

Alexey Savelyev

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7974680/publications.pdf>

Version: 2024-02-01

20
papers

1,284
citations

516710

16
h-index

839539

18
g-index

20
all docs

20
docs citations

20
times ranked

1170
citing authors

#	ARTICLE	IF	CITATIONS
1	Assessment of the DNA partial specific volume and hydration layer properties from CHARMM Drude polarizable and additive MD simulations. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 10524-10535.	2.8	4
2	Moving analytical ultracentrifugation software to a good manufacturing practices (GMP) environment. <i>PLoS Computational Biology</i> , 2020, 16, e1007942.	3.2	20
3	Moving analytical ultracentrifugation software to a good manufacturing practices (GMP) environment. , 2020, 16, e1007942.		0
4	Moving analytical ultracentrifugation software to a good manufacturing practices (GMP) environment. , 2020, 16, e1007942.		0
5	Competition among Li ⁺ , Na ⁺ , K ⁺ , and Rb ⁺ Monovalent Ions for DNA in Molecular Dynamics Simulations Using the Additive CHARMM36 and Drude Polarizable Force Fields. <i>Journal of Physical Chemistry B</i> , 2015, 119, 4428-4440.	2.6	80
6	Differential Deformability of the DNA Minor Groove and Altered BI/BII Backbone Conformational Equilibrium by the Monovalent Ions Li ⁺ , Na ⁺ , K ⁺ , and Rb ⁺ via Water-Mediated Hydrogen Bonding. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 4473-4485.	5.3	26
7	Differential Impact of the Monovalent Ions Li ⁺ , Na ⁺ , K ⁺ , and Rb ⁺ on DNA Conformational Properties. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 212-216.	4.6	51
8	All-atom polarizable force field for DNA based on the classical drude oscillator model. <i>Journal of Computational Chemistry</i> , 2014, 35, 1219-1239.	3.3	136
9	Induced Polarization Influences the Fundamental Forces in DNA Base Flipping. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 2077-2083.	4.6	59
10	Balancing the Interactions of Ions, Water, and DNA in the Drude Polarizable Force Field. <i>Journal of Physical Chemistry B</i> , 2014, 118, 6742-6757.	2.6	74
11	Recent successes in coarse-grained modeling of DNA. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2013, 3, 69-83.	14.6	81
12	Do monovalent mobile ions affect DNA's flexibility at high salt content?. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 2250.	2.8	68
13	Is DNA's Rigidity Dominated by Electrostatic or Nonelectrostatic Interactions?. <i>Journal of the American Chemical Society</i> , 2011, 133, 19290-19293.	13.7	63
14	Chemically accurate coarse graining of double-stranded DNA. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010, 107, 20340-20345.	7.1	150
15	Counterion Atmosphere and Hydration Patterns near a Nucleosome Core Particle. <i>Journal of the American Chemical Society</i> , 2009, 131, 15005-15013.	13.7	90
16	Molecular Renormalization Group Coarse-Graining of Polymer Chains: Application to Double-Stranded DNA. <i>Biophysical Journal</i> , 2009, 96, 4044-4052.	0.5	80
17	Molecular Renormalization Group Coarse-Graining of Electrolyte Solutions: Application to Aqueous NaCl and KCl. <i>Journal of Physical Chemistry B</i> , 2009, 113, 7785-7793.	2.6	76
18	Polyionic Charge Density Plays a Key Role in Differential Recognition of Mobile Ions by Biopolymers. <i>Journal of Physical Chemistry B</i> , 2008, 112, 9135-9145.	2.6	26

#	ARTICLE	IF	CITATIONS
19	Inter-DNA Electrostatics from Explicit Solvent Molecular Dynamics Simulations. Journal of the American Chemical Society, 2007, 129, 6060-6061.	13.7	62
20	Electrostatic, Steric, and Hydration Interactions Favor Na ⁺ Condensation around DNA Compared with K ⁺ . Journal of the American Chemical Society, 2006, 128, 14506-14518.	13.7	138