

## Dr Peter Yatsyshin

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Naturalised UK citizen of Ukrainian origin

### Education

- 2014 Ph.D. **Computational statistical physics and mathematical modelling**  
Imperial College London, UK  
**My Ph.D. was funded by Marie Curie fellowship**  
*Ph.D. earned on March 1, 2014*
- 2007 M.Sc. (Hons) **Theoretical Physics**  
Department of Physics and Mechanics,  
Peter the Great Saint-Petersburg Polytechnic University<sup>1</sup>, Russia  
University Web Page: <https://english.spbstu.ru>  
**Summa Cum Laude**
- 2005 B.Sc. (Hons) **Theoretical Physics**  
Department of Physics and Mechanics,  
Peter the Great Saint-Petersburg Polytechnic University<sup>1</sup>, Russia  
**Summa Cum Laude**
- 2001 High School Second Saint-Petersburg Gymnasium, Russia

### Appointments

- 2020 – **Postdoctoral Research Associate**, Data-centric Engineering group,  
The Alan Turing Institute, UK
- 2014 – 2019 **Postdoctoral Research Associate**, Complex Multiscale Systems Group,  
Department of Chemical Engineering, Imperial College London, UK
- 2012–2014 **Research Assistant**, Complex Multiphase Systems Group,  
Department of Chemical Engineering, Imperial College London, UK
- 2009–2012 **Marie-Curie Early Stage Researcher**, Complex Multiphase Systems Group,  
Department of Chemical Engineering, Imperial College London, UK
- 2007–2009 **Researcher**, Faculty of Physics and Mechanics, Saint-Petersburg State Polytechnic  
University, Russia

### Awards and recognition

- 2019 **Assistant Supervisor**. This officially recognises my supervision of PhD and MSc projects at the departments of Mathematics and Chemical Engineering at Imperial College London
- 2018 **Fellow of The Higher Education Academy (FHEA)**. Recognition reference: PR147264.  
This is a “*recognition of attainment against the UK Professional Standards Framework for teaching and learning support in higher education.*”
- 2017 **Sir William Wakeham Award** given by the Department of Chemical Engineering at Imperial College. This is the highest departmental honour, which annually recognises one or two “*early-career researchers... who have made a significant contribution to their research field.*”  
The selection is made by an independent professorial panel. Click on links below for more info.  
[Award web page with a list of winners by year](#)    [News article about the winners in 2017](#)
- 2009 **Saint-Petersburg Government Award** for outstanding post-graduate students and early stage researchers from institutions located in the Saint-Petersburg Federal District.

<sup>1</sup>This is one of the top Russian universities and research centers in the fields of physics, applied maths and engineering. It was founded in 1899 as the most advanced engineering school in Russia at the time ([Wikipedia page](#)). Currently it ranks as 11th in Russia. Unfortunately, Russian universities do not regularly participate in international rankings. Nevertheless, some ranking info is provided on the university website: <https://english.spbstu.ru/university/>.

## Research funding

- 2017 Imperial College European Partners Fund
- 2016 Imperial College “Pathways to Impact Awards”
- 2014 Travel grant to attend “Flow 14”, International Conference on Micro- and Nanofluidics

## Teaching at Imperial College London

- 2017 - 2019 Lecturer for two-semester course CE1-07 “Mathematics 1” for 1st year Chemical Engineering undergraduates (~130 students)  
*This course consists of 55 lectures, covering differential/integral calculus, linear algebra and Euclidean geometry, ODEs and introduction to linear stability. I prepared lecture materials and notes, instructed GTAs, designed/marked assessments. I provided additional support to learners in-person and online.*
- 2015 - 2018 Lecturer for one-semester course CE4-09 “Dynamical Systems” for 4th year Chemical Engineering undergraduates (~50 students)  
*This course consists of 15 lectures, covering linear stability and bifurcations, introduction to deterministic chaos theory and stochastic dynamical systems.*
- 2015 - 2016 Lecturer, GTA and tutor at the course CE1-03-4 “Scientific programming on MATLAB and Python”  
1st year Chemical Engineering undergraduates (~ 130 students)  
*The course consists of 30 lectures, covering best practices in scientific programming.*

