

Supplementary information for

**Mechanistic insight into selective hydrogenolysis of sorbitol to
propylene glycol and ethylene glycol on supported Ru catalysts**

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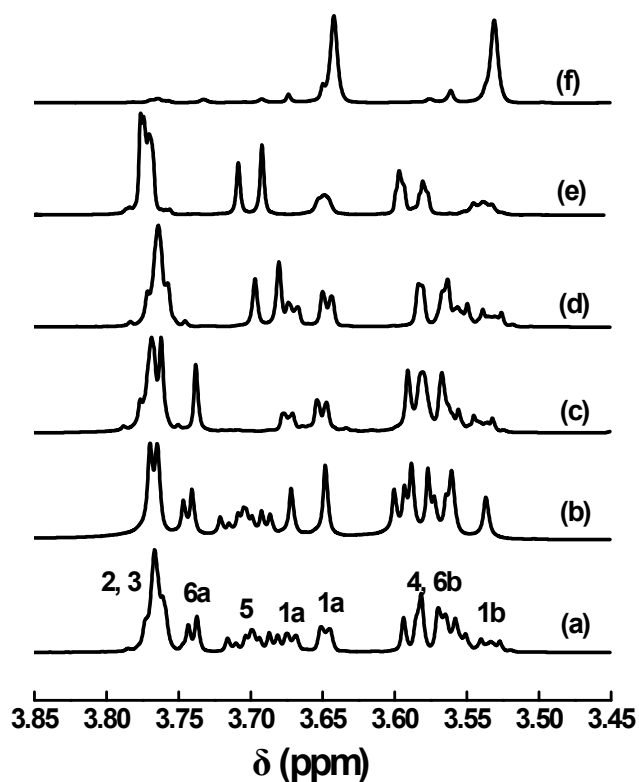


Fig. S1 ¹H NMR spectra of different D-sorbitol molecules in D₂O at room temperature.

(a) D-sorbitol; (b) D-sorbitol-2-d₁; (c) D-sorbitol-5-d₁; (d) D-sorbitol-6,6-d₂; (e) D-sorbitol-1,6,6-d₃; (f) D-sorbitol-1,2,3,4,5,6,6-d₇. All H atoms at different C positions are labeled in the spectrum of D-sorbitol (a). For D-sorbitol-1,2,3,4,5,6,6-d₇, its H atoms at C-1 position was substituted by deuterium equally at C(1)-H_a and C(1)-H_b positions.

