

Kevin Knight

List of Publications by Year in Descending Order

Source: <https://exaly.com/author-pdf/4706130/kevin-knight-publications-by-year.pdf>

Version: 2024-04-28

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

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

271
papers

7,736
citations

47
h-index

73
g-index

279
ext. papers

8,408
ext. citations

3.9
avg, IF

6.03
L-index

#	Paper	IF	Citations
271	Comprehensive determination of the high-pressure structural behaviour of BaTiO ₃ . <i>Materials Advances</i> , 2021 , 2, 6094-6103	3.3	1
270	Nuclear and magnetic structures of KMnF ₃ perovskite in the temperature interval 10 K–105 K. <i>Journal of Alloys and Compounds</i> , 2020 , 842, 155935	5.7	4
269	Low-temperature thermophysical and crystallographic properties of BaZrO ₃ perovskite. <i>Journal of Materials Science</i> , 2020 , 55, 6417-6428	4.3	5
268	Crystal and Electronic Structures of ANaIO Periodate Double Perovskites (A = Sr, Ca, Ba): Candidate Wasteforms for I-129 Immobilization. <i>Inorganic Chemistry</i> , 2020 , 59, 18407-18419	5.1	3
267	Two-dimensional spin liquid behaviour in the triangular-honeycomb antiferromagnet TbInO ₃ . <i>Nature Physics</i> , 2019 , 15, 262-268	16.2	27
266	High-Pressure Study of the Elpasolite Perovskite LaNiMnO. <i>Inorganic Chemistry</i> , 2019 , 58, 9016-9027	5.1	3
265	First-order valence transition: Neutron diffraction, inelastic neutron scattering, and x-ray absorption investigations on the double perovskite Ba ₂ PrRu _{0.9} Ir _{0.1} O ₆ . <i>Physical Review B</i> , 2019 , 99,	3.3	3
264	Crystal structures and electronic properties in 3d transition metal doped SrRuO. <i>Dalton Transactions</i> , 2019 , 48, 4730-4741	4.3	5
263	Defining an aristotype crystal structure and crystallographic distortions in leucite/pollucite-structured phases with space group ($\overline{3}d$). <i>Physics and Chemistry of Minerals</i> , 2019 , 46, 595-605	1.6	3
262	Parameterization of the crystal structure of garnet in terms of symmetry-adapted basis-vectors of the ideal tetrahedron and octahedron: Application to the pressure-dependence of the crystal structure of Y ₃ Al ₅ O ₁₂ between 0 and 126 GPa. <i>Materials Chemistry and Physics</i> , 2019 , 227, 72-82	4.4	2
261	The thermal expansion properties of halogen bond containing 1,4 dioxane halogen complexes. <i>CrystEngComm</i> , 2019 , 21, 5269-5277	3.3	1
260	Investigation of the changes in hydrogen bonding accompanying the structural reorganization at 103 K in ammonium iodate. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2019 , 75, 152-159	1.8	
259	Structure and physical properties of SeCo Mn O. <i>Journal of Physics Condensed Matter</i> , 2019 , 31, 395402	1.8	2
258	Synchrotron X-ray and neutron investigation of the structure and thermal expansion of the monoclinic Al ₁₃ Cr ₂ phase. <i>Journal of Alloys and Compounds</i> , 2019 , 781, 1198-1208	5.7	4
257	Crystal structures of NiSO ₄ ·9H ₂ O and NiSO ₄ ·8H ₂ O: magnetic properties, stability with respect to morenosite (NiSO ₄ ·7H ₂ O), the solid-solution series (Mg _x Ni _{1-x})SO ₄ ·9H ₂ O. <i>Physics and Chemistry of Minerals</i> , 2018 , 45, 695-712	1.6	3
256	A high-resolution neutron powder diffraction study of the low-temperature structural phase transitions in RbCaF ₃ perovskite. <i>Journal of Solid State Chemistry</i> , 2018 , 263, 172-181	3.3	4
255	The crystal structure of lueshite at 298 K resolved by high-resolution time-of-flight neutron powder diffraction. <i>Physics and Chemistry of Minerals</i> , 2018 , 45, 77-83	1.6	4

254	Investigation into the dehydration of selenate doped $\text{Na}_2\text{M}(\text{SO}_4)_2 \cdot 2\text{H}_2\text{O}$ (M = Mn, Fe, Co and Ni): Stabilisation of the high Na content alluaudite phases $\text{Na}_3\text{M}_{1.5}(\text{SO}_4)_3 \cdot 1.5\text{x}(\text{SeO}_4)$ (M = Mn, Co and Ni) through selenate incorporation. <i>Journal of Solid State Chemistry</i> , 2018 , 258, 64-71	3.3	17
253	Temperature-induced polymorphism in methyl stearate. <i>CrystEngComm</i> , 2018 , 20, 6885-6893	3.3	7
252	Negative 2D thermal expansion in the halogen bonded acetone bromine complex. <i>CrystEngComm</i> , 2018 , 20, 3246-3250	3.3	5
251	Thermal expansion of deuterated monoclinic natrojarosite; a combined neutron synchrotron powder diffraction study. <i>Journal of Applied Crystallography</i> , 2017 , 50, 340-348	3.8	
250	S=12 quantum critical spin ladders produced by orbital ordering in $\text{Ba}_2\text{CuTeO}_6$. <i>Physical Review B</i> , 2017 , 95,	3.3	8
249	High-resolution neutron-diffraction measurements to 8 kbar. <i>High Pressure Research</i> , 2017 , 37, 486-494	1.6	1
248	Structure, thermal expansion and incompressibility of $\text{MgSO}_4 \cdot 9\text{H}_2\text{O}$, its relationship to meridianiite ($\text{MgSO}_4 \cdot 11\text{H}_2\text{O}$) and possible natural occurrences. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2017 , 73, 47-64	1.8	13
247	Cation disorder and phase transitions in the structurally complex solar cell material $\text{Cu}_2\text{ZnSnS}_4$. <i>Journal of Materials Chemistry A</i> , 2017 , 5, 16672-16680	13	34
246	Variable stoichiometry in tectosilicates having the leucite/pollucite-type structure with particular emphasis on modelling the interframework cavity cation environment. <i>Journal of Solid State Chemistry</i> , 2017 , 251, 90-104	3.3	6
245	Structural organization in the trimethylamine iodine monochloride complex. <i>CrystEngComm</i> , 2017 , 19, 5194-5201	3.3	3
244	Low-temperature structure and the ferroelectric phase transitions in the CdTiO_3 perovskite. <i>Physical Review B</i> , 2017 , 96,	3.3	10
243	Low temperature, high pressure thermo-physical and crystallographic properties of KZnF_3 perovskite. <i>Materials Chemistry and Physics</i> , 2017 , 199, 393-407	4.4	3
242	High-pressure thermoelastic and structural properties of KCaF_3 perovskite in the low temperature Pbnm phase. <i>Journal of Alloys and Compounds</i> , 2017 , 693, 1305-1314	5.7	5
241	Phase Transition Behavior of the Layered Perovskite $\text{CsBi}_{0.6}\text{La}_{0.4}\text{Nb}_2\text{O}_7$: A Hybrid Improper Ferroelectric. <i>Crystals</i> , 2017 , 7, 135	2.3	8
240	Magnetic and structural phase diagram of the solid solution $\text{LaCo}_x\text{Mn}_{1-x}\text{O}_3$. <i>Physical Review B</i> , 2016 , 94,	3.3	14
239	Low temperature and high pressure thermoelastic and crystallographic properties of SrZrO_3 perovskite in the Pbnm phase. <i>Solid State Sciences</i> , 2016 , 62, 90-104	3.4	5
238	Orbital frustration in the S = $\frac{1}{2}$ kagome magnet vesignieite, $\text{BaCu}_3\text{V}_2\text{O}_8(\text{OH})_2$. <i>Journal of Materials Chemistry C</i> , 2016 , 4, 10315-10322	7.1	15
237	Low-temperature structural behaviour of LaCoO_3 – a high-resolution neutron study. <i>Solid State Sciences</i> , 2016 , 57, 38-43	3.4	9

