

1 Introduction

In this tutorial we will explore the ways of presenting simple kinetic data and calculating kinetic parameters from that data. Starting with the ideal case, we will then demonstrate the effect of experimental errors on the determination of the kinetic parameters using each of the methods.

1.1 The ideal model

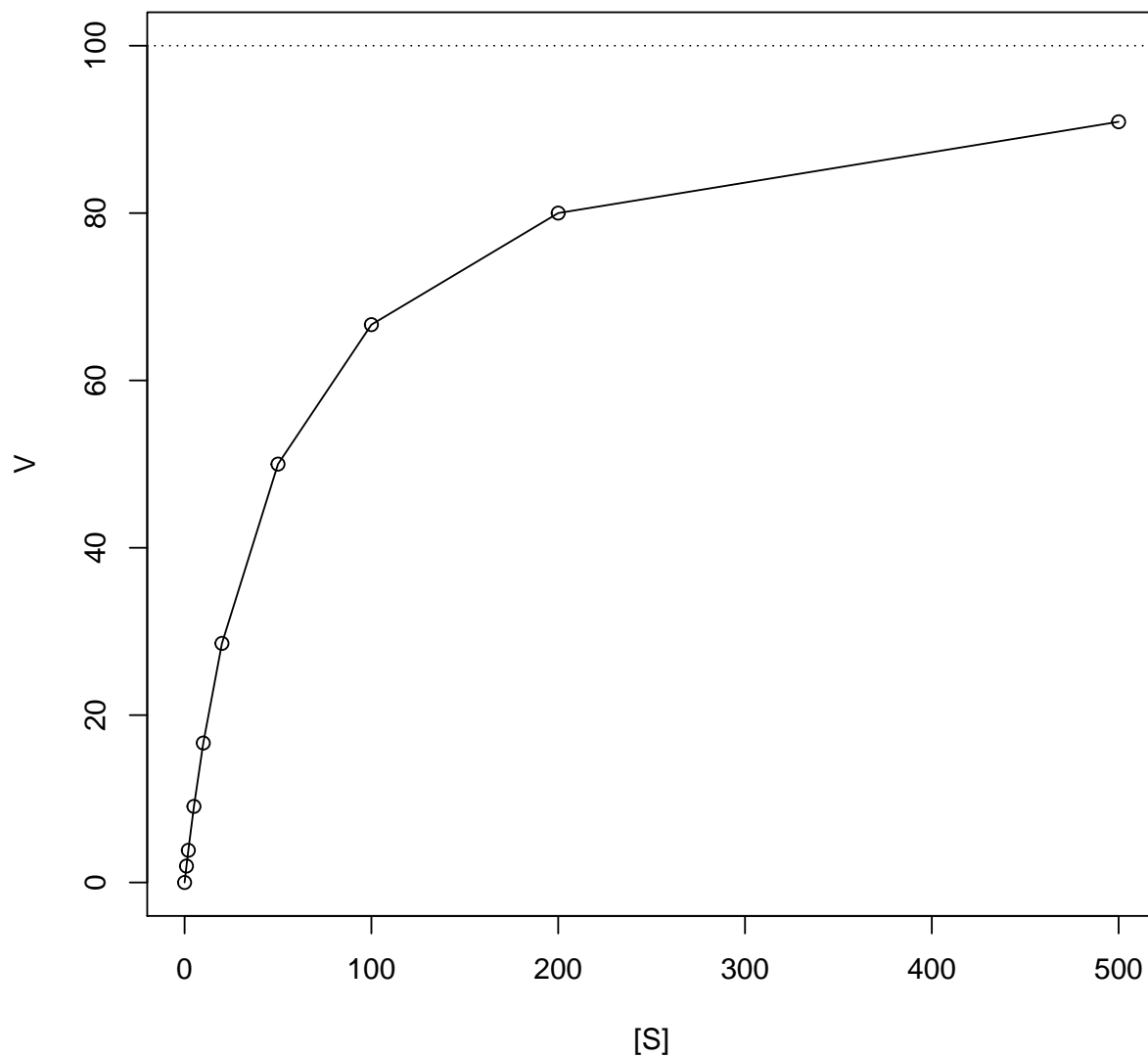
A simple kinetic reaction follows the formula $V = \frac{V_{max} \cdot [S]}{K_m + [S]}$

We can set the V_{max} and K_m to appropriate values and calculate the observed velocity V at each value of the substrate concentration $[S]$.

With V_{max} set to 100 and K_m to 50 we can calculate the effective rate for values of the substrate concentration from 0 to 500 using a standard serial dilution.

```
vmax <- 100
km <- 50
substrate <- c(0, 1, 2, 5, 10, 20, 50, 100, 200, 500)
velocity <- vmax * substrate / (substrate + km)
```

The velocity can then be plotted in several ways. The simplest way is to plot the rate against the substrate concentration. The dotted line indicates V_{max}

