

# Ecological forecasting in R

Lecture 3: latent AR and GP models

Nicholas Clark

School of Veterinary Science, University of Queensland

0900–1200 CET Tuesday 5th September, 2023

# Workflow

Press the "o" key on your keyboard to navigate among slides

Access the [tutorial html here](#)

Download the data objects and exercise  script from the html file

Complete exercises and use Slack to ask questions

Relevant open-source materials include:

[Introduction to Generalized Additive Models with !\[\]\(a870788d6ed9b8fd294b7654a8c8526b\_img.jpg\) and mgcv](#)

[Qualitative difference between stationary and non-stationary AR\(1\)s](#)

[Statistical Rethinking 2023 - 16 - Gaussian Processes](#)

# This lecture's topics

Extrapolating splines

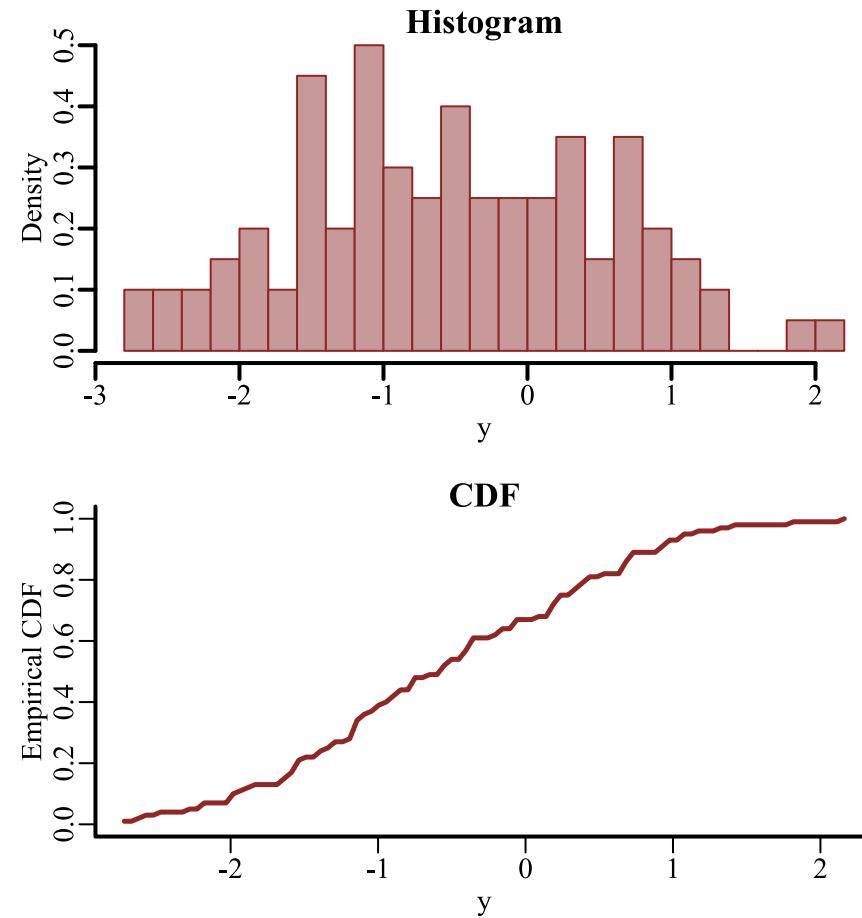
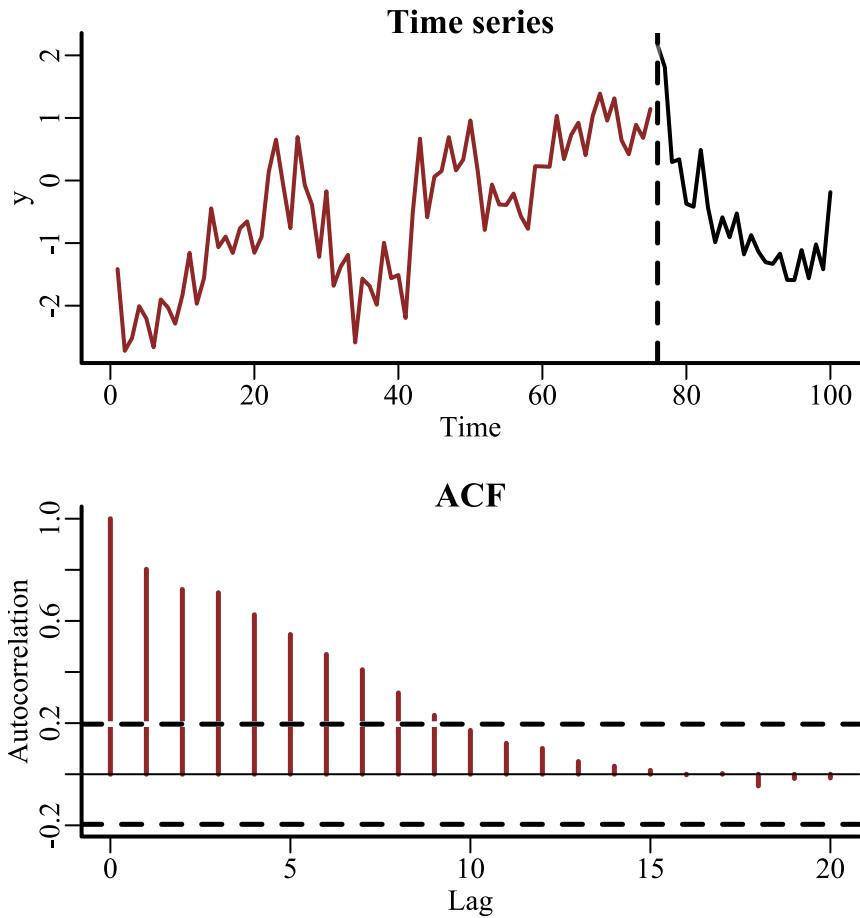
Latent autoregressive processes

Latent Gaussian Processes

Dynamic coefficient models

# Extrapolating splines

# Simulated data



# A spline of time

```
library(mvgam)
model <- mvgam(y ~
  s(time, k = 20, bs = 'bs', m = 2),
  data = data_train,
  newdata = data_test,
  family = gaussian())
```

A B-spline (`bs = 'bs'`) with `m = 2` sets the penalty on the second derivative

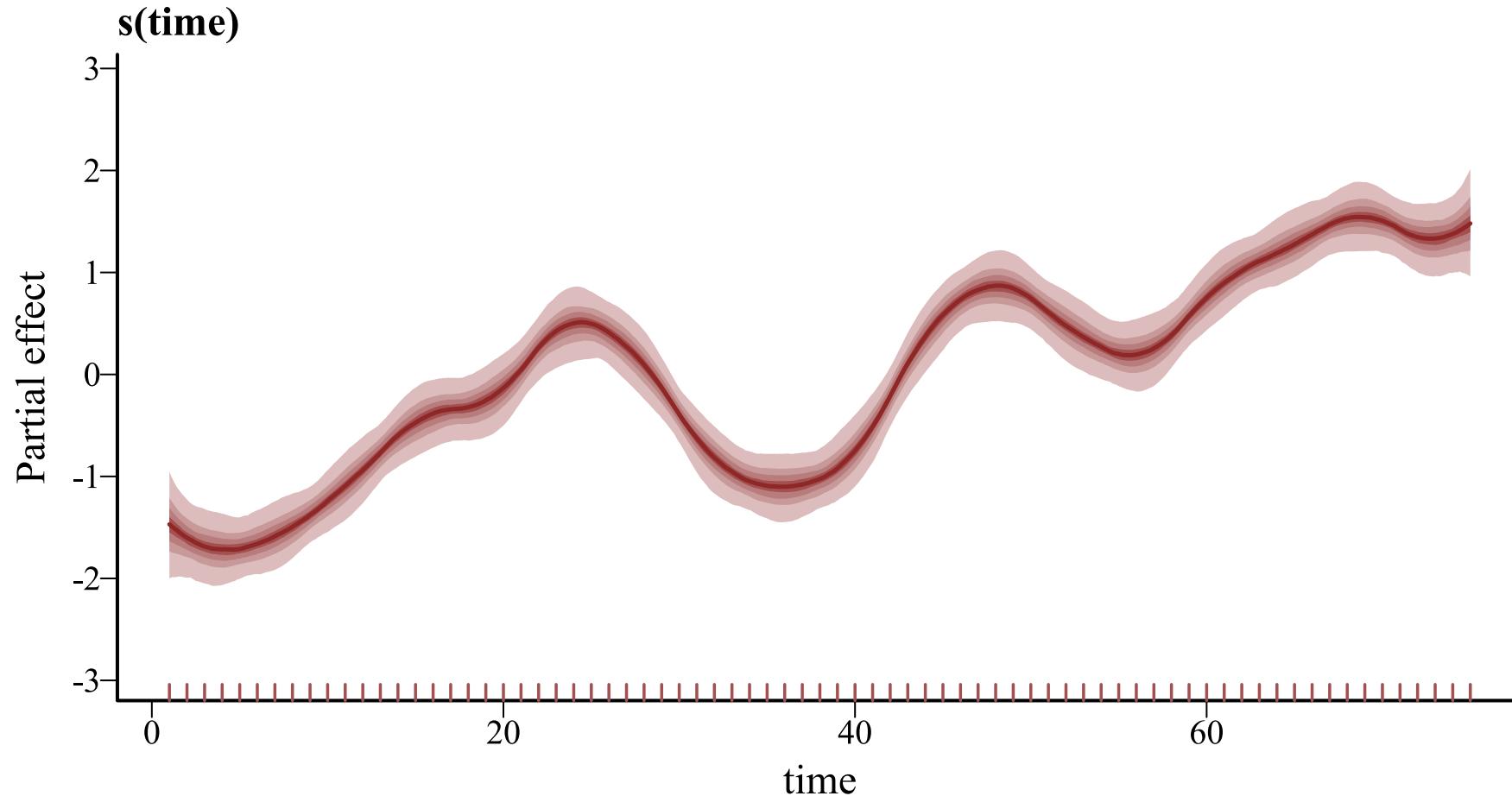
# A spline of time

```
library(mvgam)
model <- mvgam(y ~
  s(time, k = 20, bs = 'bs', m = 2),
  data = data_train,
  newdata = data_test,
  family = gaussian())
```

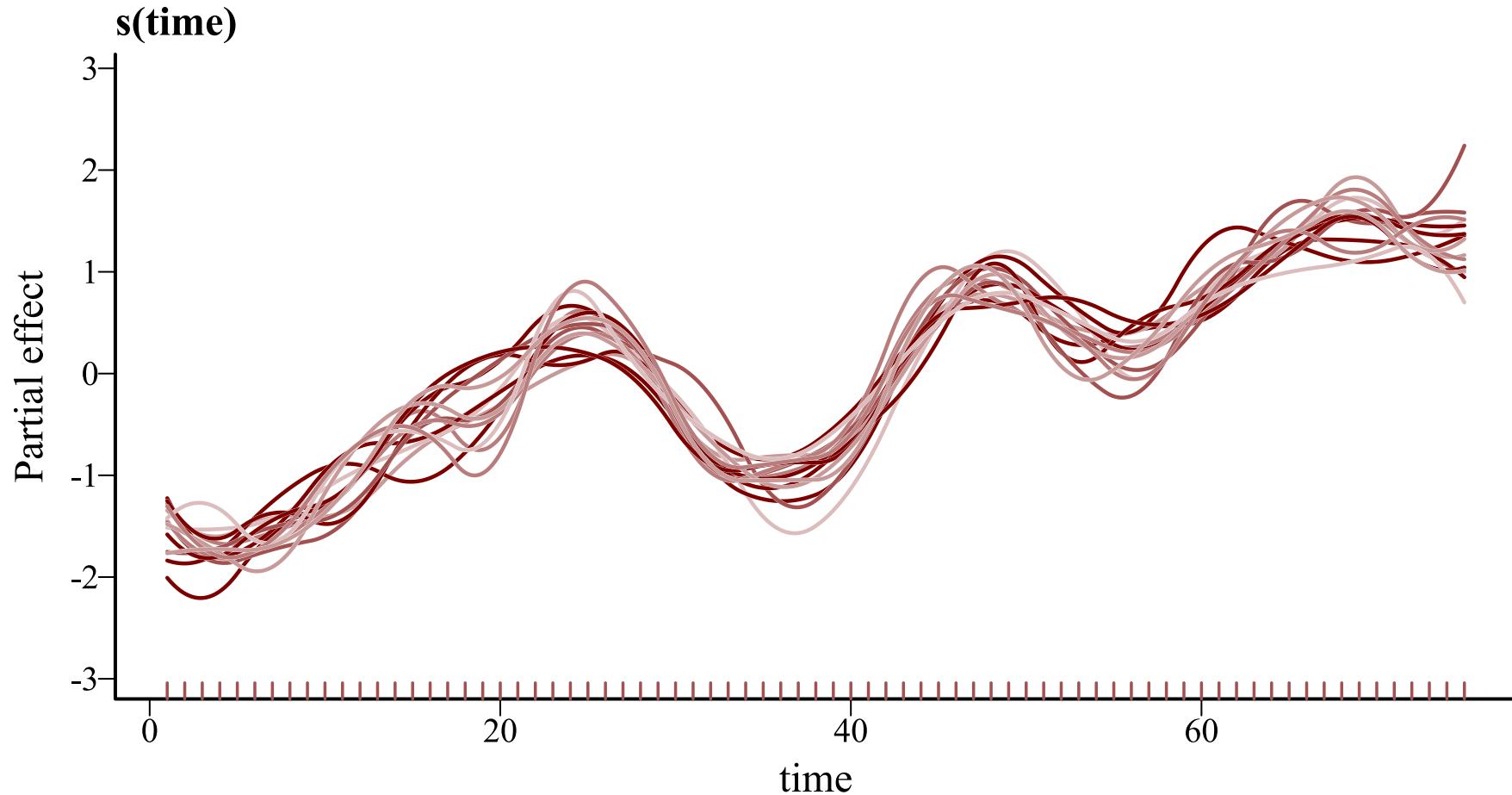
A B-spline (`bs = 'bs'`) with `m = 2` sets the penalty on the second derivative

