Correction to "Modeling the Interfacial Tension in Oil−Water-Nonionic Surfactant Mixtures Using Dissipative Particle Dynamics and Self-Consistent Field Theory"

Valeriy V. Ginzburg,* Kwanho Chang, Prasanna K. Jog, Andre B. Argenton, and Leela Rakesh

The value of χ(C−W) in Table 1 is incorrect and should read 2.47, rather than 0.47. Overall, Table 1 should read as follows.

Table 1. Flory−Huggins Parameters Used in SCFT Simulations

χ(C−E)	0.25
χ(E−E)	0.09
χ(C−W)	2.47