## **Electronic Supplementary Information**

Synthesis of multi-ion imprinted polymers based on dithizone chelation for simultaneous removal of Hg<sup>2+</sup>, Cd<sup>2+</sup>, Ni<sup>2+</sup> and Cu<sup>2+</sup> from aqueous solutions

Junqing Fu,<sup>ab</sup> Xiaoyan Wang,<sup>bc</sup> Jinhua Li,\*<sup>b</sup> Yangjun Ding<sup>a</sup> and Lingxin Chen\*<sup>ab</sup>

<sup>a</sup> Key Laboratory of Life-Organic Analysis of Shandong Province, College of Chemistry and Chemical Engineering, Qufu Normal University, Qufu 273165, China

<sup>b</sup> Key Laboratory of Coastal Environmental Processes and Ecological Remediation, Yantai Institute of Coastal Zone Research, Chinese Academy of Sciences, Yantai 264003, China

<sup>c</sup> School of Pharmacy, Binzhou Medical University, Yantai 264003, China

\*Corresponding authors.

*E-mail addresses:* jhli@yic.ac.cn (J. Li), lxchen@yic.ac.cn (L. Chen)

## Contents

**Fig. S1.** FT-IR spectra of ions contained MIIPs, MIIPs and NIPs.  $M^{2+}$  is Hg<sup>2+</sup>, Cd<sup>2+</sup>, Ni<sup>2+</sup> and Cu<sup>2+</sup> here.

Fig. S2. (A)  $N_2$  adsorption-desorption isotherms and (B) pore size distribution of MIIPs and NIPs.

Fig. S3. TG/DTG curves of MIIPs.

**Fig. S4.** (A) Static adsorption isotherm curves of MIIPs and NIPs for 4 ions in aqueous solutions. (B) Langmuir isotherm models for ions onto the MIIPs. (C) Adsorption kinetics curves of MIIPs and NIPs for ions in aqueous solution. (D) Pseudo-first-order, pseudo-second-order, Elovich and intraparticle diffusion kinetic models for ions towards MIIPs.

Table S1. Specific surface area and other related data of MIIPs and NIPs.

Table S2. Isotherm model parameters for MIIPs and NIPs.

**Table S3.** Parameters obtained of ions adsorption towards MIIPs from four kinetic models.

**Table S4.** Comparative studies on MIIPs.



Fig. S1. FT-IR spectra of ions contained MIIPs, MIIPs and NIPs.  $M^{2+}$  is  $Hg^{2+}$ ,  $Cd^{2+}$ ,  $Ni^{2+}$  and  $Cu^{2+}$  here.



Fig. S2. (A)  $N_2$  adsorption-desorption isotherms and (B) pore size distribution of MIIPs and NIPs.



Fig. S3. TG/DTG curves of MIIPs.



**Fig. S4.** (A) Static adsorption isotherm curves of MIIPs and NIPs for 4 ions in aqueous solutions. Experimental conditions: V, 10 mL; polymer, 20 mg; adsorption time, 12 h; room temperature. (B) Langmuir isotherm models for ions onto the MIIPs. (C) Adsorption kinetics curves of MIIPs and NIPs for ions in aqueous solution. Experimental conditions: the sum concentration of Hg<sup>2+</sup>, Cd<sup>2+</sup>, Ni<sup>2+</sup> and Cu<sup>2+</sup>, 100 mg/L; V, 10 mL; polymer, 20 mg; room temperature. (D) Pseudo-first-order, pseudo-second-order, Elovich and intraparticle diffusion kinetic models for ions towards MIIPs.

Table S1. Specific surface area and	d other related data of MIIPs and NIPs.	

