



Cite this: *Phys. Chem. Chem. Phys.*,  
2020, 22, 1784

DOI: 10.1039/c9cp90308a

[rsc.li/pccp](http://rsc.li/pccp)

## Correction: p-Type conductivity mechanism and defect structure of nitrogen-doped LiNbO<sub>3</sub> from first-principles calculations

Weiwei Wang,<sup>a</sup> Yang Zhong,<sup>a</sup> Dahuai Zheng,<sup>b</sup> Hongde Liu,<sup>\*a</sup> Yongfa Kong,<sup>\*ab</sup>  
Lixin Zhang,<sup>a</sup> Romano Rupp<sup>cd</sup> and Jingjun Xu<sup>\*ab</sup>

Correction for 'p-Type conductivity mechanism and defect structure of nitrogen-doped LiNbO<sub>3</sub> from first-principles calculations' by Weiwei Wang et al., *Phys. Chem. Chem. Phys.*, 2020, 22, 20–27.

The authors would like to make a correction to the author list of the published article where the name of co-author Romano Rupp was incorrectly displayed as Rupp Romano. The amended author list is as shown herein.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

<sup>a</sup> MOE Key Laboratory of Weak-Light Nonlinear Photonics and School of Physics, Nankai University, Tianjin 300071, China. E-mail: liuhd97@nankai.edu.cn, kongyf@nankai.edu.cn, jjxu@nankai.edu.cn

<sup>b</sup> TEDA Institute of Applied Physics, Nankai University, Tianjin 300457, China

<sup>c</sup> Faculty of Physics, Vienna University, Wien, A-1090, Austria

<sup>d</sup> Department of Complex Matter, Jozef Stefan Institute, Ljubljana, Slovenia