Use `newdata` argument to generate automatic probabilistic forecasts

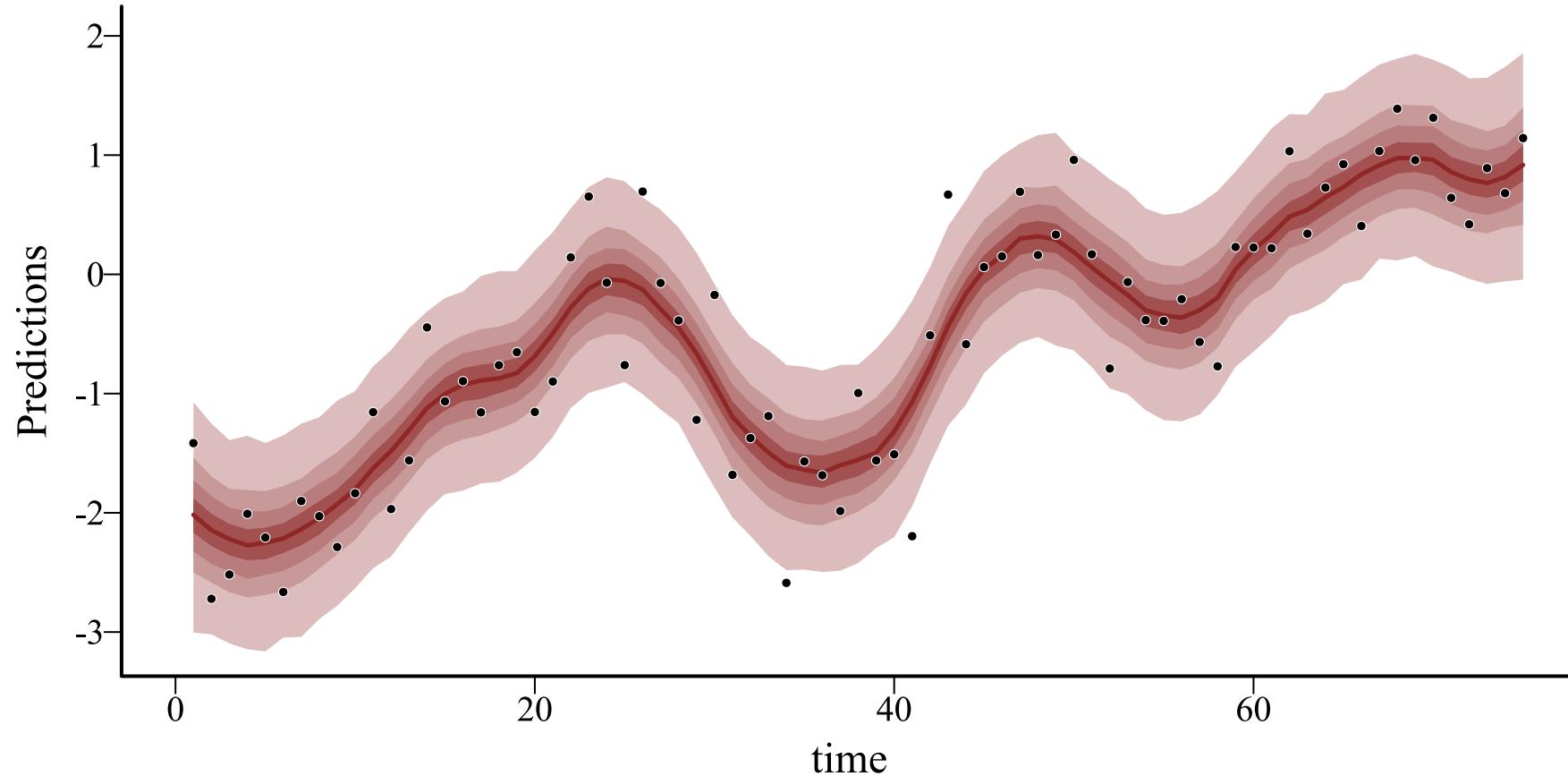
# The smooth function



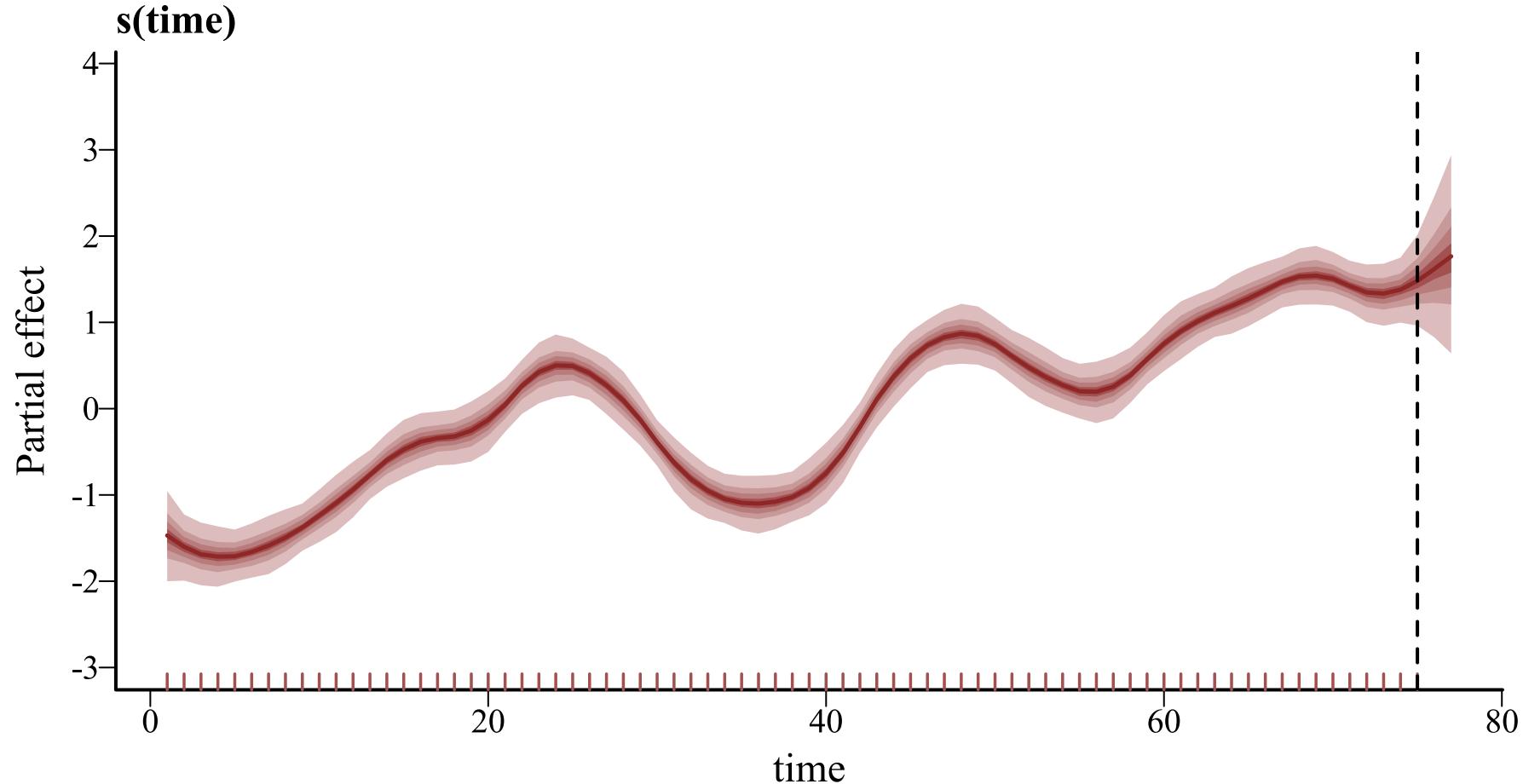
# Realizations of the function



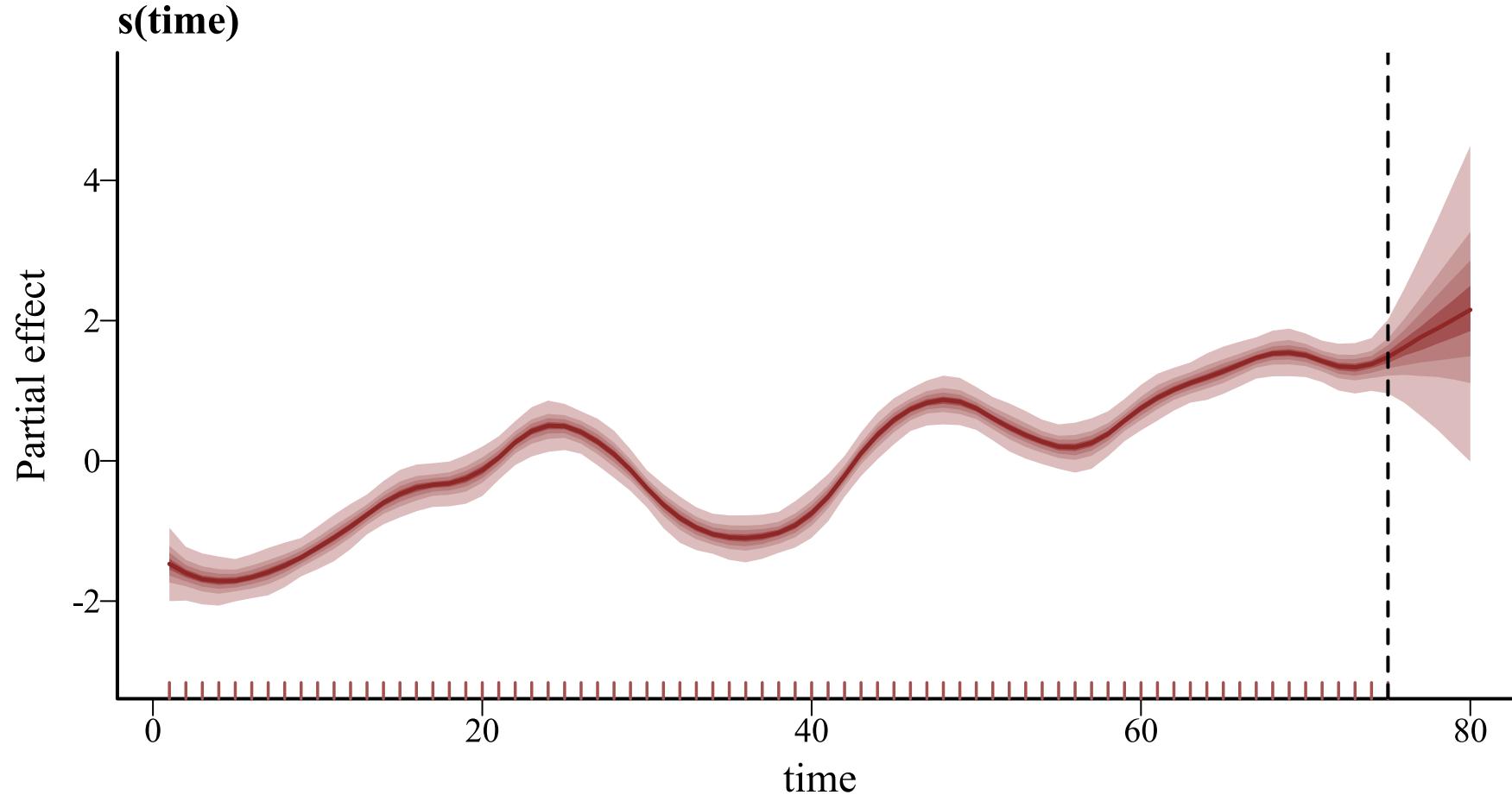
# Hindcasts 😊



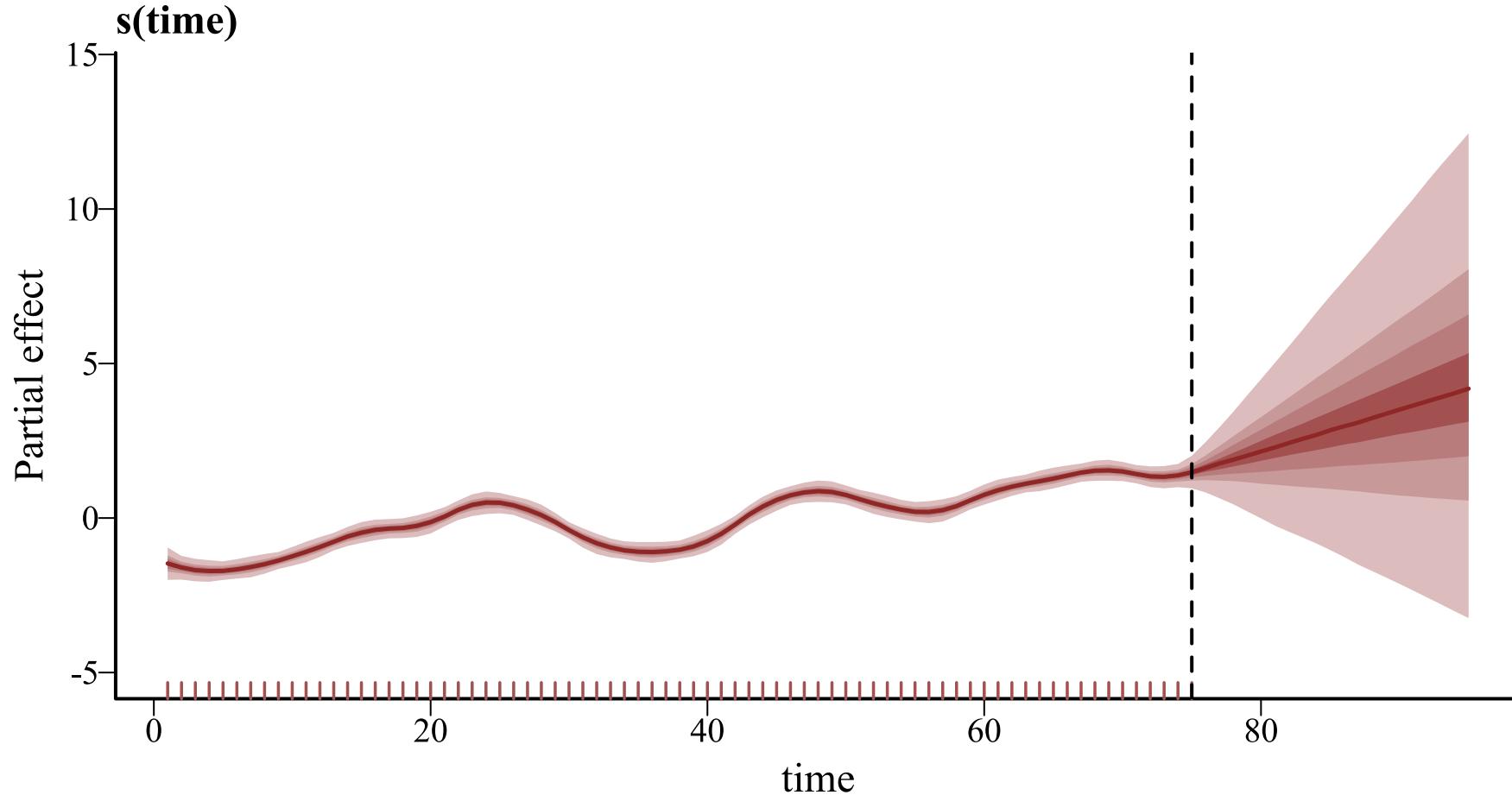
# Extrapolate 2-steps ahead 😊



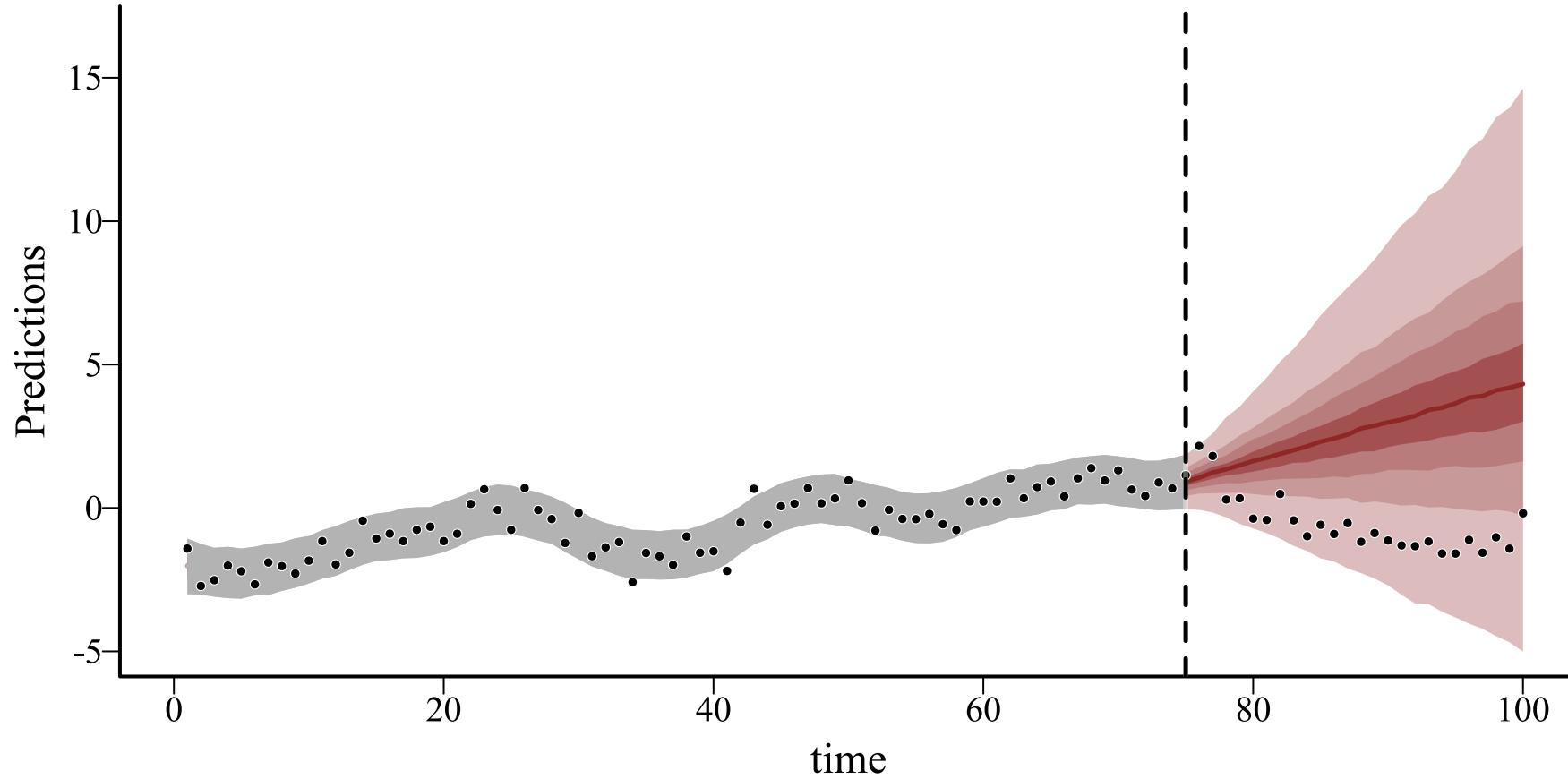
# 5-steps ahead ☹



# 20-steps ahead 😰



# Forecasts



# 2nd derivative penalty

Penalizes the overall ***curvature*** of the spline

This is default behaviour in 's `mgcv`, `brms` and `mvgam`

Provides linear extrapolations

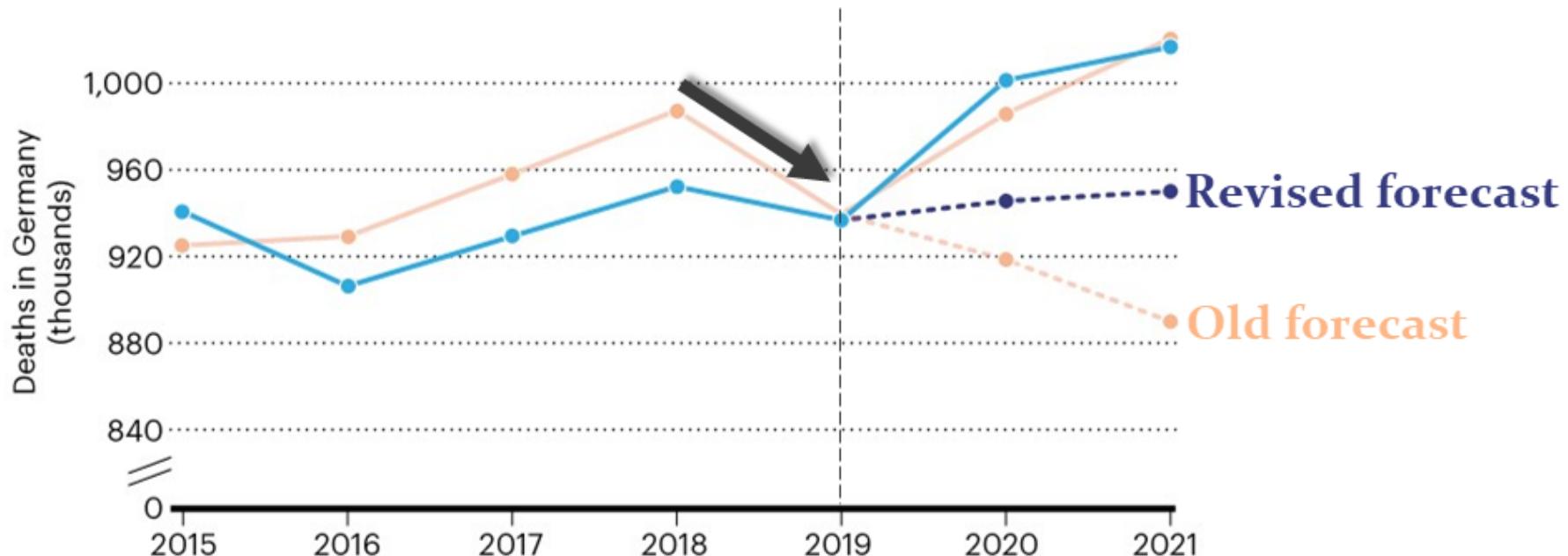