## PhD Students Supervision

- 2021– Co-supervisor for PhD project  
*“Physics-informed machine learning in statistical physics”,*  
Imperial College London, UK
- 2021– Co-supervisor for PhD project  
*“Computational multi-scale modelling of active fluids”,*  
Imperial College London, UK
- 2016–2020 Co-supervisor for PhD project  
*“Multi-phase fluids: Molecular dynamics, generalized Langevin equation and fluctuating dynamic density functional theory”,*  
Imperial College London, UK
- 2017–2019 Co-supervisor for PhD project  
*“Critical imperfect pitchfork bifurcations”,*  
Imperial College London, UK

## M.Sc. Students Supervision

- Nov 2017– Jun 2018 Co-supervisor of M.Sc. project  
*“Dynamical aspects of wetting by thin liquid films”,*  
Imperial College London, UK
- Nov 2015–Jun 2016 Co-supervisor of M.Sc. project  
*“Phase transitions and nucleation processes using density-functional theory”,*  
Imperial College London, UK
- Nov 2014–Jun 2015 Co-supervisor of M.Sc. project  
*“Nonequilibrium molecular dynamics simulations of nanoconfined fluids at solid-liquid interfaces”,*  
Imperial College London, UK

## Administrative and Conference-Related Activities

- 2015–2019 Post Doc Rep for Chemical Engineering  
2015– 2019 Co-organiser of the Annual Chemical Engineering Post-Doc Symposium  
2018 Co-organiser of minisymposium “Fluctuating Complex Dynamical Systems”, within the British Applied Mathematical Colloquium, University of St. Andrews, UK  
2016 Co-organiser of minisymposium “Statistical Mechanical and Phase Field Modelling of Inhomogeneous Fluids”, within the British Applied Mathematical Colloquium, University of Oxford, UK

## Publications

- 22 research articles in refereed scientific journals  
2 refereed book chapters  
2 refereed conference papers  
12 invited conference talks  
24 contributed conference talks

### Ph.D. thesis

**Analysis of the phase behaviour of a nano-confined Lennard-Jones fluid using pseudo-spectral approach to classical density-functional models**

P. Yatsyshin (2014). Ph.D. thesis. Imperial College London.

URL: <http://hdl.handle.net/10044/1/24122>

### a. Refereed Research Articles

Clicking on article DOI identifier will take you to the publication

- 1. Physics-constrained Bayesian inference of state functions in classical density-functional theory**  
P. Yatsyshin, S. Kalliadasis and Andrew B. Duncan (2022), *Journal of Chemical Physics* **156** 074105  
DOI: [10.1063/5.0071629](https://doi.org/10.1063/5.0071629)  
[Click for the short version, presented on NeurIPS 2020 workshop](#)
- 2. Surface nanodrops and nanobubbles: a classical density functional theory study**  
P. Yatsyshin and S. Kalliadasis (2021), *Journal of Fluid Mechanics* **913** 1  
DOI: [10.1017/jfm.2020.1167](https://doi.org/10.1017/jfm.2020.1167)
- 3. A finite-volume method for fluctuating dynamical density functional theory**  
A. Russo, S. P. Perez, M. A. Durán-Olivencia, P. Yatsyshin, J. A. Carrillo, S. Kalliadasis (2021), *Journal of Computational Physics* **428** 1  
DOI: [10.1016/j.jcp.2020.109796](https://doi.org/10.1016/j.jcp.2020.109796)
- 4. Mixing-demixing transition in polymer-grafted spherical nanoparticles**  
P. Yatsyshin, N. G. Fytas and E. Theodorakis (2020), *Soft Matter* **16** 703  
DOI: [10.1039/c9sm01639b](https://doi.org/10.1039/c9sm01639b)
- 5. Memory effects in fluctuating dynamic density-functional theory: theory and simulations**  
A. Russo, M. A. Durán-Olivencia, P. Yatsyshin and S. Kalliadasis (2020), *Journal of Physics A: Mathematical and Theoretical* **53** 445007  
DOI: [10.1088/1751-8121/ab9e8d](https://doi.org/10.1088/1751-8121/ab9e8d)

