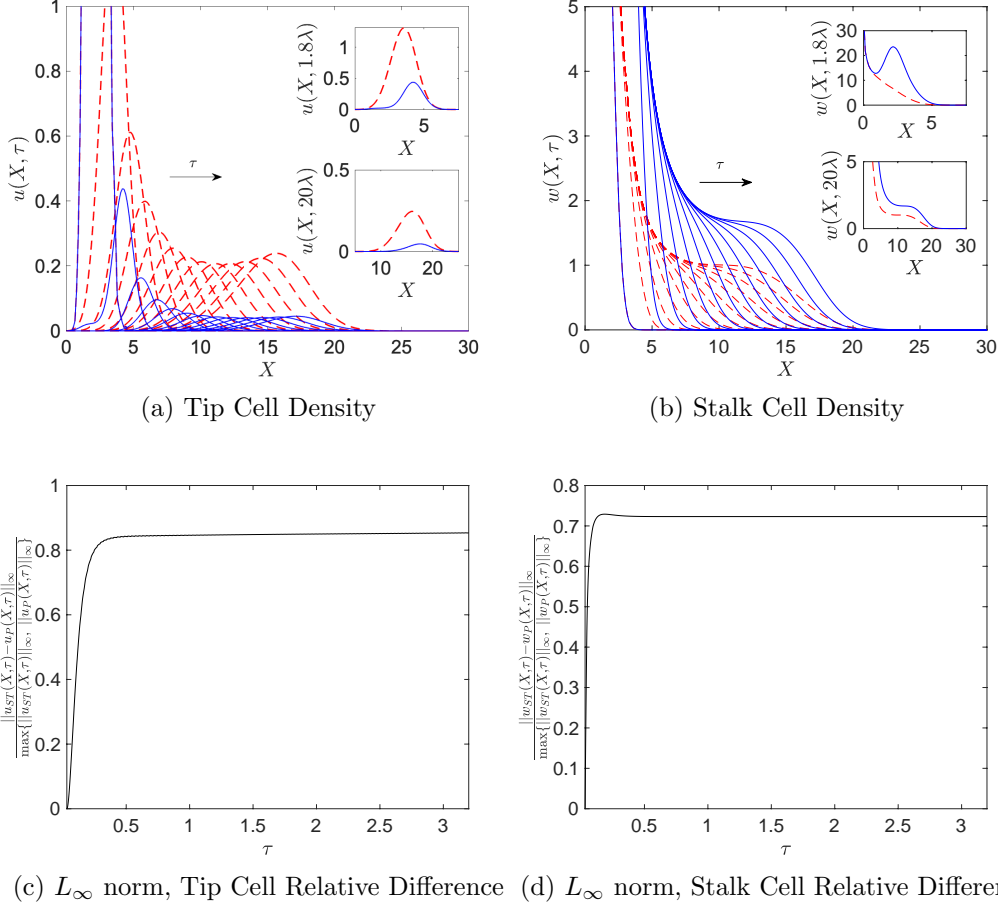


Figure 1: Numerical solutions of the ST-PDE and P-PDE systems showing how (a) the tip cell density, $u(X, \tau)$, and (b) the stalk cell density, $w(X, \tau)$, evolve over times $t = 0.2\lambda, 2.4\lambda, \dots, 2.5$, with $\lambda = 0.16$. The value of D was increased from 10^{-3} to 10^{-1} , while all other values and initial conditions are the same as those used to generate Figure 3 of the main text (this corresponds to an increase in ϵ from $10^{-3/2}$ to about 0.32). The insets in the top two panels show a zoomed-in view of the results for the times $\tau = 0.2\lambda, 2.4\lambda, 4.6\lambda$, and 6.8λ . The bottom row of figures shows the maximum relative difference between (c) tip cell densities of the ST-PDE, $u_{ST}(X, \tau)$, and P-PDE, $u_P(X, \tau)$, and (d) stalk cell densities of the ST-PDE, $w_{ST}(X, \tau)$, and P-PDE, $w_P(X, \tau)$. Key: ST-PDE (dashed red lines), P-PDE (solid blue lines). The PDEs were simulated on $\tau \in [0.2\lambda, 2.5]$, $X \in [0, 20\sqrt{\frac{\lambda}{D}}]$.