236	Neutron diffraction and multinuclear solid state NMR investigation into the structures of oxide ion conducting $\text{La}_{9.6}\text{Si}_6\text{O}_{26.4}$ and $\text{La}_8\text{Sr}_2\text{Si}_6\text{O}_{26}$, and their hydrated phases. <i>Dalton Transactions</i> , 2016 , 45, 121-33	4.3	8
235	X-ray and neutron powder diffraction analyses of $\text{Gly}[\text{MgSO}_4]_n\text{H}_2\text{O}$ and $\text{Gly}[\text{MgSO}_4]_n\text{D}_2\text{O}$, and their deuterated counterparts. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2016 , 72, 203-16	0.8	1
234	Synthesis, structural characterisation and proton conduction of two new hydrated phases of barium ferrite $\text{BaFeO}_{2.5x}(\text{OH})_{2x}$. <i>Journal of Materials Chemistry A</i> , 2016 , 4, 3415-3430	13	15
233	Phase separation in NaTaO_3 . Impact of temperature and doping. <i>Solid State Sciences</i> , 2016 , 52, 149-153	3.4	10
232	The impact of room temperature polymorphism in K doped NaTaO_3 on structural phase transition behaviour. <i>Journal of Solid State Chemistry</i> , 2016 , 238, 109-112	3.3	7
231	Thermal evolution of the crystal structure of the orthorhombic perovskite LaFeO_3 . <i>Journal of Solid State Chemistry</i> , 2015 , 230, 337-342	3.3	27
230	Structural and dielectric studies of the phase behaviour of the topological ferroelectric $\text{La}_{1-x}\text{Nd}_x\text{TaO}_4$. <i>Dalton Transactions</i> , 2015 , 44, 10673-80	4.3	25
229	A method for the monitoring of metal recrystallization based on the in-situ measurement of the elastic energy release using neutron diffraction. <i>Review of Scientific Instruments</i> , 2015 , 86, 053901	1.7	6
228	Low temperature thermoelastic properties of galena in a simple, self-consistent, two-term Debye model. <i>Physics and Chemistry of Minerals</i> , 2015 , 42, 235-242	1.6	2
227	Phase coexistence in NaTaO_3 at room temperature; a high resolution neutron powder diffraction study. <i>Solid State Sciences</i> , 2015 , 43, 15-21	3.4	15
226	New insights into the phase diagram of a magnetic perovskite, $\text{LaCo}_{1-x}\text{Mn}_x\text{O}_3$. <i>Journal of Physics Condensed Matter</i> , 2015 , 27, 165401	1.8	4
225	Low temperature structural studies of SrSnO_3 . <i>Journal of Physics Condensed Matter</i> , 2015 , 27, 365401	1.8	20
224	High-temperature structural phase transitions in neighborite: a high-resolution neutron powder diffraction investigation. <i>Physics and Chemistry of Minerals</i> , 2015 , 42, 45-52	1.6	6
223	Reply to structural and magnetic behavior of the cubic oxyfluoride SrFeO_2F studied by neutron diffraction. <i>Journal of Solid State Chemistry</i> , 2015 , 226, 326-331	3.3	10
222	Lithium insertion properties of $\text{LiTiNb}_2\text{O}_7$ investigated by neutron diffraction and first-principles modelling. <i>Journal of Solid State Chemistry</i> , 2015 , 229, 19-25	3.3	26
221	Thermoelastic and structural properties of ionically conducting cerate perovskites: (II) SrCeO_3 between 1273 K and 1723 K. <i>Dalton Transactions</i> , 2015 , 44, 10773-84	4.3	4
220	Observations on the crystal structures of lueshite. <i>Physics and Chemistry of Minerals</i> , 2014 , 41, 393-401	1.6	8
219	Introducing a large polar tetragonal distortion into Ba-doped BiFeO_3 by low-temperature fluorination. <i>Inorganic Chemistry</i> , 2014 , 53, 12572-83	5.1	21

218	Cobalt adipate, Co(C ₆ H ₈ O ₄): antiferromagnetic structure, unusual thermal expansion and magnetoelastic coupling. <i>Materials Horizons</i> , 2014 , 1, 332-337	14.4	20
217	Colossal thermal expansion and negative thermal expansion in simple halogen bonded complexes. <i>CrystEngComm</i> , 2014 , 16, 237-243	3.3	25
216	Tuning the giant magnetoelastic transition in Ba ₃ BiRu ₂ O ₉ and Ba ₃ BiRu ₂ O ₉ . <i>Journal of Physics Condensed Matter</i> , 2014 , 26, 276003	1.8	4
215	A high-pressure neutron diffraction study of the ferroelastic phase transition in RbCaF ₃ . <i>Physics and Chemistry of Minerals</i> , 2014 , 41, 461-472	1.6	9
214	From spin glass to quantum spin liquid ground states in molybdate pyrochlores. <i>Physical Review Letters</i> , 2014 , 113, 117201	7.4	30
213	Crystallographic and magnetic structure of the perovskite-type compound BaFeO _{2.5} : unrivaled complexity in oxygen vacancy ordering. <i>Inorganic Chemistry</i> , 2014 , 53, 5911-21	5.1	36
212	Equation of state and a high-pressure structural phase transition in the gillespite-structured phase Ba _{0.5} Sr _{0.5} CuSi ₄ O ₁₀ . <i>European Journal of Mineralogy</i> , 2014 , 25, 909-917	2.2	2
211	A high-resolution neutron powder diffraction investigation of galena (PbS) between 10 K and 350 K: no evidence for anomalies in the lattice parameters or atomic displacement parameters in galena or altaite (PbTe) at temperatures corresponding to the saturation of cation disorder. <i>Journal of Physics Condensed Matter</i> , 2014 , 26, 385403	1.8	15
210	Time-of-flight neutron powder diffraction with milligram samples: the crystal structures of NaCoF ₃ and NaNiF ₃ post-perovskites. <i>Journal of Applied Crystallography</i> , 2014 , 47, 1939-1947	3.8	6
209	A high-resolution powder neutron diffraction study of the crystal structure of neighborite (NaMgF ₃) between 9 and 440 K. <i>American Mineralogist</i> , 2014 , 99, 824-838	2.9	9
208	Substitution of Ti ³⁺ and Ti ⁴⁺ in hibonite (CaAl ₁₂ O ₁₉). <i>American Mineralogist</i> , 2014 , 99, 1369-1382	2.9	23
207	The temperature dependence of the volume expansivity and the thermal expansion tensor of petalite between 4.2 K and 600 K. <i>Journal of Mineralogical and Petrological Sciences</i> , 2014 , 109, 118-124	0.9	5
206	A neutron diffraction study and mode analysis of compounds of the system La _{1-x} Sr _x FeO _{3-x/2} (x=1, 0.8, 0.5, 0.2) and an investigation of their magnetic properties. <i>Journal of Solid State Chemistry</i> , 2013 , 206, 158-169	3.3	33
205	The competition between halogen bonds (Br⋯O) and C-H⋯O hydrogen bonds: the structure of the acetone-bromine complex revisited. <i>CrystEngComm</i> , 2013 , 15, 8572	3.3	15
204	High-temperature phases of multiferroic BiFe _{0.7} Mn _{0.3} O ₃ . <i>Physical Review B</i> , 2013 , 87,	3.3	1
203	Synthesis, structural and magnetic characterisation of the fluorinated compound 15R-BaFeO ₂ F. <i>Journal of Solid State Chemistry</i> , 2013 , 203, 218-226	3.3	20
202	High-temperature order-disorder transitions in the skutterudites CoGe _{1.5} Q _{1.5} (Q=S, Te). <i>Journal of Solid State Chemistry</i> , 2013 , 198, 525-531	3.3	5
201	Thermally Robust Anion-Chain Order in Oxynitride Perovskites. <i>Chemistry of Materials</i> , 2013 , 25, 5004-5016	3.6	60

