Supplementary Material (ESI) for Organic & Biomolecular Chemistry

Porphyrin-naphthodiimide interactions as a structural motif in foldamers and supramolecular assemblies

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Electronic Supplementary Information

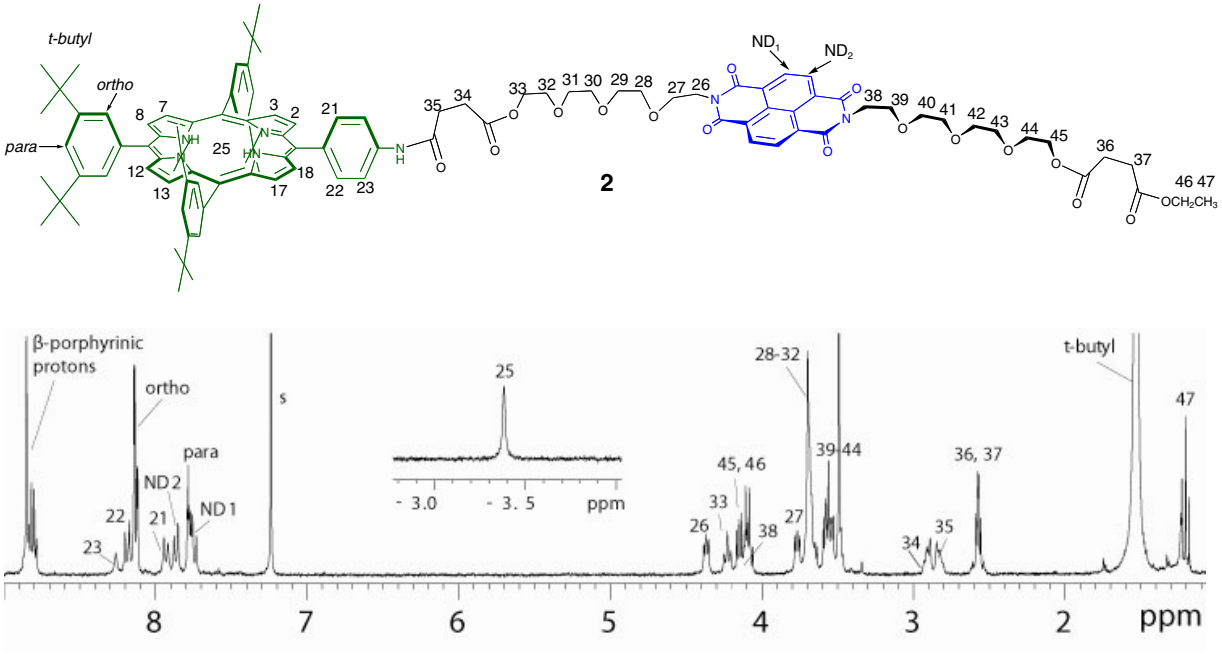


Figure S1 ^{1}H NMR spectrum and numbering scheme for **2**.

