

Supporting information for

Insights into the influence of various impurities on the stability of Cu/SiO₂ catalyst for the hydrogenation of dimethyl oxalate to ethylene glycol

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S. Wu and Y. Li are equally contributed to this work.

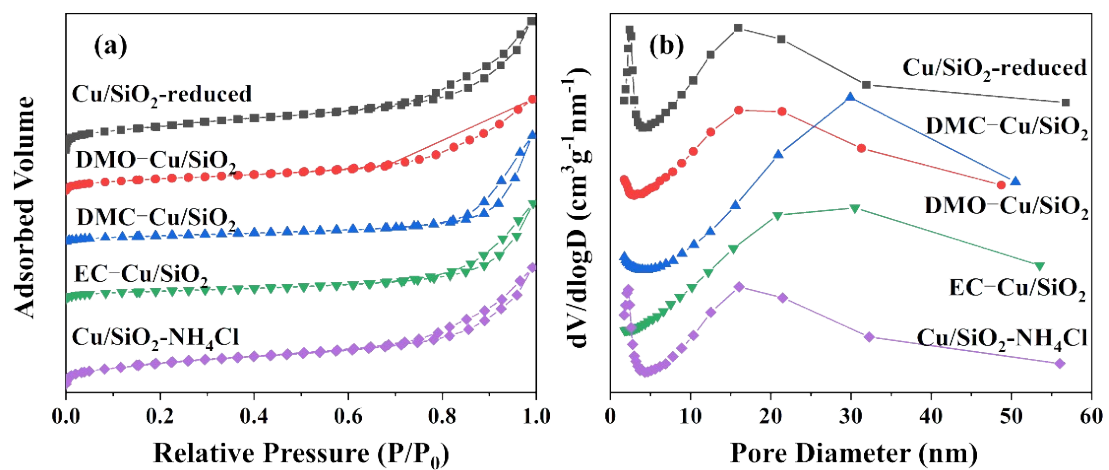


Figure S1. (a) N₂ adsorption-desorption isotherms and (b) pore size distribution curves of the reduced and used Cu/SiO₂ catalysts.

