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Erratum

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Correction for Electronic Structure Investigation and Parametrization of Biologically Relevant Iron−Sulfur Clusters

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Supporting Information

Due to an unfortunate misunderstanding, the Supporting Information for our manuscript was incomplete. Here we have added the angle bending force constants and the atomic charges for the iron−sulfur clusters. We apologize for any inconvenience this may have caused.

■ ASSOCIATED CONTENT

Supporting Information

The Supporting Information is available free of charge on the ACS Publications website at DOI: 10.1021/acs.jcim.5b00390.

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