(c) L_∞ norm, Tip Cell Relative Difference (d) L_∞ norm, Stalk Cell Relative Difference

Figure 2: Numerical solutions of the ST-PDE and P-PDE systems showing how (a) the tip cell density, $u(X, \tau)$, and (b) the stalk cell density, $w(X, \tau)$, evolve over times $t = 0.2\lambda, 1.8\lambda, \dots, 19.4\lambda$, with $\lambda = 0.16$. The parameter χ has been decreased from 0.4 to 0.04, while all other initial conditions and parameter values are the same as those used in Figure 3 of the main text (this corresponds to an increase in ϵ from $10^{-3/2}$ to about 0.32). The top insets in (a) and (b) show a zoomed-in view of the results at $\tau = 1.8\lambda$, while the bottom inset in those panels is a zoomed-in view of the results at $\tau = 20\lambda$. The bottom row of figures shows the maximum relative difference between (c) tip cell densities of the ST-PDE, $u_{ST}(X, \tau)$, and P-PDE, $u_P(X, \tau)$, and (d) stalk cell densities of the ST-PDE, $w_{ST}(X, \tau)$, and P-PDE, $w_P(X, \tau)$. Key: ST-PDE (dashed red lines), P-PDE (solid blue lines). The PDEs were simulated on $\tau \in [0.2\lambda, 20\lambda]$, $X \in [0, 20\sqrt{\frac{\lambda}{D}}]$.

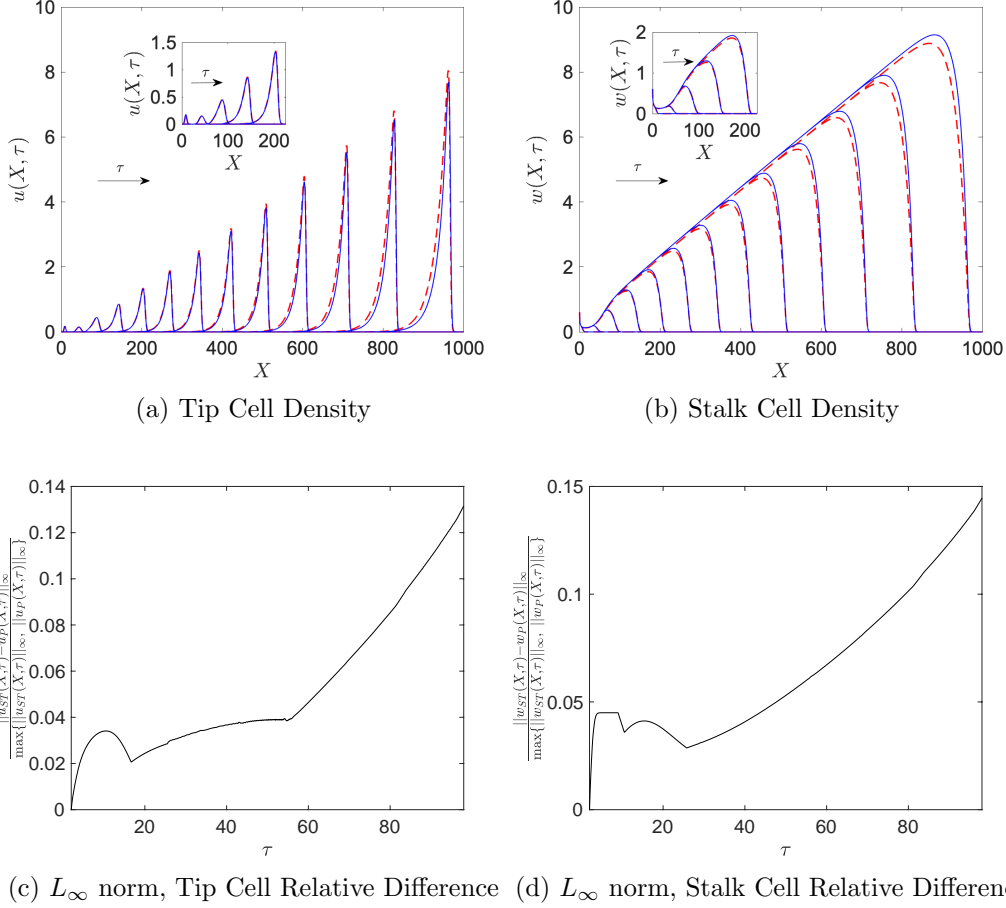
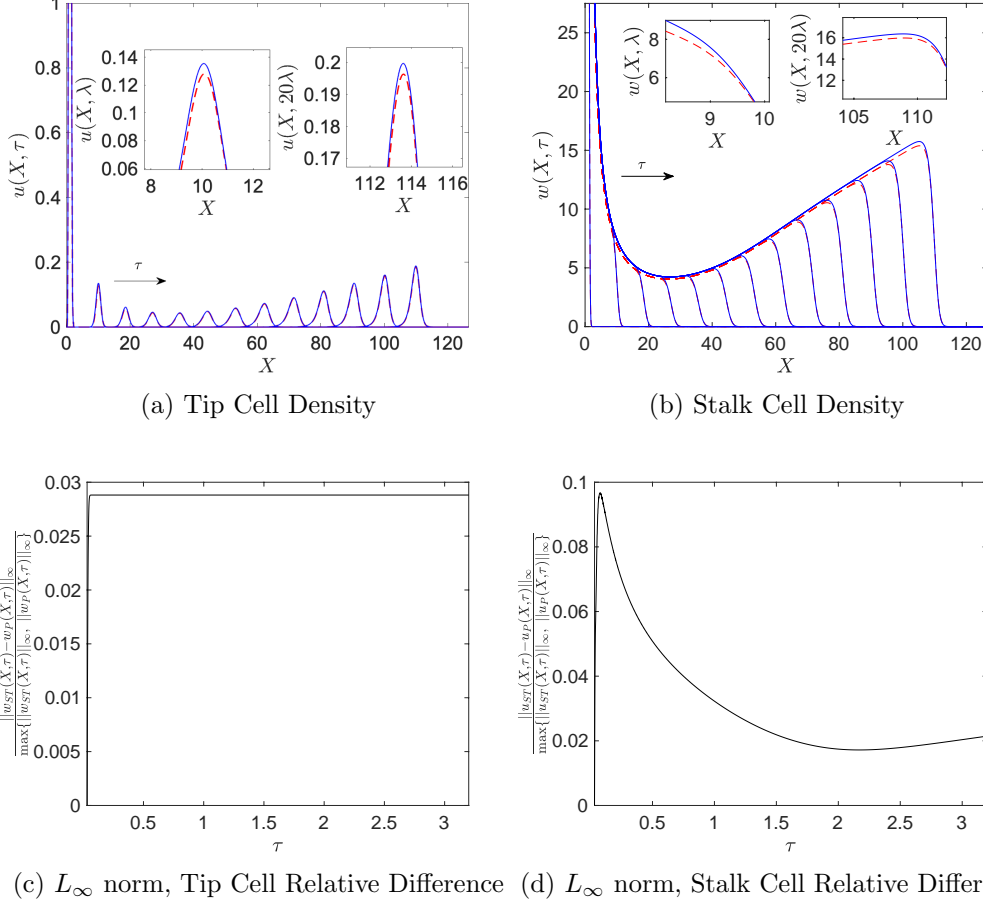


