

Andreas Kukol

List of Publications by Year in descending order

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46
papers

1,760
citations

304743

22
h-index

265206

42
g-index

48
all docs

48
docs citations

48
times ranked

2008
citing authors

#	ARTICLE	IF	CITATIONS
1	Lipid Models for United-Atom Molecular Dynamics Simulations of Proteins. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 615-626.	5.3	241
2	Experimentally based orientational refinement of membrane protein models: a structure for the Influenza A M2 H + channel 1 1 Edited by G. von Heijne. <i>Journal of Molecular Biology</i> , 1999, 286, 951-962.	4.2	141
3	Incorporation of the acetylcholine receptor dimer from <i>Torpedo californica</i> in a peptide supported lipid membrane investigated by surface plasmon and fluorescence spectroscopy. <i>Biosensors and Bioelectronics</i> , 1998, 13, 585-591.	10.1	108
4	Site-specific examination of secondary structure and orientation determination in membrane proteins: The peptidic 13C ¹⁸ O group as a novel infrared probe. <i>Biopolymers</i> , 2001, 59, 396-401.	2.4	104
5	Exploring Models of the Influenza A M2 Channel: MD Simulations in a Phospholipid Bilayer. <i>Biophysical Journal</i> , 2000, 78, 55-69.	0.5	98
6	vpu Transmembrane Peptide Structure Obtained by Site-Specific Fourier Transform Infrared Dichroism and Global Molecular Dynamics Searching. <i>Biophysical Journal</i> , 1999, 77, 1594-1601.	0.5	96
7	Model of a Putative Pore: The Pentameric α -Helical Bundle of SARS Coronavirus E Protein in Lipid Bilayers. <i>Biophysical Journal</i> , 2006, 91, 938-947.	0.5	96
8	Solid-State ¹⁷ O NMR of Amino Acids. <i>Journal of Physical Chemistry B</i> , 2004, 108, 9256-9263.	2.6	81
9	How the amyloid- β peptide and membranes affect each other: An extensive simulation study. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2013, 1828, 327-339.	2.6	66
10	Label-free electrical detection of DNA hybridization for the example of influenza virus gene sequences. <i>Analytical Biochemistry</i> , 2008, 374, 143-153.	2.4	55
11	Large-scale analysis of influenza A virus sequences reveals potential drug target sites of non-structural proteins. <i>Journal of General Virology</i> , 2009, 90, 2124-2133.	2.9	53
12	Consensus virtual screening approaches to predict protein ligands. <i>European Journal of Medicinal Chemistry</i> , 2011, 46, 4661-4664.	5.5	53
13	Use of a Single Glycine Residue to Determine the Tilt and Orientation of a Transmembrane Helix. A New Structural Label for Infrared Spectroscopy. <i>Biophysical Journal</i> , 2000, 79, 3139-3143.	0.5	52
14	The structure of the HIV-1 Vpu ion channel: modelling and simulation studies. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2001, 1512, 291-298.	2.6	49
15	A Structure for the Trimeric MHC Class II-associated Invariant Chain Transmembrane Domain. <i>Journal of Molecular Biology</i> , 2002, 320, 1109-1117.	4.2	43
16	Secondary structure, orientation, and oligomerization of phospholemman, a cardiac transmembrane protein. <i>Protein Science</i> , 2006, 15, 1127-1132.	7.6	43
17	Structure of the Influenza C Virus CM2 Protein Transmembrane Domain Obtained by Site-specific Infrared Dichroism and Global Molecular Dynamics Searching. <i>Journal of Biological Chemistry</i> , 2000, 275, 4225-4229.	3.4	36
18	Mapping the Energy Surface of Transmembrane Helix-Helix Interactions. <i>Biophysical Journal</i> , 2001, 81, 2681-2692.	0.5	29

