## Supporting slides

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### Section 1

NLME fitting algorithms

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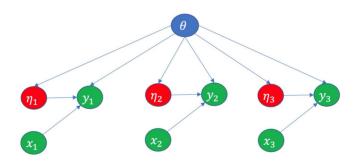
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- $x_i$ : covariates of subject i

•  $p(y=y_i \mid \theta=\theta, \eta=\eta_i, x=x_i) = p(y_i \mid \theta, \eta_i, x_i)$ : likelihood of  $(\theta, \eta_i)$  given subject i's data  $(x_i, y_i)$ . Also known as the conditional probability of  $y_i$  given  $\theta, \eta_i, x_i$ . Or just the conditional likelihood of  $\eta_i$  given  $\theta$  and  $(x_i, y_i)$ .

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- $p(y=y_i \mid \theta=\theta, x=x_i) = p(y_i \mid \theta, x_i) = \int p(y_i \mid \theta, \eta_i, x_i) \cdot p(\eta_i \mid \theta) \, d\eta_i$ : marginal likelihood of  $\theta$  given subject i's data  $y_i$ .



```
@model begin
                  @param begin
                    θ ∈ VectorDomain(4, lower = zeros(4))
                    Σ ∈ RealDomain(lower = 0.0)
                    a E RealDomain(lower = 0.0, upper = 1.0)
                  @random begin
      \eta_i \mid \theta
                    η ~ MvNormal(Ω)
                  @covariates sex wt etn
                  @pre begin
                    01 := 0[1]
                    CL = \theta[2] * ((wt / 70)^0.75) * (\theta[4]^sex) *
                    Vc = \theta[3] * exp(n[2])
                  end
y_i | \theta, \eta_i, x_i
                  @dynamics begin
                    Depot' = -Ka * Depot
                    Central' = Ka * Depot - (CL / Vc) * Central
                    Res' = Depot - Central
                  end
                  @derived begin
                    conc = @, Central / Vc
                    dv ~ @. Normal(conc. conc * Σ)
                    T max = maximum(t)
                  end
                  @observed begin
                    obs cmax = maximum(dv)
                  end
                end
```

## Marginal likelihood maximization

$$\begin{split} \theta^* &= \arg \max_{\theta} \prod_{i=1}^N p(y_i \mid \theta, x_i) \\ &= \arg \max_{\theta} \prod_{i=1}^N \int p(y_i \mid \theta, \eta_i, x_i) \cdot p(\eta_i \mid \theta) \, d\eta_i \\ \text{EBE}_i &= \eta_i^* = \arg \max_{\eta_i} \Big( p(y_i \mid \theta = \theta^*, \eta_i, x_i) \cdot p(\eta_i \mid \theta = \theta^*) \Big) \end{split}$$

## General Laplace method

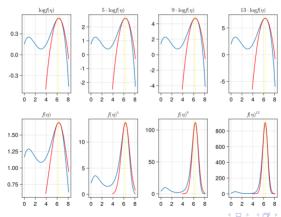
$$\int f(\eta) \, d\eta \approx f(\eta^*) \sqrt{(2\pi)^m/|-H|}$$

- ullet f: a positive scalar-valued function of  $\eta$
- $\eta$ : vector of m integration variables
- $\eta^*$  : global maximizer of  $\log f$  ,  $\frac{d\log f}{d\eta}(\eta^*)=0$
- H: second derivative matrix of  $\log f$  wrt  $\eta$  at  $\eta^*$ , must be negative definite at  $\eta = \eta^*$
- $\bullet \mid -H \mid$ : determinant of -H



## NLME Laplace method

Laplace uses a second order Taylor series approximation of  $\log f$  at  $\eta^*$ .



## NLME Laplace method

The mode is the same after multiplying  $\log f$  by a scalar. The ratio of peaks is not the same. Consider the 2 local maximizers  $\eta_1$  (lower peak) and  $\eta_2$  (higher peak).

$$\begin{split} c &= \log f(\eta_2) - \log f(\eta_1) \\ n \cdot c &= n \cdot (\log f(\eta_2) - \log f(\eta_1)) \\ e^{n \cdot c} &= f(\eta_2)^n / f(\eta_1)^n \end{split}$$

#### Summary

Approximation error of  $n \cdot \log f$  away from the mode  $\eta^*$  is not significant as n increases.



