


**Information to be requested from all CA17104 participants:**

	
<p><b>Indicate your Working Group(s) in COST Action17104:</b></p>	<p>WG2</p>
<p><b>First Name:</b></p>	<p>Patrick</p>
<p><b>Surname:</b></p>	<p>Trouillas</p>
<p><b>Department</b></p>	<p>Department of Physical Chemistry</p>
<p><b>Primary Institution</b></p>	<p>Palacký University Olomouc - Regional Centre of Advanced Technologies and Materials</p>
<p><b>Address of Primary Institution</b></p>	<p>tr. 17. listopadu 12, 771 46 Olomouc, Czech Republic</p>
<p><b>Other institutions</b></p>	<p>INSERM U1248 - Limoges University, France</p>
<p><b>Telephone:</b></p>	<p>+33631175812</p>
<p><b>e-mail:</b></p>	<p><a href="mailto:patrick.trouillas@upol.cz">patrick.trouillas@upol.cz</a> <a href="mailto:patrick.trouillas@unilim.fr">patrick.trouillas@unilim.fr</a></p>
<p><b>Link to webpage with biography:</b></p>	<p><a href="https://www.unilim.fr/ippritt/people/patrick-trouillas_new/">https://www.unilim.fr/ippritt/people/patrick-trouillas_new/</a></p>

<p><b>Link to webpage with group description:</b></p>	<p><a href="https://www.rcptm.com/groups/carbon-nanostructures-and-biomacromolecules/">https://www.rcptm.com/groups/carbon-nanostructures-and-biomacromolecules/</a></p> <p><a href="https://www.unilim.fr/ippritt/">https://www.unilim.fr/ippritt/</a></p>
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<p><b>Orcid ID or Scopus ID</b></p>	<p>orcid.org/0000-0002-8832-9606</p>
<p><b>Linkedin</b></p>	<p><i>optional</i></p>
<p><b>Expertise relevant for this COST Action:</b></p>	<p>Molecular Modelling  Molecular Dynamics  Quantum chemistry  3D models of ABC and SLC transporters  Docking of drugs or any xenobiotics in the prot patr ein chamber of these transporters  Elucidation of large conformational changes in membrane transporters  3D models of biological membranes (in silico evaluation of drug partitioning and permeation)  (In silico) Conformational analysis</p>
<p><b>Available facilities to conduct work, relevant for this COST Action:</b></p>	<p>Super-computers</p>
<p><b>Materials/Methods that could be shared with other members of this COST Action:</b></p>	<p>On-demand molecular modeling simulations</p>

NOTE: By submitting this form to the Grant Manager of CA17104, I agree that this information can be used within the scope of this COST Action (e.g. may be included on the webpage of CA17104).