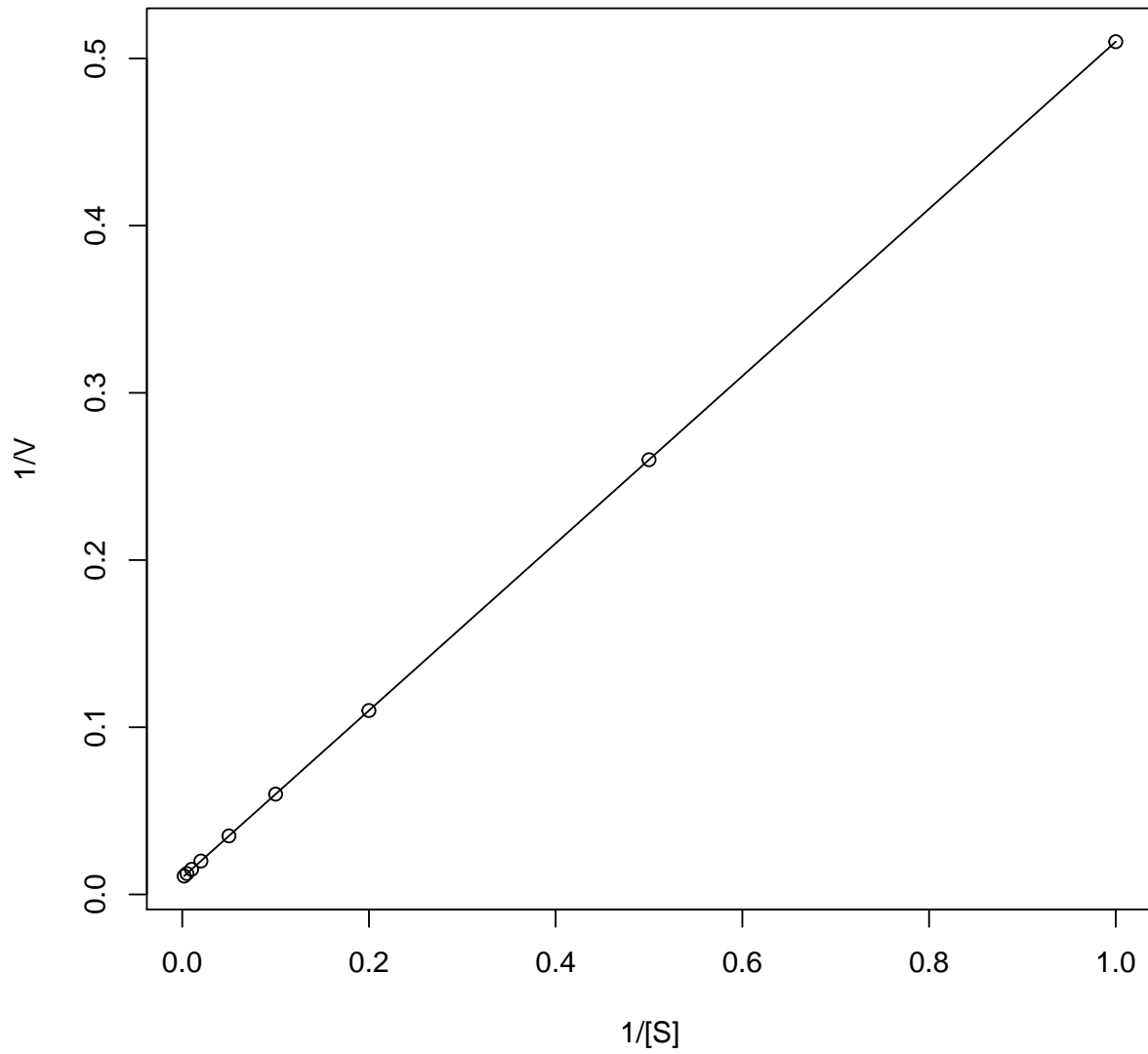


1.1.1 The Lineweaver-Burke plot

It is hard to estimate the K_m or V_{max} from such plots so typically we transform the data to a straight line, fit the straight line and use this to determine the K_m and V_{max} .

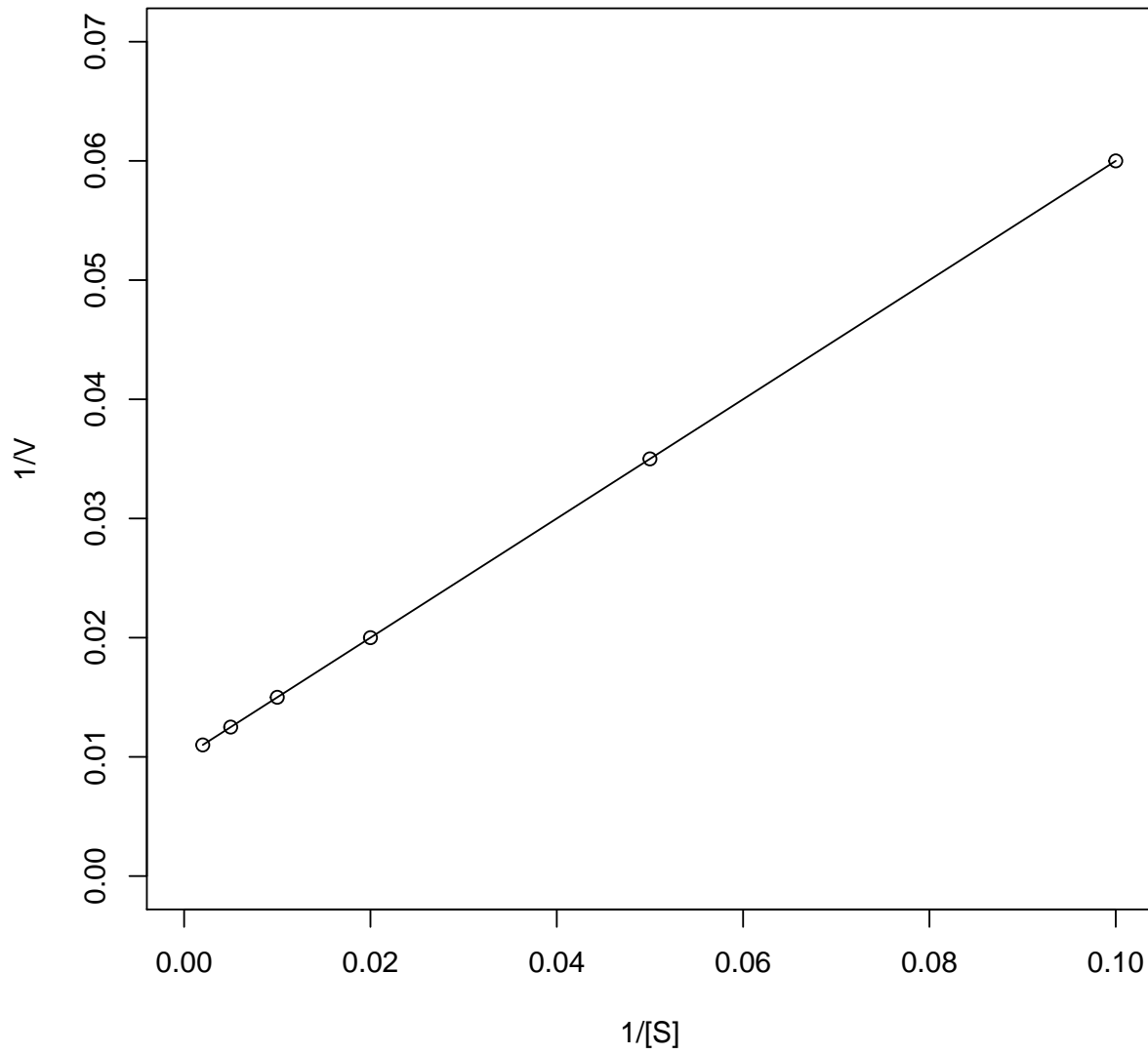
Traditionally students would use the Lineweaver-Burke plot. This is the reciprocal of the previous plot.

