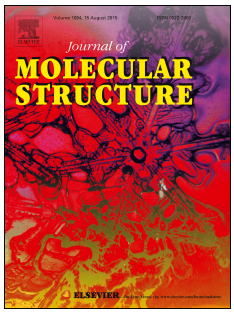
Accepted Manuscript



Quantum chemical investigation of levofloxacin-boron complexes: A computational approach

Koray Sayin, Duran Karakaş

PII: S0022-2860(18)30029-2

DOI: 10.1016/j.molstruc.2018.01.016

Reference: MOLSTR 24738

To appear in: *Journal of Molecular Structure*

Received Date: 27 October 2017

Revised Date: 9 December 2017

Accepted Date: 8 January 2018

Please cite this article as: K. Sayin, D. Karakaş, Quantum chemical investigation of levofloxacin-boron complexes: A computational approach, *Journal of Molecular Structure* (2018), doi: 10.1016/j.molstruc.2018.01.016.