Simon Billinge

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.

 220
 14,441
 58
 116

 papers
 citations
 h-index
 g-index

 241
 8
 6.65

 ext. papers
 ext. citations
 avg, IF
 L-index

#	Paper	IF	Citations
220	Controlling desolvation through polymer-assisted grinding. <i>CrystEngComm</i> , 2022 , 24, 2305-2313	3.3	O
219	Mechanistic Insight into the Precursor Chemistry of ZrO and HfO Nanocrystals; towards Size-Tunable Syntheses <i>Jacs Au</i> , 2022 , 2, 827-838		О
218	Recent advances and applications of deep learning methods in materials science. <i>Npj Computational Materials</i> , 2022 , 8,	10.9	19
217	nmfMapping: a cloud-based web application for non-negative matrix factorization of powder diffraction and pair distribution function datasets <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2022 , 78, 242-248	1.7	О
216	Hidden Local Symmetry Breaking in Silver Diamondoid Compounds is Root Cause of Ultralow Thermal Conductivity <i>Advanced Materials</i> , 2022 , e2202255	24	2
215	Powder diffraction. Nature Reviews Methods Primers, 2021, 1,		2
214	Structural Analysis of Molecular Materials Using the Pair Distribution Function. <i>Chemical Reviews</i> , 2021 ,	68.1	11
213	Toward In Situ Synchrotron Mapping of Crystal Selection Processes during Crystal Growth. <i>Chemistry of Materials</i> , 2021 , 33, 3359-3367	9.6	1
212	Superatomic solid solutions. <i>Nature Chemistry</i> , 2021 , 13, 607-613	17.6	4
211	Validation of non-negative matrix factorization for rapid assessment of large sets of atomic pair distribution function data. <i>Journal of Applied Crystallography</i> , 2021 , 54, 768-775	3.8	4
210	Dual Orbital Degeneracy Lifting in a Strongly Correlated Electron System. <i>Physical Review Letters</i> , 2021 , 126, 186402	7.4	1
209	Lowering Ternary Oxide Synthesis Temperatures by Solid-State Cometathesis Reactions. <i>Chemistry of Materials</i> , 2021 , 33, 3692-3701	9.6	3
208	A cloud platform for atomic pair distribution function analysis: PDFitc. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2021 , 77, 2-6	1.7	9
207	Ferroelectric state and polarization switching behaviour of ultrafine BaTiO3 nanoparticles with large-scale size uniformity. <i>Journal of Materials Chemistry C</i> , 2021 , 9, 5267-5276	7.1	0
206	Local and long-range atomic/magnetic structure of non-stoichiometric spinel iron oxide nanocrystallites. <i>IUCrJ</i> , 2021 , 8, 33-45	4.7	3
205	A high throughput optical method for studying compositional effects in electrocatalysts for CO reduction. <i>Nature Communications</i> , 2021 , 12, 1114	17.4	7
204	Network-Forming Liquids from Metal-Bis(acetamide) Frameworks with Low Melting Temperatures. Journal of the American Chemical Society, 2021, 143, 2801-2811	16.4	20

