

Claire E White

List of Publications by Year in descending order

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75
papers

2,657
citations

218381

26
h-index

197535

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75
all docs

75
docs citations

75
times ranked

2316
citing authors

#	ARTICLE	IF	CITATIONS
1	Chemo-mechanical properties of carbon fiber reinforced geopolymer interphase. Journal of the American Ceramic Society, 2022, 105, 1519-1532.	1.9	6
2	Time-dependent phase quantification and local structure analysis of hydroxide-activated slag via X-ray total scattering and molecular modeling. Cement and Concrete Research, 2022, 151, 106642.	4.6	8
3	Upscaling 3D Engineered Trees for Off-Grid Desalination. Environmental Science & Technology, 2022, 56, 1289-1299.	4.6	26
4	The effects of calcium hydroxide and activator chemistry on alkali-activated metakaolin pastes exposed to high temperatures. Cement and Concrete Research, 2022, 154, 106742.	4.6	11
5	Molecular Dynamics Simulations of Reverse Osmosis in Silica Nanopores. Journal of Physical Chemistry C, 2022, 126, 9161-9172.	1.5	3
6	A Roadmap for Production of Cement and Concrete with Low-CO2 Emissions. Waste and Biomass Valorization, 2021, 12, 4745-4775.	1.8	21
7	Modeling of aqueous species interaction energies prior to nucleation in cement-based gel systems. Cement and Concrete Research, 2021, 139, 106266.	4.6	20
8	A parametric study of accelerated carbonation in alkali-activated slag. Cement and Concrete Research, 2021, 145, 106454.	4.6	25
9	The effects of calcium hydroxide and activator chemistry on alkali-activated metakaolin pastes. Cement and Concrete Research, 2021, 145, 106453.	4.6	42
10	Predicting CaO-(MgO)-Al ₂ O ₃ -SiO ₂ glass reactivity in alkaline environments from force field molecular dynamics simulations. Cement and Concrete Research, 2021, 150, 106588.	4.6	16
11	Density functional modeling and total scattering analysis of the atomic structure of a quaternary $\text{CaO} \cdot 3\text{SiO}_2$ glass. https://doi.org/10.1103/PhysRevMaterials.5.053801 Physical Review Materials, 2021, 5, .	4.6	15
12	Selective Fluoride Transport in Subnanometer TiO ₂ Pores. ACS Nano, 2021, 15, 16828-16838.	7.3	16
13	Multiscale pore structure determination of cement paste via simulation and experiment: The case of alkali-activated metakaolin. Cement and Concrete Research, 2020, 137, 106212.	4.6	19
14	Assessment of Ceramic Water Filters for the Removal of Bacterial, Chemical, and Viral Contaminants. Journal of Environmental Engineering, ASCE, 2020, 146, 04020066.	0.7	2
15	Effects of magnesium content and carbonation on the multiscale pore structure of alkali-activated slags. Cement and Concrete Research, 2020, 130, 105979.	4.6	16
16	Generic and Advanced Characterization Techniques. , 2020, , 31-497.		2
17	Understanding solution processing of inorganic materials using cryo-EM. Optical Materials Express, 2020, 10, 119.	1.6	3
18	Solid Residues (Biochar, Bottom Ash, Fly Ash, etc.). , 2020, , 1307-1387.		0