200	Facile proton conduction in H ⁺ /Li ⁺ ion-exchanged garnet-type fast Li-ion conducting Li ₅ La ₃ Nb ₂ O ₁₂ . <i>Journal of Materials Chemistry A</i> , 2013 , 1, 13469	13	49
199	Synthesis, conductivity and structural aspects of Nd ₃ Zr ₂ Li ₇ B _x Al _x O ₁₂ . <i>Journal of Materials Chemistry A</i> , 2013 , 1, 14013	13	19
198	Neutron diffraction in situ monitoring of the dislocation density during martensitic transformation in a stainless steel. <i>Scripta Materialia</i> , 2013 , 68, 506-509	5.6	49
197	Thermoelastic and structural properties of ionically conducting cerate perovskites: (I) BaCeO ₃ at low temperature in the Pbnm phase. <i>Solid State Ionics</i> , 2013 , 232, 112-122	3.3	14
196	Magnetoelastic coupling and competing entropy changes in substituted CoMnSi metamagnets. <i>Physical Review B</i> , 2013 , 87,	3.3	32
195	On the soft magnetic properties of the compounds of the series Na _(x) Mn _(4.5-x/2) (VO ₄) ₃ and the magnetic structure of h.t.-Mn ₃ (VO ₄) ₂ (x = 1). <i>Dalton Transactions</i> , 2013 , 42, 7894-900	4.3	4
194	A comparison of dilatometry and in-situ neutron diffraction in tracking bulk phase transformations in a martensitic stainless steel. <i>Materials Characterization</i> , 2013 , 82, 50-57	3.9	27
193	Monitoring in situ stress/strain behaviour during plastic yielding in polymineralic rocks using neutron diffraction. <i>Journal of Structural Geology</i> , 2013 , 47, 36-51	3	5
192	Ferroelectricity and lattice distortion associated with spin orderings in a multiferroic delafossite AgFeO ₂ . <i>EPJ Web of Conferences</i> , 2013 , 40, 15008	0.3	6
191	Effect of Ga incorporation on the structure and Li ion conductivity of La ₃ Zr ₂ Li ₇ O ₁₂ . <i>Dalton Transactions</i> , 2012 , 41, 12048-53	4.3	76
190	Giant magnetoelastic effect at the opening of a spin-gap in Ba ₃ BiIr ₂ O ₉ . <i>Journal of the American Chemical Society</i> , 2012 , 134, 3265-70	16.4	32
189	Synthesis, characterization and physical properties of the skutterudites Y _b Fe ₂ Ni ₂ Sb ₁₂ (0 ≤ b ≤ 4). <i>Journal of Solid State Chemistry</i> , 2012 , 193, 36-41	3.3	15
188	Low temperature thermoelastic and structural properties of LaGaO ₃ perovskite in the Pbnm phase. <i>Journal of Solid State Chemistry</i> , 2012 , 194, 286-296	3.3	16
187	Spiral-spin-driven ferroelectricity in a multiferroic delafossite AgFeO ₂ . <i>Physical Review Letters</i> , 2012 , 109, 097203	7.4	47
186	Acentric magnetic and optical properties of chalcopyrite (CuFeS ₂). <i>Journal of Physics Condensed Matter</i> , 2012 , 24, 216001	1.8	12
185	Combined neutron and X-ray diffraction determination of disorder in doped zirconolite-2M. <i>American Mineralogist</i> , 2012 , 97, 291-298	2.9	22
184	Pressure-dependent spin fluctuations and magnetic structure in the topologically frustrated spin glass alloy Y(Mn _{0.95} Al _{0.05}) ₂ . <i>Physical Review B</i> , 2012 , 85,	3.3	2
183	Crystal structures, strain analysis, and physical properties of Sr _{0.7} Ce _{0.3} MnO ₃ . <i>Physical Review B</i> , 2012 , 85,	3.3	14

