

Erratum

Erratum to: Kinetic measurement and prediction
of the hydrogen outgassing from the polycrystalline
LiH/Li₂O/LiOH system by
L.N. Dinh et al., J. Nucl. Mater. 347 (2005) 31

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Eq. (3) of this paper needs the following clarification and correction.

The exact form of Eq. (3) should be:

$$\int_0^{\alpha} \ln \left(\frac{d\alpha}{dT/\beta} \right) d\alpha = -\frac{1}{R} \int_0^{\alpha} \frac{E d\alpha}{T} + \int_0^{\alpha} \ln(vf(\alpha)) d\alpha$$

If the activation energy barrier of the process under study is constant, the above equation can be simplified to Eq. (3) in the paper. Eq. (3) can be used as a test to see if the activation energy barrier of the process under study is constant or not. However, using Eq. (3) when E varies with α (as done in this paper) is equivalent to approximating the value of E from 0 to α with a constant E . A more accurate approach to numerically approximate values of E vs. α is as following:

$$\int_0^{\alpha_i} \ln \left(\frac{d\alpha}{dt} \right) d\alpha + \int_{\alpha_i}^{\alpha_{i+1}} \ln \left(\frac{d\alpha}{dt} \right) d\alpha + \dots = \left[-\frac{E_i}{R} \int_0^{\alpha_i} \frac{d\alpha}{T} + \int_0^{\alpha_i} \ln(vf(\alpha)) d\alpha \right] \\ + \left[-\frac{E_{i+1}}{R} \int_{\alpha_i}^{\alpha_{i+1}} \frac{d\alpha}{T} + \int_{\alpha_i}^{\alpha_{i+1}} \ln(vf(\alpha)) d\alpha \right] + \dots$$

given the α_i to α_{i+1} interval is small.

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