



## Biologia Computazionale e di Sistema

Incontro annuale del Gruppo di interesse della Società Italiana di Biochimica e Biologia Molecolare (SIB)

Co-organizzato da Progetto Bandiera InterOmics  
e Istituto di Biologia e Patologia Molecolari del CNR

**CNR, Aula Bisogno, Piazzale Aldo Moro 7 – 00185 Roma**

**10:15-10:30 – Welcome**

**10:30-13:20 – Session 1: Genome analysis, mutations and disease**

10:30-11:30 – Marco Punta (The Institute for Cancer Research, London - Centre for Evolution and Cancer)  
*Cancer genomics, cancer vaccines and cancer resistance mutations*

11:30-11:50 – Fabio Polticelli (Università di Roma Tre)  
*Identification and function prediction of an elusive HEAT repeats domain of Huntingtin*

11:50-12:10 – Giovanni Minervini (Università di Padova)  
*Genotype-phenotype relations of the von Hippel-Lindau tumor suppressor inferred from a large-scale analysis of disease mutations and interactors*

**12:10-12:40 – Coffee-Break**

12:40-13:00 – Valerio Marino (Università di Verona)  
*Evolutionary-conserved structural communication in two Neuronal Calcium Sensor proteins investigated by SNP analysis*

13:00-13:30 – Pasqualina D'Ursi – Federica Chiappori (CNR-ITB)  
*Computational studies and surface plasmon resonance in rescuing pathway impairments in CFTR: identification of key druggable targets by interaction network analysis and drug repositioning.*

**13:30-17:00 – Session 2: Protein structure, function and dynamics**

13:30-13:50 – Castrense Savojardo (Università di Bologna)  
*Computing protein-protein interaction sites from structure and sequence*

13:50-14:10 – Marco Necci (Università di Padova)  
*Integrated resources for the study of protein intrinsic disorder*

**14:10-15:30 – Lunch**

15:30-15:50 – David Sasah Staid (Università "Sapienza" di Roma)  
*Identification of point mutations related to drug resistance in bacteria*

15:50-16:10 – Giacomo Janson (Università "Sapienza" di Roma)  
*How to improve protein homology modelling within MODELLER*

16:10-16:30 – Leonardo Guidoni (Università dell'Aquila)  
*Atomistic simulations in Biochemistry*

16:30-17:00 – Veronica Morea, Allegra Via (CNR-IBPM)  
*Activities of the Bioinformatics group @IBPM-CNR*

**17:00-17:15 – Conclusions**