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**Radu Silaghi-Dumitrescu, Ioan Silaghi-Dumitrescu, Eric D. Coulter, and Donald M. Kurtz, Jr.\*:** Computational Study of the Non-Heme Iron Active Site in Superoxide Reductase and Its Reaction with Superoxide.

Pages 452 and 453. The paragraph headed “Computed Electronic Absorption Spectrum” and the caption below Figure 4 incorrectly refer to the ferric-peroxo model **4**, instead of the ferric-hydroperoxo model **6**. As stated in footnote 25, the results reported in this paragraph and shown in Figure 4 are for  $S = 1/2$  model **6**, not model **4**.

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