

Table S3. Fitting model parameters and the calculated parameters for Co released from the citrate-coated CoFe_2O_4 NPs in soil solution at different pH levels according to different dissolution kinetic models.

Fitting parameters	pH 5	pH 7	pH 8
Zero-order			
R^2	0.638	0.358	0.089
R^2 adjusted	0.625	0.335	0.056
Intercept	0.010*	0.004*	0.008*
Slope	3.307×10^{-5} *	2.038×10^{-5} *	-1.519×10^{-5}
k (mM h^{-1})	-3.307×10^{-5}	-2.038×10^{-5}	1.519×10^{-5}
M_s (mM)	0.010	0.004	0.008
First-order			
R^2	0.570	0.268	0.110
R^2 adjusted	0.554	0.242	0.079
Intercept	-4.637*	-5.584*	-4.883*
Slope	0.003*	0.004*	-0.002
k (mM h^{-1})	-2.886×10^{-3}	-3.761×10^{-3}	2.638×10^{-3}
M_s (mM)	0.010	0.004	0.008
Second-order			
R^2	0.491	0.172	0.122
R^2 adjusted	0.472	0.143	0.090
Intercept	104.136*	277.521*	140.148*
Slope	-0.258*	-0.802*	0.517
k ($\text{mM}^{-1} \text{h}^{-1}$)	-2.583×10^{-1}	-8.019×10^{-1}	5.169×10^{-1}
M_s (mM)	0.010	0.004	0.007
Pseudo-first-order			
R^2	0.048	0.176	0.195
R^2 adjusted	0.014	0.146	0.167
Intercept	-1.168*	-1.145*	-1.128*
Slope	-3.559×10^{-5}	1.827×10^{-4} *	1.115×10^{-4} *
k (h^{-1})	3.559×10^{-5}	-1.827×10^{-4}	-1.115×10^{-4}
M_s (mM)	0.311	0.318	0.324
Pseudo-second-order			
R^2	0.990	0.805	0.865
R^2 adjusted	0.989	0.798	0.861
Intercept	225.889*	1816.980*	-2179.680
Slope	69.793*	138.250*	260.020*
k ($\text{L mol}^{-1} \text{h}$)	21.564	10.519	-31.327
M_s (mM)	0.014	0.007	0.004
One-half-order			
R^2	0.605	0.316	0.101
R^2 adjusted	0.591	0.291	0.069
Intercept	0.099*	0.061*	0.088*
Slope	1.540×10^{-4}	1.363×10^{-4} *	-9.859×10^{-5}
k ($\text{mM}^{1/2} \text{h}$)	3.080×10^{-4}	-2.726×10^{-4}	-1.972×10^{-4}
M_s (mM)	0.010	0.004	0.008
Three-half-order			
R^2	0.531	0.219	0.118
R^2 adjusted	0.515	0.191	0.085
Intercept	10.183*	16.471*	11.655*
Slope	-0.014*	-0.027*	0.018
k ($\text{mM}^{-1/2} \text{h}$)	-2.721×10^{-2}	-5.380×10^{-2}	3.639×10^{-2}
M_s (mM)	0.010	0.004	0.006
Evolich			
R^2	0.641	0.188	0.043
R^2 adjusted	0.628	0.159	0.009
Intercept	0.009*	0.004*	0.007*
Slope	7.878×10^{-4} *	3.507×10^{-4} *	2.524×10^{-4}
α (mM h^{-1})	104.192	20.891	2.025×10^8

β (L mmol ⁻¹)	1269.296	2851.499	3961.965
Higuchi			
R ²	0.746	0.255	0.009
R ² adjusted	0.737	0.228	-0.027
Intercept	0.009*	0.004*	0.008*
Slope	4.652x10 ⁻⁴ *	2.235x10 ⁻⁴ *	-6.223x10 ⁻⁵
k (mM h ^{1/2})	4.652x 10 ⁻⁴	2.235 x 10 ⁻⁴	-6.223 x 10 ⁻⁵
Ms (mM)	0.009	0.004	0.008
Hixon-Crowell			
R ²	0.048	0.185	0.198
R ² adjusted	0.014	0.156	0.170
Intercept	0.104*	0.106*	0.108*
Slope	-3.703x10 ⁻⁶	1.981x10 ⁻⁵ *	1.216x10 ⁻⁵ *
k (mM ^{1/3} h ⁻¹)	-3.703 x 10 ⁻⁶	1.981 x 10 ⁻⁵	-1.981 x 10 ⁻⁵
Korsmeyer-Peppas			
n	0.070	0.075	0.113
R ²	0.664	0.175	0.019
R ² adjusted	0.652	0.146	-0.016
Intercept	-0.002*	4.032x10 ⁻⁴	0.018*
Slope	0.0306*	0.012*	3.561x10 ⁻³
k (h ⁻ⁿ)	0.0306	0.012	0.004
Baker-Lonsdale			
R ²	0.626	0.325	0.093
R ² adjusted	0.613	0.300	0.060
Intercept	0.003*	0.012*	0.023*
Slope	1.001x10 ⁻⁴ *	5.738x10 ⁻⁵ *	-4.665x10 ⁻⁵
k (h)	1.001x10 ⁻⁴	5.738x10 ⁻⁵	-4.665x10 ⁻⁵
Weibull			
R ²	0.625	0.213	0.038
R ² adjusted	0.611	0.185	0.003
Intercept	-2.034*	-2.439*	-2.192*
Slope	0.071*	0.080*	0.037
a	9.247 x 10 ⁻³	3.639 x 10 ⁻³	6.427x 10 ⁻³
b	0.071	0.080	0.037

[M]_s; theoretical saturation concentration, *Symbol indicates significance with an $\alpha=0.05$