

Supporting Information (SI):

Highly-efficient Ru/Al-SBA-15 Catalysts with Strong Lewis Acid Sites for the Water-assisted Hydrogenation of *p*-Phthalic Acid

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Table S1. Structural parameters of supports (SBA-15, AS-x).

Entry	Supports	S _{BET} (m ² /g)	D _p (nm)	V (cm ³ /g)
1	SBA-15	640	6.6	1.03
2	AS-50	827	6.5	0.89
3	AS-10	744	6.6	0.89
4	AS-3	611	6.5	0.85
5	AS-1	499	6.5	0.76
6	AS-0.5	365	6.5	0.55

Table S2. The amounts of acid sites determined by pyridine-FTIR.

Catalyst	The amounts of acid sites (umol/g)					
	150 °C			400 °C		
	B	L	Total	B	L	Total
Ru/SBA-15	2.8	35.8	38.6	0.9	2.6	3.5
Ru/AS-50	4.0	61.4	65.4	1.3	27.9	29.2
Ru/AS-10	4.1	83.4	87.5	1.4	49.4	50.8
Ru/AS-3	3.5	160.1	163.6	1.5	81.1	82.6
Ru/AS-1	2.7	92.7	95.4	1.7	33.7	35.4
Ru/AS-0.5	1.7	86.5	88.2	1.0	26.4	27.4

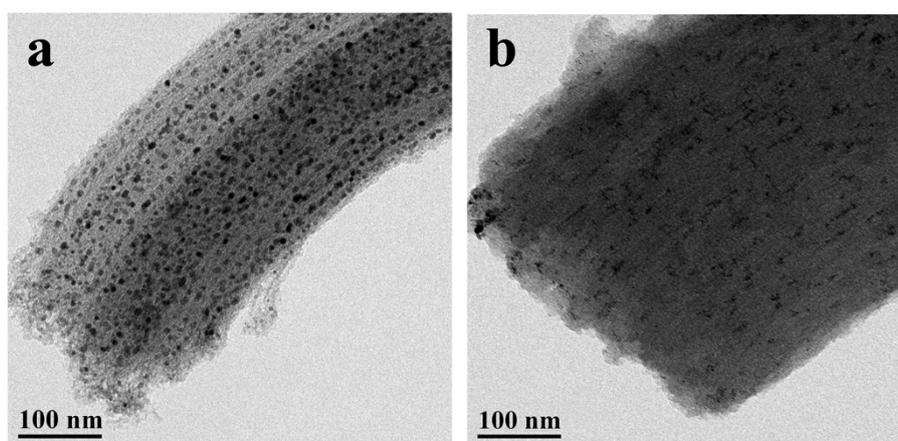


Fig. S1. TEM images of (a) Ru/AS-3 and (b) Ru/AS-3-U (U refers to used catalyst for fourteen times).