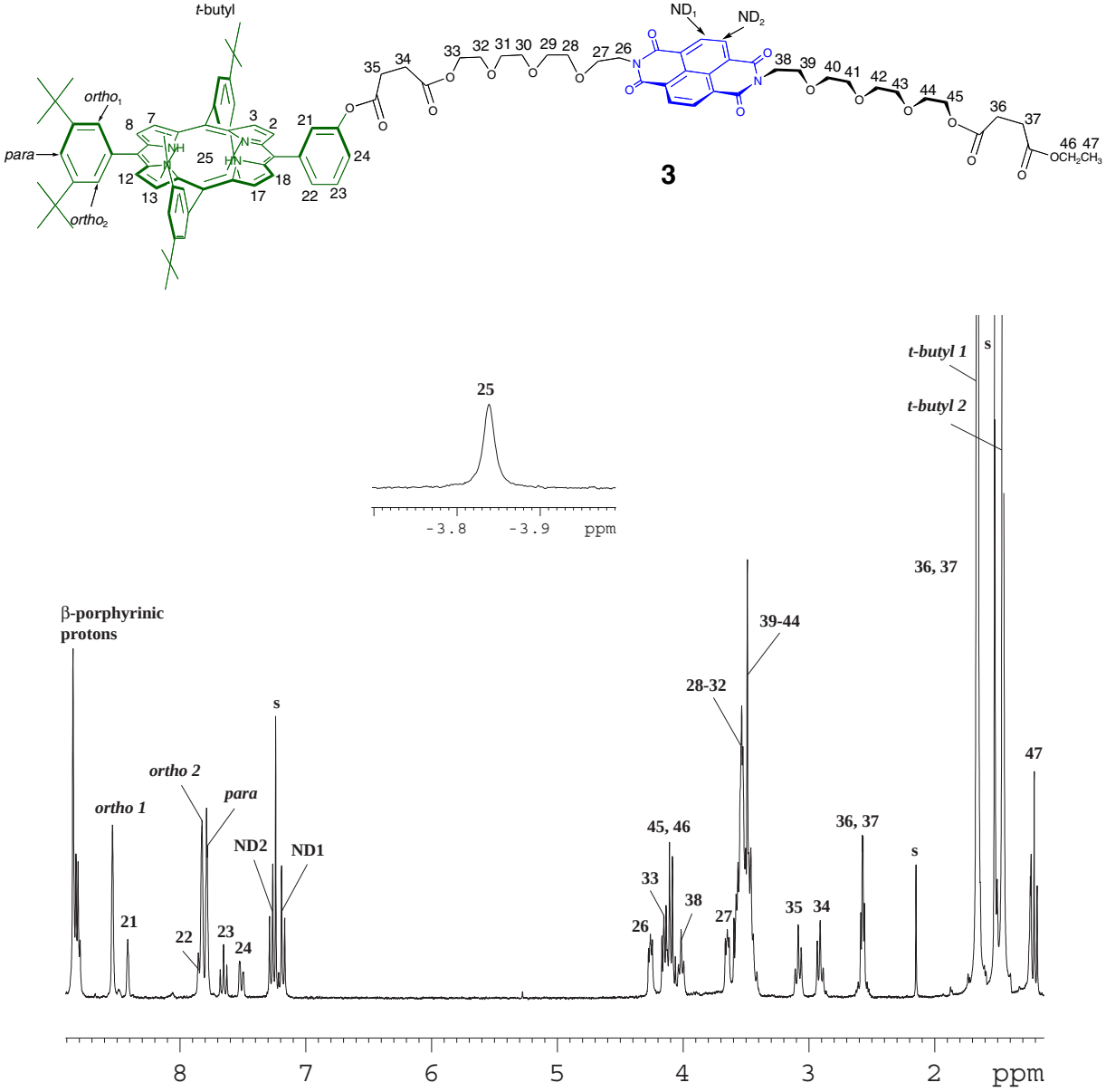


Figure S2 ^{1}H NMR spectrum and numbering scheme for **3**.

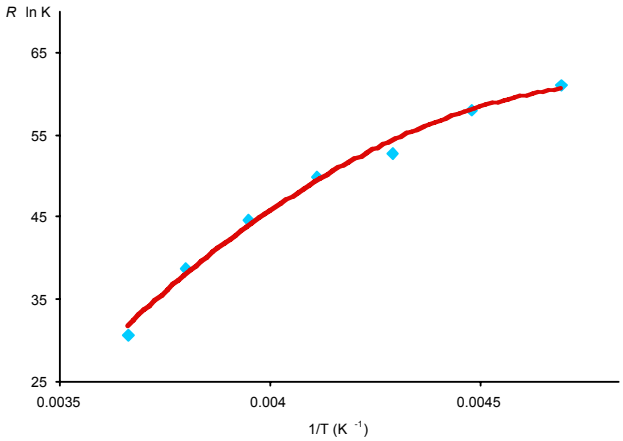


Figure S3 van't Hoff plot for the [**DN38C10.3**] system resulting from an equimolar admixture of DN38C10 and **3** in $CDCl_3$. The points are the measured data and the solid line is a guide to eye.

