

# Accepted Manuscript

Molecular dynamics simulation approach to explore atomistic molecular mechanism of peroxidase activity of apoptotic cytochrome c mutants

Gurusamy Muneeswaran, Subramanian Kartheeswaran, Muniyandi Sankaralingam, Kaliappan Muthukumar, Manickam Pandiaraj, Chandran Karunakaran

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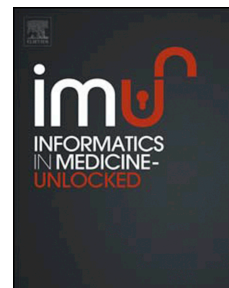
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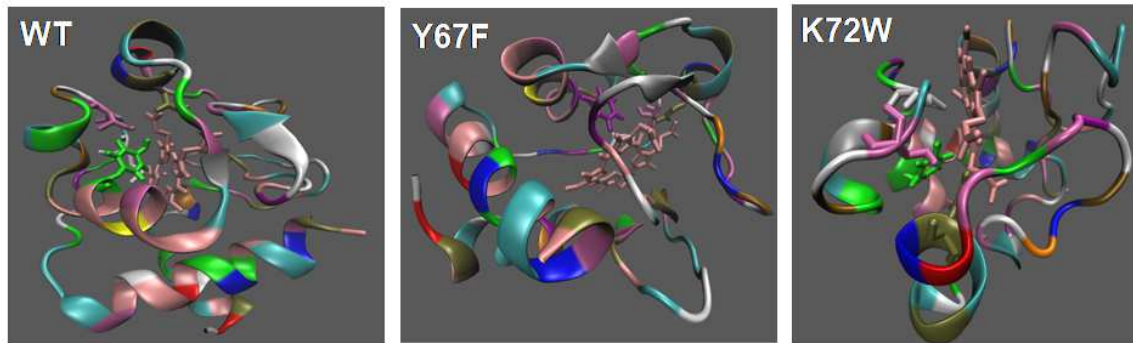
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## Graphical Abstract



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