

**Magnetic porous carbons derived from cobalt (II)-based metal-organic frameworks
for the solid-phase extraction of sulfonamides**

Sandra Yadira Mendiola Alvarez,^{1,2} Gemma Turnes Palomino,² Jorge Guzmán Mar,¹ Ma.
Aracely Hernández Ramírez,¹ Laura Hinojosa Reyes*,¹, Carlos Palomino Cabello*,²

¹ *Facultad de Ciencias Químicas, Universidad Autónoma de Nuevo León, UANL, Cd.
Universitaria, C.P. 66455 San Nicolás de los Garza, Nuevo León, México*

² *Department of Chemistry, University of the Balearic Islands, E-07122 Palma de Mallorca,
Spain*

**E-mail: carlos.palomino@uib.es *E-mail: laura.hinojosary@uanl.edu.mx*

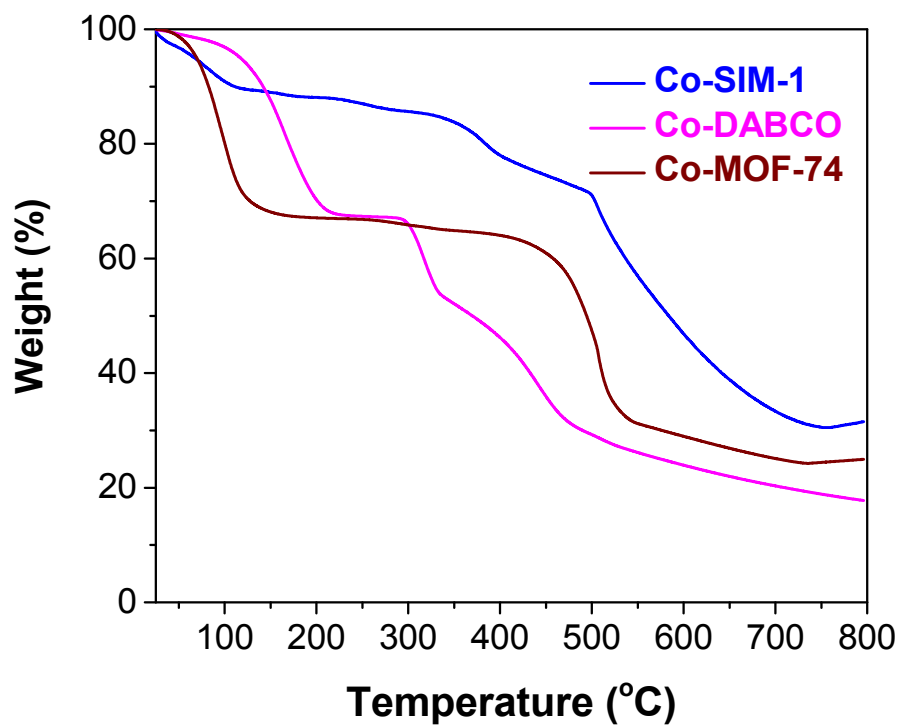


Fig. S1. TGA curves of Co-SIM-1, Co-MOF-74 and Co-DABCO.

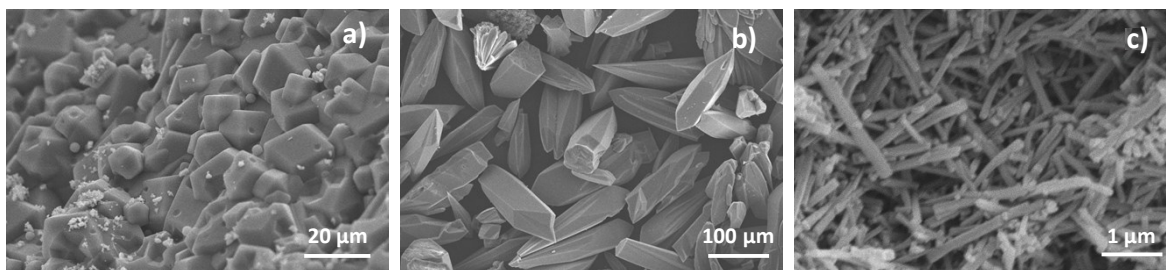


Fig. S2. SEM images of (a) Co-SIM-1, (b) Co-MOF-74 and (c) Co-DABCO metal-organic frameworks.