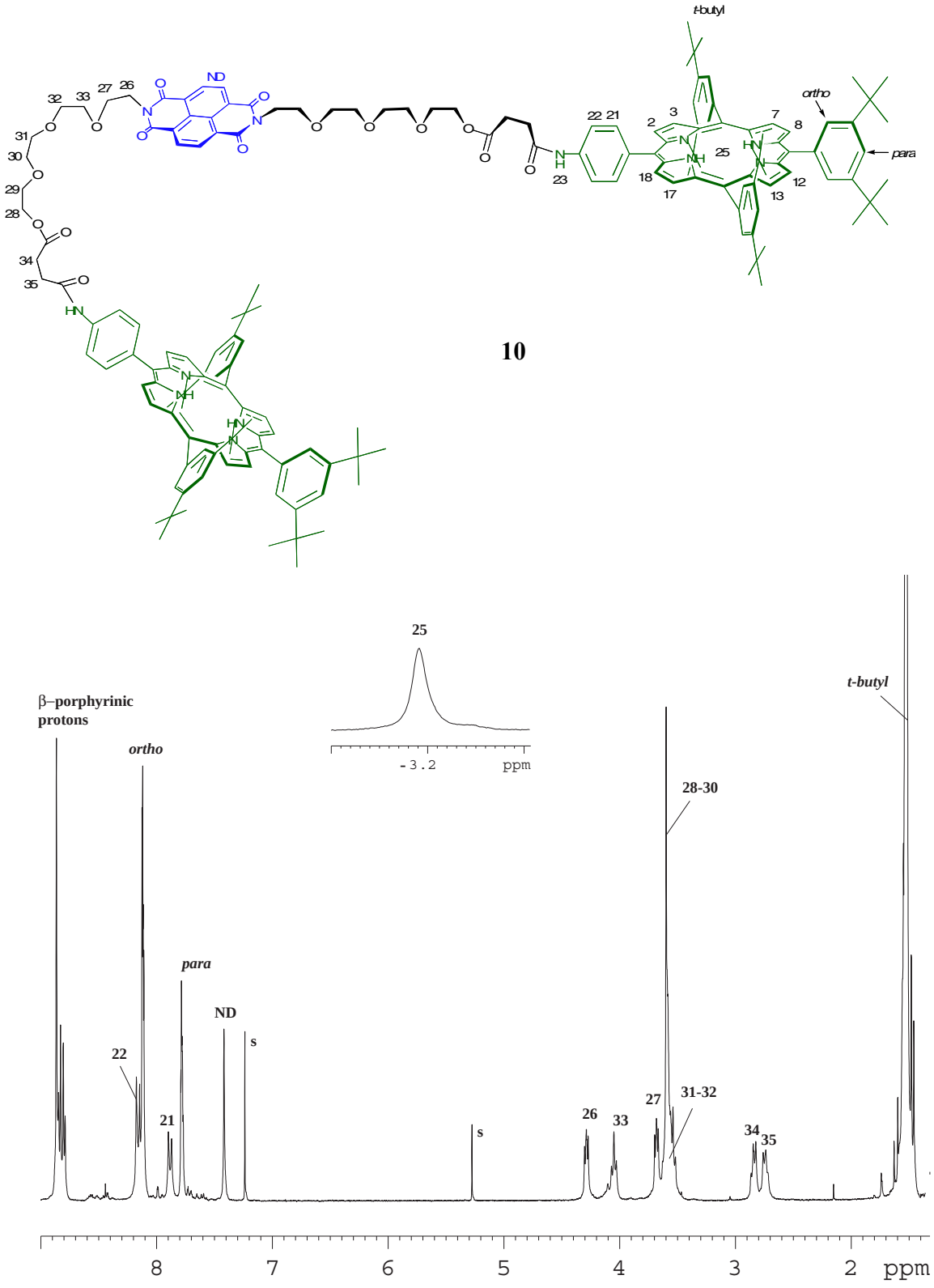
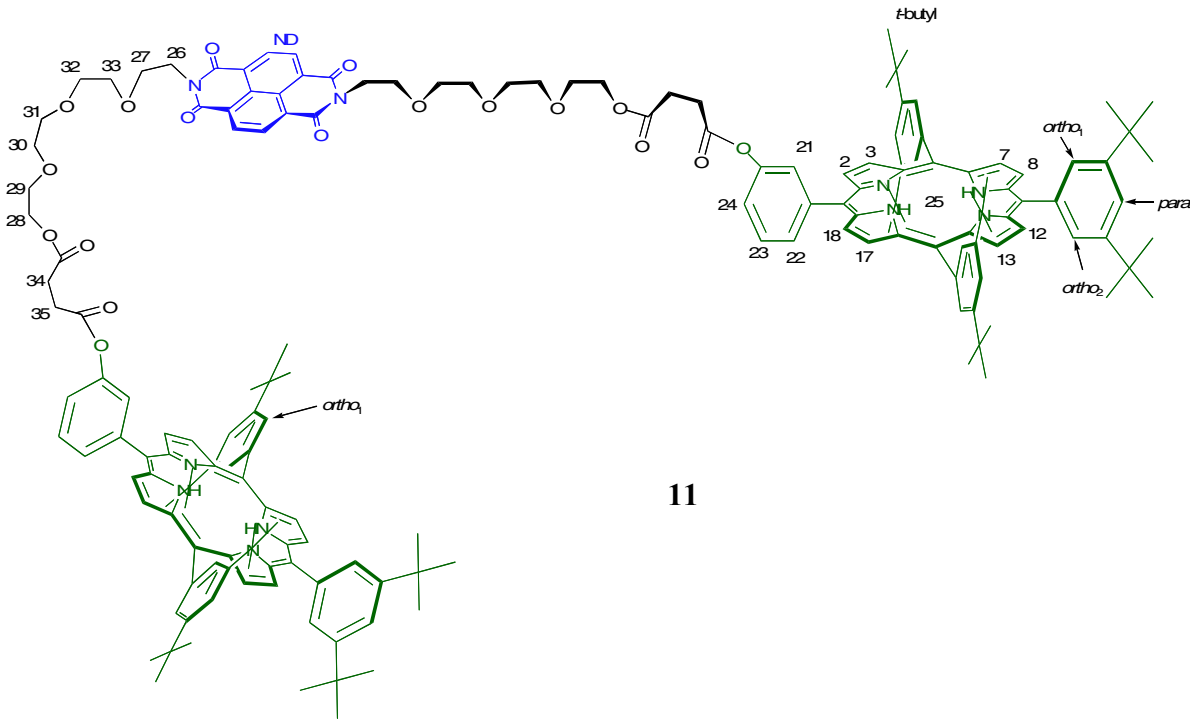


Figure S4 ^{1}H NMR spectrum and numbering scheme for **10**.



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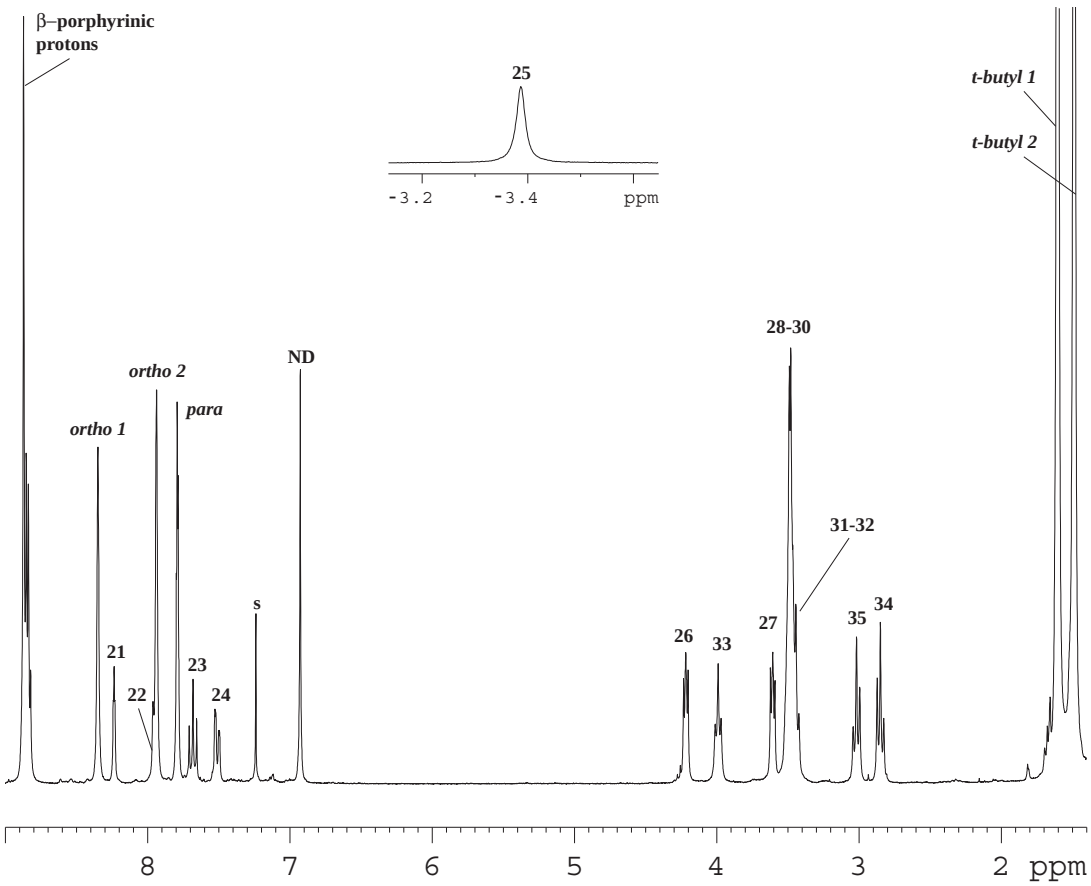

