Information to be requested from all CA17104 participants:

	
Indicate your Working Group(s) in COST Action17104:	**WG2**
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Expertise relevant for this COST Action:	Molecular Dynamics simulations, pH effects in MD simulations, Computational structural biology, Computational biophysics, Molecular docking calculations
Available facilities to conduct work, relevant for this COST Action:	Computational Cluster with ~2000 cores
Matherials/Methods that could be shared with other members of this COST Action:	Proteins and membranes structural studies, pH-dependent conformational transitions, Lipid bilayers permeation calculations

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).