

Marek Pederzoli

List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

9
papers

403
citations

7
h-index

10
g-index

10
ext. papers

435
ext. citations

3.6
avg, IF

3.26
L-index

#	Paper	IF	Citations
9	Surface Hopping Dynamics with Correlated Single-Reference Methods: 9H-Adenine as a Case Study. <i>Journal of Chemical Theory and Computation</i> , 2014 , 10, 1395-405	6.4	139
8	Non-adiabatic dynamics of pyrrole: Dependence of deactivation mechanisms on the excitation energy. <i>Chemical Physics</i> , 2010 , 375, 26-34	2.3	115
7	Nonadiabatic molecular dynamics study of the cis-trans photoisomerization of azobenzene excited to the S1 state. <i>Journal of Physical Chemistry A</i> , 2011 , 115, 11136-43	2.8	97
6	A new approach to molecular dynamics with non-adiabatic and spin-orbit effects with applications to QM/MM simulations of thiophene and selenophene. <i>Journal of Chemical Physics</i> , 2017 , 146, 114101	3.9	19
5	Orientation of Laurdan in Phospholipid Bilayers Influences Its Fluorescence: Quantum Mechanics and Classical Molecular Dynamics Study. <i>Molecules</i> , 2018 , 23,	4.8	12
4	Fluorescence of PRODAN in water: A computational QM/MM MD study. <i>Chemical Physics Letters</i> , 2014 , 597, 57-62	2.5	11
3	Photophysics of BODIPY-Based Photosensitizer for Photodynamic Therapy: Surface Hopping and Classical Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2019 , 15, 5046-5057	6.4	8
2	Cis-transphotoisomerization of azobenzene upon excitation to the S1state: an ab initio molecular dynamics and QM/MM study 2012 ,		1
1	Theoretical Investigation of the Effect of Alkylation and Bromination on Intersystem Crossing in BODIPY-Based Photosensitizers. <i>Journal of Physical Chemistry B</i> , 2021 , 125, 11617-11627	3.4	1