

# Hsiu-Wen Wang

## List of Publications by Year in descending order

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Version: 2024-02-01

38  
papers

1,279  
citations

471509

17  
h-index

345221

36  
g-index

40  
all docs

40  
docs citations

40  
times ranked

1964  
citing authors

#	ARTICLE	IF	CITATIONS
1	Resolving the Structure of $\text{Ti}_3\text{C}_2\text{Tx}$ MXenes through Multilevel Structural Modeling of the Atomic Pair Distribution Function. <i>Chemistry of Materials</i> , 2016, 28, 349-359.	6.7	374
2	Multimodality of Structural, Electrical, and Gravimetric Responses of Intercalated MXenes to Water. <i>ACS Nano</i> , 2017, 11, 11118-11126.	14.6	183
3	Boehmite and Gibbsite Nanoplates for the Synthesis of Advanced Alumina Products. <i>ACS Applied Nano Materials</i> , 2018, 1, 7115-7128.	5.0	79
4	Structure and Stability of $\text{SnO}_2$ Nanocrystals and Surface-Bound Water Species. <i>Journal of the American Chemical Society</i> , 2013, 135, 6885-6895.	13.7	67
5	Pre-Sodiated $\text{Ti}_3\text{C}_2\text{Tx}$ MXene Structure and Behavior as Electrode for Sodium-Ion Capacitors. <i>ACS Nano</i> , 2021, 15, 2994-3003.	14.6	54
6	Vibrational Density of States of Strongly H-Bonded Interfacial Water: Insights from Inelastic Neutron Scattering and Theory. <i>Journal of Physical Chemistry C</i> , 2014, 118, 10805-10813.	3.1	48
7	Multi-scale characterization of pore evolution in a combustion metamorphic complex, Hatrurim basin, Israel: Combining (ultra) small-angle neutron scattering and image analysis. <i>Geochimica Et Cosmochimica Acta</i> , 2013, 121, 339-362.	3.9	42
8	Multiscale and Multimodal Characterization of 2D Titanium Carbonitride MXene. <i>Advanced Materials Interfaces</i> , 2020, 7, 1902207.	3.7	35
9	In Situ $^{27}\text{Al}$ NMR Spectroscopy of Aluminate in Sodium Hydroxide Solutions above and below Saturation with Respect to Gibbsite. <i>Inorganic Chemistry</i> , 2018, 57, 11864-11873.	4.0	33
10	Synthesis and structure of synthetically pure and deuterated amorphous (basic) calcium carbonates. <i>Chemical Communications</i> , 2017, 53, 2942-2945.	4.1	28
11	<i>DShaper</i> : an approach for handling missing low- $Q$ data in pair distribution function analysis of nanostructured systems. <i>Journal of Applied Crystallography</i> , 2015, 48, 1651-1659.	4.5	23
12	A $\text{PH}_2\text{O}$ -dependent structural phase transition in the zeolite natrolite. <i>American Mineralogist</i> , 2008, 93, 1191-1194.	1.9	20
13	Precise determination of water exchanges on a mineral surface. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 28819-28828.	2.8	20
14	Transformation of Gibbsite to Boehmite in Caustic Aqueous Solution at Hydrothermal Conditions. <i>Crystal Growth and Design</i> , 2019, 19, 5557-5567.	3.0	19
15	Counteractions Control Local Specific Bonding Interactions and Nucleation Mechanisms in Concentrated Water-in-Salt Solutions. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 3318-3325.	4.6	19
16	Ion-ion interactions enhance aluminum solubility in alkaline suspensions of nano-gibbsite ( $\text{Al}(\text{OH})_3$ ) with sodium nitrite/nitrate. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 4368-4378.	2.8	19
17	The nano- and meso-scale structure of amorphous calcium carbonate. <i>Scientific Reports</i> , 2022, 12, 6870.	3.3	19
18	Phase transitions in natural zeolites and the importance of $\text{P}_2\text{H}_2\text{O}$ . <i>Philosophical Magazine</i> , 2010, 90, 2425-2441.	1.6	18

