

Photocatalytic Performance of Functionalized Biopolymer for Neodymium (III) Sorption and the Recovery from Leachate Solution

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Table S1. Assignment peaks of MCH, CH-POH, CH-POH+Nd and After 5 cycles of sorption desorption

38	Vibration	MCH	CH-POH	CH- POH+Nd	After 5 cycles	
39	O-H and N-H Str.	3412	3267	3274	3263	[7-1]
40	(overlapped)					
41	C-H aliph. (str.)	2928 ,2847	2917, 2848		2919, 2850	[12-8]
42	C=O (amide) and C=N-	1637	1691, 1621,	1569	1698	[13,14]
43	(overlapped) (str.)		1552			
44	1°/2° amines (bend)	1542			1575	[15,16]
45	C-H asym. (bend)	1380	1431	1396	1415	[17,18]
46	P=O (asym.)		1348		1370	[19,20]
47	O-H (in-plane) bend	1161	1261, 1236	1147	1311	[21,22]
48	P(O) (phosphate(str.))		1053	1126	1137	[19]
49	C-O, C-O-C and C-N (str.)	1019	1006	1024	1024	[23]
	P-O-C (str.)		908	946, 890	892	[19]
	-(CH ₂) _n - rocking	882	827	811	795	[24]
	P-O-C (str.)		609		596	[19,20]
	Fe-O	570	561	617	549	[25]

Table S2. Elemental analysis of MCH and CH-POH sorbents

	C	N	H	O	Fe	P
MCH [%]	27.35	4.12	5.06	30.22	33.25	0
MCH [m mol]	22.77	2.94	50.2	18.89	5.95	0
MCH-NPOH [%]	22.96	5.58	4.98	36.92	26.61	2.95
M MOL [m mol]	19.12	3.98	49.4	23.08	4.76	0.95

Table S3a. Reminder on equations used for modeling sorption isotherms [26-28].

Model	Equation	Parameters	Ref.
Langmuir	$q_{eq} = \frac{q_{m,L}C_{eq}}{1 + b_L C_{eq}}$	$q_{m,L}$ (mmol g ⁻¹): Sorption capacity at saturation of monolayer b_L (L mmol ⁻¹): Affinity coefficient	[26]
Freundlich	$q_{eq} = k_F C_{eq}^{1/n_F}$	k_F and n_F : empirical parameters of Freundlich equation	[26]
Sips	$q_{eq} = \frac{q_{m,S} b_S C_{eq}^{1/n_S}}{1 + b_S C_{eq}^{1/n_S}}$	$q_{m,L}$, b_S and n_S : empirical parameters of Sips equation (based on Langmuir and Freundlich equations)	[26]

Table S3b. Reminder on equations used for modeling uptake kinetics [26,29] [30].

Model	Equation	Parameters	Ref.
PFORE	$q(t) = q_{eq,1}(1 - e^{-k_1 t})$	$q_{eq,1}$ (mmol g ⁻¹): sorption capacity at equilibrium k_1 (min ⁻¹): apparent rate constant of PFORE	[30]
PSORE	$q(t) = \frac{q_{eq,2}^2 k_2 t}{1 + k_2 q_{eq,2} t}$	$q_{eq,2}$ (mmol g ⁻¹): sorption capacity at equilibrium k_2 (g mmol ⁻¹ min ⁻¹): apparent rate constant of PSORE	[30]
RIDE	$\frac{q(t)}{q_{eq}} = 1 - \sum_{n=1}^{\infty} \frac{6\alpha(\alpha + 1)\exp\left(\frac{-D_e q_n^2}{r^2} t\right)}{9 + 9\alpha + q_n^2 \alpha^2}$ With q_n being the non-zero roots of $\tan q_n = \frac{3 q_n}{3 + \alpha q_n^2}$ and $\frac{m q}{V C_0} = \frac{1}{1 + \alpha}$	D_e (m ² min ⁻¹): Effective diffusivity coefficient	[29]

(m (g): mass of sorbent; V (L): volume of solution; C₀ (mmol L⁻¹): initial concentration of the solution).

