

# Timothy J Fuhrer

## List of Publications by Year in descending order

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17  
papers

585  
citations

1040056

9  
h-index

1058476

14  
g-index

17  
all docs

17  
docs citations

17  
times ranked

438  
citing authors

#	ARTICLE	IF	CITATIONS
1	Faculty Professional Development on Inclusive Pedagogy Yields Chemistry Curriculum Transformation, Equity Awareness, and Community. <i>Journal of Chemical Education</i> , 2022, 99, 291-300.	2.3	4
2	Thiourea-Catalyzed Amidation of Esters: A New Method for the Preparation of Amides. <i>Letters in Organic Chemistry</i> , 2022, 19, .	0.5	0
3	Toward Taming the Chemical Reversibility of Perfluoropyridine through Molecular Design with Applications to Pre- and Postmodifiable Polymer Architectures. <i>Macromolecules</i> , 2021, 54, 5586-5594.	4.8	11
4	Fluoromaticity: The Molecular Orbital Contributions of Fluorine Substituents to the $\pi$ -Systems of Aromatic Rings. <i>ACS Omega</i> , 2021, 6, 32607-32617.	3.5	6
5	Theoretical Prediction and Explanation of Reaction Site Selectivity in the Addition of a Phenoxy Group to Perfluoropyrimidine, Perfluoropyridazine, and Perfluoropyrazine. <i>Molecules</i> , 2021, 26, 7637.	3.8	1
6	Density Functional Theory Investigation of Fulvene-Derivatized Fullerenes as Candidates for Organic Solar Cells. <i>Journal of Physical Chemistry A</i> , 2020, 124, 10324-10329.	2.5	1
7	Temperature Dependent Stabilities of the C <sub>32</sub> and K@C <sub>32</sub> , Ca@C <sub>32</sub> and Sc@C <sub>32</sub> Endohedral Metallofullerene Isomeric Set. <i>Chemical Data Collections</i> , 2020, 28, 100409.	2.3	0
8	Theoretical Explanation of Reaction Site Selectivity in the Addition of a Phenoxy Group to Perfluoropyridine. <i>Journal of Physical Chemistry A</i> , 2019, 123, 9450-9455.	2.5	18
9	Computational study of synergistic effects of electron withdrawing groups as catalysts for fullerene formation. <i>Chemical Data Collections</i> , 2018, 17-18, 415-418.	2.3	0
10	Isolated pentagon rule violating endohedral metallofullerenes explained using the $H_{1/4}$ rule: A statistical mechanical study of the C <sub>84</sub> Isomeric Set. <i>Journal of Computational Chemistry</i> , 2015, 36, 146-150.	3.3	5
11	Enhanced Dipole Moments in Trimetallic Nitride Template Endohedral Metallofullerenes with the Pentalene Motif. <i>Journal of the American Chemical Society</i> , 2013, 135, 3351-3354.	13.7	28
12	A missing link in the transformation from asymmetric to symmetric metallofullerene cages implies a top-down fullerene formation mechanism. <i>Nature Chemistry</i> , 2013, 5, 880-885.	13.6	138
13	Nanoscale Fullerene Compression of an Yttrium Carbide Cluster. <i>Journal of the American Chemical Society</i> , 2012, 134, 8487-8493.	13.7	92
14	Electronic Properties and <sup>13</sup> C NMR Structural Study of Y <sub>3</sub> N@C <sub>88</sub> . <i>Inorganic Chemistry</i> , 2011, 50, 4256-4259.	4.0	24
15	Gd <sub>2</sub> @C <sub>79</sub> N: Isolation, Characterization, and Monoadduct Formation of a Very Stable Heterofullerene with a Magnetic Spin State of $S = 15/2$ . <i>Journal of the American Chemical Society</i> , 2011, 133, 9741-9750.	13.7	104
16	<sup>89</sup> Y and <sup>13</sup> C NMR Cluster and Carbon Cage Studies of an Yttrium Metallofullerene Family, Y <sub>3</sub> N@C <sub>2n</sub> (n) <small>Tj ETQq0 0.0.rgBT /Overlock 1</small>	13.7	81
17	Highly Regioselective Derivatization of Trimetallic Nitride Templated Endohedral Metallofullerenes via a Facile Photochemical Reaction. <i>Journal of the American Chemical Society</i> , 2008, 130, 17755-17760.	13.7	72