182	Impact of Jahn-Teller active Mn ³⁺ on strain effects and phase transitions in Sr _{0.65} Pr _{0.35} MnO ₃ . <i>Physical Review B</i> , 2012 , 85,	3.3	20
181	Structural and thermoelastic properties of CaTiO ₃ perovskite between 7 K and 400 K. <i>Journal of Alloys and Compounds</i> , 2011 , 509, 6337-6345	5.7	23
180	THE LOW-TEMPERATURE AND HIGH-PRESSURE THERMOELASTIC AND STRUCTURAL PROPERTIES OF CHALCOPYRITE, CuFeS ₂ . <i>Canadian Mineralogist</i> , 2011 , 49, 1015-1034	0.7	30
179	Structural and thermoelastic study of the protonic conducting perovskite SrCe _{0.95} Yb _{0.05} O _{3-δ} between 373 K and 1273 K. <i>Journal of Electroceramics</i> , 2011 , 27, 143-153	1.5	6
178	Thermoelastic properties and crystal structure of CaPtO ₃ post-perovskite from 0 to 9 GPa and from 2 to 973 K. <i>Journal of Applied Crystallography</i> , 2011 , 44, 999-1016	3.8	8
177	Negative linear compressibility and massive anisotropic thermal expansion in methanol monohydrate. <i>Science</i> , 2011 , 331, 742-6	33.3	178
176	CENTROSYMMETRIC PEROVSKITE CRYSTAL STRUCTURES WITH SPACE GROUP Pbnm: CRYSTALLOGRAPHIC PARAMETERIZATION OF KCaF ₃ BETWEEN 100 AND 400 K IN TERMS OF THE AMPLITUDES OF SYMMETRY-ADAPTED BASIS VECTORS OF THE CUBIC ARISTOTYPE PHASE. <i>Canadian Mineralogist</i> , 2011 , 49, 793-808	0.7	16
175	High-temperature phase transitions of hexagonal YMnO ₃ . <i>Physical Review B</i> , 2011 , 83,	3.3	158
174	Determination of structural chirality of berlinite and quartz using resonant x-ray diffraction with circularly polarized x-rays. <i>Physical Review B</i> , 2010 , 81,	3.3	18
173	Experimental evidence of anapolar moments in the antiferromagnetic insulating phase of V ₂ O ₃ obtained from x-ray resonant Bragg diffraction. <i>Physical Review B</i> , 2010 , 81,	3.3	25
172	Symmetry and strain analysis of structural phase transitions in Pr _{0.48} Ca _{0.52} MnO ₃ . <i>Physical Review B</i> , 2010 , 82,	3.3	22
171	Structural variations in the wesselsiteeffenbergerite (Sr _{1-x} Ba _x CuSi ₄ O ₁₀) solid solution. <i>European Journal of Mineralogy</i> , 2010 , 22, 411-423	2.2	9
170	X-ray absorption and neutron diffraction studies of (Sr(1-x)Ce(x))MnO ₃ : transition from coherent to incoherent static Jahn-Teller distortions. <i>Journal of Physics Condensed Matter</i> , 2010 , 22, 445401	1.8	8
169	Giant magnetoelastic coupling in a metallic helical metamagnet. <i>Physical Review Letters</i> , 2010 , 104, 247202	7.0	70
168	Zigzag ladders with staggered magnetic chirality in the S=3/2 compound BaCaCr ₂ O ₄ . <i>Physical Review B</i> , 2010 , 81,	3.3	38
167	The polar phase of NaNbO ₃ : a combined study by powder diffraction, solid-state NMR, and first-principles calculations. <i>Journal of the American Chemical Society</i> , 2010 , 132, 8732-46	16.4	137
166	Combined experimental and modelling studies of proton conducting La _{1-x} Ba _x O _{4-δ} : proton location and dopant site selectivity. <i>Journal of Materials Chemistry</i> , 2010 , 20, 10412		12
165	Analytical expressions to determine the isothermal compressibility tensor and the isobaric thermal expansion tensor for monoclinic crystals: application to determine the direction of maximum compressibility in jadeite. <i>Physics and Chemistry of Minerals</i> , 2010 , 37, 529-533	1.6	33

164	The Γ - Γ transition in BiFeO ₃ : A Powder Neutron Diffraction Study. <i>Advanced Functional Materials</i> , 2010 , 20, 2116-2123	15.6	81
163	The crystal structure of perdeuterated methanol hemiammoniate (CD ₃ OD _{0.5} ND ₃) determined from neutron powder diffraction data at 4.2 and 180 K. <i>Journal of Applied Crystallography</i> , 2010 , 43, 328-336	3.8	8
162	Polysomatic apatites. <i>Acta Crystallographica Section B: Structural Science</i> , 2010 , 66, 1-16		28
161	Revision of the structure of Cs ₂ CuSi ₅ O ₁₂ leucite as orthorhombic Pbc _a . <i>Acta Crystallographica Section B: Structural Science</i> , 2010 , 66, 51-9		15
160	PARAMETERIZATION OF CENTROSYMMETRIC ELPASOLITE-TYPE CRYSTAL STRUCTURES IN TERMS OF SYMMETRY-ADAPTED BASIS-VECTORS OF THE PRIMITIVE CUBIC ARISTOTYPE PHASE. <i>Canadian Mineralogist</i> , 2009 , 47, 401-420	0.7	5
159	Ambi-site substitution of Mn in lanthanum germanate apatites. <i>Materials Research Bulletin</i> , 2009 , 44, 1806-1809	5.1	15
158	The thermal expansion and crystal structure of mirabilite (Na ₂ SO ₄ ·10D ₂ O) from 4.2 to 300 K, determined by time-of-flight neutron powder diffraction. <i>Physics and Chemistry of Minerals</i> , 2009 , 36, 29-46	1.6	35
157	Phase behaviour and thermoelastic properties of perdeuterated ammonia hydrate and ice polymorphs from 0 to 2 GPa. <i>Journal of Applied Crystallography</i> , 2009 , 42, 846-866	3.8	27
156	The crystal structure of perdeuterated methanol monoammoniate (CD ₃ OD ₁ ND ₃) determined from neutron powder diffraction data at 4.2 and 180 K. <i>Journal of Applied Crystallography</i> , 2009 , 42, 1054-1061	3.8	10
155	Structure, crystal chemistry and thermal evolution of the Bi ₂ O ₃ -related phase Bi ₉ ReO ₁₇ . <i>Journal of Solid State Chemistry</i> , 2009 , 182, 2468-2474	3.3	5
154	Synchrotron X-ray absorption spectroscopy and X-ray powder diffraction studies of the structure of johnbaumite [Ca ₁₀ (AsO ₄) ₆ (OH,F) ₂] and synthetic Pb-, Sr- and Ba-arsenate apatites and some comments on the crystal chemistry of the apatite structure type in general. <i>Mineralogical Magazine</i> , 2009 , 73, 433-455	1.7	17
153	PARAMETERIZATION OF THE CRYSTAL STRUCTURES OF CENTROSYMMETRIC ZONE-BOUNDARY-TILTED PEROVSKITES: AN ANALYSIS IN TERMS OF SYMMETRY-ADAPTED BASIS-VECTORS OF THE CUBIC ARISTOTYPE PHASE. <i>Canadian Mineralogist</i> , 2009 , 47, 381-400	0.7	29
152	Ferroelectric-paraelectric transition in BiFeO ₃ : crystal structure of the orthorhombic beta phase. <i>Physical Review Letters</i> , 2009 , 102, 027602	7.4	261
151	Temperature- and pressure-induced proton transfer in the 1:1 adduct formed between squaric acid and 4,4'-bipyridine. <i>Journal of the American Chemical Society</i> , 2009 , 131, 3884-93	16.4	72
150	PROTON LOCATION AND HYDROGEN BONDING IN THE HYDROUS LEAD COPPER SULFATES LINARITE, PbCu(SO ₄)(OH) ₂ , AND CALEDONITE, Pb ₅ Cu ₂ (SO ₄) ₃ CO ₃ (OH) ₆ . <i>Canadian Mineralogist</i> , 2009 , 47, 649-662	0.7	8
149	Characteristic length scale for strain fields around impurity cations in perovskites. <i>Physical Review B</i> , 2009 , 80,	3.3	25
148	Right handed or left handed? Forbidden x-ray diffraction reveals chirality. <i>Physical Review Letters</i> , 2008 , 100, 145502	7.4	47
147	(Ca _{0.37} Sr _{0.63})TiO ₃ perovskite: an example of an unusual class of tilted perovskites. <i>Journal of Physics Condensed Matter</i> , 2008 , 20, 135202	1.8	17

