## Liang Wu

## List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/1790597/publications.pdf

Version: 2024-02-01

39	816	18	27
papers	citations	h-index	g-index
41	41	41	823
all docs	docs citations	times ranked	citing authors

#	Article	IF	CITATIONS
1	Enantio―and Diastereodivergent Synthesis of Spirocycles through Dualâ€Metalâ€Catalyzed [3+2] Annulation of 2â€Vinyloxiranes with Nucleophilic Dipoles. Angewandte Chemie - International Edition, 2021, 60, 24941-24949.	13.8	110
2	Pd(II)-Catalyzed Asymmetric Addition of Arylboronic Acids to Isatin-Derived Ketimines. Organic Letters, 2016, 18, 288-291.	4.6	74
3	Ni(II)-catalyzed asymmetric alkenylations of ketimines. Nature Communications, 2018, 9, 2258.	12.8	60
4	Palladium-Catalyzed Aerobic Aminooxygenation of Alkenes for Preparation of Isoindolinones. Organic Letters, 2015, 17, 5566-5569.	4.6	59
5	Cobaltâ€Catalyzed Asymmetric Allylation of Cyclic Ketimines. Chemistry - A European Journal, 2018, 24, 1241-1245.	3.3	42
6	Woven Polymer Networks via the Topological Transformation of a [2]Catenane. Journal of the American Chemical Society, 2020, 142, 14343-14349.	13.7	37
7	Palladiumâ€Catalyzed Addition of Arylboronic Acids to <i>para</i> ―Quinone Methides for Preparation of Diarylacetates. Advanced Synthesis and Catalysis, 2017, 359, 1028-1036.	4.3	29
8	Enantio―and Diastereodivergent Synthesis of Spirocycles through Dualâ€Metal atalyzed [3+2] Annulation of 2â€Vinyloxiranes with Nucleophilic Dipoles. Angewandte Chemie, 2021, 133, 25145-25153.	2.0	28
9	Thiophene Derivative as a High Electrochemical Active Anode Material for Sodium-Ion Batteries: The Effect of Backbone Sulfur. Chemistry of Materials, 2018, 30, 8426-8430.	6.7	25
10	Copper (II)/RuPHOXâ€Catalyzed Enantioselective Mannichâ€Type Reaction of Glycine Schiff Bases with Cyclic Ketimines. Advanced Synthesis and Catalysis, 2018, 360, 4625-4633.	4.3	25
11	Enthalpy-change driven synthesis of high-entropy perovskite nanoparticles. Nano Research, 2022, 15, 4867-4872.	10.4	25
12	Predicting Thermodynamic Properties of Alkanes by High-Throughput Force Field Simulation and Machine Learning. Journal of Chemical Information and Modeling, 2018, 58, 2502-2516.	5.4	23
13	A Co(ii)-catalyzed asymmetric ring opening reaction of spiro-epoxyoxindoles with allylboron. Organic Chemistry Frontiers, 2020, 7, 862-867.	4.5	22
14	Orientational ordering and phase behaviour of binary mixtures of hard spheres and hard spherocylinders. Journal of Chemical Physics, 2015, 143, 044906.	3.0	21
15	A Transferrable Coarse-Grained Force Field for Simulations of Polyethers and Polyether Blends. Macromolecules, 2019, 52, 249-261.	4.8	21
16	Ni-Catalyzed Enantioconvergent Coupling of Epoxides with Alkenylboronic Acids: Construction of Oxindoles Bearing Quaternary Carbons. CCS Chemistry, 2020, 2, 623-631.	7.8	21
17	Extracting the mechanisms and kinetic models of complex reactions from atomistic simulation data. Journal of Computational Chemistry, 2019, 40, 1586-1592.	3.3	20
18	Reshaping the Cathodic Catalyst Layer for Anion Exchange Membrane Fuel Cells: From Heterogeneous Catalysis to Homogeneous Catalysis. Angewandte Chemie - International Edition, 2021, 60, 4049-4054.	13.8	19

