

**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  $j$  ( $\chi_j$ ) as:

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

where  $L_{p,j(37)}$  is the lattice parameter, at 37 °C, of lipid  $j$ ,  $L_{p,PE(37)}$  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:

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