

Keiko Takano

List of Publications by Year in descending order

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Version: 2024-02-01

12
papers

105
citations

1478505

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1281871

11
g-index

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all docs

12
docs citations

12
times ranked

165
citing authors

#	ARTICLE	IF	CITATIONS
1	Reversibility of 1,4-Metal Migration in Cp*Rh ^{III} and Cp*Ir ^{III} Complexes. <i>Organometallics</i> , 2014, 33, 2142-2145.	2.3	32
2	Theoretical Study on Internal Alkyne/Vinylidene Isomerization in Group 8 Transition-Metal Complexes. <i>Organometallics</i> , 2015, 34, 3934-3943.	2.3	26
3	Affinity of HIV-1 antibody 2G12 with monosaccharides: A theoretical study based on explicit and implicit water models. <i>Computational Biology and Chemistry</i> , 2014, 49, 36-44.	2.3	8
4	Ruthenium Vinylidene Complexes Generated by Selective 1,2-Migration of P- and S-Substituents: Synthesis, Structures, and Dichromism Arising from an Intramolecular CH ^{δ+} ⋯O Hydrogen Bond. <i>Organometallics</i> , 2020, 39, 711-718.	2.3	8
5	Substituent effects on the photophysical properties of 2,9-substituted phenanthroline copper(I) complexes: a theoretical investigation. <i>ChemPhysChem</i> , 2021, 22, 509-515.	2.1	7
6	Interaction analysis of HIV-1 antibody 2G12 and Man9GlcNAc2 ligand: Theoretical calculations by fragment molecular orbital and MD methods. <i>Chemical Physics Letters</i> , 2013, 578, 144-149.	2.6	6
7	Theoretical study on photophysical properties of 3-hydroxyechinenone and the effects of interactions with orange carotenoid protein. <i>Chemical Physics Letters</i> , 2016, 647, 95-102.	2.6	5
8	CHEMICAL DESCRIPTION OF THE INTERACTION BETWEEN GLYCAN LIGAND AND SIGLEC-7 USING AB INITIO FMO METHOD AND CLASSICAL MD SIMULATION. <i>Journal of Theoretical and Computational Chemistry</i> , 2013, 12, 1350060.	1.8	4
9	Water molecules inside protein structure affect binding of monosaccharides with HIV-1 antibody 2G12. <i>Journal of Computational Chemistry</i> , 2016, 37, 2341-2348.	3.3	3
10	R/X exchange reactions in cis-[M(R) ₂ {P(X)(NMeCH ₂) ₂ } ₂] (M = Pd, Pt), via a phosphonium intermediate. <i>Dalton Transactions</i> , 2016, 45, 19216-19220.	3.3	3
11	Systematic Interaction Analysis of Anti-Human Immunodeficiency Virus Type-1 Neutralizing Antibodies with High Mannose Glycans Using Fragment Molecular Orbital and Molecular Dynamics Methods. <i>Journal of Computational Chemistry</i> , 2020, 41, 31-42.	3.3	3
12	Theoretical Study on Adjacent Agostic Interaction in Ruthenium Complexes. <i>Journal of Computer Chemistry Japan</i> , 2019, 18, 162-163.	0.1	0