



Linear scaling DFT calculations for large tungsten systems using an optimized local basis

URL: <https://www.sciencedirect.com/science/article/pii/S235217911730090X?via%3Dihub>

Authors: Mohr, Stephan / Eixarch, Marc / Amsler, Maximilian / Mantsinen, Mervi / Genovese, Luigi

Research Lines: [Ab-Initio Electronic Structure Methods](#) / [Computational Modeling for Fusion](#)

Publication: Nuclear Materials and Energy

Volume / Pagination: 15 / 64-70

Barcelona Supercomputing Center - Centro Nacional de Supercomputación

Source URL (retrieved on 23 Abr 2024 - 10:06): <https://www.bsc.es/es/research-and-development/publications/linear-scaling-dft-calculations-large-tungsten-systems-using>