Figure S5 ^{1}H NMR spectrum and numbering scheme for **11**

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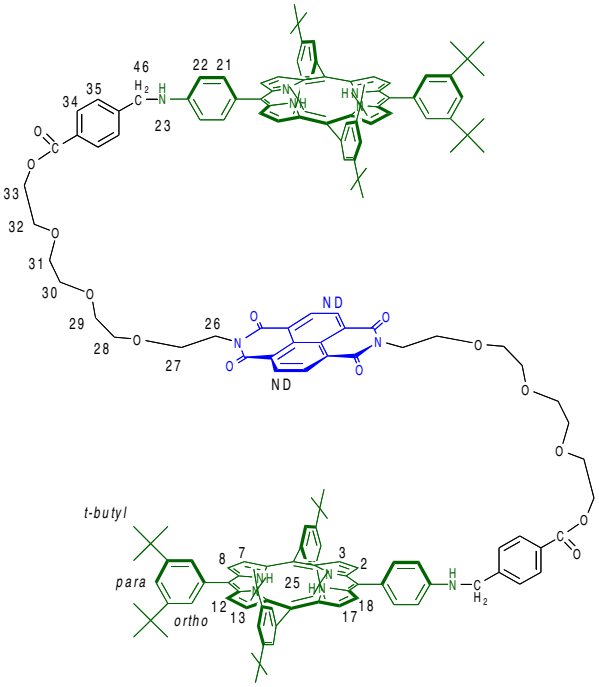


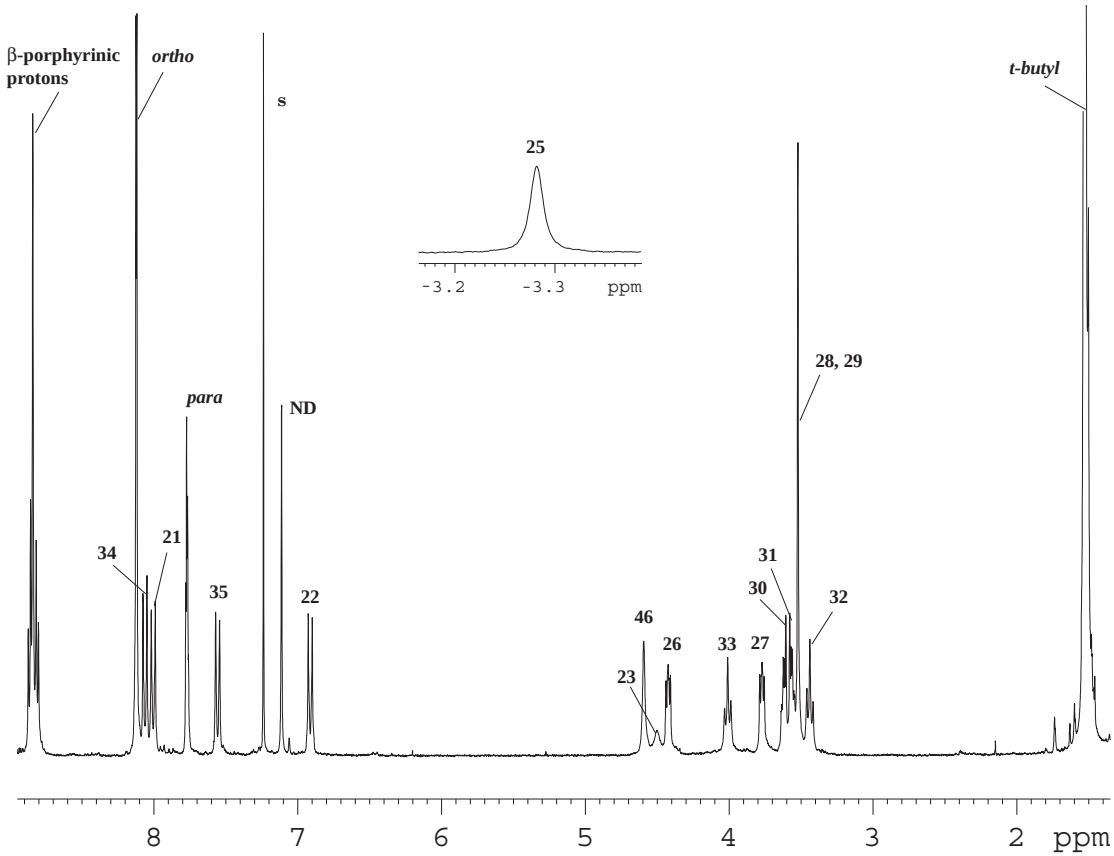


Figure S6 ^{1}H NMR spectrum and numbering scheme for **12**.