6. **Dynamics of the Desai-Zwanzig model in multiwell and random energy landscapes**  
S. N. Gomes, S. Kalliadasis, G. A. Pavliotis and P. Yatsyshin (2019), *Phys. Rev. E* **99** 032109  
DOI: [10.1103/PhysRevE.99.032109](https://doi.org/10.1103/PhysRevE.99.032109)
7. **General framework for nonclassical nucleation**  
M. A. Durán-Olivencia, P. Yatsyshin, S. Kalliadasis and J. F. Lutsko (2018), *New J. Phys.* **20** 083019  
DOI: [10.1088/1367-2630/aad170](https://doi.org/10.1088/1367-2630/aad170)
8. **Microscopic aspects of wetting using classical density-functional theory**  
P. Yatsyshin, M. A. Durán-Olivencia and S. Kalliadasis (2018), *J. Phys.: Condens. Matter* **30** 274003  
DOI: [10.1088/1361-648X/aac6fa](https://doi.org/10.1088/1361-648X/aac6fa)
9. **Wetting of a plane with a narrow solvophobic stripe**  
P. Yatsyshin, A. O. Parry, C. Rascón and S. Kalliadasis (2018), *Mol. Phys.* **116** 1990  
DOI: [10.1080/00268976.2018.1473648](https://doi.org/10.1080/00268976.2018.1473648)
10. **General framework for fluctuating dynamic density functional theory**  
M. A. Durán-Olivencia, P. Yatsyshin, B. D. Goddard and S. Kalliadasis (2017), *New J. Phys.* **19** 123022  
DOI: [10.1088/1367-2630/aa9041](https://doi.org/10.1088/1367-2630/aa9041)
11. **Nonequilibrium molecular dynamics simulations of nanoconfined fluids at solid-liquid interfaces**  
M. Morciano, M. Fasano, A. Nold, C. C. Braga, P. Yatsyshin, D. N. Sibley, B. D. Goddard, E. Chiavazzo, P. Asinari, and S. Kalliadasis (2017), *J. Chem. Phys.* **146** 244507  
DOI: [1.4986904](https://doi.org/10.1063/1.4986904)
12. **Classical density functional study of wetting transitions on nanopatterned surfaces**  
P. Yatsyshin, A. O. Parry, C. Rascón and S. Kalliadasis (2017), *J. Phys.: Condens. Matter* **29** 094001  
DOI: [1361-648X/aa4fd7](https://doi.org/10.1088/1361-648X/aa4fd7)
13. **Pseudospectral methods for density functional theory in bounded and unbounded domains**  
A. Nold, B. D. Goddard, P. Yatsyshin, N. Savva and S. Kalliadasis (2016), *J. Comp. Phys.* **334** 639  
DOI: [10.1016/j.jcp.2016.12.023](https://doi.org/10.1016/j.jcp.2016.12.023)
14. **Mean-field phenomenology of wetting in nanogrooves**  
P. Yatsyshin and S. Kalliadasis (2016), *Mol. Phys.* **114** 2688  
DOI: [10.1080/00268976.2016.1224393](https://doi.org/10.1080/00268976.2016.1224393)
15. **Complete Prewetting**  
P. Yatsyshin, A. O. Parry and S. Kalliadasis (2016), *J. Phys.: Condens. Matter* **28** 275001  
DOI: [10.1088/0953-8984/28/27/275001](https://doi.org/10.1088/0953-8984/28/27/275001)  

[Article was highlighted by reviewers as being “particularly significant to the community” and is featured at the journal's news blog **JPhys+**. This is a blog for articles that have wider appeal and interest as well as scientific importance and covers research news from across the *Journal of Physics* series.]

