

Effective Isolation of Succinic Acid from Aqueous Media with the Use of Anion Exchange Resins

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SUPPORTING INFORMATION

Table S1. The general properties of anion exchange resins used in the study.

General properties	Lewatit MP-62	Lewatit MP-64	Lewatit M-500	Lewatit M-600
Type	weak basic	weak basic	strong basic	strong basic
Ionic form	free base	free base	Cl ⁻	Cl ⁻
Functional group	tertiary amine	tertiary & quaternary amine	quaternary amine (Type I)	quaternary amine (Type II)
Matrix structure	cross-linked polystyrene	cross-linked polystyrene	cross-linked polystyrene	cross-linked polystyrene
Appearance	beige, opaque	beige, opaque	yellow, translucent	white, translucent
Maximum temperature (°C)	130	70	70	30
Operating pH range	0-8	0-7	0-12	0-11
Density (g/mL)	1.02	1.04	1.08	1.10
Total capacity (eq/L)	1.7	1.3	1.3	1.3

Table S2. Non-linear and linear equations of the kinetic models studied.

Kinetic model	Non-linear equation	Linear equation	Plot
Pseudo-first-order	$dq_t/dt = k_1 \cdot (q_e - q_t)$	$\log(q_e - q_t) = \log q_e - k_1 t$	$\log(q_e - q_t) vs t$
Pseudo-second-order	$dq_t/dt = k_2 \cdot (q_e - q_t)$	$\frac{t}{q_t} = \frac{1}{k_2 q_e^2} + \frac{1}{q_e} \cdot t$	$t/q_t vs t$
Elovich	$dq_t/dt = \alpha \cdot exp^{[\frac{\beta}{\alpha}]}$	$q_t = \frac{\ln(\alpha \cdot \beta)}{\beta} + \frac{\ln t}{\beta}$	$q_t vs \ln t$
Intraparticle diffusion	---	$q_t = k_{id} \cdot t^{0.5} + I$	$q_t vs t^{0.5}$

Table S3. The linear equations of the isotherm models studied.

Model name	Linear equation	Plot
Langmuir	$\frac{1}{q_e} = \frac{1}{q_{\max} K_L} \cdot \frac{1}{C_e} + \frac{1}{q_{\max}}$ $R_L = 1/(1 + (K_L \cdot C_o))$	$1/q_e$ vs $1/C_e$
Freundlich	$\log q_e = \frac{1}{n} \cdot \log C_e + \log K$	$\log q_e$ vs $\log C_e$
Temkin	$q_e = B \cdot \ln K_T + B \cdot \ln C$	q_e vs $\ln C_e$
Scatchard	$\frac{q_e}{C_e} = q_{\max} \cdot K_S - K_S \cdot q_e$	q_e/C_e vs q_e

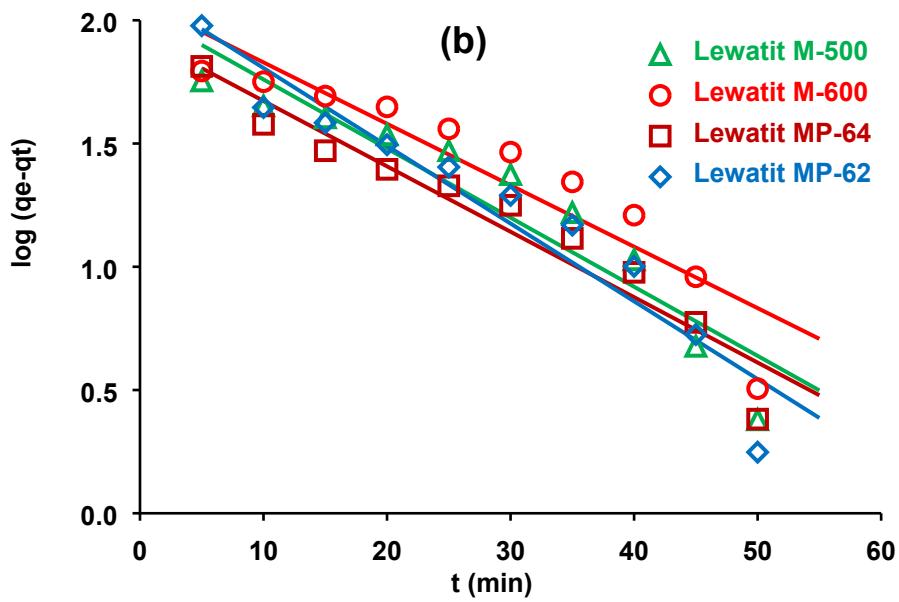
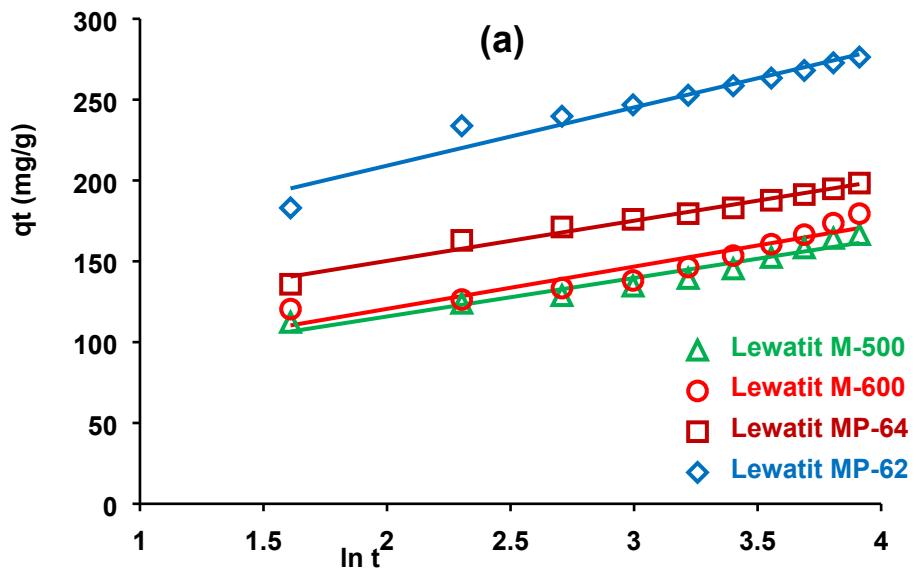


Fig. S1 The plots of the kinetic analysis for the isolation of SA using anion exchange resins
a) Elovich kinetic model b) Pseudo-first order kinetic model