

Andris Anspoks

List of Publications by Year in descending order

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papers

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687220

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times ranked

951
citing authors

#	ARTICLE	IF	CITATIONS
1	Neural Network Approach for Characterizing Structural Transformations by X-Ray Absorption Fine Structure Spectroscopy. <i>Physical Review Letters</i> , 2018, 120, 225502.	2.9	85
2	Atomic structure relaxation in nanocrystalline NiO studied by EXAFS spectroscopy: Role of nickel vacancies. <i>Physical Review B</i> , 2012, 86, .	1.1	60
3	Interpretation of the Ni K-edge EXAFS in nanocrystalline nickel oxide using molecular dynamics simulations. <i>Journal of Non-Crystalline Solids</i> , 2011, 357, 2604-2610.	1.5	52
4	Temperature dependence of the local structure and lattice dynamics of wurtzite-type ZnO. <i>Acta Materialia</i> , 2014, 79, 194-202.	3.8	30
5	Probing NiO nanocrystals by EXAFS spectroscopy. <i>Solid State Communications</i> , 2010, 150, 2270-2274.	0.9	24
6	Local structure relaxation in nanosized tungstates. <i>Solid State Communications</i> , 2014, 183, 22-26.	0.9	23
7	Thermal disorder and correlation effects in anti-perovskite-type copper nitride. <i>Acta Materialia</i> , 2017, 129, 61-71.	3.8	21
8	Thermal holograms in doped ferroelectric SBN crystals. <i>Ferroelectrics</i> , 1988, 80, 277-280.	0.3	20
9	External pressure and composition effects on the atomic and electronic structure of SnWO ₄ . <i>Solar Energy Materials and Solar Cells</i> , 2015, 143, 627-634.	3.0	19
10	Local structural investigation of hafnia-zirconia polymorphs in powders and thin films by X-ray absorption spectroscopy. <i>Acta Materialia</i> , 2019, 180, 158-169.	3.8	19
11	Analysis of extended x-ray absorption fine structure data from copper tungstate by the reverse Monte Carlo method. <i>Physica Scripta</i> , 2014, 89, 044006.	1.2	18
12	Treatment of disorder effects in X-ray absorption spectra beyond the conventional approach. <i>Radiation Physics and Chemistry</i> , 2020, 175, 108112.	1.4	16
13	Local structure of nanosized tungstates revealed by evolutionary algorithm. <i>Physica Status Solidi (A) Applications and Materials Science</i> , 2015, 212, 265-273.	0.8	15
14	The role of yttrium and titanium during the development of ODS ferritic steels obtained through the STARS route: TEM and XAS study. <i>Journal of Nuclear Materials</i> , 2018, 504, 8-22.	1.3	14
15	ODS ferritic steels obtained from gas atomized powders through the STARS processing route: Reactive synthesis as an alternative to mechanical alloying. <i>Nuclear Materials and Energy</i> , 2018, 17, 1-8.	0.6	14
16	The Use of X-ray Absorption Spectra for Validation of Classical Force-Field Models. <i>Zeitschrift Fur Physikalische Chemie</i> , 2016, 230, 537-549.	1.4	13
17	Investigation of precipitate in an austenitic ODS steel containing a carbon-rich process control agent. <i>Nuclear Materials and Energy</i> , 2018, 15, 237-243.	0.6	13
18	High-temperature X-ray absorption spectroscopy study of thermochromic copper molybdate. <i>Acta Materialia</i> , 2019, 179, 26-35.	3.8	13

