

## [pyDockCG: new coarse-grained potential for protein-protein docking.](#)

**Authors:** [Solernou, Albert](#) / [Fernández-Recio, Juan](#)

**Publication:** The journal of physical chemistry. B

**Volume / Pagination:** 115 / 6032-9

**Palabras clave:** [Algorithms](#), [Computer Simulation](#), [Multiprotein Complexes](#), [Protein Binding](#), [Protein Interaction Mapping](#), [Solvents](#), [Static Electricity](#), [Thermodynamics](#)

Barcelona Supercomputing Center - Centro Nacional de Supercomputación

---

**Source URL (retrieved on 14 Dic 2024 - 13:17):** <https://www.bsc.es/es/research-and-development/publications/pydockcg-new-coarse-grained-potential-protein-protein-docking>