Lineweaver–Burke plot



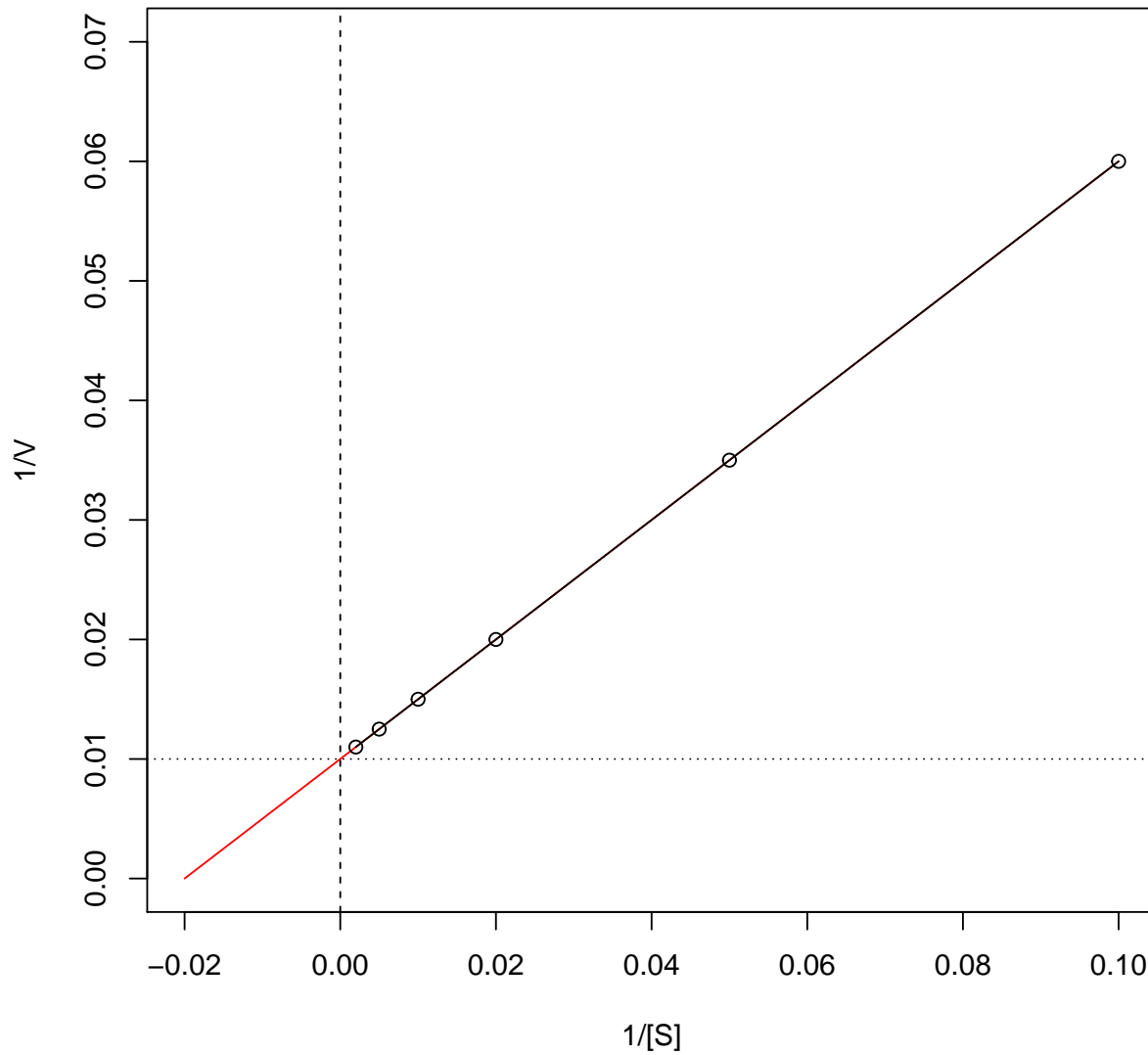
Zooming in on the Y axis:

Lineweaver–Burke plot



Extending the line to intercept the X-axis gives $1/V_{max}$ as the Y intercept and $-1/K_m$ as the X intercept.

