

Omar Baig

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Assessing Configurational Sampling in the Quantum Mechanics/Molecular Mechanics Calculation of Temoporfin Absorption Spectrum and Triplet Density of States. <i>Molecules</i> , 2018, 23, 2932.	3.8	10
2	Molecular dynamics simulation of aqueous 1-dodecyl-3-methylimidazolium chloride: Emerging micelles. <i>Journal of Molecular Liquids</i> , 2018, 272, 766-777.	4.9	10
3	The effect of N-heterocyclic carbene units on the absorption spectra of Fe(II) complexes: a challenge for theory. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 27605-27616.	2.8	8