203	Autonomous experimentation systems for materials development: A community perspective. <i>Matter</i> , 2021 , 4, 2702-2726	12.7	26	
202	Modern crystallography and its foundations. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2021 , 77, 1	1.7	O	
201	Linking far-from-equilibrium defect structures in ceramics to electromagnetic driving forces. Journal of Materials Chemistry A, 2021 , 9, 8425-8434	13	О	
200	Understanding electronic peculiarities in tetragonal FeSe as local structural symmetry breaking. <i>Physical Review B</i> , 2020 , 102,	3.3	10	
199	sasPDF: pair distribution function analysis of nanoparticle assemblies from small-angle scattering data. <i>Journal of Applied Crystallography</i> , 2020 , 53, 699-709	3.8	6	
198	Ultrafast x-ray diffraction study of melt-front dynamics in polycrystalline thin films. <i>Science Advances</i> , 2020 , 6, eaax2445	14.3	9	
197	A thermal-gradient approach to variable-temperature measurements resolved in space. <i>Journal of Applied Crystallography</i> , 2020 , 53, 662-670	3.8	8	
196	Cluster-mining: an approach for determining core structures of metallic nanoparticles from atomic pair distribution function data. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2020 , 76, 24-31	1.7	17	
195	Structure-mining: screening structure models by automated fitting to the atomic pair distribution function over large numbers of models. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2020 , 76, 395-409	1.7	9	
194	Rapid desolvation-triggered domino lattice rearrangement in a metal-organic framework. <i>Nature Chemistry</i> , 2020 , 12, 90-97	17.6	60	
193	Active Reaction Control of Cu Redox State Based on Real-Time Feedback from In Situ Synchrotron Measurements. <i>Journal of the American Chemical Society</i> , 2020 , 142, 18758-18762	16.4	6	
192	Defect-Accommodating Intermediates Yield Selective Low-Temperature Synthesis of YMnO Polymorphs. <i>Inorganic Chemistry</i> , 2020 , 59, 13639-13650	5.1	7	
191	Complete Strain Mapping of Nanosheets of Tantalum Disulfide. <i>ACS Applied Materials & Materials & Interfaces</i> , 2020 , 12, 43173-43179	9.5	3	
190	Two-orbital degeneracy lifted local precursor to a metal-insulator transition in MgTi2O4. <i>Physical Review B</i> , 2020 , 102,	3.3	3	
189	Local Structural Effects Due to Micronization and Amorphization on an HIV Treatment Active Pharmaceutical Ingredient. <i>Molecular Pharmaceutics</i> , 2020 , 17, 2370-2389	5.6	8	
188	Nanometre-scale structure from powder diffraction: total scattering and atomic pair distribution function analysis 2019 , 649-672		8	
187	Local orbital degeneracy lifting as a precursor to an orbital-selective Peierls transition. <i>Nature Communications</i> , 2019 , 10, 3638	17.4	21	
186	Quantitative Structural Characterization of Catalytically Active TiO2 Nanoparticles. <i>ACS Applied Nano Materials</i> , 2019 , 2, 6268-6276	5.6	9	

185	Two-Dimensional Arrays of Transition Metal Nitride Nanocrystals. <i>Advanced Materials</i> , 2019 , 31, e19023	39 <u>-3</u> 4	59
184	Photoinduced dynamics of nematic order parameter in FeSe. <i>Physical Review B</i> , 2019 , 99,	3.3	10
183	Stabilization of reactive CoO cubane oxygen-evolution catalysts within porous frameworks. Proceedings of the National Academy of Sciences of the United States of America, 2019, 116, 11630-11639.	9 ^{11.5}	29
182	The rise of the X-ray atomic pair distribution function method: a series of fortunate events. Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences, 2019, 377, 201804	13	49
181	Microporous Battery Electrodes from Molecular Cluster Precursors. <i>ACS Applied Materials & amp; Interfaces</i> , 2019 , 11, 11292-11297	9.5	4
180	Scalable Synthesis of Ultrathin Mn3N2 Exhibiting Room-Temperature Antiferromagnetism. <i>Advanced Functional Materials</i> , 2019 , 29, 1809001	15.6	37
179	ProtonElectron Conductivity in Thin Films of a CobaltDxygen Evolving Catalyst. <i>ACS Applied Energy Materials</i> , 2019 , 2, 3-12	6.1	30
178	Room temperature local nematicity in FeSe superconductor. <i>Physical Review B</i> , 2019 , 100,	3.3	19
177	Using a machine learning approach to determine the space group of a structure from the atomic pair distribution function. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019 , 75, 633-643	3 ^{1.7}	26
176	Zirconium Phosphate: The Pathway from Turbostratic Disorder to Crystallinity. <i>Inorganic Chemistry</i> , 2019 , 58, 14260-14274	5.1	8
175	Algorithm for distance list extraction from pair distribution functions. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019 , 75, 658-668	1.7	5
174	Size-Dependent Lattice Dynamics of Atomically Precise Cadmium Selenide Quantum Dots. <i>Physical Review Letters</i> , 2019 , 122, 026101	7.4	6
173	Stoichiometric Control over Ferroic Behavior in Ba(Ti1NFex)O3 Nanocrystals. <i>Chemistry of Materials</i> , 2019 , 31, 1318-1335	9.6	21
172	Synthesis and Properties of Plasmonic Boron-Hyperdoped Silicon Nanoparticles. <i>Advanced Functional Materials</i> , 2019 , 29, 1807788	15.6	17
171	Anthracene as a Launchpad for a Phosphinidene Sulfide and for Generation of a Phosphorus-Sulfur Material Having the Composition PS, a Vulcanized Red Phosphorus That Is Yellow. <i>Journal of the American Chemical Society</i> , 2019 , 141, 431-440	16.4	17
170	Early stage structural development of prototypical zeolitic imidazolate framework (ZIF) in solution. <i>Nanoscale</i> , 2018 , 10, 4291-4300	7.7	40
169	PDFgetN3: atomic pair distribution functions from neutron powder diffraction data using adlhoc corrections. <i>Journal of Applied Crystallography</i> , 2018 , 51, 1492-1497	3.8	16
168	Correlated local dipoles in PbTe. <i>Physical Review Materials</i> , 2018 , 2,	3.2	25

