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Additions and Corrections

Analyzing the $n \rightarrow \pi^*$ Electronic Transition of Formaldehyde in Water. A Sequential Monte Carlo/Time-Dependent Density Functional Theory

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In printed version of volume 19, number 2, page 306, column 1, the first sentence is missing, related to page 305:

results are not clear cut for this low intense $n \rightarrow \pi^*$ transition.

The on-line version is correct. The whole comprehensive sentence is:

However, because of the great ability of formaldehyde to form aggregates in water the results are not clear cut for this low intense $n \rightarrow \pi^*$ transition.

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