

# Gerard T Barkema

## List of Publications by Year in descending order

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Version: 2024-02-01

31

papers

456

citations

687220

13

h-index

713332

21

g-index

31

all docs

31

docs citations

31

times ranked

490

citing authors

#	ARTICLE	IF	CITATIONS
1	Comparison of cluster algorithms for the bond-diluted Ising model. <i>Physical Review E</i> , 2022, 105, 015313.	0.8	2
2	Structural dynamics of polycrystalline graphene. <i>Physical Review E</i> , 2022, 105, 044116.	0.8	3
3	Rupture of amorphous graphene via void formation. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 16966-16972.	1.3	2
4	Efficient simulation of semiflexible polymers with stiff bonds. <i>Physical Review E</i> , 2017, 95, 012502.	0.8	2
5	Structural characterization of carbon nanotubes via the vibrational density of states. <i>Carbon</i> , 2017, 118, 58-65.	5.4	15
6	An introduction to Monte Carlo methods. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2015, 418, 78-87.	1.2	48
7	Monte Carlo methods beyond detailed balance. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2015, 418, 88-93.	1.2	9
8	SAWdoubler: A program for counting self-avoiding walks. <i>Computer Physics Communications</i> , 2013, 184, 891-898.	3.0	11
9	A model for the dynamics of extensible semiflexible polymers. <i>Journal of Statistical Mechanics: Theory and Experiment</i> , 2012, 2012, P12019.	0.9	8
10	Critical exponents of the pair contact process with diffusion. <i>Journal of Statistical Mechanics: Theory and Experiment</i> , 2012, 2012, P03009.	0.9	8
11	Saturation of front propagation in a reaction diffusion process describing plasma damage in porous low- $\langle \text{mml:math} \text{ xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline" } \rangle \langle \text{mml:mrow} \langle \text{mml:mi} k \rangle \langle \text{mml:mi} \rangle \rangle \langle \text{mml:mrow} \rangle \langle / \text{mml:math} \rangle$ materials. <i>Physical Review B</i> , 2011, 83, .	1.1	4
12	The equilibrium winding angle of a polymer around a bar. <i>Journal of Statistical Mechanics: Theory and Experiment</i> , 2011, 2011, P10020.	0.9	12
13	Exact enumeration of self-avoiding walks. <i>Journal of Statistical Mechanics: Theory and Experiment</i> , 2011, 2011, P06019.	0.9	36
14	Structural modes of a polymer in the repton model. <i>Journal of Chemical Physics</i> , 2011, 134, 154901.	1.2	8
15	Simulations of color development in tinted paints. <i>Journal of Colloid and Interface Science</i> , 2010, 344, 256-260.	5.0	0
16	Elastic lattice polymers. <i>Physical Review E</i> , 2010, 81, 061801.	0.8	9
17	Frequency-dependent stiffening of semiflexible networks: A dynamical nonaffine to affine transition. <i>Physical Review E</i> , 2010, 82, 061902.	0.8	29
18	Non-Markovian dynamics of clusters during nucleation. <i>Physical Review E</i> , 2009, 79, 062101.	0.8	7

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19	Inverse Langmuir method for oligonucleotide microarray analysis. <i>BMC Bioinformatics</i> , 2009, 10, 64.	1.2	13
20	Analysis of the variability of the axial dipole moment of a numerical geodynamo model. <i>Physics of the Earth and Planetary Interiors</i> , 2009, 173, 228-232.	0.7	14
21	Linear model for fast background subtraction in oligonucleotide microarrays. <i>Algorithms for Molecular Biology</i> , 2009, 4, 15.	0.3	1
22	Monte Carlo study of multiply crosslinked semiflexible polymer networks. <i>Physical Review E</i> , 2008, 78, 051801.	0.8	44
23	Universality class of the pair contact process with diffusion. <i>Physical Review E</i> , 2008, 78, 031129.	0.8	15
24	Modeling background intensity in DNA microarrays. <i>Physical Review E</i> , 2008, 77, 061915.	0.8	10
25	Physical-Chemistry-Based Analysis of Affymetrix Microarray Data. <i>Journal of Physical Chemistry B</i> , 2006, 110, 22786-22795.	1.2	13
26	Spinodal decomposition via surface diffusion in polymer mixtures. <i>Physical Review E</i> , 2006, 74, 011804.	0.8	10
27	Comment on "Solving the riddle of the bright mismatches: Labeling and effective binding in oligonucleotide arrays". <i>Physical Review E</i> , 2006, 73, 063901; author reply 063902.	0.8	11
28	Phase separation driven by surface diffusion: A Monte Carlo study. <i>Physical Review E</i> , 2005, 72, 046131.	0.8	37
29	Nucleation times in the two-dimensional Ising model. <i>Physical Review E</i> , 2005, 71, 031601.	0.8	36
30	Magnetization reversal times in the two-dimensional Ising model. <i>Physical Review E</i> , 2003, 67, 026119.	0.8	14
31	Universality in the pair contact process with diffusion. <i>Physical Review E</i> , 2003, 68, 036113.	0.8	25