

# Christopher A O'keefe

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8672273/publications.pdf>

Version: 2024-02-01

16  
papers

892  
citations

687220

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996849

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16  
docs citations

16  
times ranked

998  
citing authors

#	ARTICLE	IF	CITATIONS
1	A molecular shuttle that operates inside a metal-organic framework. <i>Nature Chemistry</i> , 2015, 7, 514-519.	6.6	247
2	Metal-Organic Frameworks with Mechanically Interlocked Pillars: Controlling Ring Dynamics in the Solid-State via a Reversible Phase Change. <i>Journal of the American Chemical Society</i> , 2014, 136, 7403-7409.	6.6	127
3	Mechanically Interlocked Linkers inside Metal-Organic Frameworks: Effect of Ring Size on Rotational Dynamics. <i>Journal of the American Chemical Society</i> , 2015, 137, 9643-9651.	6.6	98
4	Formulation of Metal-Organic Framework-Based Drug Carriers by Controlled Coordination of Methoxy PEG Phosphate: Boosting Colloidal Stability and Redispersibility. <i>Journal of the American Chemical Society</i> , 2021, 143, 13557-13572.	6.6	88
5	Toward an Understanding of SEI Formation and Lithium Plating on Copper in Anode-Free Batteries. <i>Journal of Physical Chemistry C</i> , 2021, 125, 16719-16732.	1.5	55
6	Structural and Dynamic Disorder, Not Ionic Trapping, Controls Charge Transport in Highly Doped Conducting Polymers. <i>Journal of the American Chemical Society</i> , 2022, 144, 3005-3019.	6.6	45
7	Electrolyte Reactivity at the Charged Ni-Rich Cathode Interface and Degradation in Li-Ion Batteries. <i>ACS Applied Materials &amp; Interfaces</i> , 2022, 14, 13206-13222.	4.0	45
8	A Study of Transition-Metal Organometallic Complexes Combining <sup>35</sup> Cl Solid-State NMR Spectroscopy and <sup>35</sup> Cl...NQR Spectroscopy and First-Principles DFT Calculations. <i>Chemistry - A European Journal</i> , 2013, 19, 12396-12414.	1.7	34
9	Thermally Driven Dynamics of a Rotaxane Wheel about an Imidazolium Axle inside a Metal-Organic Framework. <i>ChemPlusChem</i> , 2016, 81, 836-841.	1.3	33
10	New Route to Battery Grade NaPF <sub>6</sub> for Na-Ion Batteries: Expanding the Accessible Concentration. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 24882-24887.	7.2	31
11	An Investigation of Chlorine Ligands in Transition-Metal Complexes via <sup>35</sup> Cl Solid-State NMR and Density Functional Theory Calculations. <i>Inorganic Chemistry</i> , 2014, 53, 9581-9597.	1.9	28
12	Exploring the dynamics of Zr-based metal-organic frameworks containing mechanically interlocked molecular shuttles. <i>Faraday Discussions</i> , 2021, 225, 358-370.	1.6	24
13	Sodium Borates: Expanding the Electrolyte Selection for Sodium-Ion Batteries. <i>Angewandte Chemie - International Edition</i> , 2022, 61, .	7.2	15
14	New Route to Battery Grade NaPF <sub>6</sub> for Na-Ion Batteries: Expanding the Accessible Concentration. <i>Angewandte Chemie</i> , 0, .	1.6	8
15	Dynamics of a [2]rotaxane wheel in a crystalline molecular solid. <i>Chemical Communications</i> , 2021, 57, 8210-8213.	2.2	8
16	Sodium Borates: Expanding the Electrolyte Selection for Sodium-Ion Batteries. <i>Angewandte Chemie</i> , 2022, 134, .	1.6	6