

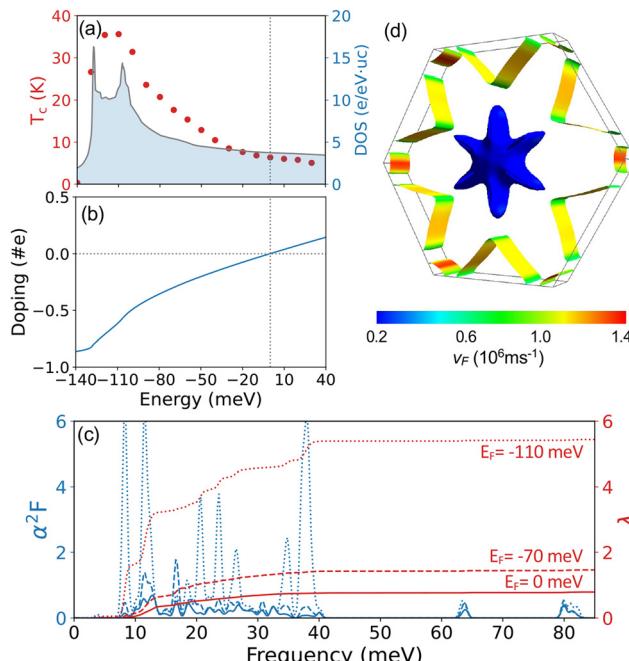
## CORRECTION

[View Article Online](#)  
[View Journal](#) | [View Issue](#)
Cite this: *Nanoscale*, 2023, **15**, 11727DOI: 10.1039/d3nr90127k  
rsc.li/nanoscale

## Correction: Superconductivity in functionalized niobium-carbide MXenes

Cem Sevik,<sup>a,b</sup> Jonas Bekaert<sup>a</sup> and Milorad V. Milošević<sup>a,c</sup>Correction for 'Superconductivity in functionalized niobium-carbide MXenes' by Cem Sevik *et al.*, *Nanoscale*, 2023, **15**, 8792–8799, <https://doi.org/10.1039/D3NR00347G>.

In Fig. 4d and 5d of the published article, the colour bar legend (representing the Fermi velocities on the Fermi surface) should be rotated 180 degrees from the original depiction. The correct colour bar orientation is shown below.



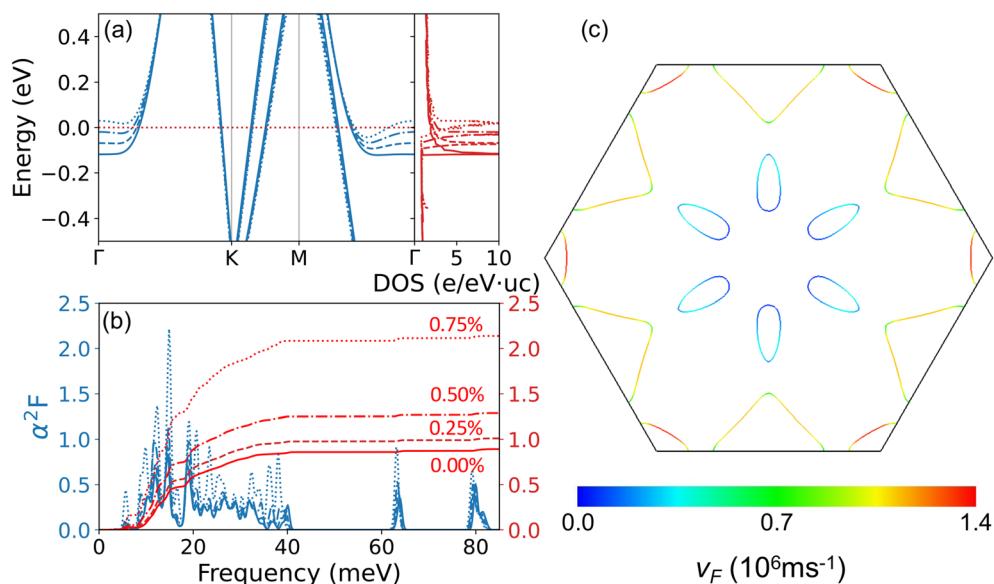
**Fig. 4** The effect of gating on the superconducting properties of bulk-layered  $\text{Nb}_2\text{CCl}_2$ . (a)  $T_c$  along with the electronic density of states, as a function of the electronic energy level. (b) Carrier doping levels with the Fermi level shift – negative values correspond to hole doping. (c) Eliashberg function,  $\alpha^2F$ , and the electron–phonon coupling constant,  $\lambda$ , for different values of the gating-shifted Fermi level. (d) Fermi surface, along with Fermi velocities, for the Fermi level shifted by gating to  $\sim 80$  meV below the intrinsic value.

<sup>a</sup>Department of Physics & NANOLab Center of Excellence, University of Antwerp, Groenenborgerlaan 171, B-2020 Antwerp, Belgium. E-mail: cem.sevik@uantwerpen.be, jonas.bekaert@uantwerpen.be, milorad.milosevic@uantwerpen.be

<sup>b</sup>Department of Mechanical Engineering, Faculty of Engineering, Eskisehir Technical University, 26555 Eskisehir, Turkey

<sup>c</sup>Instituto de Física, Universidade Federal de Mato Grosso, Cuiabá, Mato Grosso 78060-900, Brazil





**Fig. 5** The calculated (a) electronic band structure, and (b) the Eliashberg function,  $\alpha^2F$ , and the electron–phonon coupling constant,  $\lambda$ , for two-dimensional  $\text{Nb}_2\text{CCl}_2$  under applied tensile strain. The Fermi surface corresponding to the case with 0.75% applied tensile strain is presented in panel (c).

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