146	Neutron powder diffraction studies of sulfuric acid hydrates. II. The structure, thermal expansion, incompressibility, and polymorphism of sulfuric acid tetrahydrate (D2SO4.4D2O). <i>Journal of Chemical Physics</i> , 2008 , 128, 054506	3.9	14
145	Local structure and disorder in crystalline Pb9Al8O21. <i>Journal of Solid State Chemistry</i> , 2008 , 181, 1087-1102	3.9	11
144	The crystal structure and thermal expansion tensor of MgSO4·1D2O(meridianiite) determined by neutron powder diffraction. <i>Physics and Chemistry of Minerals</i> , 2008 , 35, 207-221	1.6	62
143	POWDER NEUTRON-DIFFRACTION STUDIES OF CLINOPYROXENES. I. THE CRYSTAL STRUCTURE AND THERMOELASTIC PROPERTIES OF JADEITE BETWEEN 1.5 AND 270 K. <i>Canadian Mineralogist</i> , 2008 , 46, 1593-1622	0.7	13
142	Cation Substitution in Defect Thiospinels: Structural and Magnetic Properties of GaV4-xMoxS8 (0 ≤ x ≤ 4). <i>Chemistry of Materials</i> , 2007 , 19, 5035-5044	9.6	16
141	Thermal expansion of deuterated hopeite, Zn3(PO4)2·4D2O. <i>American Mineralogist</i> , 2007 , 92, 1038-1047.	2.9	2
140	Structural anomalies at the magnetic transition in centrosymmetric BiMnO3. <i>Physical Review B</i> , 2007 , 75,	3.3	70
139	Structures and phase diagram for the system CaTiO3-xBa2/3TiO3. <i>Journal of Solid State Chemistry</i> , 2007 , 180, 1083-1092	3.3	30
138	High-resolution neutron powder diffraction study on the phase transitions in BaPbO3. <i>Journal of Solid State Chemistry</i> , 2007 , 180, 1559-1565	3.3	16
137	Crystal structure of Ln1/3NbO3 (Ln=Nd, Pr) and phase transition in Nd1/3NbO3. <i>Journal of Solid State Chemistry</i> , 2007 , 180, 1846-1851	3.3	20
136	Crystal structures and thermal expansion of ϵ -MgSO4 and δ -MgSO4 from 4.2 to 300 K by neutron powder diffraction. <i>Journal of Applied Crystallography</i> , 2007 , 40, 761-770	3.8	30
135	Cooperative mechanisms of fast-ion conduction in gallium-based oxides with tetrahedral moieties. <i>Nature Materials</i> , 2007 , 6, 871-5	27	164
134	High temperature structural and thermoelastic behaviour of mantle orthopyroxene: an in situ neutron powder diffraction study. <i>Physics and Chemistry of Minerals</i> , 2007 , 34, 185-200	1.6	14
133	Structural studies of the proton conducting perovskite δ -Ba0.6Ba0.4ScO2.8. <i>Solid State Ionics</i> , 2007 , 178, 943-949	3.3	24
132	Polar and magnetoelectric multipoles in gallium ferrate inferred from optical and x-ray measurements. <i>Journal of Physics Condensed Matter</i> , 2007 , 19, 376205	1.8	10
131	Observation of two spin gap energies in the filled skutterudite compound CeOs4Sb12. <i>Physical Review B</i> , 2007 , 75,	3.3	19
130	Vanadium magnetoelectric multipoles in V2O3. <i>Physical Review B</i> , 2007 , 75,	3.3	22
129	Structural phase transition and magnetism in hexagonal SrMnO3 by magnetization measurements and by electron, x-ray, and neutron diffraction studies. <i>Physical Review B</i> , 2007 , 75,	3.3	34

128	Structural basis for the anomalous low-temperature thermal expansion behaviour of the gillespite-structured phase Ba _{0.5} Sr _{0.5} CuSi ₄ O ₁₀ . <i>European Journal of Mineralogy</i> , 2007 , 19, 189-200	2.2	7
127	The high-pressure phase diagram of ammonia dihydrate. <i>High Pressure Research</i> , 2007 , 27, 201-212	1.6	50
126	Suppression of strain coupling in perovskite La _{0.6} Sr _{0.1} TiO ₃ by cation disorder. <i>Physical Review B</i> , 2007 , 76,	3.3	25
125	Cation ordering in MgTi ₂ O ₅ (karrooite): Probing temperature dependent effects with neutrons. <i>American Mineralogist</i> , 2007 , 92, 1165-1180	2.9	18
124	Unusual high-temperature structural behaviour in ferroelectric Bi ₂ WO ₆ . <i>Chemistry - A European Journal</i> , 2006 , 12, 1493-9	4.8	129
123	Neutron powder diffraction studies of sulfuric acid hydrates. I. The structure of sulfuric acid hemitrisikaidekahydrate D ₂ SO ₄ .6(1/2)D ₂ O. <i>Journal of Chemical Physics</i> , 2006 , 125, 144510	3.9	14
122	Rhombohedral to cubic phase transition in the relaxor ferroelectric PZN. <i>Journal of Physics Condensed Matter</i> , 2006 , 18, L233-L240	1.8	19
121	Structural relationships and a phase diagram for (Ca,Sr)TiO ₃ perovskites. <i>Journal of Physics Condensed Matter</i> , 2006 , 18, 10725-10749	1.8	64
120	Enantioselective surface chemistry of R-2-bromobutane on Cu(643)R&S and Cu(531)R&S. <i>Journal of Physical Chemistry B</i> , 2006 , 110, 10411-20	3.4	31
119	The thermoelastic properties of MgSO ₄ .7D ₂ O (epsomite) from powder neutron diffraction and ab initio calculation. <i>European Journal of Mineralogy</i> , 2006 , 18, 449-462	2.2	41
118	PbZn(1/3)Nb(2/3)O ₃ at 4.2 and 295 K. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2006 , 62, i46-8		6
117	Structures of the cation-deficient perovskite Nd(0.7)Ti(0.9)Al(0.1)O ₃ from high-resolution neutron powder diffraction in combination with group-theoretical analysis. <i>Acta Crystallographica Section B: Structural Science</i> , 2006 , 62, 60-7		17
116	Structures and phase transitions in the ordered double perovskites Ba ₂ BiIII BiVO ₆ and Ba ₂ BiIIISbVO ₆ . <i>Acta Crystallographica Section B: Structural Science</i> , 2006 , 62, 537-46		96
115	Phases II and IV of 1,3,5-trichloro-2,4,6-trimethylbenzene: Ab initio crystal structure determination by high-resolution powder diffraction. <i>Journal of Molecular Structure</i> , 2006 , 791, 41-52	3.4	10
114	Neutron diffraction and inelastic neutron scattering investigations of the ordered double perovskite Ba ₂ PrIrO ₆ . <i>Physica B: Condensed Matter</i> , 2006 , 378-380, 543-545	2.8	13
113	Phase transitions in PZN ₈₅ PT ₁₅ in the range 4.2 $\bar{5}$ 0 K. <i>Physica B: Condensed Matter</i> , 2006 , 385-386, 160-162	2.8	4
112	Structural and magnetic properties of the Kagom \bar{a} ntiferromagnet YbBaCo ₄ O ₇ . <i>Journal of Solid State Chemistry</i> , 2006 , 179, 1136-1145	3.3	128
111	Structure and thermoelectric properties of the ordered skutterudite CoGe _{1.5} Te _{1.5} . <i>Journal of Solid State Chemistry</i> , 2006 , 179, 2047-2053	3.3	41