Slope remains unchanged from the last boundary of training data

Uncertainty grows but has no probabilistic understanding of time

This behaviour is widely known; ***but spline extrapolation is still commonplace***

# COVID death tolls: scientists acknowledge errors in WHO estimates

Researchers with the World Health Organization explain mistakes in high-profile mortality estimates for Germany and Sweden.

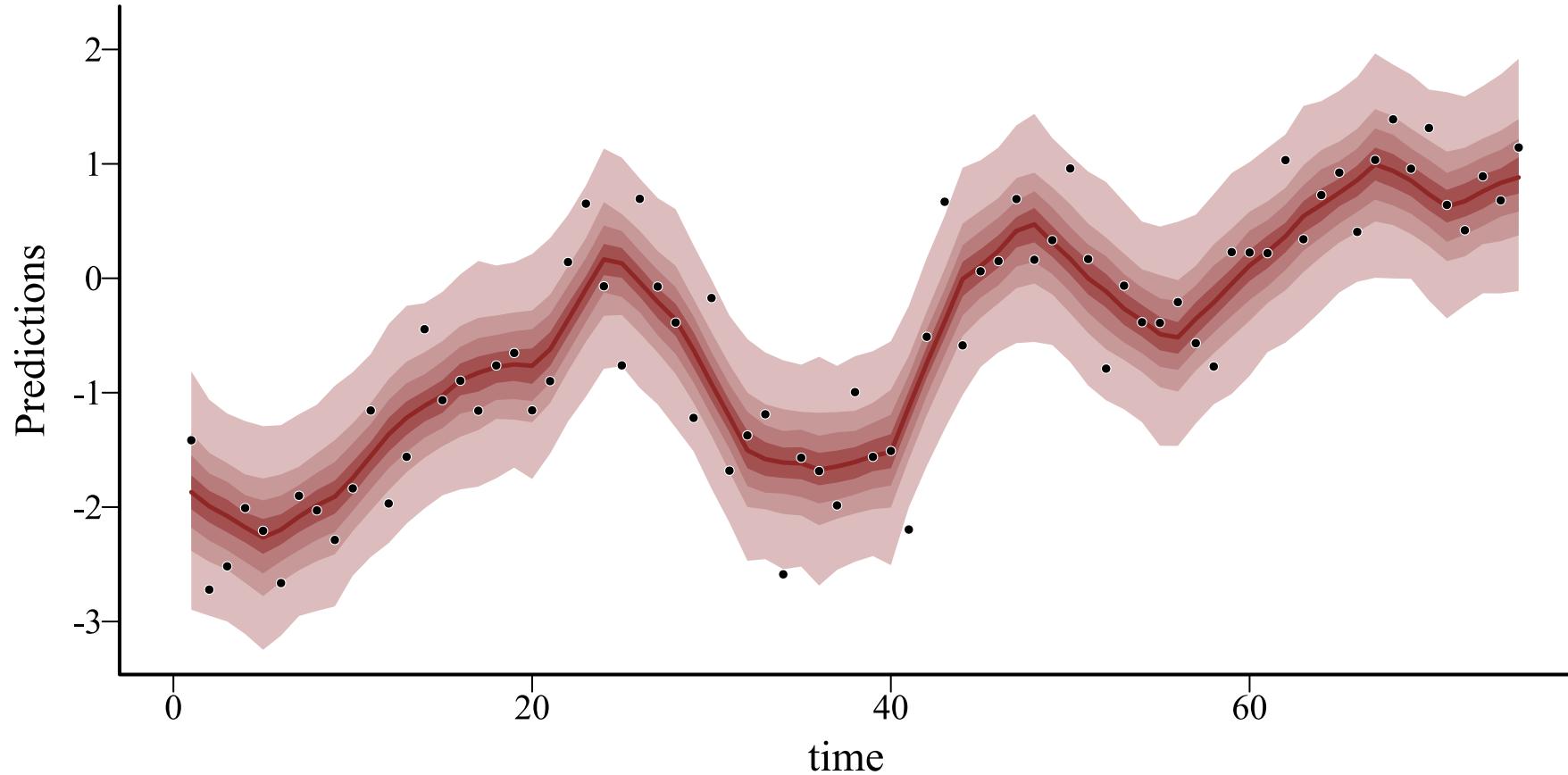


# 1st derivative penalty?

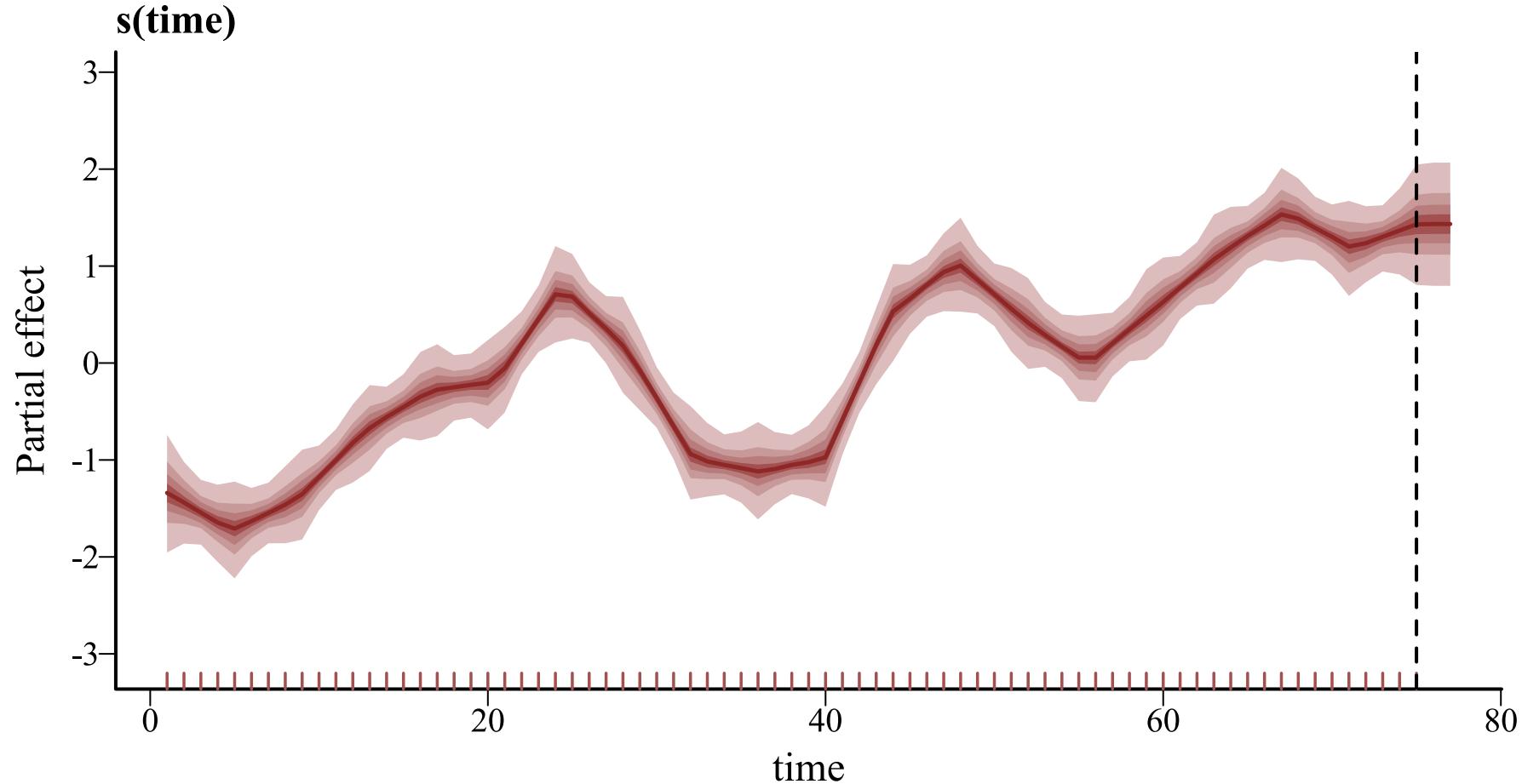
```
model ← mvgam(y ~  
    s(time, k = 20, bs = 'bs', m = 1),  
    data = data_train,  
    newdata = data_test,  
    family = gaussian())
```

