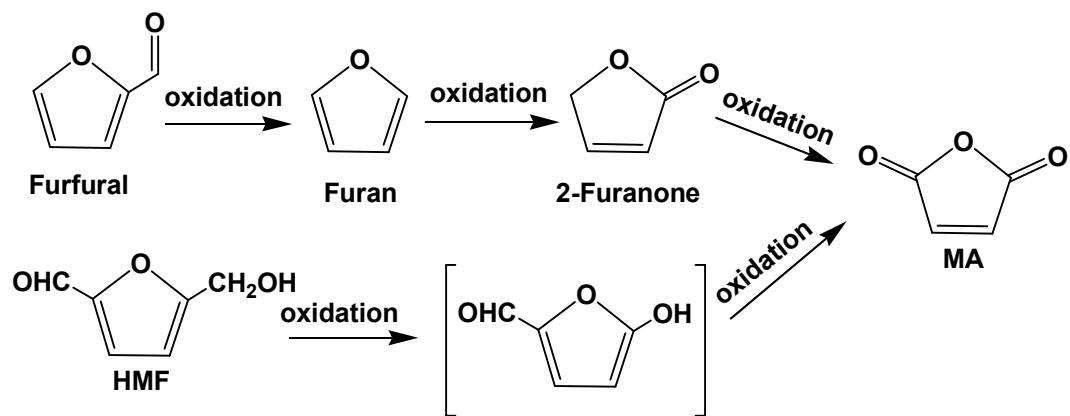


**Supporting Information**

For

**One-pot synthesis of 1,4-butanediol via the deep hydrogenation of  
maleic anhydride over Cu-xMo/SiO<sub>2</sub> catalysts**

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**Scheme S1** Selective oxidation of furfural or HMF to maleic anhydride.

**Table S1** Quantitative acidity according to NH<sub>3</sub>-TPD for various Cu-xMo/SiO<sub>2</sub> catalysts.

Catalyst	Acid amount (mmol·g <sup>-1</sup> )			
	Weak	Moderate	Strong	Total acidity
Cu/SiO <sub>2</sub>	0.15	0.14	0.08	0.37
Cu-0.02Mo/SiO <sub>2</sub>	0.22	0.39	0	0.61
Cu-0.03Mo/SiO <sub>2</sub>	0.22	0.28	0	0.50
Cu-0.1Mo/SiO <sub>2</sub>	0.35	0.36	0	0.71
Cu-0.25Mo/SiO <sub>2</sub>	0.22	0.37	0	0.59
Mo/SiO <sub>2</sub>	0	0	0	0

**Table S2** Quantitative hydrogen uptake for various Cu-xMo/SiO<sub>2</sub> catalysts.

Sample	Cu content (wt%)	Amount of H <sub>2</sub> consumption (mmol/g <sub>Cu</sub> )
Cu/SiO <sub>2</sub>	37.7	4.38
Cu-0.03Mo/SiO <sub>2</sub>	32.1	5.54
Cu-0.25Mo/SiO <sub>2</sub>	31.7	5.84