- parameters	specific surface area (m <sup>2</sup> /g)		cumulative	cumulative	average
	BET method	Langmuir method	pore area (m <sup>2</sup> /g)	pore volume (mL/g)	pore diameter (nm)
MIIPs	15.59	27.44	43.82	0.2404	21.94
NIPs	4.62	7.37	38.26	0.0955	9.98

Langmuir model			Freundlich model			
Sorbent	$Q_{\max}$ (mg g <sup>-1</sup> )	$K_1(\text{L mg}^{-1})$	R	$K_{\rm f} ({\rm mg^{1-(1/n)}}{\rm L^{1/n}}{\rm g^{-1}})$	1/ <i>n</i>	R
MIIP	18.178	0.006	0.990	0.006	4.066	0.988
NIP	3.516	0.011	0.998	0.007	3.301	0.998

Table S2. Isotherm model parameters for MIIPs and NIPs.

**Table S3.** Parameters obtained of ions adsorption towards MIIPs from four kinetic models.

Kinematic model	Model parameter		
Pseudo-first-order	$k_1^a$ (min <sup>-1</sup> )	$Q_{e}^{b}(\mu mol/g)$	r
$\ln(Q_{\rm e}-Q_{\rm t}) = \ln Q_{\rm e} - k_{\rm l}t$	0.066	16.345	0.990
Pseudo-second-order	$k_2^c(g/(\mu mol \cdot min))$	$Q_{\rm e}$ (µmol/g)	r
$\frac{t}{Q_{\rm t}} = \frac{1}{k_{\rm ad}Q_{\rm e}^2} + \frac{t}{Q_{\rm e}}$	0.01	17.271	0.996
Elovich	$\alpha^d (\mu mol/g)$	β <sup>d</sup> (min∙g/µmol)	r
$Q_{t} = \frac{1}{\beta} ln(\alpha\beta) + \frac{1}{\beta} ln(t)$	4.627	0.290	0.954
Intraparticle diffusion	$k_{\mathrm{p}}{}^{e}(\mu\mathrm{mol}/(\mathrm{gmin}^{0.5}))$	$C^{f}(\mu mol/g)$	r
$Q_{\rm t} = k_{\rm p} t^{1/2} + C$	2.183	0.644	0.658
	0.144	14.992	0.959

 $^{a} k_{1}$  is the rate constant of adsorption in pseudo-first-order model.

 ${}^{b} Q_{e}$  is the final adsorption amount at equilibrium,  $Q_{t}$  is the instantaneous adsorption amount in the adsorbent at time t.

 $^{c}$   $k_{2}$  is the rate constant of adsorption in pseudo-second-order model.

 $^{d} \alpha$  and  $\beta$  represent the initial adsorption rate and desorption constant in Elovich model.

<sup>*e*</sup>  $k_p$  indicates the intraparticle diffusion rate constant and relates to the particle size *R* with the  $k_p = 6Q_e \sqrt{\frac{D}{D}}$ 

equation  $k_p = \frac{6Q_e}{R} \sqrt{\frac{D}{\pi}}$ .

 ${}^{f}C$  provides information about the thickness of the boundary layer.

 Table S4. Comparative studies on MIIPs.

Template	Polymerization method	Monomer	Application	Ref.	
Ce(IV)-Gd(III)	Thermal polymerization	But-2-enedioic acid bis-[(2-	Screen-printed carbon	16	
		amino-ethyl)-amide	electrodes	10	
Cu(II)-Cd(II)	Precipitation polymerization	DEM <sup>a</sup>	Electrochemistry	17	
Cd(II)-Pb(II)	Precipitation polymerization	DEM	Solid-phase extraction	18	
As(V)-Cr(III)	Emulsion polymerization	$DDDPA^b$ & 4- $VP^c$		19	
Hg(II)-Cd(II)-Ni(II)-Cu(II)	Sol-gel process	Dithizone & APTES	Solid-phase extraction	This work	

<sup>*a*</sup> DEM: 2-(diethylamino) ethylmetacrilate.

<sup>b</sup> DDDPA: 1,12-dodecanediol-O,O'-diphenyl-phosphonic acid.

<sup>*c*</sup> 4-VP: 4-vinylpyridine.