110	Phase relations and crystal structures in the systems (Bi,Ln) ₂ WO ₆ and (Bi,Ln) ₂ MoO ₆ (Ln=lanthanide). <i>Journal of Solid State Chemistry</i> , 2006 , 179, 3437-3444	3.3	27
109	Methodology and recent developments for using neutron diffraction to characterize the mechanical properties of rocks. <i>Physica B: Condensed Matter</i> , 2006 , 385-386, 938-941	2.8	4
108	Does the modulated magnetic structure of BiFeO ₃ change at low temperatures?. <i>Journal of Physics Condensed Matter</i> , 2006 , 18, 2069-2075	1.8	52
107	Pressure dependence of the crystal structure of SrCeO ₃ perovskite. <i>Journal of Alloys and Compounds</i> , 2005 , 394, 131-137	5.7	16
106	High temperature structural phase transitions in SrSnO ₃ perovskite. <i>Materials Research Bulletin</i> , 2005 , 40, 507-520	5.1	69
105	Crystal chemistry, cation ordering and thermoelastic behaviour of CoMgSiO ₄ olivine at high temperature as studied by in situ neutron powder diffraction. <i>Physics and Chemistry of Minerals</i> , 2005 , 32, 655-664	1.6	11
104	The incompressibility and thermal expansivity of D ₂ O ice II determined by powder neutron diffraction. <i>Journal of Applied Crystallography</i> , 2005 , 38, 612-618	3.8	37
103	Temperature dependence of structural parameters in the perovskite Na _{0.74} WO ₃ . <i>Physica Status Solidi (B): Basic Research</i> , 2005 , 242, 854-868	1.3	2
102	Strain mechanism for order-parameter coupling through successive phase transitions in PrAlO ₃ . <i>Physical Review B</i> , 2005 , 72,	3.3	62
101	Transformation processes in LaAlO ₃ : Neutron diffraction, dielectric, thermal, optical, and Raman studies. <i>Physical Review B</i> , 2005 , 72,	3.3	183
100	Ferroelectric-paraelectric phase transition in the n=2 Aurivillius phase Bi ₃ Ti _{1.5} W _{0.5} O ₉ : A neutron powder diffraction study. <i>Physical Review B</i> , 2005 , 71,	3.3	17
99	The crystal structure of KCaF ₃ at 4.2 and 300 K: A re-evaluation using high-resolution powder neutron diffraction. <i>Powder Diffraction</i> , 2005 , 20, 7-13	1.8	15
98	Quadrupole and hexadecapole ordering in DyB ₂ C ₂ : Direct observation with resonant x-ray diffraction. <i>Physical Review B</i> , 2004 , 69,	3.3	33
97	Structure and magnetism in synthetic pyrrhotite Fe ₇ S ₈ : A powder neutron-diffraction study. <i>Physical Review B</i> , 2004 , 70,	3.3	93
96	The role of hydrogen bonding in the thermal expansion and dehydration of brushite, di-calcium phosphate dihydrate. <i>Physics and Chemistry of Minerals</i> , 2004 , 31, 606-624	1.6	57
95	Thermal expansion and crystal structure of cementite, Fe ₃ C, between 4 and 600 K determined by time-of-flight neutron powder diffraction. <i>Journal of Applied Crystallography</i> , 2004 , 37, 82-90	3.8	153
94	A reduced moment antiferromagnetic Kondo lattice compound: Ce ₈ Pd ₂₄ Ga. <i>Journal of Magnetism and Magnetic Materials</i> , 2004 , 272-276, 622-624	2.8	5
93	Temperature-induced phase transitions in BaTbO ₃ . <i>Journal of Solid State Chemistry</i> , 2004 , 177, 1667-1673	3.3	33

92	Neutron powder diffraction study of phase transitions in Sr ₂ SnO ₄ . <i>Journal of Solid State Chemistry</i> , 2004 , 177, 4081-4086	3.3	17
91	No evidence for large-scale proton ordering in Antarctic ice from powder neutron diffraction. <i>Journal of Chemical Physics</i> , 2004 , 120, 11376-9	3.9	39
90	Using neutron diffraction measurements to characterize the mechanical properties of polymineralic rocks. <i>Mineralogical Magazine</i> , 2003 , 67, 967-987	1.7	16
89	Spontaneous strain variations through the low temperature phase transitions of deuterated lawsonite. <i>American Mineralogist</i> , 2003 , 88, 534-546	2.9	36
88	Structural distortions in the layered perovskites CsANb ₂ O ₇ (A=Nd, Bi). <i>Journal of Solid State Chemistry</i> , 2003 , 173, 309-313	3.3	44
87	Structures of Na _{0.74} WO ₃ . <i>Acta Crystallographica Section B: Structural Science</i> , 2003 , 59, 584-7		8
86	Using neutron diffraction to investigate the elastic properties of anisotropic rocks: Results from an olivine + orthopyroxene mylonite. <i>Journal of Geophysical Research</i> , 2003 , 108,		8
85	Determination of B-site ordering and structural transformations in the mixed transition metal perovskites La ₂ CoMnO ₆ and La ₂ NiMnO ₆ . <i>Journal of Physics Condensed Matter</i> , 2003 , 15, 4927-4936	1.8	215
84	A high-resolution neutron powder diffraction study of ammonia dihydrate (ND ₃ ?2D ₂ O) phase I. <i>Journal of Chemical Physics</i> , 2003 , 119, 10806-10813	3.9	33
83	Reduced moment magnetic ordering in a Kondo lattice compound: Ce ₈ Pd ₂₄ Ga. <i>Physical Review B</i> , 2003 , 67,	3.3	6
82	Displacive Phase Transitions in and Strain Analysis of Fe-Doped CaTiO ₃ Perovskites at High Temperatures by Neutron Diffraction. <i>Journal of Solid State Chemistry</i> , 2002 , 167, 459-471	3.3	13
81	Thermal expansion and crystal structure of FeSi between 4 and 1173 K determined by time-of-flight neutron powder diffraction. <i>Physics and Chemistry of Minerals</i> , 2002 , 29, 132-139	1.6	88
80	An interpretation of two X-ray studies of vanadium sesquioxide (V ₂ O ₃). <i>Physica B: Condensed Matter</i> , 2002 , 318, 261-266	2.8	
79	Thermal expansion and atomic displacement parameters of cubic KMgF ₃ perovskite determined by high-resolution neutron powder diffraction. <i>Journal of Applied Crystallography</i> , 2002 , 35, 291-295	3.8	51
78	Contrasting Structural Behavior in the Aurivillius Phase Ferroelectrics Bi ₄ Ti ₃ O ₁₂ , BaBi ₄ Ti ₄ O ₁₅ and Ba ₂ Bi ₄ Ti ₅ O ₁₈ . <i>Materials Research Society Symposia Proceedings</i> , 2002 , 755, 1		
77	Accurate quantification of the modal mineralogy of rocks when image analysis is difficult. <i>Mineralogical Magazine</i> , 2002 , 66, 189-200	1.7	35
76	Orbital magnetization of a Mott insulator, V ₂ O ₃ , revealed by resonant x-ray Bragg diffraction. <i>Physical Review B</i> , 2002 , 65,	3.3	20
75	High-temperature phase transitions in tungsten trioxide - the last word?. <i>Journal of Physics Condensed Matter</i> , 2002 , 14, 377-387	1.8	73