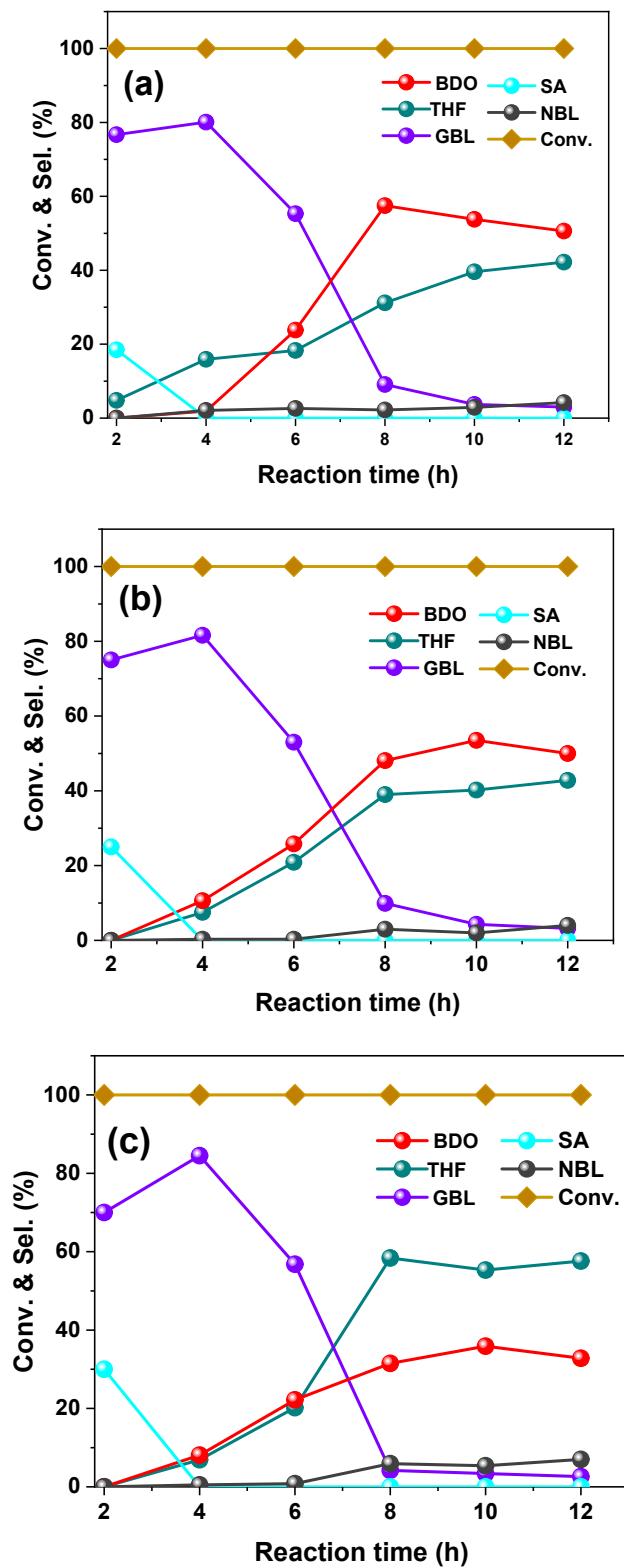
**Table S3** Physicochemical properties of the spent Cu-0.03Mo/SiO<sub>2</sub> catalyst.

Catalyst	S <sub>BET</sub> <sup>a</sup> (m <sup>2</sup> ·g <sup>-1</sup> )	Vp <sup>a</sup> (cm <sup>3</sup> ·g <sup>-1</sup> )	Cu loading <sup>b</sup> (wt%)	d <sub>Cu</sub> <sup>c</sup> (nm)
Cu-0.03Mo/SiO <sub>2</sub> -used	291	0.50	31.4	14.1
Cu-0.03Mo/SiO <sub>2</sub> -fresh	357	0.61	32.1	3.8

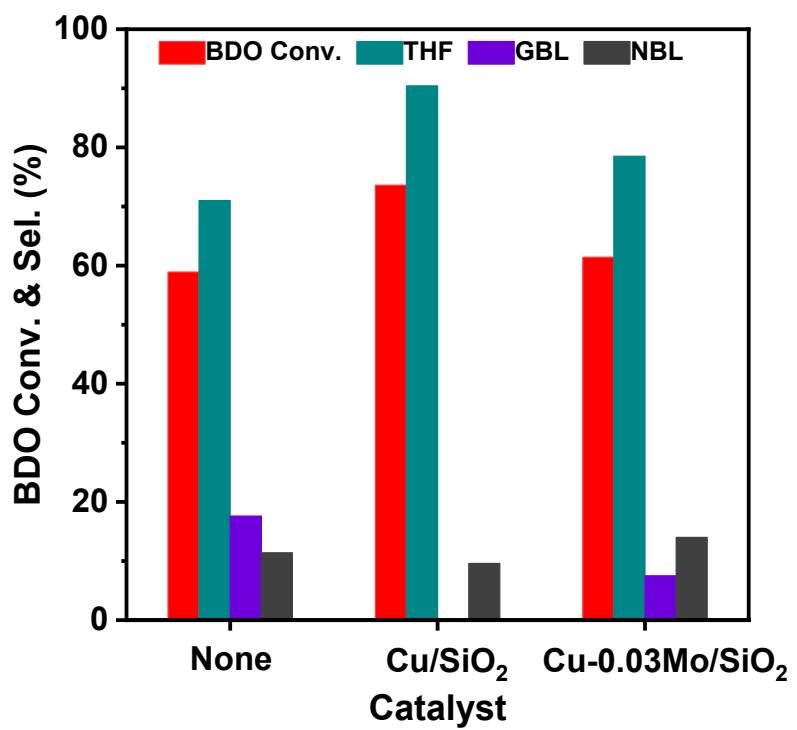
<sup>a</sup> Determined by N<sub>2</sub> physisorption.

<sup>b</sup> Determined by ICP-OES analysis.

