

# Alfredo E Cardenas

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7031282/publications.pdf>

Version: 2024-02-01

18  
papers

526  
citations

623574

14  
h-index

839398

18  
g-index

19  
all docs

19  
docs citations

19  
times ranked

496  
citing authors

#	ARTICLE	IF	CITATIONS
1	Unassisted Transport of <i>N</i> -Acetyl-L-tryptophanamide through Membrane: Experiment and Simulation of Kinetics. <i>Journal of Physical Chemistry B</i> , 2012, 116, 2739-2750.	1.2	59
2	Kinetics of cytochrome C folding: Atomically detailed simulations. <i>Proteins: Structure, Function and Bioinformatics</i> , 2003, 51, 245-257.	1.5	57
3	Calculating Iso-Committer Surfaces as Optimal Reaction Coordinates with Milestoning. <i>Entropy</i> , 2017, 19, 219.	1.1	51
4	MOIL-opt: Energy-Conserving Molecular Dynamics on a GPU/CPU System. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 3072-3082.	2.3	47
5	Membrane Permeation of a Peptide: It Is Better to be Positive. <i>Journal of Physical Chemistry B</i> , 2015, 119, 6412-6420.	1.2	44
6	Bridging the Gap between Long Time Trajectories and Reaction Pathways. <i>Advances in Chemical Physics</i> , 2003, , 93-129.	0.3	37
7	Computational study of peptide permeation through membrane: searching for hidden slow variables. <i>Molecular Physics</i> , 2013, 111, 3565-3578.	0.8	37
8	Modeling kinetics and equilibrium of membranes with fields: Milestoning analysis and implication to permeation. <i>Journal of Chemical Physics</i> , 2014, 141, 054101.	1.2	33
9	The Impact of Protonation on Early Translocation of Anthrax Lethal Factor: Kinetics from Molecular Dynamics Simulations and Milestoning Theory. <i>Journal of the American Chemical Society</i> , 2017, 139, 14837-14840.	6.6	30
10	Physiological Calcium Concentrations Slow Dynamics at the Lipid-Water Interface. <i>Biophysical Journal</i> , 2018, 115, 1541-1551.	0.2	30
11	Atomically Detailed Simulations of Helix Formation with the Stochastic Difference Equation. <i>Biophysical Journal</i> , 2003, 85, 2919-2939.	0.2	28
12	Peptide Permeation across a Phosphocholine Membrane: An Atomically Detailed Mechanism Determined through Simulations and Supported by Experimentation. <i>Journal of Physical Chemistry B</i> , 2022, 126, 2834-2849.	1.2	17
13	Markovian and Non-Markovian Modeling of Membrane Dynamics with Milestoning. <i>Journal of Physical Chemistry B</i> , 2016, 120, 8208-8216.	1.2	16
14	Direct Measurement of the Effect of Cholesterol and 6-Ketocholestanol on the Membrane Dipole Electric Field Using Vibrational Stark Effect Spectroscopy Coupled with Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2017, 121, 3424-3436.	1.2	15
15	A peptide-derived strategy for specifically targeting the mitochondria and ER of cancer cells: a new approach in fighting cancer. <i>Chemical Science</i> , 2022, 13, 6929-6941.	3.7	11
16	Calcium-Lipid Interactions Observed with Isotope-Edited Infrared Spectroscopy. <i>Biophysical Journal</i> , 2020, 118, 2694-2702.	0.2	9
17	Dramatic Shape Changes Occur as Cytochrome <i>c</i> Folds. <i>Journal of Physical Chemistry B</i> , 2020, 124, 8240-8248.	1.2	4
18	Molecular Dynamics at Extended Timescales. <i>Israel Journal of Chemistry</i> , 2014, 54, 1302-1310.	1.0	1