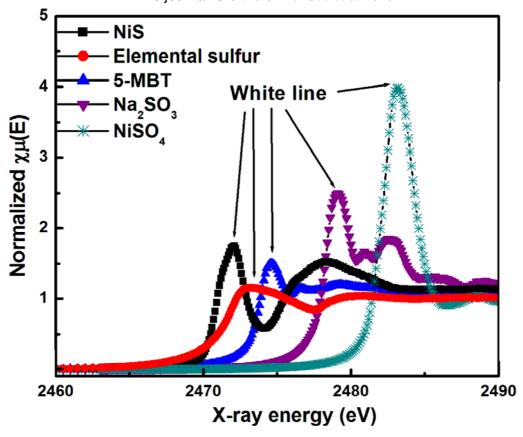
Sulfur poisoning mechanism of steam reforming catalysts: an x-ray absorption near edge structure (XANES) spectroscopic study

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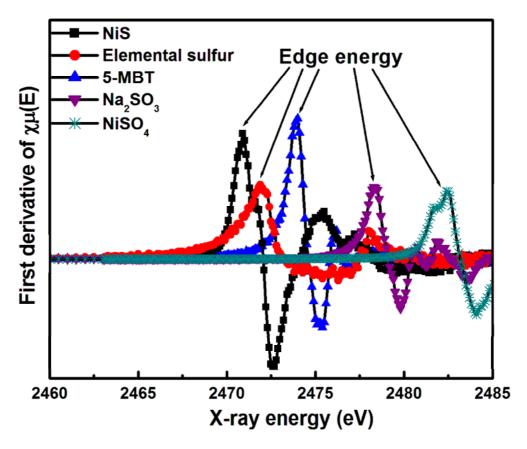


Figure S1. Sulfur standards. Top: absorption spectra (showing white line); Bottom: first derivative of the absorption spectra (showing edge energy).

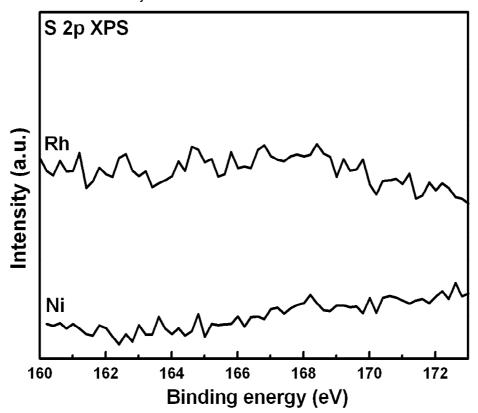


Figure S2 S 2p XPS of the used 2wt%Rh/CeO₂-Al₂O₃ and 10wt%Ni/CeO₂-Al₂O₃ after steam reforming reaction with sulfur at 800 °C for 55 h. Sulfur peaks would appear somewhere between 161-169 eV depending on their oxidation states and chemical environment.

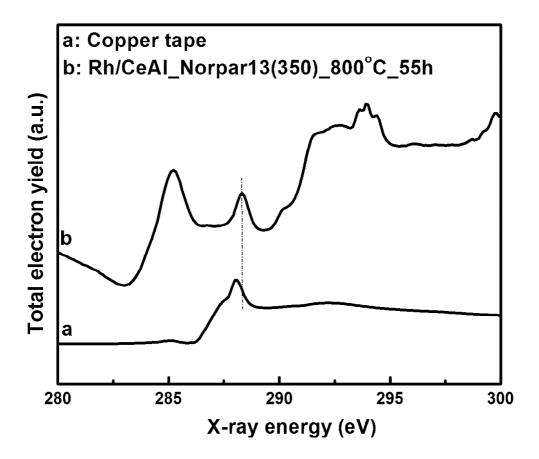


Figure S3. Carbon K-edge XANES spectrum of the copper tape used for sample preparation together with the spectrum of the Rh catalyst tested in the steam reforming reaction at 800 °C for 55 hours.

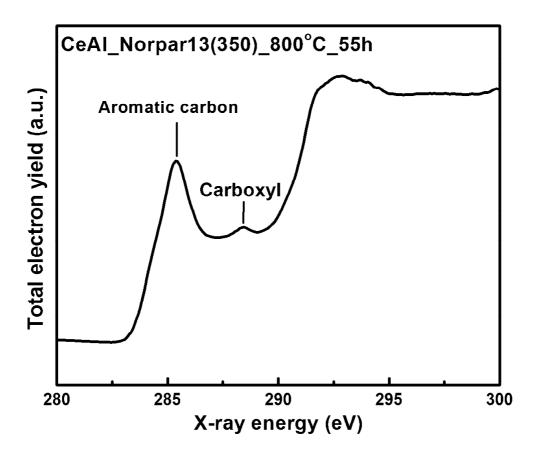


Figure S4. Carbon K-edge XANES spectrum of the used CeO₂-Al₂O₃ support after steam reforming of Norpar13 with 350 ppmw sulfur at 800 °C for 55 h. The reaction conditions are exactly the same as those used for testing of the Ni and Rh catalysts.

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Table S1 Contents of metal sulfide, organic sulfide, and sulfate on the used 10wt%Ni/CeO₂-Al₂O₃ catalysts after steam reforming of Norpar13 with 350 ppmw sulfur at 800 °C for different reaction time

Reaction time (h)	Metal sulfide (M-S)	Organic sulfide (C-S-C)	Sulfate (SO ₄ ²⁻)
	mg/g·cat	mg/g·cat	mg/g*cat
5	0.85	0	0.05
15	0.69	0.07	0.17
30	0.22	0.44	0.20
55	0.54	0.22	0.20

Table S2 Molar ratio of S: C: Me (Me = Rh or Ni) for the used Rh and Ni catalysts after steam reforming of Norpar13 with 350 ppmw sulfur at 800 °C for 55 h

	S	С	Rh	Ni
2wt%Rh/CeO ₂ -Al ₂ O ₃	0.45	107.3	1	
10wt%Ni/CeO ₂ -Al ₂ O ₃	0.23	814.9		1