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167	Synthesis, characterization, and growth mechanism of motifs of ultrathin cobalt-substituted NaFeSi2O6 nanowires. <i>CrystEngComm</i> , 2018 , 20, 223-236	3.3	3
166	StructureBroperty insights into nanostructured electrodes for Li-ion batteries from local structural and diffusional probes. <i>Journal of Materials Chemistry A</i> , 2018 , 6, 127-137	13	17
165	Magnetism in semiconducting molybdenum dichalcogenides. Science Advances, 2018, 4, eaat3672	14.3	56
164	Improved Models for Metallic Nanoparticle Cores from Atomic Pair Distribution Function (PDF) Analysis. <i>Journal of Physical Chemistry C</i> , 2018 , 122, 29498-29506	3.8	28
163	Emphanitic anharmonicity in PbSe at high temperature and anomalous electronic properties in the PbQ(Q=S,Se,Te) system. <i>Physical Review B</i> , 2018 , 98,	3.3	17
162	Recent results on assigned and unassigned distance geometry with applications to protein molecules and nanostructures. <i>Annals of Operations Research</i> , 2018 , 271, 161-203	3.2	17
161	Barium titanate nanoparticles: Short-range lattice distortions with long-range cubic order. <i>Physical Review B</i> , 2018 , 98,	3.3	8
160	Pair Distribution Function Analysis of ZrO[Nanocrystals and Insights in the Formation of ZrOE/Ba[LuD[Nanocomposites. <i>Materials</i> , 2018 , 11,	3.5	13
159	Hollow organic capsules assemble into cellular semiconductors. <i>Nature Communications</i> , 2018 , 9, 1957	17.4	20
158	Coherent Nanotwins and Dynamic Disorder in Cesium Lead Halide Perovskite Nanocrystals. <i>ACS Nano</i> , 2017 , 11, 3819-3831	16.7	181
157	Modelling and validation of particle size distributions of supported nanoparticles using the pair distribution function technique. <i>Journal of Applied Crystallography</i> , 2017 , 50, 741-748	3.8	15
156	Superconducting order from disorder in 2H-TaSe 2lk S x. <i>Npj Quantum Materials</i> , 2017 , 2,	5	41
155	Signatures of the topological s superconducting order parameter in the type-II Weyl semimetal T -MoTe. <i>Nature Communications</i> , 2017 , 8, 1082	17.4	62
154	Robust Nanostructure from High Throughput Powder Diffraction Data. <i>Microscopy and Microanalysis</i> , 2017 , 23, 172-173	0.5	
153	Unlocking the structure of mixed amorphous-crystalline ceramic oxide films synthesized under low temperature electromagnetic excitation. <i>Journal of Materials Chemistry A</i> , 2017 , 5, 18434-18441	13	16
152	Cation Exchange Induced Transformation of InP Magic-Sized Clusters. <i>Chemistry of Materials</i> , 2017 , 29, 7984-7992	9.6	49
151	Pressure tuning of structure, superconductivity, and novel magnetic order in the Ce-underdoped electron-doped cuprate T? P r1.3\(\text{La0.7CexCuO4 (x=0.1)}. \(\text{Physical Review B}, \(\text{2017}, 96, \)	3.3	6
150	2D molybdenum and vanadium nitrides synthesized by ammoniation of 2D transition metal carbides (MXenes). <i>Nanoscale</i> , 2017 , 9, 17722-17730	7.7	192

