

# Gunaretnam Rajagopal

## List of Publications by Year in descending order

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47  
papers

4,333  
citations

172457

29  
h-index

223800

46  
g-index

48  
all docs

48  
docs citations

48  
times ranked

3512  
citing authors

#	ARTICLE	IF	CITATIONS
1	Quantum Monte Carlo simulations of solids. <i>Reviews of Modern Physics</i> , 2001, 73, 33-83.	45.6	1,813
2	Finite-size effects and Coulomb interactions in quantum Monte Carlo calculations for homogeneous systems with periodic boundary conditions. <i>Physical Review B</i> , 1996, 53, 1814-1832.	3.2	265
3	Calculations of Silicon Self-Interstitial Defects. <i>Physical Review Letters</i> , 1999, 83, 2351-2354.	7.8	166
4	Patterns and barriers for fission of charged small metal clusters. <i>Physical Review Letters</i> , 1991, 67, 3058-3061.	7.8	126
5	Exchange and correlation in silicon. <i>Physical Review B</i> , 1998, 57, 8972-8982.	3.2	117
6	Optical spectra of localized excess electrons in alkali halide clusters. <i>Physical Review Letters</i> , 1990, 64, 2933-2936.	7.8	106
7	The tumor suppressor p53: Cancer and aging. <i>Cell Cycle</i> , 2008, 7, 842-847.	2.6	106
8	Metallization of ionic clusters. <i>Physical Review Letters</i> , 1991, 67, 727-730.	7.8	104
9	Elimination of Coulomb finite-size effects in quantum many-body simulations. <i>Physical Review B</i> , 1997, 55, R4851-R4854.	3.2	99
10	Finite-size errors in quantum many-body simulations of extended systems. <i>Physical Review B</i> , 1999, 59, 1917-1929.	3.2	98
11	Monte Carlo energy and variance-minimization techniques for optimizing many-body wave functions. <i>Physical Review B</i> , 1999, 59, 12344-12351.	3.2	94
12	Born-Oppenheimer dynamics using density-functional theory: Equilibrium and fragmentation of small sodium clusters. <i>Journal of Chemical Physics</i> , 1991, 94, 608-616.	3.0	85
13	Carbon clusters near the crossover to fullerene stability. <i>Physical Review B</i> , 2000, 62, 15394-15397.	3.2	83
14	Identifying microRNA/mRNA dysregulations in ovarian cancer. <i>BMC Research Notes</i> , 2012, 5, 164.	1.4	82
15	Diffusion quantum Monte Carlo calculations of the excited states of silicon. <i>Physical Review B</i> , 1998, 57, 12140-12144.	3.2	81
16	Variational and diffusion quantum Monte Carlo calculations at nonzero wave vectors: Theory and application to diamond-structure germanium. <i>Physical Review B</i> , 1995, 51, 10591-10600.	3.2	79
17	Quantum Monte Carlo Calculations for Solids Using SpecialkPoints Methods. <i>Physical Review Letters</i> , 1994, 73, 1959-1962.	7.8	72
18	Quantum Monte Carlo Investigation of Exchange and Correlation in Silicon. <i>Physical Review Letters</i> , 1997, 78, 3350-3353.	7.8	65

