

# Chang Woo Kim

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/101369/publications.pdf>

Version: 2024-02-01

10  
papers

201  
citations

1307594

7  
h-index

1372567

10  
g-index

10  
all docs

10  
docs citations

10  
times ranked

307  
citing authors

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