149	Local Environment of Terbium(III) Ions in Layered Nanocrystalline Zirconium(IV) Phosphonate-Phosphate Ion Exchange Materials. <i>Inorganic Chemistry</i> , 2017 , 56, 8837-8846	5.1	24
148	Real-space investigation of short-range magnetic correlations in fluoride pyrochlores NaCaCo2F7 and NaSrCo2F7 with magnetic pair distribution function analysis. <i>Physical Review Materials</i> , 2017 , 1,	3.2	9
147	Verification of Anderson Superexchange in MnO via Magnetic Pair Distribution Function Analysis and ablinitio Theory. <i>Physical Review Letters</i> , 2016 , 116, 197204	7.4	23
146	Towards atomistic understanding of polymorphism in the solvothermal synthesis of ZrO nanoparticles. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2016 , 72, 645-650	1.7	28
145	Polymorphism in magic-sized Au144(SR)60 clusters. <i>Nature Communications</i> , 2016 , 7, 11859	17.4	126
144	Towards solution and refinement of organic crystal structures by fitting to the atomic pair distribution function. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2016 , 72, 62-72	1.7	17
143	Control of electronic properties of 2D carbides (MXenes) by manipulating their transition metal layers. <i>Nanoscale Horizons</i> , 2016 , 1, 227-234	10.8	242
142	Recrystallization, Phase Composition, and Local Structure of Amorphous Lactose from the Total Scattering Pair Distribution Function. <i>Crystal Growth and Design</i> , 2016 , 16, 210-220	3.5	25
141	Celebrating 100 years of the Debye scattering equation. <i>Acta Crystallographica Section A:</i> Foundations and Advances, 2016 , 72, 589-590	1.7	16
140	Cooperative coupling of static magnetism and bulk superconductivity in the stripe phase of La2\(\text{BaxCuO4}: \text{ Pressure- and doping-dependent studies. } \text{ Physical Review B, 2016, 94,}	3.3	8
139	Volume-wise destruction of the antiferromagnetic Mott insulating state through quantum tuning. <i>Nature Communications</i> , 2016 , 7, 12519	17.4	22
138	Assigned and unassigned distance geometry: applications to biological molecules and nanostructures. <i>4or</i> , 2016 , 14, 337-376	1.4	34
137	Local atomic and magnetic structure of dilute magnetic semiconductor (Ba,K)(Zn,Mn)2As2. <i>Physical Review B</i> , 2016 , 94,	3.3	23
136	Structures of Hard Phases in Thermoplastic Polyurethanes. <i>Macromolecules</i> , 2016 , 49, 7350-7358	5.5	23
135	Atomic electron tomography: 3D structures without crystals. Science, 2016, 353,	33.3	129
134	Direct Observation of Dynamic Symmetry Breaking above Room Temperature in Methylammonium Lead Iodide Perovskite. <i>ACS Energy Letters</i> , 2016 , 1, 880-887	20.1	177
133	Investigating short-range magnetic correlations in real space with the magnetic pair distribution function (mPDF). <i>Neutron News</i> , 2016 , 27, 14-16	0.4	1
132	Novel trends in pair distribution function approaches on bulk systems with nanoscale heterogeneities. <i>Neutron News</i> , 2016 , 27, 27-31	0.4	1

