

## REACTION KINETICS

In this lecture we present an example from “reaction kinetics”. The purpose is to describe how to set up the kinetic equations, to solve them with MATLAB, and to draw a conclusion about the reaction from the computation. Perhaps this will be useful in you own [project](#).

## 1. OXIDATION OF NO

Inspired by: P. Atkins and L. Jones, *Chemical Principles. The Quest for Insight*. Freeman, New York, first edition, 1999, pp. 624–625.

We consider the oxidation of NO to  $\text{N}_2\text{O}_2$ . The following rate of formation of  $\text{N}_2\text{O}_2$  has been observed experimentally:

$$(1) \quad \frac{d}{dt}[\text{N}_2\text{O}_2] = k[\text{O}_2][\text{NO}]^2. \quad (\text{mol}/(\text{L s}))$$

This corresponds to the formula



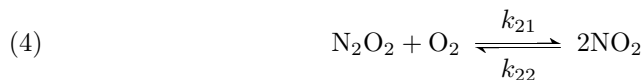
Here  $[\text{N}_2\text{O}_2]$ , measured in  $\text{mol}/\text{L}$ , is the concentration of  $\text{N}_2\text{O}_2$ . The rate constant  $k$  is measured in  $\text{L}^2/(\text{mol}^2\text{s})$ . This is a *third order reaction*, because the rate is proportional to the product of three concentrations.

In order to explain this empirical formula one has proposed the following two-step reaction mechanism.

**Step 1.** NO is decomposed into an intermediate product  $\text{N}_2\text{O}_2$  in a fast reaction:



**Step 2.** The intermediate product reacts with  $\text{O}_2$  in a slow reaction:



We now write down the reaction rates for the four reactions ( $\text{mol}/(\text{L s})$ ):

$$(5) \quad \begin{aligned} r_{11} &= k_{11}[\text{NO}]^2 &&= k_{11}u_1^2, \\ r_{12} &= k_{12}[\text{N}_2\text{O}_2] &&= k_{12}u_3, \\ r_{21} &= k_{21}[\text{N}_2\text{O}_2][\text{O}_2] &&= k_{21}u_3u_2, \\ r_{22} &= k_{22}[\text{NO}_2]^2 &&= k_{22}u_4^2. \end{aligned}$$

Here we introduced the variables

$$(6) \quad u_1 = [\text{NO}], \quad u_2 = [\text{O}_2], \quad u_3 = [\text{N}_2\text{O}_2], \quad u_4 = [\text{NO}_2]. \quad (\text{mol}/\text{L})$$

What are the units of the rate constants  $k_{11}, k_{12}, k_{21}, k_{22}$ ?

Finally, we write down the differential equations for the concentrations:

$$(7) \quad \begin{aligned} \dot{u}_1 &= -2r_{11} + 2r_{12}, \\ \dot{u}_2 &= -r_{21} + r_{22}, \\ \dot{u}_3 &= r_{11} - r_{12} - r_{21} + r_{22}, \\ \dot{u}_4 &= 2r_{21} - 2r_{22}. \end{aligned}$$

The numbers  $\pm 1, \pm 2$  that occur in front of the rates are called *stoichiometric numbers*. For example, the stoichiometric numbers of  $\text{NO}_2$  in reactions 21 and 22 are 2 and -2, respectively.

The equations are implemented and solved in the MATLAB programs [no2.m](#), [no2a.m](#), [no2test.m](#). The file `no2test.m` is a *script file*, which starts the computations and plots the solutions.

We use the data  $k_{11} = 10$ ,  $k_{12} = 10$  (fast reaction),  $k_{21} = 0.01$ ,  $k_{22} = 0$  (slow reaction), and initial values  $u_{10} = 0.5$ ,  $u_{20} = 1$ ,  $u_{30} = 0$ ,  $u_{40} = 0$ . Download the programs and compute!

In order to test if the proposed two-step mechanism explains the empirical third order rate law (1), we also compute the quotient

$$(8) \quad \frac{\dot{u}_4}{u_2 u_1^2}.$$

If (1) holds, then this quotient should be constant  $= k$ . A calculation in Atkins and Jones shows that  $k = 2k_{21}k_{11}/k_{12}$ . The quotient (8) is plotted with black dots, and we see that it quickly becomes constant  $= 2k_{21}k_{11}/k_{12}$ .

This verifies that the two-step mechanism explains the empirical rate law (1).

You can download the MATLAB programs from

<http://www.phil.chalmers.se/education/courses/2001/ala-b/matlab/facit/>