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Addition/Correction

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Correction to "Conformational Distribution of *trans*-Stilbene in Solution Investigated by Liquid Crystal NMR Spectroscopy and Compared with In Vacuo Theoretical Predictions"

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We realized that, due to a mistake in the printed output of our computer program, in Table 2 of the published article the values corresponding to the interaction tensor components $\varepsilon^{R}_{2,0}$, $\varepsilon^{R}_{2,2}$, and $\varepsilon^{(H_6—C=C—H_7)}_{2,0}$ are listed in the wrong order and are given in the wrong units (kJ mol^{-1} units instead of RT units, the right ones). The right complete Table 2 is given here.

Table 2. Optimized Values of the Iteration Parameters Required by the AP-DPD Approach

	t-St in ZLI1132
$\phi^{M}_{1} = \pm\phi^{M}_{2}$/deg	16.80 ± 0.02
$h_1 = h_2$/deg	10 (after parametrization)
$w_{(C2)}$	0.59 ± 0.01
$\varepsilon^{R}_{2,0}/RT$	1.349 ± 0.001
$\varepsilon^{R}_{2,2}/RT$	0.650 ± 0.002
$\varepsilon^{(H_6—C=C—H_7)}_{2,0}/RT$	0.473 ± 0.002

We want anyway emphasize that the mistake was only in the output, so that the right values have been fortunately used during calculations and the results of the work are not affected at all.

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