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19	Ceramic Water Filters for the Removal of Bacterial, Chemical, and Viral Contaminants. Journal of Environmental Engineering, ASCE, 2019, 145, .	0.7	7
20	Physical Properties of Interfacial Layers Developed on Weathered Silicates: A Case Study Based on Labradorite Feldspar. Journal of Physical Chemistry C, 2019, 123, 24520-24532.	1.5	12
21	Nanoscale Ordering and Depolymerization of Calcium Silicate Hydrates in the Presence of Alkalis. Journal of Physical Chemistry C, 2019, 123, 24873-24883.	1.5	30
22	Symmetry-Induced Stability in Alkali-Doped Calcium Silicate Hydrate. Journal of Physical Chemistry C, 2019, 123, 14081-14088.	1.5	13
23	Alkali-activation of CaO-FeOx-SiO2 slag: Formation mechanism from in-situ X-ray total scattering. Cement and Concrete Research, 2019, 122, 179-188.	4.6	46
24	<i>In situ</i> quasi-elastic neutron scattering study on the water dynamics and reaction mechanisms in alkali-activated slags. Physical Chemistry Chemical Physics, 2019, 21, 10277-10292.	1.3	20
25	Spray drying OZ439 nanoparticles to form stable, water-dispersible powders for oral malaria therapy. Journal of Translational Medicine, 2019, 17, 97.	1.8	24
26	Amorphous nanoparticles by self-assembly: processing for controlled release of hydrophobic molecules. Soft Matter, 2019, 15, 2400-2410.	1.2	29
27	Nanoscale Chemical Degradation Mechanisms of Sulfate Attack in Alkali-activated Slag. Journal of Physical Chemistry C, 2018, 122, 5992-6004.	1.5	37
28	Rapid Recovery of Clofazimine-Loaded Nanoparticles with Long-Term Storage Stability as Anti-Cryptosporidium Therapy. ACS Applied Nano Materials, 2018, 1, 2184-2194.	2.4	20
29	Highly Surface-Active Ca(OH) ₂ Monolayer as a CO ₂ Capture Material. Nano Letters, 2018, 18, 1786-1793.	4.5	24
30	Drying-induced atomic structural rearrangements in sodium-based calcium-alumino-silicate-hydrate gel and the mitigating effects of ZrO ₂ nanoparticles. Physical Chemistry Chemical Physics, 2018, 20, 8593-8606.	1.3	11
31	A uniaxial load frame for in situ neutron studies of stress-induced changes in cementitious materials and related systems. Review of Scientific Instruments, 2018, 89, 092903.	0.6	4
32	Removal Mechanisms of Contaminants in Ceramic Water Filters. Journal of Environmental Engineering, ASCE, 2018, 144, .	0.7	6
33	Equation of state of the commensurate phases of $\text{CaO} \cdot x\text{FeO} \cdot \text{SiO}_2$ glassy slags and resultant inorganic polymer binders. Journal of the American Ceramic Society, 2018, 101, 5846-5857.	1.9	40
34	Accumulators for the Capture of Heavy Metals in Thermal Conversion Systems. Journal of Environmental Engineering, ASCE, 2018, 144, 04018118.	0.7	0
35	Molecular structure of $\text{CaO} \cdot x\text{FeO} \cdot \text{SiO}_2$ glassy slags and resultant inorganic polymer binders. Journal of the American Ceramic Society, 2018, 101, 5846-5857.	1.9	40
36	Anisotropic crystallization in solution processed chalcogenide thin film by linearly polarized laser. Applied Physics Letters, 2017, 110, .	1.5	11

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37	Mechanism of zinc oxide retardation in alkali-activated materials: an in situ X-ray pair distribution function investigation. <i>Journal of Materials Chemistry A</i> , 2017, 5, 11794-11804.	5.2	89
38	Evolution of the pore structure during the early stages of the alkali-activation reaction: an in situ small-angle neutron scattering investigation. <i>Journal of Applied Crystallography</i> , 2017, 50, 61-75.	1.9	20
39	Structure and properties of clay ceramics for thermal energy storage. <i>Journal of the American Ceramic Society</i> , 2017, 100, 4748-4759.	1.9	10
40	Impact of activator chemistry on permeability of alkali-activated slags. <i>Journal of the American Ceramic Society</i> , 2017, 100, 4848-4859.	1.9	20
41	Environmental stability of 2D anisotropic tellurium containing nanomaterials: anisotropic to isotropic transition. <i>Nanoscale</i> , 2017, 9, 12288-12294.	2.8	41
42	Novel Surface Molecular Functionalization Route To Enhance Environmental Stability of Tellurium-Containing 2D Layers. <i>ACS Applied Materials & Interfaces</i> , 2017, 9, 44625-44631.	4.0	15
43	Nanoscale Charge-Balancing Mechanism in Alkali-Substituted Calcium Silicate Hydrate Gels. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 5266-5272.	2.1	47
44	Effects of temperature on the atomic structure of synthetic calcium silicate deuterate gels: A neutron pair distribution function investigation. <i>Cement and Concrete Research</i> , 2016, 79, 93-100.	4.6	26
45	Impact of chemical variability of ground granulated blast-furnace slag on the phase formation in alkali-activated slag pastes. <i>Cement and Concrete Research</i> , 2016, 89, 310-319.	4.6	82
46	High-pressure polymorphism of PbF_2 to 75 GPa. <i>Physical Review B</i> , 2016, 94, .	1.1	9
47	Modeling the Formation of Alkali Aluminosilicate Gels at the Mesoscale Using Coarse-Grained Monte Carlo. <i>Langmuir</i> , 2016, 32, 11580-11590.	1.6	16
48	Elucidating the atomic structures of different sources of fly ash using X-ray and neutron PDF analysis. <i>Fuel</i> , 2016, 177, 148-156.	3.4	18
49	Nanoscale heterogeneities in a fractured alkali-activated slag binder: A helium ion microscopy analysis. <i>Cement and Concrete Research</i> , 2016, 79, 45-48.	4.6	7
50	Role of Magnesium-Stabilized Amorphous Calcium Carbonate in Mitigating the Extent of Carbonation in Alkali-Activated Slag. <i>Chemistry of Materials</i> , 2015, 27, 6625-6634.	3.2	52
51	In situ X-ray pair distribution function analysis of accelerated carbonation of a synthetic calcium silicate hydrate gel. <i>Journal of Materials Chemistry A</i> , 2015, 3, 8597-8605.	5.2	78
52	Intrinsic differences in atomic ordering of calcium (alumino)silicate hydrates in conventional and alkali-activated cements. <i>Cement and Concrete Research</i> , 2015, 67, 66-73.	4.6	72
53	Uncovering the True Atomic Structure of Disordered Materials: The Structure of a Hydrated Amorphous Magnesium Carbonate ($MgCO_3 \cdot 3D_2O$). <i>Chemistry of Materials</i> , 2014, 26, 2693-2702.	3.2	26
54	Structure of kaolinite and influence of stacking faults: Reconciling theory and experiment using inelastic neutron scattering analysis. <i>Journal of Chemical Physics</i> , 2013, 138, 194501.	1.2	12