Akaike Information Criterion, AIC:

$$AIC = N \ln \left(\frac{\sum_{i=0}^N (y_{i,exp.} - y_{i,model})^2}{N} \right) + 2N_p + \frac{2N_p(N_p + 1)}{N - N_p - 1}$$

Where N is the number of experimental points, N_p the number of model parameters, y_{i,exp.} and y_{i,model} the experimental and calculated values of the tested variable.

Table S4. XRF analysis of the study ore materials.

Major oxides (%)	Wt. (%)	Trace elements	ppm
SiO ₂	76.51	U	1117
Al ₂ O ₃	8.56	REE	1296
TiO ₂	0.95	Mn	35.4
Fe ₂ O ₃ ^{total}	3.26	Zn	86.4
CaO	1.19		
MgO	0.45		
Na ₂ O	0.43		
K ₂ O	1.18		
P ₂ O ₅	0.49		
L.O.I	5.68		
Total	98.7		

Table S5. Chemical composition of the prepared carbonate leach liquor at (pH=0.3)

Constituent	Conc. (mg/L)	Constituent	Conc. (mg/L)
Fe	2690	Ca	448
Al ₂ O ₃	2538	U	388
Mg	350	REE	451

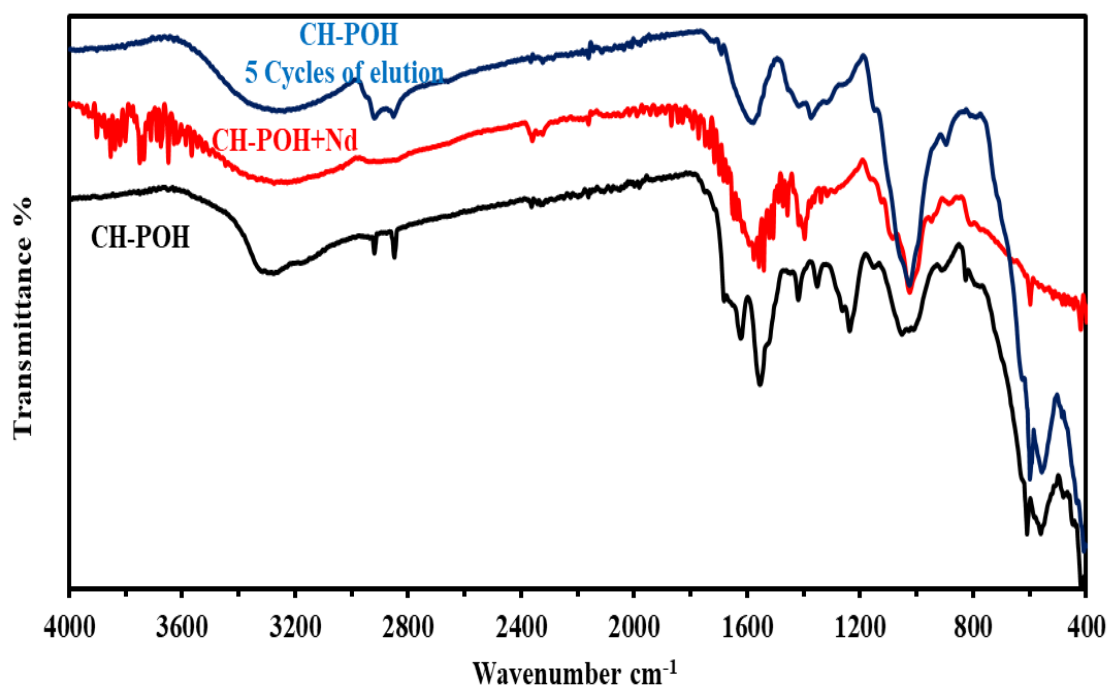


Figure S1. The full range of the FTIR spectra.

