

# Cassandra D M Churchill

## List of Publications by Year in descending order

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20  
papers

439  
citations

759055

12  
h-index

752573

20  
g-index

20  
all docs

20  
docs citations

20  
times ranked

680  
citing authors

#	ARTICLE	IF	CITATIONS
1	Noncovalent Interactions Involving Histidine: The Effect of Charge on $\pi$ - $\pi$ Stacking and T-Shaped Interactions with the DNA Nucleobases. <i>Journal of Physical Chemistry B</i> , 2009, 113, 16046-16058.	1.2	80
2	Effects of the biological backbone on stacking interactions at DNA-protein interfaces: the interplay between the backbone $\pi$ and $\sigma$ components. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 14515.	1.3	43
3	Developing a computational model that accurately reproduces the structural features of a dinucleoside monophosphate unit within B-DNA. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 16373.	1.3	43
4	A Preliminary Investigation of the Additivity of $\pi$ - $\pi$ or $\pi$ - $\sigma$ Stacking and T-Shaped Interactions between Natural or Damaged DNA Nucleobases and Histidine. <i>Journal of Physical Chemistry B</i> , 2010, 114, 3355-3367.	1.2	33
5	The Unique Binding Mode of Laulimalide to Two Tubulin Protofilaments. <i>Chemical Biology and Drug Design</i> , 2015, 86, 190-199.	1.5	27
6	Effects of the biological backbone on DNA-protein stacking interactions. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 10657.	1.3	25
7	Evaluating How Discrete Water Molecules Affect Protein-DNA $\pi$ - $\pi$ and $\pi$ - $\sigma$ Stacking and T-Shaped Interactions: The Case of Histidine-Adenine Dimers. <i>Journal of Physical Chemistry B</i> , 2011, 115, 10990-11003.	1.2	24
8	Elucidating the Mechanism of Action of the Clinically Approved Taxanes: A Comprehensive Comparison of Local and Allosteric Effects. <i>Chemical Biology and Drug Design</i> , 2015, 86, 1253-1266.	1.5	22
9	A new antiproliferative noscapine analogue: chemical synthesis and biological evaluation. <i>Oncotarget</i> , 2016, 7, 40518-40530.	0.8	21
10	Effect of Watson-Crick and Hoogsteen Base Pairing on the Conformational Stability of C8-Phenoxy-2-deoxyguanosine Adducts. <i>Journal of Physical Chemistry B</i> , 2010, 114, 12995-13004.	1.2	17
11	Analysis of the binding mode of laulimalide to microtubules: Establishing a laulimalide-tubulin pharmacophore. <i>Journal of Biomolecular Structure and Dynamics</i> , 2016, 34, 1455-1469.	2.0	16
12	Formation Mechanism and Structure of a Guanine-Uracil DNA Intrastrand Cross-Link. <i>Chemical Research in Toxicology</i> , 2011, 24, 2189-2199.	1.7	15
13	Mathematical and computational modeling in biology at multiple scales. <i>Theoretical Biology and Medical Modelling</i> , 2014, 11, 52.	2.1	12
14	Inactivation of Protein Tyrosine Phosphatases by Peracids Correlates with the Hydrocarbon Chain Length. <i>Cellular Physiology and Biochemistry</i> , 2015, 36, 1069-1083.	1.1	12
15	Probing the Basis of $\alpha$ -Synuclein Aggregation by Comparing Simulations to Single-Molecule Experiments. <i>Biophysical Journal</i> , 2019, 117, 1125-1135.	0.2	11
16	Designing and Testing of Novel Taxanes to Probe the Highly Complex Mechanisms by Which Taxanes Bind to Microtubules and Cause Cytotoxicity to Cancer Cells. <i>PLoS ONE</i> , 2015, 10, e0129168.	1.1	11
17	DNA Distortion Caused by Uracil-Containing Intrastrand Cross-Links. <i>Journal of Physical Chemistry B</i> , 2016, 120, 1195-1204.	1.2	10
18	Synthesis and biological evaluation of structurally simplified noscapine analogues as microtubule binding agents. <i>Canadian Journal of Chemistry</i> , 2017, 95, 649-655.	0.6	10

#	ARTICLE	IF	CITATIONS
19	The fragment molecular orbital method and understanding monomer polarization. Chemical Physics Letters, 2012, 554, 185-189.	1.2	5
20	Molecular Dynamics and Related Computational Methods with Applications to Drug Discovery. Springer Proceedings in Mathematics and Statistics, 2018, , 267-285.	0.1	2