

ONETEP: Linear Scaling DFT with Plane-Waves – Methods and Applications –

Arash A. Mostofi¹, Peter D. Haynes², Chris-Kriton Skylaris³
and Mike C. Payne²

¹*Department of Materials Science and Engineering,
Massachusetts Institute of Technology, Cambridge MA, USA*

²*Cavendish Laboratory, University of Cambridge, Cambridge UK*

³*Physical and Theoretical Chemistry Laboratory,
University of Oxford, Oxford UK*

We present ONETEP[1] (Order- N Electronic Total Energy Package), a density functional method, based on plane-waves, whose computational cost scales only linearly with the number of atoms.

Although at first sight the extended nature of plane-waves makes them appear unsuitable for representing the localised orbitals used in linear scaling methods – which is why localised basis sets such as Gaussians, numerical atomic orbitals, spherical waves and B-splines have been used commonly instead – we have developed a linear-scaling method based on plane-waves which overcomes this difficulty and which can achieve the same accuracy and convergence rate as traditional plane-wave calculations[2].

We describe the main features of our method which result in its success, and present recent developments and results for realistic applications.

[1] *J. Chem. Phys.* **122**, 084119 (2005).

[2] *J. Chem. Phys.* **119**, 8842 (2003).