Information to be requested from all CA17104 participants:

	**PHOTO**
Indicate your Working Group(s) in COST Action17104:	(WG4)
First Name:	Ilza
Surname:	Pajeva
Department	QSAR & Molecular Modelling
Primary Institution	Institute of Biophysics and Biomedical Engineering, Bulgarian Academy of Sciences
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Link to webpage with biography:	*optional*

Link to webpage with group description:	http://biomed.bas.bg/en/departments/qsar-and-molecular-modelling/

Orcid ID or Scopus ID	https://orcid.org/0000-0001-5412-1586
Linkedin	*optional*
Expertise relevant for this COST Action:	Application of ligand- and structure-based methods for in silico drug design and computational toxicology
Available facilities to conduct work, relevant for this COST Action:	HPC cluster and specialized software for molecular modelling (e.g. MOE, GOLD, Schroedinger, free software)
Matherials/Methods that could be shared with other members of this COST Action:	In silico models /predictions of pharmacological and toxicological effects of bioactive molecules

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).