CORRECTIONS

Tatsuya Yamaue,* Hiroto Mukai, Kinji Asaka, and Masao Doi: Electrostress Diffusion Coupling Model for Polyelectrolyte Gels. Volume 38, Number 4, February 22, 2005, pp 1349−1356.

The following equations should appear as follows:

$$\mathbf{j}_s = -\kappa\nabla p - \lambda\nabla\psi \tag{2}$$

$$\nabla\cdot\mathbf{j}_e = 0 \tag{9}$$

$$\nabla\cdot(c_p w_p \mathbf{v}_p + c_s w_s \mathbf{v}_s + \sum_i c_i w_i \mathbf{v}_i) = 0 \tag{10}$$

$$\nabla\cdot(\sigma - p\mathbf{I}) = 0 \tag{11}$$

$$\lambda_p = -\frac{q_p}{\zeta_p}(1-\phi_p) \tag{19}$$

$$\lambda_i = \sum_i \frac{c_i q_i w_i}{\zeta_i} \tag{20}$$

$$(\sigma - p\mathbf{I})\cdot\mathbf{n} = \mathbf{f}_{\mathrm{bc}} \quad \text{for a deformable boundary} \tag{21}$$

$$(\sigma - (k_{\mathrm{B}}T\sum_i c_i + p)\mathbf{I})\cdot\mathbf{n} = \mathbf{f}_{\mathrm{bc}} \tag{23}$$

$$(\sigma - (k_{\mathrm{B}}T\sum_i [[c_i]] + [[p]])\mathbf{I})\cdot\mathbf{n} = 0 \tag{24}$$

$$\mathbf{j}_s\cdot\mathbf{n} = 0 \quad \text{for an impermeable wall} \tag{26}$$

$$\mathbf{j}_s\cdot\mathbf{n} = \kappa_b[[p]] \tag{27}$$

$$\mathbf{j}_e\cdot\mathbf{n} = \dot{\Phi}/C \tag{29}$$

$$\zeta_p = \frac{6\pi\eta}{\xi_b{}^2 c_p} \tag{31}$$

$$u_y = \frac{xy}{R(t)}. \tag{39}$$

$$u_z = \frac{xz}{R(t)} \tag{40}$$

$$f(x,t) = \frac{\partial u_x}{\partial x} + \frac{2x}{R(t)} \tag{41}$$

$$p(x,t) = \sigma_{xx}(x,t) \tag{45}$$

$$D' = \left(\kappa - \frac{\lambda^2}{\sigma_e}\right)\left(K + \frac{4}{3}G\right) \tag{52}$$

On p 1352, right side, lines 12−13, the sentence should read: For typical swollen polyelectrolyte gel of fluorocarbon, we can take $\phi_p \sim 0.5$, $\xi_b \sim 1.0$ [nm], and $c_i \sim 1.0/z$ [mol/L].

On p 1352, right side, lines 42−43, the sentence should read: For ϕ_p and c_i, we use the value $\phi_p \sim 0.7$ and $c_i \sim 1.0/z$ [mol/L], which are reported in ref 15.

$f(x,t) \equiv \nabla\mathbf{u}(\mathbf{r},t)$ is modifed to $f(x,t) \equiv \nabla\cdot\mathbf{u}(\mathbf{r},t)$.

In the caption of Figure 7, $\tau_{\mathrm{relax}}/\tau_{\mathrm{relax}}$, H^+ is modified to $\tau_{\mathrm{relax}}/\tau_{\mathrm{relax},H^+}$.

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