Lineweaver–Burke plot

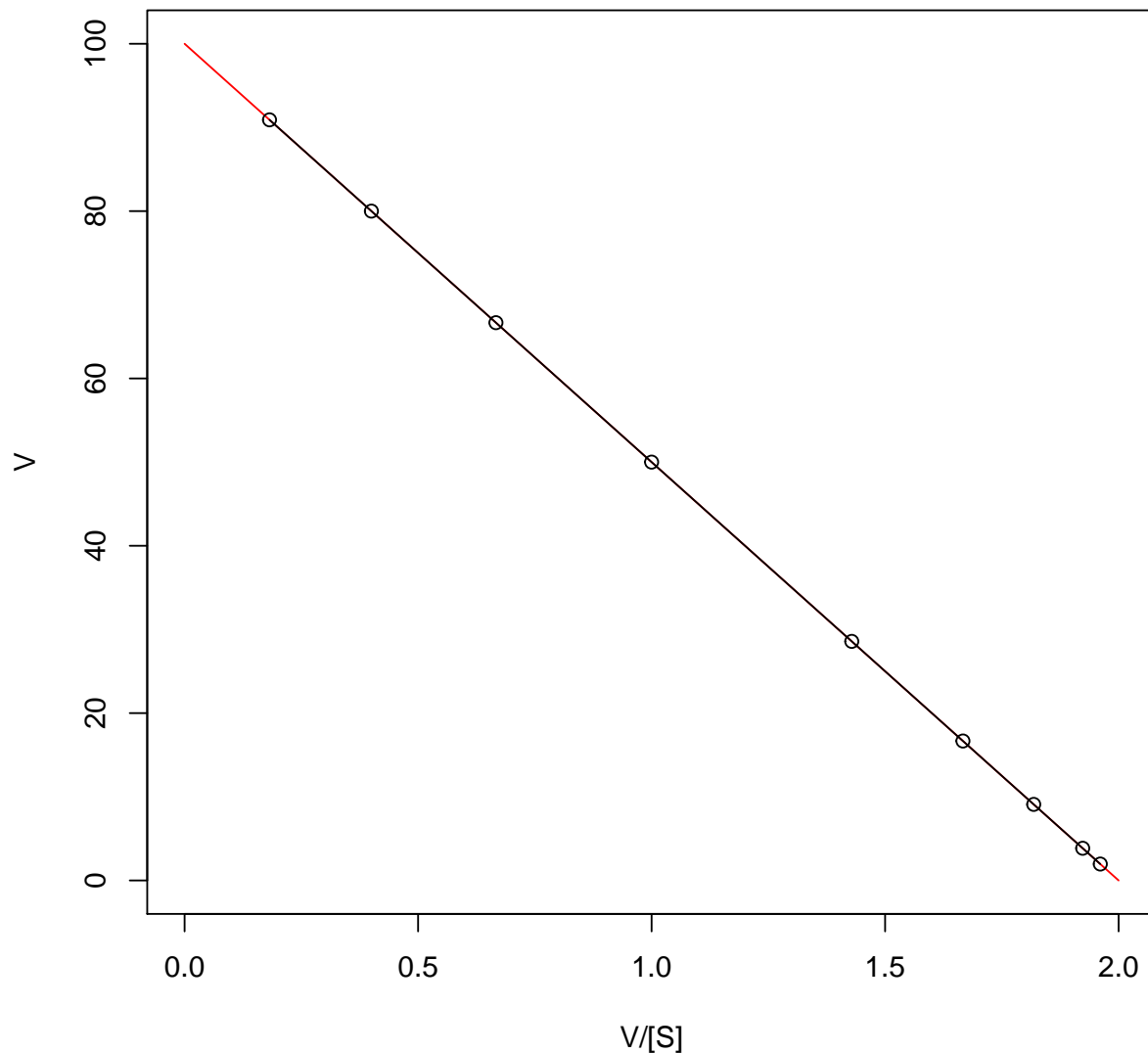


1.1.2 The Eadie-Hofstee plot

Alternative transformations which are less susceptible to errors are the Eadie-Hofstee plot and the Haynes-Woolf plot.

Eadie-Hofstee plots the velocity against the velocity divided by the substrate concentration. This gives a straight line of the form $v = -K_m \frac{v}{[S]} + V_{max}$ where v is the y-axis, $\frac{v}{[S]}$ is the x-axis, the intercept is V_{max} and the gradient is $-K_m$. Again, the line can be extrapolated to give the intercept as indicated by the red line.

Eadie-Hofstee plot

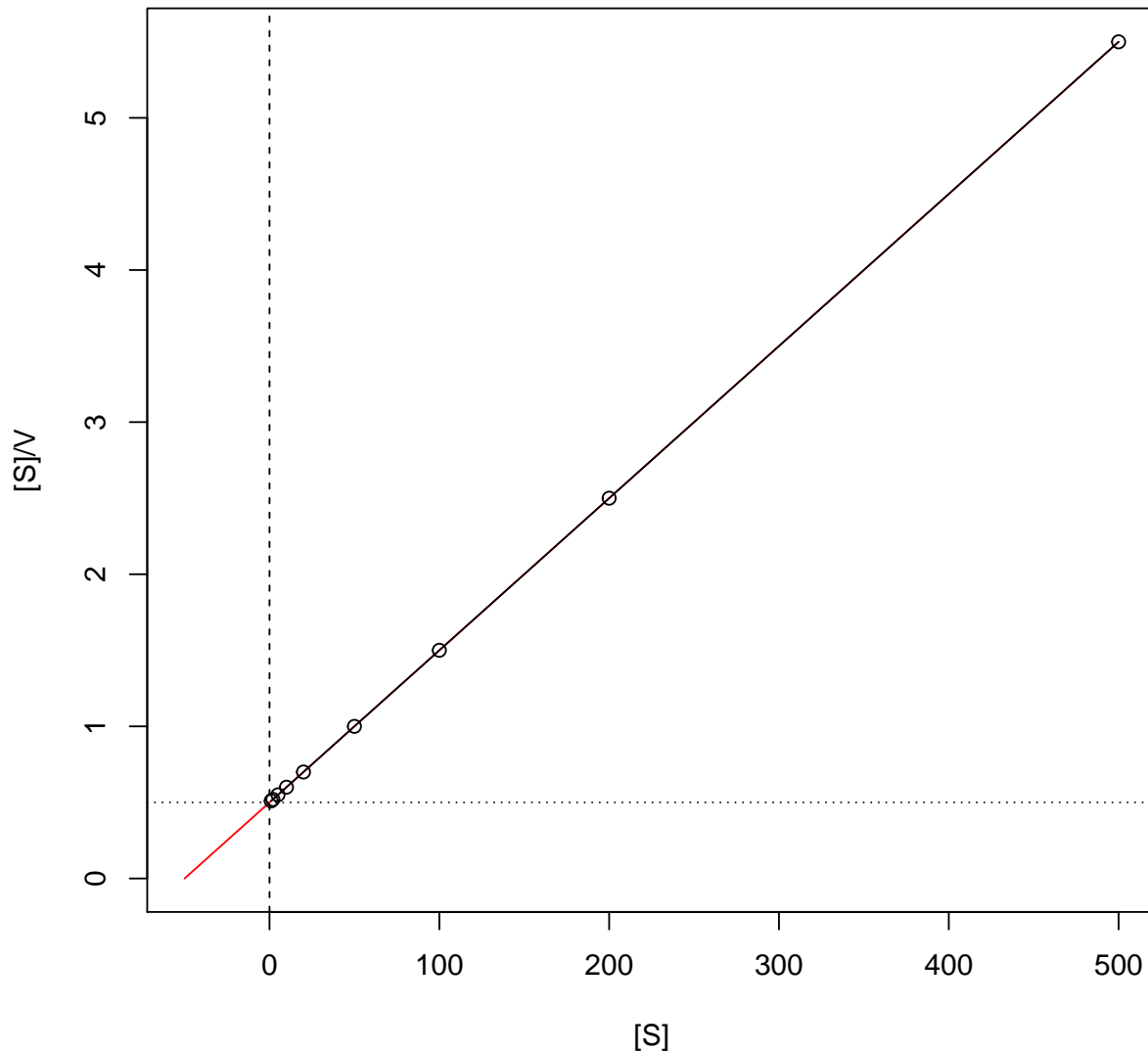


1.1.3 The Haynes-Woolf plot

The Haynes-Woolf plot is possibly the most accurate of the linear transformations. Dividing the substrate concentration by the velocity, and plotting this against the substrate concentration gives a plot which looks superficially like a Lineweaver-Burke plot but does not fall prey to the same reciprocal error scaling.