Using `m = 1` sets the penalty on the first derivative

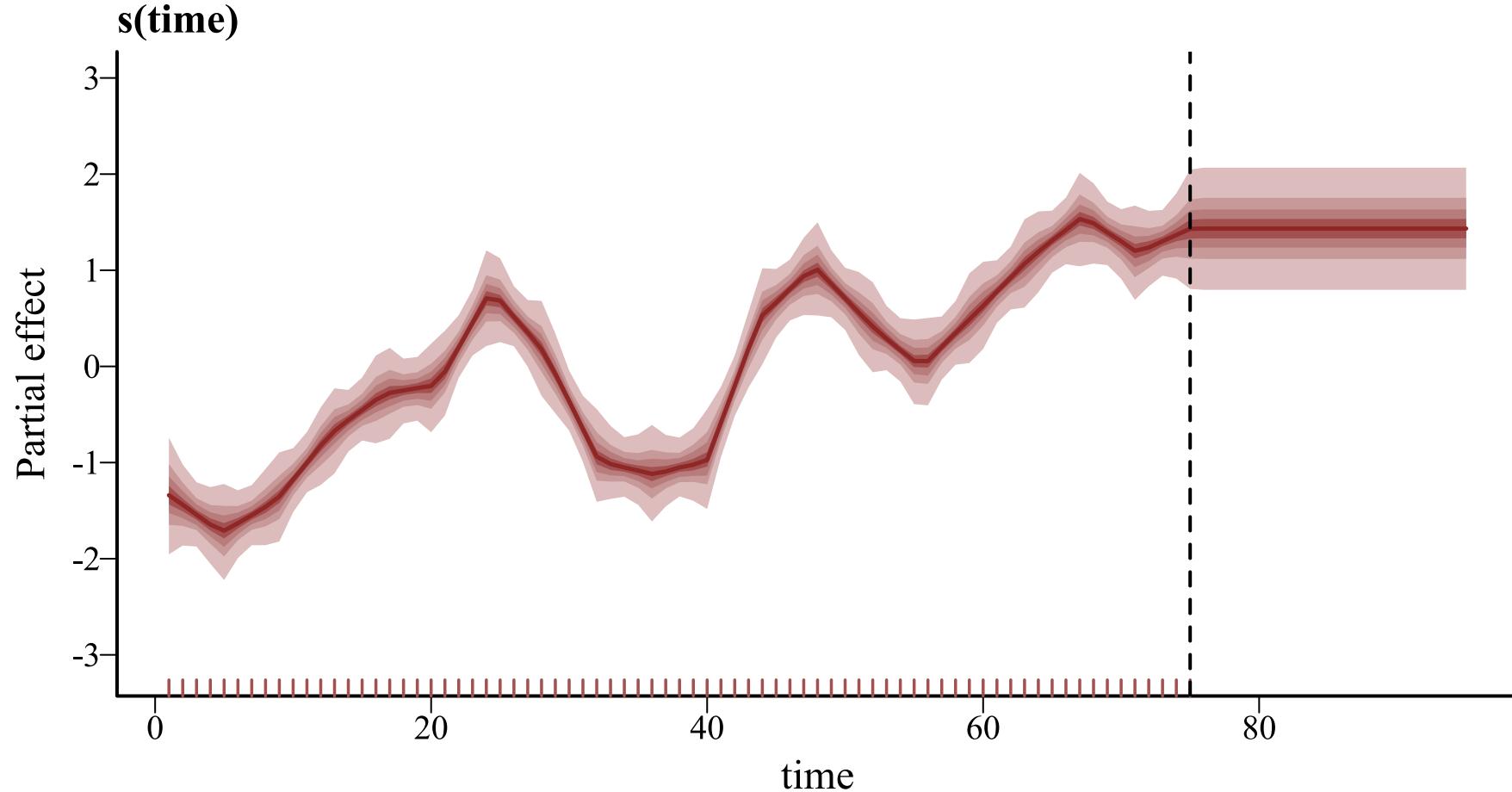
# Hindcasts 😊



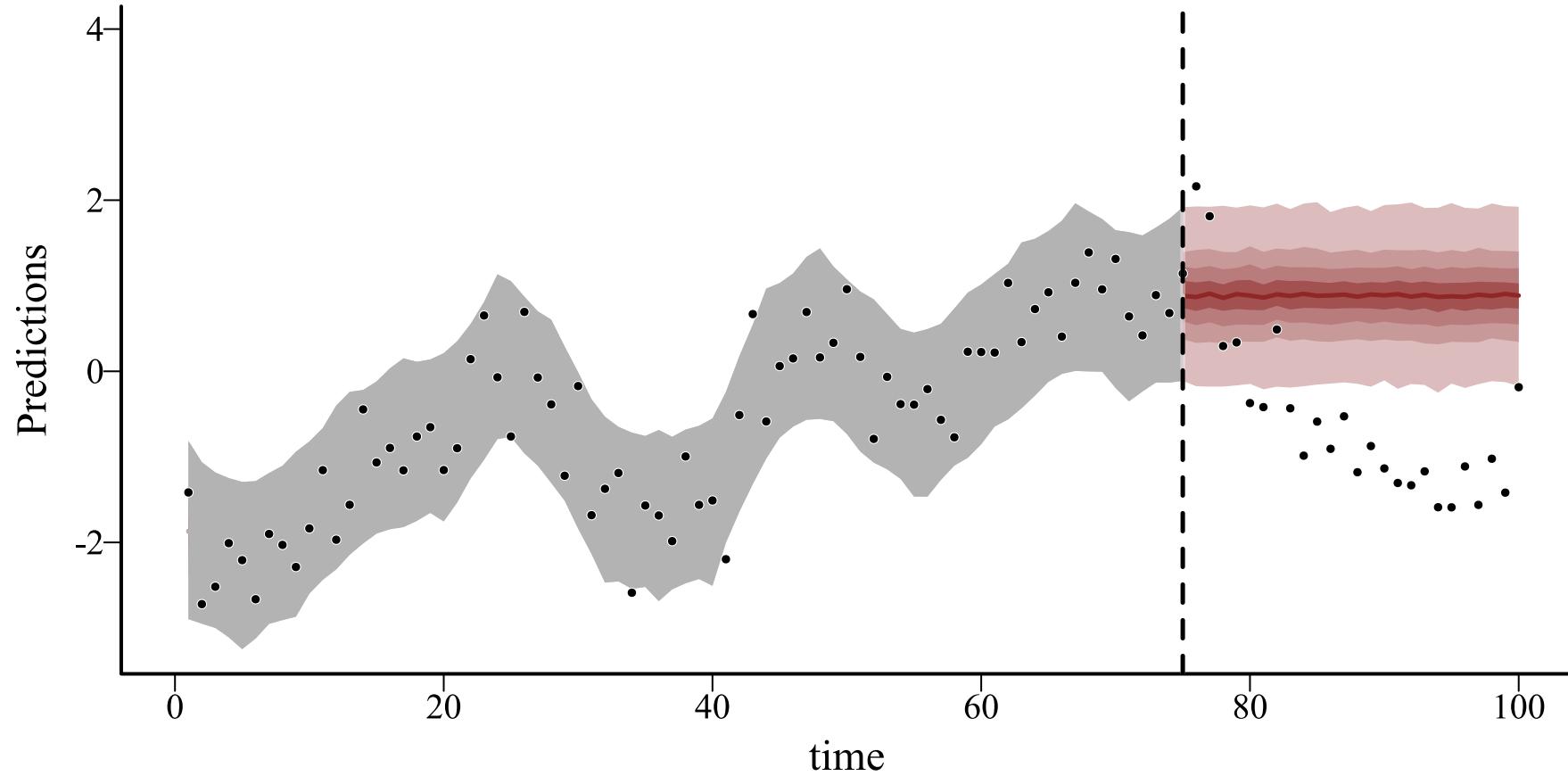
# 2-step ahead prediction 😊



# 20-steps ahead



# Forecasts



# 1st derivative penalty

Penalizes deviations from a flat function

Provides flat extrapolations

Mean remains unchanged from last boundary of the training data

Uncertainty remains unrealistically narrow

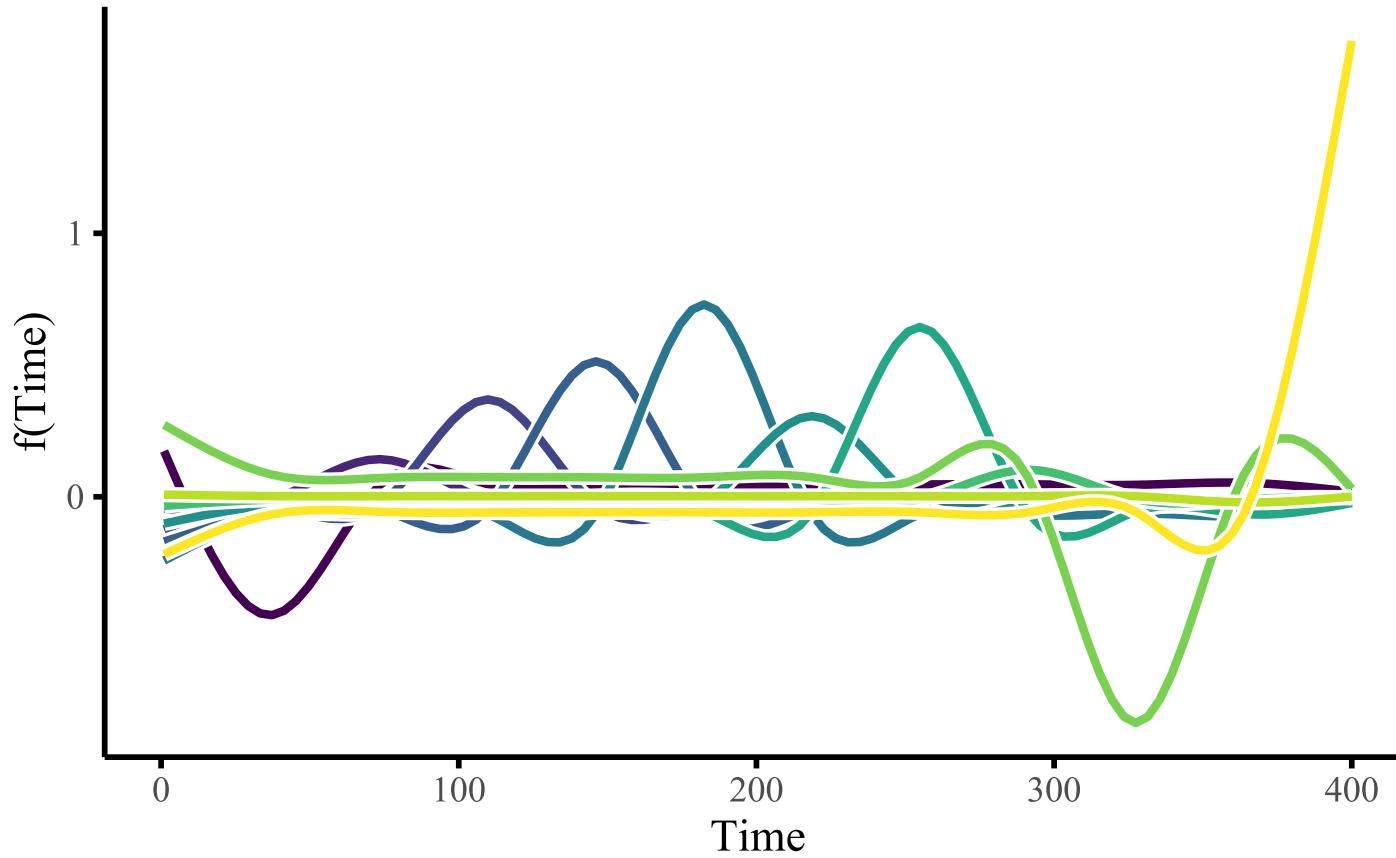
Not commonly used, though there are exceptions

