

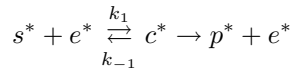
# Enzyme kinetics ala Lin & Segel

Harald Hanche-Olsen

hanche@math.ntnu.no

## Reaction and basic model

The reaction looks like:



which leads to the model

$$\begin{aligned}\frac{ds^*}{dt^*} &= -k_1 s^* e^* + k_{-1} c^* \\ \frac{dc^*}{dt^*} &= k_1 s^* e^* - k_{-1} c^* - k_2 c^* \\ \frac{de^*}{dt^*} &= -k_1 s^* e^* + k_{-1} c^* + k_2 c^* \\ \frac{dp^*}{dt^*} &= k_2 c^*\end{aligned}$$

We shall use the initial conditions

$$s^*(0) = \bar{s}, \quad c^*(0) = 0, \quad e^*(0) = \bar{e}, \quad p^*(0) = 0.$$

**Immediate consequences.** We get  $c^* + e^* = \bar{e}$  and  $s^* + c^* + p^* = \bar{s}$ . Hence we need only solve for  $S^*$  and  $c^*$ . Substitute  $e^* = \bar{e} - c^*$  into the first two equations and get our final, non-scaled model:

$$\begin{aligned}(1) \quad \frac{ds^*}{dt^*} &= -k_1 \bar{e} s^* + (k_1 s^* + k_{-1}) c^* \\ (2) \quad \frac{dc^*}{dt^*} &= k_1 \bar{e} s^* - (k_1 s^* + k_{-1} + k_2) c^* \\ (3) \quad s^*(0) &= \bar{s}, \quad c^*(0) = 0\end{aligned}$$

## Scaling

Put

$$s^* = \bar{s}s, \quad c^* = \bar{e}c, \quad t^* = \frac{t}{k_1\bar{e}}$$

and, with the nondimensional parameters

$$\kappa = \frac{k_{-1} + k_2}{k_1\bar{s}}, \quad \lambda = \frac{k_2}{k_1\bar{s}}, \quad \varepsilon = \frac{\bar{e}}{\bar{s}}$$

we have the problem on its non-dimensional form:

$$\begin{aligned} \dot{s} &= -s + (s + \kappa - \lambda)c \\ \varepsilon \dot{c} &= s - (s + \kappa)c \\ s(0) &= 1, \quad c(0) = 0 \end{aligned}$$

For later reference, it will be useful to remember that  $0 < \lambda < \kappa$ , and we shall assume  $0 < \varepsilon \ll 1$ .

## Solution by perturbation

**Outer solution.** First, put  $\varepsilon = 0$  in the differential equations. The second equation becomes the algebraic equation  $s - (s + \kappa)c = 0$ . We solve this for  $c$  and substitute in the first equation, which becomes  $\dot{s} = -\lambda s / (s + \kappa)$ . We call the solution  $(s_0, c_0)$ :

$$(4) \quad s_0 + \kappa \ln s_0 = Q - \lambda t, \quad c_0 = \frac{s_0}{s_0 + \kappa}$$

The integration constant  $Q$  might be determined by using  $s_0(0) = 1$ , so  $Q = 1$ . But then the other initial condition is not satisfied, and so we are not quite sure whether even the first one is satisfied. So presumably,  $Q$  must be determined by matching.

**Inner solution.** Initially we expect  $s \approx 1$ , in which case  $\varepsilon \dot{c} \approx 1 - (1 + \kappa)c$ . Thus  $c$  seems to tend towards  $1 / (1 + \kappa)$ . Introduce this time constant as a new time scale, and define the *inner dimensionless time*  $\tau$  by<sup>1</sup>

$$t = \frac{\varepsilon}{1 + \kappa} \tau$$

---

<sup>1</sup>Here we depart from Lin & Segel in a small way. It simplifies a few formulae, but changes nothing important.

With  $s(t) = S(\tau)$  and  $c(t) = C(\tau)$  the equations become

$$\begin{aligned}(1 + \kappa)S' &= \varepsilon(-S + (S + \kappa - \lambda)C) \\ (1 + \kappa)C' &= S - (S + \kappa)C\end{aligned}$$

and, of course, we can use the initial conditions as well, so  $S(0) = 1$  and  $C(0) = 0$ .