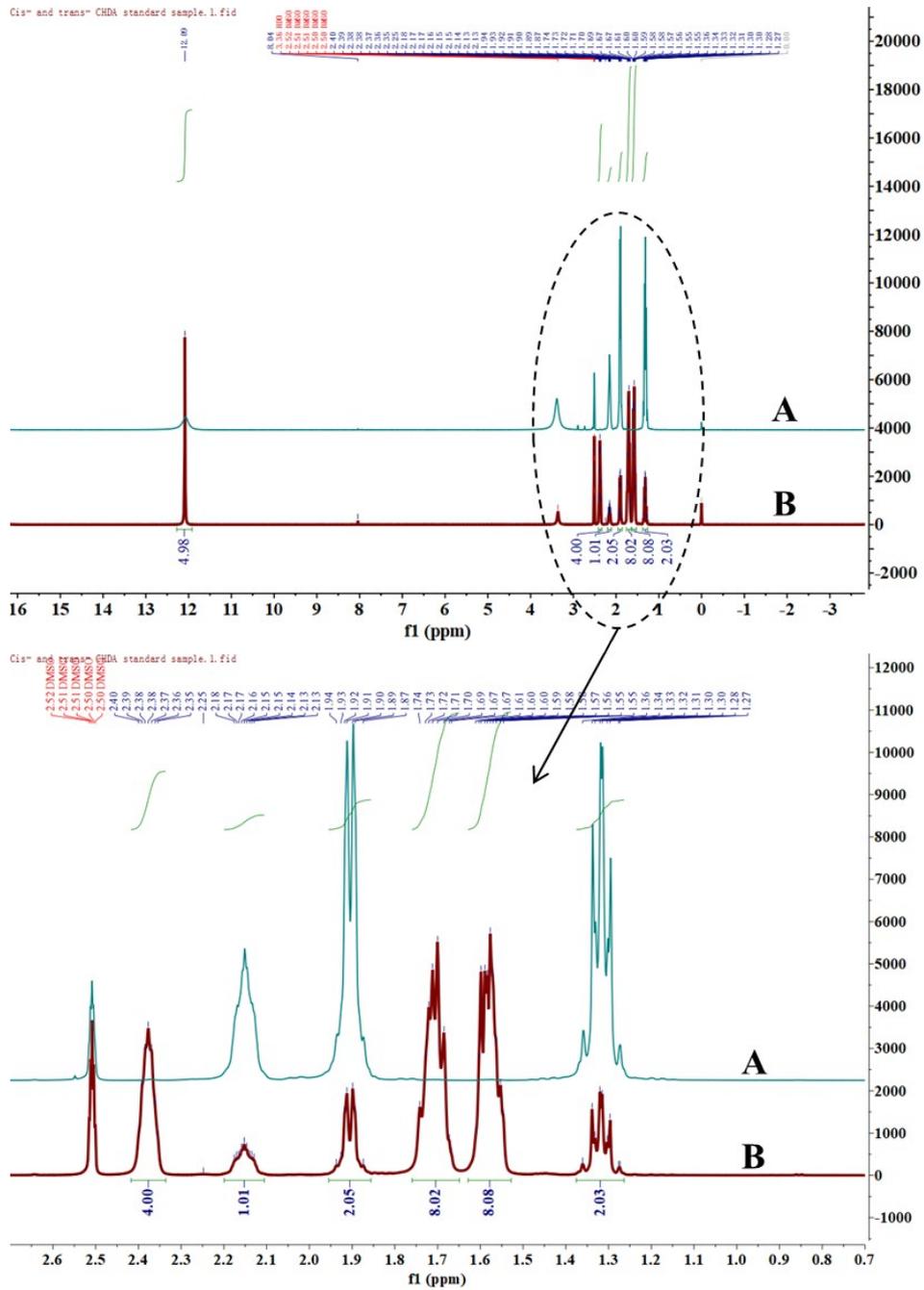


Fig. S2. ^1H NMR spectra of CHDA standard samples. (A) *trans*-CHDA and (B) *cis*- and *trans*-CHDA mixture. $\text{DMSO-}d_6$ was used to dissolve the samples.

