Simon Billinge

List of Publications by Citations

220 papers	14,441 citations	58 h-index	116 g-index
241 ext. papers	16,486 ext. citations	8 avg, IF	6.65 L-index

#	Paper	IF	Citations
220	PDFfit2 and PDFgui: computer programs for studying nanostructure in crystals. *Journal of Physics Condensed Matter*, **2007**, 19, 335219	1.8	962
219	PDFgetX2: a GUI-driven program to obtain the pair distribution function from X-ray powder diffraction data. *Journal of Applied Crystallography*, **2004**, 37, 678-678	3.8	765
218	Structures of the ferroelectric phases of barium titanate. *The Journal of Physical Chemistry*, **1993**, 97, 2368-2377		529
217	The problem with determining atomic structure at the nanoscale. *Science*, **2007**, 316, 561-5	33.3	525
216	Rapid-acquisition pair distribution function (RA-PDF) analysis. *Journal of Applied Crystallography*, **2003**, 36, 1342-1347	3.8	433
215	Direct Observation of Lattice Polaron Formation in the Local Structure of La1-xCaxMnO3. *Physical Review Letters*, **1996**, 77, 715-718	7.4	388
214	Beyond crystallography: the study of disorder, nanocrystallinity and crystallographically challenged materials with pair distribution functions. *Chemical Communications*, **2004**, 749-60	5.8	364
213	Structure of V(2)O(5)*nH(2)O xerogel solved by the atomic pair distribution function technique. *Journal of the American Chemical Society*, **2002**, 124, 10157-62	16.4	358
212	PDFFIT, a program for full profile structural refinement of the atomic pair distribution function. *Journal of Applied Crystallography*, **1999**, 32, 572-575	3.8	357
211	Synthesis and characterization of two-dimensional Nb4C3 (MXene). *Chemical Communications*, **2014** , 50, 9517-20	5.8	321
210	Textural Mesoporosity and the Catalytic Activity of Mesoporous Molecular Sieves with Wormhole Framework Structures. *Journal of the American Chemical Society*, **1999**, 121, 8835-8842	16.4	308
209	Nature of Activated Manganese Oxide for Oxygen Evolution. *Journal of the American Chemical Society*, **2015**, 137, 14887-904	16.4	295
208	Structure of methylammonium lead iodide within mesoporous titanium dioxide: active material in high-performance perovskite solar cells. *Nano Letters*, **2014**, 14, 127-33	11.5	258
207	Entropically stabilized local dipole formation in lead chalcogenides. *Science*, **2010**, 330, 1660-3	33.3	254
206	Control of electronic properties of 2D carbides (MXenes) by manipulating their transition metal layers. *Nanoscale Horizons*, **2016**, 1, 227-234	10.8	242
205	PDFgetN: a user-friendly program to extract the total scattering structure factor and the pair distribution function from neutron powder diffraction data. *Journal of Applied Crystallography*, **2000**, 33, 1192-1192	3.8	233
204	2D molybdenum and vanadium nitrides synthesized by ammoniation of 2D transition metal carbides (MXenes). *Nanoscale*, **2017**, 9, 17722-17730	7.7	192

