```
client199-on-tre:~ grass$ Espresso
0: Script directory: /Users/grass/workspace/eclipse/programs/Espresso/scripts

    Espresso -

                                A MPI Parallel Molecular Dynamics Program
* (c) 2002-2006
* Max-Planck-Institute for Polymer Research
* Mainz, Germany
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>code_info
ESPResSo: 2.0.4s, Last Change: October 11th, 2007
{ Compilation status { MPI fake } { FFTW3 } { PARTIAL_PERIODIC } { ELECTROSTATICS } { EXTE
RNAL_FORCES } { CONSTRAINTS } { LENNARD_JONES } { BOND_ANGLE_COSINE } { LB } }
{ Debug status { } }
```