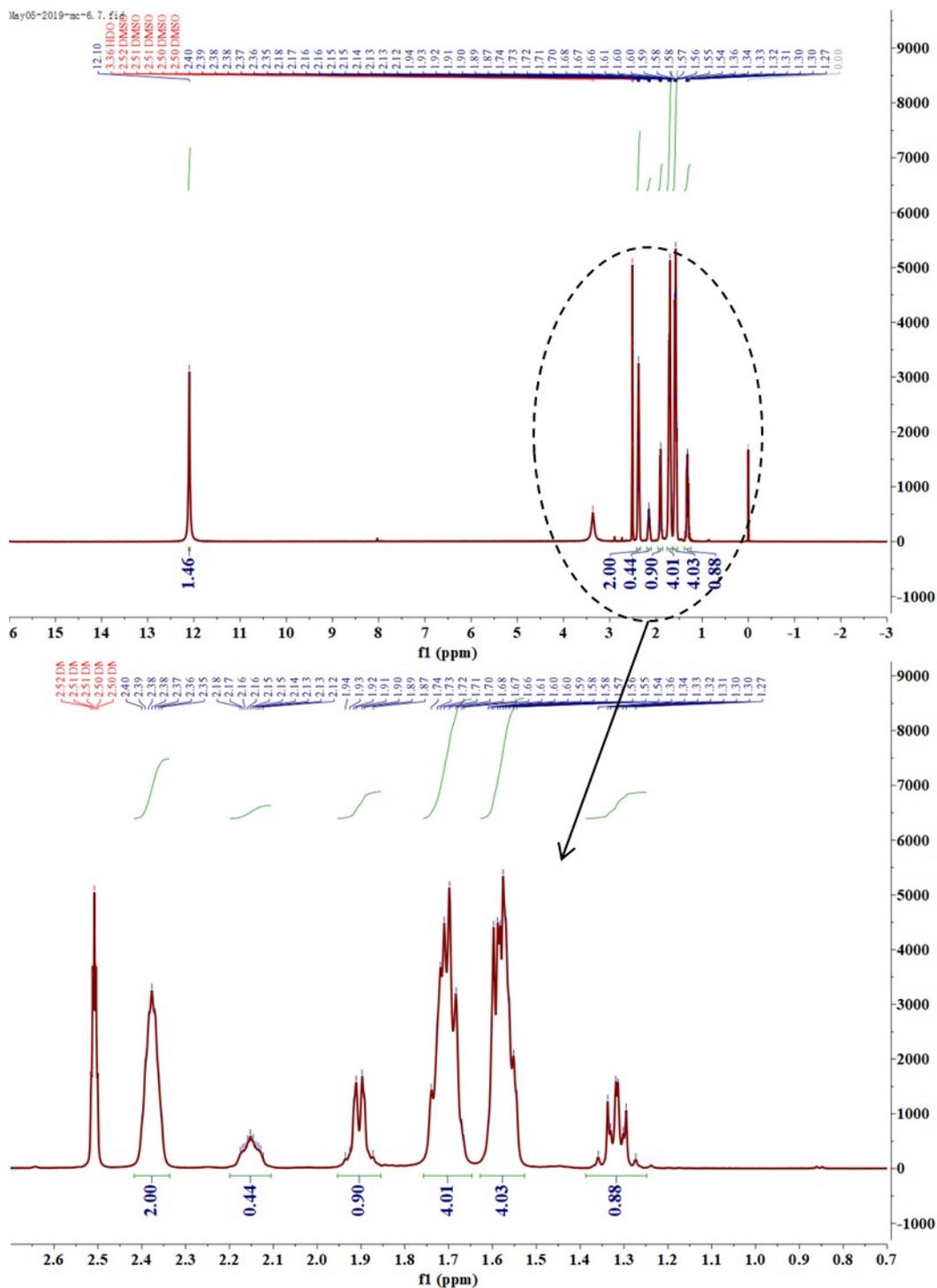


Fig. S3. ^1H NMR spectra of the hydrogenation product (*p*-phthalic acid was hydrogenated in H_2O solvent for 2h). $\text{DMSO-}d_6$ was used to dissolve the product.

