



Correction to: Molecular Interactions and Mechanisms in the ^1H NMR Relaxation of Residual CHCl_3 in Deuteriochloroform Solution of a Two-Chain Ionic Surfactant

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A few typographical errors have been noticed in the published article.

On page 1262, Eq. 7, the Planck's constant term, $\hbar = h/2\pi$, should have 2 as an exponent, and in the denominator the hydrogen atom–chlorine atom internuclear separation distance, r_{HCl} , should have 6 as an exponent, and not 6 as a coefficient. The correct equation is:

$$\frac{1}{T_1} = \frac{15 \gamma_{\text{H}}^2 \hbar^2 (0.75 \gamma_{^{35}\text{Cl}}^2 + 0.25 \gamma_{^{37}\text{Cl}}^2) \tau_c}{r_{\text{HCl}}^6} \quad (7)$$

On page 1262, Eq. 8, the Planck's constant term, $\hbar = h/2\pi$, should have 2 as an exponent. The correct equation is:

$$\frac{1}{T_1} = \frac{1}{T_2} = \frac{2 \gamma_{\text{H}}^2 \gamma_{\text{H}}^2 \hbar^2 I(I+1) \tau_c}{r_{\text{HH}}^6} \quad (8)$$

The correct equations were used in the analysis and, therefore, all calculations and results in the published article remain unchanged.

The original article can be found online at <https://doi.org/10.1007/s10953-018-0789-x>.

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