

Liam Wilbraham

List of Publications by Year in descending order

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Version: 2024-02-01

32
papers

2,242
citations

331670

21
h-index

377865

34
g-index

61
all docs

61
docs citations

61
times ranked

2646
citing authors

#	ARTICLE	IF	CITATIONS
1	Current understanding and challenges of solar-driven hydrogen generation using polymeric photocatalysts. <i>Nature Energy</i> , 2019, 4, 746-760.	39.5	638
2	Accelerated Discovery of Organic Polymer Photocatalysts for Hydrogen Evolution from Water through the Integration of Experiment and Theory. <i>Journal of the American Chemical Society</i> , 2019, 141, 9063-9071.	13.7	264
3	Photocatalytic Hydrogen Evolution from Water Using Fluorene and Dibenzothiophene Sulfone-Conjugated Microporous and Linear Polymers. <i>Chemistry of Materials</i> , 2019, 31, 305-313.	6.7	173
4	Exploring excited states using Time Dependent Density Functional Theory and density-based indexes. <i>Coordination Chemistry Reviews</i> , 2015, 304-305, 166-178.	18.8	118
5	Maximising the hydrogen evolution activity in organic photocatalysts by co-polymerisation. <i>Journal of Materials Chemistry A</i> , 2018, 6, 11994-12003.	10.3	93
6	Side-chain tuning in conjugated polymer photocatalysts for improved hydrogen production from water. <i>Energy and Environmental Science</i> , 2020, 13, 1843-1855.	30.8	92
7	Nitrogen Containing Linear Poly(phenylene) Derivatives for Photo-catalytic Hydrogen Evolution from Water. <i>Chemistry of Materials</i> , 2018, 30, 5733-5742.	6.7	88
8	Revealing the Origins of Mechanically Induced Fluorescence Changes in Organic Molecular Crystals. <i>Advanced Materials</i> , 2018, 30, e1800817.	21.0	82
9	Multiconfiguration Pair-Density Functional Theory Predicts Spin-State Ordering in Iron Complexes with the Same Accuracy as Complete Active Space Second-Order Perturbation Theory at a Significantly Reduced Computational Cost. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 2026-2030.	4.6	57
10	High-Throughput Screening Approach for the Optoelectronic Properties of Conjugated Polymers. <i>Journal of Chemical Information and Modeling</i> , 2018, 58, 2450-2459.	5.4	57
11	Understanding Aggregation-Induced Emission in Molecular Crystals: Insights from Theory. <i>Journal of Physical Chemistry C</i> , 2017, 121, 5747-5752.	3.1	52
12	Hydrogen evolution from water using heteroatom substituted fluorene conjugated co-polymers. <i>Journal of Materials Chemistry A</i> , 2020, 8, 8700-8705.	10.3	47
13	Photocatalytic proton reduction by a computationally identified, molecular hydrogen-bonded framework. <i>Journal of Materials Chemistry A</i> , 2020, 8, 7158-7170.	10.3	45
14	Comparison of structural dynamics and coherence of π - π^* and MLCT light-induced spin state trapping. <i>Chemical Science</i> , 2017, 8, 4978-4986.	7.4	43
15	Structure-activity relationships in well-defined conjugated oligomer photocatalysts for hydrogen production from water. <i>Chemical Science</i> , 2020, 11, 8744-8756.	7.4	41
16	Describing Excited State Intramolecular Proton Transfer in Dual Emissive Systems: A Density Functional Theory Based Analysis. <i>Journal of Physical Chemistry B</i> , 2015, 119, 2459-2466.	2.6	38
17	Mapping binary copolymer property space with neural networks. <i>Chemical Science</i> , 2019, 10, 4973-4984.	7.4	36
18	Photocatalytic polymers of intrinsic microporosity for hydrogen production from water. <i>Journal of Materials Chemistry A</i> , 2021, 9, 19958-19964.	10.3	36

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19	Electrostatic Embedding To Model the Impact of Environment on Photophysical Properties of Molecular Crystals: A Self-Consistent Charge Adjustment Procedure. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 3316-3324.	5.3	28
20	Communication: Evaluating non-empirical double hybrid functionals for spin-state energetics in transition-metal complexes. <i>Journal of Chemical Physics</i> , 2018, 148, 041103.	3.0	28
21	Modelling photophysical properties of metal-organic frameworks: a density functional theory based approach. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 25176-25182.	2.8	27
22	Solvent tuned single molecule dual emission in protic solvents: effect of polarity and H-bonding. <i>Faraday Discussions</i> , 2015, 185, 285-297.	3.2	21
23	Photocatalytically active ladder polymers. <i>Faraday Discussions</i> , 2019, 215, 84-97.	3.2	20
24	Computational high-throughput screening of polymeric photocatalysts: exploring the effect of composition, sequence isomerism and conformational degrees of freedom. <i>Faraday Discussions</i> , 2019, 215, 98-110.	3.2	17
25	Electron Storage System Based on a Two-Way Inversion of Redox Potentials. <i>Journal of the American Chemical Society</i> , 2020, 142, 5162-5176.	13.7	17
26	Using high-throughput virtual screening to explore the optoelectronic property space of organic dyes; finding diketopyrrolopyrrole dyes for dye-sensitized water splitting and solar cells. <i>Sustainable Energy and Fuels</i> , 2021, 5, 704-719.	4.9	15
27	Emissive Azobenzenes Delivered on a Silver Coordination Polymer. <i>Inorganic Chemistry</i> , 2018, 57, 15009-15022.	4.0	14
28	Mapping the optoelectronic property space of small aromatic molecules. <i>Communications Chemistry</i> , 2020, 3, .	4.5	14
29	Insight into the self-assembly of water-soluble perylene bisimide derivatives through a combined computational and experimental approach. <i>Nanoscale</i> , 2019, 11, 15917-15928.	5.6	13
30	Exploring and mapping chemical space with molecular assembly trees. <i>Science Advances</i> , 2021, 7, eabj2465.	10.3	13
31	Modeling UV-Vis spectra of low dimensional materials using electrostatic embedding: The case of CdSe. <i>Journal of Computational Chemistry</i> , 2021, 42, 1212-1224.	3.3	3
32	The potential scarcity, or not, of polymeric overall water splitting photocatalysts. <i>Sustainable Energy and Fuels</i> , 2022, 6, 2233-2242.	4.9	2