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

2006 J. Phys.: Condens. Matter 18 10529

(http://iopscience.iop.org/0953-8984/18/46/C01)

J. Phys.: Condens. Matter **18** (2006) 10529 doi:10.1088/0953-8984/18/46/C01

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.

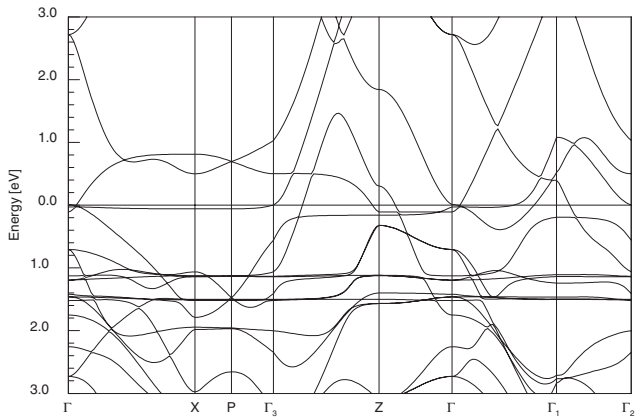

Figure 2(b). The band structure of $YbRh_2Si_2$ for LDA+U calculation.

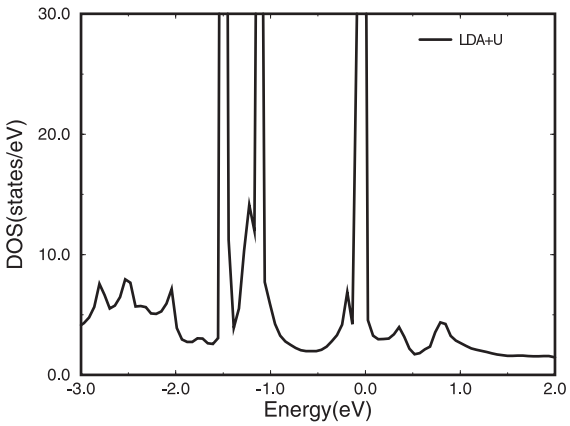

Figure 3(b). The density of states of $YbRh_2Si_2$ for LDA+U calculation.