Figure 3: Numerical solutions of the ST-PDE and P-PDE systems showing how (a) the tip cell density, $u(X, \tau)$, and (b) the stalk cell density, $w(X, \tau)$, evolve over times $\tau = 0.2\lambda, 1.8\lambda, \dots, 9.8\lambda$, when λ is increased from 0.16 to 10 (this corresponds to increases in ϵ from $10^{-3/2}$ to about 0.32 and α from 10^{-3} to 0.0625). Other parameter values: as in Figure 3 of the main text. The insets in (a) and (b) show a zoomed-in view of the results at times $\tau = 0.2\lambda, 1.8\lambda, 2.4\lambda, 4\lambda$, and 5.6λ . The bottom row shows the maximum relative difference between (c) tip cell densities of the ST-PDE, $u_{ST}(X, \tau)$, and P-PDE, $u_P(X, \tau)$, and (d) stalk cell densities of the ST-PDE, $w_{ST}(X, \tau)$, and P-PDE, $w_P(X, \tau)$. Key: ST-PDE (dashed red lines), P-PDE (solid blue lines). The PDEs were simulated on $\tau \in [0.2\lambda, 9.8\lambda]$, $X \in [0, 1000]$.



(c) L_∞ norm, Tip Cell Relative Difference (d) L_∞ norm, Stalk Cell Relative Difference

Figure 4: Numerical solutions of the ST-PDE and P-PDE systems showing how (a) the tip cell density, $u(X, \tau)$, and (b) the stalk cell density, $w(X, \tau)$, evolve over times $\tau = 0.2\lambda, 1.8\lambda, \dots, 19.4\lambda$, with $\lambda = 0.16$. The parameter a_e was increased from 0.0391 to 1, while keeping all other initial conditions and parameter values the same as those used to generate Figure 3 of the main text (this corresponds to an increase in β from 39.1 to 1000, while $\epsilon = 10^{-3/2}$ and $\alpha = 10^{-3}$ remain constant). The left inset in (a) and (b) shows a zoomed-in view of the solution at $\tau = \lambda$, while the right inset in those panels shows a zoomed-in view of the solution at $\tau = 20\lambda$. The bottom row of figures shows the maximum relative difference between (c) tip cell densities of the ST-PDE, $u_{ST}(X, \tau)$, and P-PDE, $u_P(X, \tau)$, and (d) stalk cell densities of the ST-PDE, $w_{ST}(X, \tau)$, and P-PDE, $w_P(X, \tau)$. Key: ST-PDE (dashed red lines), P-PDE (solid blue lines). The PDEs were simulated on $\tau \in [0.2\lambda, 20\lambda]$, $X \in [0, 10\sqrt{\frac{\lambda}{D}}]$.