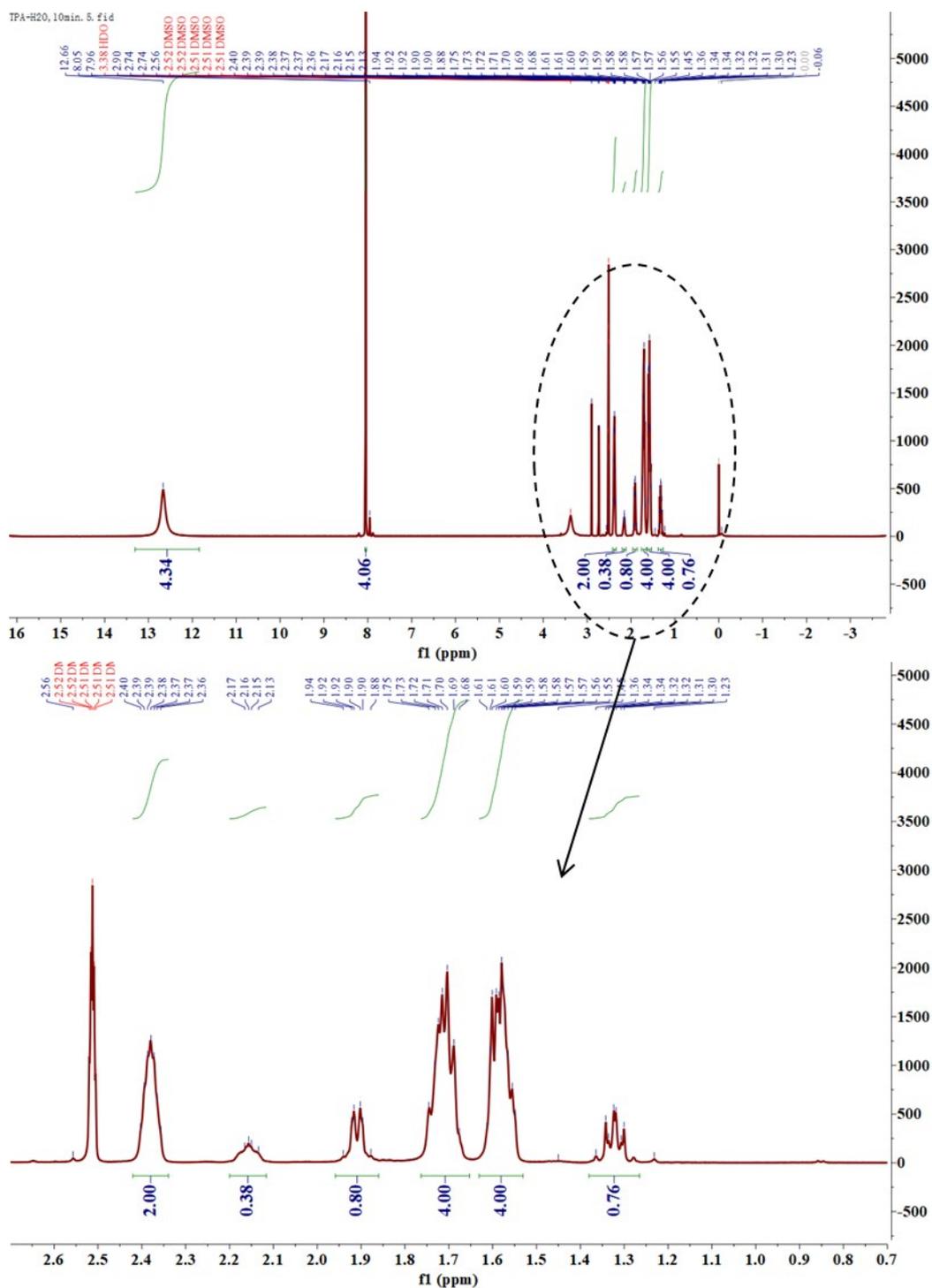


Fig. S5. ^1H NMR spectra of the hydrogenation product (*p*-phthalic acid was hydrogenated in H_2O solvent for 30min). $\text{DMSO-}d_6$ was used to dissolve the product.

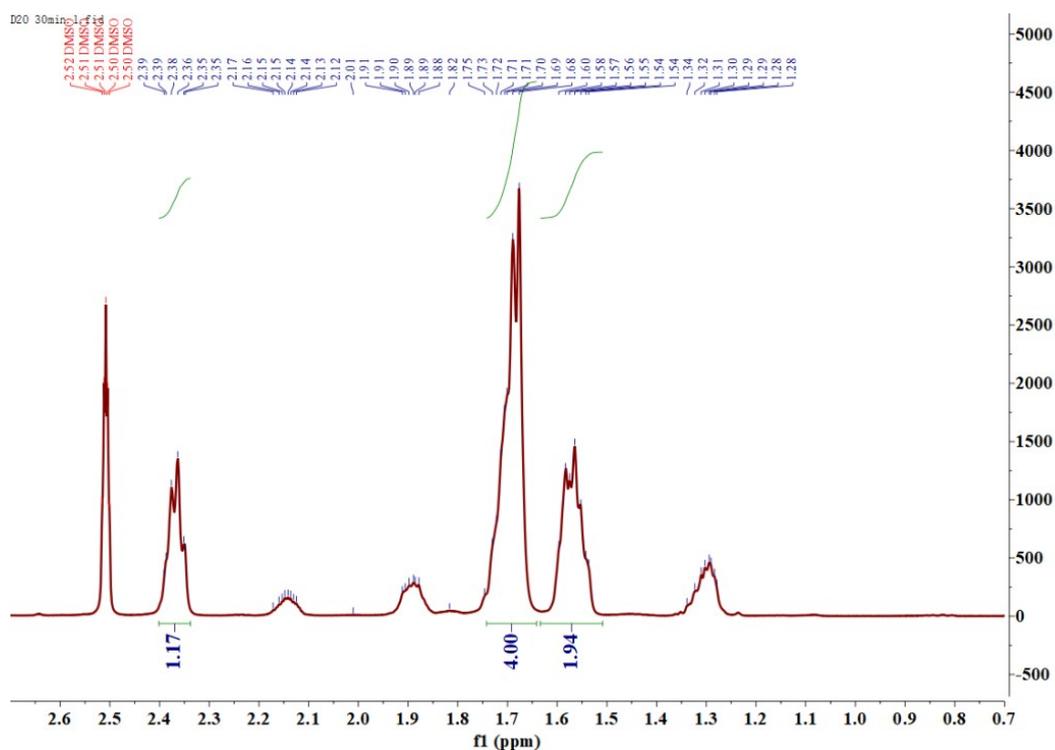


Fig. S6. ^1H NMR spectra of the hydrogenation product (*p*-phthalic acid was hydrogenated in D_2O solvent for 2h). DMSO-d_6 was used to dissolve the product.

