Ansuman Lahiri

List of Publications by Year in descending order

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41 papers 337 citations

933447 10 h-index 17 g-index

44 all docs 44 docs citations

44 times ranked 473 citing authors

#	Article	IF	CITATIONS
1	LNA-induced dynamic stability in a therapeutic aptamer: insights from molecular dynamics simulations. Journal of Biomolecular Structure and Dynamics, 2023, 41, 2221-2230.	3.5	6
2	Inosine and its methyl derivatives: Occurrence, biogenesis, and function in RNA. Progress in Biophysics and Molecular Biology, 2022, 169-170, 21-52.	2.9	12
3	Data-informed reparameterization of modified RNA and the effect of explicit water models: application to pseudouridine and derivatives. Journal of Computer-Aided Molecular Design, 2022, 36, 205-224.	2.9	4
4	Molecular Dynamics Simulation of the Conformational Preferences of Pseudouridine Derivatives: Improving the Distribution in the Glycosidic Torsion Space. Journal of Chemical Information and Modeling, 2020, 60, 4995-5002.	5.4	5
5	Ensemble Allosteric Model for the Modified Wobble Hypothesis. Journal of Physical Chemistry Letters, 2020, 11, 6337-6343.	4.6	2
6	Comparative study of the SBP-box gene family in rice siblings. Journal of Biosciences, 2020, 45, 1.	1.1	2
7	Probing the functional conformations of an atypical proline-rich fusion peptide. Physical Chemistry Chemical Physics, 2019, 21, 20727-20742.	2.8	1
8	Structural Stability of the Anticodon Stem Loop Domains of the Unmodified Yeast and <i>Escherichia coli</i> tRNA ^{Phe} : Differing Views from Different Force Fields. ACS Omega, 2019, 4, 3029-3044.	3.5	1
9	Computational and NMR studies of RNA duplexes with an internal pseudouridine-adenosine base pair. Scientific Reports, 2019, 9, 16278.	3.3	30
10	Dynamical Features of Cognate Site Recognition in bZIP–DNA Interaction. ACS Omega, 2019, 4, 292-308.	3.5	1
11	Plant Polypeptide Hormone Systemin Prefers Polyproline II Conformation in Solution. ACS Omega, 2017, 2, 6831-6843.	3.5	O
12	Reparameterizations of the <i>I‡</i> Torsion and Lennard-Jones <i>If</i> Parameters Improve the Conformational Characteristics of Modified Uridines. Journal of Computational Chemistry, 2016, 37, 1576-1588.	3.3	12
13	Effect of Inactivating Mutations on Peptide Conformational Ensembles: The Plant Polypeptide Hormone Systemin. Journal of Chemical Information and Modeling, 2016, 56, 1267-1281.	5.4	3
14	Role of tryptophan 135 of Chandipura virus phosphoprotein P in dimerization and complex formation with leader RNA: structural aspect using time resolved anisotropy and simulation. RSC Advances, 2015, 5, 104582-104593.	3.6	0
15	Rapid communication capturing the destabilizing effect of dihydrouridine through molecular	2.4	4
	simulations. Biopolymers, 2014, 101, 985-991.		
16	Conformational Preferences of Modified Uridines: Comparison of AMBER Derived Force Fields. Journal of Chemical Information and Modeling, 2014, 54, 1129-1142.	5.4	19
16	Conformational Preferences of Modified Uridines: Comparison of AMBER Derived Force Fields. Journal	5.4 2.3	19

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19	Comparative analysis of abscisic acidâ€regulated transcriptomes in <i>Arabidopsis</i> . Plant Biology, 2011, 13, 28-35.	3.8	11
20	Genome wide gene expression regulation by HIP1 Protein Interactor, HIPPI: Prediction and validation. BMC Genomics, 2011, 12, 463.	2.8	13
21	Inclusion of chrysin in \hat{l}^2 -cyclodextrin nanocavity and its effect on antioxidant potential of chrysin: A spectroscopic and molecular modeling approach. Journal of Molecular Structure, 2010, 977, 180-188.	3.6	84
22	Arabidopsis thaliana regulatory element analyzer. Bioinformatics, 2008, 24, 2263-2264.	4.1	2
23	TRABAS: a database for transcription regulation by ABA signaling. In Silico Biology, 2008, 8, 511-6.	0.9	3
24	DYNAMICS OF LEUCINE-RICH REPEAT PROTEINS. Biophysical Reviews and Letters, 2007, 02, 207-219.	0.8	2
25	Interactions of HIPPI, a molecular partner of Huntingtin interacting protein HIP1, with the specific motif present at the putative promoter sequence of the caspaseâ€1, caspaseâ€8 and caspaseâ€10 genes. FEBS Journal, 2007, 274, 3886-3899.	4.7	10
26	Molecular dynamics simulation of the preferred conformations of 2-thiouridine in aqueous solution. Theoretical Chemistry Accounts, 2007, 117, 267-273.	1.4	9
27	Theoretical Analysis of the Excited State Properties of Wybutine: A Natural Probe for Transfer RNA Dynamics. International Journal of Molecular Sciences, 2004, 5, 75-83.	4.1	6
28	Exploring the idea of self-guided dynamics. Journal of Chemical Physics, 2001, 114, 5993-5999.	3.0	9
29	Molecular Dynamics of the Anticodon Domain of Yeast tRNAPhe:Codon-Anticodon Interaction. Biophysical Journal, 2000, 79, 2276-2289.	0.5	25
30	Examining the characteristics of chaos in biomolecular dynamics: a random matrix approximation. Chemical Physics Letters, 1999, 311, 459-466.	2.6	7
31	Denaturation of supercoiled DNA: a Monte Carlo study. Biophysical Chemistry, 1998, 75, 177-186.	2.8	7
32	Properties of dianionic oxyphosphorane intermediates from hybrid QM/MM simulation: implications for ribozyme reactions. Computational and Theoretical Chemistry, 1997, 419, 51-55.	1.5	5
33	Ligand binding isotherm for DNA in the presence of supercoil-induced non-B form: a theoretical analysis. Biophysical Chemistry, 1996, 58, 239-243.	2.8	2
34	Structure and energetics of plectonemically supercoiled DNA. Biopolymers, 1994, 34, 799-804.	2.4	3
35	Computational approach to the study of supercoil-induced structural polymorphism in DNA. Computational and Theoretical Chemistry, 1993, 286, 211-218.	1.5	1
36	Theoretical analysis of gel electrophoretic data for interaction of lysine rich histone with-supercoiled DNA. Biophysical Chemistry, 1992, 42, 223-228.	2.8	0

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37	Melting characteristics of highly supercoiled DNA. Biophysical Chemistry, 1992, 42, 229-234.	2.8	5
38	A semiempirical expression for the gel electrophoretic mobility of supercoiled DNA. Biopolymers, 1992, 32, 893-896.	2.4	0
39	Influencing the B-Z switch in supercoiled DNA. Biophysical Chemistry, 1991, 39, 85-90.	2.8	1
40	Effect of supercoiling on the melting characteristics of heteropolynucleotides. Biophysical Chemistry, 1991, 40, 33-41.	2.8	5
41	Theory of a supercoil-induced B-Z transition in closed circular DNA. Computational and Theoretical Chemistry, 1991, 230, 431-435.	1.5	2