203	Structural analysis of complex materials using the atomic pair distribution function – a practical guide. *Zeitschrift Fur Kristallographie - Crystalline Materials*, **2003**, 218, 132-143	1	182
202	Coherent Nanotwins and Dynamic Disorder in Cesium Lead Halide Perovskite Nanocrystals. *ACS Nano*, **2017**, 11, 3819-3831	16.7	181
201	Direct Observation of Dynamic Symmetry Breaking above Room Temperature in Methylammonium Lead Iodide Perovskite. *ACS Energy Letters*, **2016**, 1, 880-887	20.1	177
200	Two-Step Nucleation and Growth of InP Quantum Dots via Magic-Sized Cluster Intermediates. *Chemistry of Materials*, **2015**, 27, 1432-1441	9.6	176
199	Quantitative size-dependent structure and strain determination of CdSe nanoparticles using atomic pair distribution function analysis. *Physical Review B*, **2007**, 76,	3.3	160
198	Complex modeling: a strategy and software program for combining multiple information sources to solve ill posed structure and nanostructure inverse problems. *Acta Crystallographica Section A: Foundations and Advances*, **2015**, 71, 562-8	1.7	152
197	Revealing the mechanisms behind SnO2 nanoparticle formation and growth during hydrothermal synthesis: an in situ total scattering study. *Journal of the American Chemical Society*, **2012**, 134, 6785-92	16.4	150
196	Imaging Dirac-mass disorder from magnetic dopant atoms in the ferromagnetic topological insulator Crx(Bi0.1Sb0.9)2-xTe3. *Proceedings of the National Academy of Sciences of the United States of America*, **2015**, 112, 1316-21	11.5	147
195	Ab initio determination of solid-state nanostructure. *Nature*, **2006**, 440, 655-8	50.4	146
194	Atomic structures and gram scale synthesis of three tetrahedral quantum dots. *Journal of the American Chemical Society*, **2014**, 136, 10645-53	16.4	139
193	Structure of nanocrystalline materials using atomic pair distribution function analysis: Study of LiMoS2. *Physical Review B*, **2002**, 65,	3.3	139
192	Intermediate-range structure of self-assembled cobalt-based oxygen-evolving catalyst. *Journal of the American Chemical Society*, **2013**, 135, 6403-6	16.4	138
191	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, 232-9		138
190	Polyhedral units and network connectivity in calcium aluminosilicate glasses from high-energy X-Ray diffraction. *Physical Review Letters*, **2000**, 85, 3436-9	7.4	138
189	Neutron diffraction evidence of microscopic charge inhomogeneities in the CuO2 plane of superconducting La2-xSrxCuO4 (0. *Physical Review Letters*, **2000**, 84, 5856-9	7.4	138
188	Orbital correlations in the pseudocubic O and rhombohedral R phases of LaMnO3. *Physical Review Letters*, **2005**, 94, 177203	7.4	135
187	Structure of nanocrystalline Ti3C2 MXene using atomic pair distribution function. *Physical Review Letters*, **2014**, 112, 125501	7.4	129
186	Atomic electron tomography: 3D structures without crystals. *Science*, **2016**, 353,	33.3	129

185	Polymorphism in magic-sized Au144(SR)60 clusters. *Nature Communications*, **2016**, 7, 11859	17.4	126
184	Mercury binding sites in thiol-functionalized mesostructured silica. *Journal of the American Chemical Society*, **2005**, 127, 8492-8	16.4	118
183	Measuring Correlated Atomic Motion Using X-ray Diffraction. *Journal of Physical Chemistry A*, **1999**, 103, 921-924	2.8	114
182	High Real-Space Resolution Measurement of the Local Structure of Ga1□InxAs Using X-Ray Diffraction. *Physical Review Letters*, **1999**, 83, 4089-4092	7.4	113
181	Lattice dynamics and correlated atomic motion from the atomic pair distribution function. *Physical Review B*, **2003**, 67,	3.3	105
180	Local structure of nanoporous carbons. *The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties*, **1999**, 79, 1519-1530		96
179	Fine-Scale Nanostructure in □Al2O3. *Chemistry of Materials*, **2006**, 18, 3242-3248	9.6	88
178	Improved measures of quality for the atomic pair distribution function. *Journal of Applied Crystallography*, **2003**, 36, 53-64	3.8	84
177	Role of framework sodium versus local framework structure in determining the hydrothermal stability of MCM-41 mesostructures. *Journal of the American Chemical Society*, **2002**, 124, 97-103	16.4	82
176	Application of Atomic Pair Distribution Function Analysis to Materials with Intrinsic Disorder. Three-Dimensional Structure of Exfoliated-Restacked WS2: Not Just a Random Turbostratic Assembly of Layers. *Journal of the American Chemical Society*, **2000**, 122, 11571-11576	16.4	82
175	Understanding the formation and evolution of ceria nanoparticles under hydrothermal conditions. *Angewandte Chemie - International Edition*, **2012**, 51, 9030-3	16.4	78
174	Reciprocal-space instrumental effects on the real-space neutron atomic pair distribution function. *Journal of Applied Crystallography*, **2004**, 37, 110-116	3.8	78
173	Simulation of nanoporous carbons: A chemically constrained structure. *The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties*, **1999**, 79, 1499-1518		76
172	Local vibrations and negative thermal expansion in ZrW2O8. *Physical Review Letters*, **2014**, 112, 045505	7.4	75
171	Pair distribution function computed tomography. *Nature Communications*, **2013**, 4, 2536	17.4	75
170	Characterisation of amorphous and nanocrystalline molecular materials by total scattering. *CrystEngComm*, **2010**, 12, 1366-1368	3.3	66
169	Structure of Sulfate Adsorption Complexes on Ferrihydrite. *Environmental Science and Technology Letters*, **2014**, 1, 97-101	11	65
168	Mechanisms for iron oxide formation under hydrothermal conditions: an in situ total scattering study. *ACS Nano*, **2014**, 8, 10704-14	16.7	65