131	Enhanced thermoelectric power and electronic correlations in RuSe2. APL Materials, 2015, 3, 041513	5.7	10
130	Calibration and data collection protocols for reliable lattice parameter values in electron pair distribution function studies. <i>Journal of Applied Crystallography</i> , 2015 , 48, 244-251	3.8	17
129	Algorithm for systematic peak extraction from atomic pair distribution functions. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2015 , 71, 392-409	1.7	9
128	Structural Evolution of Iron Antimonides from Amorphous Precursors to Crystalline Products Studied by Total Scattering Techniques. <i>Journal of the American Chemical Society</i> , 2015 , 137, 9652-8	16.4	16
127	Correlating Size and Composition-Dependent Effects with Magnetic, MBsbauer, and Pair Distribution Function Measurements in a Family of Catalytically Active Ferrite Nanoparticles. <i>Chemistry of Materials</i> , 2015 , 27, 3572-3592	9.6	61
126	X-Ray Diffraction Computed Tomography for Structural Analysis of Electrode Materials in Batteries. <i>Journal of the Electrochemical Society</i> , 2015 , 162, A1310-A1314	3.9	39
125	Complex modeling: a strategy and software program for combining multiple information sources to solve ill posed structure and nanostructure inverse problems. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2015 , 71, 562-8	1.7	152
124	Nature of Activated Manganese Oxide for Oxygen Evolution. <i>Journal of the American Chemical Society</i> , 2015 , 137, 14887-904	16.4	295
123	Resonant spin tunneling in randomly oriented nanospheres of Mn12 acetate. <i>Physical Review B</i> , 2015 , 91,	3.3	3
122	Total-scattering pair-distribution function of organic material from powder electron diffraction data. <i>Microscopy and Microanalysis</i> , 2015 , 21, 459-71	0.5	13
121	Demonstration of thin film pair distribution function analysis (tfPDF) for the study of local structure in amorphous and crystalline thin films. <i>IUCrJ</i> , 2015 , 2, 481-9	4.7	41
120	Reconciliation of local and long-range tilt correlations in underdoped La2\BaxCuO4(0\D.155). <i>Physical Review B</i> , 2015 , 91,	3.3	25
119	Magnetic structure determination from the magnetic pair distribution function (mPDF): ground state of MnO. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2015 , 71, 325-34	1.7	29
118	Imaging Dirac-mass disorder from magnetic dopant atoms in the ferromagnetic topological insulator Crx(Bi0.1Sb0.9)2-xTe3. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, 1316-21	11.5	147
117	Two-Step Nucleation and Growth of InP Quantum Dots via Magic-Sized Cluster Intermediates. <i>Chemistry of Materials</i> , 2015 , 27, 1432-1441	9.6	176
116	Detection and characterization of nanoparticles in suspension at low concentrations using the X-ray total scattering pair distribution function technique. <i>Nanoscale</i> , 2015 , 7, 5480-7	7.7	30
115	Modelling pair distribution functions (PDFs) of organic compounds: describing both intra- and intermolecular correlation functions in calculated PDFs. <i>Journal of Applied Crystallography</i> , 2015 , 48, 171-178	3.8	37
114	Hollandites as a new class of multiferroics. <i>Scientific Reports</i> , 2014 , 4, 6203	4.9	26

113	Structure of Sulfate Adsorption Complexes on Ferrihydrite. <i>Environmental Science and Technology Letters</i> , 2014 , 1, 97-101	11	65
112	Structure of methylammonium lead iodide within mesoporous titanium dioxide: active material in high-performance perovskite solar cells. <i>Nano Letters</i> , 2014 , 14, 127-33	11.5	258
111	Local off-centering symmetry breaking in the high-temperature regime of SnTe. <i>Physical Review B</i> , 2014 , 89,	3.3	58
110	Atomic structures and gram scale synthesis of three tetrahedral quantum dots. <i>Journal of the American Chemical Society</i> , 2014 , 136, 10645-53	16.4	139
109	Bulk metallic glass-like scattering signal in small metallic nanoparticles. ACS Nano, 2014 , 8, 6163-70	16.7	23
108	Mechanisms for iron oxide formation under hydrothermal conditions: an in situ total scattering study. <i>ACS Nano</i> , 2014 , 8, 10704-14	16.7	65
107	Synthesis and characterization of two-dimensional Nb4C3 (MXene). <i>Chemical Communications</i> , 2014 , 50, 9517-20	5.8	321
106	Local vibrations and negative thermal expansion in ZrW2O8. <i>Physical Review Letters</i> , 2014 , 112, 045505	7.4	75
105	Robust structure and morphology parameters for CdS nanoparticles by combining small-angle X-ray scattering and atomic pair distribution function data in a complex modeling framework. Journal of Applied Crystallography, 2014 , 47, 561-565	3.8	16
104	Quantitative Structural Analysis of Nanoparticles Using Electron Pair Distribution Function (ePDF). <i>Microscopy and Microanalysis</i> , 2014 , 20, 630-631	0.5	1
103	Cu(Ir xCrx)BD model system for studying nanoscale phase coexistence at the metal-insulator transition. <i>Scientific Reports</i> , 2014 , 4, 4081	4.9	12
102	On the estimation of statistical uncertainties on powder diffraction and small-angle scattering data from two-dimensional X-ray detectors. <i>Journal of Applied Crystallography</i> , 2014 , 47, 1273-1283	3.8	15
101	Structure of nanocrystalline Ti3C2 MXene using atomic pair distribution function. <i>Physical Review Letters</i> , 2014 , 112, 125501	7.4	129
100	Nanoscale coherent intergrowthlike defects in a crystal of La1.9Ca1.1Cu2O6+Imade superconducting by high-pressure oxygen annealing. <i>Physical Review B</i> , 2014 , 90,	3.3	3
99	Intra-unit-cell nematic charge order in the titanium-oxypnictide family of superconductors. <i>Nature Communications</i> , 2014 , 5, 5761	17.4	22
98	Evolution of atomic structure during nanoparticle formation. <i>IUCrJ</i> , 2014 , 1, 165-71	4.7	39
97	Magnetic pair distribution function analysis of local magnetic correlations. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2014 , 70, 3-11	1.7	35
96	Pair distribution function computed tomography. <i>Nature Communications</i> , 2013 , 4, 2536	17.4	75