With  $\varepsilon = 0$  we get  $S' = 0$ , so the initial conditions imply  $S = 1$ . We plug that into the second equation and get  $C' = \frac{1}{1+\kappa} - C$ , which is easy to solve with the initial condition  $C = 0$ . Renaming this solution  $(S_0, C_0)$  we thus have

$$S_0(\tau) = 1, \quad C_0(\tau) = \frac{1 - e^{-\tau}}{1 + \kappa}$$

**Matching.** To this order, we can match inner and outer solutions by simply requiring that  $\lim_{\tau \rightarrow \infty} S(\tau) = \lim_{t \rightarrow 0} s(t)$ . Thus  $Q = 1$ , as we guessed before. The similar equation  $\lim_{\tau \rightarrow \infty} C(\tau) = \lim_{t \rightarrow 0} c(t)$  turns out to be automatically satisfied.

**To the next order.** We try to substitute power series  $S = S_0 + \varepsilon S_1 + \dots$ ,  $C = C_0 + \varepsilon C_1 + \dots$ ,  $s = s_0 + \varepsilon s_1 + \dots$ , and  $c = c_0 + \varepsilon c_1 + \dots$  into the inner and outer equations. For  $S_0$ ,  $C_0$ ,  $s_0$  and  $c_0$  we find the equations and solutions we have already discovered. Next, we find

$$\begin{aligned}S_1 &= -\frac{1}{(1 + \kappa)^2}(\lambda\tau + (1 + \kappa - \lambda)(1 - e^{-\tau})) \\ C_1 &= -\frac{1}{(1 + \kappa)^4}(\lambda\kappa\tau + \kappa(1 + \kappa - 2\lambda) + (1 + \kappa - \lambda)e^{-2\tau} \\ &\quad + [\frac{1}{2}\lambda\tau^2 + (1 + \kappa - \lambda)(1 - \kappa)\tau - ((1 + \kappa)^2 - \lambda - 2\kappa\lambda)]e^{-\tau})\end{aligned}$$

while  $s_1$  and  $c_1$  are given by

$$\dot{s}_1 = (c_0 - 1)s_1 + (\kappa - \lambda + s_0)c_1, \quad \dot{c}_0 = s_1(1 - c_0) - (\kappa + s_0)c_1$$

which requires an initial condition for  $s_1$ .<sup>2</sup> So we introduce the intermediate time scale  $\Psi$ , put  $t = \Psi\tau_i$ , and note that

$$\tau = \frac{1 + \kappa}{\varepsilon}t = \frac{(1 + \kappa)\Psi}{\varepsilon}\tau_i$$

---

<sup>2</sup>At this stage, we do not need an initial condition for  $c_1$ , as this system is actually just one differential equation and one algebraic equation,  $c_0$  being a known function.

We get, therefore

$$\begin{aligned} S(\tau) &= S_0 \left( \frac{(1 + \kappa)\Psi}{\varepsilon} \tau_i \right) + \varepsilon S_1 \left( \frac{(1 + \kappa)\Psi}{\varepsilon} \tau_i \right) + \dots \\ &= 1 - \frac{\lambda \tau_i \Psi}{1 + \kappa} - \frac{(1 + \kappa - \lambda)\varepsilon}{(1 + \kappa)^2} + \text{TST} + \dots \end{aligned}$$

which needs to be matched to  $s(t) = s_0(t) + \varepsilon s_1(t) + \dots$  with  $t = \Psi \tau_i$ .

