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Geometric and electronic properties in a series of phosphorescent heteroleptic Cu(I) complexes:

Crystallographic and computational studies

Katharina Kubi ek^{a,b,*}, Sreevidya Thekku Veedu^{a,b}, Darina Storozhuk^{a,b,c}, Reza Kia^{a,b,l,*}, and Simone Techert^{a,b,c,*}

^aFS-SCS, Deutsches Elektronen-Synchrotron (DESY), Notkestraße 85, 22607 Hamburg, Germany

^bMax Planck Institute for Biophysical Chemistry, Am Faßberg 11, 37077 Göttingen, Germany

^cInstitute of X-ray Physics, Georg-August-Universität Göttingen, Friedrich-Hund-Platz 1, 37077 Göttingen, Germany

*Corresponding authors. *E-mail addresses*: katharina.kubicek@desy.de (K.K.);

rkia@sharif.edu (R.K.); simone.techert@desy.de (S.T.).

¹Present address: Chemistry Department, Sharif University of Technology, P.O. Box 11155-3516, Tehran, Iran Download English Version:

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