Evaluation of a Porous Imine-Based Covalent Organic Framework for Solid-Phase Extraction of Nitroimidazoles

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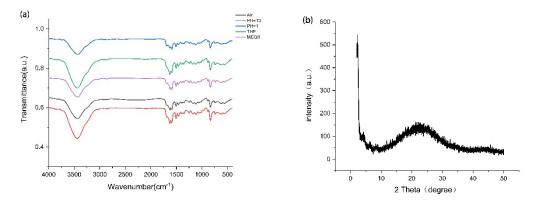


Fig.S1 (a) The FTIR spectrum of BP-COF in different conditions; (b) The PXRD pattern of BP-COF

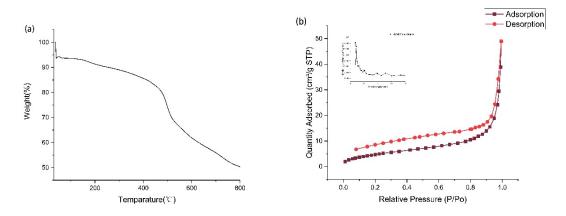


Fig.S2 (a)The TGA analysis of BP-COF; (b) N2 adsorption desorption isotherms of BP-COF.

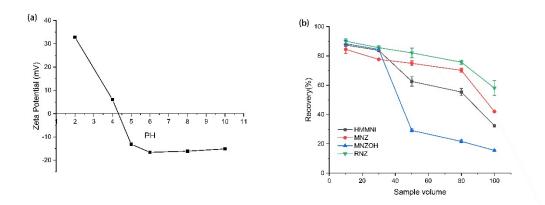


Fig.S3(a) Zeta potential of BP-COF; (b) Effect of the sample volume on the recoveries of NDZs (conditions: amount of sorbent, 60 mg; washing solvent, 1 mL water; elution solvent, 2 mL MeOH with 5% NH₃; sample pH,4; Spiked level, 10 ng ml⁻¹)

| the four 3-INDES. | | | | | |
|-------------------|------------------|--------------------|---------------------|--------------------------|----------|
| Analyte s | Parent ion (m/z) | Daughter ion (m/z) | Cone voltage (V) | Collision energy (eV) | RT (min) |
| HMM NI | 158.06 | 140.05*94.05 | 8 | 10-22 | 1.43 |
| MNZ | 172.07 | 128.05*82.05 | 18 | 12-22 | 1.5 |
| MNZO H | 188.07 | 126.03*123.06 | 16 | 10-14 | 1.12 |
| RNZ | 201.06 | 140.05*55.05 | 14 | 12-20 | 1.94 |

Table S1 Monitored LC-MS/MS transitions, sample cone voltages, collision energies and retention time of the four 5-NDZs.