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55	In situ synchrotron X-ray pair distribution function analysis of the early stages of gel formation in metakaolin-based geopolymers. <i>Applied Clay Science</i> , 2013, 73, 17-25.	2.6	82
56	Inelastic neutron scattering analysis of the thermal decomposition of kaolinite to metakaolin. <i>Chemical Physics</i> , 2013, 427, 82-86.	0.9	14
57	In situ X-ray pair distribution function analysis of geopolymer gel nanostructure formation kinetics. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 8573.	1.3	60
58	Nanostructural characterization of geopolymers by advanced beamline techniques. <i>Cement and Concrete Composites</i> , 2013, 36, 56-64.	4.6	33
59	Pair distribution function analysis of amorphous geopolymer precursors and binders: the importance of complementary molecular simulations. <i>Zeitschrift für Kristallographie</i> , 2012, 227, 304-312.	1.1	11
60	Comment on "Structure-Directing Role of Counterions in the Initial Stage of Zeolite Synthesis". <i>Journal of Physical Chemistry C</i> , 2012, 116, 1619-1621.	1.5	10
61	Molecular mechanisms responsible for the structural changes occurring during geopolymerization: Multiscale simulation. <i>AIChE Journal</i> , 2012, 58, 2241-2253.	1.8	60
62	X-ray microtomography shows pore structure and tortuosity in alkali-activated binders. <i>Cement and Concrete Research</i> , 2012, 42, 855-864.	4.6	394
63	Density functional modelling of silicate and aluminosilicate dimerisation solution chemistry. <i>Dalton Transactions</i> , 2011, 40, 1348-1355.	1.6	66
64	The use of XANES to clarify issues related to bonding environments in metakaolin: a discussion of the paper S. Sperinck et al., "Dehydroxylation of kaolinite to metakaolin-a molecular dynamics study", <i>J. Mater. Chem.</i> , 2011, 21, 2118-2125. <i>Journal of Materials Chemistry</i> , 2011, 21, 7007.	6.7	23
65	Effect of Temperature on the Local Structure of Kaolinite Intercalated with Potassium Acetate. <i>Chemistry of Materials</i> , 2011, 23, 188-199.	3.2	33
66	Quantitative Mechanistic Modeling of Silica Solubility and Precipitation during the Initial Period of Zeolite Synthesis. <i>Journal of Physical Chemistry C</i> , 2011, 115, 9879-9888.	1.5	28
67	Evolution of Local Structure in Geopolymer Gels: An <i>In Situ</i> Neutron Pair Distribution Function Analysis. <i>Journal of the American Ceramic Society</i> , 2011, 94, 3532-3539.	1.9	110
68	Treatment of hydrogen background in bulk and nanocrystalline neutron total scattering experiments. <i>Journal of Applied Crystallography</i> , 2011, 44, 532-539.	1.9	24
69	Discussion of Y. Zhang et al., "Study of ion cluster reorientation process of geopolymerisation reaction using semi-empirical AM1 calculations", <i>Cem Concr Res</i> 39(12): 1174-1179; 2009. <i>Cement and Concrete Research</i> , 2010, 40, 827-828.	4.6	4
70	Extracting differential pair distribution functions using <i>MIXSCAT</i> . <i>Journal of Applied Crystallography</i> , 2010, 43, 635-638.	1.9	9
71	The Effects of Temperature on the Local Structure of Metakaolin-Based Geopolymer Binder: A Neutron Pair Distribution Function Investigation. <i>Journal of the American Ceramic Society</i> , 2010, 93, 3486-3492.	1.9	135
72	Combining density functional theory (DFT) and pair distribution function (PDF) analysis to solve the structure of metastable materials: the case of metakaolin. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 3239.	1.3	137

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73	Density Functional Modeling of the Local Structure of Kaolinite Subjected to Thermal Dehydroxylation. <i>Journal of Physical Chemistry A</i> , 2010, 114, 4988-4996.	1.1	113
74	What Is the Structure of Kaolinite? Reconciling Theory and Experiment. <i>Journal of Physical Chemistry B</i> , 2009, 113, 6756-6765.	1.2	63
75	Modeling Silica Nanoparticle Dissolution in TPAOH~TEOS~H2O Solutions. <i>Journal of Physical Chemistry C</i> , 2008, 112, 14769-14775.	1.5	16