

# Accepted Manuscript

The acidic pH-induced structural changes in apo-CP43 by spectral methodologies and molecular dynamics simulations

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PII: S0022-2860(17)31280-2

DOI: [10.1016/j.molstruc.2017.09.082](https://doi.org/10.1016/j.molstruc.2017.09.082)

Reference: MOLSTR 24331

To appear in: *Journal of Molecular Structure*

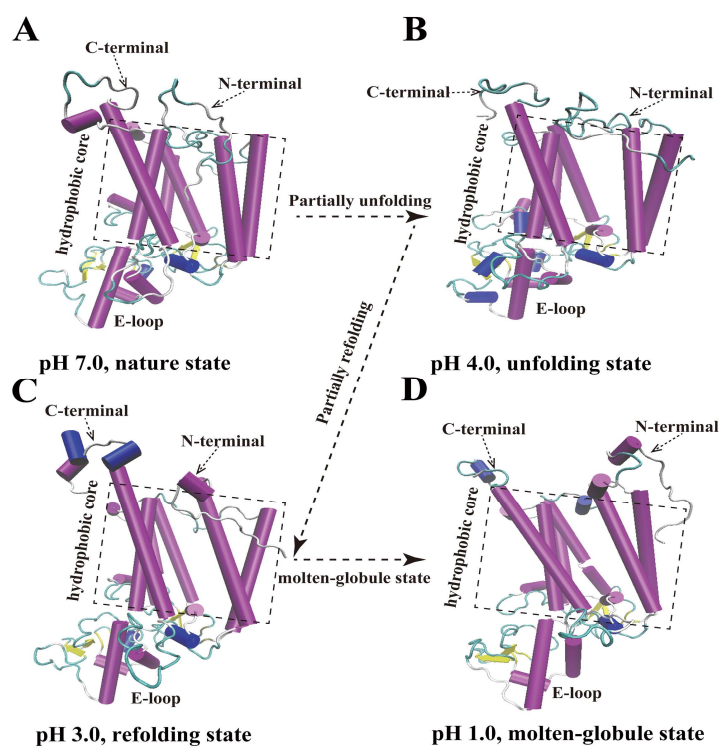
Received Date: 18 July 2017

Revised Date: 21 September 2017

Accepted Date: 21 September 2017

Please cite this article as: W. Wang, X. Li, Q. Wang, X. Zhu, Q. Zhang, L. Du, The acidic pH-induced structural changes in apo-CP43 by spectral methodologies and molecular dynamics simulations, *Journal of Molecular Structure* (2017), doi: 10.1016/j.molstruc.2017.09.082.

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