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19	Local structure of copper nitride revealed by EXAFS spectroscopy and a reverse Monte Carlo/evolutionary algorithm approach. <i>Physica Scripta</i> , 2016, 91, 054003.	1.2	12
20	Advanced approach to the local structure reconstruction and theory validation on the example of the W L ₃ -edge extended x-ray absorption fine structure of tungsten. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2018, 26, 025004.	0.8	12
21	High efficiency angular deflection of the laser beam: PLZT/intracavity array. <i>Ferroelectrics</i> , 1992, 131, 301-306.	0.3	11
22	X-ray absorption spectroscopy of Cu-doped WO ₃ films for use in electrochemical metallization cell memory. <i>Journal of Non-Crystalline Solids</i> , 2014, 401, 87-91.	1.5	10
23	Changes in structure and conduction type upon addition of Ir to ZnO thin films. <i>Thin Solid Films</i> , 2017, 636, 694-701.	0.8	10
24	Local structure and dynamics of wurtzite-type ZnO from simulation-based EXAFS analysis. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2014, 11, 1472-1475.	0.8	9
25	Extended x-ray absorption fine structure spectroscopy and first-principles study of SnWO ₄ . <i>Physica Scripta</i> , 2014, 89, 044005.	1.2	9
26	Local structure relaxation in nanocrystalline Ni _{1-x} O thin films. <i>Thin Solid Films</i> , 2014, 553, 58-62.	0.8	7
27	Local Structure of Cobalt Tungstate Revealed by EXAFS Spectroscopy and Reverse Monte Carlo/Evolutionary Algorithm Simulations. <i>Zeitschrift Fur Physikalische Chemie</i> , 2016, 230, 551-568.	1.4	7
28	Pressure-induced structural changes in Ir^{\pm} -MoO ₃ probed by X-ray absorption spectroscopy. <i>IOP Conference Series: Materials Science and Engineering</i> , 0, 503, 012018.	0.3	7
29	The local atomic structure and thermoelectric properties of Ir-doped ZnO: hybrid DFT calculations and XAS experiments. <i>Journal of Materials Chemistry C</i> , 2021, 9, 4948-4960.	2.7	7
30	Local Structure of Multiferroic MnWO ₄ and Mn _{0.7} Co _{0.3} WO ₄ Revealed by the Evolutionary Algorithm. <i>Ferroelectrics</i> , 2015, 483, 68-74.	0.3	6
31	Effect of Pressure and Temperature on the Local Structure and Lattice Dynamics of Copper(II) Oxide. <i>Physics Procedia</i> , 2016, 85, 27-35.	1.2	6
32	Local structure of A-atom in ABO ₃ perovskites studies by RMC-EXAFS. <i>Radiation Physics and Chemistry</i> , 2020, 175, 108072.	1.4	6
33	Probing vacancies in NiO nanoparticles by EXAFS and molecular dynamics simulations. <i>Journal of Physics: Conference Series</i> , 2013, 430, 012027.	0.3	5
34	ODS steel raw material local structure analysis using X-ray absorption spectroscopy. <i>IOP Conference Series: Materials Science and Engineering</i> , 2015, 77, 012029.	0.3	5
35	Interpretation of the Cu K-edge EXAFS spectra of Cu ₃ N using ab initio molecular dynamics. <i>Radiation Physics and Chemistry</i> , 2020, 175, 108100.	1.4	5
36	Controlling of visible and infrared radiation by use of holographic gratings in PLZT ceramics. <i>Ferroelectrics</i> , 1989, 90, 203-207.	0.3	4

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37	Local structure studies of SrTi ₁₆ O ₃ and SrTi ₁₈ O ₃ . Physica Scripta, 2014, 89, 044002.	1.2	4
38	Origin of Pressure-Induced Metallization in Cu ₃ N: An X-ray Absorption Spectroscopy Study. Physica Status Solidi (B): Basic Research, 2018, 255, 1800073.	0.7	4
39	Recent progress in high pressure X-ray absorption spectroscopy studies at the ODE beamline. High Pressure Research, 2020, 40, 82-87.	0.4	4
40	Study of Copper Nitride Thin Film Structure. Latvian Journal of Physics and Technical Sciences, 2016, 53, 31-37.	0.4	4
41	Effect of cobalt doping on the local structure and dynamics of multiferroic MnWO ₄ and Mn _{0.7} Co _{0.3} WO ₄ . Journal of Physics: Conference Series, 2013, 430, 012109.	0.3	3
42	Local Structure Studies of Ti for SrTi ₁₆ O ₃ and SrTi ₁₈ O ₃ by Advanced X-ray Absorption Spectroscopy Data Analysis. Ferroelectrics, 2015, 485, 42-52.	0.3	3
43	Pressure-induced insulator-to-metal transition in \pm -SnWO ₄ . Journal of Physics: Conference Series, 2016, 712, 012122.	0.3	3
44	Local dynamics and phase transition in quantum paraelectric SrTiO ₃ studied by Ti K-edge x-ray absorption spectroscopy. Journal of Physics: Conference Series, 2016, 712, 012101.	0.3	3
45	The influence of Zn ²⁺ ions on the local structure and thermochromic properties of Cu _{1-x} Zn _x MoO ₄ solid solutions. Materials Today Communications, 2021, 28, 102607. Revealing the local structure of CuMo _{1-x} Zn _x MoO ₄ solid solutions. Journal of Physics: Conference Series, 2021, 2021, 012001.	0.9	2
46	Revealing the local structure of CuMo _{1-x} Zn _x MoO ₄ solid solutions. Journal of Physics: Conference Series, 2021, 2021, 012001.	2.7	2
47	High-pressure x-ray absorption spectroscopy study of tin tungstates. Physica Scripta, 2015, 90, 094003.	1.2	1
48	Origin of Pressure-Induced Metallization in Cu ₃ N: An X-ray Absorption Spectroscopy Study (Phys.) Tj ETQq0 0 0 rgBT/Overlock 10 Tf 50	0.7	1
49	Influence of Pressure and Temperature on X-Ray Induced Photoreduction of Nanocrystalline CuO. Latvian Journal of Physics and Technical Sciences, 2018, 55, 13-19.	0.4	1
50	Photoinduced light scattering in electrooptical PLZT ceramics. Ferroelectrics, 1993, 145, 259-269.	0.3	0
51	Kinetic Monte Carlo modeling of Y ₂ O ₃ nano-cluster formation in radiation resistant matrices. Nuclear Instruments & Methods in Physics Research B, 2018, 434, 13-22.	0.6	0