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## Correction to "Energy dissipation in multifrequency atomic force microscopy"

Valentina Pukhova<sup>1,2</sup>, Francesco Banfi<sup>2</sup> and Gabriele Ferrini<sup>\*2</sup>

Correction

Address:

<sup>1</sup>Dipartimento di Fisica, Università degli Studi di Milano, I-20122 Milano, Italy and <sup>2</sup>Interdisciplinary Laboratories for Advanced Materials Physics (i-LAMP) and Dipartimento di Matematica e Fisica, Università Cattolica, I-25121 Brescia, Italy

Email:

Gabriele Ferrini\* - gabriele.ferrini@unicatt.it

\* Corresponding author

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In the section "Energy dissipation" of the above manuscript, there is a typesetting error in the mathematical expressions after Equation 5. The correct form must be:

The energy balance of each decaying mode obtained from Equation 4 in the time window  $0 < t < \tau = 200 \mu s$  (see Figure 1) can be written as

$$E_{b_i} = E_{\gamma_i} \tag{5}$$

where

$$E_{b_i} = \Delta K_i + \Delta U_i$$
  
$$E_{\gamma_i} = \int_0^{\tau} m_{eq} \gamma_i v_i^2 dt$$

*i* is the index of the mode,  $\Delta K_i = 1/2 \ m_{\rm eq}(v_i(0)^2 - v_i(\tau)^2)$  is the variation of kinetic energy, and  $\Delta U_i = 1/2 \ k_i(z_i(0)^2 - z_i(\tau)^2)$  is the variation of elastic potential energy.

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