Structures for a new era of drug discovery

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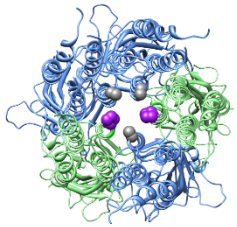
11th February 2016



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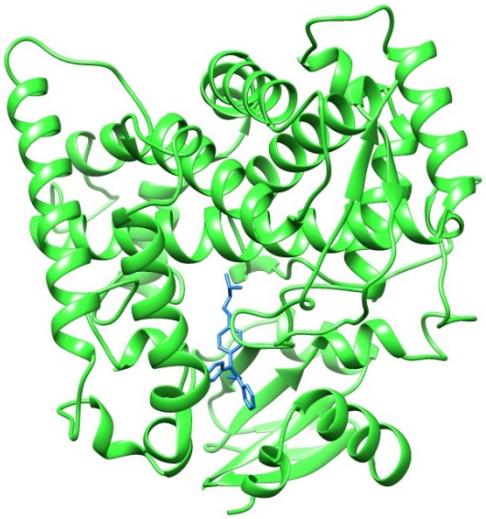
working molecular knowledge across genomes

What We Do



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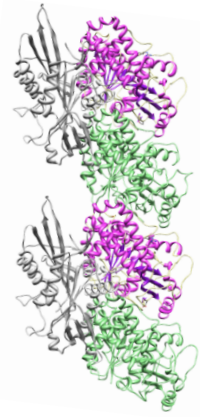
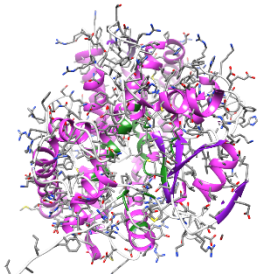
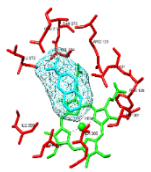
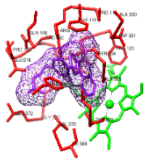


Genome / Multi-Genome Scale :

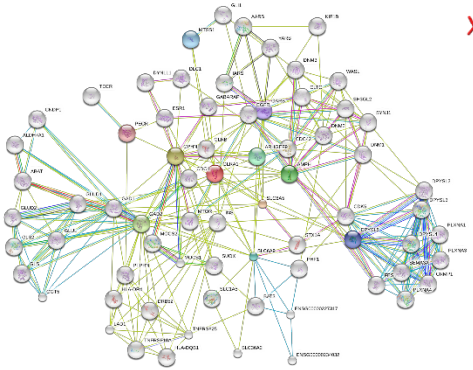




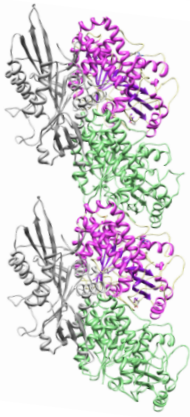
- Structural modelling at vast scales
- Affinity docking of vast compound libraries “Rapid *in silico* exploration”
- Pharmacological profiling and toxicity screening, precision medicine



- Prediction of protein-protein interactions and networks



- *In silico* pathway analysis
- Comparative structural proteomics



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Partners - Projects - Products

TargetPath - a lead discovery platform for antibiotics

Host-Pathogen Interactomes

In silico pharmacological profiling panel

In silico toxicity screening panel

[dstl]

Centre for Defence Enterprise

Unilever

NC 3Rs

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Partners - Projects - Products

In silico analysis of pipeline drug compounds

Pharmacological profiling of Brazilian natural compounds

Tumour-specific cancer drug targets based on exome sequences

Institute NEO MED

Universidade Federal Fluminense

International Cancer Genome Consortium

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What Makes Us Different……

- R&D based on a track record of 50+ journal publications
- Validated workflows in partnership with target industries

"Moleculomics is the only company offering structurally-based open-ended lead discovery, toxicity screening and protein network identification, that is completely unlimited in reach, applied across the entire proteome."

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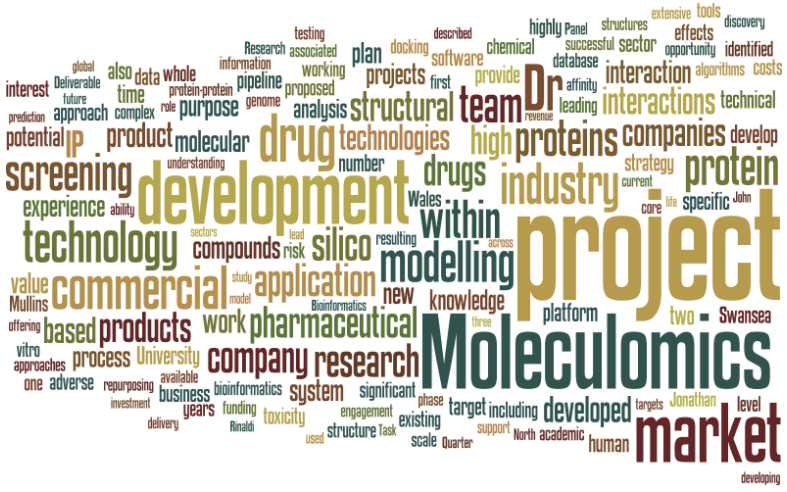
Swansea University
Prifysgol Abertawe

THE UNIVERSITY RESEARCH UNDERGRADUATE POSTGRADUATE

Genome and Structural Bioinformatics



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Welsh Government