

CORRECTION

Correction: Molecular dynamic simulations to investigate the structural impact of known drug resistance mutations on HIV-1C Integrase-Dolutegravir binding

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Reference

1. Chitongo R, Obasa AE, Mikasi SG, Jacobs GB, Cloete R (2020) Molecular dynamic simulations to investigate the structural impact of known drug resistance mutations on HIV-1C Integrase-Dolutegravir binding. PLoS ONE 15(5): e0223464. <https://doi.org/10.1371/journal.pone.0223464> PMID: 32379830



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