74	Static and dynamic structures of $\text{CD}_3\text{ND}_3\text{GeCl}_3$ studied by TOF high resolution neutron powder diffraction and solid state NMR. <i>Dalton Transactions RSC</i> , 2002 , 2112-2118		34
73	The effect of ferromagnetism on the equation of state of Fe_3C studied by first-principles calculations. <i>Earth and Planetary Science Letters</i> , 2002 , 203, 567-575	5.3	102
72	Displacive components of the low-temperature phase transitions in lawsonite. <i>American Mineralogist</i> , 2001 , 86, 566-577	2.9	31
71	Structural phase transitions, oxygen vacancy ordering and protonation in doped BaCeO_3 : results from time-of-flight neutron powder diffraction investigations. <i>Solid State Ionics</i> , 2001 , 145, 275-294	3.3	106
70	Neutron scattering studies of an antiferromagnetic Kondo compound: $\text{Ce}_8\text{Pd}_{24}\text{Ga}$. <i>Journal of Physics Condensed Matter</i> , 2001 , 13, 459-473	1.8	4
69	Composition and temperature dependence of cation ordering in Ni-Mg olivine solid solutions: a time-of-flight neutron powder diffraction and EXAFS study. <i>American Mineralogist</i> , 2001 , 86, 1170-1187	2.9	21
68	Resonant $1s$ - β x-ray Bragg diffraction and structure factors for transition-metal compounds. <i>Physical Review B</i> , 2001 , 64,	3.3	3
67	Calculated x-ray dichroic signals and resonant Bragg diffraction structure factors for DyB_2C_2 . <i>Physical Review B</i> , 2001 , 64,	3.3	20
66	The microscopic origin of thermal cracking in rocks: An investigation by simultaneous time-of-flight neutron diffraction and acoustic emission monitoring. <i>Geophysical Research Letters</i> , 2001 , 28, 2105-2108	4.9	24
65	Single-crystal X-ray diffraction analysis of pyrene II at 93K. <i>Journal of Molecular Structure</i> , 2000 , 520, 29-32	3.4	42
64	Neutron powder diffraction studies of the thermal behaviour of deuterated chalcantinite. <i>Physica B: Condensed Matter</i> , 2000 , 276-278, 897-898	2.8	4
63	Powder neutron diffraction studies of $\text{BaCe}_{0.9}\text{Y}_{0.1}\text{O}_{2.95}$ and BaCeO_3 at 4.2 K: a possible structural site for the proton. <i>Solid State Ionics</i> , 2000 , 127, 43-48	3.3	49
62	Residual stress measurements in an SiC continuous fiber reinforced Ti matrix composite. <i>Scripta Materialia</i> , 2000 , 42, 775-779	5.6	7
61	Octahedral cation ordering in olivine at high temperature. II: an in situ neutron powder diffraction study on synthetic MgFeSiO_4 (Fa50). <i>Physics and Chemistry of Minerals</i> , 2000 , 27, 630-637	1.6	65
60	A high temperature structural phase transition in crocoite (PbCrO_4) at 1068 K: crystal structure refinement at 1073 K and thermal expansion tensor determination at 1000 K. <i>Mineralogical Magazine</i> , 2000 , 64, 291-300	1.7	16
59	Resonant x-ray Bragg diffraction from orbital moments in vanadium sesquioxide (V_2O_3) and haematite ($\alpha\text{Fe}_2\text{O}_3$). <i>Journal of Physics Condensed Matter</i> , 2000 , 12, L367-L372	1.8	19
58	The structural phase transitions in strontium zirconate revisited. <i>Journal of Physics Condensed Matter</i> , 2000 , 12, L677-L683	1.8	160
57	Temperature related structural variation of the hydrous components in gypsum. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2000 , 215,	1	9

56	In situ high-t neutron diffraction studies of non-convergent order/disorder in minerals: From simple oxides to complex silicates. <i>Phase Transitions</i> , 1999 , 69, 17-34	1.3	4
55	Oxygen vacancy ordering in neodymium-doped barium cerate. <i>Solid State Communications</i> , 1999 , 112, 73-78	1.6	13
54	High-resolution neutron diffraction evidence for strain-induced microdomains during the phase transition in UNi ₂ Sn. <i>Physica B: Condensed Matter</i> , 1999 , 262, 312-316	2.8	1
53	On the lattice parameters of sodium niobate at room temperature and above. <i>Physica B: Condensed Matter</i> , 1999 , 266, 368-372	2.8	77
52	High-temperature phases of NaNbO ₃ and NaTaO ₃ . <i>Acta Crystallographica Section B: Structural Science</i> , 1999 , 55, 24-30		110
51	Structure and dynamics of methyl groups in the deuterated microporous organic/organic hybrid, aluminium methylphosphonate- α . <i>Chemical Physics Letters</i> , 1999 , 313, 505-513	2.5	14
50	Temperature evolution between 50 K and 320 K of the thermal expansion tensor of gypsum derived from neutron powder diffraction data. <i>Physics and Chemistry of Minerals</i> , 1999 , 26, 477-483	1.6	20
49	A neutron powder diffraction study of cation ordering in high-temperature synthetic amphiboles. <i>European Journal of Mineralogy</i> , 1999 , 11, 321-332	2.2	32
48	In-situ neutron diffraction study of non-convergent cation ordering in the (Fe ₃ O ₄) _{1-x} (MgAl ₂ O ₄) _x spinel solid solution. <i>American Mineralogist</i> , 1999 , 84, 555-563	2.9	9
47	Muon sites and diffusion in barium cerate. <i>Solid State Ionics</i> , 1998 , 113-115, 341-345	3.3	2
46	Structural Study of the Proton Conductor Cs ₃ H(SeO ₄) ₂ by High Resolution Neutron Powder Diffraction. <i>Materials Science Forum</i> , 1998 , 278-281, 726-731	0.4	1
45	Structures and phase transitions of in the temperature range. <i>Journal of Physics Condensed Matter</i> , 1998 , 10, 9975-9989	1.8	8
44	Structural order parameters of the transition - a soft-Ising-system. <i>Journal of Physics Condensed Matter</i> , 1998 , 10, 8799-8813	1.8	
43	Fe-Mn cation ordering in fayalite β ephroite (Fe _x Mn _{1-x}) ₂ SiO ₄ olivines: a neutron diffraction study. <i>Mineralogical Magazine</i> , 1998 , 62, 607-615	1.7	17
42	The influence of thermal vibrations on the average structure of cubic perovskite: a combined molecular dynamics and neutron diffraction study. <i>Journal of Physics Condensed Matter</i> , 1997 , 9, L647-L655	1.8	16
41	Structural aspects of the phase transition in. <i>Journal of Physics Condensed Matter</i> , 1997 , 9, 3503-3519	1.8	30
40	The structure of Bi ₂ Sn ₂ O ₇ at 725 °C by high-resolution neutron diffraction: implications for bismuth(III)-containing pyrochlores. <i>Journal of the Chemical Society Dalton Transactions</i> , 1997 , 2551-2556		21
39	Cooperative Jahn-Teller Effect in Titanium Alum. <i>Journal of the American Chemical Society</i> , 1997 , 119, 3324-3332	16.4	21