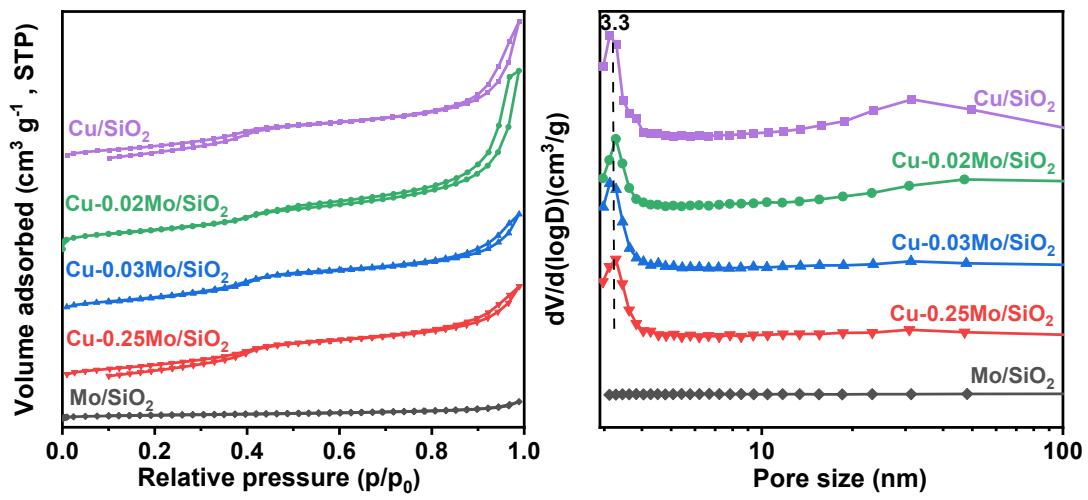
<sup>c</sup> Cu particle size was calculated by the Scherrer equation.



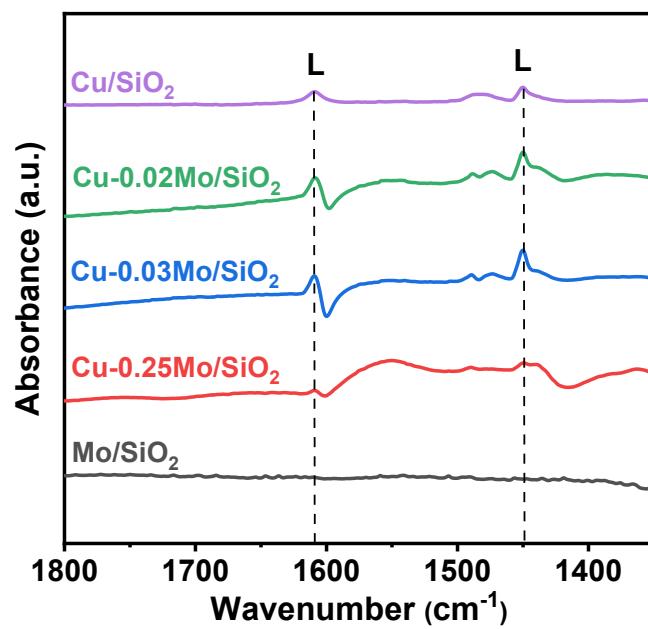
**Fig. S1** Kinetic profiles for the hydrogenation of MA with (a) Cu-0.02Mo/SiO<sub>2</sub>, (b) Cu-0.1Mo/SiO<sub>2</sub>, and (c) Cu-0.25Mo/SiO<sub>2</sub> catalysts. General conditions: 0.2 g of catalyst, 0.2 g of MA, 10 mL of 1,4-dioxane, 5 MPa H<sub>2</sub>, 200 °C.



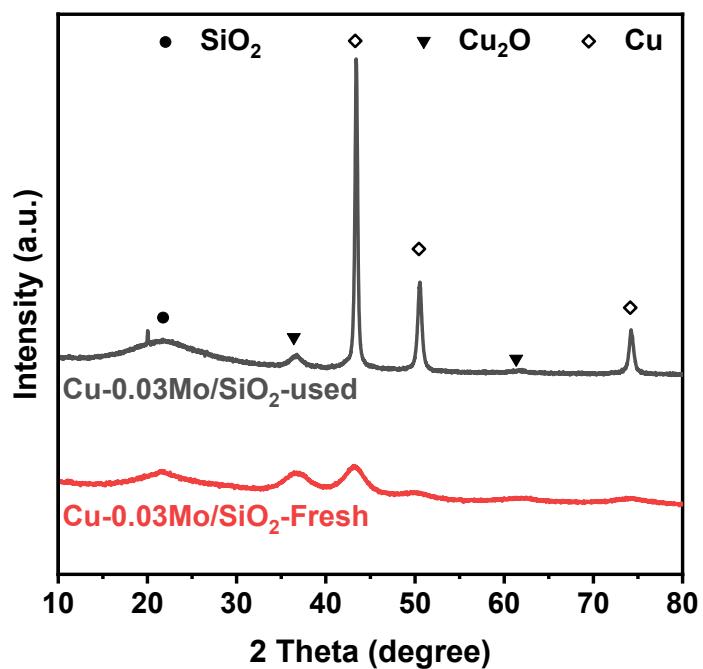
**Fig. S2** Product distribution in the hydrogenation of BDO. Reaction conditions: 0.2 g of catalyst, 0.2 g of BDO, 10 mL of 1,4-dioxane, 200 °C, 5 MPa H<sub>2</sub>, 10 h.



