715.	3.0	9
20	Monte Carlo simulations of systems with complex energy landscapes. <i>Computer Physics Communications</i> , 2009, 180, 475-479.	7.5	6
21	Growth-induced polarity formation in two-component crystals of organic molecules: A statistical analysis. <i>Journal of Physics and Chemistry of Solids</i> , 2006, 67, 2517-2527.	4.0	5
22	Effect of reduced cooperativity on growth-induced polarity formation: a comparison between different growth models. <i>Philosophical Magazine</i> , 2007, 87, 1683-1703.	1.6	4
23	WANG'S LANDAU SIMULATIONS OF ADSORBED AND CONFINED LATTICE PROTEINS. <i>International Journal of Modern Physics C</i> , 2012, 23, 1240008.	1.7	4
24	Surface adsorption of lattice HP proteins: Thermodynamics and structural transitions using Wang-Landau sampling. <i>Journal of Physics: Conference Series</i> , 2012, 402, 012046.	0.4	3
25	Wang's Landau sampling of the interplay between surface adsorption and folding of HP lattice proteins. <i>Molecular Simulation</i> , 2014, 40, 640-655.	2.0	3
26	Effects of an external electrical field on the polarization of growing organic crystals: a theoretical study. <i>Chemical Physics Letters</i> , 2003, 377, 340-346.	2.6	2
27	Universality behaviour for polarity formation in channel-type inclusion compounds. <i>Journal of Mathematical Chemistry</i> , 2009, 45, 869-881.	1.5	2
28	Thermodynamics and structural properties of a confined HP protein determined by Wang-Landau simulation. <i>Journal of Physics: Conference Series</i> , 2013, 454, 012071.	0.4	2
29	MONTE CARLO SIMULATIONS OF PROTEIN MODELS: AT THE INTERFACE BETWEEN STATISTICAL PHYSICS AND BIOLOGY. , 2010, , .		0