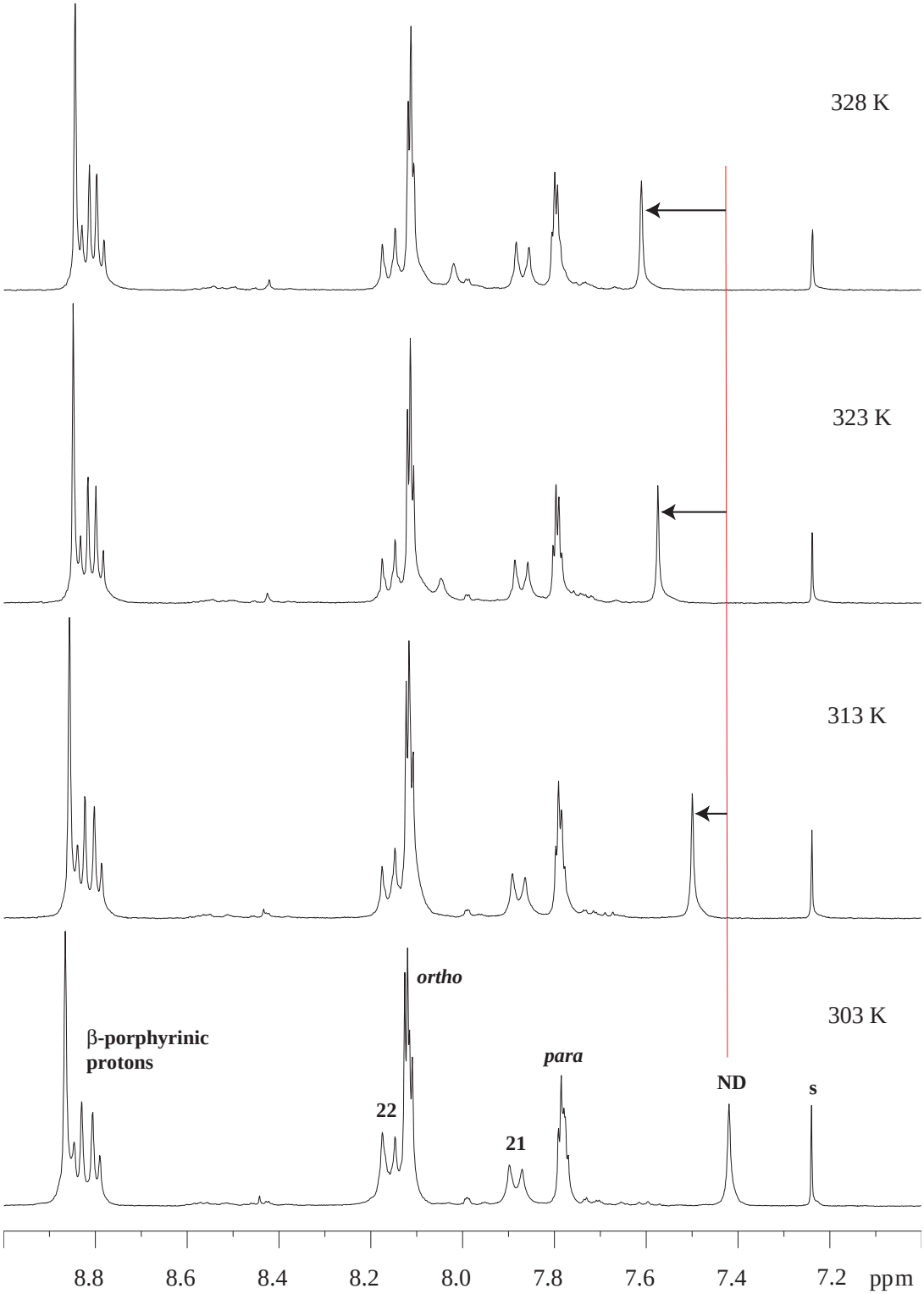


Figure S7 Variable high temperature ^{1}H NMR experiments for **10**. As the temperature is increased, the naphthalene protons ND are shifted downfield. Selected peaks are identified by the non-systematic numbering shown above (s is $CDCl_3$).