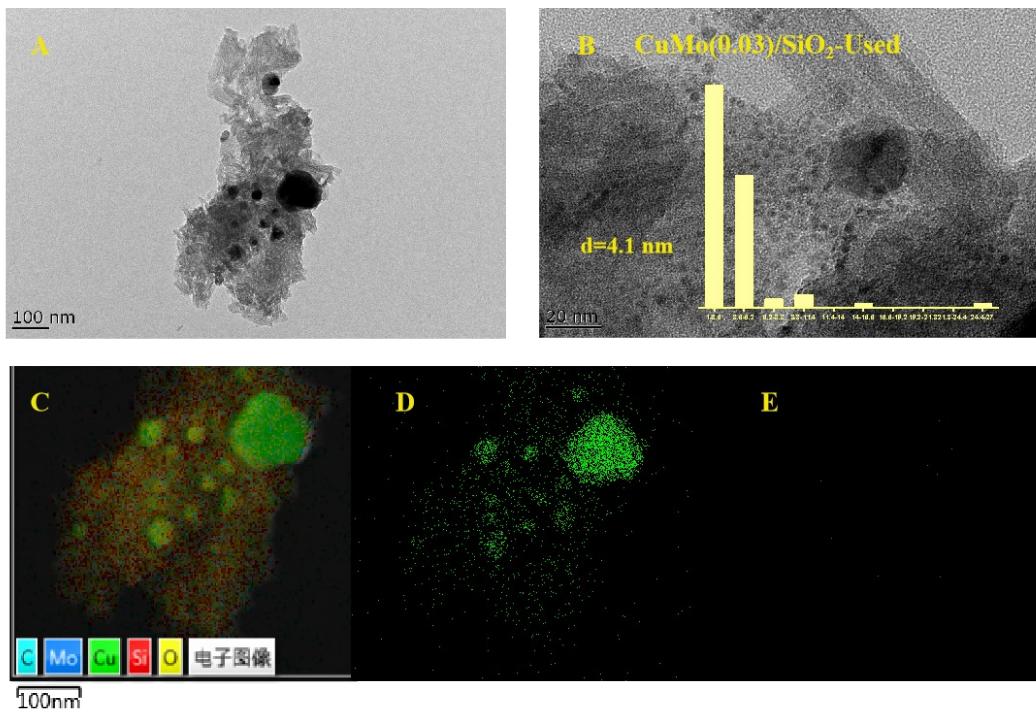
**Fig. S3** (a) N<sub>2</sub> adsorption-desorption isotherms and (b) pore size distributions of various Cu-xMo/SiO<sub>2</sub> catalysts.



