

# Yafei Dai

## List of Publications by Year in descending order

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17  
papers

890  
citations

840119

11  
h-index

887659

17  
g-index

17  
all docs

17  
docs citations

17  
times ranked

1372  
citing authors

#	ARTICLE	IF	CITATIONS
1	Cationic Covalent Organic Framework Nanosheets for Fast Li-Ion Conduction. <i>Journal of the American Chemical Society</i> , 2018, 140, 896-899.	6.6	331
2	Super Moisture-Absorbent Gels for All-Weather Atmospheric Water Harvesting. <i>Advanced Materials</i> , 2019, 31, e1806446.	11.1	281
3	Density functional theory study of the structure and energetics of negatively charged oligopyrroles. <i>International Journal of Quantum Chemistry</i> , 2011, 111, 2295-2305.	1.0	56
4	Polar Ketone-Functionalized Metal-Organic Framework Showing a High CO <sub>2</sub> Adsorption Performance. <i>Inorganic Chemistry</i> , 2017, 56, 2363-2366.	1.9	44
5	Energetics, structure, and charge distribution of reduced and oxidized n-pyrrole oligomers: A density functional approach. <i>Journal of Chemical Physics</i> , 2008, 129, 164903.	1.2	27
6	Density Functional Study of Nonlinear Optical Properties of Grossly Warped Nanographene C <sub>80</sub> H <sub>30</sub> . <i>Journal of Physical Chemistry C</i> , 2014, 118, 3313-3318.	1.5	27
7	Diffusion and desorption of oxygen atoms on graphene. <i>Journal of Physics Condensed Matter</i> , 2013, 25, 405301.	0.7	24
8	Fiber-Shaped Electrochemical Capacitors Based on Plasma-Engraved Graphene Fibers with Oxygen Vacancies for Alternating Current Line Filtering Performance. <i>ACS Applied Energy Materials</i> , 2019, 2, 993-999.	2.5	16
9	Monte Carlo study of oligopyrroles in condensed phases. <i>Journal of Chemical Physics</i> , 2010, 133, 034905.	1.2	15
10	Pressure distribution based optimization of phase-coded acoustical vortices. <i>Journal of Applied Physics</i> , 2014, 115, .	1.1	15
11	Porous germanene as a highly efficient gas separation membrane. <i>Nanoscale</i> , 2017, 9, 17505-17512.	2.8	12
12	A short review of nanographenes: structures, properties and applications. <i>Molecular Physics</i> , 2018, 116, 987-1002.	0.8	10
13	Density functional theory study of neutral and oxidized thiophene oligomers. <i>Journal of Chemical Physics</i> , 2013, 139, 184905.	1.2	9
14	Energetics, Structure, and Electron Detachment Spectra of Calcium and Zinc Neutral and Anion Clusters: A Density Functional Theory Study. <i>Journal of Physical Chemistry A</i> , 2008, 112, 11052-11060.	1.1	7
15	Electronic and optical properties of pyrrole and thiophene oligomers: A density functional theory study. <i>International Journal of Quantum Chemistry</i> , 2017, 117, e25453.	1.0	7
16	Coronoid nanographene C <sub>216</sub> as hydrogen purification membrane: A density functional theory study. <i>Carbon</i> , 2018, 135, 112-117.	5.4	6
17	Comment on "Electrical-Driven Transport of Endohedral Fullerene Encapsulating a Single Water Molecule". <i>Physical Review Letters</i> , 2014, 113, 119601.	2.9	3