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19	Combinatorial appraisal of transition states for <i>in situ</i> pair distribution function analysis. <i>Journal of Applied Crystallography</i> , 2017, 50, 1744-1753.	4.5	18
20	Structure determination of the 2.5 hydrate MgSO <sub>4</sub> phase by simulated annealing. <i>American Mineralogist</i> , 2009, 94, 1071-1074.	1.9	17
21	Decoding Oxyanion Aqueous Solvation Structure: A Potassium Nitrate Example at Saturation. <i>Journal of Physical Chemistry B</i> , 2018, 122, 7584-7589.	2.6	14
22	Resolving local configurational contributions to X-ray and neutron radial distribution functions within solutions of concentrated electrolytes – a case study of concentrated NaOH. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 6828-6838.	2.8	14
23	An Atomistic Carbide-Derived Carbon Model Generated Using ReaxFF-Based Quenched Molecular Dynamics. <i>Journal of Carbon Research</i> , 2017, 3, 32.	2.7	13
24	Coupled Multimodal Dynamics of Hydrogen-Containing Ion Networks in Water-Deficient, Sodium Hydroxide-Aluminate Solutions. <i>Journal of Physical Chemistry B</i> , 2018, 122, 12097-12106.	2.6	12
25	X-ray diffraction study of the zeolite natrolite: TP <sub>2</sub> H <sub>2</sub> O phase diagram and phase transitions during dehydration/rehydration. <i>European Journal of Mineralogy</i> , 2010, 22, 271-284.	1.3	10
26	Solvothermal Synthesis and Surface Chemistry To Control the Size and Morphology of Nanoquartz. <i>Crystal Growth and Design</i> , 2015, 15, 5327-5331.	3.0	10
27	Structure and dynamics of water on the forsterite surface. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 27822-27829.	2.8	10
28	Hydroxide promotes ion pairing in the NaNO <sub>2</sub> –NaOH–H <sub>2</sub> O system. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 112-122.	2.8	8
29	Local molecular environment drives speciation and reactivity of ion complexes in concentrated salt solution. <i>Journal of Molecular Liquids</i> , 2021, 340, 116898.	4.9	8
30	Solution and Interface Structure and Dynamics in Geochemistry: Gateway to Link Elementary Processes to Mineral Nucleation and Growth. <i>Crystal Growth and Design</i> , 2022, 22, 853-870.	3.0	8
31	Local structural distortion and electrical transport properties of Bi(Ni <sub>1/2</sub> Ti <sub>1/2</sub> )O <sub>3</sub> perovskite under high pressure. <i>Scientific Reports</i> , 2016, 5, 18229.	3.3	7
32	The “good,” the “bad,” and the “hidden” in neutron scattering and molecular dynamics of ionic aqueous solutions. <i>Journal of Chemical Physics</i> , 2022, 156, .	3.0	6
33	Cluster defects in gibbsite nanoplates grown at acidic to neutral pH. <i>Nanoscale</i> , 2021, 13, 17373-17385.	5.6	5
34	Pressure/temperature fluid cell apparatus for the neutron powder diffractometer instrument: Probing atomic structure in situ. <i>Review of Scientific Instruments</i> , 2014, 85, 125116.	1.3	4
35	Pair distribution function analysis applied to decahedral gold nanoparticles. <i>Physica Scripta</i> , 2017, 92, 114002.	2.5	4
36	Theory-Guided Inelastic Neutron Scattering of Crystalline Alkaline Aluminate Salts Bearing Principal Motifs of Solution-State Species. <i>Inorganic Chemistry</i> , 2021, 60, 16223-16232.	4.0	4

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37	PH <sub>2</sub> O-dependent structural phase transitions in the zeolite mesolite: Real- and reciprocal-space crystal structure refinements. <i>American Mineralogist</i> , 2010, 95, 686-698.	1.9	3
38	Infrared spectroscopic characterization of dehydration and accompanying phase transition behaviors in NAT-topology zeolites. <i>Physics and Chemistry of Minerals</i> , 2012, 39, 277-293.	0.8	3