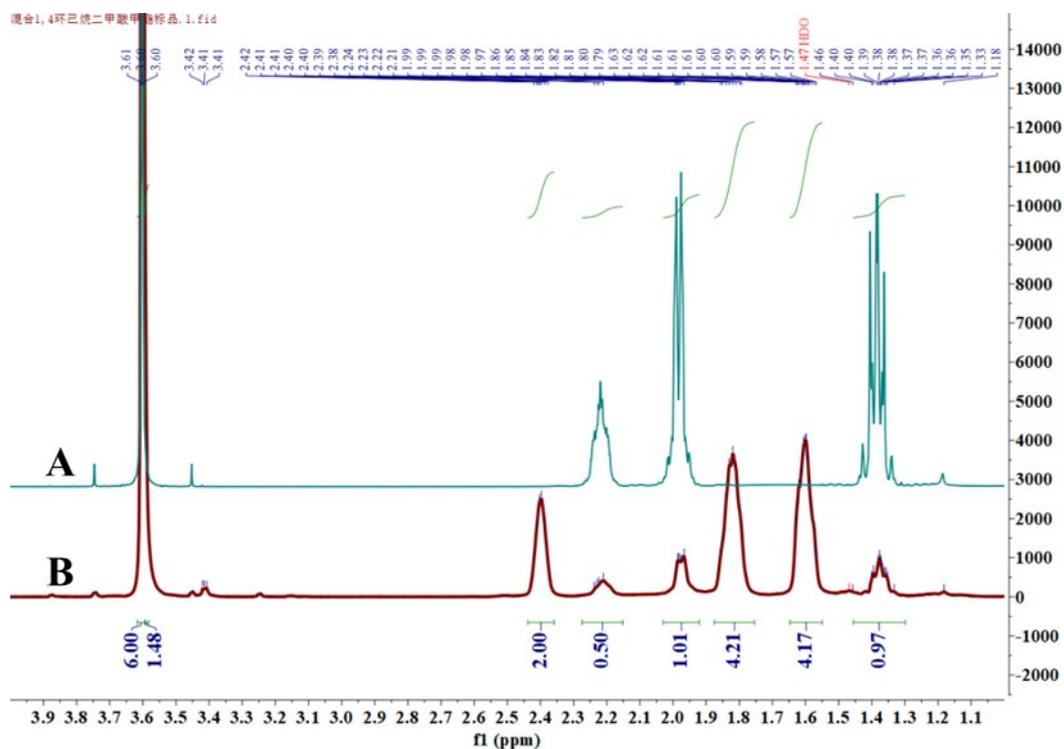


Fig. S7. ^1H NMR spectrum of standard dimethyl 1,4-cyclohexanedicarboxylate samples. (A) *trans*-dimethyl 1,4-cyclohexanedicarboxylate and (B) *trans*- and *cis*-dimethyl 1,4-cyclohexanedicarboxylate mixture. CDCl_3 was used to dissolve the product.

