# Fitting Mathematical Models to Biological Data using Non-Linear Least-Squares (NLLS) 

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## Outline

- Why NLLS?
- The NLLS fitting method
- Practicals (in R) overview

Why NLLS?

## LINEAR MODELS

- These are all good Linear Models (really?!):



- The data can be modelled (aka "a mathematical model fitted to them") as a linear combination of variables and coefficients
- Easily fitted using Ordinary Least Squares (OLS)
- Linear models can include curved responses (e.g. Polynomial regression)


## WHAT MAKES A MODEL NON-LINEAR?

- OLS can be used to fit (model) equations that are intrinsically linear, e.g.,
- Straight line: $y_{i}=\beta_{0}+\beta_{1} x_{i}+\varepsilon_{i}$
- Polynomial (quadratic): $y_{i}=\beta_{0}+\beta_{1} x_{i}+\beta_{2} x_{i}^{2}+\varepsilon_{i}$
- Another quadratic: $y_{i}=e^{\beta_{0}}+\beta_{1} x_{i}+\beta_{2} x_{i}^{2}+\varepsilon_{i}$
- What is intrinsic linearity? - the equation of the model to be fitted should be linear in the parameters (the $\beta^{\prime}$ s)
- Some non-linear models:
- $y_{i}=\beta_{0} x_{i}^{\beta_{1}}+\varepsilon_{i}$
- $y_{i}=\beta_{0}+\beta_{1} x_{i}^{\beta_{2}}+\varepsilon_{i}$
- $y_{i}=\beta_{0} e^{\beta_{2} x_{i}}+\varepsilon_{i}$
- $y_{i}=\frac{\beta_{0} x_{i}}{\beta_{1}+x_{i}}+\varepsilon_{i}$

In all of these, at least one parameter (a $\beta$ ) is non-linear (e.g., $x_{i}^{\beta_{2}}$, $e^{\beta_{2} x_{i}}$, etc.)

## The Least-Squares Solution

## Recall what the Least Squares method does:

- Consider data on a response variable $y$, a predictor (independent) variable $x$, and $n$ observations.
- Say we want to fit a model to these data: $f\left(x_{i}, \boldsymbol{\beta}\right)+\varepsilon_{i}$ ( $\boldsymbol{\beta}=\left(\beta_{0}, \beta_{1}, \ldots, \beta_{k}\right)$ are the model's $k+1$ parameters $)$
- An example of $f\left(x_{i}, \boldsymbol{\beta}\right)+\varepsilon_{i}$ could be: $y_{i}=\beta_{0}+\beta_{1} x_{i}+\varepsilon_{i}$ (linear regression)
- The objective of any least squares method is to find estimates of values of the parameters $\left(\hat{\beta}_{j}\right)$ that minimize the sum $(S)$ of squared residuals $\left(r_{i}\right)$ (AKA RSS):

$$
\mathrm{RSS}=S=\sum_{i=1}^{n}\left[y_{i}-f\left(x_{i}, \boldsymbol{\beta}\right)\right]^{2}=\sum_{i=1}^{n} r_{i}^{2}
$$

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- Let's picture this using a simple (OLS) example; fitting the model $y_{i}=\beta_{0}+\beta_{1} x_{i}+\varepsilon_{i} \ldots$




























































































## If THE MODEL IS LINEAR, THE LEAST-SQUARE SOLUTION IS EXACT



$$
\begin{aligned}
y_{i} & =\beta_{0}+\beta_{1} x_{i}+\varepsilon_{i} \\
9.50 & =5+4 \times 1+0.50 \\
11.00 & =5+4 \times 2-2.00 \\
19.58 & =5+4 \times 3+2.58 \\
20.00 & =5+4 \times 4-1.00
\end{aligned}
$$

The least squares solution here is:
$\beta_{0}=5 ; \beta_{1}=4$

- This system of (linear) equations can be compactly represented (and solved using matrix algebra) as $\mathbf{Y}=\mathbf{X} \boldsymbol{\beta}+\boldsymbol{\varepsilon}$


## INTRINSIC NON-LINEARITY MAKES LEAST-AQUARES MODEL FITTING DIFFICULT

- In an intrinsically non-linear model such as $y_{i}=\beta_{0} e^{\beta_{2} x_{i}}+\varepsilon_{i}$, the nice trick of solving $\mathbf{Y}=\mathbf{X} \boldsymbol{\beta}+\varepsilon$ exactly is impossible

Linear Least-Squares Minimization


Non-Linear Least-Squares Minimization


## OK, FINE, WHY WOULD I EVER NEED NLLS?

- Many observations in biology are just not well-fitted by a linear model
- That is, the underlying biological phenomena/phenomenon are not well-described by a linear equation
- Examples:
- Michaelis-Menten biochemical (reaction) kinetics
- Allometric growth
- Responses of metabolic rates to changing temperature
- Consumer-Resource (e.g., predator-prey) functional responses
- Individual growth
- Population growth
- Time-series data (e.g., fitting a sinusoidal function)
- Can you think of some examples?


