Accepted Manuscript

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

Koray Sayin, Duran Karakas

PII: S0022-2860(18)30029-2

DOI: 10.1016/j.molstruc.2018.01.016

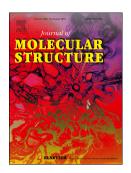
Reference: MOLSTR 24738

To appear in: Journal of Molecular Structure

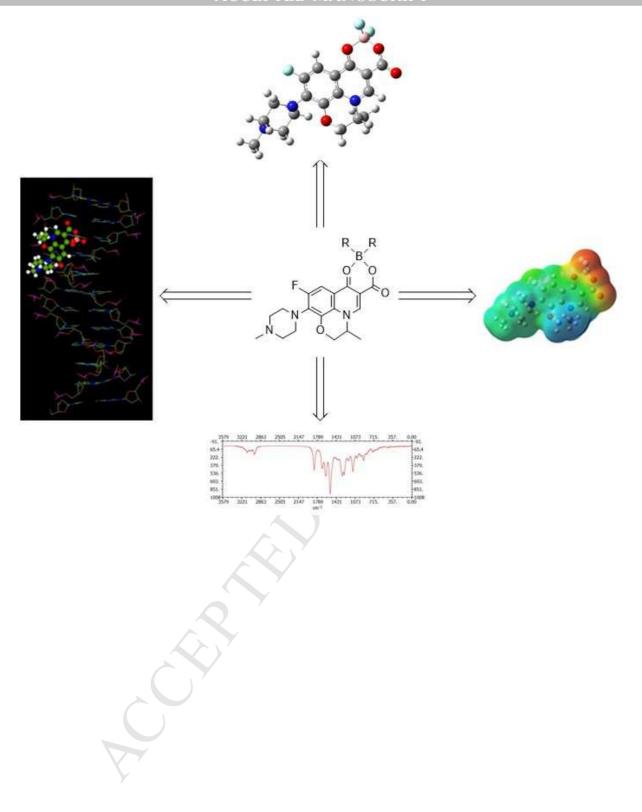
Received Date: 27 October 2017
Revised Date: 9 December 2017
Accepted Date: 8 January 2018

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

This is a PDF file of an unedited manuscript that has been accepted for publication. As a service to our customers we are providing this early version of the manuscript. The manuscript will undergo copyediting, typesetting, and review of the resulting proof before it is published in its final form. Please note that during the production process errors may be discovered which could affect the content, and all legal disclaimers that apply to the journal pertain.



ACCEPTED MANUSCRIPT



Download English Version:

https://daneshyari.com/en/article/7808035

Download Persian Version:

https://daneshyari.com/article/7808035

<u>Daneshyari.com</u>