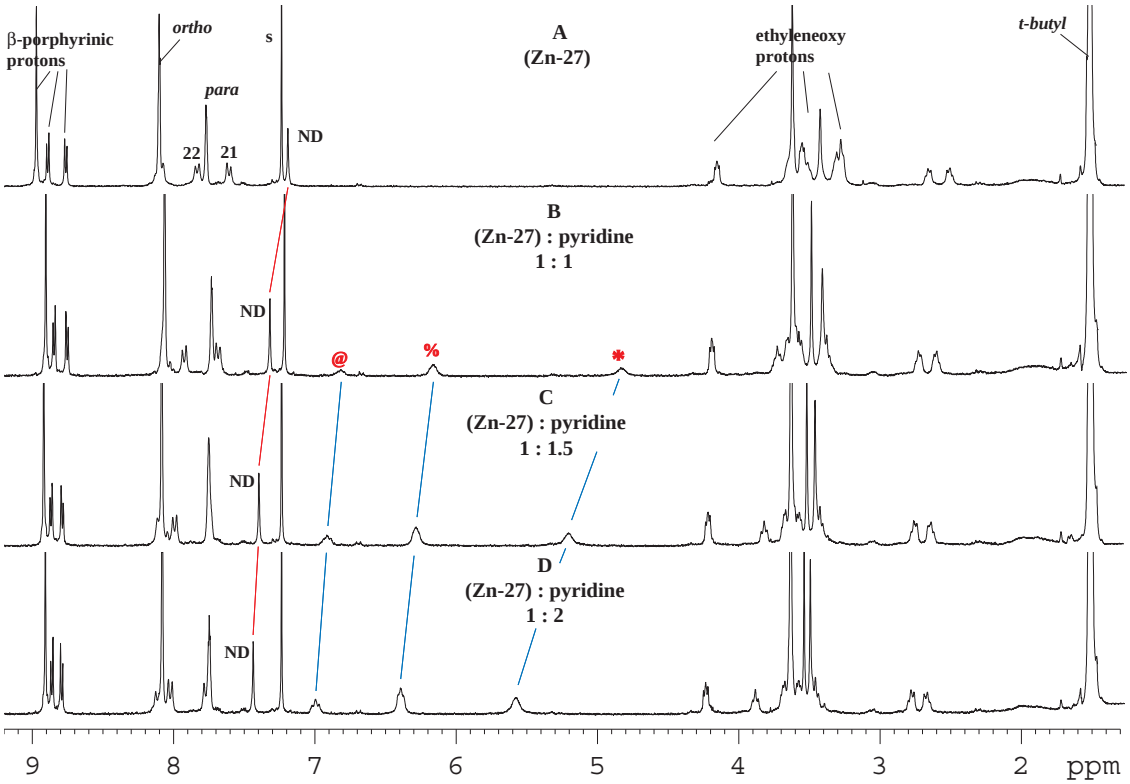


Figure S8 ^{1}H NMR spectra ($CDCl_3$) showing gradual downfield shifts of the diimide protons upon binding of pyridine (py) to **Zn10.**

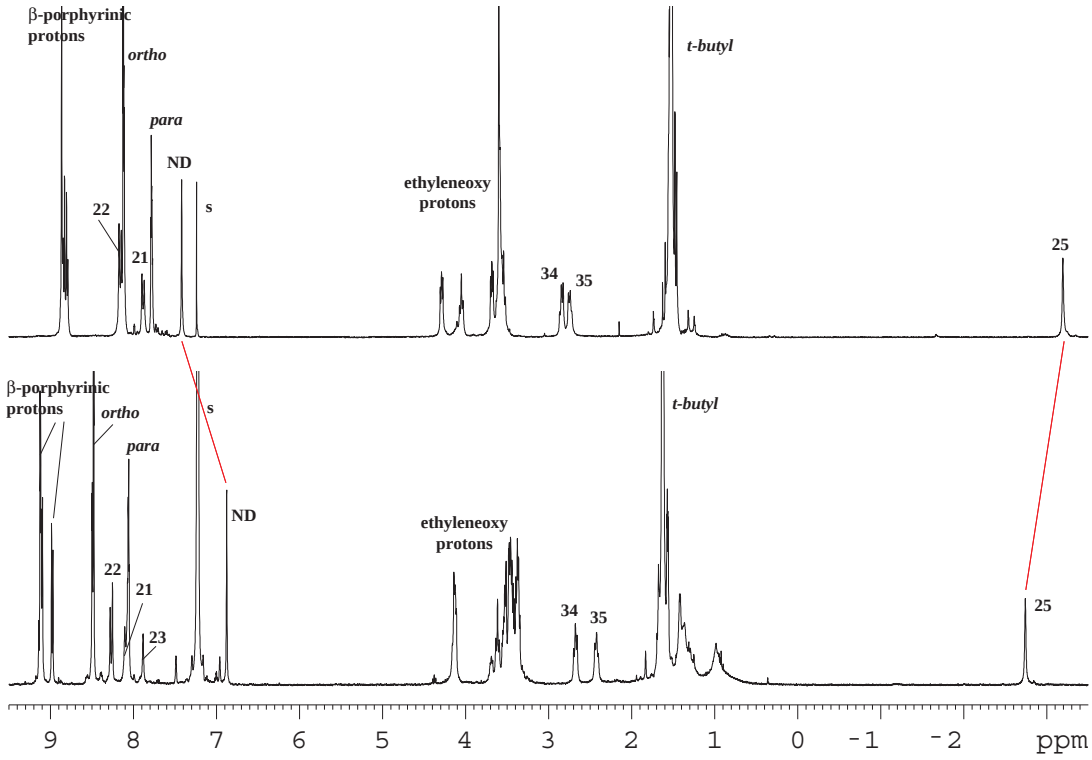


Figure S9 ^{1}H NMR spectra of **10** in: $CDCl_3$ (top) and d_6-benzene (bottom).