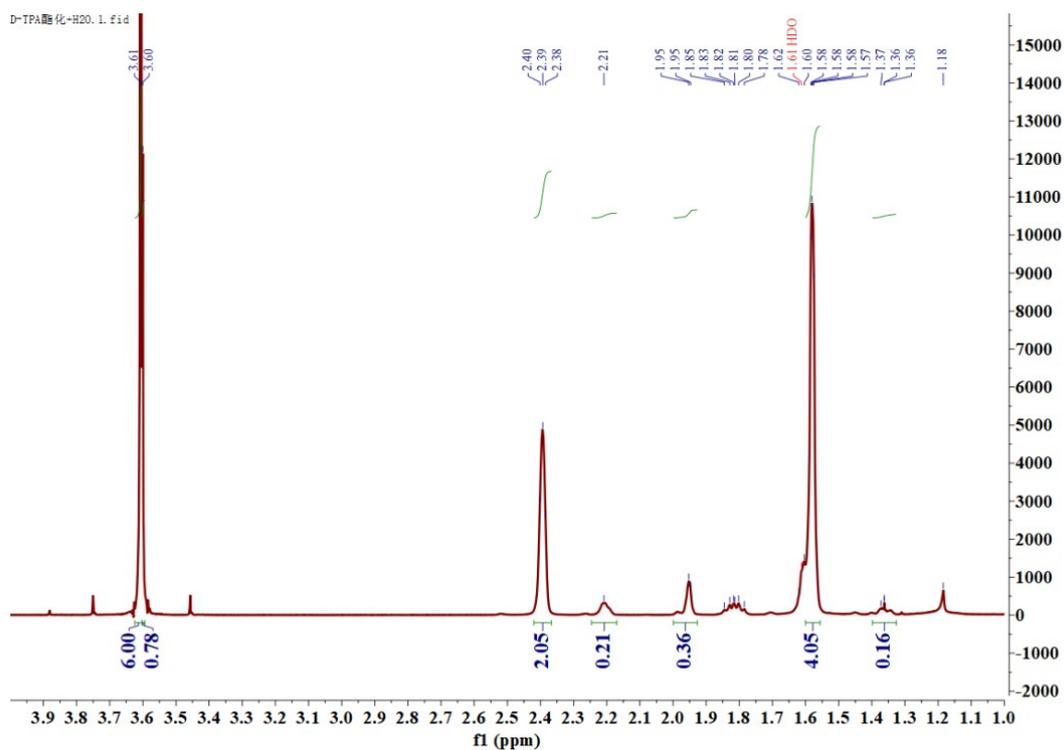


Fig. S8. ^1H NMR spectrum of product (*p*-phthalic acid- d_4 was hydrogenated in H_2O following by esterifying to dimethyl 1, 4-cyclohexanedicarboxylate). CDCl_3 was used to dissolve the product.

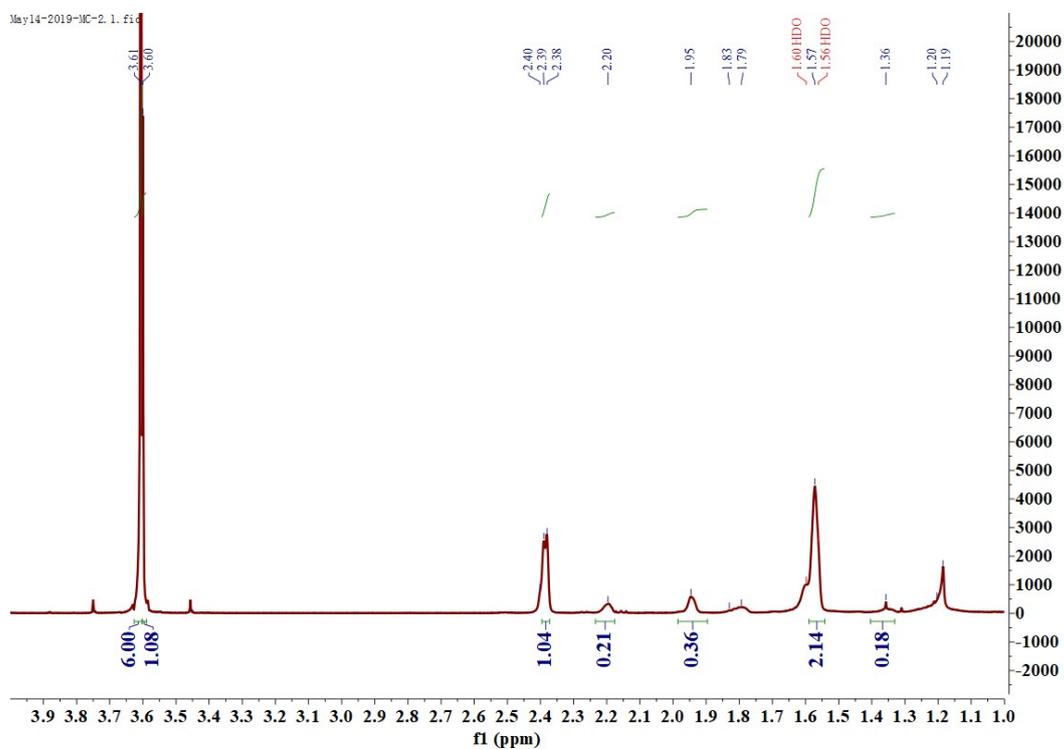


Fig. S9. ^1H NMR spectrum of product (*p*-phthalic acid- d_4 was hydrogenated in D_2O following by esterifying to dimethyl 1, 4-cyclohexanedicarboxylate). CDCl_3 was used to dissolve the product.