

# Cuihong Wang

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8255386/publications.pdf>

Version: 2024-02-01

8  
papers

70  
citations

1937685

4  
h-index

1872680

6  
g-index

8  
all docs

8  
docs citations

8  
times ranked

117  
citing authors

#	ARTICLE	IF	CITATIONS
1	Adsorption and properties of aromatic amino acids on single-walled carbon nanotubes. <i>Nanoscale</i> , 2012, 4, 1146-1153.	5.6	45
2	Conformers, properties, and docking mechanism of the anticancer drug docetaxel: DFT and molecular dynamics studies. <i>Journal of Computational Chemistry</i> , 2018, 39, 889-900.	3.3	10
3	Binding modes of cabazitaxel with the different human $\beta$ -tubulin isotypes: DFT and MD studies. <i>Journal of Molecular Modeling</i> , 2020, 26, 162.	1.8	6
4	A DFT study on the high-density assembly of doxorubicin drug delivery by single-walled carbon nanotubes. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2021, 134, 114892.	2.7	6
5	Design and interaction mechanism of ligand targeted with folate receptor $\beta$ 1 and $\beta$ 2. <i>Journal of Physical Organic Chemistry</i> , 2018, 31, e3719.	1.9	2
6	Endohedral and exohedral complexes of 1,2-benzene with carbon nanotubes and high-density assembly of multiple benzenes inside of a carbon nanotube. <i>International Journal of Quantum Chemistry</i> , 2019, 119, e25936.	2.0	1
7	Conformers, Properties of the Anticancer Drug Plocabulin, and its Binding Mechanism with p-Glycoprotein: DFT and MD Studies. <i>Australian Journal of Chemistry</i> , 2021, 74, 529.	0.9	0
8	Interaction mechanism of novel fluorescent antifolates targeted with folate receptors $\beta$ 1 and $\beta$ 2 via molecular docking and molecular dynamic simulations. <i>Journal of Molecular Modeling</i> , 2022, 28, .	1.8	0