

# Alexander J White

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6114386/publications.pdf>

Version: 2024-02-01

31  
papers

847  
citations

471509

17  
h-index

477307

29  
g-index

32  
all docs

32  
docs citations

32  
times ranked

759  
citing authors

#	ARTICLE	IF	CITATIONS
1	Mixed Stochastic-Deterministic Time-Dependent Density Functional Theory: Application to Stopping Power of Warm Dense Carbon. <i>Journal of Physics Condensed Matter</i> , 2022, , .	1.8	6
2	Proton stopping measurements at low velocity in warm dense carbon. <i>Nature Communications</i> , 2022, 13, .	12.8	13
3	Nonadiabatic Excited-State Molecular Dynamics Methodologies: Comparison and Convergence. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 2970-2982.	4.6	20
4	An Ab Initio Multiple Cloning Method for Non-Adiabatic Excited-State Molecular Dynamics in NWChem. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 3629-3643.	5.3	15
5	Review of the first charged-particle transport coefficient comparison workshop. <i>High Energy Density Physics</i> , 2020, 37, 100905.	1.5	42
6	Fast and Universal Kohn-Sham Density Functional Theory Algorithm for Warm Dense Matter to Hot Dense Plasma. <i>Physical Review Letters</i> , 2020, 125, 055002.	7.8	27
7	Mixed quantum-classical approach to model non-adiabatic electron-nuclear dynamics: Detailed balance and improved surface hopping method. <i>Journal of Chemical Physics</i> , 2020, 153, 074116.	3.0	2
8	Static and dynamic properties of multi-ionic plasma mixtures. <i>Physical Review E</i> , 2020, 101, 033207.	2.1	13
9	NEXMD Software Package for Nonadiabatic Excited State Molecular Dynamics Simulations. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 5771-5783.	5.3	56
10	Non-adiabatic Excited-State Molecular Dynamics: Theory and Applications for Modeling Photophysics in Extended Molecular Materials. <i>Chemical Reviews</i> , 2020, 120, 2215-2287.	47.7	231
11	Multicomponent mutual diffusion in the warm, dense matter regime. <i>Physical Review E</i> , 2019, 100, 033213.	2.1	13
12	Enhancement factor of nuclear reactions in partially ionized plasmas and asymmetric mixtures. <i>Contributions To Plasma Physics</i> , 2019, 59, e201800155.	1.1	3
13	Checking the Salpeter enhancement of nuclear reactions at intermediate coupling in asymmetric mixtures. <i>Physics of Plasmas</i> , 2019, 26, 012702.	1.9	3
14	Improved Ehrenfest Approach to Model Correlated Electron-Nuclear Dynamics. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 433-440.	4.6	10
15	Ab Initio Studies on the Stopping Power of Warm Dense Matter with Time-Dependent Orbital-Free Density Functional Theory. <i>Physical Review Letters</i> , 2018, 121, 145001.	7.8	44
16	Time-dependent orbital-free density functional theory for electronic stopping power: Comparison to the Mermin-Kohn-Sham theory at high temperatures. <i>Physical Review B</i> , 2018, 98, .	3.2	27
17	Photoexcited Nonadiabatic Dynamics of Solvated Push-Pull $\pi$ -Conjugated Oligomers with the NEXMD Software. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 3955-3966.	5.3	39
18	Correlation and transport properties for mixtures at constant pressure and temperature. <i>Physical Review E</i> , 2017, 95, 063202.	2.1	24

#	ARTICLE	IF	CITATIONS
19	Enhancement of nuclear reaction rates in asymmetric binary ionic mixtures. <i>Contributions To Plasma Physics</i> , 2017, 57, 512-517.	1.1	5
20	Coupled wave-packets for non-adiabatic molecular dynamics: a generalization of Gaussian wave-packet dynamics to multiple potential energy surfaces. <i>Chemical Science</i> , 2016, 7, 4905-4911.	7.4	18
21	Non-adiabatic molecular dynamics by accelerated semiclassical Monte Carlo. <i>Journal of Chemical Physics</i> , 2015, 143, 014115.	3.0	13
22	Semiclassical Monte Carlo: A first principles approach to non-adiabatic molecular dynamics. <i>Journal of Chemical Physics</i> , 2014, 141, 184101.	3.0	17
23	Nonequilibrium Atomic Limit for Transport and Optical Response of Molecular Junctions. <i>Journal of Physical Chemistry C</i> , 2014, 118, 11159-11173.	3.1	21
24	Raman Scattering in Molecular Junctions: A Pseudoparticle Formulation. <i>Nano Letters</i> , 2014, 14, 699-703.	9.1	30
25	Effects of Electromagnetic Coupling on Conductance Switching of a Gated Tunnel Junction. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 3545-3550.	4.6	6
26	Coherence in charge and energy transfer in molecular junctions. <i>Physical Review B</i> , 2013, 88, .	3.2	26
27	Quantum transport with two interacting conduction channels. <i>Journal of Chemical Physics</i> , 2013, 138, 174111.	3.0	14
28	Non-Markovian theory of collective plasmon-molecule excitations in nanojunctions combined with classical electrodynamic simulations. , 2013, , .		1
29	Inelastic transport: a pseudoparticle approach. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 13809.	2.8	37
30	Molecular nanoplasmonics: Self-consistent electrodynamics in current-carrying junctions. <i>Physical Review B</i> , 2012, 86, .	3.2	20
31	Collective Plasmon-Molecule Excitations in Nanojunctions: Quantum Consideration. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2738-2743.	4.6	51