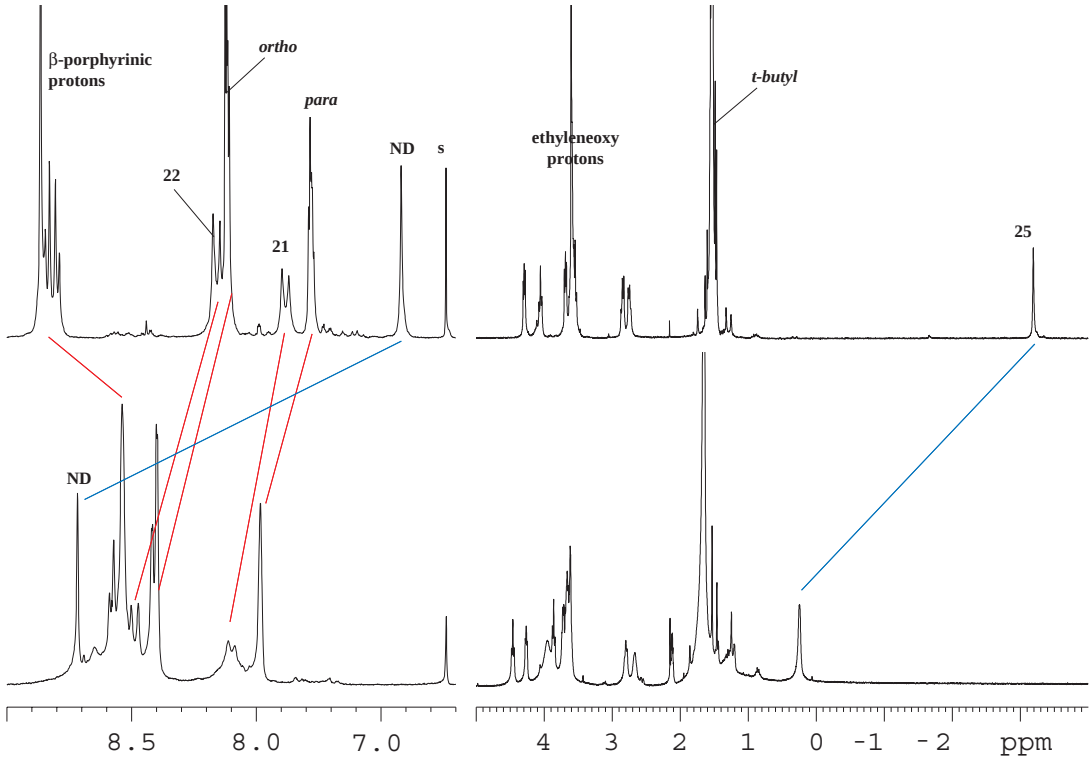
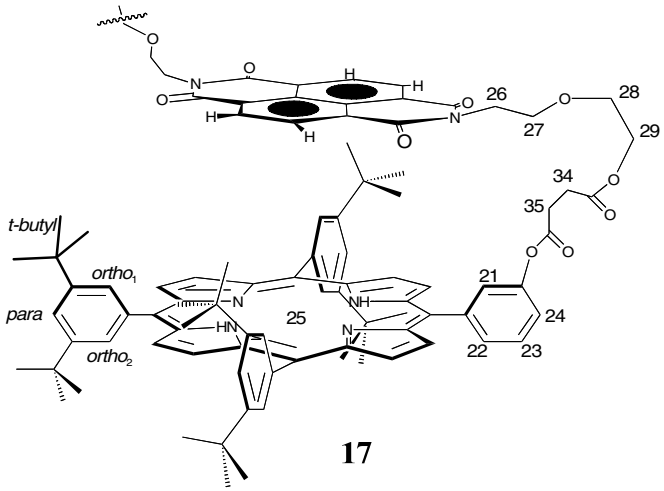


Figure S10 ^{1}H NMR spectra ($CDCl_3$) showing the shift of naphthodiimide protons ND to about its original frequency upon protonation (bottom), relative to that of the free-base derivative **10** (top).



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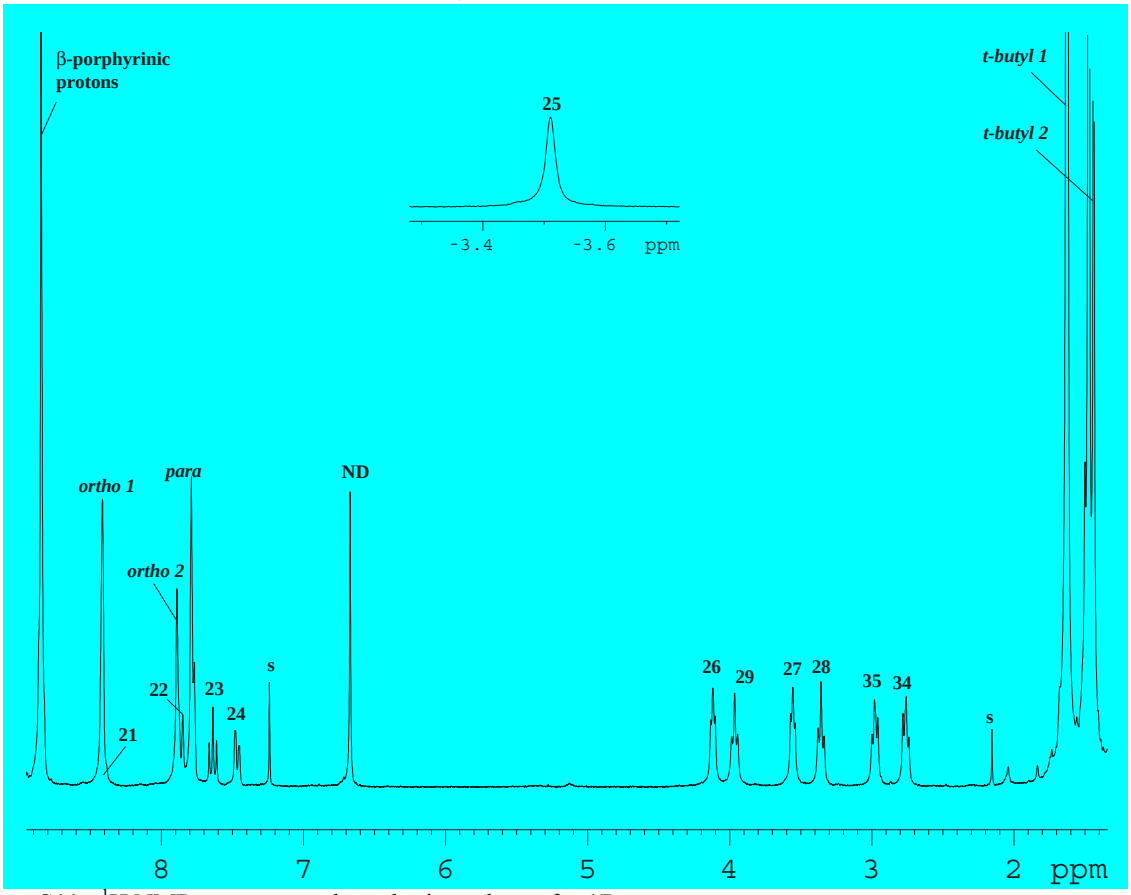