167	Signatures of the topological s superconducting order parameter in the type-II Weyl semimetal T -MoTe. *Nature Communications*, **2017**, 8, 1082	17.4	62
166	Correlating Size and Composition-Dependent Effects with Magnetic, Mössbauer, and Pair Distribution Function Measurements in a Family of Catalytically Active Ferrite Nanoparticles. *Chemistry of Materials*, **2015**, 27, 3572-3592	9.6	61
165	Data requirements for the reliable use of atomic pair distribution functions in amorphous pharmaceutical fingerprinting. *Pharmaceutical Research*, **2011**, 28, 1041-8	4.5	60
164	Rapid desolvation-triggered domino lattice rearrangement in a metal-organic framework. *Nature Chemistry*, **2020**, 12, 90-97	17.6	60
163	Two-Dimensional Arrays of Transition Metal Nitride Nanocrystals. *Advanced Materials*, **2019**, 31, e1902393	24	59
162	Local off-centering symmetry breaking in the high-temperature regime of SnTe. *Physical Review B*, **2014**, 89,	3.3	58
161	Structure of intercalated Cs in zeolite ITQ-4: an array of metal ions and correlated electrons confined in a pseudo-1D nanoporous host. *Physical Review Letters*, **2002**, 89, 075502	7.4	58
160	Study of Local Structure in Selected Organic–Inorganic Perovskites in the Pm3 m Phase. *Chemistry of Materials*, **2008**, 20, 1272-1277	9.6	57
159	Tuning the Surface Structure and Optical Properties of CdSe Clusters Using Coordination Chemistry. *Journal of Physical Chemistry Letters*, **2011**, 2, 3075-3080	6.4	56
158	Search for a structural response to the intermediate phase in GexSe1−x glasses. *Physical Review B*, **2008**, 77,	3.3	56
157	Magnetism in semiconducting molybdenum dichalcogenides. *Science Advances*, **2018**, 4, eaat3672	14.3	56
156	Synthesis, crystal structure, and magnetism of Fe1.00(2)Se1.00(3) single crystals. *Physical Review B*, **2011**, 83,	3.3	55
155	Confirmation of disordered structure of ultrasmall CdSe nanoparticles from X-ray atomic pair distribution function analysis. *Physical Chemistry Chemical Physics*, **2013**, 15, 8480-6	3.6	54
154	Probing local and long-range structure simultaneously: an in situ study of the high-temperature phase transition of alpha-AlF3. *Journal of the American Chemical Society*, **2004**, 126, 4756-7	16.4	50
153	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, 20180413	3	49
152	Cation Exchange Induced Transformation of InP Magic-Sized Clusters. *Chemistry of Materials*, **2017**, 29, 7984-7992	9.6	49
151	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. *Physical Review B*, **2010**, 81,	3.3	49
150	The atomic pair distribution function: past and present. *Zeitschrift Fur Kristallographie - Crystalline Materials*, **2004**, 219,	1	49

