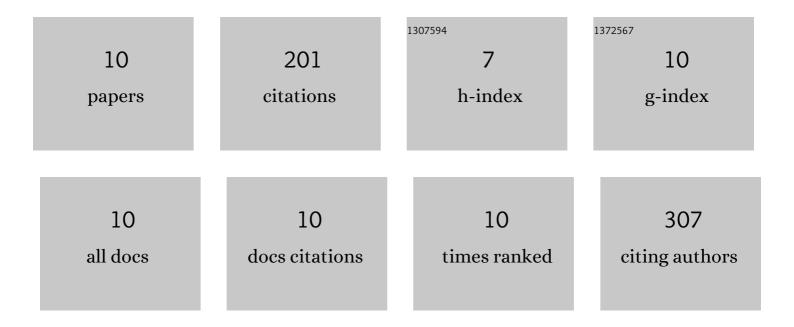
## Chang Woo Kim

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Theory of dissipation pathways in open quantum systems. Journal of Chemical Physics, 2021, 154, 084109.	3.0	7
2	Uncovering the Conformational Distribution of a Small Protein with Nanoparticle-Aided Cryo-Electron Microscopy Sampling. Journal of Physical Chemistry Letters, 2021, 12, 6565-6573.	4.6	4
3	Protein folding from heterogeneous unfolded state revealed by time-resolved X-ray solution scattering. Proceedings of the National Academy of Sciences of the United States of America, 2020, 117, 14996-15005.	7.1	33
4	Toward monitoring the dissipative vibrational energy flows in open quantum systems by mixed quantum–classical simulations. Journal of Chemical Physics, 2020, 152, 244109.	3.0	5
5	Effect of Underdamped Vibration on Excitation Energy Transfer: Direct Comparison between Two Different Partitioning Schemes. Journal of Physical Chemistry A, 2019, 123, 1186-1197.	2.5	14
6	Excited state energy fluctuations in the Fenna–Matthews–Olson complex from molecular dynamics simulations with interpolated chromophore potentials. Physical Chemistry Chemical Physics, 2018, 20, 3310-3319.	2.8	31
7	Constructing an Interpolated Potential Energy Surface of a Large Molecule: A Case Study with Bacteriochlorophyll <i>a</i> Model in the Fenna–Matthews–Olson Complex. Journal of Chemical Theory and Computation, 2016, 12, 5235-5246.	5.3	23
8	Effect of Chromophore Potential Model on the Description of Exciton–Phonon Interactions. Journal of Physical Chemistry Letters, 2015, 6, 2875-2880.	4.6	27
9	Phase Behavior and Conductivity of Sulfonated Block Copolymers Containing Heterocyclic Diazole-Based Ionic Liquids. Macromolecules, 2012, 45, 8702-8713.	4.8	46
10	Comparison of Arsenic Acid with Phosphoric Acid in the Interaction with a Water Molecule and an Alkali/Alkaline-Earth Metal Cation. Journal of Physical Chemistry A, 2011, 115, 11355-11361.	2.5	11