URL: <https://jphysplus.iop.org/2016/06/23/derjaguin-in-flatland-prewetting-spreads-out/> ]
16. **Density functional study of condensation in capped capillaries**  
P. Yatsyshin, N. Savva and S. Kalliadasis (2015), *J. Phys.: Condens. Matter* **27** 275104  
DOI: [10.1088/0953-8984/27/27/275104](https://doi.org/10.1088/0953-8984/27/27/275104)

17. **Wetting of prototypical one- and two-dimensional systems: thermodynamics and density functional theory**  
P. Yatsyshin, N. Savva and S. Kalliadasis (2015), *J. Chem. Phys.* **142** 034708  
DOI: [10.1063/1.4905605](https://doi.org/10.1063/1.4905605)
18. **Unification of dynamic density functional theory for colloidal fluids to include inertia and hydrodynamic interactions: derivation and numerical experiments**  
B. Goddard, A. Nold, N. Savva, P. Yatsyshin, and S. Kalliadasis (2013), *J. Phys.: Condens. Matter* **25** 035101  
DOI: [10.1088/0953-8984/25/3/035101](https://doi.org/10.1088/0953-8984/25/3/035101)  
[Article featured as a news item on the Imperial College website  
URL: <http://tinyurl.com/z9mm4rt>]
19. **Geometry-induced phase transition in fluids: Capillary prewetting**  
P. Yatsyshin, N. Savva, and S. Kalliadasis (2013), *Phys. Rev. E (Rapid Comm)*, **87**, 020402(R).  
DOI: [10.1103/PhysRevE.87.020402](https://doi.org/10.1103/PhysRevE.87.020402)
20. **Spectral methods for the equations of classical density-functional theory: relaxation dynamics of microscopic films**  
P. Yatsyshin, N. Savva, and S. Kalliadasis (2012), *J. Chem. Phys.*, **136**, 124113  
DOI: [10.1063/1.3697471](https://doi.org/10.1063/1.3697471)
21. **Structure of electron-positron clusters: Hartree-Fock approximation**  
P. Yatsyshin, R. G. Polozkov, V. K. Ivanov, and A. V. Solovyov (2009), *Phys. Scr.*, **80**, 048126  
DOI: [10.1088/0031-8949/80/04/048126](https://doi.org/10.1088/0031-8949/80/04/048126)
22. **Resonances in the cross section of photodetachment of  $2p$  electrons from negative ions  $Na^-$**   
V. K. Ivanov and P. I. Yatsyshin (2009), *Technical Physics*, **54**, 7  
DOI: [10.1134/S1063784209010022](https://doi.org/10.1134/S1063784209010022)

## b. Book Chapters

23. **Classical Density Functional Theory and Nanofluidics: Adsorption and the Interface Binding Potential**  
P. Yatsyshin, M.-A. Durán-Olivencia and S. Kalliadasis  
In: Klaus Sattler (ed) *21st Century Nanoscience. A Handbook* Chapter 14 CRC Press (2020)  
DOI: [10.1201/9780429347313-14](https://doi.org/10.1201/9780429347313-14)
24. **Classical density-functional theory studies of fluid adsorption on nanopatterned planar surfaces**  
P. Yatsyshin and S. Kalliadasis (2018)  
In: L.L. Bonilla, E. Kaxiras and R. Melnik (eds) *Coupled Mathematical Models for Physical and Biological Nanoscale Systems and Their Applications*. BIRS-16w5069 2016. Springer Proceedings in Mathematics & Statistics, vol 232. Springer, Cham

## c. Conference Papers

25. **Data-driven density functional theory: a case for physics-informed learning**  
P. Yatsyshin, S. Kalliadasis and A. B. Duncan (2021)  
*Machine Learning and the Physical Sciences Workshop at the 35th Conference on Neural Information Processing Systems (NeurIPS)*

Workshop website

Conference Article

Video Introduction

## 26. Equilibrium fluid structures in prototypical nanosystems

P. Yatsyshin, N. Savva and S. Kalliadasis (2016),

*8th GRACM International Congress on Computational Mechanics*, University of Thessaly, Volos (Greece), July 2015