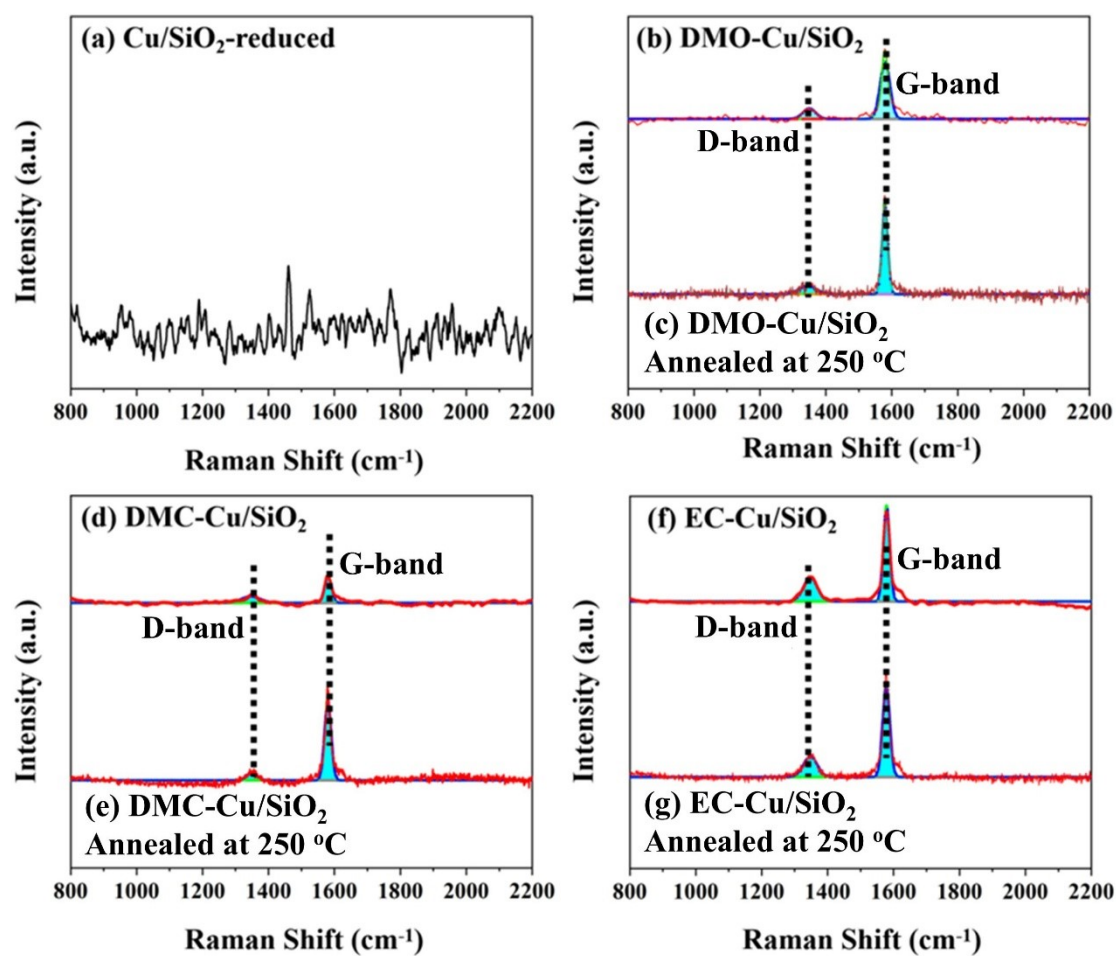


Figure S2. Raman spectra of the reduced and used Cu/SiO₂ catalysts with different feedstocks. (a) reduced, (b, d, f) after reaction, (c, e, g) after reaction and annealed at 250 °C.

Table S1. Actual weight loss of the used Cu/SiO₂ catalysts calculated from TG analysis.

Catalyst	Total loss (%)	Soft coke (%)	Heavy coke (%)
DMO-Cu/SiO ₂	2.79	2.19	0.60
DMC-Cu/SiO ₂	4.44	3.33	1.11
EC-Cu/SiO ₂	7.13	6.42	0.71
Cu/SiO ₂ -NH ₄ Cl	1.97	1.35	0.62

Table S2. I_D/I_G values derived from Raman spectra of the used Cu/SiO₂ catalysts with different feedstocks.

	DMO-Cu/SiO ₂	DMC-Cu/SiO ₂	EC-Cu/SiO ₂
I_D/I_G (before annealing)	0.18	0.24	0.26
I_D/I_G (after annealing)	0.09	0.16	0.22

Table S3. Main carbonaceous deposit species in the used Cu/SiO₂ catalysts and the properties.

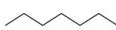



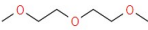


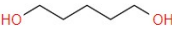
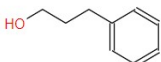
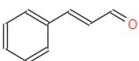
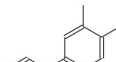
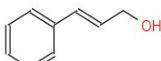
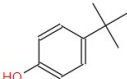
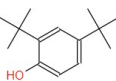

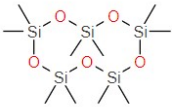

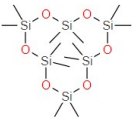
Name	Structural formula	Molecular formula	Relative molecular mass	Boiling point (°C)
n-heptane		C ₇ H ₁₆	100.2	98
toluene		C ₇ H ₈	92.1	111
p-xylene		C ₈ H ₁₀	106.2	138
1,2,4-trimethylbenzene		C ₉ H ₁₂	120.2	168
diethylene glycol dimethyl ether		C ₆ H ₁₄ O ₃	134.2	159
1,2-dimethyl-3-ethylbenzene		C ₁₀ H ₁₄	134.2	194
1,4-butanediol		C ₄ H ₁₀ O ₂	90.1	228
1,5-pentanediol		C ₅ H ₁₂ O ₂	104.1	239
3-phenylpropanol		C ₉ H ₁₂ O	136.2	238
trans-cinnamaldehyde		C ₉ H ₈ O	132.2	250
3,4-xylylaldehyde		C ₁₀ H ₁₂ O	148.2	245
cinnamyl alcohol		C ₉ H ₁₀ O	134.2	250
4-tert-butylphenol		C ₁₀ H ₁₄ O	150.2	238
2,4-di-tert-butylphenol		C ₁₄ H ₂₂ O	206.3	265

Table S3. Main carbonaceous deposit species in the used Cu/SiO₂ catalysts and the

properties (continued).

Name	Structural formula	Molecular formula	Relative molecular mass	Boiling point (°C)
18-crown-6		C ₁₂ H ₂₄ O ₆	264.3	395
decamethylcyclopentasiloxane		C ₁₀ H ₃₀ O ₅ Si ₅	370.8	90
dodecane		C ₁₂ H ₂₆	170.3	215
dodecamethylcyclohexasiloxane		C ₁₂ H ₃₆ O ₆ Si ₆	444.9	245