

## Correction to Gadolinium Acetylacetonate Tetraphenyl Monoporphyrinate Complex and Some of Its Derivatives: EXAFS Study and Molecular Dynamics Simulation

J. H. Agondanou, I. Nicolis, E. Curis, J. Purans, G. A. Spyroulias, A. G. Coutsolelos,\* and S. Bénazeth\*

*Inorg. Chem.* **2007**, *46* (17), 6871–6879. DOI: 10.1021/ic061861x

### **S** Supporting Information

Page 6879: Because of a production error, the original Supporting Information file was not published. The file is published with this Addition and Correction and contains the information presented below.

### **■ ASSOCIATED CONTENT**

#### **S** Supporting Information

Structural data obtained from the best-fit analysis on the raw data of the monomers Sm(tpp)(acac) and Gd(tpyp)(acac) (Table S1) and the dimer GdH(tpyp)<sub>2</sub> (Table S2) and figure showing the EXAFS of the single- and multiple-scattering paths that produce the largest contribution to the total EXAFS signal (Figure S1). This material is available free of charge via the Internet at <http://pubs.acs.org>.

Published: October 2, 2013