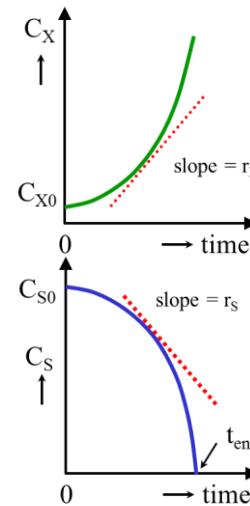


Bioprocesses → Bioreactors

Batch

$$C_x = C_{x0} \exp(\mu^{\max} t)$$

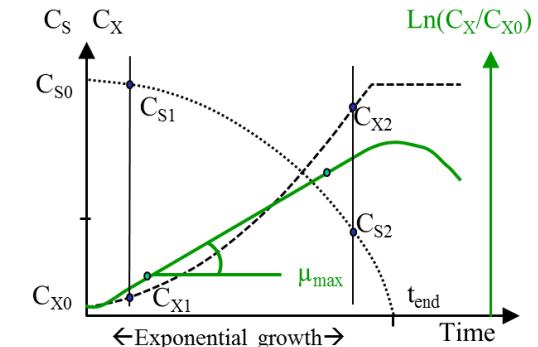
Parameters		state Variables
reactor	operator	
V	C_{S0} C_{X0}	q_s^{\max} μ^{\max} Y_{SX}^{\max} m_s, K_s



$$C_{s0} - C_s = \frac{q_s^{\max}}{\mu^{\max}} C_{X0} [\exp(\mu^{\max} t) - 1]$$

$$\mu^{\max} = \text{slope} \ln\left(\frac{C_x}{C_{x0}}\right)$$

$$Y_{SX} \approx Y_{SX}^{\max} = \frac{C_{X2} - C_{X1}}{C_{s1} - C_{s2}}$$

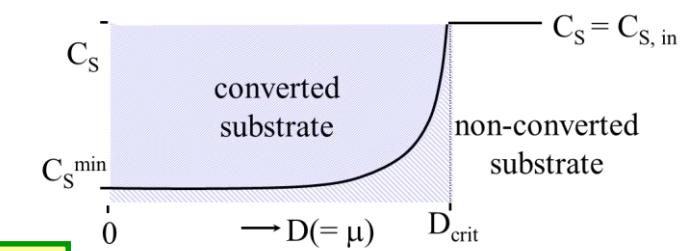


Chemostat

Parameters		state Variables
reactor	operator	
V	$\phi_{L,in}$ $\phi_{L,out}$	C_s C_x

$$C_s = \frac{C_s^{\min} + K_s \frac{D}{\mu^{\max}}}{1 - \frac{D}{\mu^{\max}}}$$

$$C_x = Y_{SX} (\alpha \cdot C_{s,in} - C_s)$$



$$\mu = \frac{r_x}{C_x} = \frac{Q_{L,out}}{V} \stackrel{\text{def}}{=} D : \text{Dilution rate}$$

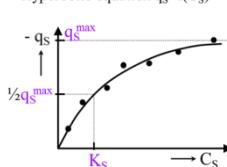
Lineweaver-Burk linearisation

$$\frac{1}{q_s} = \frac{k_s}{q_s^{\max}} \frac{1}{C_s} + \frac{1}{q_s^{\max}}$$

Hanes-Woolf linearisation

$$\frac{C_s}{q_s} = \frac{1}{q_s^{\max}} C_s + \frac{k_s}{q_s^{\max}}$$

Hyperbolic equation $q_s = f(C_s)$



$$D_{crit} = \mu^{\max} \cdot \frac{(C_{s,in} - C_s^{\min})}{K_s + C_{s,in}} \approx \mu^{\max}$$