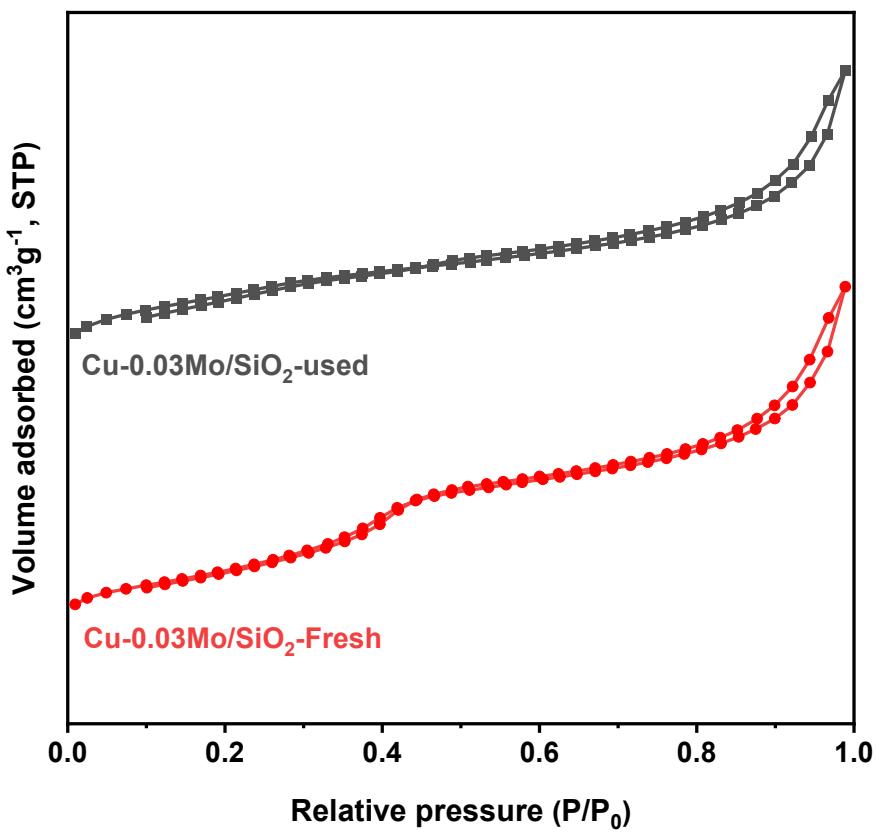
**Fig. S4** FTIR spectra of pyridine adsorbed on  $\text{Cu-xMo/SiO}_2$  catalysts at 150 °C.



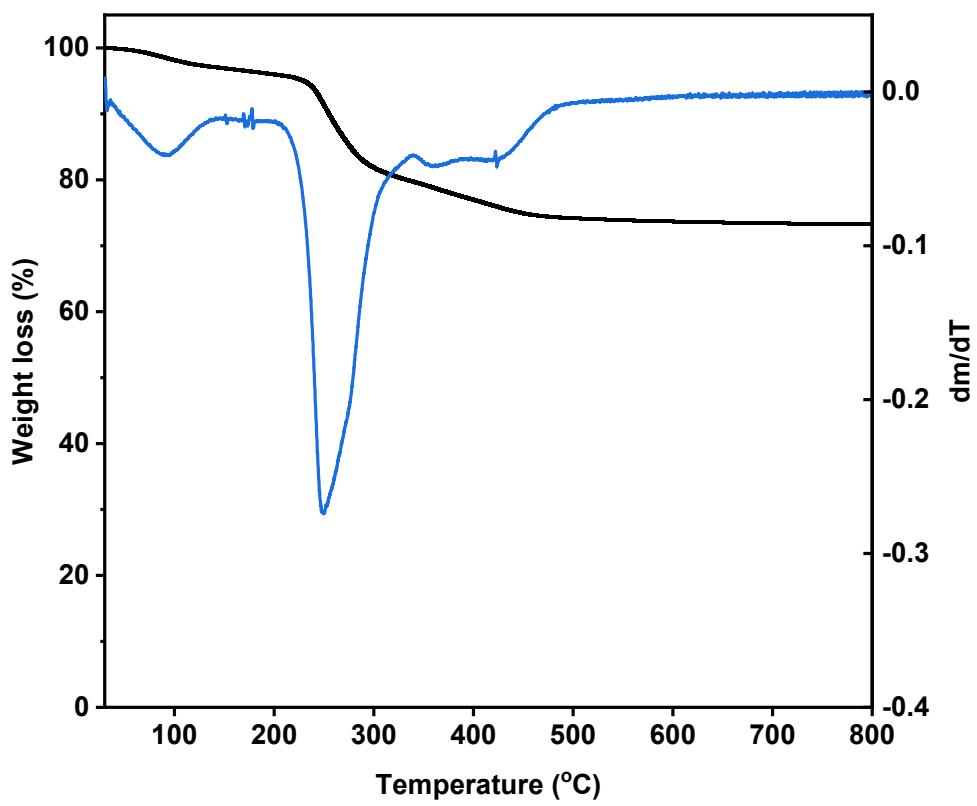
**Fig. S5** XRD patterns of Cu-0.03Mo/SiO<sub>2</sub> before and after long-term stability test.



**Fig. S6** (A) TEM image and (B) high-resolution TEM image of used Cu-0.03Mo/SiO<sub>2</sub>. (C) HAADF-STEM-EDX mapping of used Cu-0.03Mo/SiO<sub>2</sub> and (D) Cu element, (E) Mo element.



**Fig. S7** N<sub>2</sub> sorption isotherms of the Cu-0.03Mo/SiO<sub>2</sub> catalyst before and after long-term stability test.



**Fig. S8** TG curve of the used Cu-0.03Mo/SiO<sub>2</sub> catalyst.