**Changing penalties when using splines will impact how they extrapolate**

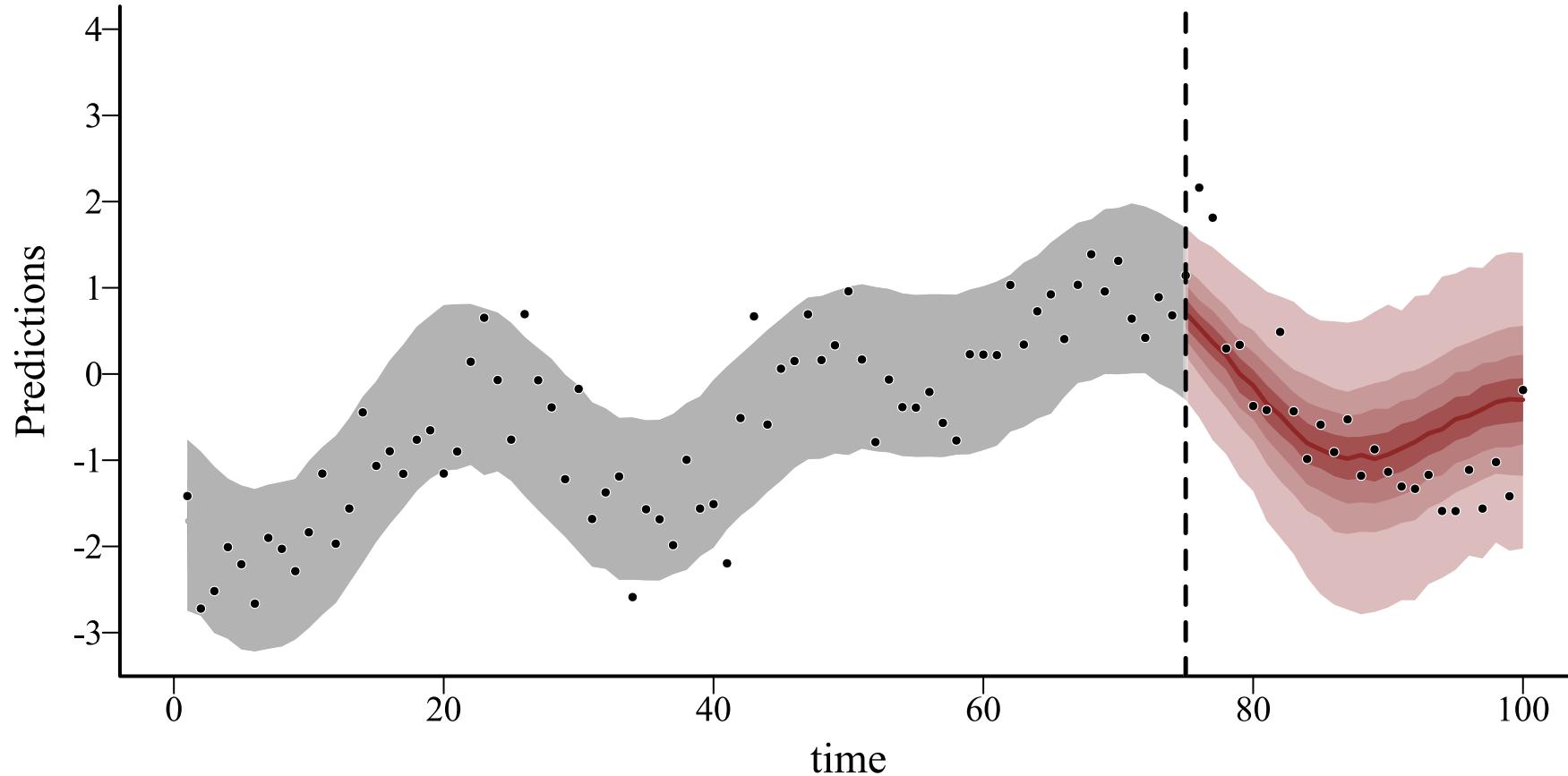
**Extrapolation also reacts *strongly* to what the spline is doing at the boundaries**

**This is because splines only have *local knowledge***

# Basis functions $\Rightarrow$ local knowledge



# We need *global knowledge*



# First, a few other pitfalls

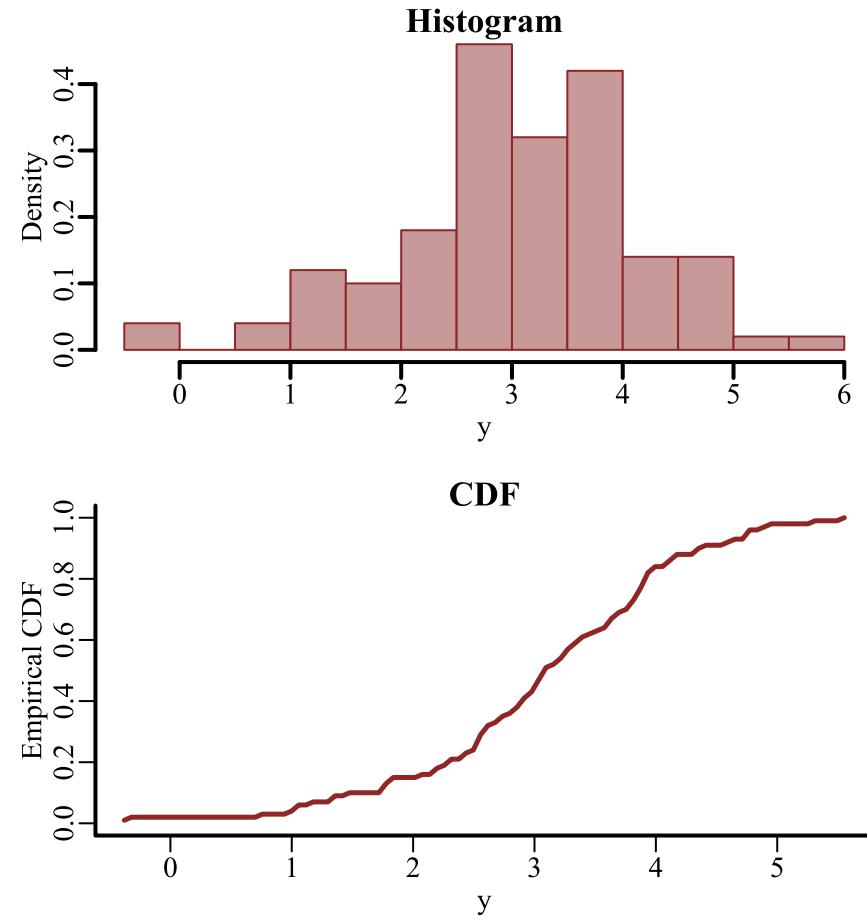
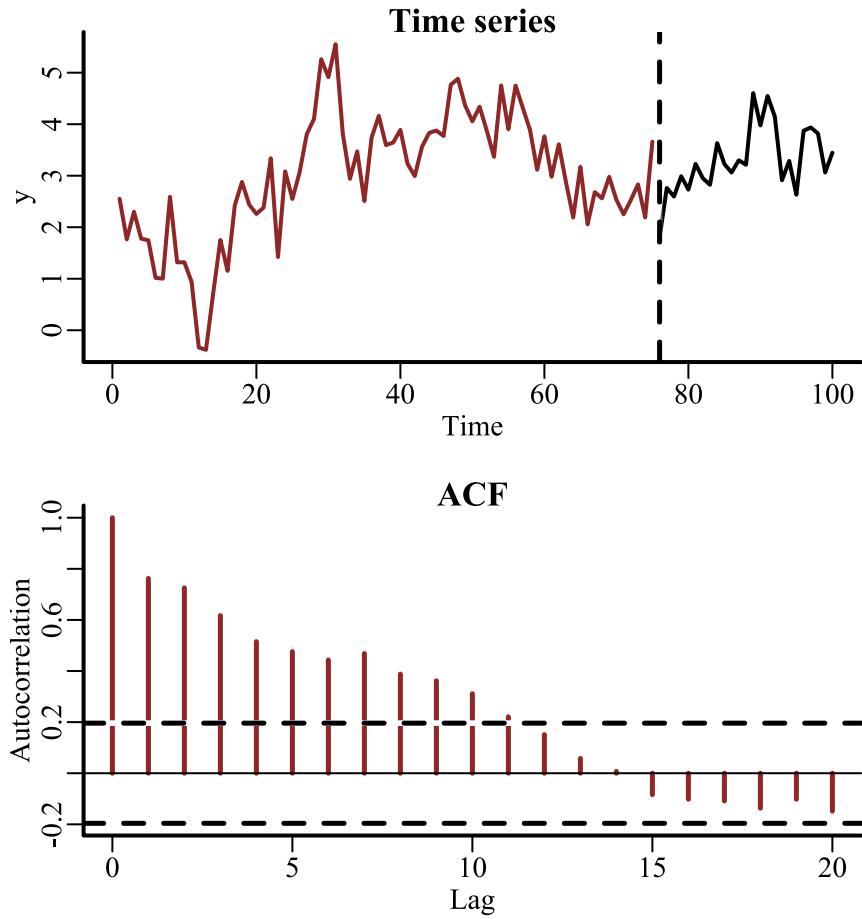
`mgcv`  has a heuristic checking function (`gam.check`) to inform whether a spline is *wiggly enough*

Can be useful to understand if your functions are complex enough to capture patterns in observed data

But can also be misleading when dealing with time series

`mvgam`  includes an underlying object of class `gam` that can be checked with `gam.check`

# Simulated data



# Restricted smooth of time

```
model <- mvgam(y ~ s(time, k = 6),  
                 family = gaussian(),  
                 data = data_train,  
                 newdata = data_test)
```

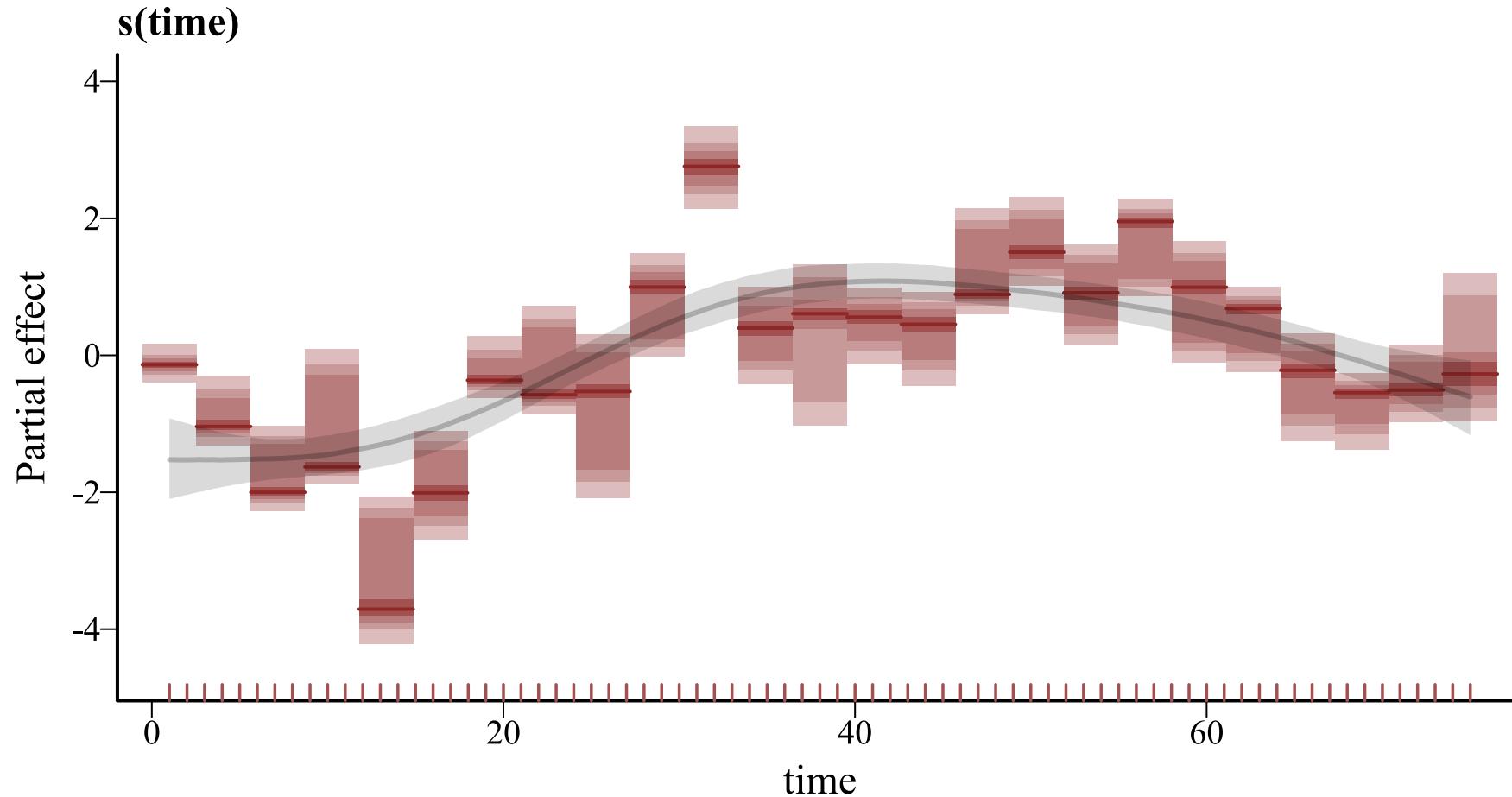
Using a thin plate spline with low maximum complexity (**k = 6**)

# Check basis complexity

```
gam.check(model$mgcv_model)
```

```
##  
## Method: REML  Optimizer: outer newton  
## full convergence after 6 iterations.  
## Gradient range [-2.516432e-07,-8.657903e-09]  
## (score 94.00124 & scale 0.590227).  
## Hessian positive definite, eigenvalue range [1.028413,36.58177].  
## Model rank =  6 / 6  
##  
## Basis dimension (k) checking results. Low p-value (k-index<1) may  
## indicate that k is too low, especially if edf is close to k'.  
##  
##          k'  edf k-index p-value  
## s(time) 5.00 4.41    0.55  <2e-16 ***  
## ---  
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
```

