

Yusuke Kanematsu

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	PyDISH: database and analysis tools for heme porphyrin distortion in heme proteins. Database: the Journal of Biological Databases and Curation, 2023, 2023, .	3.0	12
2	Hydrogen/Deuterium Transfer from Anisole to Methoxy Radicals: A Theoretical Study of a Deuterium-Labeled Drug Model. Journal of Physical Chemistry A, 2022, 126, 155-163.	2.5	3
3	Global Analysis of Heme Proteins Elucidates the Correlation between Heme Distortion and the Heme-Binding Pocket. Journal of Chemical Information and Modeling, 2022, 62, 775-784.	5.4	10
4	Structure of Heme-binding Pocket in Heme Protein is Generally Rigid and can be Predicted by AlphaFold2. Chemistry Letters, 2022, 51, 704-708.	1.3	8
5	Short-lived intermediate in N ₂ O generation by P450 NO reductase captured by time-resolved IR spectroscopy and XFEL crystallography. Proceedings of the National Academy of Sciences of the United States of America, 2021, 118, .	7.1	21
6	H/D isotope effect between adsorbed water (H ₂ O, D ₂ O, and HDO) and H ₂ O- and D ₂ O-ice Ih(0001) basal surfaces based on the combined plane wave and localized basis set method. Applied Surface Science, 2021, 561, 150100.	6.1	1
7	A computational examination of the electric-field-induced proton transfer along the interface hydrogen bond between proton donating and accepting self-assembled monolayers. Chemical Physics Letters, 2020, 741, 137091.	2.6	5
8	Computational study of distortion effect of Fe-porphyrin found as a biological active site. Japanese Journal of Applied Physics, 2020, 59, 010502.	1.5	14
9	Strong Hydrogen Bonds at the Interface between Proton-Donating and -Accepting Self-Assembled Monolayers on Au(111). Langmuir, 2018, 34, 2189-2197.	3.5	16
10	Statistical and quantum-chemical analysis of the effect of heme porphyrin distortion in heme proteins: Differences between oxidoreductases and oxygen carrier proteins. Chemical Physics Letters, 2018, 710, 108-112.	2.6	13
11	Applicability of Density Functional Tight Binding Method with Dispersion Correction to Investigate the Adsorption of Porphyrin/Porphycene Metal Complexes on Graphene. Chemistry Letters, 2017, 46, 51-52.	1.3	8
12	Multicomponent DFT study of geometrical H/D isotope effect on hydrogen-bonded organic conductor, $\text{I}^{\text{9}}\text{-H}_3$ (Cat EDT-ST) 2. Chemical Physics Letters, 2017, 674, 168-172.	2.6	12
13	Capturing an initial intermediate during the P450 _{nor} enzymatic reaction using time-resolved XFEL crystallography and caged-substrate. Nature Communications, 2017, 8, 1585.	12.8	74
14	Inverse Ubbelohde effect in the short hydrogen bond of photosystem II: Relation between H/D isotope effect and symmetry in potential energy profile. Journal of Computational Chemistry, 2016, 37, 2140-2145.	3.3	6
15	Quantum Mechanics/Molecular Mechanics Study of the Sialyltransferase Reaction Mechanism. Biochemistry, 2016, 55, 5764-5771.	2.5	4
16	Theoretical study of the H/D isotope effect on phase transition of hydrogen-bonded organic conductor $\text{I}^{\text{9}}\text{-H}_3$ (Cat-EDT-TTF) 2. Physical Chemistry Chemical Physics, 2016, 18, 29673-29680.	2.8	26
17	Vibrational analysis on the revised potential energy curve of the low-barrier hydrogen bond in photoactive yellow protein. Computational and Structural Biotechnology Journal, 2016, 14, 16-19.	4.1	6
18	Isotope effect on the circular dichroism spectrum of methyl $\text{I}^{\text{9}}\text{-D}$ -glucopyranoside in aqueous solution. Scientific Reports, 2016, 5, 17900.	3.3	9

#	ARTICLE	IF	CITATIONS
19	Performance Test of Multicomponent Quantum Mechanical Calculation with Polarizable Continuum Model for Proton Chemical Shift. <i>Journal of Physical Chemistry A</i> , 2015, 119, 4933-4938.	2.5	11
20	Theoretical analysis of geometry and NMR isotope shift in hydrogen-bonding center of photoactive yellow protein by combination of multicomponent quantum mechanics and ONIOM scheme. <i>Journal of Chemical Physics</i> , 2014, 141, 185101.	3.0	22
21	Development of multicomponent hybrid density functional theory with polarizable continuum model for the analysis of nuclear quantum effect and solvent effect on NMR chemical shift. <i>Journal of Chemical Physics</i> , 2014, 140, 164111.	3.0	16
22	Substrate shielding and hydrolytic reaction in hydrolases. <i>Proteins: Structure, Function and Bioinformatics</i> , 2013, 81, 926-932.	2.6	1
23	Nuclear quantum effect and H/D isotope effect in excited state intramolecular proton transfer and electron-induced intramolecular proton transfer reactions in 8-hydroxyquinoline. <i>International Journal of Quantum Chemistry</i> , 0, , .	2.0	1