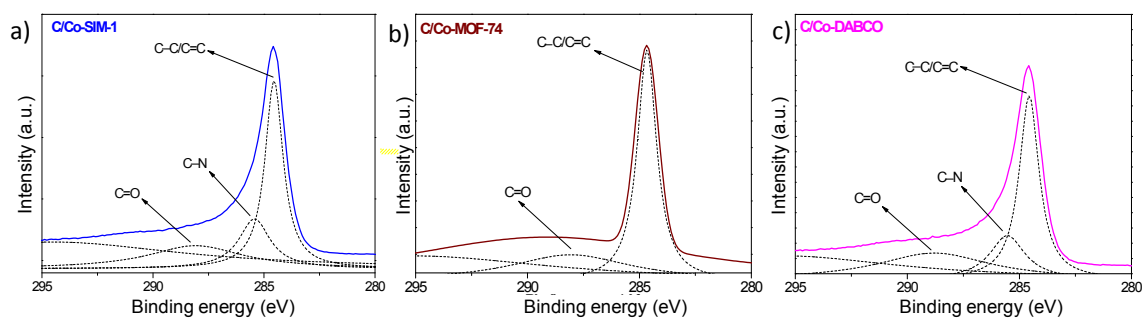


Fig. S3. High resolution C 1s XPS spectra of (a) C/Co-SIM-1, (b) C/Co-MOF-74 and (c) C/Co-DABCO samples.

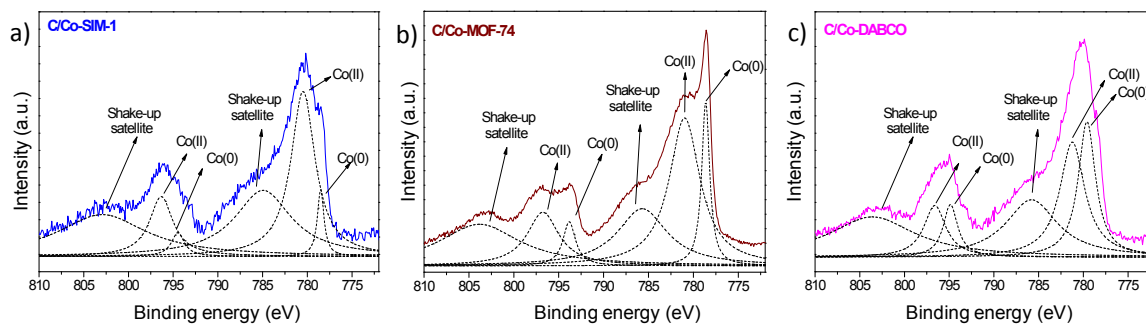


Fig. S4. High resolution Co 2p XPS spectra of (a) C/Co-SIM-1, (b) C/Co-MOF-74 and (c) C/Co-DABCO samples.

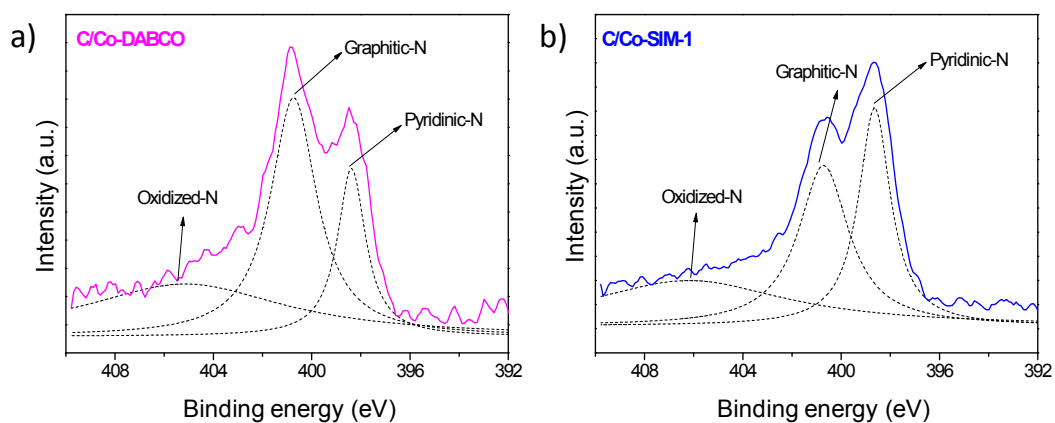


Fig. S5. High resolution N 1s XPS spectra of (a) C/Co-SIM-1 and (b) C/Co-DABCO samples.