Figure S11 ^{1}H NMR spectrum and numbering scheme for **17**.

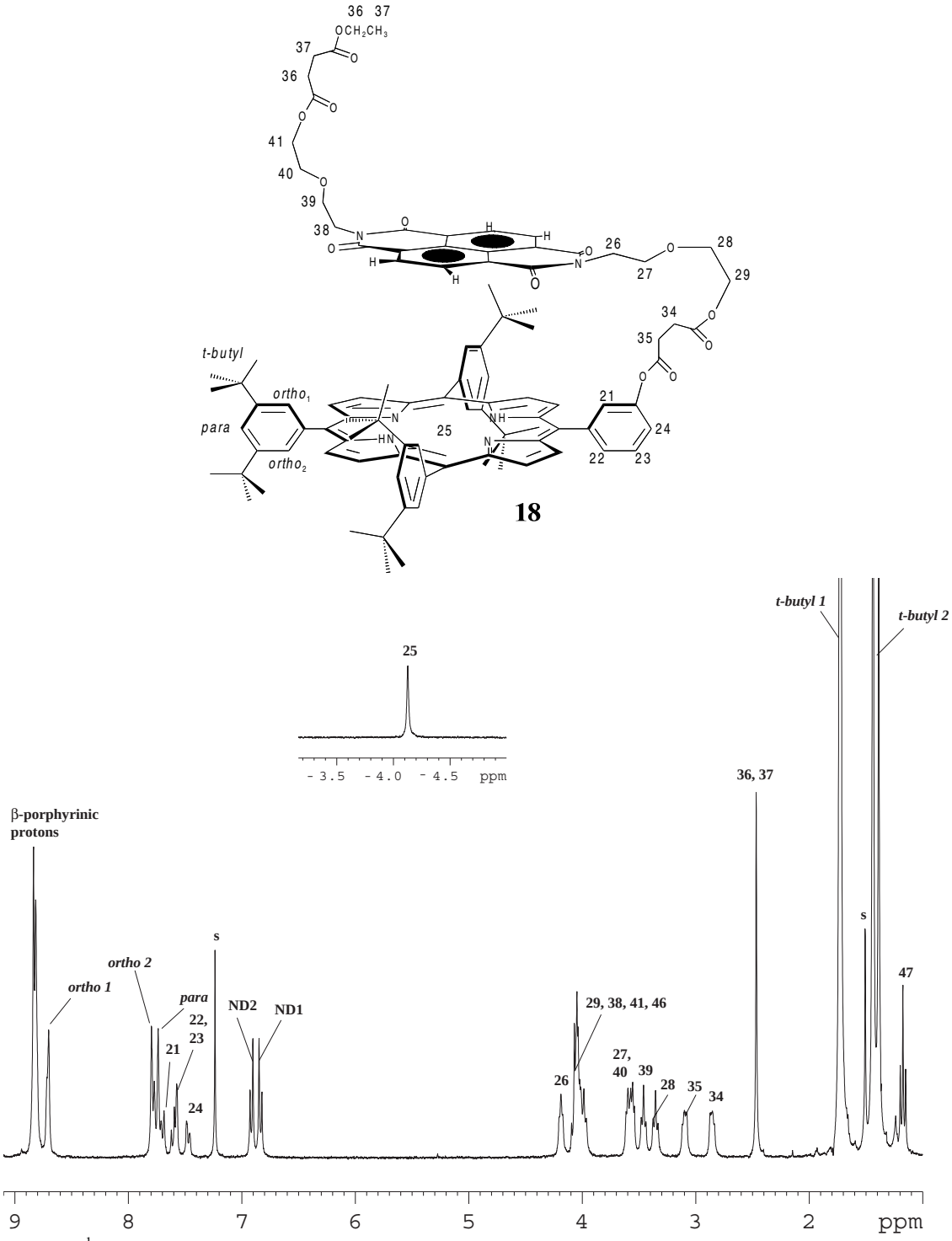

Figure S12 ^{1}H NMR spectrum and numbering scheme for **18**.