

Leonor Saiz

List of Publications by Year in descending order

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48
papers

2,445
citations

257450

24
h-index

206112

48
g-index

53
all docs

53
docs citations

53
times ranked

2800
citing authors

#	ARTICLE	IF	CITATIONS
1	Hydrogen bonding in liquid alcohols: a computer simulation study. <i>Journal of Molecular Structure</i> , 1997, 416, 243-248.	3.6	211
2	Structure and Dynamics of Liquid Ethanol. <i>Journal of Physical Chemistry B</i> , 1997, 101, 78-86.	2.6	194
3	Predicting human olfactory perception from chemical features of odor molecules. <i>Science</i> , 2017, 355, 820-826.	12.6	194
4	Computer Simulation Studies of Model Biological Membranes. <i>Accounts of Chemical Research</i> , 2002, 35, 482-489.	15.6	181
5	Structural Properties of a Highly Polyunsaturated Lipid Bilayer from Molecular Dynamics Simulations. <i>Biophysical Journal</i> , 2001, 81, 204-216.	0.5	136
6	DNA looping in gene regulation: from the assembly of macromolecular complexes to the control of transcriptional noise. <i>Current Opinion in Genetics and Development</i> , 2005, 15, 136-144.	3.3	129
7	Three-dimensional modeling of single stranded DNA hairpins for aptamer-based biosensors. <i>Scientific Reports</i> , 2017, 7, 1178.	3.3	113
8	Dynamics in hydrogen bonded liquids: water and alcohols. <i>Journal of Molecular Liquids</i> , 2002, 96-97, 3-17.	4.9	99
9	Structure of liquid ethylene glycol: A molecular dynamics simulation study with different force fields. <i>Journal of Chemical Physics</i> , 2001, 114, 3187-3199.	3.0	89
10	Electrostatic interactions in a neutral model phospholipid bilayer by molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2002, 116, 3052-3057.	3.0	88
11	Towards an Understanding of Complex Biological Membranes from Atomistic Molecular Dynamics Simulations. <i>Bioscience Reports</i> , 2002, 22, 151-173.	2.4	83
12	DNA looping: the consequences and its control. <i>Current Opinion in Structural Biology</i> , 2006, 16, 344-350.	5.7	79
13	Dielectric properties of liquid ethanol. A computer simulation study. <i>Journal of Chemical Physics</i> , 2000, 113, 2814-2822.	3.0	76
14	Partitioning of Anesthetics into a Lipid Bilayer and their Interaction with Membrane-Bound Peptide Bundles. <i>Biophysical Journal</i> , 2006, 91, 2815-2825.	0.5	67
15	Concentration Effects of Volatile Anesthetics on the Properties of Model Membranes: A Coarse-Grain Approach. <i>Biophysical Journal</i> , 2005, 88, 1524-1534.	0.5	65
16	Inferring the in vivo looping properties of DNA. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 17642-17645.	7.1	54
17	Dynamics and hydrogen bonding in liquid ethanol. <i>Molecular Physics</i> , 1999, 97, 897-905.	1.7	47
18	Stochastic dynamics of macromolecular assembly networks. <i>Molecular Systems Biology</i> , 2006, 2, 2006.0024.	7.2	47

