

**Guillaume Chevrot**

140 Allée de Fakenham, 45160 Olivet (FRANCE)  
 +33 6 25 42 76 56 (cell)  
 guillaume.chevrot@gmail.com  
<https://gchevrot.github.io/home/>

**Computational scientist****Professional background**

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- 09/2015 – Present** Research Engineer  
 MISC, Complex Systems Institute  
 University of Orléans  
 • Combining ActivePapers and Exec&Share.  
 • Statistical analysis and modeling neural signal.  
 • Text mining.
- 04/2014 – 08/2015** Postdoctoral Research Associate  
 MEMPHYS, Center for Biomembrane Physics  
 University of Southern Denmark (SDU)  
 Khandelia group  
 • Enhanced stability of proteins in 3<sup>rd</sup> generation Ionic liquids.  
 • Behavior of the motor-protein Kinesin-5 with an inhibitor.  
 • Lipid droplet genesis.
- 01/2011 – 02/2014** Postdoctoral Research Associate  
 CBM-CNRS / SOLEIL synchrotron  
 Kneller group  
 • Extraction of internal motions from molecular dynamics simulations.  
 • Impact of anisotropic atomic motions on incoherent neutron scattering intensities.  
 • New approach to molecular diffusion tensors.
- 02/2008 – 08/2010** Postdoctoral Research Associate  
 CEA-DAM Ile de France  
 Soulard group  
 • Molecular dynamics simulations of carbon nanoparticles under detonation conditions.  
 • Modeling the kinetics of carbon nanoparticles coalescence.  
 • Development of programs to analyze molecular dynamics simulations.
- 10/2004 – 01/2008** Ph.D. in computational chemistry  
 University of Strasbourg, France  
 Modeling and Molecular Simulations laboratory  
 Wipff group  
 • Molecular dynamics simulations of dicarbolides anions, BTP and 18C6 extractants.  
 • Simulations in pure liquids and at liquid-liquid interfaces.  
 • Study at classical interfaces and at [BMI][PF<sub>6</sub>] ionic liquids interfaces.
- 01/2004 – 06/2004** Intern (M2 - computational chemistry)  
 University of Strasbourg, France  
 Modeling and Molecular Simulations laboratory  
 Wipff group  
 • Surfactant behavior of dicarbolides anions at interfaces: a molecular dynamics study.
- 04/2003 – 08/2003** Intern (M1 – chemistry)  
 CEA Valduc. Dijon Area, France  
 Hubinois group  
 • Gas analysis by mass spectrometry and gas chromatography.
- 06/2002 – 09/2002** Laboratory assistant. Dijon Area, France.  
 Sanofi pharmaceutical group.  
 • Water, drug and raw materials analysis.

## Educational background

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<b>2004 – 2008</b>	Ph.D. in computational chemistry. University of Strasbourg, France.
<b>2003 – 2004</b>	Master degree (M2) in computational chemistry. University of Strasbourg, France.
<b>2002 – 2003</b>	Master degree (M1) in chemistry. University of Dijon, France.
<b>07/2009</b>	Methods in molecular simulation summer school 2009 (5 - 14 <sup>th</sup> july). University of Sheffield, United Kingdom.

## Skills

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<ul style="list-style-type: none"> <li>- Extensive experience in molecular simulations.</li> <li>- Programming experience in Python, R, C, Shell scripting, Mathematica, Django, HTML, CSS, Git, Drupal.</li> <li>- Experience working with large data sets (HDF5, NetCDF).</li> <li>- HPC experience.</li> <li>- English: full professional proficiency TOEIC test (11-28-06): 885/990.</li> </ul>	<ul style="list-style-type: none"> <li>- 11 articles published in international peer-reviewed journals.</li> <li>- Oral and poster communications in international and national congress.</li> <li>- Contribution to the european project « europart ».</li> <li>- Technological monitoring.</li> <li>- Preparing and delivering reports, presentations.</li> </ul>
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## Publications

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**G. Chevrot**, E.E. Fileti, and V.V. Chaban. *Enhanced stability of the model mini-protein in amino acid ionic liquids and their aqueous solutions*. J. Comput. Chem., **2015**, 36, 2044.

**G. Chevrot**, K. Hinsén, and G.R. Kneller. *Model-free simulation approach to molecular diffusion tensors*. J. Chem. Phys., **2013**, 139, 154110.

G.R. Kneller and **G. Chevrot**. *Impact of anisotropic atomic motions in proteins on powder-averaged incoherent neutron scattering intensities*. J. Chem. Phys., **2012**, 137, 225101.

**G. Chevrot**, A. Sollier, and N. Pineau. *Molecular dynamics and kinetic study of carbon coagulation in the release wave of detonation products*. J. Chem. Phys., **2012**, 136, 084506.

**G. Chevrot**, P. Calligari, K. Hinsén, and G.R. Kneller. *Least constraint approach to the extraction of internal motions from molecular dynamics trajectories of flexible macromolecules*. J. Chem. Phys., **2011**, 135, 084110.

J.H. Los, N. Pineau, **G. Chevrot**, G. Vignoles and J.-M. Leyssale. *Formation of multiwall fullerenes from nanodiamonds studied by atomistic simulations*. Phys. Rev. B, **2009**, 155420.

**G. Chevrot**, E. Bourasseau, N. Pineau and J.-B. Maillet. *Molecular dynamics simulations of nanocarbons at high pressure and temperature*. Carbon, **2009**, 47, 3392-3402.

**G. Chevrot**, R. Schurhammer and G. Wipff. *Molecular dynamics study of dicarbolide anions in nitrobenzene solution and at its aqueous interface. Synergistic effect in the Eu(III) assisted extraction*. Phys. Chem. Chem. Phys., **2007**, 9, 5928-5938.

**G. Chevrot**, R. Schurhammer and G. Wipff. *Synergistic effect of dicarbolide anions in liquid - liquid extraction: a molecular dynamics study at the octanol – water interface*. Phys. Chem. Chem. Phys., **2007**, 9, 1991-2003.

**G. Chevrot**, R. Schurhammer and G. Wipff. *Molecular dynamics simulations of the aqueous interface with the [BMI][PF6] ionic liquid: comparison of different solvent models*. Phys. Chem. Chem. Phys., **2006**, 8, 4166-4174.

**G. Chevrot**, R. Schurhammer and G. Wipff. *Surfactant behavior of ‘ellipsoidal’ dicarbolide anions: a molecular dynamics study*. J. Phys. Chem. B, **2006**, 110, 9488-9498.