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19	Understanding and Describing the Liquid-Crystalline States of Polypeptide Solutions: A Coarse-Grained Model of PBLG in DMF. Macromolecules, 2014, 47, 1482-1493.	4.8	15
20	Metal–Tannin Coordination Assembly Route to Nanostructured High-Entropy Oxide Perovskites with Abundant Defects. Chemistry of Materials, 2022, 34, 1746-1755.	6.7	14
21	Demixing, surface nematization, and competing adsorption in binary mixtures of hard rods and hard spheres under confinement. Journal of Chemical Physics, 2018, 148, 164701.	3.0	13
22	Multifunctional Catalyst CuS for Nonaqueous Rechargeable Lithium–Oxygen Batteries. ACS Applied Materials & Samp; Interfaces, 2021, 13, 50065-50075.	8.0	13
23	Monte Carlo simulation of vapor-liquid equilibrium and critical asymmetry of square-well dimer fluid. Journal of Chemical Physics, 2012, 136, 214508.	3.0	11
24	Improved renormalization group theory for critical asymmetry of fluids. Journal of Chemical Physics, 2013, 139, 124103.	3.0	11
25	On accuracy of predicting densities and solubility parameters of polymers using atomistic simulations. Molecular Simulation, 2017, 43, 510-518.	2.0	10
26	Cholesteric ordering predicted using a coarse-grained polymeric model with helical interactions. Soft Matter, 2018, 14, 344-353.	2.7	10
27	Liquid Crystal Phase Behaviour of Attractive Disc-Like Particles. International Journal of Molecular Sciences, 2013, 14, 16414-16442.	4.1	8
28	All-atom and coarse-grained force fields for polydimethylsiloxane. Molecular Simulation, 2017, 43, 1513-1522.	2.0	8
29	Manipulation of cholesteric liquid crystal phase behavior and molecular assembly by molecular chirality. Physical Review E, 2019, 100, 022703.	2.1	8
30	Solvent-free synthesis of N-doped carbon-based catalyst for high-efficient reduction of 4-nitrophenol. Journal of Environmental Chemical Engineering, 2021, 9, 105649.	6.7	8
31	Critical asymmetry in renormalization group theory for fluids. Journal of Chemical Physics, 2013, 138, 234502.	3.0	7
32	Deciphering Single-Bacterium Adhesion Behavior Modulated by Extracellular Electron Transfer. Nano Letters, 2021, 21, 5105-5115.	9.1	5
33	Efficient Electrocatalytic Upgradation of Furan-Based Biomass: Key Roles of a Two-Dimensional Mesoporous Poly(m-phenylenediamine)-Graphene Heterostructure and a Ternary Electrolyte. Macromolecules, 0, , .	4.8	5
34	Ni-Catalyzed Enantioconvergent Coupling of Epoxides with Alkenylboronic Acids: Construction of Oxindoles Bearing Quaternary Carbons. CCS Chemistry, 2020, 2, 623-631.	7.8	4
35	Prediction of self-assemblies of sodium dodecyl sulfate and fragrance additives using coarse-grained force fields. Journal of Molecular Modeling, 2017, 23, 211.	1.8	3
36	Reshaping the Cathodic Catalyst Layer for Anion Exchange Membrane Fuel Cells: From Heterogeneous Catalysis to Homogeneous Catalysis. Angewandte Chemie, 2021, 133, 4095-4100.	2.0	2

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37	Frontispiece: Cobaltâ€Catalyzed Asymmetric Allylation of Cyclic Ketimines. Chemistry - A European Journal, 2018, 24, .	3.3	0
38	Coarse-Grained Force Fields Built on Atomistic Force Fields. Molecular Modeling and Simulation, 2021, , 143-180.	0.2	0
39	A New Parameterization of an All-Atom Force Field for Cellulose. Jom, 2021, 73, 2335-2346.	1.9	0