# Unmodelled variation



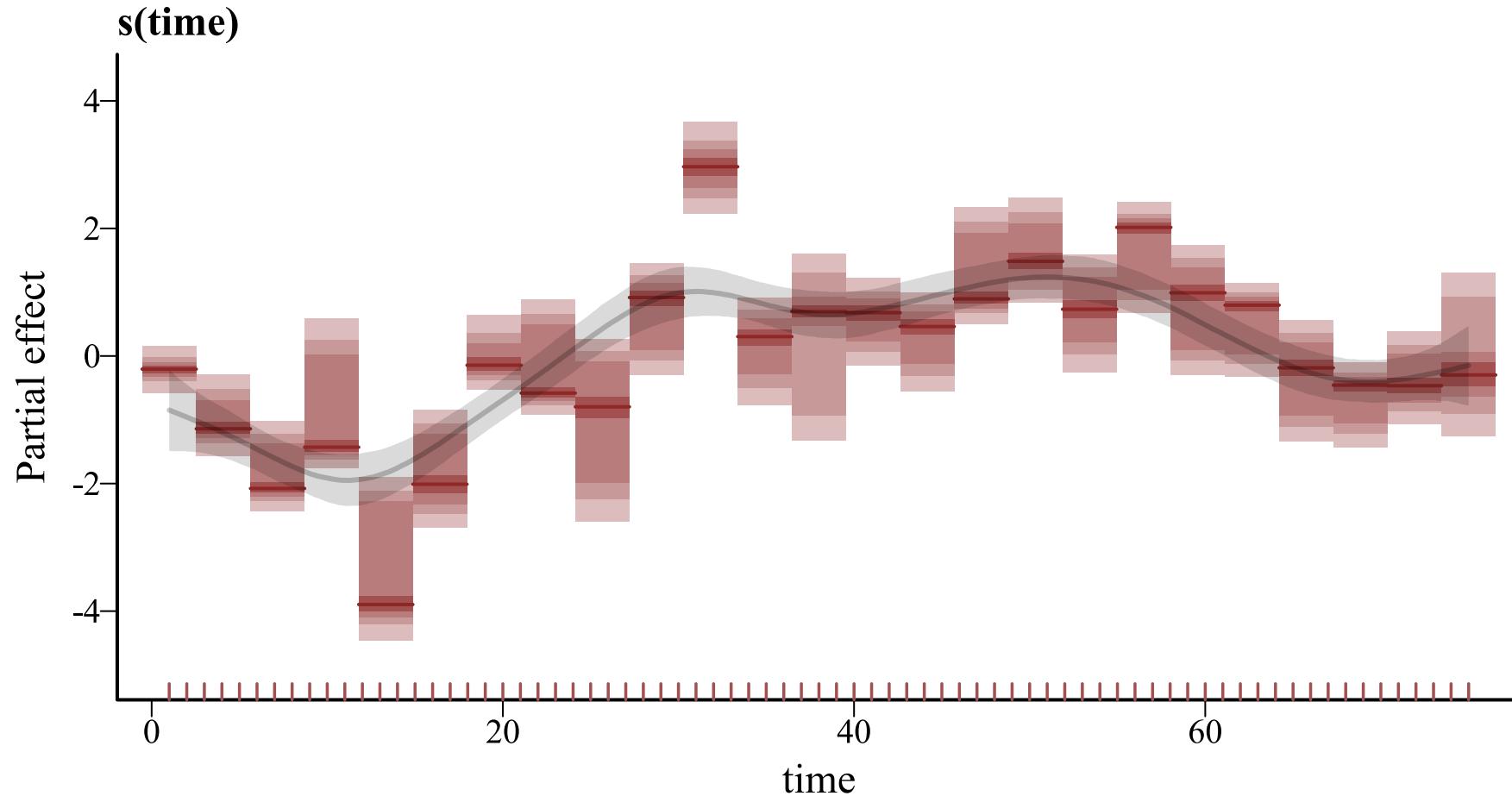
# Increase complexity?

```
model ← mvgam(y ~ s(time, k = 15), family = gaussian(),
               data = data_train, newdata = data_test)

gam.check(model$mgcv_model)
```

```
## 
## Method: REML   Optimizer: outer newton
## full convergence after 5 iterations.
## Gradient range [-1.64113e-07,3.260311e-08]
## (score 86.84541 & scale 0.4085855).
## Hessian positive definite, eigenvalue range [1.993815,36.91466].
## Model rank =  15 / 15
##
## Basis dimension (k) checking results. Low p-value (k-index<1) may
## indicate that k is too low, especially if edf is close to k'.
##
##          k'    edf k-index p-value
## s(time) 14.00   8.57    0.82   0.045 *
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
```

# Not wiggly enough



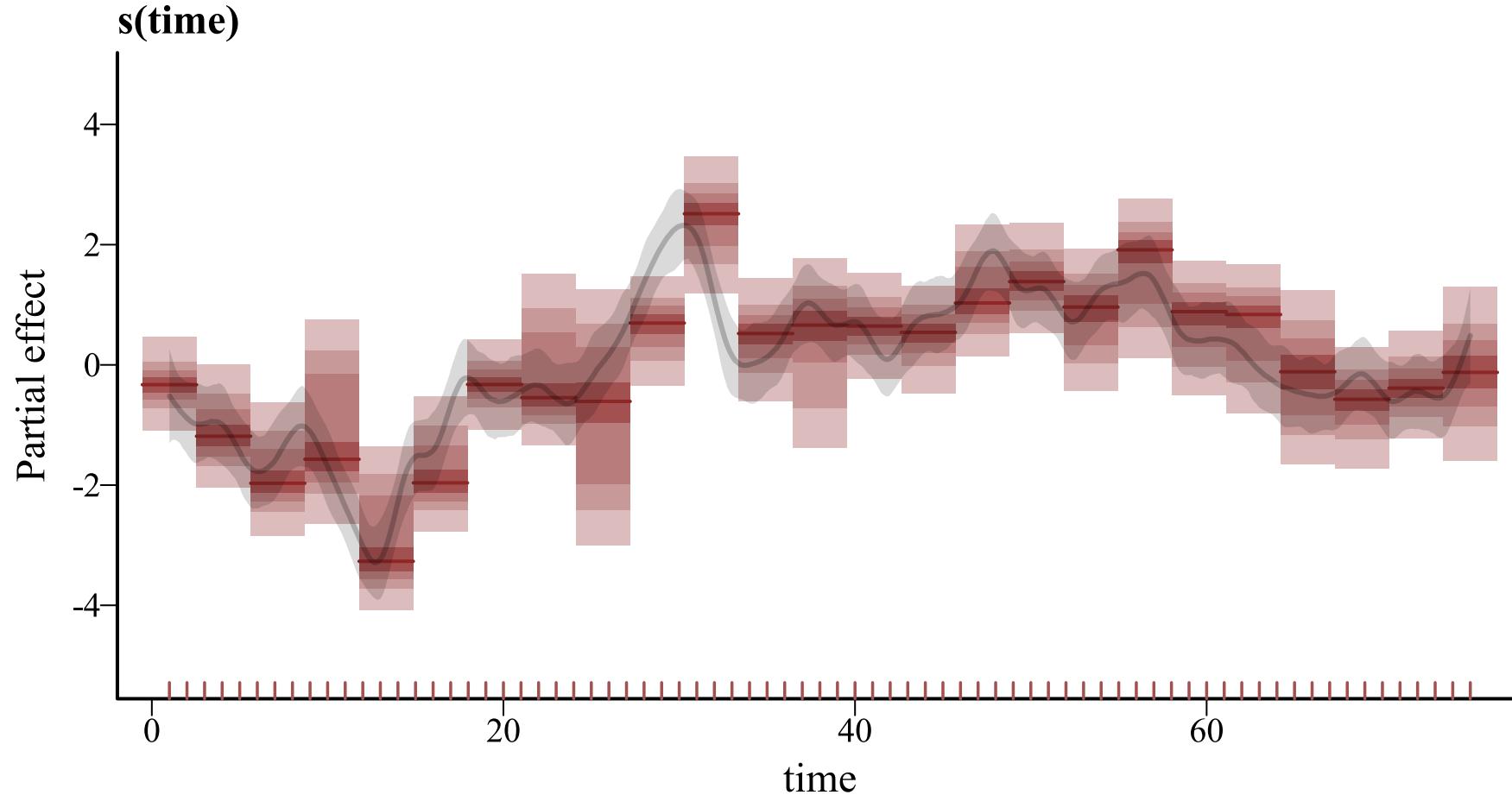
# Even more complex?

```
model ← mvgam(y ~ s(time, k = 50), family = gaussian(),
               data = data_train, newdata = data_test)

gam.check(model$mgcv_model)
```

```
##  
## Method: REML  Optimizer: outer newton  
## full convergence after 6 iterations.  
## Gradient range [1.752317e-07,4.46345e-07]  
## (score 84.14271 & scale 0.372791).  
## Hessian positive definite, eigenvalue range [0.883367,37.11506].  
## Model rank = 50 / 50  
##  
## Basis dimension (k) checking results. Low p-value (k-index<1) may  
## indicate that k is too low, especially if edf is close to k'.  
##  
##          k'   edf k-index p-value  
## s(time) 49.0 10.4     0.9    0.2
```

# Finally wiggly enough

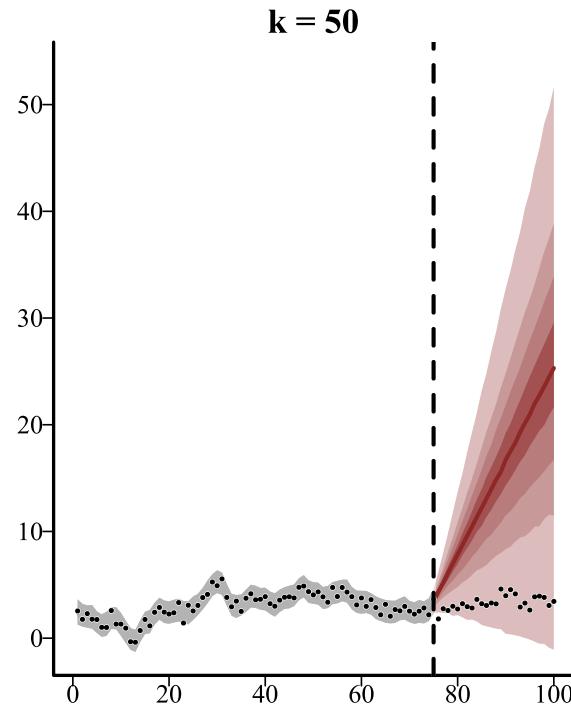
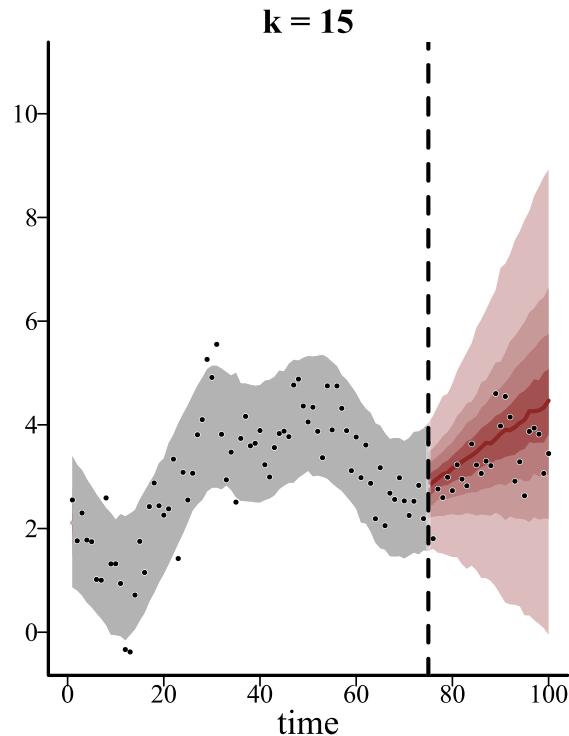
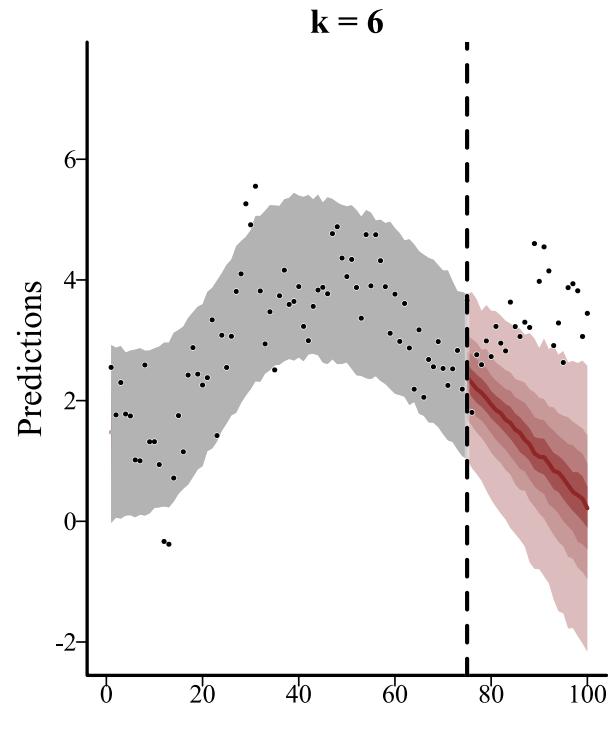


**Capturing this autocorrelation is important**

**Improves inferences on other parts of the model, while also giving more appropriate p-values, confidence intervals etc... in frequentist paradigms**

**But what effect does this variation in wiggliness have on forecasts?**

# Forecasts vary *hugely*





`gam.check` is sensitive to unmodelled autocorrelation

Raising `k` to satisfy warnings may improve inference on historical patterns, but leads to even more unpredictable extrapolation behaviour

If the goal is to produce predictions (i.e. to forecast), we can do better with appropriate *time series models*

# Ok. Can we just do this?

A linear model with an autoregressive term

$$\begin{aligned}\mathbf{Y}_t &\sim \text{Normal}(\mu_t, \sigma) \\ \mu_t &= \alpha + \beta_1 \mathbf{Y}_{t-1} + \dots\end{aligned}$$

Where:

$\alpha$  is an intercept coefficient

$\beta_1$  is a ***first-order autoregressive coefficient***

Can sometimes work because of identity link; but missingness, measurement error will still cause problems

# What about Poisson?

A Poisson GLM with an autoregressive term

$$\mathbf{Y}_t \sim \text{Poisson}(\lambda_t)$$

$$\log(\lambda_t) = \alpha + \beta_1 \mathbf{Y}_{t-1} + \dots$$

Where:

$\alpha$  is an intercept coefficient

$\beta_1$  is a ***first-order autoregressive coefficient***

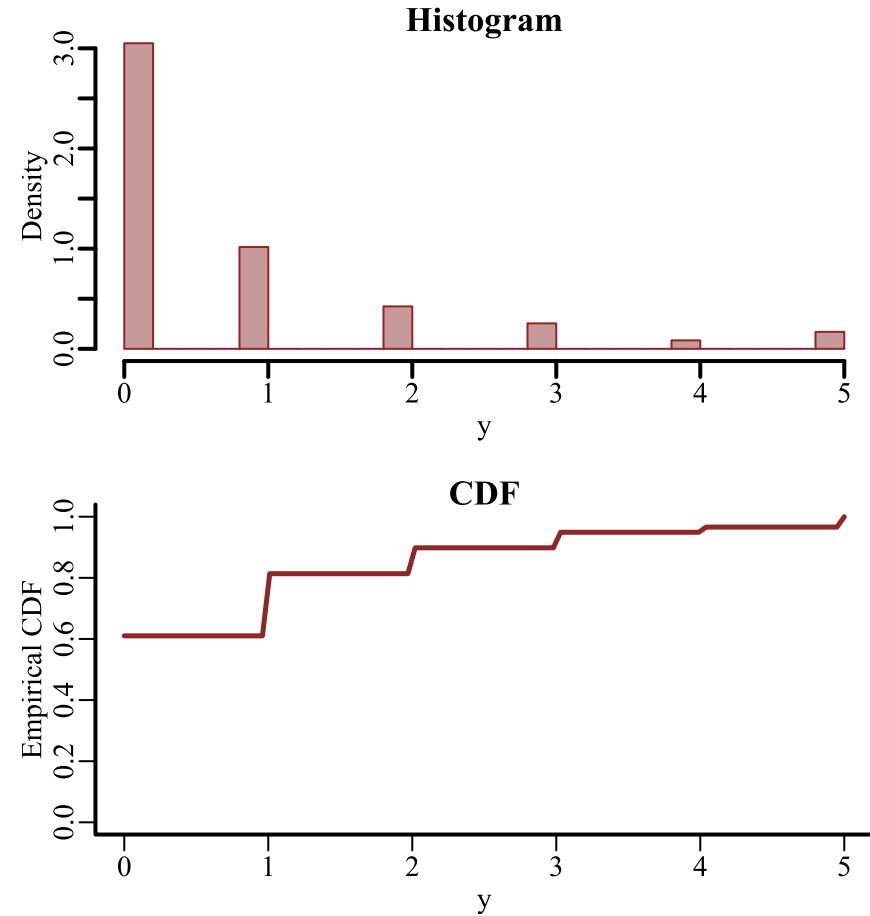
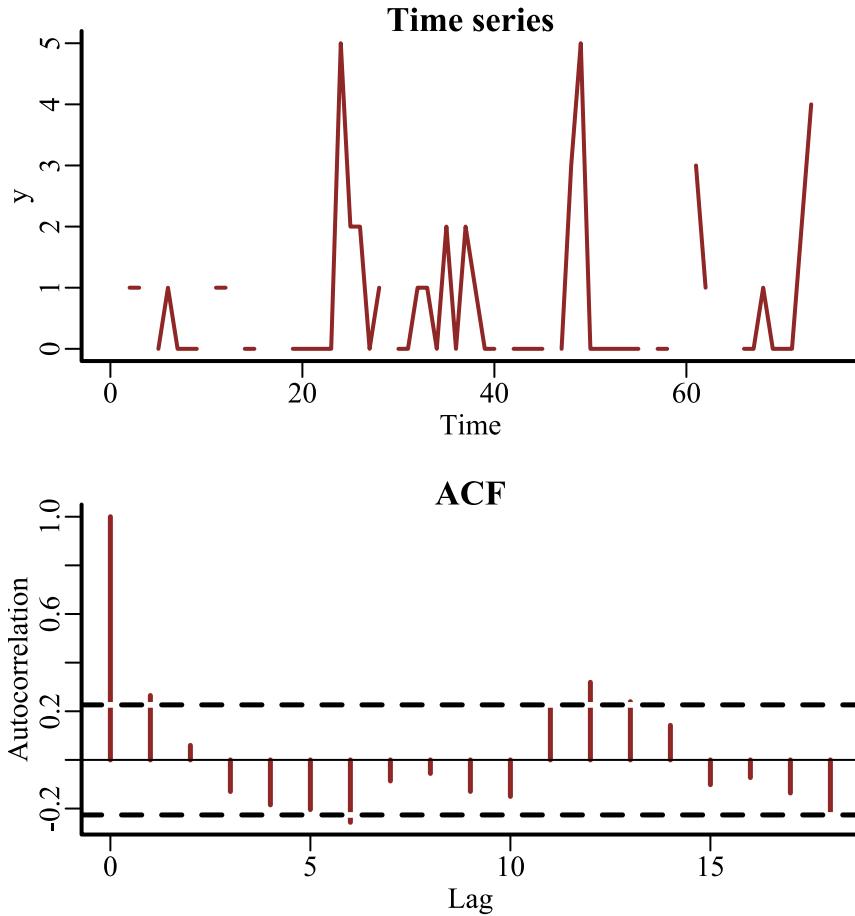
# A motivating example

```
# set seed for reproducibility
set.seed(222)

# simulate an integer-valued time series with some missing observations
sim_data ← sim_mvgam(T = 100, n_series = 1,
                      trend_model = 'RW',
                      prop_missing = 0.2)
```

y	season	year	series	time
NA		1	1	series_1
1	2	1	series_1	2
1	3	1	series_1	3
NA	4	1	series_1	4

# The simulated data



# Use the **tscount** ?

```
# attempt a tscount time series model
# which can fit autoregressive models for count time series
library(tscount)

# use the tsglm function for AR modelling
tsglm(sim_data$data_train$y,

      # model using outcome at lag 1 as the predictor
      model = list(past_obs = 1))
```

```
## Error in tsglm.meanfit(ts = ts, model = model, xreg = xreg, link = link, :
# Cannot make estimation with missing values in time series or covariates
```

NAs cause big problems in autoregressive models

# NAs compound over lags

time	y	y_lag1	y_lag2	season	year	series
1	NA	NA	NA	1	1	series_1
2	1	NA	NA	2	1	series_1
3	1	1	NA	3	1	series_1
4	NA	1	1	4	1	series_1
5	0	NA	1	5	1	series_1
6	1	0	NA	6	1	series_1
7	0	1	0	7	1	series_1
8	0	0	1	8	1	series_1

# Only 2/8 rows complete

time	y	y_lag1	y_lag2	season	year	series
1	NA	NA	NA	1	1	series_1
2	1	NA	NA	2	1	series_1
3	1	1	NA	3	1	series_1
4	NA	1	1	4	1	series_1
5	0	NA	1	5	1	series_1
6	1	0	NA	6	1	series_1
7	0	1	0	7	1	series_1
8	0	0	1	8	1	series_1

# Other problems of AR observations

Measurement errors also compound

Difficult / impossible to ensure stability of forecasts

Can use  $\log(Y_{t-lag})$  as predictors, but this doesn't always work

Challenging to link dynamics across multiple series

Not extendable to other types of dynamics

Smooth temporal evolution

Changepoint models

Stochastic variance / volatility

etc...

Latent  
autoregressive  
processes

# Dynamic Poisson GLM

A dynamic Poisson GLM can use *autocorrelated latent residuals*

$$Y_t \sim \text{Poisson}(\lambda_t)$$

$$\log(\lambda_t) = \alpha + \dots + z_t$$

$$z_t \sim \text{Normal}(z_{t-1}, \sigma)$$

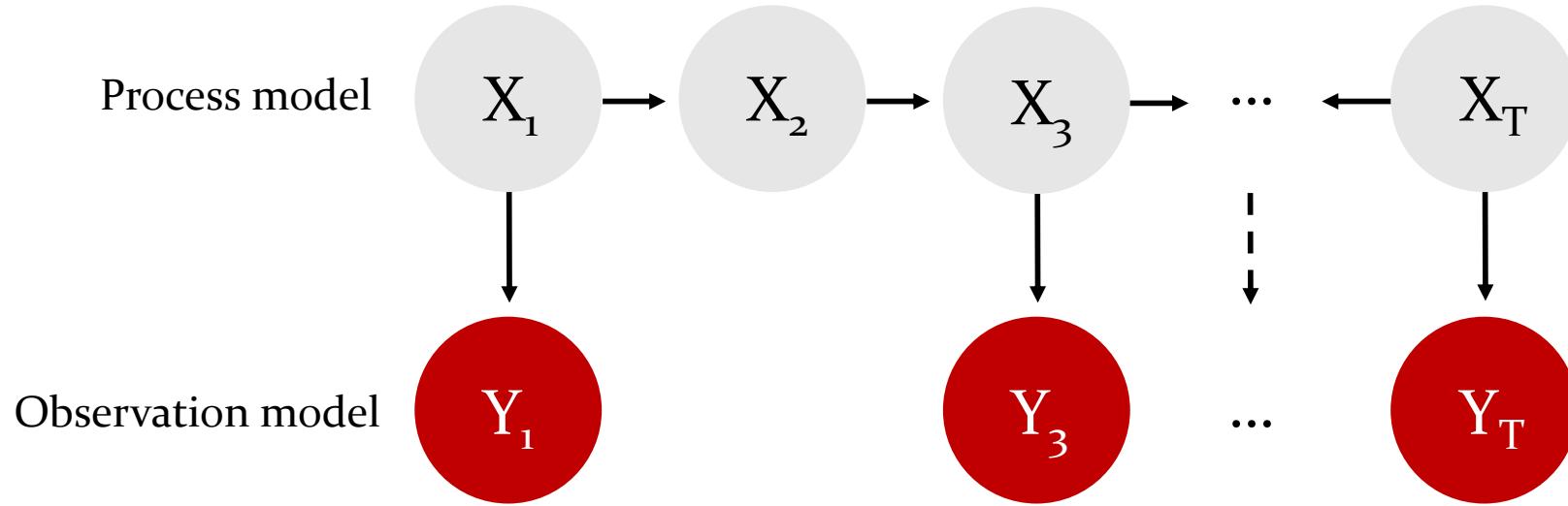
$$\sigma \sim \text{Exponential}(2)$$

Where:

$z_t$  is the value of the latent residual at time  $t$

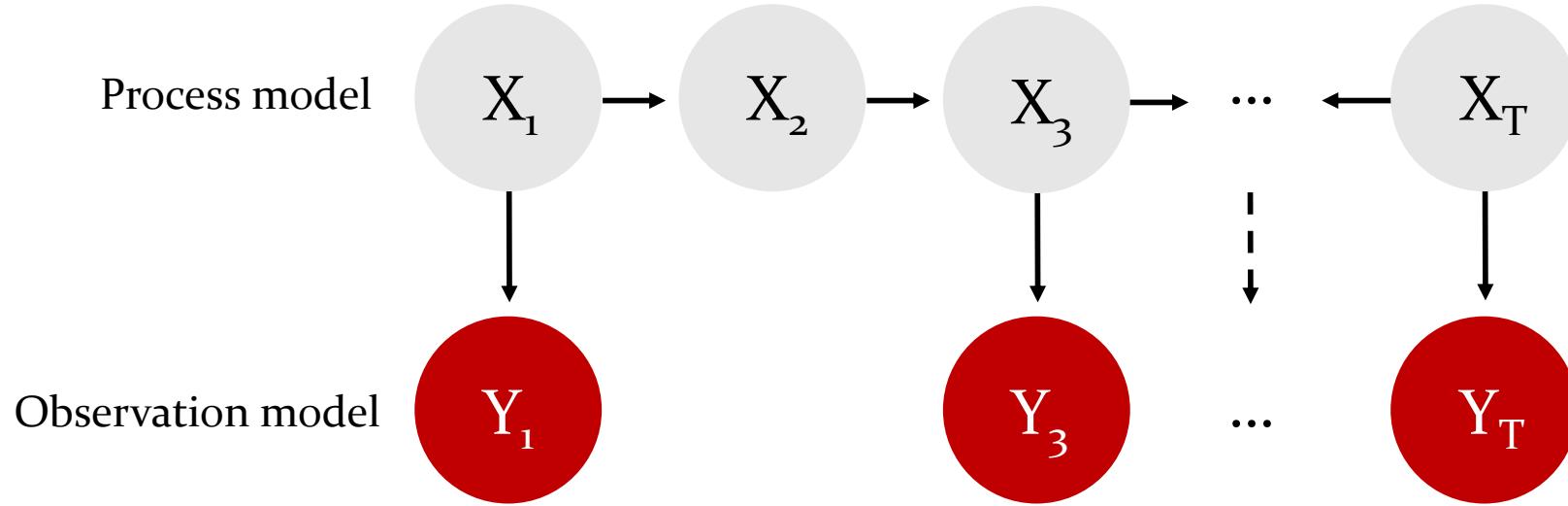
$\sigma$  captures variation in the latent dynamic process

# Evolves *independently*



Missing observations do not impede evolution of the *latent* process

# Evolves *independently*



The latent process model can take on a *huge variety* of forms

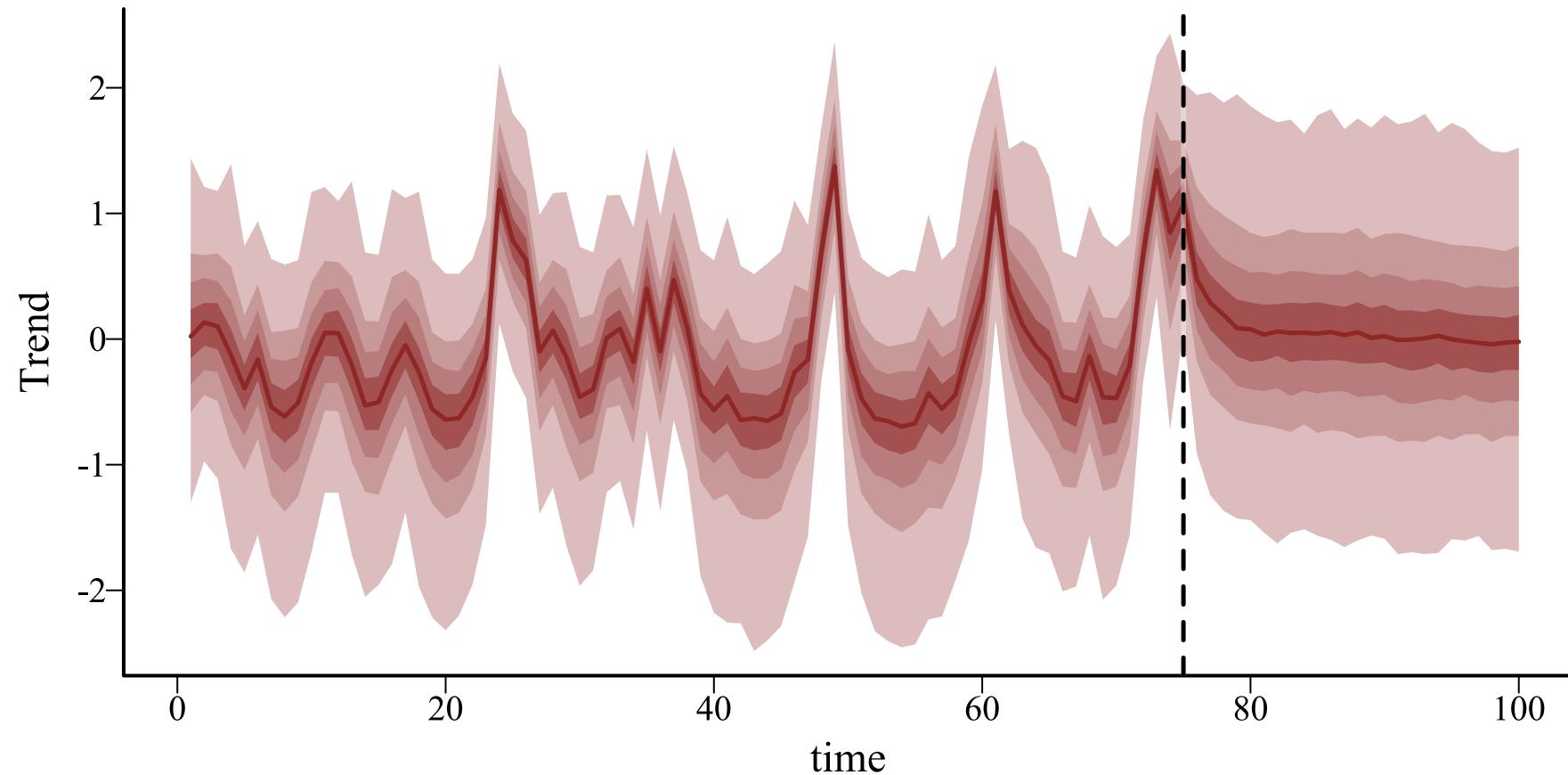
# Back to the example

```
mod_example <- mvgam(y ~ 1,  
                      trend_model = 'AR1',  
                      data = sim_data$data_train,  
                      newdata = sim_data$data_test,  
                      family = poisson())
```

