Erratum: First-principles study of the elasticity, piezoelectricity, and vibrational modes in $LiGaO_2$ compared with ZnO and GaN [Phys. Rev. B 81, 235214 (2010)]

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In Fig. 3 the y-axis label should be Im$[\varepsilon(\omega)]$, Im$[-\varepsilon^{-1}(\omega)]$.

Figure 5 (which was identical to Fig. 6 in the original) should be replaced by the Fig. 5 presented here.

The caption in Fig. 6 should say $b_2(y)$ instead of $b_2(x)$. The text in the last paragraph of Sec. III E 2 should state: For b_2 symmetry, the visual agreement between the spectra is excellent. Again, the spectrum can roughly be divided into three regions, a low-frequency mode near 300 cm^{-1} followed by a three peak region between 400 and 550 cm^{-1}, and a high-frequency band between 600 and 800 with a sharp dip in the middle. We note that the features in the middle range (350–550 cm^{-1}) are somewhat shifted up relative to the experiment while for the high-frequency range, the theory seems to underestimate slightly the experimental frequencies. Note, for example, the position of the sharp dip in reflectivity. The spectrum for b_1 symmetry, on the other hand, shows a series of weak features below 450 cm^{-1} and two main peaks between 450–600 cm^{-1} and 600–800 cm^{-1}.

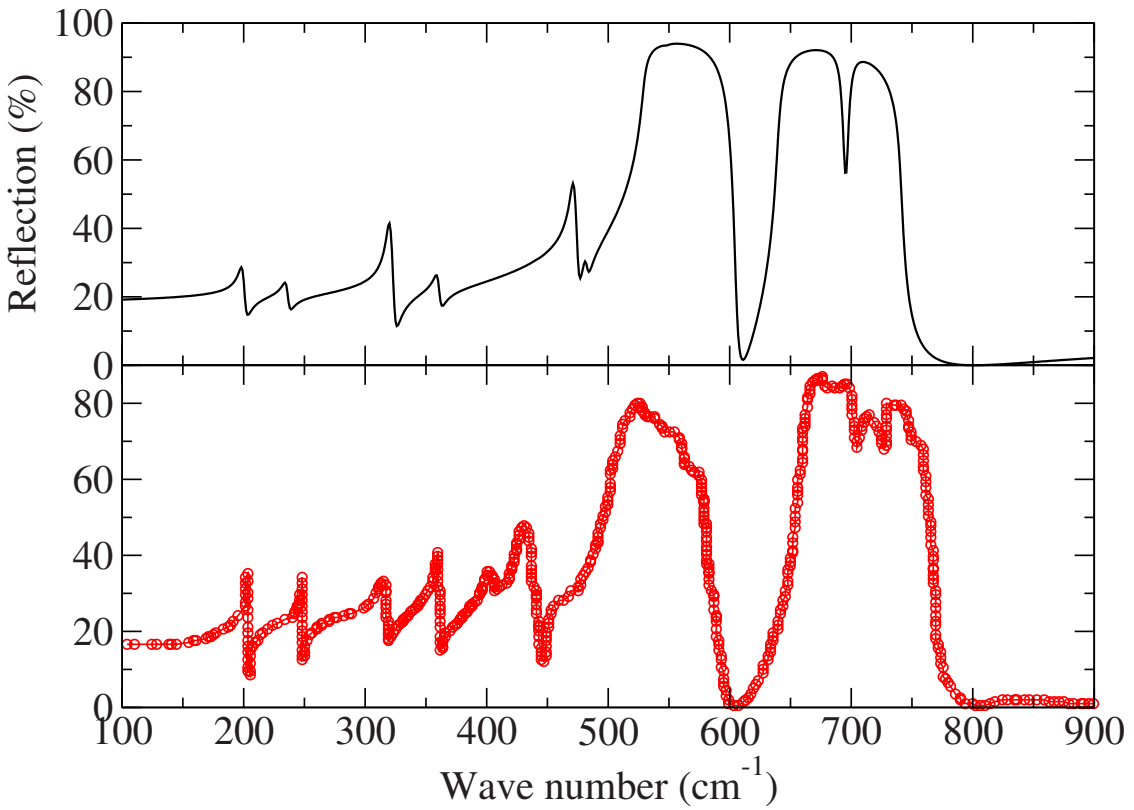

FIG. 5. (Color online) Comparison of experimental (lower panel) infrared reflectivity of $b_1(x)$ with calculated (upper panel) reflectivity.

