## Synthesis, Characterization and Electrochemistry of the Rhodium(III) Complexes of *meso*-Substituted [14]Tribenzotriphyrin(2.1.1)

Zhaoli Xue,<sup>a\*</sup> Yemei Wang,<sup>a</sup> John Mack,<sup>b\*</sup> Scebi Mkhize,<sup>b</sup> Tebello Nyokong,<sup>b</sup> Yuanyuan Fang,<sup>a</sup> Zhongping Ou<sup>a</sup> and Karl M. Kadish <sup>c\*</sup>

<sup>a</sup>School of Chemistry and Chemical Engineering, Jiangsu University, Zhenjiang 212013, China <sup>b</sup>Department of Chemistry, Rhodes University Grahamstown 6140, South Africa <sup>c</sup>Department of Chemistry, University of Houston, Houston, Texas 77204-5003, United States



Table S1. Dihedral angles between *meso*-phenyl rings and the triphyrin mean plane (deg)<sup>a</sup>.

	2b		2b
$\Phi_1{}^a$	88.21	C9-C29	1.502(8)
$\Phi_2$	78.53	C18-C35	1.498(9)
$\Phi_3$	86.71	C19-C41	1.500(8)
$\Phi_4$	85.69	C28-C47	1.501(9)

<sup>*a*</sup> Angles  $\Phi_i$  (i = 1, 2, 3, 4) are defined by the dihedral angles between the *meso*-phenyl rings and the triphyrin mean plane.



**Figure S1.** Packing diagram of **2b** and the distance between the two triphyrin mean plan (blue line) and the two Rh(III) ions (red line). The solvent molecules are omitted for clarity.



**Figure S2.** <sup>1</sup>H NMR spectra of a) **1a** and b) **2a** in CDCl<sub>3</sub>. An asterisk is used to denote the presence of a solvent peak.





Figure S3. <sup>1</sup>H NMR spectra of 2a in CDCl<sub>3</sub> at room temperature.



Figure S4. <sup>1</sup>H NMR spectra of 2b in CDCl<sub>3</sub> at room temperature.



Figure S5. <sup>1</sup>H NMR spectra of 2c in CDCl<sub>3</sub> at room temperature.

$$-8.82$$
  
 $-8.72$   
 $-8.13$   
 $7.146$   
 $-7.41$   
 $-7.416$   
 $-6.66$   
 $-6.66$ 



Figure S6. <sup>1</sup>H NMR spectra of 2d in CDCl<sub>3</sub> at room temperature.