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Correction to "Lipid spontaneous curvatures estimated from temperature-dependent changes in inverse hexagonal phase lattice parameters: the effects of metal cations"

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Correction:

It has come to our attention that there are two typos in the original manuscript concerning equations 7 and 8. These should read as;

We define the 'curvature power 'of lipid j (χ_j) as:

$$\chi_j = -(L_{p,\mathrm{PE}(37)} - L_{p,\mathrm{j}(37)}) \, / x_j \qquad (7).$$

where $Lp_{,\mathrm{j}(37)}$ is the lattice parameter, at 37 °C, of lipid j, $Lp_{,\mathrm{PE}(37)}$ is the lattice parameter of DOPE and x_j is the mole fraction of the lipid j.

The universal curve can be fitted with the linear equation:

$$c_0 = 0.0731 \, \chi - 0.3694 \, . \qquad (8)$$