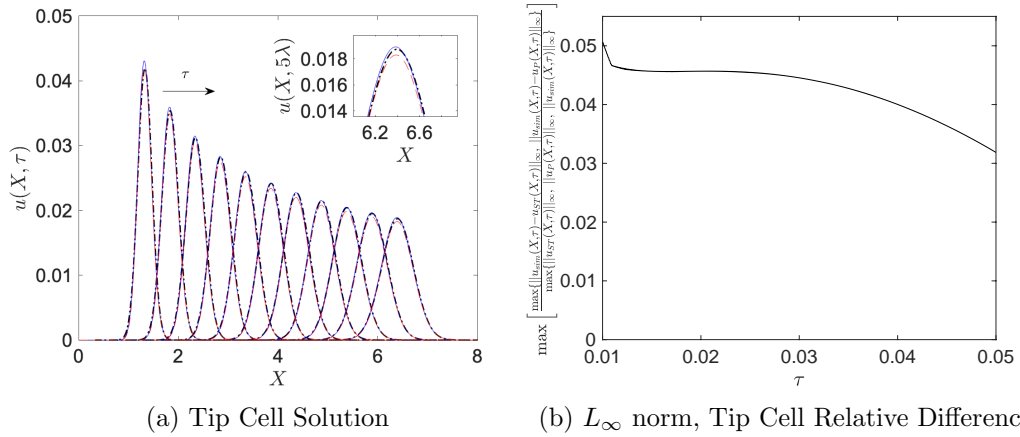


Figure 5: Self-similar solution given by equation (20) of the main text in (a) the independent variable X for $\tau = 0.2\lambda, 0.6\lambda, \dots, 5\lambda$, where $\lambda = 0.01$. In order to better compare the self-similar solution to the ST-PDE and P-PDE results, we set the maximum value of the self-similar solution at $\tau = 0.2\lambda$ equal to that of the ST-PDE and P-PDE tip cell densities at $\tau = 0.2\lambda$. The inset in (a) shows a zoomed-in view of the results at $\tau = 5\lambda$. In (b), we have plotted the relative error between the similarity solution and the P-PDE solution. Key for (a): P-PDE solutions (solid blue lines), ST-PDE solutions (dashed red lines), solution to equation (19) of the main text (dashed-dot black line).

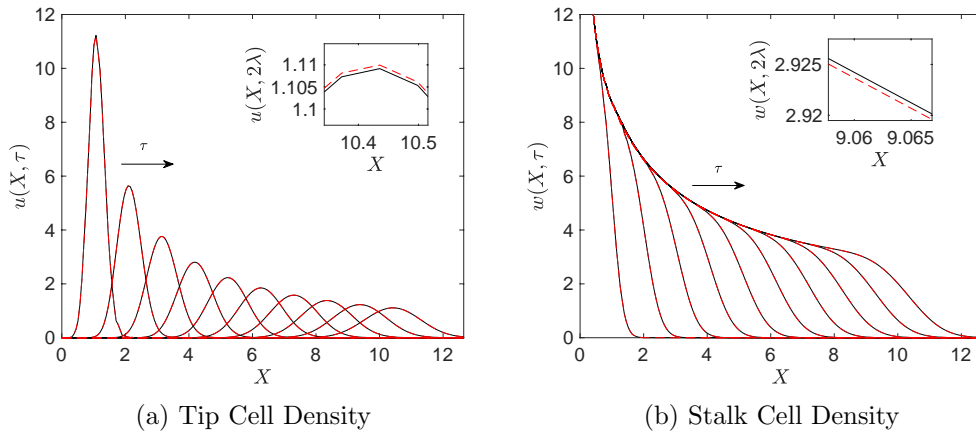


Figure 6: Numerical solution of the leading order dynamics showing how (a) the tip cell density, $u(X, \tau)$, and (b) the stalk cell density, $w(X, \tau)$, evolve over times $\tau = 0.2\lambda, 0.4\lambda, \dots, 2\lambda$, with $\lambda = 0.16$, along with results from the original P-PDE (the independent variables of the inner solution have been transformed back into the unscaled variables X and τ for ease of comparison). The inset in both panels shows a zoomed-in view of the results at time $\tau = 2\lambda$. Key: P-PDE solution (solid black lines), leading order solutions (dashed red lines). Parameter values: $D = 10^{-3}$, $\chi = 0.4$, $a_n = 1$, $a_e = 0.0391$, $\mu = 160$ (this corresponds to $\epsilon = 10^{-3/2}$, $\alpha = 10^{-3}$, and $\beta = 39.1$). The PDEs were simulated on $\tau \in [0.2\lambda, 2\lambda]$, $X \in [0, \sqrt{\frac{\lambda}{D}}]$. P-ABM solutions at $\tau = 0.2\lambda$, column averaged in the y -direction, were used as initial conditions.