

Corrigendum

Corrigendum to: Molecular dynamics simulations revealed Ca^{2+} -dependent conformational change of Calmodulin (FEBS 26172)

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An unfortunate mistake occurred in Fig. 3B. The correct figure is given below. Note that one of the lines is dotted.

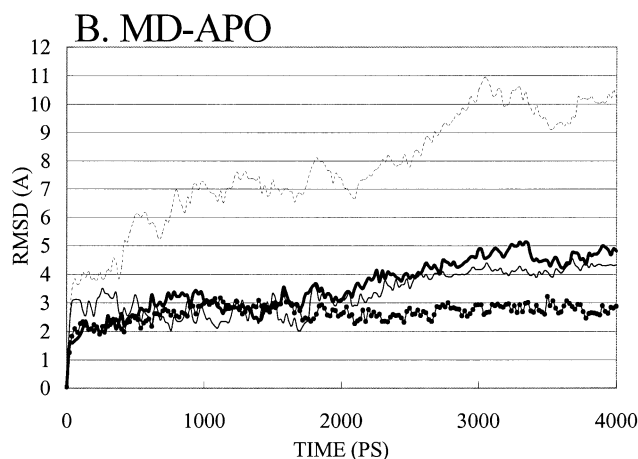


Fig. 3B.

The legends to Figs. 5 and 6 were misplaced. The correct legends are given below.

Fig. 5. Time evolution of the VDA in MD-holo. See Fig. 1B for a visual definition of the VDA.

Fig. 6. Time evolution of the bending angle of the central helix. The angle between helices D and E was plotted.

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