

# Florian Richter

## List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

18  
papers

3,480  
citations

15  
h-index

18  
g-index

18  
ext. papers

4,198  
ext. citations

9.7  
avg, IF

4.38  
L-index

#	Paper	IF	Citations
18	A Light-Oxygen-Voltage Receptor Integrates Light and Temperature. <i>Journal of Molecular Biology</i> , <b>2021</b> , 433, 167107	6.5	5
17	Computational Aminoacyl-tRNA Synthetase Library Design for Photocaged Tyrosine. <i>International Journal of Molecular Sciences</i> , <b>2019</b> , 20,	6.3	16
16	Optogenetic Control by Pulsed Illumination. <i>ChemBioChem</i> , <b>2018</b> , 19, 1296-1304	3.8	15
15	Switchable Cas9. <i>Current Opinion in Biotechnology</i> , <b>2017</b> , 48, 119-126	11.4	32
14	Photoactivatable Mussel-Based Underwater Adhesive Proteins by an Expanded Genetic Code. <i>ChemBioChem</i> , <b>2017</b> , 18, 1819-1823	3.8	39
13	Engineering of temperature- and light-switchable Cas9 variants. <i>Nucleic Acids Research</i> , <b>2016</b> , 44, 10003-10014	10.1470	70
12	Library-Aided Probing of Linker Determinants in Hybrid Photoreceptors. <i>ACS Synthetic Biology</i> , <b>2016</b> , 5, 1117-1126	5.7	32
11	Upgrading a microplate reader for photobiology and all-optical experiments. <i>Photochemical and Photobiological Sciences</i> , <b>2015</b> , 14, 270-9	4.2	27
10	Light-Induced Rearrangement of the $\beta$ Strand in the BLUF Photoreceptor SyPixD (Slr1694). <i>Journal of Physical Chemistry Letters</i> , <b>2015</b> , 6, 4749-53	6.4	11
9	Computational design of a red fluorophore ligase for site-specific protein labeling in living cells. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2014</b> , 111, E4551-9	11.5	57
8	Design of activated serine-containing catalytic triads with atomic-level accuracy. <i>Nature Chemical Biology</i> , <b>2014</b> , 10, 386-91	11.7	52
7	Computational Protein Design for Synthetic Biology <b>2013</b> , 101-122		3
6	Computational design of catalytic dyads and oxyanion holes for ester hydrolysis. <i>Journal of the American Chemical Society</i> , <b>2012</b> , 134, 16197-206	16.4	114
5	ROSETTA3: an object-oriented software suite for the simulation and design of macromolecules. <i>Methods in Enzymology</i> , <b>2011</b> , 487, 545-74	1.7	1216
4	De novo enzyme design using Rosetta3. <i>PLoS ONE</i> , <b>2011</b> , 6, e19230	3.7	224
3	RosettaScripts: a scripting language interface to the Rosetta macromolecular modeling suite. <i>PLoS ONE</i> , <b>2011</b> , 6, e20161	3.7	311
2	RosettaRemodel: a generalized framework for flexible backbone protein design. <i>PLoS ONE</i> , <b>2011</b> , 6, e24169	3.7	200

- 1 Quantitative reactivity profiling predicts functional cysteines in proteomes. *Nature*, **2010**, 468, 790-5 50.4 1056