

Erratum: *Ab initio* modeling of the two-dimensional energy landscape of screw dislocations in bcc transition metals [Phys. Rev. B **89**, 024104 (2014)]

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The units and values of the dislocation core energies in Table II were incorrect due to a conversion unit error. Here are the correct values, given in $\text{meV}/\text{\AA}$.

This correction consequently shifts all the normalized values in Fig. 4, but all the conclusions written in the paper remain unchanged.

	E_{core} ($\text{meV}/\text{\AA}$)
V	116
Nb	171
Ta	153
Mo	416
W	501
Fe	206

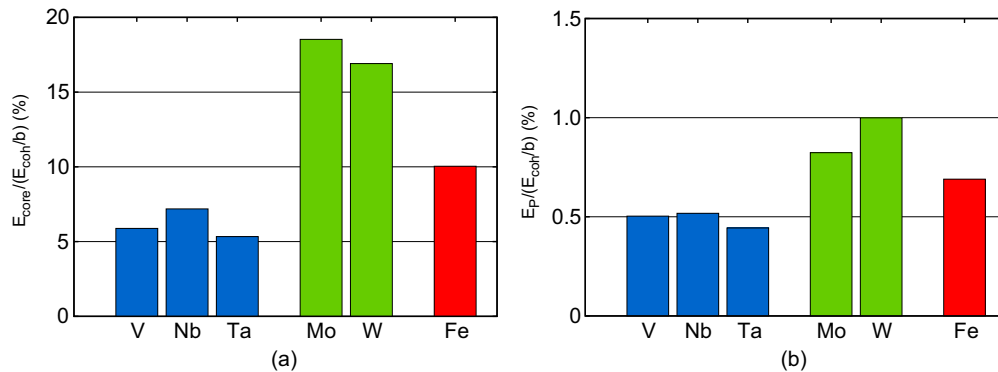


FIG. 4. (Color online) Group dependence of the normalized energy for (a) easy core energy and (b) Peierls energy.