149	Nyquist-Shannon sampling theorem applied to refinements of the atomic pair distribution function. *Physical Review B*, **2011**, 84,	3.3	46
148	Tiopronin gold nanoparticle precursor forms aurophilic ring tetramer. *Inorganic Chemistry*, **2010**, 49, 10858-66	5.1	42
147	Local and average structures of the spin-glass pyrochlore Y2Mo2O7 from neutron diffraction and neutron pair distribution function analysis. *Physical Review B*, **2009**, 79,	3.3	42
146	Superconducting order from disorder in 2H-TaSe 2−x S x. *Npj Quantum Materials*, **2017**, 2,	5	41
145	Demonstration of thin film pair distribution function analysis (tfPDF) for the study of local structure in amorphous and crystalline thin films. *IUCrJ*, **2015**, 2, 481-9	4.7	41
144	Early stage structural development of prototypical zeolitic imidazolate framework (ZIF) in solution. *Nanoscale*, **2018**, 10, 4291-4300	7.7	40
143	Lattice dynamics reveals a local symmetry breaking in the emergent dipole phase of PbTe. *Physical Review B*, **2012**, 86,	3.3	40
142	X-Ray Diffraction Computed Tomography for Structural Analysis of Electrode Materials in Batteries. *Journal of the Electrochemical Society*, **2015**, 162, A1310-A1314	3.9	39
141	Evolution of atomic structure during nanoparticle formation. *IUCrJ*, **2014**, 1, 165-71	4.7	39
140	Structures of alkali metals in silica gel nanopores: new materials for chemical reductions and hydrogen production. *Journal of the American Chemical Society*, **2007**, 129, 1386-92	16.4	39
139	Charge Density Wave Caused by Reducing ThSe3 by One Electron. Superstructure and Short-Range Order in ATh2Se6 (A = K, Rb) Studied by X-ray Diffraction, Electron Diffraction, and Diffuse Scattering. *Journal of the American Chemical Society*, **1998**, 120, 10706-10714	16.4	38
138	Scalable Synthesis of Ultrathin Mn3N2 Exhibiting Room-Temperature Antiferromagnetism. *Advanced Functional Materials*, **2019**, 29, 1809001	15.6	37
137	Modelling pair distribution functions (PDFs) of organic compounds: describing both intra- and intermolecular correlation functions in calculated PDFs. *Journal of Applied Crystallography*, **2015**, 48, 171-178	3.8	37
136	Seeking Supersolidity in Helium Layers. *Physics Magazine*, **2010**, 3,	1.1	36
135	Mesostructured Non-Oxidic Solids with Adjustable Worm-hole Shaped Pores: M-Ge-Q (Q=S, Se) Frameworks Based on Tetrahedral [Ge4Q10]4− Clusters. *Advanced Materials*, **2000**, 12, 85-91	24	36
134	Neutron Scattering Studies of Compositional Heterogeneity in Sol−Gel Processed Lead Zirconate Titanates. *Chemistry of Materials*, **1998**, 10, 3611-3619	9.6	36
133	Magnetic pair distribution function analysis of local magnetic correlations. *Acta Crystallographica Section A: Foundations and Advances*, **2014**, 70, 3-11	1.7	35
132	Assigned and unassigned distance geometry: applications to biological molecules and nanostructures. *4or*, **2016**, 14, 337-376	1.4	34