38	Crystal structure of BaMo ₃ O ₁₀ solved from powder diffraction data. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 1997 , 212,	1	3
37	The static compressibility of gypsum. <i>Geophysical Research Letters</i> , 1997 , 24, 1267-1270	4.9	11
36	Structural phase transitions in germanate analogues of investigated by high-resolution neutron powder diffraction. <i>Journal of Physics Condensed Matter</i> , 1997 , 9, 3833-3851	1.8	23
35	Crystal structure and paramagnetic behaviour of. <i>Journal of Physics Condensed Matter</i> , 1997 , 9, 6563-6577	7.8	85
34	Determination of the hydrogen position in Cs ₃ H(SeO ₄) ₂ at 483 K. <i>Physica B: Condensed Matter</i> , 1997 , 234-236, 89-91	2.8	3
33	First-order valence change in CeNi _{1-x} CoxSn (0.35 ≤ x ≤ 0.4). <i>Physica B: Condensed Matter</i> , 1997 , 234-236, 872-874	2.8	3
32	Powder neutron diffraction studies of the thermal expansion, compressibility and dehydration of deuterated gypsum. <i>Physica B: Condensed Matter</i> , 1997 , 234-236, 942-944	2.8	3
31	Nuclear and magnetic structures of the distorted hexagonal perovskites ND ₄ FeCl ₃ and NH ₄ FeCl ₃ . <i>Physica B: Condensed Matter</i> , 1997 , 241-243, 385-386	2.8	
30	Crystal Structure Solution from Neutron Powder Diffraction Data by a new Monte Carlo Approach Incorporating Restrained Relaxation of the Molecular Geometry. <i>Journal of Applied Crystallography</i> , 1997 , 30, 968-974	3.8	15
29	Distortion Characteristics Across the Structural Phase Transition in (Cu _{1-x} Zn _x)WO ₄ . <i>Acta Crystallographica Section B: Structural Science</i> , 1997 , 53, 102-112		48
28	High-Temperature Study of Octahedral Cation Exchange in Olivine by Neutron Powder Diffraction. <i>Science</i> , 1996 , 271, 1713-1715	33.3	54
27	Constrained Rietveld refinement of [D ₁ H ₁]decadeuteriodopamine deuteriobromide using powder neutron diffraction data. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1996 , 92, 4555-4559		4
26	Thermal expansion of gypsum investigated by neutron powder diffraction. <i>American Mineralogist</i> , 1996 , 81, 847-851	2.9	63
25	A neutron powder diffraction determination of the thermal expansion tensor of crocoite (PbCrO ₄) between 60 K and 290 K. <i>Mineralogical Magazine</i> , 1996 , 60, 963-972	1.7	23
24	Some observations on the crystal structure of (R,S)-propranolol hydrochloride. <i>International Journal of Pharmaceutics</i> , 1996 , 137, 255-259	6.5	5
23	Neutron diffraction of a complex of 1,8-bis(dimethylamino)naphthalene with 1,2-dichloromaleic acid. <i>Acta Crystallographica Section B: Structural Science</i> , 1996 , 52, 691-696		31
22	The crystal structure of perdeuterated pyrene II at 4.2 K. <i>Chemical Physics Letters</i> , 1996 , 258, 490-494	2.5	10
21	Assignment of tunnelling lines by single crystal neutron spectroscopy. <i>Physica B: Condensed Matter</i> , 1996 , 226, 238-240	2.8	10

20	Rb ₃ H(SeO ₄) ₂ at 4K by Neutron Powder Diffraction. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 1996 , 52, 1061-1063		10
19	Neutron powder diffraction study of the scintillator material ZnWO ₄ . <i>Journal of Materials Science</i> , 1996 , 31, 2873-2877	4.3	62
18	Determination of olivine cooling rates from metal-cation ordering. <i>Nature</i> , 1996 , 381, 407-409	50.4	49
17	Neutron Diffraction Pole-Figure Measurements Using a Pulsed White Beam and the Linear Julios-Detector. <i>Materials Science Forum</i> , 1996 , 228-231, 259-264	0.4	
16	The low-temperature behaviour of analcime. 1: high-resolution neutron powder diffraction. <i>Mineralogical Magazine</i> , 1996 , 60, 499-507	1.7	8
15	The crystal structures of some doped and undoped alkaline earth cerate perovskites. <i>Materials Research Bulletin</i> , 1995 , 30, 347-356	5.1	82
14	First-order valence phase transition in CeNi _{1-x} CoxSn alloys. <i>Physical Review B</i> , 1995 , 52, 12790-12797	3.3	32
13	A new uranium based compound with a cubic structure: UNi _{0.8} Pt _{0.2} . <i>Solid State Communications</i> , 1994 , 92, 389-392	1.6	2
12	Structural study of barium titanate between 150 and 425 K. <i>Phase Transitions</i> , 1994 , 48, 217-236	1.3	20
11	Space group and lattice constants for barium cerate and minor corrections to the crystal structures of BaCe _{0.9} Y _{0.1} O _{2.95} and BaCe _{0.9} Gd _{0.1} O _{2.95} . <i>Journal of Materials Chemistry</i> , 1994 , 4, 899		38
10	Structural study of potassium niobate between 200 and 823 K. <i>Phase Transitions</i> , 1994 , 52, 261-275	1.3	10
9	The crystal structure of ferroelectric Bi ₂ WO ₆ at 961 K. <i>Ferroelectrics</i> , 1993 , 150, 319-330	0.6	24
8	Koordinationszahl 4 oder 6 für Lithium??: Die Kristallstruktur von wasserfreiem Lithiumpermanganat, Li[MnO ₄]. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 1993 , 619, 1419-1425	1.3	7
7	An introduction to Bayesian model selection. <i>Physica D: Nonlinear Phenomena</i> , 1993 , 66, 234-242	3.3	25
6	Crystal structures of gadolinium- and yttrium-doped barium cerate. <i>Journal of Materials Chemistry</i> , 1992 , 2, 709		33
5	The crystal structure of russellite; a re-determination using neutron powder diffraction of synthetic Bi ₂ WO ₆ . <i>Mineralogical Magazine</i> , 1992 , 56, 399-409	1.7	104
4	Enantiospecific preparation of [(2 <i>r</i> ,6 <i>s</i>)-endo]-5-aza-1,10,10-trimethyl-3-oxatricyclo[5.2.1.0 ^{2,6}]decan-4-one by a nitrene-mediated route from [(1 <i>s</i>)-endo]-1-borneol and its utility as a chiral auxiliary in some asymmetric transformations. <i>Tetrahedron</i> , 1992 , 48, 7979-8006	2.4	43
3	Time-of-flight powder diffraction test experiments at a pulsed neutron beam using the position-sensitive detector JULIOS. <i>Nuclear Instruments and Methods in Physics Research, Section A: Accelerators, Spectrometers, Detectors and Associated Equipment</i> , 1992 , 317, 202-212	1.2	6

2	The crystal structures of CuInSe ₂ and CuInTe ₂ . <i>Materials Research Bulletin</i> , 1992 , 27, 161-167	5.1	90
1	Shallow ion implantation in gallium arsenide. <i>Vacuum</i> , 1984 , 34, 199-201	3.7	1