

Matthew D Kundrat

List of Publications by Year in descending order

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8
papers

574
citations

1307594
7
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1588992
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g-index

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all docs

8
docs citations

8
times ranked

574
citing authors

#	ARTICLE	IF	CITATIONS
1	Modeling of the Chiroptical Response of Chiral Amino Acids in Solution Using Explicit Solvation and Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 1051-1060.	5.3	46
2	Ab Initio and Density Functional Theory Modeling of the Chiroptical Response of Glycine and Alanine in Solution Using Explicit Solvation and Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2008, 4, 1902-1914.	5.3	47
3	Computational Modeling of the Optical Rotation of Amino Acids: A New Look at an Old Rule for pH Dependence of Optical Rotation. <i>Journal of the American Chemical Society</i> , 2008, 130, 4404-4414.	13.7	38
4	Time Dependent Density Functional Theory Modeling of Chiroptical Properties of Small Amino Acids in Solution. <i>Journal of Physical Chemistry A</i> , 2006, 110, 4115-4123.	2.5	47
5	Calculation of circular dichroism spectra from optical rotatory dispersion, and vice versa, as complementary tools for theoretical studies of optical activity using time-dependent density functional theory. <i>Journal of Chemical Physics</i> , 2006, 125, 194110.	3.0	57
6	Time Dependent Density Functional Theory Modeling of Specific Rotation and Optical Rotatory Dispersion of the Aromatic Amino Acids in Solution. <i>Journal of Physical Chemistry A</i> , 2006, 110, 12908-12917.	2.5	30
7	Hydrogen Desorption Exceeding Ten Weight Percent from the New Quaternary Hydride Li ₃ BN ₂ H ₈ . <i>ChemInform</i> , 2005, 36, no.	0.0	2
8	Hydrogen Desorption Exceeding Ten Weight Percent from the New Quaternary Hydride Li ₃ BN ₂ H ₈ . <i>Journal of Physical Chemistry B</i> , 2005, 109, 6-8.	2.6	307