

# Rhys Heffernan

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7155417/publications.pdf>

Version: 2024-02-01

12  
papers

1,565  
citations

759055

12  
h-index

1199470

12  
g-index

12  
all docs

12  
docs citations

12  
times ranked

1628  
citing authors

#	ARTICLE	IF	CITATIONS
1	Capturing non-local interactions by long short-term memory bidirectional recurrent neural networks for improving prediction of protein secondary structure, backbone angles, contact numbers and solvent accessibility. <i>Bioinformatics</i> , 2017, 33, 2842-2849.	1.8	300
2	Improving prediction of secondary structure, local backbone angles and solvent accessible surface area of proteins by iterative deep learning. <i>Scientific Reports</i> , 2015, 5, 11476.	1.6	290
3	Gram-positive and Gram-negative protein subcellular localization by incorporating evolutionary-based descriptors into Chou's general PseAAC. <i>Journal of Theoretical Biology</i> , 2015, 364, 284-294.	0.8	232
4	Sixty-five years of the long march in protein secondary structure prediction: the final stretch?. <i>Briefings in Bioinformatics</i> , 2018, 19, bbw129.	3.2	168
5	SPIDER2: A Package to Predict Secondary Structure, Accessible Surface Area, and Main-Chain Torsional Angles by Deep Neural Networks. <i>Methods in Molecular Biology</i> , 2017, 1484, 55-63.	0.4	137
6	Predicting backbone $\phi/\psi$ angles and dihedrals from protein sequences by stacked sparse auto-encoder deep neural network. <i>Journal of Computational Chemistry</i> , 2014, 35, 2040-2046.	1.5	133
7	Single-sequence-based prediction of protein secondary structures and solvent accessibility by deep whole-sequence learning. <i>Journal of Computational Chemistry</i> , 2018, 39, 2210-2216.	1.5	84
8	Highly accurate sequence-based prediction of half-sphere exposures of amino acid residues in proteins. <i>Bioinformatics</i> , 2016, 32, 843-849.	1.8	79
9	SPIN2: Predicting sequence profiles from protein structures using deep neural networks. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018, 86, 629-633.	1.5	62
10	Advancing the Accuracy of Protein Fold Recognition by Utilizing Profiles From Hidden Markov Models. <i>IEEE Transactions on Nanobioscience</i> , 2015, 14, 761-772.	2.2	34
11	Protein fold recognition using HMM-HMM alignment and dynamic programming. <i>Journal of Theoretical Biology</i> , 2016, 393, 67-74.	0.8	33
12	Detecting Proline and Non-Proline Cis Isomers in Protein Structures from Sequences Using Deep Residual Ensemble Learning. <i>Journal of Chemical Information and Modeling</i> , 2018, 58, 2033-2042.	2.5	13