

Correction to Efficient Resolution of Enantiomers by Coupling Preferential Crystallization and Dissolution. Part 2. A Parametric Simulation Study To Identify Suitable Process Conditions

Matthias J. Eicke,* Guillaume Levilain, and Andreas Seidel-Morgenstern

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In the original article, the authors unfortunately did not include the considered batch idle time, t_{idle} , in eq 22. The correct equation for the productivity is thus as follows:

$$\text{Pr}_k^j(t) = \frac{m_{kS}^j(t) - m_{k,\text{seeds}}^j}{V_{\text{cryst}}(t + t_{\text{idle}})} ee_S^j, m_{k,\text{seeds}}^j = 0 \text{ for } j = \text{T2},$$
$$k = \begin{cases} \text{E1(L-Thr)} & \text{for } j = \text{T1} \\ \text{E2(D-Thr)} & \text{for } j = \text{T2} \end{cases} \quad (22)$$

The time t_{idle} , needed to prepare a subsequent batch, was set to be 1 h. Higher productivities are possible, if this idle time can be reduced.

The mistake has been kindly pointed out by a reader, and the authors apologize for any inconvenience this might have caused.