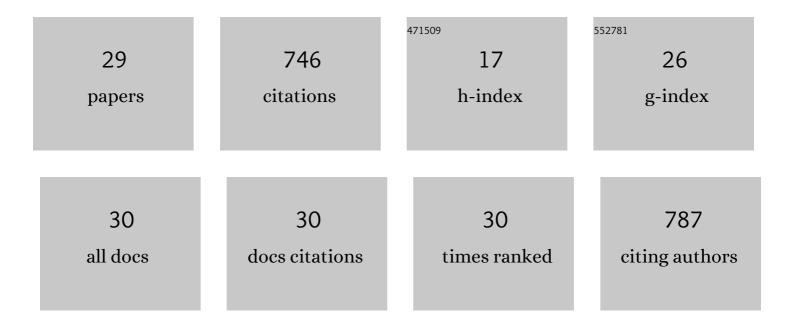
Dorit Shemesh

List of Publications by Year in descending order

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| # | Article | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Absorption spectra of pyruvic acid in water: insights from calculations for small hydrates and comparison to experiment. Physical Chemistry Chemical Physics, 2020, 22, 12658-12670. | 2.8 | 19 |
| 2 | Impact of pH and NaCl and CaCl ₂ Salts on the Speciation and Photochemistry of Pyruvic Acid in the Aqueous Phase. Journal of Physical Chemistry A, 2020, 124, 5071-5080. | 2.5 | 18 |
| 3 | Experimental and Theoretical Studies of the Environmental Sensitivity of the Absorption Spectra and Photochemistry of Nitenpyram and Analogs. ACS Earth and Space Chemistry, 2019, 3, 2063-2075. | 2.7 | 8 |
| 4 | Molecular Dynamics of Photoinduced Reactions of Acrylic Acid: Products, Mechanisms, and Comparison with Experiment. Journal of Physical Chemistry Letters, 2018, 9, 527-533. | 4.6 | 15 |
| 5 | Adjacent keto and enol groups in photochemistry of a cyclic molecule: Products, mechanisms and dynamics. Chemical Physics, 2018, 515, 177-186. | 1.9 | 3 |
| 6 | Photochemistry of Thin Solid Films of the Neonicotinoid Imidacloprid on Surfaces. Environmental Science & Technology, 2017, 51, 2660-2668. | 10.0 | 37 |
| 7 | DMAP-assisted sulfonylation as an efficient step for the methylation of primary amine motifs on solid support. Beilstein Journal of Organic Chemistry, 2017, 13, 806-816. | 2.2 | 9 |
| 8 | Dynamics of Photochemical Reactions of Organic Carbonyls and their Clusters. Advances in Chemical Physics, 2016, , 1-22. | 0.3 | 0 |
| 9 | Photochemical Reactions of Cyclohexanone: Mechanisms and Dynamics. Journal of Physical Chemistry A, 2016, 120, 7112-7120. | 2.5 | 17 |
| 10 | Absorption spectra and aqueous photochemistry of β-hydroxyalkyl nitrates of atmospheric interest. Molecular Physics, 2015, 113, 2179-2190. | 1.7 | 22 |
| 11 | Computational Studies of Atmospherically-Relevant Chemical Reactions in Water Clusters and on Liquid Water and Ice Surfaces. Accounts of Chemical Research, 2015, 48, 399-406. | 15.6 | 89 |
| 12 | Photochemistry of aldehyde clusters: cross-molecular versus unimolecular reaction dynamics. Physical Chemistry Chemical Physics, 2014, 16, 23861-23868. | 2.8 | 21 |
| 13 | Ab initio and semi-empirical Molecular Dynamics simulations of chemical reactions in isolated molecules and in clusters. Physical Chemistry Chemical Physics, 2014, 16, 9760-9775. | 2.8 | 35 |
| 14 | Dynamics of Triplet-State Photochemistry of Pentanal: Mechanisms of Norrish I, Norrish II, and H Abstraction Reactions. Journal of Physical Chemistry A, 2013, 117, 11711-11724. | 2.5 | 26 |
| 15 | Experimental and Theoretical Study of Aqueous <i>cis</i> -Pinonic Acid Photolysis. Journal of Physical Chemistry A, 2013, 117, 12930-12945. | 2.5 | 60 |
| 16 | Femtosecond timescale deactivation of electronically excited peroxides at ice surfaces. Molecular Physics, 2012, 110, 605-617. | 1.7 | 21 |
| 17 | Absorption Spectra and Photolysis of Methyl Peroxide in Liquid and Frozen Water. Journal of Physical Chemistry A, 2012, 116, 6068-6077. | 2.5 | 49 |
| 18 | Effect of the Chirality of Residues and γâ€Turns on the Electronic Excitation Spectra, Excitedâ€6tate Reaction Paths and Conical Intersections of Capped Phenylalanine–Alanine Dipeptides. ChemPhysChem, 2011, 12, 1833-1840. | 2.1 | 14 |

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| # | Article | IF | CITATIONS |
|----|---|------|-----------|
| 19 | Role of excited-state hydrogen detachment and hydrogen-transfer processes for the excited-state deactivation of an aromatic dipeptide: N-acetyl tryptophan methyl amide. Physical Chemistry Chemical Physics, 2010, 12, 4899. | 2.8 | 29 |
| 20 | Photophysics of the Trp-Gly Dipeptide: Role ofÂElectron and Proton Transfer Processes forÂEfficient Excited-State Deactivation. , 2010, , 641-649. | | 0 |
| 21 | Photophysics of the Trp-Gly dipeptide: Role of electron and proton transfer processes for efficient excited-state deactivation. Chemical Physics Letters, 2009, 482, 38-43. | 2.6 | 27 |
| 22 | Efficient Excited-State Deactivation of the Gly-Phe-Ala Tripeptide via an Electron-Driven Proton-Transfer Process. Journal of the American Chemical Society, 2009, 131, 1374-1375. | 13.7 | 65 |
| 23 | Computational Studies of the Photophysics of Neutral and Zwitterionic Amino Acids in an Aqueous Environment: Tyrosineâ^'(H2O)2 and Tryptophanâ^'(H2O)2 Clusters. Journal of Physical Chemistry A, 2009, 113, 542-550. | 2.5 | 57 |
| 24 | Vibrational spectroscopy for glycine adsorbed on silicon clusters: Harmonic and anharmonic calculations for models of the Si(100)-2×1 surface. Chemical Physics, 2008, 347, 218-228. | 1.9 | 26 |
| 25 | Dynamical Simulations of Photoionization of Small Biological Molecules. , 2006, , 213-237. | | 0 |
| 26 | Classical Trajectory Simulations of Photoionization Dynamics of Tryptophan:Â Intramolecular Energy Flow, Hydrogen-Transfer Processes and Conformational Transitionsâ€. Journal of Physical Chemistry A, 2006, 110, 8401-8408. | 2.5 | 18 |
| 27 | Photoionization dynamics of glycine adsorbed on a silicon cluster: "On-the-fly―simulations. Journal of Chemical Physics, 2005, 122, 184704. | 3.0 | 15 |
| 28 | Different chemical dynamics for different conformers of biological molecules: Photoionization of glycine. Journal of Chemical Physics, 2005, 122, 241104. | 3.0 | 24 |
| 29 | Photoionization Dynamics of Glycine:  The First 10 Picoseconds. Journal of Physical Chemistry A, 2004, 108, 11477-11484. | 2.5 | 22 |