The graph forms a straight line according to the equation $\frac{[S]}{v} = \frac{1}{V_{max}}[S] + \frac{K_m}{V_{max}}$. Like the Lineweaver-Burke plot, the line can be extended through the y-axis to intercept the x-axis at $-K_m$, the y-axis intercept being at $\frac{K_m}{V_{max}}$.

Haynes–Woolf plot



With modern computers, fitting the experimental data directly is a good idea as the transformations amplify errors in the data.

1.2 Introducing experimental errors

Errors can be introduced into an experiment in many ways and fall into several categories. We will include fixed errors (which may be found in for example machine reading variation) and proportional errors such as those found in pipetting or errors in fixed time measurement when measuring rates.

We will implement these using the `jitter()` command in a function that will return a set of experimental results with error added.

```
fixederror <- velocity - jitter(velocity, factor = 2)
variableerror <- velocity * (velocity - jitter(velocity, factor = 0.2))
```

We can combine these errors and build a table of experimental values. Each time `jitter()` is called it produces a new random distribution of errors. These values will be saved in a table of substrate (column 1)

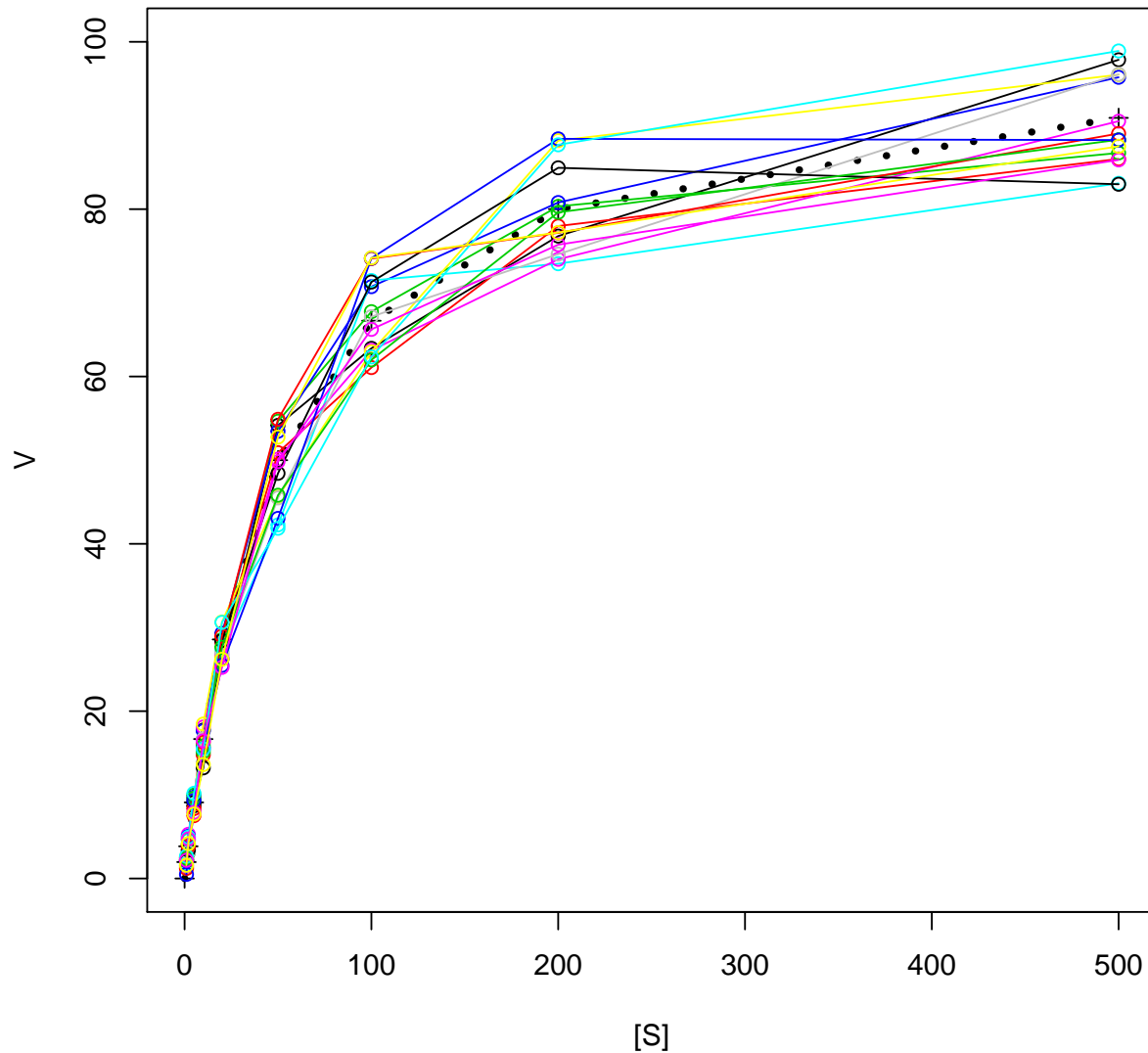
and velocity measurement (column 2) with the experiment number in column 3. The velocity calculations will be performed in a function `doexperiment` which takes as parameters V_{max} , K_m , and the substrate concentrations.

```
doexperiment <- function(vm, k, subs) {  
  # add some proportional error to the substrate for concentration errors.  
  errsubs <- subs + subs * (subs - jitter(subs))  
  # calculate the velocity that should be observed  
  vobs <- vm * (errsubs/(k + errsubs))  
  # add measurement (fixed) and time errors (proportional) to the observed  
  # value  
  errvobs <- jitter(vobs, factor = 2) + vobs * (vobs - jitter(vobs, factor = 0.2))  
  # return the experimental values  
  errvobs  
}
```

```
experiment <- cbind(substrate[-1], doexperiment(vmax, km, substrate[-1]), 1)  
for (e in 2:15) {  
  experiment <- rbind(experiment, cbind(substrate[-1], doexperiment(vmax,  
    km, substrate[-1]), e))  
}
```

The latter line is a loop that gives us many replicates as we want. In this case we repeat 15 times.

