

# Chunyi Zhang

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9166710/publications.pdf>

Version: 2024-02-01

11  
papers

190  
citations

1307594

7  
h-index

1372567

10  
g-index

11  
all docs

11  
docs citations

11  
times ranked

111  
citing authors

| #  | ARTICLE  | IF   | CITATIONS |
|----|--|------|-----------|
| 1  | Finite-temperature phonon dispersion and vibrational dynamics of BaTiO <sub>3</sub> from first-principles molecular dynamics. <i>Physical Review B</i> , 2022, 105, .                                  | 3.2  | 3         |
| 2  | Dissolving salt is not equivalent to applying a pressure on water. <i>Nature Communications</i> , 2022, 13, 822.   | 12.8 | 41        |
| 3  | Many-body effects in the X-ray absorption spectra of liquid water. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022, 119, e2201258119.                    | 7.1  | 11        |
| 4  | Importance of nuclear quantum effects on the hydration of chloride ion. <i>Physical Review Materials</i> , 2021, 5, .  | 2.4  | 11        |
| 5  | Modeling Liquid Water by Climbing up Jacob's Ladder in Density Functional Theory Facilitated by Using Deep Neural Network Potentials. <i>Journal of Physical Chemistry B</i> , 2021, 125, 11444-11456. | 2.6  | 40        |
| 6  | Isotope effects in x-ray absorption spectra of liquid water. <i>Physical Review B</i> , 2020, 102, .   | 3.2  | 6         |
| 7  | Isotope effects in molecular structures and electronic properties of liquid water via deep potential molecular dynamics based on the SCAN functional. <i>Physical Review B</i> , 2020, 102, .          | 3.2  | 22        |
| 8  | Nano-architected metamaterials: Carbon nanotube-based nanotrusses. <i>Carbon</i> , 2018, 131, 38-46.   | 10.3 | 29        |
| 9  | Finite-temperature infrared and Raman spectra of high-pressure hydrogen from first-principles molecular dynamics. <i>Physical Review B</i> , 2018, 98, .   | 3.2  | 14        |
| 10 | Thermal conductance of one-dimensional materials calculated with typical lattice models. <i>Physical Review E</i> , 2016, 94, 052131.  | 2.1  | 5         |
| 11 | Structural and Dynamic Properties of Solvated Hydroxide and Hydronium Ions in Water from Ab Initio Modeling. <i>Journal of Chemical Physics</i> , 0, , .   | 3.0  | 8         |