

First-principles study of the electronic structure of heavy fermion YbRh₂Si₂

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2006 J. Phys.: Condens. Matter 18 10529

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Erratum

First-principles study of the electronic structure of heavy fermion YbRh_2Si_2 T Jeong 2006 *J. Phys.: Condens. Matter* **18** 6289–6297

The band structure in figure 2(b) and density of states in figure 3(b) for LDA +U calculation are incorrect and should be replaced by the new figures below. This change does not affect the conclusions of the paper. W E Pickett is removed from the list of authors.

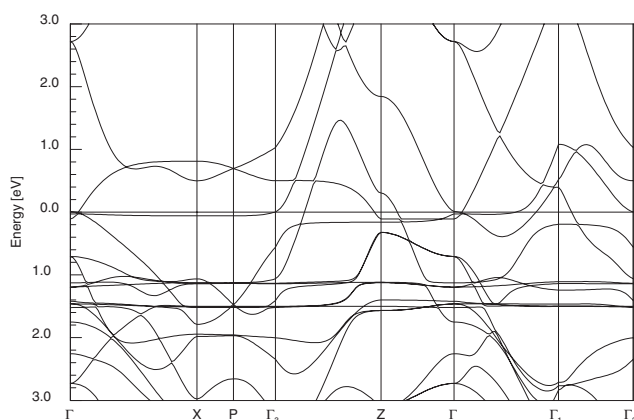


Figure 2(b). The band structure of YbRh_2Si_2 for LDA+U calculation.

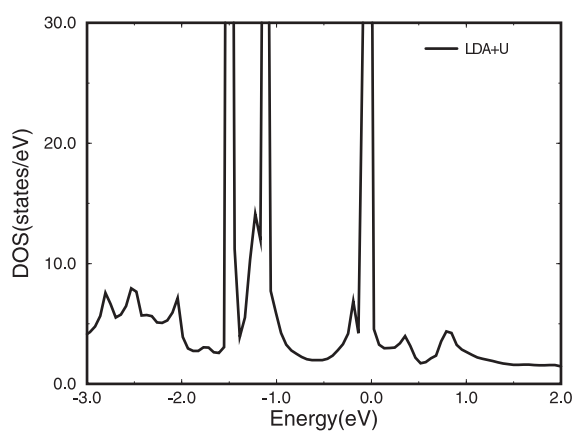


Figure 3(b). The density of states of YbRh_2Si_2 for LDA+U calculation.