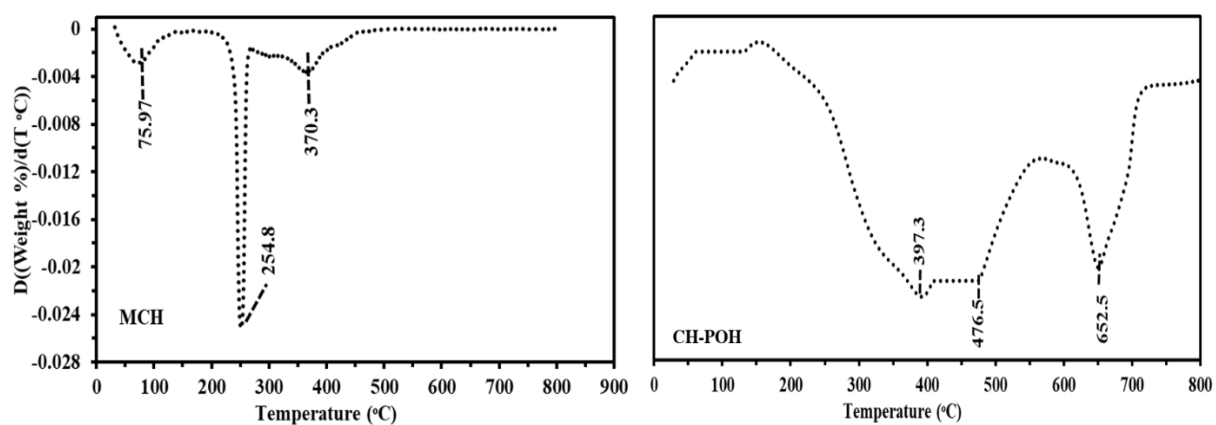


Figure S2. DrTG of chitosan magnetite (MCH) and functionalized sorbent (CH-POH)

