

# Charlie Tsai

## List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

31  
papers

7,571  
citations

28  
h-index

34  
g-index

34  
ext. papers

8,991  
ext. citations

12.9  
avg, IF

6.23  
L-index

#	Paper	IF	Citations
31	Enhancing Electrocatalytic Water Splitting by Strain Engineering. <i>Advanced Materials</i> , <b>2019</b> , 31, e18070014	14	240
30	Computational Design of Active Site Structures with Improved Transition-State Scaling for Ammonia Synthesis. <i>ACS Catalysis</i> , <b>2018</b> , 8, 4017-4024	13.1	52
29	A Highly Active Molybdenum Phosphide Catalyst for Methanol Synthesis from CO and CO <sub>2</sub> . <i>Angewandte Chemie</i> , <b>2018</b> , 130, 15265-15270	3.6	12
28	A Highly Active Molybdenum Phosphide Catalyst for Methanol Synthesis from CO and CO <sub>2</sub> . <i>Angewandte Chemie - International Edition</i> , <b>2018</b> , 57, 15045-15050	16.4	46
27	Rapid flame doping of Co to WS <sub>2</sub> for efficient hydrogen evolution. <i>Energy and Environmental Science</i> , <b>2018</b> , 11, 2270-2277	35.4	45
26	The Predominance of Hydrogen Evolution on Transition Metal Sulfides and Phosphides under CO <sub>2</sub> Reduction Conditions: An Experimental and Theoretical Study. <i>ACS Energy Letters</i> , <b>2018</b> , 3, 1450-1457	20.1	48
25	Mechanistic insights into heterogeneous methane activation. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 19, 3575-3581	3.6	72
24	Scaling Relations for Adsorption Energies on Doped Molybdenum Phosphide Surfaces. <i>ACS Catalysis</i> , <b>2017</b> , 7, 2528-2534	13.1	30
23	Electrochemical generation of sulfur vacancies in the basal plane of MoS <sub>2</sub> for hydrogen evolution. <i>Nature Communications</i> , <b>2017</b> , 8, 15113	17.4	396
22	Investigating Catalyst-Support Interactions To Improve the Hydrogen Evolution Reaction Activity of Thiomolybdate [Mo <sub>3</sub> S <sub>13</sub> ] <sub>2</sub> Nanoclusters. <i>ACS Catalysis</i> , <b>2017</b> , 7, 7126-7130	13.1	55
21	Promoter Effects of Alkali Metal Cations on the Electrochemical Reduction of Carbon Dioxide. <i>Journal of the American Chemical Society</i> , <b>2017</b> , 139, 11277-11287	16.4	381
20	Understanding trends in C-H bond activation in heterogeneous catalysis. <i>Nature Materials</i> , <b>2017</b> , 16, 225-229	27.9	276
19	Two-Dimensional Materials as Catalysts for Energy Conversion. <i>Catalysis Letters</i> , <b>2016</b> , 146, 1917-1921	2.8	39
18	Automated Discovery and Construction of Surface Phase Diagrams Using Machine Learning. <i>Journal of Physical Chemistry Letters</i> , <b>2016</b> , 7, 3931-3935	6.4	56
17	Direct and continuous strain control of catalysts with tunable battery electrode materials. <i>Science</i> , <b>2016</b> , 354, 1031-1036	33.3	369
16	How Doped MoS <sub>2</sub> Breaks Transition-Metal Scaling Relations for CO <sub>2</sub> Electrochemical Reduction. <i>ACS Catalysis</i> , <b>2016</b> , 6, 4428-4437	13.1	193
15	Chemical and Phase Evolution of Amorphous Molybdenum Sulfide Catalysts for Electrochemical Hydrogen Production. <i>ACS Nano</i> , <b>2016</b> , 10, 624-32	16.7	86

14	Direct Water Decomposition on Transition Metal Surfaces: Structural Dependence and Catalytic Screening. <i>Catalysis Letters</i> , <b>2016</b> , 146, 718-724	2.8	11
13	Scaling Relationships for Binding Energies of Transition Metal Complexes. <i>Catalysis Letters</i> , <b>2016</b> , 146, 304-308	2.8	18
12	Activating and optimizing MoS <sub>2</sub> basal planes for hydrogen evolution through the formation of strained sulphur vacancies. <i>Nature Materials</i> , <b>2016</b> , 15, 48-53	27	1563
11	Predicting Promoter-Induced Bond Activation on Solid Catalysts Using Elementary Bond Orders. <i>Journal of Physical Chemistry Letters</i> , <b>2015</b> , 6, 3670-4	6.4	12
10	Designing an improved transition metal phosphide catalyst for hydrogen evolution using experimental and theoretical trends. <i>Energy and Environmental Science</i> , <b>2015</b> , 8, 3022-3029	35.4	671
9	Rational design of MoS <sub>2</sub> catalysts: tuning the structure and activity via transition metal doping. <i>Catalysis Science and Technology</i> , <b>2015</b> , 5, 246-253	5.5	128
8	The Challenge of Electrochemical Ammonia Synthesis: A New Perspective on the Role of Nitrogen Scaling Relations. <i>ChemSusChem</i> , <b>2015</b> , 8, 2180-6	8.3	666
7	Theoretical insights into the hydrogen evolution activity of layered transition metal dichalcogenides. <i>Surface Science</i> , <b>2015</b> , 640, 133-140	1.8	256
6	Transition-metal doped edge sites in vertically aligned MoS <sub>2</sub> catalysts for enhanced hydrogen evolution. <i>Nano Research</i> , <b>2015</b> , 8, 566-575	10	478
5	Tuning the MoS <sub>2</sub> edge-site activity for hydrogen evolution via support interactions. <i>Nano Letters</i> , <b>2014</b> , 14, 1381-7	11.5	533
4	Understanding the Reactivity of Layered Transition-Metal Sulfides: A Single Electronic Descriptor for Structure and Adsorption. <i>Journal of Physical Chemistry Letters</i> , <b>2014</b> , 5, 3884-9	6.4	57
3	Active edge sites in MoSe <sub>2</sub> and WSe <sub>2</sub> catalysts for the hydrogen evolution reaction: a density functional study. <i>Physical Chemistry Chemical Physics</i> , <b>2014</b> , 16, 13156-64	3.6	364
2	Molybdenum Sulfides and Selenides as Possible Electrocatalysts for CO <sub>2</sub> Reduction. <i>ChemCatChem</i> , <b>2014</b> , 6, 1899-1905	5.2	194
1	Operando Characterization of an Amorphous Molybdenum Sulfide Nanoparticle Catalyst during the Hydrogen Evolution Reaction. <i>Journal of Physical Chemistry C</i> , <b>2014</b> , 118, 29252-29259	3.8	66