Experimental results



1.2.1 Fitting the data

To fit the data we can use the drc package. (Note: these commands do not run automatically. You will need to run the install once and the library command each time you start R.)

```
install.packages("drc")
```

```
library(drc)
```

The drm method can be used to fit the data to a 2-parameter Michaelis-Menten equation. It takes a data frame as argument and fits velocity against substrate concentration. First we can test how well it works against ideal data:

```
ideal <- as.data.frame(cbind(substrate, velocity))  
# create the data frame
```

```
colnames(ideal) <- c("substrate", "velocity")
# set the column names
fit <- drm(velocity ~ substrate, data = ideal, fct = MM.2())
# fit the data to the equation
```

The results are found in `fit$coefficients`. The first value is the fit to V_{max} and the second is K_m .

```
fit$coefficients

## d:(Intercept) e:(Intercept)
##           100           50
```

The ideal data has, as expected given a perfect fit.

We can fit the data for each experiment and see how the results compare to the real result.

```
# first convert our experimental data to a data frame
experiment <- as.data.frame(experiment)
colnames(experiment) <- c("substrate", "velocity", "expr")
data <- subset(experiment, expr == 1)
fit <- drm(velocity ~ substrate, data = data, fct = MM.2())
mmres <- as.data.frame(cbind(fit$coefficients[1], fit$coefficients[2]))
mmres

##           V1    V2
## d:(Intercept) 104.2 55.8
```

This is not as good a fit. We can complete the rest of the fits with a loop and see how they match to the ideal fit.

```
for (e in 2:15) {
  data <- subset(experiment, expr == e)
  fit <- drm(velocity ~ substrate, data = data, fct = MM.2())
  mmres <- rbind(mmres, fit$coefficients)
}
colnames(mmres) <- c("Vmax", "Km")
mean(mmres[, "Vmax"])

## [1] 100.4

sd(mmres[, "Vmax"])

## [1] 5.933

mean(mmres[, "Km"])

## [1] 51.87

sd(mmres[, "Km"])

## [1] 7.993
```

1.2.2 Fitting to the Lineweaver-Burke plot

To fit the Lineweaver-Burke plot (and any of the other linear plots) we first need to calculate the x and y values for each datapoint.

```
x <- 1/substrate[-1]
y <- 1/velocity[-1]
```

These can be fitted to a straight line with the `lm` function.

```
fit <- lm(y ~ x)
fit

##
## Call:
## lm(formula = y ~ x)
##
## Coefficients:
## (Intercept)          x
##      0.01      0.50
```

The coefficients are the intercept and gradient. V_{max} is the reciprocal of the intercept. K_m is the gradient divided by the intercept.

```
# calculate Vmax
1/fit$coefficients[1]

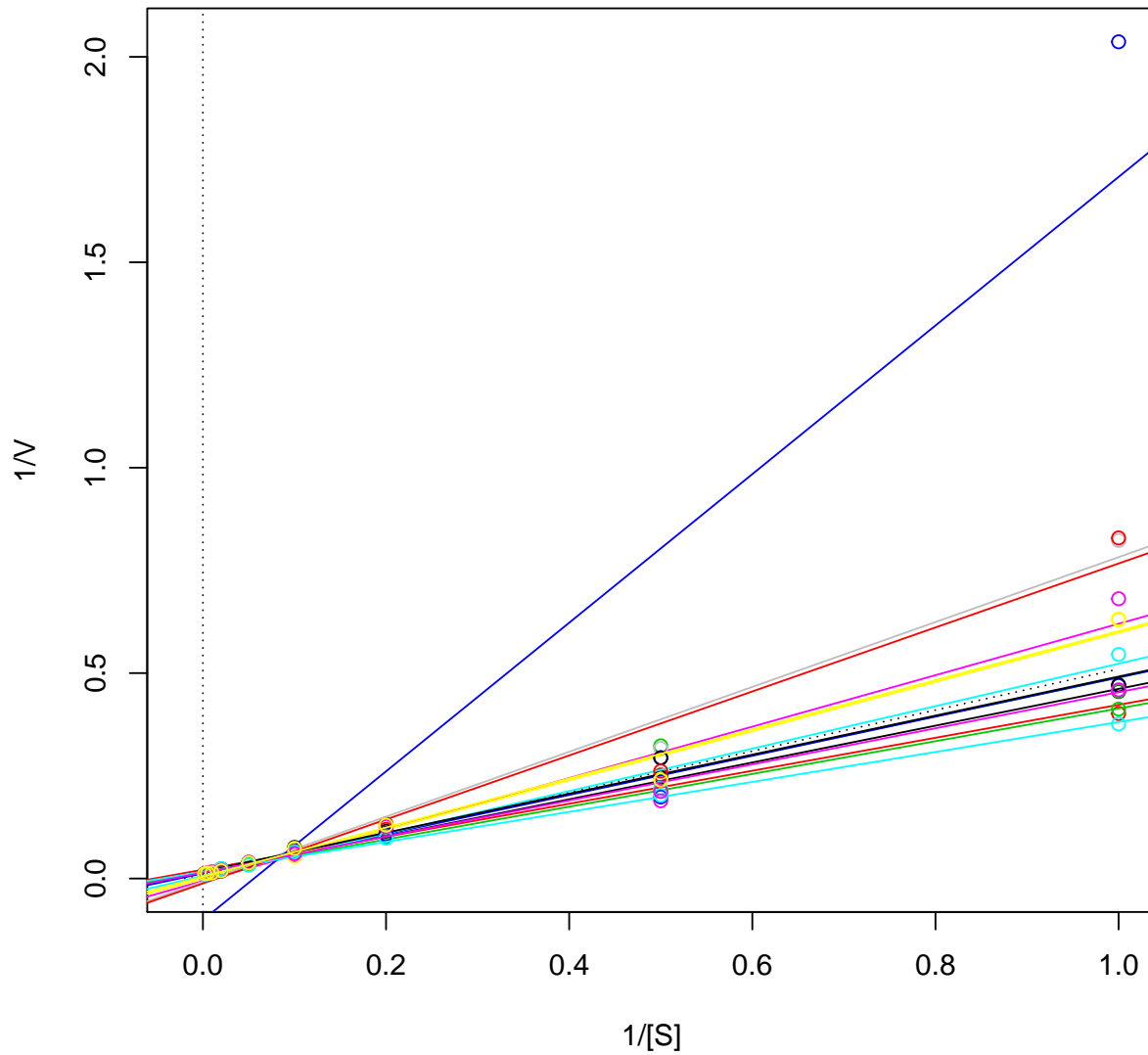
## (Intercept)
##      100

# calculate Km
fit$coefficients[2]/fit$coefficients[1]

## x
## 50
```

Again, we will repeat this for the experimental data. First we will plot the transformed data, then fit a straight line to each data set.