(2011-2013)

95	Toward Phase Quantification at the Nanoscale Using the Total Scattering Pair Distribution Function (TSPDF) Method: Recrystallization of Cryomilled Sulfamerazine. <i>Crystal Growth and Design</i> , 2013 , 13, 4239-4244	3.5	27
94	Evidence for Anomalous Bond Softening and Disorder Below 2 nm Diameter in Carbon-Supported Platinum Nanoparticles from the Temperature-Dependent Peak Width of the Atomic Pair Distribution Function. <i>Journal of Physical Chemistry C</i> , 2013 , 117, 7226-7230	3.8	18
93	Confirmation of disordered structure of ultrasmall CdSe nanoparticles from X-ray atomic pair distribution function analysis. <i>Physical Chemistry Chemical Physics</i> , 2013 , 15, 8480-6	3.6	54
92	Intermediate-range structure of self-assembled cobalt-based oxygen-evolving catalyst. <i>Journal of the American Chemical Society</i> , 2013 , 135, 6403-6	16.4	138
91	Towards a robust ad hoc data correction approach that yields reliable atomic pair distribution functions from powder diffraction data. <i>Journal of Physics Condensed Matter</i> , 2013 , 25, 454202	1.8	17
90	Local structural evidence for strong electronic correlations in spinel LiRh2O4. <i>Physical Review B</i> , 2013 , 88,	3.3	15
89	Evidence for short-range-ordered charge stripes far above the charge-ordering transition in La1.67Sr0.33NiO4. <i>Physical Review Letters</i> , 2013 , 111, 096404	7.4	25
88	Lattice dynamics reveals a local symmetry breaking in the emergent dipole phase of PbTe. <i>Physical Review B</i> , 2012 , 86,	3.3	40
87	Quantitative nanostructure characterization using atomic pair distribution functions obtained from laboratory electron microscopes. <i>Zeitschrift Fil Kristallographie</i> , 2012 , 227, 248-256		32
86	Local structure of ReO3 at ambient pressure from neutron total-scattering study. <i>Physical Review B</i> , 2012 , 86,	3.3	12
85	Understanding the Formation and Evolution of Ceria Nanoparticles Under Hydrothermal Conditions. <i>Angewandte Chemie</i> , 2012 , 124, 9164-9167	3.6	10
84	Understanding the formation and evolution of ceria nanoparticles under hydrothermal conditions. <i>Angewandte Chemie - International Edition</i> , 2012 , 51, 9030-3	16.4	78
83	Revealing the mechanisms behind SnO2 nanoparticle formation and growth during hydrothermal synthesis: an in situ total scattering study. <i>Journal of the American Chemical Society</i> , 2012 , 134, 6785-92	16.4	150
82	In-Situ Monitoring of Particle Growth at PEMFC Cathode under Accelerated Cycling Conditions. <i>Electrochemical and Solid-State Letters</i> , 2012 , 15, B72		25
81	Pair Distribution Function Technique: Principles and Methods. <i>NATO Science for Peace and Security Series B: Physics and Biophysics</i> , 2012 , 183-193	0.2	1
80	Diverse structural and magnetic properties of differently prepared MnAs nanoparticles. <i>ACS Nano</i> , 2011 , 5, 2970-8	16.7	15
79	Tuning the Surface Structure and Optical Properties of CdSe Clusters Using Coordination Chemistry. <i>Journal of Physical Chemistry Letters</i> , 2011 , 2, 3075-3080	6.4	56
78	Synthesis, crystal structure, and magnetism of Fe1.00(2)Se1.00(3) single crystals. <i>Physical Review B</i> , 2011 , 83,	3.3	55

