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CORRECTION

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Correction: Assembly and relaxation behaviours of phosphatidylethanolamine monolayers investigated by polarization and frequency resolved SFG-VS

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Correction for 'Assembly and relaxation behaviours of phosphatidylethanolamine monolayers investigated by polarization and frequency resolved SFG-VS' by Feng Wei *et al.*, *Phys. Chem. Chem. Phys.*, 2015, **17**, 25114–25122.

The authors would like to make some corrections to their article based on some recent calculations:

The spectral resolution of IR-scanning sum frequency generation vibrational spectroscopy (SFG-VS) system can be calculated using eqn (13) in ref. 1:

$$\Delta\nu_{\rm Instr}^{2} = \frac{8\ln(2)}{(2\pi c)^{2}} \Delta\omega_{\rm SF}^{2} = \frac{8\ln(2)}{(2\pi c)^{2}} (\Delta\omega_{\rm IR}^{2} + \Delta\omega_{\rm Vis}^{2})$$
(13)

where $\Delta\omega_{\rm IR}$ and $\Delta\omega_{\rm Vis}$ are the Guassian bandwidths/linewidths of incident IR beam and visible beam, respectively. So based on our latest experiments and calculation results:

- 1. "Bandwidth $\approx 1 \text{ cm}^{-1}$ " in the abstract (line 3) should be corrected into "Linewidth $\approx 1.5 \text{ cm}^{-1}$ ".
- 2. "The bandwidth of the final IR beam is calculated to be $\approx 1~\text{cm}^{-1}$, about twice the transform-limited width of a 30 ps Gaussian pulse (0.48 cm⁻¹)." in "Experimental" (page 2, left column, last third line) should be corrected into "The linewidths of the final IR beams are calculated to be $\approx 1.5~\text{cm}^{-1}$."
- 3. "The spectral resolution of the current SFG system is $\approx 1 \text{ cm}^{-1}$." in "Introduction" (page 2, left column, paragraph 2, line 11–12) should be corrected into "The calculated spectral resolution $\Delta \nu_{\text{Instr}}$ of the current SFG system is about 4.6 cm⁻¹."

Reference

1 F. Wei, W. X. Xia, Z. J. Hu, W. H. Li, J. Y. Zhang and W. Q. Zheng, *Chin. J. Chem. Phys.*, 2016, **29**, 171–178. The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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