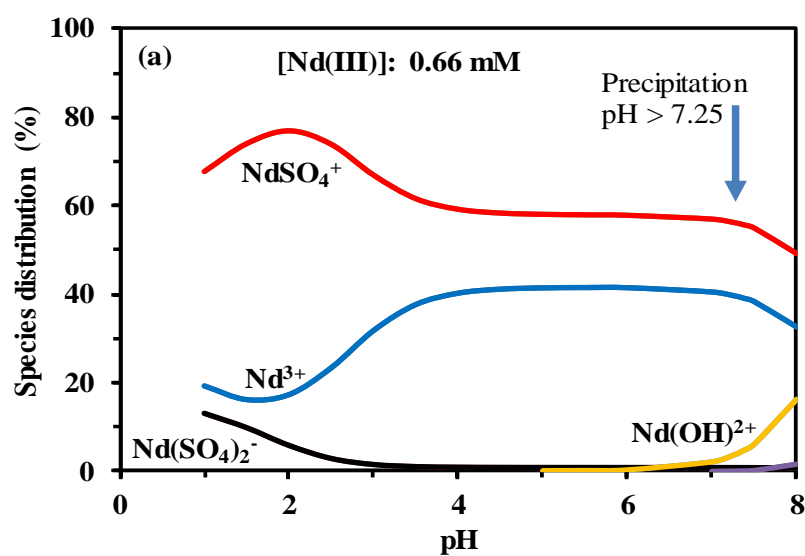


Figure S3. Speciation diagrams for Nd(III) under the experimental conditions

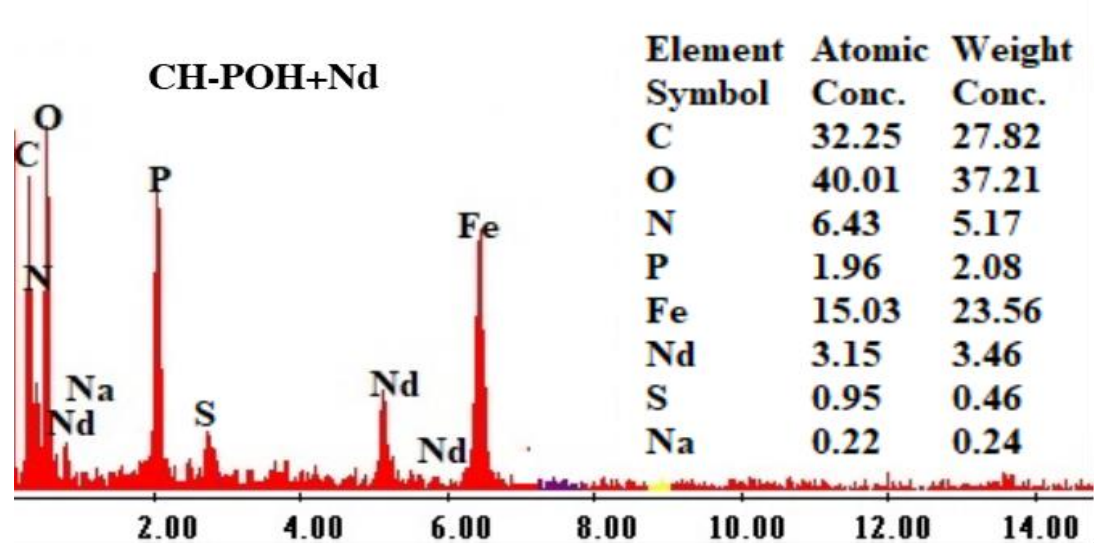


Figure S4. EDX analysis of the CH-POH after Nd(III) sorption

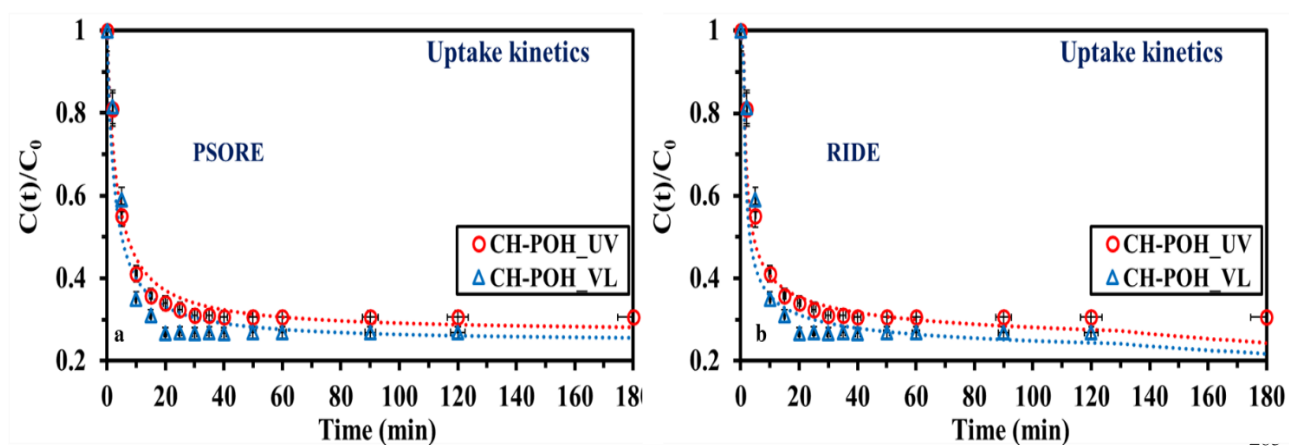


Figure S5. The unfitted profiles of the PSORE and RIDE for CH-POH sorbent, (pH_0 : 4; C_0 : 0.36 mmol Cd L⁻¹; SD: 0.66 g L⁻¹; T: 21 ± 1 °C; v: 210 rpm).

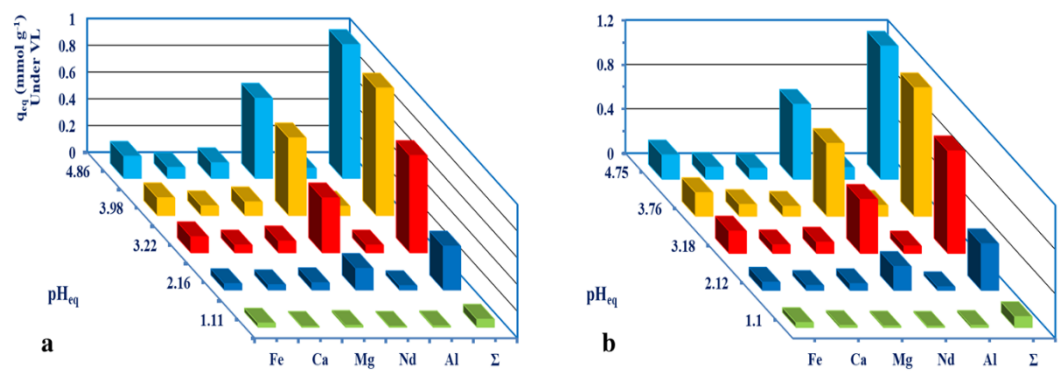


Figure S6. Total sorption capacity of CH-POH in polymetallic equimolar solution under VL(a) and (UV) conditions

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