77	Data requirements for the reliable use of atomic pair distribution functions in amorphous pharmaceutical fingerprinting. <i>Pharmaceutical Research</i> , 2011 , 28, 1041-8	4.5	60
76	Nyquist-Shannon sampling theorem applied to refinements of the atomic pair distribution function. <i>Physical Review B</i> , 2011 , 84,	3.3	46
75	Detailed mapping of the local Ir4+ dimers through the metal-insulator transitions of CuIr2S4 thiospinel by X-ray atomic pair distribution function measurements. <i>Physical Review Letters</i> , 2011 , 106, 045501	7∙4	22
74	Local structural investigation of SmFeAsOExF(x) high temperature superconductors. <i>Journal of Physics Condensed Matter</i> , 2011 , 23, 272201	1.8	7
73	Quantitative nanoparticle structures from electron crystallography data. <i>Physical Review B</i> , 2010 , 81,	3.3	5
72	Magnetic phase transition in V2O3 nanocrystals. <i>Physical Review B</i> , 2010 , 82,	3.3	18
71	Nanoscale disorder and local electronic properties of CaCu3Ti4O12: An integrated study of electron, neutron, and x-ray diffraction, x-ray absorption fine structure, and first-principles calculations. <i>Physical Review B</i> , 2010 , 81,	3.3	49
70	Advances in Scattering Probes for Materials. MRS Bulletin, 2010, 35, 495-503	3.2	4
69	Tiopronin gold nanoparticle precursor forms aurophilic ring tetramer. <i>Inorganic Chemistry</i> , 2010 , 49, 10	08 5 8-6€	j 42
68	Entropically stabilized local dipole formation in lead chalcogenides. <i>Science</i> , 2010 , 330, 1660-3	33.3	254
67	Characterisation of amorphous and nanocrystalline molecular materials by total scattering. <i>CrystEngComm</i> , 2010 , 12, 1366-1368	3.3	66
66	Seeking Supersolidity in Helium Layers. <i>Physics Magazine</i> , 2010 , 3,		
	Seeking Supersolidity in Mediam Layers. Thysics Magazine, 2010, 3,	1.1	36
65	Crystal structure solution from experimentally determined atomic pair distribution functions. Journal of Applied Crystallography, 2010 , 43, 623-629	3.8	21
65 64	Crystal structure solution from experimentally determined atomic pair distribution functions.	3.8	
	Crystal structure solution from experimentally determined atomic pair distribution functions. Journal of Applied Crystallography, 2010, 43, 623-629 Relationship between the atomic pair distribution function and small-angle scattering: implications	3.8	21
64	Crystal structure solution from experimentally determined atomic pair distribution functions. Journal of Applied Crystallography, 2010, 43, 623-629 Relationship between the atomic pair distribution function and small-angle scattering: implications for modeling of nanoparticles. Acta Crystallographica Section A: Foundations and Advances, 2009, 65, 2 Local and average structures of the spin-glass pyrochlore Y2Mo2O7 from neutron diffraction and	3.8 32-9	21
64	Crystal structure solution from experimentally determined atomic pair distribution functions. <i>Journal of Applied Crystallography</i> , 2010 , 43, 623-629 Relationship between the atomic pair distribution function and small-angle scattering: implications for modeling of nanoparticles. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2009 , 65, 2 Local and average structures of the spin-glass pyrochlore Y2Mo2O7 from neutron diffraction and neutron pair distribution function analysis. <i>Physical Review B</i> , 2009 , 79, Study of Local Structure in Selected OrganicIhorganic Perovskites in the Pm3 m Phase. <i>Chemistry</i>	3.8 32-9 3.3	21 138 42

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