131	Quantitative nanostructure characterization using atomic pair distribution functions obtained from laboratory electron microscopes. *Zeitschrift F Kristallographie*, **2012**, 227, 248-256		32
130	Proton Electron Conductivity in Thin Films of a Cobalt Oxygen Evolving Catalyst. *ACS Applied Energy Materials*, **2019**, 2, 3-12	6.1	30
129	Detection and characterization of nanoparticles in suspension at low concentrations using the X-ray total scattering pair distribution function technique. *Nanoscale*, **2015**, 7, 5480-7	7.7	30
128	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	11.5	29
127	Magnetic structure determination from the magnetic pair distribution function (mPDF): ground state of MnO. *Acta Crystallographica Section A: Foundations and Advances*, **2015**, 71, 325-34	1.7	29
126	Symmetry Breaking in Nanostructure Development of Carbogenic Molecular Sieves: Effects of Morphological Pattern Formation on Oxygen and Nitrogen Transport. *Chemistry of Materials*, **1996**, 8, 2159-2171	9.6	29
125	Towards atomistic understanding of polymorphism in the solvothermal synthesis of ZrO nanoparticles. *Acta Crystallographica Section A: Foundations and Advances*, **2016**, 72, 645-650	1.7	28
124	Nanoscale structural domains in the phonon-glass thermoelectric material Zn4Sb3. *Physical Review B*, **2007**, 75,	3.3	28
123	Improved Models for Metallic Nanoparticle Cores from Atomic Pair Distribution Function (PDF) Analysis. *Journal of Physical Chemistry C*, **2018**, 122, 29498-29506	3.8	28
122	Toward Phase Quantification at the Nanoscale Using the Total Scattering Pair Distribution Function (TSPDF) Method: Recrystallization of Cryomilled Sulfamerazine. *Crystal Growth and Design*, **2013**, 13, 4239-4244	3.5	27
121	Hollandites as a new class of multiferroics. *Scientific Reports*, **2014**, 4, 6203	4.9	26
120	Using a machine learning approach to determine the space group of a structure from the atomic pair distribution function. *Acta Crystallographica Section A: Foundations and Advances*, **2019**, 75, 633-643	1.7	26
119	Autonomous experimentation systems for materials development: A community perspective. *Matter*, **2021**, 4, 2702-2726	12.7	26
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117	Reconciliation of local and long-range tilt correlations in underdoped La2BaxCuO4(0.155). *Physical Review B*, **2015**, 91,	3.3	25
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115	In-Situ Monitoring of Particle Growth at PEMFC Cathode under Accelerated Cycling Conditions. *Electrochemical and Solid-State Letters*, **2012**, 15, B72		25
114	Correlated local dipoles in PbTe. *Physical Review Materials*, **2018**, 2,	3.2	25

113	Local Environment of Terbium(III) Ions in Layered Nanocrystalline Zirconium(IV) Phosphonate-Phosphate Ion Exchange Materials. *Inorganic Chemistry*, **2017**, 56, 8837-8846	5.1	24
112	Verification of Anderson Superexchange in MnO via Magnetic Pair Distribution Function Analysis and ab□nitio Theory. *Physical Review Letters*, **2016**, 116, 197204	7.4	23
111	Bulk metallic glass-like scattering signal in small metallic nanoparticles. *ACS Nano*, **2014**, 8, 6163-70	16.7	23
110	Cu(x)UTe3: stabilization of UTe3 in the ZrSe3 structure type via copper insertion. The artifact of Te-Te chains and evidence for distortions due to long range modulations. *Journal of the American Chemical Society*, **2001**, 123, 4755-62	16.4	23
109	Local atomic and magnetic structure of dilute magnetic semiconductor (Ba,K)(Zn,Mn)2As2. *Physical Review B*, **2016**, 94,	3.3	23
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107	Intra-unit-cell nematic charge order in the titanium-oxypnictide family of superconductors. *Nature Communications*, **2014**, 5, 5761	17.4	22
106	Detailed mapping of the local Ir4+ dimers through the metal-insulator transitions of CuIr2S4 thiospinel by X-ray atomic pair distribution function measurements. *Physical Review Letters*, **2011**, 106, 045501	7.4	22
105	Accurate Structure Determination of Mo6SyIz Nanowires from Atomic Pair Distribution Function (PDF) Analysis. *Chemistry of Materials*, **2006**, 18, 100-106	9.6	22
104	Volume-wise destruction of the antiferromagnetic Mott insulating state through quantum tuning. *Nature Communications*, **2016**, 7, 12519	17.4	22
103	Local orbital degeneracy lifting as a precursor to an orbital-selective Peierls transition. *Nature Communications*, **2019**, 10, 3638	17.4	21
102	Crystal structure solution from experimentally determined atomic pair distribution functions. *Journal of Applied Crystallography*, **2010**, 43, 623-629	3.8	21
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99	Hollow organic capsules assemble into cellular semiconductors. *Nature Communications*, **2018**, 9, 1957	17.4	20
98	Room temperature local nematicity in FeSe superconductor. *Physical Review B*, **2019**, 100,	3.3	19
97	Recent advances and applications of deep learning methods in materials science. *Npj Computational Materials*, **2022**, 8,	10.9	19
96	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. *Journal of Physical Chemistry C*, **2013**, 117, 7226-7230	3.8	18