#	ARTICLE	IF	CITATIONS
19	Optimized wave functions for quantum Monte Carlo studies of atoms and solids. <i>Physical Review B</i> , 1996, 53, 9640-9648.	3.2	57
20	Ab initio calculations of the cohesive energy and the bulk modulus of aluminium. <i>Journal of Physics Condensed Matter</i> , 2002, 14, 8787-8793.	1.8	53
21	Negative Feedback Governs Gonadotrope Frequency-Decoding of Gonadotropin Releasing Hormone Pulse-Frequency. <i>PLoS ONE</i> , 2009, 4, e7244.	2.5	47
22	Homology-based hydrogen bond information improves crystallographic structures in the <scp>PDB</scp>. <i>Protein Science</i> , 2018, 27, 798-808.	7.6	41
23	The p53 knowledgebase: an integrated information resource for p53 research. <i>Oncogene</i> , 2007, 26, 1517-1521.	5.9	40
24	Quantum Monte Carlo calculations of the one-body density matrix and excitation energies of silicon. <i>Physical Review B</i> , 1998, 57, 15293-15302.	3.2	39
25	The path from big data to precision medicine. <i>Expert Review of Precision Medicine and Drug Development</i> , 2016, 1, 129-143.	0.7	37
26	An Optimized Ewald Method for Long-Ranged Potentials. <i>Journal of Computational Physics</i> , 1994, 115, 399-405.	3.8	34
27	Relativistic corrections to atomic energies from quantum Monte Carlo calculations. <i>Physical Review A</i> , 1995, 51, 1898-1904.	2.5	33
28	Hybrid simulations of stochastic reaction-diffusion processes for modeling intracellular signaling pathways. <i>Physical Review E</i> , 2006, 74, 051910.	2.1	33
29	ArhGAP9, a novel MAP kinase docking protein, inhibits Erk and p38 activation through WW domain binding. <i>Journal of Molecular Signaling</i> , 2007, 2, 1.	0.5	33
30	Quantifying Forces Mediated by Integral Tight Junction Proteins in Cell-Cell Adhesion. <i>Experimental Mechanics</i> , 2009, 49, 3-9.	2.0	29
31	Whole-genome sequencing analysis of phenotypic heterogeneity and anticipation in Li-Fraumeni cancer predisposition syndrome. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 15497-15501.	7.1	29
32	Pseudopotentials for correlated-electron calculations. <i>Physical Review B</i> , 2000, 62, 13347-13355.	3.2	27
33	Dynamics, Spectra, and Relaxation Phenomena of Excess Electrons in Clusters. <i>Israel Journal of Chemistry</i> , 1990, 30, 85-105.	2.3	22
34	Quantum Monte Carlo Calculations of the Energy of the Relativistic Homogeneous Electron Gas. <i>Physical Review Letters</i> , 1996, 77, 1099-1102.	7.8	16
35	Comment on "Quantum Monte Carlo study of the dipole moment of CO". <i>J. Chem. Phys.</i> 110, 11700 (1999). <i>Journal of Chemical Physics</i> , 2000, 112, 4419-4420.	3.0	16
36	Facilitating arrhythmia simulation: the method of quantitative cellular automata modeling and parallel running. <i>BioMedical Engineering OnLine</i> , 2004, 3, 29.	2.7	16

#	ARTICLE	IF	CITATIONS
37	3D Meshless Prostate Segmentation and Registration in Image Guided Radiotherapy. Lecture Notes in Computer Science, 2009, 12, 43-50.	1.3	15
38	Group-based variant calling leveraging next-generation supercomputing for large-scale whole-genome sequencing studies. BMC Bioinformatics, 2015, 16, 304.	2.6	12
39	An exact calculation of the flavour-changing quark-photon vertex. Physics Letters, Section B: Nuclear, Elementary Particle and High-Energy Physics, 1985, 156, 405-410.	4.1	11
40	Quantum Monte Carlo calculations for ground and excited states. International Journal of Quantum Chemistry, 2002, 86, 218-225.	2.0	11
41	A Quantum Monte Carlo Approach to the Adiabatic Connection Method. Advances in Quantum Chemistry, 1998, 33, 189-207.	0.8	10
42	Oscillations in intracellular signaling cascades. Physical Review E, 2007, 75, 061901.	2.1	10
43	First-principles calculations of the adsorbate-induced surface stress of. Surface Science, 1997, 372, 179-184.	1.9	9
44	Molecular force spectroscopy of homophilic nectin-1 interactions. Biochemical and Biophysical Research Communications, 2007, 362, 886-892.	2.1	6
45	Quantum Monte Carlo Study of Silicon Self-interstitial Defects. VLSI Design, 2001, 13, 229-235.	0.5	5
46	Gastric Cancer (Biomarkers) Knowledgebase (GCBKB): A Curated and Fully Integrated Knowledgebase of Putative Biomarkers Related to Gastric Cancer. Biomarker Insights, 2006, 1, 117727190600100.	2.5	1
47	Identification and Modeling of TEL-AML1 (ETV6-RUNX1) Molecular Signature In Acute Lymphoblastic Leukemia. Blood, 2010, 116, 2500-2500.	1.4	0