

# Andrei Kryjevski

## List of Publications by Year in descending order

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12  
papers

138  
citations

1307594

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1720034

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g-index

12  
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12  
docs citations

12  
times ranked

126  
citing authors

#	ARTICLE	IF	CITATIONS
1	Photoinduced Single- and Multiple-Electron Dynamics Processes Enhanced by Quantum Confinement in Lead Halide Perovskite Quantum Dots. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 3032-3039.	4.6	52
2	Multiple exciton generation in silicon quantum dot arrays: density functional perturbation theory computation. <i>Molecular Physics</i> , 2014, 112, 430-440.	1.7	21
3	Theoretical predictions on efficiency of bi-exciton formation and dissociation in chiral carbon nanotubes. <i>Journal of Chemical Physics</i> , 2016, 145, 154112.	3.0	14
4	Amorphous silicon nanomaterials: Quantum dots versus nanowires. <i>Journal of Renewable and Sustainable Energy</i> , 2013, 5, .	2.0	12
5	Multiple exciton generation in chiral carbon nanotubes: Density functional theory based computation. <i>Journal of Chemical Physics</i> , 2017, 147, 154106.	3.0	12
6	Singlet fission in chiral carbon nanotubes: Density functional theory based computation. <i>Journal of Chemical Physics</i> , 2017, 147, 034106.	3.0	11
7	Dynamics of Charge Transfer and Multiple Exciton Generation in the Doped Silicon Quantum Dot-Carbon Nanotube System: Density Functional Theory-Based Computation. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 5759-5764.	4.6	9
8	Enhanced multiple exciton generation in amorphous silicon nanowires and films. <i>Molecular Physics</i> , 2013, 151, 1-15.	1.7	7
9	Spatially non-uniform field response in arrays of silicon quantum dots: DFT computation. , 2013, , .		0
10	Toward First-Principles Description of Carrier Relaxation in Nanoparticles. <i>ACS Symposium Series</i> , 2015, , 201-213.	0.5	0
11	Comprehensive Study of Multiple Exciton Generation in Chiral Carbon Nanotubes Using Many-Body Perturbation Theory Based on Density Functional Theory Simulations. <i>ACS Symposium Series</i> , 2019, , 157-179.	0.5	0
12	Electronic structure of semiconductor nanoparticles from stochastic evaluation of imaginary-time path integral. <i>Physical Review Research</i> , 2021, 3, .	3.6	0