URL: <http://8gracm.mie.uth.gr/Papers/Session%20D1-B2/P.%20Yatsyshin.pdf>

## d. Conference Presentation

### Invited talks and seminars

1. *Data-Driven Classical Density Functional Theory: A Case for Physics Informed Learning*, New directions in classical density functional theory workshop, International Center for Mathematical Sciences, UK, April 2021  
URL: <https://www.icms.org.uk/events/workshops/cdft>  
Video of presentation
2. *Machine-learning the DFT of a classical statistical-mechanical system: A case for physics-informed learning*, Density Functional Days in Tübingen, Tübingen University, Germany, September 2020
3. *Wetting on striped walls: interplay between pre-wetting and interface unbending*, Density Functional Days in Tübingen, Tübingen University, Germany, September 2019
4. *Statistical Mechanics of Wetting*, Open Statistical Physics, School of Mathematics and Statistics, Open University, UK, March 2018
5. *Computational statistical mechanical framework for soft condensed matter. Mean-field description of wetting at the nanoscale*, Chemical Engineering Postdoc Symposium, Department of Chemical Engineering, Imperial College London, UK, April 2017.
6. *A computational statistical-mechanical framework: Classical density functional theory with applications to phase transition in nano-confined fluids*, Department of Mathematics & Statistics, The Open University, UK, November 2016.
7. *Statistical Mechanics of classical fluids: Density functional theory and equilibrium and dynamics of wetting*, Physics of Fluids group, University of Twente, Enschede, The Netherlands, in October 2016.
8. *Wetting at the nanoscale. Equilibrium and dynamics*  
Video available at URL: (the link below is clickable in the .pdf file)  
<http://www.birs.ca/events/2016/5-day-workshops/16w5069/videos/watch/201608301028-Yatsyshin.html>,  
2016 BIRS Workshop on Coupled Mathematical Models for Physical and Biological Nanoscale Systems and Their Applications, in Banff, Alberta, Canada, August 2016.
9. *Implementing computations with equilibrium and dynamic classical density functional theory*, Department of Mathematical Sciences, Loughborough University, Leicestershire, UK, in May 2016.
10. *Classical density functional theory for Lennard-Jones fluids. Applications to wetting on planar and sculpted substrates*, Molecular Systems Engineering group, Imperial College London, UK, in April 2016.
11. *Density functional theory for surface tensions and more*, presented within the four-part seminar "Exploration of fluids", Imperial College London, October 2014, chaired by Prof. S. H. Davis, Northwestern University, Royal Academy of Engineering Visiting Professor to the Department of Chemical Engineering, Imperial College London.



12. *Geometry-induced phase transitions*, Chemical Engineering Ph.D. symposium, Imperial College London, UK, June 2013.

