## **Text S4** Reaction set for stochastic simulations

We developed a set of reactions that, using the law of mass action and the quasi-steady-state approximation, would exactly match our set of PDE equations. The complete reaction set can be found in Table S11 below. For a Hill coefficient of exactly two, we assumed that each promoter had a single binding site to which only the dimerized form of the appropriate activator or inhibitor could bind. The table of reactions follows the species and variable naming scheme used in the PDE set found in (6)-(16) with the addition of dimers  $p_{x2}$  and promoters  $Pr_x$ . The dissociation constants  $K_x$  are combinations of the on and off dimerization and binding constants. In general,  $K_x = \sqrt{\frac{k_{\text{doff}_x} k_{\text{off}_x}}{k_{\text{don}_x} k_{\text{on}_x}}}$ .

**Table S11:** Full reaction set for stochastic simulations. Kinetic rate constants chosen to validate quasi-steady-state approximation and to match Parameter Set 2 of Table S1.

Reaction Type	Reactions
Dimerization	$p_x + p_x \xrightarrow[k_{\text{doff}_x}]{k_{\text{doff}_x}} p_{x2} \text{ for } x \in \{C, L, RA, T\}$
Transcription factor-promoter binding	$Pr_x + p_{x2} = \frac{k_{\text{On}_x}}{k_{\text{Off}_x}} Pr_x p_{x2} \text{ for } x \in \{C, L, RA\}$
	$Pr_{x} + p_{T2} \xrightarrow{k_{\text{on}_{T}}} Pr_{x}p_{T2} \text{ for } x \in \{TO, TQ\}$ $Pr_{C} \xrightarrow{V_{PR}} Pr_{C} + m_{L}$
Maximal transcription	$Pr_C \xrightarrow{V_{P_R}} Pr_C + m_L$
	$Pr_L \xrightarrow{r_{LiacO-1}} Pr_L + m_{TO}$
	$Pr_{TO} \xrightarrow{V_{P_{LtetO-1}}} Pr_{TO} + m_C$
	$Pr_{TQ} \xrightarrow{V_{P_{LtetO}-1}} Pr_{TQ} + m_{I}$
	$Pr_{RA}p_{RA2} \xrightarrow{V_{P_{Lux}I}} Pr_{RA}p_{RA2} + m_{TQ}$ $Pr_{C}p_{C2} \xrightarrow{\ell_{P_{R}}V_{P_{R}}} Pr_{C}p_{C2} + m_{L}$
Leaky transcription	$Pr_C p_{C2} \xrightarrow{\ell_{P_R} v_{P_R}} Pr_C p_{C2} + m_L$
	$Pr_{L}p_{L2} \xrightarrow{\ell_{P_{LlacO-1}}V_{P_{LlacO-1}}} Pr_{L}p_{L2} + m_{TO}$
	$Pr_{TO}p_{T2} \xrightarrow{\ell_{P_{LtetO-1}}V_{P_{LtetO-1}}} Pr_{TO}p_{T2} + m_{C}$
	$Pr_{TO}p_{T2} \xrightarrow{\ell_{P_{LtetO-1}}V_{P_{LtetO-1}}} Pr_{TO}p_{T2} + m_I$
	$Pr_{RA} \xrightarrow{\ell_{P_{LuxI}} V_{P_{LuxI}}} Pr_{RA} + m_{TQ}$
Translation	$m_x \xrightarrow{\epsilon_x} p_x \text{ for } x \in \{C, I, L\}$
AHL production	$m_{Tx} \xrightarrow{\epsilon_{Tx}} p_T \text{ for } x \in \{O, Q\}$ $p_I \xrightarrow{v_3} p_I + A$
ATTL production	$p_I \longrightarrow p_I + A$ $k_{\text{bon}_{AB}}$
LuxR-AHL binding	$p_R + A \xrightarrow{k_{\text{bon}_{AR}}} p_{RA}$
Degradation	$p_x \xrightarrow{\gamma_x} \varnothing \text{ for } x \in \{C, T, L, I\}$
	$m_x \xrightarrow{\gamma_{mO}} \varnothing \text{ for } x \in \{C, L, TO\}$
	$m_x \xrightarrow{\gamma_{mQ}} \varnothing \text{ for } x \in \{I, TQ\}$ $A \xrightarrow{\gamma_A} \varnothing$
Diffusion	$A \xrightarrow{A} \varnothing$ $A \text{ in cell } x \xrightarrow{d_{AHL}} A \text{ in cell } x \pm 1$
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