Corrections

Prediction of Ionic Liquid Properties. I. Volumetric Properties as a Function of Temperature at 0.1 MPa. Johan Jacquemin,* Rile Ge, Paul Nancarrow, David W. Rooney, Margarida F. Costa Gomes, Agilio A. H. Pádua, and Christopher Hardacre,* *J. Chem. Eng. Data* **2008,** *53,* 716–726.

Page 722. Equations 6 and 7 are wrong, and the following equations must be used in order to obtain a comparable relationship with eq 4. These errors are typographical and do not influence the calculations and conclusions claimed in this article.

$$V^*_{[C_4mim]^+}(\delta T) = V^*_{[C_4mim]^+}(T^*) + \Phi_{[C_4mim]^+} \cdot \left(\sum_{i=1}^{2} A_i \cdot (\delta T)^i \right) \quad (6)$$

$$V^*_{[NTf_2]^-}(\delta T) = V^*_{[NTf_2]^-}(T^*) + \Phi_{[NTf_2]^-} \cdot \left(\sum_{i=1}^{2} A_i \cdot (\delta T)^i \right) \quad (7)$$

where A_i are the coefficients obtained by eq 4 and $\Phi_{[C_4mim]^+}$ and $\Phi_{[NTf_2]^-}$ are the effective molar volume fractions at 298.15 K of ions constituting $[C_4mim][NTf_2]$.

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