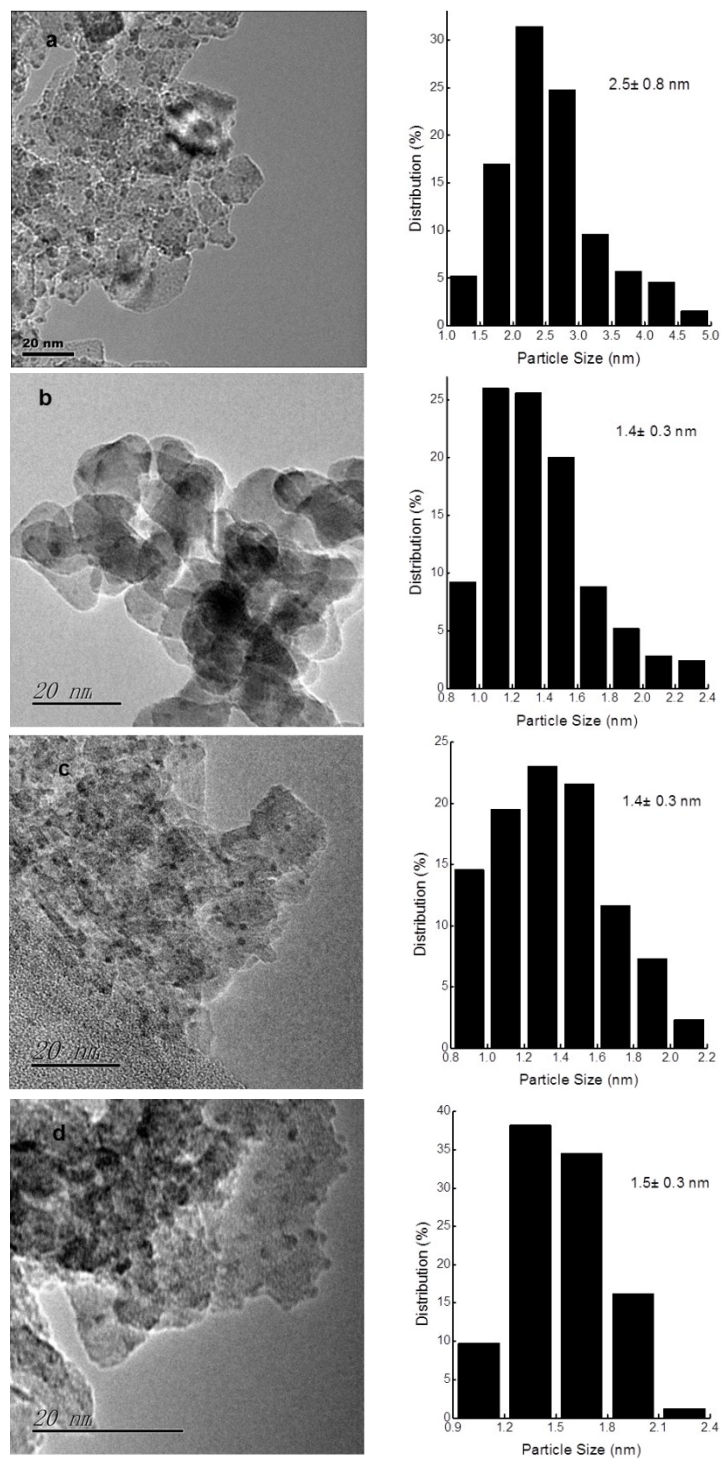


Fig. S2 TEM micrographs and histograms of Ru particle size distribution for different Ru catalysts (scale bar=20 nm). (a) Ru/MgO, 2.5 ± 0.8 nm; (b) Ru/m-ZrO₂, 1.4 ± 0.3 nm; (c) Ru/Al₂O₃, 1.4 ± 0.3 nm; (d) Ru/TiO₂, 1.5 ± 0.3 nm.

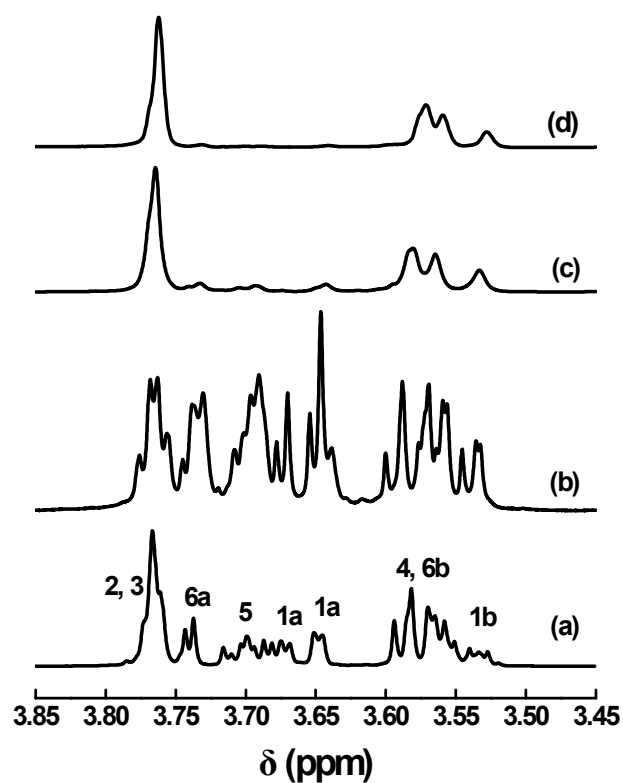


Fig. S3 ^1H NMR spectra of different D-sorbitol molecules in D_2O at room temperature.