Lineweaver–Burke plot



The estimates from the experimental data are significantly worse than a direct fit.

```
mean(lbres$Vmax)
## [1] -59.61
sd(lbres$Vmax)
## [1] 382.6
mean(lbres$Km)
## [1] -46.02
sd(lbres$Km)
## [1] 227.2
```

1.2.3 Fitting to the Eadie-Hofstee plot

To fit the Eadie-Hofstee we first need to calculate the x and y values for each datapoint.

```
x <- velocity[-1]/substrate[-1]
y <- velocity[-1]
```

These can be fitted to a straight line with the `lm` function.

```
fit <- lm(y ~ x)
fit

##
## Call:
## lm(formula = y ~ x)
##
## Coefficients:
## (Intercept)          x
##          100         -50
```

The coefficients are the intercept and gradient. V_{max} is the intercept. K_m is the gradient.

```
# calculate Vmax
fit$coefficients[1]

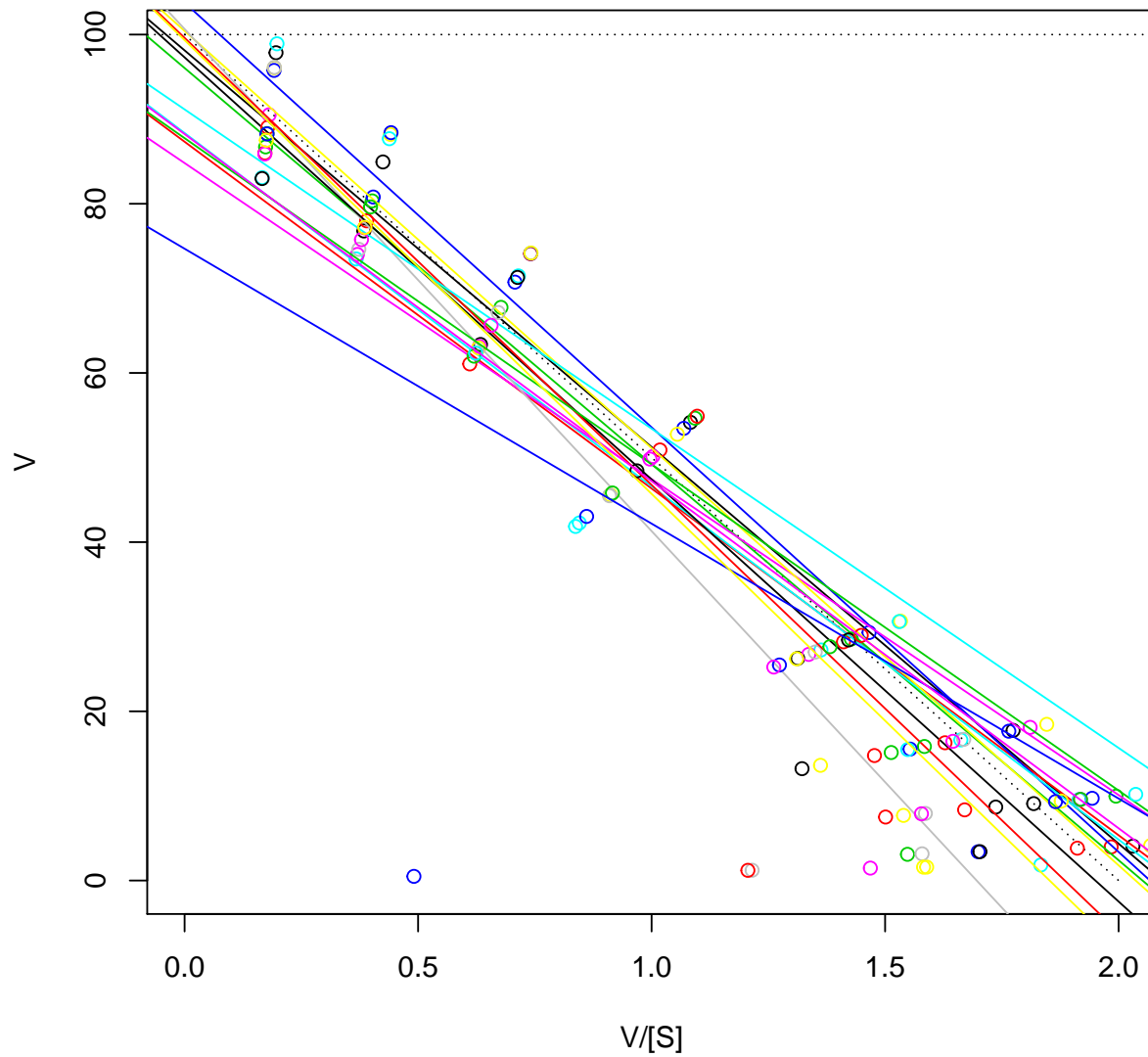
## (Intercept)
##          100

# calculate Km
-fit$coefficients[2]

## x
## 50
```

Again, we will repeat this for the experimental data. First we will plot the transformed data, then fit a straight line to each data set.