## Non-Linear Model Example: Temperature AND METABOLISM


$B=B_{0} e^{-\frac{E}{k T}} f\left(T, T_{p k}, E_{D}\right)$
$T$ = temperature (K)
$k=$ Boltzmann constant $\left(\mathrm{eV} \mathrm{K}^{-1}\right)$
$E=$ Activation energy (eV)
$T_{p k}=$ Temperature of peak performance
$E_{D}=$ Deactivation energy (eV)
(J H van't Hoff 1884, S Arrhenius 1889)

## The NLLS Fitting Method

## THE NLLS METHOD: OVERVIEW

- OK, so we cannot find an exact, simple solution to the least-squares problem for non-linear models
- But we can use a computer to find a approximate but close-to-optimal least-squares solution as follows:
- Choose starting (initial values for the parameters we want to estimate ( $\beta_{j}$ 's)
- Then, adjust the parameters iteratively (using a specific "algorithm" that is better than searching randomly) such that the RSS is gradually decreased
- Eventually, if it all goes well, a combination of $\beta_{j}$ 's that is very close to the desired solution (where the RSS is approximately minimized) can be found


## THE NLLS FITTING / OPTIMIZATION PROCESS

Linear Least-Squares Minimization


Non-Linear Least-Squares Minimization


## THE NLLS FITTING / OPTIMIZATION PROCESS

The general procedure / algorithm is:
(1) Start with an initial value for each parameter in the model
(2) Generate the curve defined by the initial values
(3) Calculate the residual sum-of-squares (RSS)
(9) Adjust the parameters to make the curve come closer to the data points. This the tricky part - more on this in the next slide
(5) Adjust the parameters again so that the curve comes even closer to the points (RSS decreases)
(6) Repeat 4-5
(3) Stop simulations when the adjustments make virtually no difference to the RSS

## NLLS FITTING / OPTIMIZATION ALGORITHMS

The tricky part - adjust parameters to make curve come closer to the data points (step 4) - has two main algorithms that are generally used:

- The Gauss-Newton algorithm is often used, but doesn't work very well if the model to be fitted is mathematically complicated (the parameter search "landscape" is difficult) and the starting values for parameters are far-off-optimal
- The Levenberg-Marquardt algorithm switches between Gauss-Newton and "gradient descent" and is more robust against starting values that are far-off-optimal and is more reliable in most scenarios.


## NLLS FITS - ASSESSMENT AND REPORTING

- Once the NLLS fitting is done, you need to get the goodness of fit measures
- First, of course, examine the fits visually
- Report the goodness-fit results:
- Sums of deviations of the data points from the final model fit (final RSS)
- Estimated coefficients
- For each coefficient, standard error (can be used for CI's), t-statistic and corresponding (two-tailed) p-value
- You will learn to calculate all these in the practicals
- You may also want to compare and select between multiple competing models
- Unlike in Linear Models, $\mathrm{R}^{2}$ values should not be used to interpret the quality of a NLLS fit (more on this in the practicals).


## NLLS Assumptions

NLLS-regression has all the assumptions of OLS-regression:

- No (in practice, minimal) measurement error in explanatory variable ( $x$-axis variable)
- Data have constant normal variance - errors in the $y$-axis are homogeneously distributed over the $x$-axis range
- The measurement/observation errors are Normally distributed (Gaussian)
- What if the errors are not normal? - Interpret results cautiously, and use Maximum Likelihood or Bayesian fitting methods instead


## Practicals Overview

## NLLS Fitting Practicals

- We will use R
- For fitting simple non-linear models, the nl s function in R is sufficient
- It uses the Gauss-Newton algorithm by default
- The command is nls()
- It is part of the stats base package (so no extra installation and loading of package necessary)
- For fitting complex non-linear models the Levenberg-Marquardt (LM) algorithm is better
- The command is nlsLM()
- It is available through the the minpack.lm package http://cran.r-project.org/web/packages/minpack.lm
- It offers additional features like the ability to "bound" parameters to realistic values


## NLLS Fitting Practicals

- We will start with NLLS fitting of the Michaelis-Menten model of biochemical reaction kinetics:

$$
V=\frac{V_{\max }[S]}{K_{m}+[S]}
$$

- $S=$ Substrate density
- $V_{\max }=$ Maximum reaction rate (at saturating substrate concentration)
- $K_{M}=$ Half-saturation constant; the $S$ at which reaction rate reaches half of possible maximum $\left(=\frac{1}{2} V_{\max }\right)$


Substrate abundance (S)

- You will use NLLS fitting to obtain estimates of $V_{\max }$ and $K_{M}$
- Note that $V_{\max } \leq 0$ and $K_{M} \leq 0$ are physically impossible (useful fir picking starting values)


## READINGS

- Motulsky, Harvey, and Arthur Christopoulos. Fitting models to biological data using linear and nonlinear regression: a practical guide to curve fitting. OUP USA, 2004.