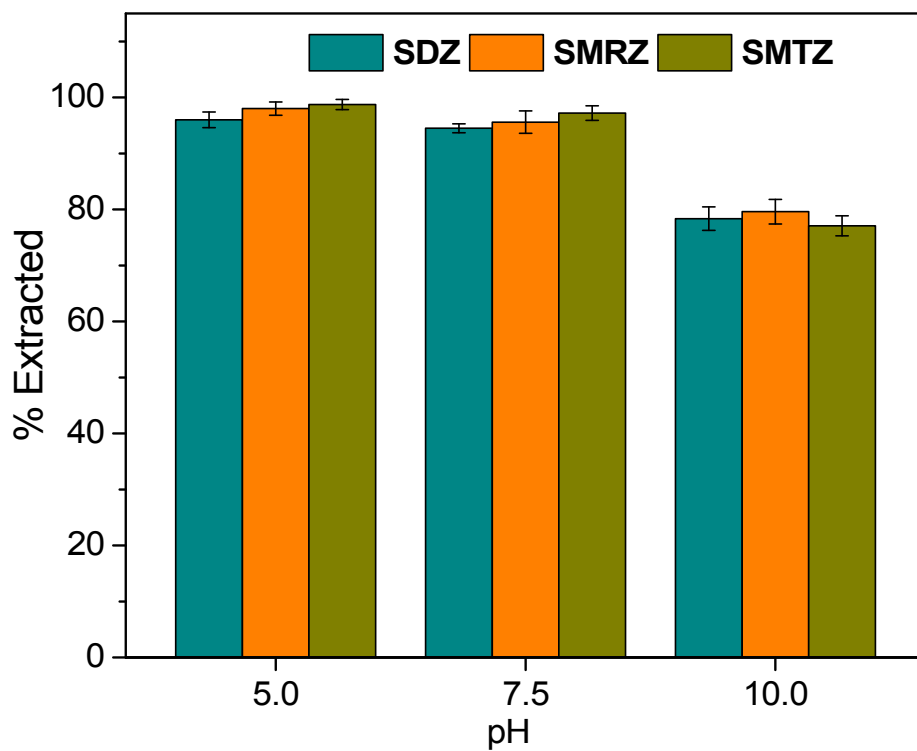


Fig. S6. Effect of pH on the adsorption of sulfonamides (5 ppm, each) on C/Co-SIM-1.

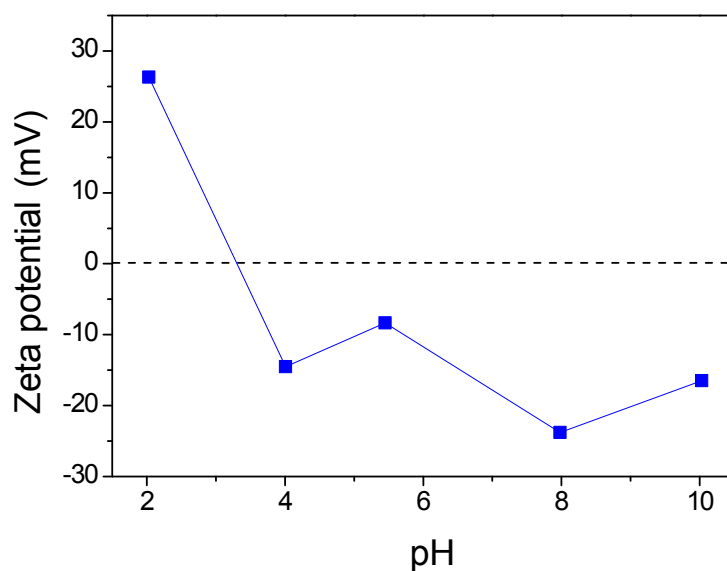


Fig. S7. Zeta potential values of C/Co-SIM-1 at different pH values.

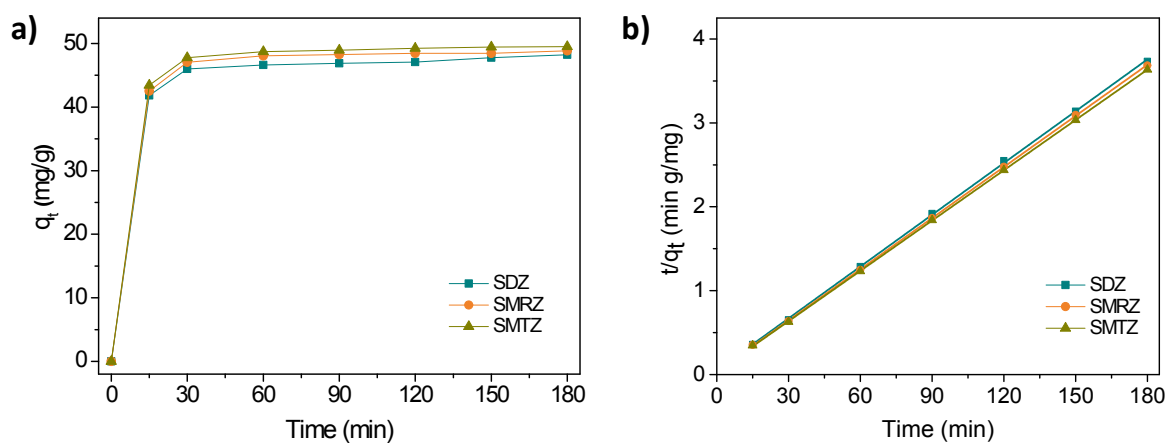


Fig. S8. (a) SNs adsorbed (mg/g) versus contact time (min) using C/Co-SIM-1 as adsorbent ($C_{\text{Sulfonamide}}$ 5 ppm, each). (b) Linear fit of pseudo-second order kinetics model for the adsorption of sulfonamides on C/Co-SIM-1.