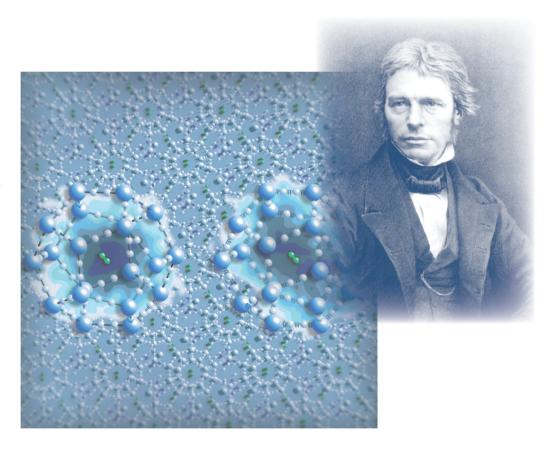
Hydrogen Storage Materials

Rutherford Appleton Laboratory, Didcot, Oxon, United Kingdom 18–20 April 2011



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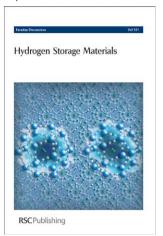
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See Peterson et al., Faraday Discuss., 2011, **151**, 37-46.

The potential-energy surfaces, calculated from ab-initio molecular-dynamics, for the center of mass of a H2 guest in the H2/THFclathrate composite, which differ for flexible (right) and rigid (left) H2O-cage calculations.

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