

# Adrian W Lange

## List of Publications by Citations

**Source:** <https://exaly.com/author-pdf/11650033/adrian-w-lange-publications-by-citations.pdf>

**Version:** 2024-04-28

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

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

12  
papers

2,836  
citations

9  
h-index

12  
g-index

12  
ext. papers

3,329  
ext. citations

4.9  
avg, IF

4.26  
L-index

#	Paper	IF	Citations
12	Advances in molecular quantum chemistry contained in the Q-Chem 4 program package. <i>Molecular Physics</i> , <b>2015</b> , 113, 184-215	1.7	2068
11	Both intra- and interstrand charge-transfer excited states in aqueous B-DNA are present at energies comparable to, or just above, the (1)ππ* excitonic bright states. <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 3913-22	16.4	168
10	Charge-transfer excited states in a π-stacked adenine dimer, as predicted using long-range-corrected time-dependent density functional theory. <i>Journal of Physical Chemistry B</i> , <b>2008</b> , 112, 6304-8	3.4	150
9	A smooth, nonsingular, and faithful discretization scheme for polarizable continuum models: the switching/Gaussian approach. <i>Journal of Chemical Physics</i> , <b>2010</b> , 133, 244111	3.9	137
8	Software for the frontiers of quantum chemistry: An overview of developments in the Q-Chem 5 package. <i>Journal of Chemical Physics</i> , <b>2021</b> , 155, 084801	3.9	115
7	Polarizable Continuum Reaction-Field Solvation Models Affording Smooth Potential Energy Surfaces. <i>Journal of Physical Chemistry Letters</i> , <b>2010</b> , 1, 556-561	6.4	89
6	Symmetric versus asymmetric discretization of the integral equations in polarizable continuum solvation models. <i>Chemical Physics Letters</i> , <b>2011</b> , 509, 77-87	2.5	58
5	A simple polarizable continuum solvation model for electrolyte solutions. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 204110	3.9	16
4	Improving Generalized Born Models by Exploiting Connections to Polarizable Continuum Models. I. An Improved Effective Coulomb Operator. <i>Journal of Chemical Theory and Computation</i> , <b>2012</b> , 8, 1999-2014	6.4	13
3	Intrinsically smooth discretisation of Connolly's solvent-excluded molecular surface. <i>Molecular Physics</i> , <b>2020</b> , 118, e1644384	1.7	9
2	Improving Generalized Born Models by Exploiting Connections to Polarizable Continuum Models. II. Corrections for Salt Effects. <i>Journal of Chemical Theory and Computation</i> , <b>2012</b> , 8, 4381-92	6.4	8
1	Response to [Comment on A smooth, nonsingular, and faithful discretization scheme for polarizable continuum models: The switching/Gaussian approach[J. Chem. Phys. 134, 117101 (2011)]. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 117102	3.9	5