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19	Large-scale analysis of influenza A virus nucleoprotein sequence conservation reveals potential drug-target sites. <i>Virology</i> , 2014, 454-455, 40-47.	2.4	29
20	Enhancing resolution and sensitivity of ¹⁷ O solid-state NMR through combining double rotation, ¹ H decoupling and satellite modulation for biomolecular applications. <i>Chemical Physics Letters</i> , 2006, 421, 42-46.	2.6	27
21	Conformational Flexibility of the Peptide Hormone Ghrelin in Solution and Lipid Membrane Bound: A Molecular Dynamics Study. <i>Journal of Biomolecular Structure and Dynamics</i> , 2006, 23, 357-363.	3.5	25
22	The Transmembrane Domain of the Oncogenic Mutant ErbB-2 Receptor: A Structure Obtained from Site-specific Infrared Dichroism and Molecular Dynamics. <i>Journal of Molecular Biology</i> , 2006, 361, 945-953.	4.2	23
23	Methyl group effect on the proton affinity of methylated acetophenones studied by two mass spectrometric techniques. <i>Organic Mass Spectrometry</i> , 1993, 28, 1107-1110.	1.3	22
24	Systematic molecular dynamics searching in a lipid bilayer: Application to the glycoporphin A and oncogenic ErbB-2 transmembrane domains. <i>Journal of Molecular Graphics and Modelling</i> , 2006, 25, 226-233.	2.4	22
25	Solid-state ¹⁷ O NMR spectroscopy of a phospholemman transmembrane domain protein: Implications for the limits of detecting dilute ¹⁷ O sites in biomaterials. <i>Solid State Nuclear Magnetic Resonance</i> , 2008, 33, 72-75.	2.3	22
26	Evolutionary conservation of influenza A PB2 sequences reveals potential target sites for small molecule inhibitors. <i>Virology</i> , 2017, 509, 112-120.	2.4	16
27	Phospholemman Transmembrane Structure Reveals Potential Interactions with Na ⁺ /K ⁺ -ATPase. <i>Journal of Biological Chemistry</i> , 2007, 282, 32742-32748.	3.4	14
28	A cellulose-based bioassay for the colorimetric detection of pathogen DNA. <i>Analytical and Bioanalytical Chemistry</i> , 2014, 406, 7887-7898.	3.7	14
29	Genomic evidence for genes encoding leucine-rich repeat receptors linked to resistance against the eukaryotic extra- and intracellular <i>Brassica napus</i> pathogens <i>Leptosphaeria maculans</i> and <i>Plasmodiophora brassicae</i> . <i>PLoS ONE</i> , 2018, 13, e0198201.	2.5	13
30	Site-specific IR spectroscopy and molecular modelling combined towards solving transmembrane protein structure. <i>Spectroscopy</i> , 2005, 19, 1-16.	0.8	11
31	Structure and dynamics of the kinase IKK- β " A key regulator of the NF-kappa B transcription factor. <i>Journal of Structural Biology</i> , 2011, 176, 133-142.	2.8	11
32	Integrating molecular modelling methods to advance influenza A virus drug discovery. <i>Drug Discovery Today</i> , 2021, 26, 503-510.	6.4	11
33	G-quadruplex formation of FXVD1 pre-mRNA indicates the possibility of regulating expression of its protein product. <i>Archives of Biochemistry and Biophysics</i> , 2014, 560, 52-58.	3.0	8
34	Prediction of ligands to universally conserved binding sites of the influenza a virus nuclear export protein. <i>Virology</i> , 2019, 537, 97-103.	2.4	8
35	Electrostatic determinants of the ion channel control of the nicotinic acetylcholine receptor of <i>Torpedo californica</i> . <i>European Biophysics Journal</i> , 1998, 27, 618-625.	2.2	7
36	The Structure of Ghrelin. <i>Vitamins and Hormones</i> , 2007, 77, 1-12.	1.7	6

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37	L30A Mutation of Phospholemman Mimics Effects of Cardiac Glycosides in Isolated Cardiomyocytes. <i>Biochemistry</i> , 2016, 55, 6196-6204.	2.5	5
38	Evaluation of a novel virtual screening strategy using receptor decoy binding sites. <i>Journal of Negative Results in BioMedicine</i> , 2016, 15, 15.	1.4	5
39	Site-specific examination of secondary structure and orientation determination in membrane proteins: The peptidic ^{13}C / ^{18}O group as a novel infrared probe. <i>Biopolymers</i> , 2001, 59, 396-401.	2.4	5
40	Recent discoveries of influenza A drug target sites to combat virus replication. <i>Biochemical Society Transactions</i> , 2016, 44, 932-936.	3.4	4
41	Lipid Membranes for Membrane Proteins. <i>Methods in Molecular Biology</i> , 2015, 1215, 73-90.	0.9	4
42	Ion channels' analysis and simulation embedded in a flexible environment. <i>Computers & Chemistry</i> , 1997, 21, 181-184.	1.2	1
43	Lipid models for united-atom molecular dynamics simulations of protein. <i>Nature Precedings</i> , 2011, , .	0.1	1
44	Influenza A nucleoprotein binding sites for antivirals: current research and future potential. <i>Future Virology</i> , 2014, 9, 625-627.	1.8	1
45	Transmembrane Protein Models Based on High-Throughput Molecular Dynamics Simulations with Experimental Constraints. <i>Methods in Molecular Biology</i> , 2008, 443, 213-227.	0.9	1
46	Consensus virtual screening approaches to predict protein ligands. <i>Nature Precedings</i> , 2011, , .	0.1	0