THE JOURNAL OF PHYSICAL CHEMISTRY B

pubs.acs.org/JPCB

Addition/Correction

Correction to "The Amide I Spectrum of Proteins—Optimization of Transition Dipole Coupling Parameters Using Density Functional Theory Calculations"

Cesare M. Baronio and Andreas Barth*

J. Phys. Chem. B **2020**, *124* (9), 1703−1714. DOI: 10.1021/acs.jpcb.9b11793

Cite This: *J. Phys. Chem. B* 2020, 124, 2730−2730

Read Online

ACCESS | Metrics & More | Article Recommendations

The file names given in the SI text are no longer correct:

jp9b11793_si_002.xlsx is the table with the results of the optimization for all six structural models together.

jp9b11793_si_003.xlsx contains the results of the optimization for each secondary structure separately.

Published: March 24, 2020