Biophysical *Journal*
Correction

Biophysical Society

Dynamics of Membrane-Bound G12V-KRAS from Simulations and Single-Molecule FRET in Native Nanodiscs

Priyanka Prakash, Douglas Litwin, Hong Liang, Suparna Sarkar-Banerjee, Drew Dolino, Yong Zhou, John F. Hancock, Vasanthi Jayaraman, and Alemayehu A. Gorfe*

(Biophysical Journal *116*, 179–183; January 22, 2019)
In Materials and Methods, the last sentence under the sub-heading MD Simulation should read:

The trajectory was analyzed primarily based on the distance between Cα atoms of residues 132 and 183 (ζ) and the angle between a vector from residue 5 to residue 9 Cα atoms and the membrane normal (**Θ**) (see Supporting Materials and Methods for details).

*Correspondence: alemayehu.g.abebe@uth.tmc.edu
https://doi.org/10.1016/j.bpj.2019.12.032