

# Tamar Stein

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8691415/publications.pdf>

Version: 2024-02-01

24  
papers

5,639  
citations

471061

17  
h-index

642321

23  
g-index

25  
all docs

25  
docs citations

25  
times ranked

5887  
citing authors

#	ARTICLE	IF	CITATIONS
1	Isomerization of hydrogen cyanide and hydrogen isocyanide in a cluster environment: quantum chemical study. <i>Journal of Chemical Physics</i> , 2022, 156, 054307.	1.2	3
2	Stereoselective Diels–Alder Reactions of <i>gem</i> -Diborylalkenes: Toward the Synthesis of <i>gem</i> -Diboron-Based Polymers. <i>Journal of the American Chemical Society</i> , 2021, 143, 6211-6220.	6.6	19
3	Molecular dynamics reveals formation path of benzonitrile and other molecules in conditions relevant to the interstellar medium. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	3.3	6
4	The Effect of Cluster Size on the Intra-Cluster Ionic Polymerization Process. <i>Molecules</i> , 2021, 26, 4782.	1.7	2
5	Electrochemical deposition of N-heterocyclic carbene monolayers on metal surfaces. <i>Nature Communications</i> , 2020, 11, 5714.	5.8	32
6	Molecular growth upon ionization of van der Waals clusters containing HCCH and HCN is a pathway to prebiotic molecules. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 20337-20348.	1.3	8
7	Photoredox-Mediated Reaction of <i>gem</i> -Diborylalkenes: Reactivity Toward Diverse 1,1-Bisborylalkanes. <i>Chemistry - A European Journal</i> , 2020, 26, 5360-5364.	1.7	24
8	Molecular Formation upon Ionization of van der Waals Clusters and Implication to Astrochemistry. <i>Israel Journal of Chemistry</i> , 2020, 60, 842-849.	1.0	10
9	Probing solvation and reactivity in ionized polycyclic aromatic hydrocarbon-water clusters with photoionization mass spectrometry and electronic structure calculations. <i>Faraday Discussions</i> , 2019, 217, 414-433.	1.6	18
10	Ab initio dynamics and photoionization mass spectrometry reveal ion-molecule pathways from ionized acetylene clusters to benzene cation. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E4125-E4133.	3.3	24
11	Mechanisms of the Formation of Adenine, Guanine, and Their Analogues in UV-Irradiated Mixed NH <sub>3</sub> :H <sub>2</sub> O Molecular Ices Containing Purine. <i>Astrobiology</i> , 2017, 17, 771-785.	1.5	25
12	Probing Ionic Complexes of Ethylene and Acetylene with Vacuum-Ultraviolet Radiation. <i>Journal of Physical Chemistry A</i> , 2016, 120, 5053-5064.	1.1	11
13	Advances in molecular quantum chemistry contained in the Q-Chem 4 program package. <i>Molecular Physics</i> , 2015, 113, 184-215.	0.8	2,561
14	Seniority-based coupled cluster theory. <i>Journal of Chemical Physics</i> , 2014, 141, 244104.	1.2	110
15	Seniority zero pair coupled cluster doubles theory. <i>Journal of Chemical Physics</i> , 2014, 140, 214113.	1.2	147
16	Stability of Hemi-Bonded vs Proton-Transferred Structures of (H <sub>2</sub> O) <sub>2</sub> <sup>+</sup> , (H <sub>2</sub> S) <sub>2</sub> <sup>+</sup> , and (H <sub>2</sub> Se) <sub>2</sub> <sup>+</sup> Studied with Projected Hartree–Fock Methods. <i>Journal of Physical Chemistry A</i> , 2014, 118, 7261-7266.	1.1	26
17	Curvature and Frontier Orbital Energies in Density Functional Theory. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 3740-3744.	2.1	145
18	Calculation of transition dipole moment in fluorescent proteins towards efficient energy transfer. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 4109.	1.3	60

#	ARTICLE	IF	CITATIONS
19	Excitation Gaps of Finite-Sized Systems from Optimally Tuned Range-Separated Hybrid Functionals. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 1515-1531.	2.3	765
20	Charge-Transfer-Like $\tilde{\epsilon}^+$ Excitations in Time-Dependent Density Functional Theory: A Conundrum and Its Solution. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 2408-2415.	2.3	221
21	Fundamental Gaps in Finite Systems from Eigenvalues of a Generalized Kohn-Sham Method. <i>Physical Review Letters</i> , 2010, 105, 266802.	2.9	377
22	Prediction of charge-transfer excitations in coumarin-based dyes using a range-separated functional tuned from first principles. <i>Journal of Chemical Physics</i> , 2009, 131, 244119.	1.2	313
23	Reliable Prediction of Charge Transfer Excitations in Molecular Complexes Using Time-Dependent Density Functional Theory. <i>Journal of the American Chemical Society</i> , 2009, 131, 2818-2820.	6.6	729
24	Accelerated Multiphosphorylated Peptide Synthesis. <i>Organic Process Research and Development</i> , 0, , .	1.3	3