## NLME Laplace method

- $\bullet$  There are N functions  $\{f_i: i=1\dots N\}$  to be integrated, one for each subject
- $\bullet \ f_i(\eta_i) = p(y_i \mid \theta, \eta_i, x_i) \cdot p(\eta_i \mid \theta)$
- Laplace method

$$\int f_i(\eta_i)\,d\eta_i = f_i(\mathsf{EBE}_i)\sqrt{(2\pi)^m/|-H_i|}$$

•  $H_i$ : second derivative matrix of  $\log f_i$  wrt  $\eta_i$  at  $\mathsf{EBE}_i$ , must be negative definite at  $\eta_i = \mathsf{EBE}_i$ 



 $\ensuremath{\mathsf{FOCE}}(\ensuremath{\mathsf{I}})$  approximates the Hessian  $H_i$  for each subject i. Assume the following:

$$\log f_i(\eta_i) = \log p(y_i \mid \theta, \eta_i, x_i) + \log p(\eta_i \mid \theta)$$
$$= L_i(g_i(\eta_i)) + \log p(\eta_i \mid \theta)$$

#### where:

- $g_i$  returns the vector of IPREDs  $\mu_i$  and the residual standard deviations  $\sigma_i$  (constant in the additive error model case), at all observed time points, and
- ullet  $L_i$  is the log probability of  $y_i$  given  $\mu_i$  and  $\sigma_i$



 $g_i$  is usually the most expensive component of  $\log f_i$ , because it often involves solving a differential equation. So let's approximate it!

#### First order Taylor series expansion

FO

$$g_i(\eta_i) = g_i(0) + \frac{dg_i}{d\eta_i}(0) \cdot \eta_i$$

FOCE(I)

$$g_i(\eta_i) = g_i(\mathsf{EBE}_i) + \frac{dg_i}{d\eta_i}(\mathsf{EBE}_i) \cdot (\eta_i - \mathsf{EBE}_i)$$

- FOCE(I) ensures that the approximation error in  $g_i$  (and  $\log f_i$  by extension) is low in the proximity of  $\mathsf{EBE}_i$ .
- FO does not ensure that so it only works well if:
  - EBE<sub>i</sub> is not far from 0, or
  - $g_i$  is close to linear in the interval  $[0, \mathsf{EBE}_i]$ .
- FO requires a correction term in the Laplace method because the gradient of  $\log f_i$  wrt  $\eta_i$  at  $\eta_i=0$  is not 0.

#### Chain rule for Hessians

$$(L_i \cdot g_i)''(\eta_i) = \frac{dg_i}{d\eta_i}^T \cdot \frac{\partial^2 L_i}{\partial g_i \cdot \partial g_i^T} \cdot \frac{dg_i}{d\eta_i} + \sum_{t=1}^d \left( \frac{\partial L_i}{\partial g_{i,t}} \cdot \underbrace{\frac{\partial^2 g_{i,t}}{\partial \eta_i \cdot \partial \eta_i^T}} \right)$$

where d is twice the number of observed time points (corresponding to  $\mu_i$  and  $\sigma_i$ ) and  $g_{i,t}$  is the  $t^{th}$  component of  $g_i$ .

#### Summary

If  $g_i$  is linear in  $\eta_i$ :

$$\bullet \ \frac{\partial^2 g_{i,t}}{\partial \eta_i \cdot \partial \eta_i^T} = 0$$

• 
$$J_i = \frac{\partial g_i}{\partial \eta_i}$$
 is constant

The Hessian simplifies to:

$$(L_i \cdot g_i)''(\eta_i) = J_i^T \cdot \frac{\partial^2 L_i}{\partial q_i \cdot \partial q_i^T} \cdot J_i$$

- ullet  $J_i$  can be computed for each subject using finite difference at
  - $\eta_i = 0$  for FO, or
  - $\eta_i = \mathsf{EBE}_i$  for  $\mathsf{FOCE}(\mathsf{I})$
- For many data distributions,  $\frac{\partial^2 L_i}{\partial g_i \cdot \partial g_i^T}$  is both diagonal and has a closed form. Doesn't have to be Gaussian!

Recall

$$\begin{split} \log f_i(\eta_i) &= \log p(y_i \mid \theta, \eta_i, x_i) + \log p(\eta_i \mid \theta) \\ &= L_i(\underbrace{g_i(\eta_i)}_{\text{approx}}) + \log p(\eta_i \mid \theta) \end{split}$$

For many random effects distributions, the Hessian of  $\log p(\eta_i \mid \theta)$  wrt  $\eta_i$  has a closed form. Doesn't have to be Gaussian!

- In Pumas, FOCE is always "with interaction".
- Pumas FOCE supports a number of data distributions:
  - Continuous: Normal, LogNormal, Gamma, Exponential, Beta
  - Discrete: NegativeBinomial, Bernoulli, Binomial, Poisson, Categorical

- Pumas supports a number of random effect distributions:
  - Unbounded: Cauchy, Gumbel, Laplace, Logistic, Normal, NormalCanon, NormalInverseGaussian, PGeneralizedGaussian, TDist
  - Positive: BetaPrime, Chi, Chisq, Erlang, Exponential, Frechet, Gamma, InverseGamma, InverseGaussian, Kolmogorov, LogNormal, NoncentralChisq, Rayleigh, Weibull
  - Between 0 and 1: Beta, LogitNormal
  - Other bounded: Uniform, Arcsine, Biweight, Cosine, Epanechnikov, LogUniform, Semicircle, SymTriangularDist, Triweight

- Use FOCE if supported, otherwise use Laplace.
- Avoid FO.

