Electronic Supplementary Information

Structure-chiroptical property relationship of kinetically labile camphor-derivative β-diketone Yb(III) complexes: The adducts coexist as diastereomers or not?

Yiji Lin,^a Shigang Wan,^a Fang Zou,^a Yuekui Wang^{*b} and Hui Zhang^{*a}

^a State Key Laboratory of Physical Chemistry of Solid Surface and Department of Chemistry, College of Chemistry and Chemical Engineering, Xiamen University, Xiamen 361005, China.

^b Key Laboratory of Chemical Biology and Molecular Engineering of the Education Ministry, Institute of Molecular Science, Shanxi University, Taiyuan 030006, China.

Figure S1. The coordination polyhedra of complexes I and II.

Figure S2. The observed and simulated X-ray diffraction patterns of complex I.

Figure S3. The observed and simulated X-ray diffraction patterns of complex III.

Figure S4. The bulk sample solid-state CD spectra of complexes I and II.

Figure S5. Solid-state electronic absorption spectrum of complex I.

Figure S6. Solid-state electronic absorption spectrum of complex III.

Figure S7. Solution electronic absorption spectrum of complex I.

Figure S8. Solution electronic absorption spectrum of complex III.

Figure S9. ORTEP structures of complexes I and II.

Figure S10. ORTEP structures of complexes III and IV.



Figure S1. The coordination polyhedra of complexes I and II.

Figure S2. The observed (red) and simulated (black) X-ray powder diffraction spectra of complex I.



Figure S3. The observed (red) and simulated (black) X-ray powder diffraction spectra of complex III.



Figure S4. Solid-state CD spectra of the bulk samples I and II (black: I; red: II).



Figure S5. Solid-state absorption spectrum of complex I.



Figure S6. Solid-state absorption spectrum of complex III.



Figure S7. Solution absorption spectrum of complex I.



Figure S8. Solution absorption spectrum of complex III.



Figure S9. ORTEP structure of complexes I and II (- C_3F_7 groups are omitted for clarity). Thermal ellipsoids are shown at the 30% probability level.



Figure S10. ORTEP structure of complexes **III** (left) and **IV** (right) ($-C_3F_7$ groups are omitted for clarity). Thermal ellipsoids are shown at the 30% probability level.

