

Antonio Caliri

List of Publications by Year in descending order

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33
papers

242
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1163117

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1058476

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times ranked

339
citing authors

#	ARTICLE	IF	CITATIONS
1	Phospholipase A2 Isolated from the Venom of <i>Crotalus durissus terrificus</i> Inactivates Dengue virus and Other Enveloped Viruses by Disrupting the Viral Envelope. <i>PLoS ONE</i> , 2014, 9, e112351.	2.5	53
2	LMProt: An Efficient Algorithm for Monte Carlo Sampling of Protein Conformational Space. <i>Biophysical Journal</i> , 2004, 87, 1567-1577.	0.5	15
3	Membrane vesiculation induced by proteins of the dengue virus envelope studied by molecular dynamics simulations. <i>Journal of Physics Condensed Matter</i> , 2017, 29, 504002.	1.8	14
4	The predictive power of $r(0)$ in an epidemic probabilistic system. <i>Journal of Biological Physics</i> , 2003, 29, 63-75.	1.5	12
5	Solution of deterministic stochastic epidemic models by dynamical Monte Carlo method. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2000, 282, 546-558.	2.6	11
6	Partition function of long-ranged ferromagnet and of a long-ranged fully frustrated model. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1984, 106, 74-76.	2.1	10
7	Phase Transition in a System of Hard Disks by Monte Carlo Simulation. <i>Physical Review Letters</i> , 1987, 58, 2312-2314.	7.8	10
8	Deterministic folding: The role of entropic forces and steric specificities. <i>Journal of Chemical Physics</i> , 2001, 114, 4235-4242.	3.0	9
9	Extensive structural change of the envelope protein of dengue virus induced by a tuned ionic strength: conformational and energetic analyses. <i>Journal of Computer-Aided Molecular Design</i> , 2012, 26, 1311-1325.	2.9	9
10	In silico assessment of S100A12 monomer and dimer structural dynamics: implications for the understanding of its metal-induced conformational changes. <i>Journal of Biological Inorganic Chemistry</i> , 2014, 19, 1113-1120.	2.6	9
11	The 3-phenylcoumarin derivative 6,7-dihydroxy-3-[3,4-methylenedioxyphenyl]-coumarin downmodulates the Fc γ R- and CR-mediated oxidative metabolism and elastase release in human neutrophils: Possible mechanisms underlying inhibition of the formation and release of neutrophil extracellular traps. <i>Free Radical Biology and Medicine</i> , 2018, 115, 421-435.	2.9	9
12	The water factor in the protein-folding problem. <i>Brazilian Journal of Physics</i> , 2004, 34, 90-101.	1.4	7
13	Entropic force and folding of macromolecules. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1996, 220, 178-182.	2.1	6
14	Temporal duration and event size distribution at the epidemic threshold. <i>Journal of Biological Physics</i> , 1999, 25, 309-324.	1.5	6
15	Steric constraints as folding coadjuvant. <i>Physical Review E</i> , 2003, 67, 031901.	2.1	6
16	Topology-dependent protein folding rates analyzed by a stereochemical model. <i>Journal of Chemical Physics</i> , 2005, 123, 154906.	3.0	6
17	A new technique for grand canonical Monte Carlo simulation: Application in a hard disk system. <i>Journal of Chemical Physics</i> , 1989, 91, 6328-6336.	3.0	5
18	Geometrical effects on folding of macromolecules. <i>Journal of Chemical Physics</i> , 1997, 106, 7856-7861.	3.0	5

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19	Epidemic phase and the site percolation with distant-neighbor interactions. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1998, 238, 54-58.	2.1	5
20	Distinct conformational properties determined by implicit and explicit representation of protein-solvent interactions. An analytical and computer simulation study. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2009, 388, 4097-4104.	2.6	5
21	The role of disulfide bridges in the 3-D structures of the antimicrobial peptides gomesin and protegrin-1: a molecular dynamics study. <i>Genetics and Molecular Research</i> , 2008, 7, 1070-1088.	0.2	5
22	Effect of local thermal fluctuations on folding kinetics: A study from the perspective of nonextensive statistical mechanics. <i>Physical Review E</i> , 2011, 84, 041903.	2.1	4
23	Nonextensive statistical mechanics applied to protein folding problem: kinetics aspects. <i>Brazilian Journal of Physics</i> , 2009, 39, .	1.4	4
24	Two-dimensional chain folding-random energy interaction. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1993, 183, 327-331.	2.1	3
25	Geometric constraints in polymer chains: analysis on the pearl-necklace model by Monte Carlo simulation. <i>Computational and Theoretical Chemistry</i> , 1995, 335, 123-127.	1.5	3
26	Extended secondary structures in proteins. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2014, 1844, 384-388.	2.3	3
27	Free energy of a spin glass. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1985, 108, 111-114.	2.1	2
28	The Early Events of the Protein Folding Process. <i>Current Physical Chemistry</i> , 2013, 3, 69-76.	0.2	2
29	Theory and calculations for a spin glass. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1985, 109, 282-288.	2.1	1
30	Thermal properties of the $\hat{A} \pm J$ infinite-range Sherrington-Kirkpatrick model by analysis of the partition function zeros. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 1991, 154, 287-292.	2.1	1
31	Entropic formulation for the protein folding process: Hydrophobic stability correlates with folding rates. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2018, 490, 1111-1124.	2.6	1
32	Percolao e o fenmeno epidmico: uma abordagem temporal e espacial da difuso de doenas. <i>Scientia Agricola</i> , 1998, 55, 418-427.	1.2	1
33	Bacillus subtilis Engagement Induced via Sporulation: a Case of Bacterial Communication. <i>Brazilian Journal of Physics</i> , 2022, 52, 1.	1.4	0