

Osz, Judit, Gabriella Bodó, Rui Miguel Mamede Branca, and Csaba Bagyinka. 2005. Theoretical calculations on hydrogenase kinetics: explanation of the lag phase and the enzyme concentration dependence of the activity of hydrogenase uptake. *Biophys. J.* 89:1957–1964.

Equation 3 printed incorrectly. The correct version is below.

$$\begin{aligned}\dot{E}_1 &= -aE_1 \\ \dot{E}_2 &= aE_1 - bE_2E_3 + dE_4H_2M_0^2 \\ \dot{E}_3 &= +bE_2E_3 - cE_3 \\ \dot{E}_4 &= cE_3 - dE_4H_2M_0^2 \\ \dot{H}_2 &= -dE_4H_2M_0^2 \\ \dot{M}_0 &= -2dE_4H_2M_0^2 \\ \dot{M}_r &= 2dE_4H_2M_0^2\end{aligned}$$

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doi: 10.1529/biophysj.105.900129

Bicho, Ana, and Christof Grewer. 2005. Rapid substrate-induced charge movements of the GABA transporter GAT1. *Biophys. J.* 89:211–231.

In the Fig. 8 legend, some of the kinetic parameters were not correct. We thank Dr. Donald D. F. Loo for pointing this problem out to us. The correct values are:

$k_{120} = 0.5 \text{ M}^{-1} \text{ ms}^{-1}$ ,  $k_{210} = 0.01 \text{ ms}^{-1}$ ,  $z_{12} = 0.9$ ,  $k_{230} = 10 \text{ M}^{-1} \text{ ms}^{-1}$ ,  $k_{320} = 0.1 \text{ ms}^{-1}$ ,  $z_{23} = 0.3$ ,  $k_{340} = 10,000 \text{ M}^{-1} \text{ ms}^{-1}$ ,  $k_{430} = 1 \text{ ms}^{-1}$ ,  $z_{34} = 0$ ,  $k_{450} = 1 \text{ ms}^{-1}$ ,  $k_{540} = 1 \text{ ms}^{-1}$ ,  $z_{45} = 0.2$ ,  $k_{560} = 0.3 \text{ ms}^{-1}$ ,  $k_{650} = 1.5 \cdot 10^7 \text{ M}^{-3} \text{ ms}^{-1}$ ,  $z_{56} = 0.2$ ,  $k_{160} = 0.005 \text{ ms}^{-1}$ ,  $k_{610} = 0.005 \text{ ms}^{-1}$ ,  $z_{61} = -0.8$ ,  $k_{670} = 200 \text{ M}^{-1} \text{ ms}^{-1}$ ,  $k_{760} = 20 \text{ ms}^{-1}$ ,  $z_{67} = 0.01$ ,  $k_{780} = 0.4 \text{ ms}^{-1}$ ,  $k_{870} = 0.02 \text{ ms}^{-1}$ ,  $z_{78} = 0.005$ ,  $k_{810} = 50 \text{ ms}^{-1}$ ,  $k_{180} = 10,000 \text{ M}^{-1} \text{ ms}^{-1}$ , and  $z_{81} = -0.215$ .

These revised kinetic parameters lead to simulations indistinguishable from those shown in Figs. 9 and 10. However, in Fig. 11, the y axis scale has to be multiplied by a factor of  $\sim 10$ . Finally, it should be noted that the background currents in the simulations before GABA application were subtracted to allow for a better comparison of the simulated currents with the experimental traces.

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doi: 10.1529/biophysj.105.900130

Jayachithra, Kandaswamy, Thallampuranam Krishnaswamy Suresh Kumar, Ta-Jung Lu, Chin Yu, and Der-Hang Chin. 2005. Cold instability of aponeocarcinostatin and its stabilization by labile chromophore. *Biophys. J.* 88:4252–4261.

On page 4254, the third sentence of the section “pH stability of apoNCS” in Results should read:

Without urea, the  $T_m$  of the protein is almost unchanged in the pH range of 4.0–10.0.

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Alves, I. D., G. F. J. Salgado, Z. Salamon, M. F. Brown, G. Tollin, and V. J. Hruby. 2005. Phosphatidylethanolamine enhances rhodopsin photoactivation and transducin binding in a solid supported lipid bilayer as determined using plasmon-waveguide resonance spectroscopy. *Biophys. J.* 88:198–210.

On page 200, line 18 of section “Formation of solid-supported lipid bilayers”, butanol/methanol (0.05:0.95:0.5, v/v) should read: butanol/methanol (0.05:9.5:0.5, v/v).

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