Eadie-Hofstee plot



The estimates from the experimental data are significantly better than the Lineweaver-Burke plot.

```
mean(ehres$Vmax)
## [1] 93.17
sd(ehres$Vmax)
## [1] 7.914
mean(ehres$Km)
## [1] 45.26
sd(ehres$Km)
## [1] 7.41
```

1.2.4 Fitting to the Haynes-Woolf plot

To fit the Haynes-Woolf plot we first need to calculate the x and y values for each datapoint.

```
x <- substrate[-1]
y <- substrate[-1]/velocity[-1]
```

These can be fitted to a straight line with the `lm` function.

```
fit <- lm(y ~ x)
fit

##
## Call:
## lm(formula = y ~ x)
##
## Coefficients:
## (Intercept)          x
##          0.50          0.01
```

The coefficients are the intercept and gradient. $\frac{K_m}{V_{max}}$ is the y-intercept. $-K_m$ is the x-intercept.

```
# calculate Vmax
1/fit$coefficients[2]

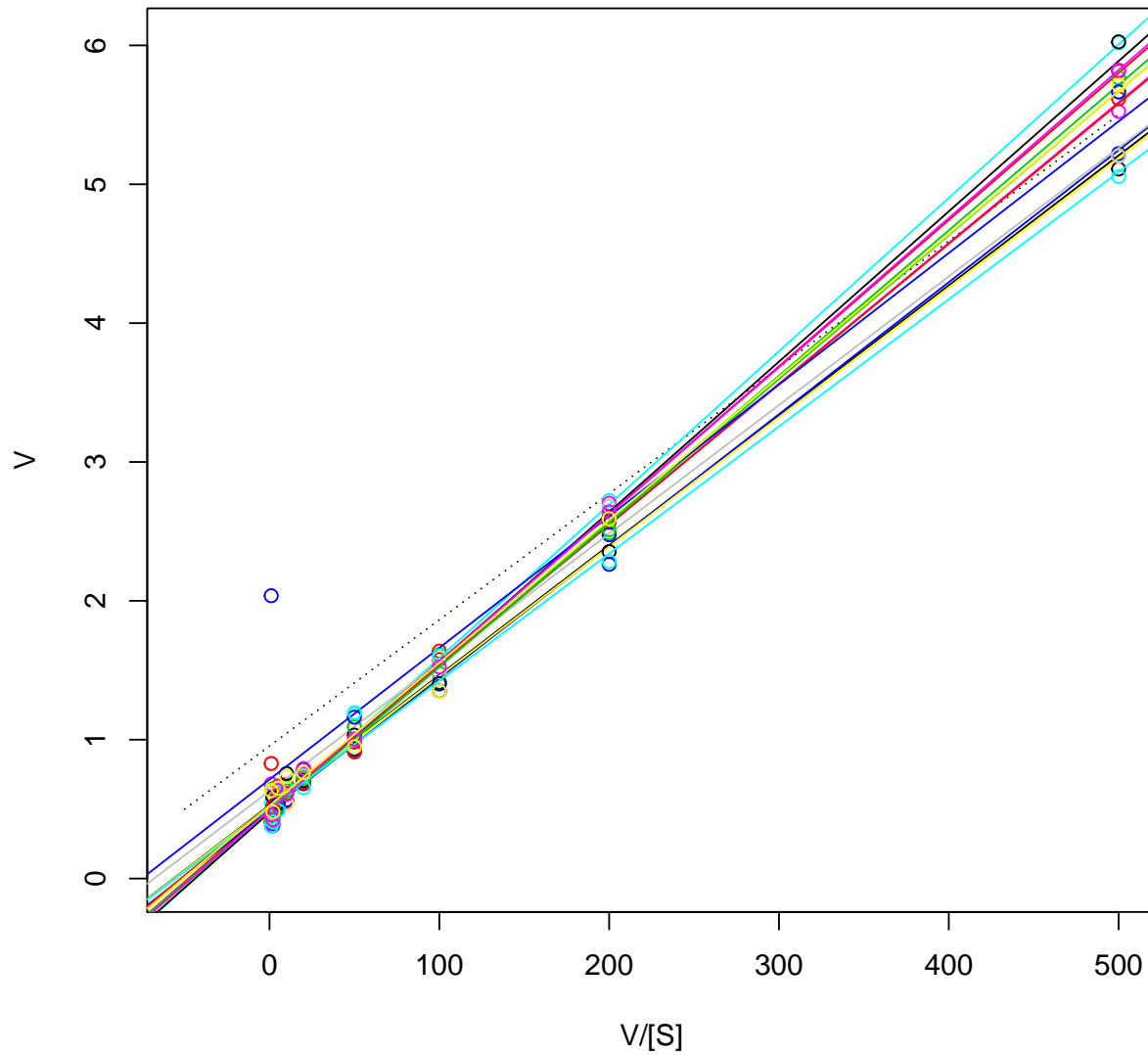
##      x
## 100

# calculate Km
fit$coefficient[1]/fit$coefficients[2]

## (Intercept)
##          50
```

Again, we will repeat this for the experimental data. First we will plot the transformed data, then fit a straight line to each data set.

Haynes-Woolf plot



The estimates from the experimental data are significantly better than the Lineweaver-Burke plot.

```
mean(hwres$Vmax)
## [1] 99.97
sd(hwres$Vmax)
## [1] 6.345
mean(hwres$Km)
## [1] 52.79
sd(hwres$Km)
## [1] 8.804
```


1.3 Conclusions

We can compare the four methods described here - three linear and one direct fitting.

```
mmres$method <- "Direct"
lbres$method <- "Lineweaver-Burke"
ehres$method <- "Eadie-Hoftzee"
hwres$method <- "Haynes-Woolf"
results <- rbind(mmres, lbres, ehres, hwres)
tapply(results$Vmax, results$method, mean)

##          Direct      Eadie-Hoftzee      Haynes-Woolf Lineweaver-Burke
##          100.38           93.17           99.97           -59.61

tapply(results$Vmax, results$method, sd)

##          Direct      Eadie-Hoftzee      Haynes-Woolf Lineweaver-Burke
##           5.933           7.914           6.345           382.611

tapply(results$Km, results$method, mean)

##          Direct      Eadie-Hoftzee      Haynes-Woolf Lineweaver-Burke
##           51.87           45.26           52.79           -46.02

tapply(results$Km, results$method, sd)

##          Direct      Eadie-Hoftzee      Haynes-Woolf Lineweaver-Burke
##           7.993           7.410           8.804           227.160
```

From these results we can see that the Haynes-Woolf plot gives the closest results to the real data, very similar to the direct fitting. Eadie-Hoftzee is close. Lineweaver-Burke should not be used.