### Section 2

# Diagnostics

### Standard error estimation

#### Goal

Estimate the covariance matrix of the estimator  $\theta^*$ :

$$\theta^* = \arg \max_{\theta} \prod_{i=1}^{N} p(y_i \mid \theta, x_i)$$

#### Asymptotic covariance

$$\begin{split} \theta^* &\sim \mathcal{N}(\theta_0, V) = \mathcal{N}(\theta_0, A^{-1}BA^{-1}) \\ A &= \sum_{i=1}^N \frac{\partial^2 \log p(y_i \mid \theta, x_i)}{\partial \theta \cdot \partial \theta^T}(\theta_0) \\ B &= \sum_{i=1}^N \frac{\partial \log p(y_i \mid \theta, x_i)}{\partial \theta}(\theta_0) \times \frac{\partial \log p(y_i \mid \theta, x_i)}{\partial \theta}(\theta_0)^T \end{split}$$

where  $\theta_0$  is the set of unknown true parameters.

#### Estimated covariance

$$\begin{split} & \theta^* \overset{a}{\sim} \mathcal{N}(\theta_0, \hat{V}) = \mathcal{N}(\theta_0, \hat{A}^{-1} \hat{B} \hat{A}^{-1}) \\ & \hat{A} = \sum_{i=1}^N \frac{\partial^2 \log p(y_i \mid \theta, x_i)}{\partial \theta \cdot \partial \theta^T} (\theta^*) \\ & \hat{B} = \sum_{i=1}^N \frac{\partial \log p(y_i \mid \theta, x_i)}{\partial \theta} (\theta^*) \times \frac{\partial \log p(y_i \mid \theta, x_i)}{\partial \theta} (\theta^*)^T \end{split}$$

#### Standard error estiamtes

The square root of the diagonal elements of  $\hat{V}$  are the standard error estimates of  $\theta^*$ .

## Computing $\hat{V}$

- ullet  $\hat{A}$  and  $\hat{B}$  can be approximated with finite difference
- $\log p(y_i \mid \theta, x_i)$  itself needs to be approximated with Laplace/FOCE/FO
- $\hat{A}^{-1} \cdot \hat{B} \cdot \hat{A}^{-1}$  is computed using a generalized eigenvalue problem.

## Computing $\hat{V}$

Consider the following generalized eigenvalue problem:

$$\begin{split} \hat{B} \cdot U &= \hat{A} \cdot U \cdot \Lambda \\ I &= U^T \cdot \hat{A} \cdot U \end{split}$$

where U is the matrix of generalized eigenvectors and  $\Lambda$  is the diagonal matrix of generalized eigenvalues.

## Computing $\hat{V}$

The inverse of the matrix of eigenvectors U is obtained from the following constraint on U:

$$\begin{split} (U^T \cdot \hat{A}) \cdot U &= I \\ U^{-1} &= U^T \cdot \hat{A} \end{split}$$

## Computing $\hat{V}$

The following identity is true:

$$\begin{split} \hat{B} \cdot U &= \hat{A} \cdot U \cdot \Lambda \\ \hat{A}^{-1} \cdot \hat{B} &= U \cdot \Lambda \cdot U^{-1} \\ \hat{A}^{-1} \cdot \hat{B} &= U \cdot \Lambda \cdot U^T \cdot \hat{A} \\ \hat{V} &= \hat{A}^{-1} \cdot \hat{B} \cdot \hat{A}^{-1} &= U \cdot \Lambda \cdot U^T \end{split}$$

#### Failed estimator

- If the computed  $\hat{A}$  is: a) singular, b) near singular, or c) has negative eigenvalues, the sandwich estimator will fail.
- This is a sign of poor identifiability of at least 1 parameter and/or significant numerical errors.
- $\bullet$  Even if a single (IIV) parameter is not identifiable given the data,  $\hat{A}$  will be singular.

#### Failed estimator

• Numerical errors in the finite difference or Laplace/FOCE/FO can also cause the computed approximate  $\hat{A}$  to be singular (or have small negative eigenvalues) even when the exact matrix  $\hat{A}$  may be only near singular and positive definite.

## Weighted residuals

## Visual predictive check

## Section 3

Time to event

## **Definitions**

Quantity	Formula
Instantaneous hazard	$\lambda(t) > 0$
Cumulative hazard	$\Lambda(t) = \int_0^t \lambda(t')  dt'$
Survival function: probability of survival up to time $t$	$S(t) = \exp(-\Lambda(t))$
Failure function: probability of death/failure before time $t$	F(t) = 1 - S(t)
Probability density function of time of death $t$	$f(t) = \frac{dF}{dt} = \lambda(t) \cdot exp(-\Lambda(t))$
Expected time of death ${\cal E}[t]$	$E[t] = \int_0^\infty t \cdot f(t)  dt = \int_0^\infty S(t)  dt$

## Log likelihood

The log likelihood for censored survival data is given by the following 2 formulas:

ullet For censored subjects at time t (patient survived until time t)

$$\log \operatorname{likelihood} = \log S(t) = -\Lambda(t)$$

For subjects dead at time t

$$\log \mathsf{likelihood} = \log f(t) = \log \lambda(t) - \Lambda(t)$$