#	ARTICLE	IF	CITATIONS
19	Influence of Anesthetic and Nonimmobilizer Molecules on the Physical Properties of a Polyunsaturated Lipid Bilayer. <i>Journal of Physical Chemistry B</i> , 2003, 107, 14500-14508.	2.6	41
20	Ab initio thermodynamic modeling of distal multisite transcription regulation. <i>Nucleic Acids Research</i> , 2007, 36, 726-731.	14.5	38
21	CplexA: a <i>Mathematica</i> package to study macromolecular-assembly control of gene expression. <i>Bioinformatics</i> , 2010, 26, 2060-2061.	4.1	37
22	Influence of Highly Polyunsaturated Lipid Acyl Chains of Biomembranes on the NMR Order Parameters. <i>Journal of the American Chemical Society</i> , 2001, 123, 7381-7387.	13.7	34
23	Reliable Prediction of Complex Phenotypes from a Modular Design in Free Energy Space: An Extensive Exploration of the <i>lac</i> Operon. <i>ACS Synthetic Biology</i> , 2013, 2, 576-586.	3.8	31
24	Multilevel Deconstruction of the In Vivo Behavior of Looped DNA-Protein Complexes. <i>PLoS ONE</i> , 2007, 2, e355.	2.5	27
25	Multiprotein DNA Looping. <i>Physical Review Letters</i> , 2006, 96, 238103.	7.8	26
26	Systems Biophysics of Gene Expression. <i>Biophysical Journal</i> , 2013, 104, 2574-2585.	0.5	26
27	Computational modelling of Smad-mediated negative feedback and crosstalk in the TGF- β superfamily network. <i>Journal of the Royal Society Interface</i> , 2013, 10, 20130363.	3.4	26
28	The Transmembrane Domain of the Acetylcholine Receptor: Insights from Simulations on Synthetic Peptide Models. <i>Biophysical Journal</i> , 2005, 88, 959-970.	0.5	23
29	Effect of the Pore Region of a Transmembrane Ion Channel on the Physical Properties of a Simple Membrane. <i>Journal of Physical Chemistry B</i> , 2004, 108, 2608-2613.	2.6	21
30	The physics of protein-DNA interaction networks in the control of gene expression. <i>Journal of Physics Condensed Matter</i> , 2012, 24, 193102.	1.8	20
31	Control of gene expression by modulated self-assembly. <i>Nucleic Acids Research</i> , 2011, 39, 6854-6863.	14.5	15
32	Characterization of Negative Feedback Network Motifs in the TGF- β Signaling Pathway. <i>PLoS ONE</i> , 2013, 8, e83531.	2.5	15
33	All-or-none amyloid disassembly via chaperone-triggered fibril unzipping favors clearance of β -synuclein toxic species. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	7.1	15
34	Trafficking Coordinate Description of Intracellular Transport Control of Signaling Networks. <i>Biophysical Journal</i> , 2011, 101, 2315-2323.	0.5	13
35	Protein-protein/DNA interaction networks: versatile macromolecular structures for the control of gene expression. <i>IET Systems Biology</i> , 2008, 2, 247-255.	1.5	12
36	Suppression and enhancement of transcriptional noise by DNA looping. <i>Physical Review E</i> , 2014, 89, 062703.	2.1	11

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37	Computational design of single-stranded DNA hairpin aptamers immobilized on a biosensor substrate. <i>Scientific Reports</i> , 2021, 11, 10984.	3.3	9
38	Cell-to-cell and type-to-type heterogeneity of signaling networks: insights from the crowd. <i>Molecular Systems Biology</i> , 2021, 17, e10402.	7.2	9
39	In silico identification of potential therapeutic targets in the TGF- β 2 signal transduction pathway. <i>Molecular BioSystems</i> , 2014, 10, 537.	2.9	6
40	Clearly Detectable, Kinetically Restricted Solid-Solid Phase Transition in cis-Ceramide Monolayers. <i>Langmuir</i> , 2018, 34, 11749-11758.	3.5	6
41	Reliably quantifying the evolving worldwide dynamic state of the COVID-19 outbreak from death records, clinical parametrization, and demographic data. <i>Scientific Reports</i> , 2021, 11, 19952.	3.3	6
42	Ascertaining the initiation of epidemic resurgences: an application to the COVID-19 second surges in Europe and the Northeast United States. <i>Royal Society Open Science</i> , 2021, 8, 210773.	2.4	6
43	Literature-Based Automated Reconstruction, Expansion, and Refinement of the TGF- β 2 Superfamily Ligand-Receptor Network. <i>Journal of Membrane Biology</i> , 2014, 247, 381-386.	2.1	5
44	Determinants of protein-ligand complex formation in the thyroid hormone receptor β : A molecular dynamics simulation study. <i>Computational and Theoretical Chemistry</i> , 2014, 1038, 57-66.	2.5	4
45	Computational Analysis of the TGF-Beta and BMP Signal Transduction Pathways. <i>Biophysical Journal</i> , 2011, 100, 164a.	0.5	3
46	Insights into Signaling and the Functional Complexity of Biological Membranes. <i>Journal of Membrane Biology</i> , 2017, 250, 335-336.	2.1	2
47	Field-induced force-suppression in ferromagnetic colloids. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2001, 293, 51-58.	2.6	1
48	Computing at the Front-End by Receptor Networks. <i>Cell Systems</i> , 2017, 5, 316-318.	6.2	1