95	Magnetic phase transition in V2O3 nanocrystals. *Physical Review B*, **2010**, 82,	3.3	18
94	Determination of standard uncertainties in fits to pair distribution functions. *Acta Crystallographica Section A: Foundations and Advances*, **2004**, 60, 315-7		18
93	Calibration and data collection protocols for reliable lattice parameter values in electron pair distribution function studies. *Journal of Applied Crystallography*, **2015**, 48, 244-251	3.8	17
92	Towards solution and refinement of organic crystal structures by fitting to the atomic pair distribution function. *Acta Crystallographica Section A: Foundations and Advances*, **2016**, 72, 62-72	1.7	17
91	Towards a robust ad hoc data correction approach that yields reliable atomic pair distribution functions from powder diffraction data. *Journal of Physics Condensed Matter*, **2013**, 25, 454202	1.8	17
90	Planar nets of Ti atoms comprising squares and rhombs in the new binary antimonide Ti2Sb. *Journal of the American Chemical Society*, **2004**, 126, 8295-302	16.4	17
89	Cluster-mining: an approach for determining core structures of metallic nanoparticles from atomic pair distribution function data. *Acta Crystallographica Section A: Foundations and Advances*, **2020**, 76, 24-31	1.7	17
88	Synthesis and Properties of Plasmonic Boron-Hyperdoped Silicon Nanoparticles. *Advanced Functional Materials*, **2019**, 29, 1807788	15.6	17
87	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. *Journal of the American Chemical Society*, **2019**, 141, 431-440	16.4	17
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80	Unlocking the structure of mixed amorphous-crystalline ceramic oxide films synthesized under low temperature electromagnetic excitation. *Journal of Materials Chemistry A*, **2017**, 5, 18434-18441	13	16
79	LATTICE EFFECTS IN HIGH-TC SUPERCONDUCTORS **1996**, 265-373		16
78	Celebrating 100 years of the Debye scattering equation. *Acta Crystallographica Section A: Foundations and Advances*, **2016**, 72, 589-590	1.7	16

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73	Total-scattering pair-distribution function of organic material from powder electron diffraction data. *Microscopy and Microanalysis*, **2015**, 21, 459-71	0.5	13
72	The Liga algorithm for ab initio determination of nanostructure. *Acta Crystallographica Section A: Foundations and Advances*, **2008**, 64, 631-40		13
71	PDF from X-ray powder diffraction for nanometer-scale atomic structure analysis of quasicrystalline alloys. *Zeitschrift Fur Kristallographie - Crystalline Materials*, **2005**, 220, 962-967	1	13
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69	Cu(Ir□ xCrx)S□a model system for studying nanoscale phase coexistence at the metal-insulator transition. *Scientific Reports*, **2014**, 4, 4081	4.9	12
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67	Structural compliance, misfit strain, and stripe nanostructures in cuprate superconductors. *Physical Review B*, **2002**, 66,	3.3	11
66	Structural Analysis of Molecular Materials Using the Pair Distribution Function. *Chemical Reviews*, **2021**,	68.1	11
65	Photoinduced dynamics of nematic order parameter in FeSe. *Physical Review B*, **2019**, 99,	3.3	10
64	Enhanced thermoelectric power and electronic correlations in RuSe2. *APL Materials*, **2015**, 3, 041513	5.7	10
63	Understanding electronic peculiarities in tetragonal FeSe as local structural symmetry breaking. *Physical Review B*, **2020**, 102,	3.3	10
62	Understanding the Formation and Evolution of Ceria Nanoparticles Under Hydrothermal Conditions. *Angewandte Chemie*, **2012**, 124, 9164-9167	3.6	10
61	Quantitative Structural Characterization of Catalytically Active TiO2 Nanoparticles. *ACS Applied Nano Materials*, **2019**, 2, 6268-6276	5.6	9
60	Algorithm for systematic peak extraction from atomic pair distribution functions. *Acta Crystallographica Section A: Foundations and Advances*, **2015**, 71, 392-409	1.7	9

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56	A cloud platform for atomic pair distribution function analysis: PDFitc. *Acta Crystallographica Section A: Foundations and Advances*, **2021**, 77, 2-6	1.7	9
55	Nanometre-scale structure from powder diffraction: total scattering and atomic pair distribution function analysis **2019**, 649-672		8
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