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# TENTH INTERNATIONAL WORKSHOP ON COMPUTATIONAL CONDENSED MATTER PHYSICS: TOTAL ENERGY AND FORCE METHODS | (smr 1300)

## Saturday 13 January 2001

### First-principle molecular-dynamics simulation of minerals at deep Earth's temperatures and pressures - Main Building Main Lecture Hall (15:00-15:30)

| time | title | presenter |
|---|---|---|
| 15:00 | First-principle molecular-dynamics simulation of minerals at deep Earth's temperatures and pressures | S. SCANDOLO |