(a) D-sorbitol; (b) The sorbitol molecules separated from the solution of D-sorbitol-1,2,3,4,5,6,6- d_7 hydrogenolysis in H_2O and H_2 ; (c) The sorbitol molecules separated from the solution of D-sorbitol hydrogenolysis in D_2O and H_2 ; (d) The sorbitol molecules separated from the solution of D-sorbitol hydrogenolysis in D_2O and D_2 . All H atoms at different C positions are labeled in the spectrum of D-sorbitol (a).

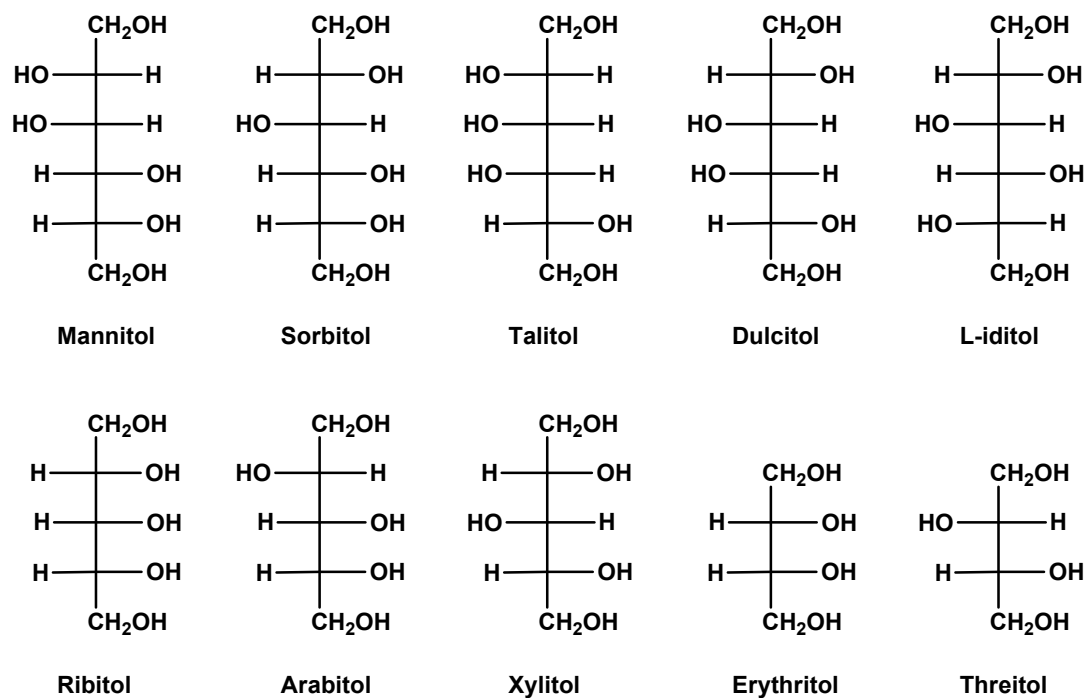


Fig. S4 Fischer projections of hexitols, pentitols and tetritols

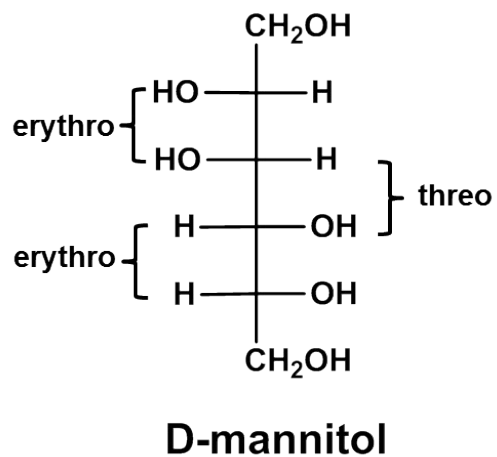


Fig. S5 Erythro and threo sequences of hydroxyl groups of D-mannitol.