Thus we need a power series representation for  $s_0$  in terms of  $t$ : Write  $s_0(t) = 1 + \gamma t + \dots$  and insert into the first part of (4):  $1 + \gamma t + \kappa \gamma t + \dots = 1 - \lambda t$ , so  $\gamma = -\lambda/(1 + \kappa)$ , and we can write

$$s(t) = s_0(\tau_i \Psi) + \varepsilon s_1(\tau_i \Psi) + \dots = 1 - \frac{\lambda \tau_i \Psi}{1 + \kappa} + s_1(0)\varepsilon + \dots$$

and we see that the required matching implies

$$s_1(0) = -\frac{(1 + \kappa - \lambda)\varepsilon}{(1 + \kappa)^2}.$$

Note: the next term in the approximation for  $s(t)$  would contain a factor  $\Psi^2$ . In order for our argument to remain valid, we need  $\Psi^2 \ll \varepsilon$  as  $\varepsilon \rightarrow 0$ ; thus we must have

$$\lim_{\varepsilon \rightarrow 0} \frac{\Psi}{\varepsilon} = \infty, \quad \lim_{\varepsilon \rightarrow 0} \frac{\Psi}{\sqrt{\varepsilon}} = 0.$$

## A better scaling

The above analysis is due to Lin & Segel. An improved scaling was introduced by Segel & Slemrod (see References, below).

They find the proper *time scales* first. (The choice of  $\bar{s}$  as the scale for  $s^*$  is pretty obvious even at this stage, though.)

After the initial transient, in the quasi-steady state  $dc^*/dt^* \approx 0$ . Assuming this is exact, we solve (2) for  $c^*$  to get

$$c^* = \frac{s^*}{s^* + K} \bar{e}, \quad \text{where } K = \frac{k_{-1} + k_2}{k_1}.$$

Substitute into (1) and simplify, to get

$$\frac{ds^*}{dt^*} = -\frac{k_2 \bar{e} s^*}{s^* + K}$$

(which is easily solved, but never mind that for now). Clearly the maximal value of  $s^*$  is  $\bar{s}$ , and one good way to get a time scale is

$$(5) \quad T = \frac{\max |s^*|}{\max \left| \frac{ds^*}{dt^*} \right|} = \frac{\bar{s}}{\frac{k_2 \bar{e} \bar{s}}{\bar{s} + K}} = \frac{\bar{s} + K}{k_2 \bar{e}}.$$

This, then is the *long time scale* of the problem.

On the other hand, during the initial transient  $s^* \approx \bar{s}$  and so (2) becomes

$$\frac{dc^*}{dt^*} \approx k_1 \bar{e} \bar{s} - k_1 (\bar{s} + K) c^*.$$

The equilibrium of this equation is at

$$c^* = \frac{\bar{e} \bar{s}}{\bar{s} + K},$$

so this value is actually a good scale for  $c^*$ . Moreover, the approach to this equilibrium is exponential with a time constant

$$\varepsilon T = \frac{1}{k_1 (\bar{s} + K)}.$$

This is the appropriate choice of the *short time constant*. The ratio between the two time constants is

$$\varepsilon = \frac{k_2 \bar{e}}{k_1 (\bar{s} + K)^2}.$$

We must have  $\varepsilon \ll 1$ , or else our assumptions of very different time scales is wrong, and our analysis becomes suspect. (But in practice, we can get surprisingly good results even for  $\varepsilon \approx 1$ .)

We thus end up with the following scalings, for the outer solution:

$$s^* = \bar{s} s, \quad c^* = \frac{\bar{e} \bar{s}}{\bar{s} + K} c, \quad t^* = \frac{\bar{s} + K}{k_2 \bar{e}} t.$$

For the inner solution the rescaled dimensionless time  $\tau$  is given by

$$t = \varepsilon \tau$$

where  $\varepsilon$  is given above.

## Some principles

Segel & Slemrod used these general principles to derive their scaling:

1. Dependent variables ( $s^*$  and  $c^*$  in our example) should be scaled according to their maximal value, so that their dimensionless versions vary between 0 and 1.
2. Independent variables ( $t^*$  in our example) should be scaled so that the dependent variables vary considerably over the chosen scale. In other words, the derivatives of the dependent variables ought to have a maximum value of 1 in the scaled model. (In our example, this was achieved by equation (5).)

## References

C.C LIN & L.A. SEGEL: *Mathematics applied to deterministic problems in the natural sciences*. ISBN 0898712297. Chapter 10 (pp. 302–320).

LEE A. SEGEL & MARSHALL SLEMROD: The quasi-steady state assumption: A case study in perturbation. *SIAM Review* **31** #3 (1989), pp. 446–477 (available at the mathematics library).