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Jonathan McMaster, Michael D. Carducci, Yi-Shan Yang, Edward I. Solomon,* and John H. Enemark*: Electronic Spectral Studies of Molybdenyl Complexes. 2. MCD Spectroscopy of $[MoOS_4]^-$ Centers.

Page 694. Figure 7 as published did not show the relative phases of the contour plot of the HOMO. The correct version of Figure 7 appears below.

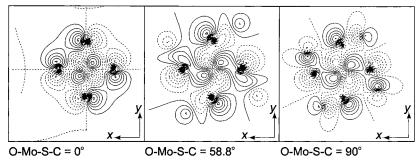


Figure 7. DFT calculation contour plot of the highest occupied molecular orbital (HOMO) of $[MoO(SMe)_4]^-$ at $O-Mo-S-C(\alpha)$ dihedral angles of 0°, 58.8°, and 90°.

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