### Contributed talks

13. M. A. Durán-Olivencia, A. Russo, Yatsyshin, P. and S. Kalliadasis *Memory effects in fluctuating dynamic density-functional theory: theory and simulations*, 74th Annual Meeting of the APS Division of Fluid Dynamics, Phoenix (ARI, USA), November 2021
14. Yatsyshin, P., S. Kalliadasis, and A. B. Duncan *Data-Driven Classical Density Functional Theory: A Case for Physics Informed Learning*, APS March Meeting, Virtual meeting due to Covid-19 (USA), March 2021
15. Yatsyshin, P., A. O. Parry, C. Rascón, M. A. Durán-Olivencia, and S. Kalliadasis *Phase transitions at the interfaces and in the bulk. Equilibrium and dynamics studies using classical density functional theory*, Liblice 2018 (Czech Republic), June 2018
16. Yatsyshin, P., M. A. Durán-Olivencia, A. O. Parry, C. Rascón and S. Kalliadasis *Understanding interfacial wetting transitions with classical density functional theory*, APS March Meeting, Los Angeles (CA, USA), March 2018
17. Yatsyshin, P., A. O. Parry, C. Rascón and S. Kalliadasis *Wetting of chemically nanopatterned walls*, Workshop of the Fundamental Theoretical Approaches to the Equation of State, Manchester (UK), January 2018
18. Yatsyshin, P., A. O. Parry, C. Rascón, M. A. Durán-Olivencia, and S. Kalliadasis *Wetting of heterogeneous substrates. A classical density-functional-theory approach*, 70th Annual Meeting of the APS Division of Fluid Dynamics, Denver (CO, USA), November 2017
19. Yatsyshin, P., M. A. Durán-Olivencia, A. O. Parry, C. Rascón and S. Kalliadasis *Mean-field treatment of wetting at the nanoscale*, to be given at the summer school Complex Motion in Fluids 2017, Cambridge (UK), September 2017
20. Yatsyshin, P., A. O. Parry, C. Rascón and S. Kalliadasis *Classical density functional study of wetting transitions on nanopatterned surfaces*, Thermodynamics 2017, Edinburgh (UK), August 2017
21. Yatsyshin, P., M. A. Durán-Olivencia, A. O. Parry, C. Rascón and S. Kalliadasis *Wetting in flatland: Complex interfacial transitions at inhomogeneous solid-gas interfaces*, APS March Meeting, New Orleans (LA, USA), March 2017
22. Yatsyshin, P., D. N. Sibley, M. A. Durán-Olivencia and S. Kalliadasis, *Dynamics of two-phase interfaces and surface tensions: A density-functional theory perspective*, 69th Annual Meeting of the APS Division of Fluid Dynamics, Portland (OR, USA), November 2016
23. Yatsyshin, P., N. Savva, A. Nold, B. D. Goddard and S. Kalliadasis, *Mapped-Chebyshev spectral collocation approach to the integral and integral-differential equations of the classical equilibrium and dynamic density functional theory*, British Applied Mathematical Colloquium, University of Oxford, Oxford (UK), April 2016
24. Yatsyshin, P., A. O. Parry and S. Kalliadasis, *A computational DFT study of structural transitions in textured solid-fluid interfaces*, 68th Annual Meeting of the APS Division of Fluid Dynamics, Boston (MA, USA), November 2015
25. Yatsyshin, P., D. N. Sibley, N. Savva and S. Kalliadasis, *Molecular-level description of nano-drops: contact angles, dynamics of wetting and coalescence*, Droplets 2015, University of Twente, Enschede (Netherlands), October 2015
26. Yatsyshin, P., N. Savva and S. Kalliadasis, *Thermodynamics and statistical mechanics of wetting transitions: fluid phase behavior in prototypical nanostructured substrates*, Thermodynamics 2015, Technical University of Denmark, Copenhagen (Denmark), September 2015

27. Yatsyshin, P., N. Savva and S. Kalliadasis, *Equilibrium fluid structures in prototypical nanosystems*, 8th GRACM International Congress on Computational Mechanics, University of Thessaly, Volos (Greece), July 2015
28. Yatsyshin, P., D. N. Sibley, N. Savva and S. Kalliadasis, *Droplets and the three-phase contact line at the nano-scale. Statics and dynamics*, 67th Annual Meeting of the APS Division of Fluid Dynamics, San Francisco (CA, USA), November 2014
29. Yatsyshin, P., N. Savva and S. Kalliadasis, *Capillary Condensation Revisited: Wetting on a Capped Capillary*, Thermodynamics 2013, University of Manchester, Manchester (UK), September 2013
30. Yatsyshin, P., N. Savva and S. Kalliadasis, *Micro confined inhomogeneous fluids. Wetting on a capped capillary*, British Applied Mathematical Colloquium, University of Leeds, Leeds (UK), April 2013
31. Yatsyshin, P., N. Savva and S. Kalliadasis, *Relaxational Dynamics of Microscopic films. Spectral Methods for the Equations of Classical Density Functional Theory*, Summer School on Wave Patterns and Interactions in Advection-Dominated Flows, University of Thessaly, Volos (Greece), July 2012
32. Yatsyshin, P., N. Savva and S. Kalliadasis, *Integral and Integral-Differential equations of classical density functional theory. A novel numerical approach*, British Applied Mathematical Colloquium, University College London, London (UK), March 2012
33. Yatsyshin, P., N. Savva and S. Kalliadasis, *Dynamics of a thin film layer on planar substrate obtained from a dynamic density functional theory approach*, International Conference on Multiscale Complex Fluid Flows and Interfacial Phenomena, Free University of Brussels, Brussels (Belgium), November 2010