

Curriculum vitae

Personal data	<i>Aleksejs Gopejenko</i> <i>19.10.1983</i> <i>Latvian</i> <i>agopejen@inbox.lv</i>
Education	<i>10.2017 – today</i> PhD student, University of Latvia, physics. <i>09.2005 - 06.2007</i> Master degree, University of Latvia, physics, diploma MD C Nr. 2973 <i>09.2001 – 07.2005</i> Professional higher, Information System Management Institute (ISMA), information systems, qualification – system analytic, diploma PD A Nr. 0466. <i>09.1990 – 06.2001</i> Riga secondary school №40
Additional education	<i>01.09.2008 – 12.09.2008</i> <i>2. Karlsruhe international summer school on fusion technologies</i> <i>10.2004 – 02.2005</i> <i>Education in the University of Osnabrück, faculty of physics</i>
Work experience	<i>11.2013 – today</i> researcher (Institute of Solid State Physics) <i>11.2007 – 11.2013</i> assistant (Institute of Solid State Physics) <i>05.2007 – 10.2007</i> engineer (Institute of Solid State Physics) <i>09.2002 – 10.2005</i> engineer (A/S Izglītības nams)
Scientific works	<i>3 researches</i> , 9 scientific publications Research projects <i>Yttrium oxide precipitates theoretical modelling in iron in support of oxide dispersed strengthened steels formation</i> <i>Lead zirconate titanate atomic and electronic structure ab initio calculations</i> <i>Lead zirconate atomic and electronic structure ab initio calculations</i>

- Publications** A. Gopejenko, S. Piskunov, and Yu.F. Zhukovskii. Ab initio modelling of the effects of varying Zr (Ti) concentrations on the atomic and electronic properties of stoichiometric PZT solid solutions. *Comput. Theor. Chem.*, 2017, **1104**, p. 56-60
- A. Gopejenko, Yu.F. Zhukovskii, E.A. Kotomin, Yu.A. Mastrikov, P.V. Vladimirov, V.A. Borodin, and A. Möslang, Ab initio modelling of Y–O cluster formation in fcc-Fe lattice. - *Phys. Stat. Sol. B*, 2016, **253**, p. 2136-2143.
- A. Gopejenko, Yu.F. Zhukovskii, P.V. Vladimirov, E.A. Kotomin, and A. Möslang, Interaction between oxygen and yttrium impurity atoms as well as vacancies in fcc iron lattice: Ab initio modeling. - Proc. NATO ARW „Nanodevices and Nanomaterials for Ecological Security” (Eds. Yuri N. Shunin and Arnold E. Kiv; Springer: Dordrecht, 2012), p. 149-160.
- A. Gopejenko, Yu.F. Zhukovskii, P.V. Vladimirov, E.A. Kotomin, and A. Möslang, Modeling of yttrium, oxygen atoms and vacancies in γ -iron lattice. - *J. Nucl. Mater.*, 2011, **416**, p. 40-44.
- A. Gopejenko, Yu.F. Zhukovskii, P.V. Vladimirov, E.A. Kotomin, and A. Möslang, Ab initio simulation of yttrium oxide nanocluster formation on fcc Fe lattice. - *J. Nucl. Mater.*, 2010, **406**, p. 345–350.
- E.A. Kotomin, S. Piskunov, Yu.F. Zhukovskii, R.I. Eglitis, A. Gopejenko, and D.E. Ellis, The electronic properties of an oxygen vacancy at ZrO₂-terminated (001) surfaces of cubic PbZrO₃: Computer simulations from the first principles. - *Phys. Chem. & Chem. Phys.*, 2008, **10**, 4258-4263.
- S. Piskunov, A. Gopeyenko, E.A. Kotomin, Yu.F. Zhukovskii, D.E. Ellis, Atomic and electronic structure of perfect and defective PbZrO₃ perovskite: hybrid DFT calculations of cubic and orthorhombic phases, *Comput. Mater. Sci.*, 2007, **41**, p. 195-201.
- A. Gopeyenko, S. Piskunov, and Yu.N. Shunin, The atomic and electronic structure of pure and defective PbZrO₃, *Computer Modelling and New Technologies (Latvia)*, 2006, 10(4), p. 7-16.
- Yu.N. Shunin, A.V. Gopejenko, Phase-shift functions method for nanoclusters electronic structure calculations in solids, *Computer Modelling and New Technologies (Latvia)*, 2006, 10(4), p. 15-31.

Place, date

Signature