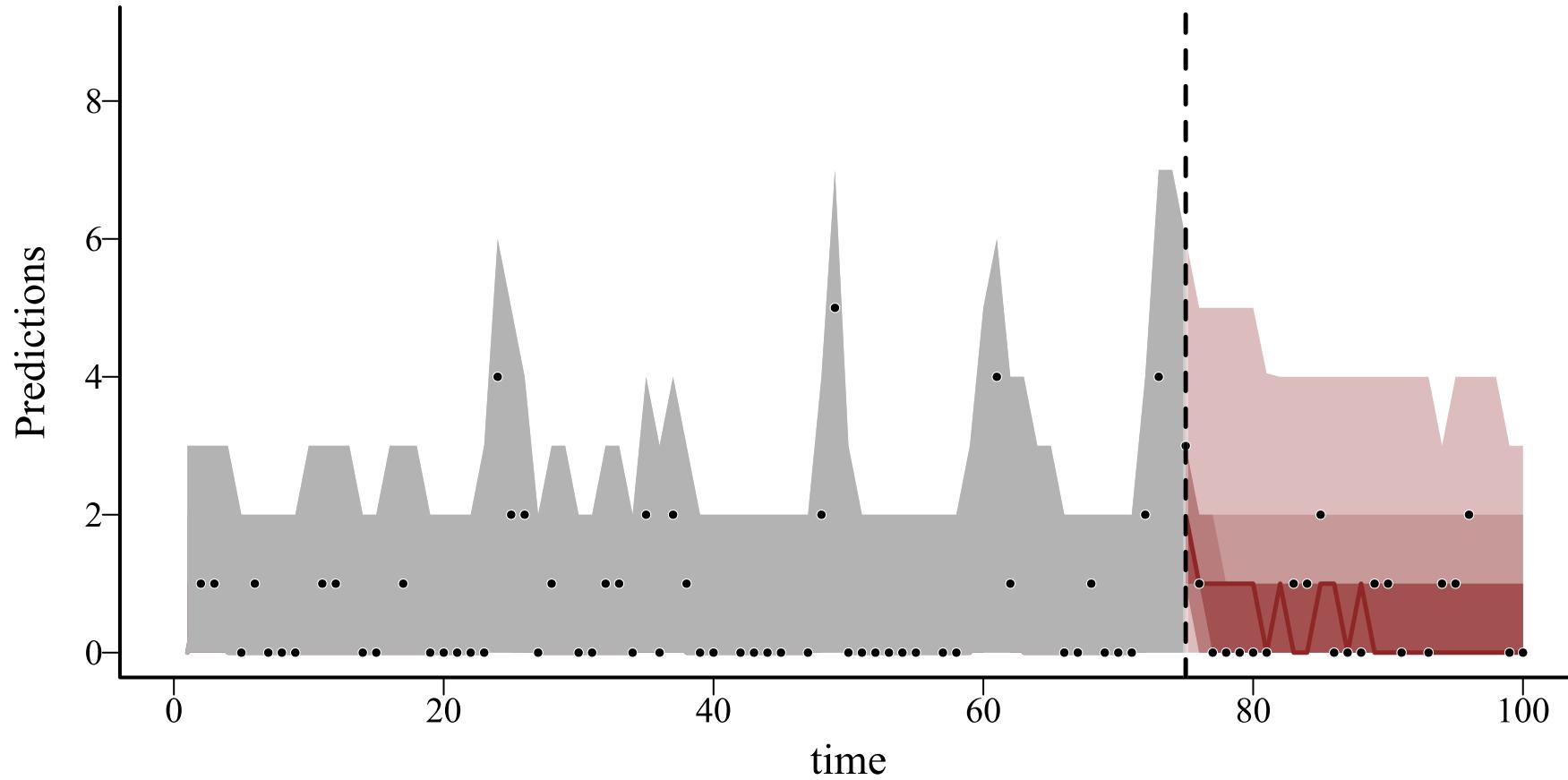
`mvgam`  has no problem with these observations

Fit a model with latent AR1 dynamics and just an intercept in the observation model

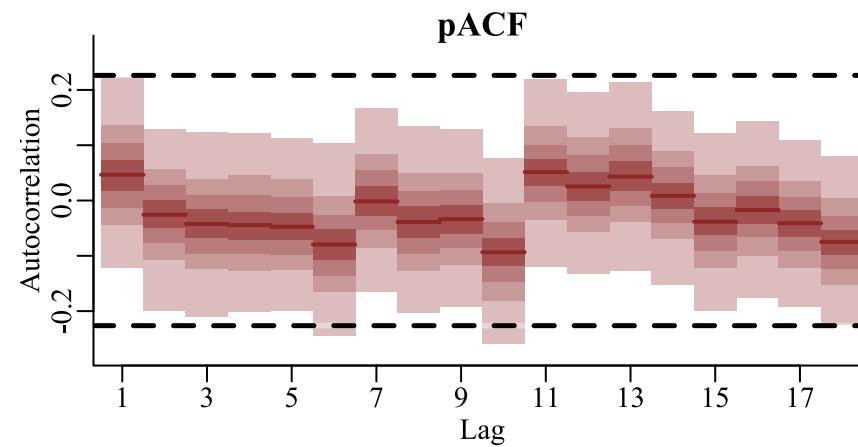
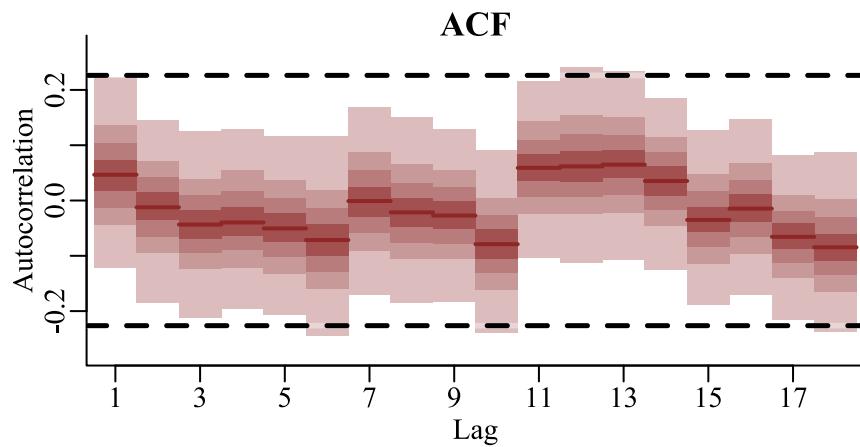
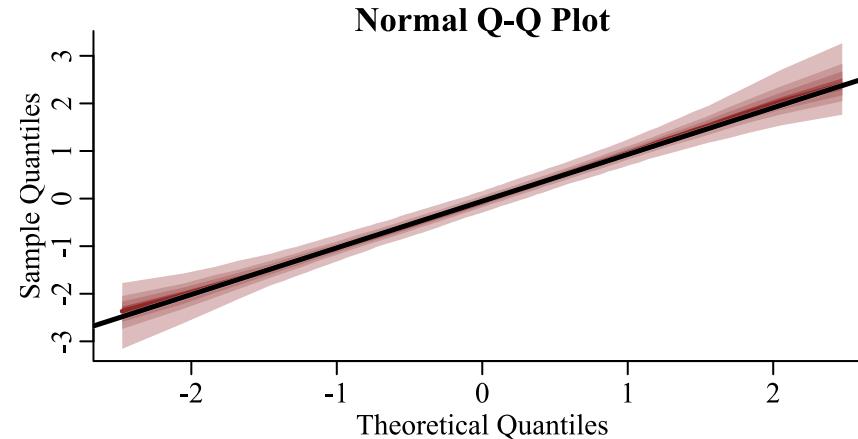
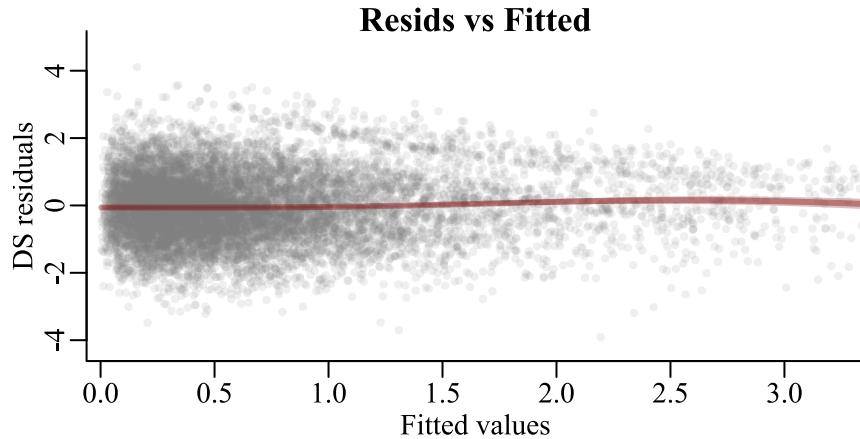
# The latent trend



# Forecasts



# Residuals



# A tougher example?

```
# set seed for reproducibility
set.seed(100)

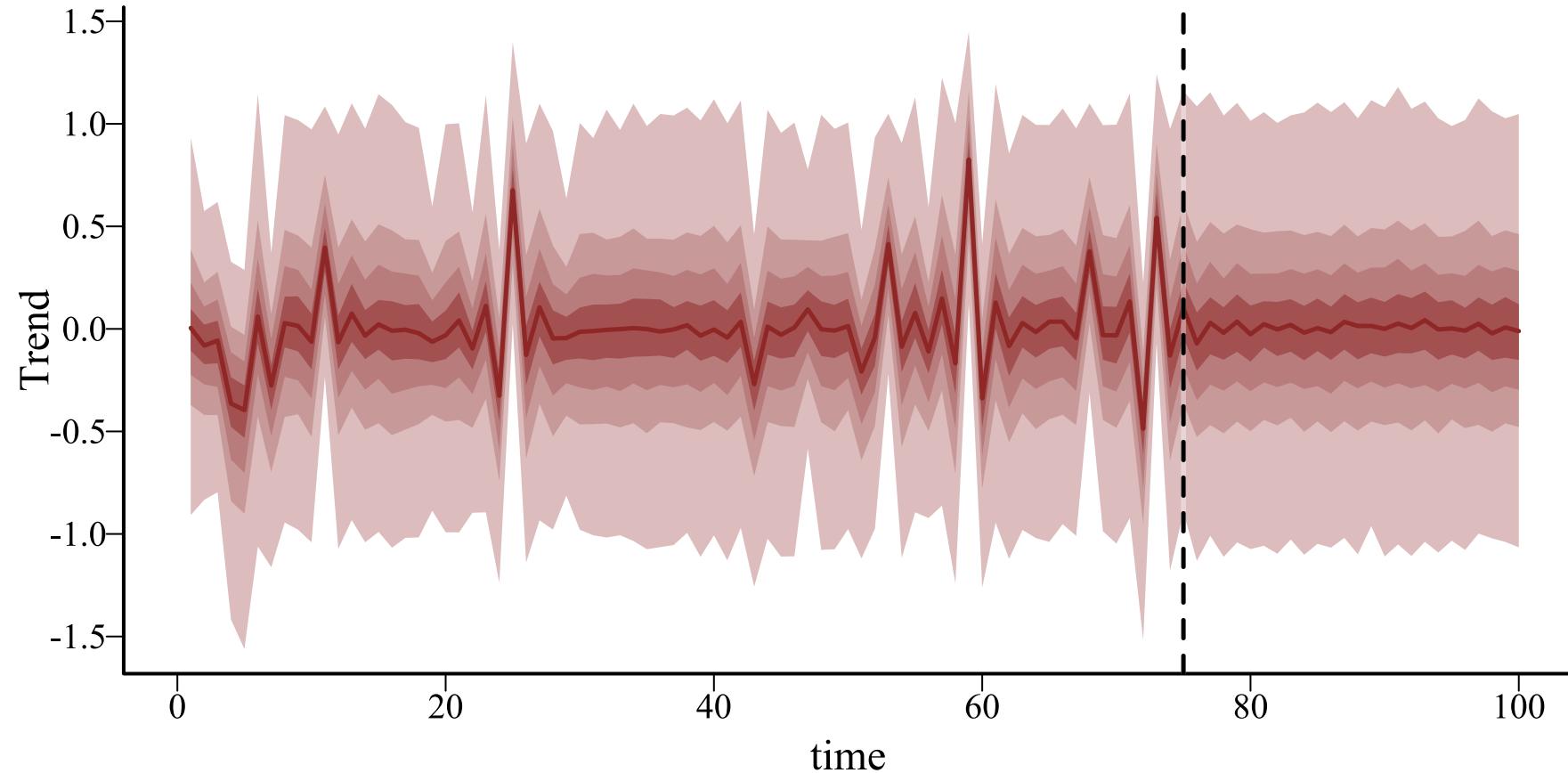
# simulate an integer-valued time series with some missing observations
sim_data2 ← sim_mvgam(T = 100, n_series = 1,
                       mu = 1,
                       trend_model = 'RW',
                       prop_missing = 0.75)
```

75% of observations missing!

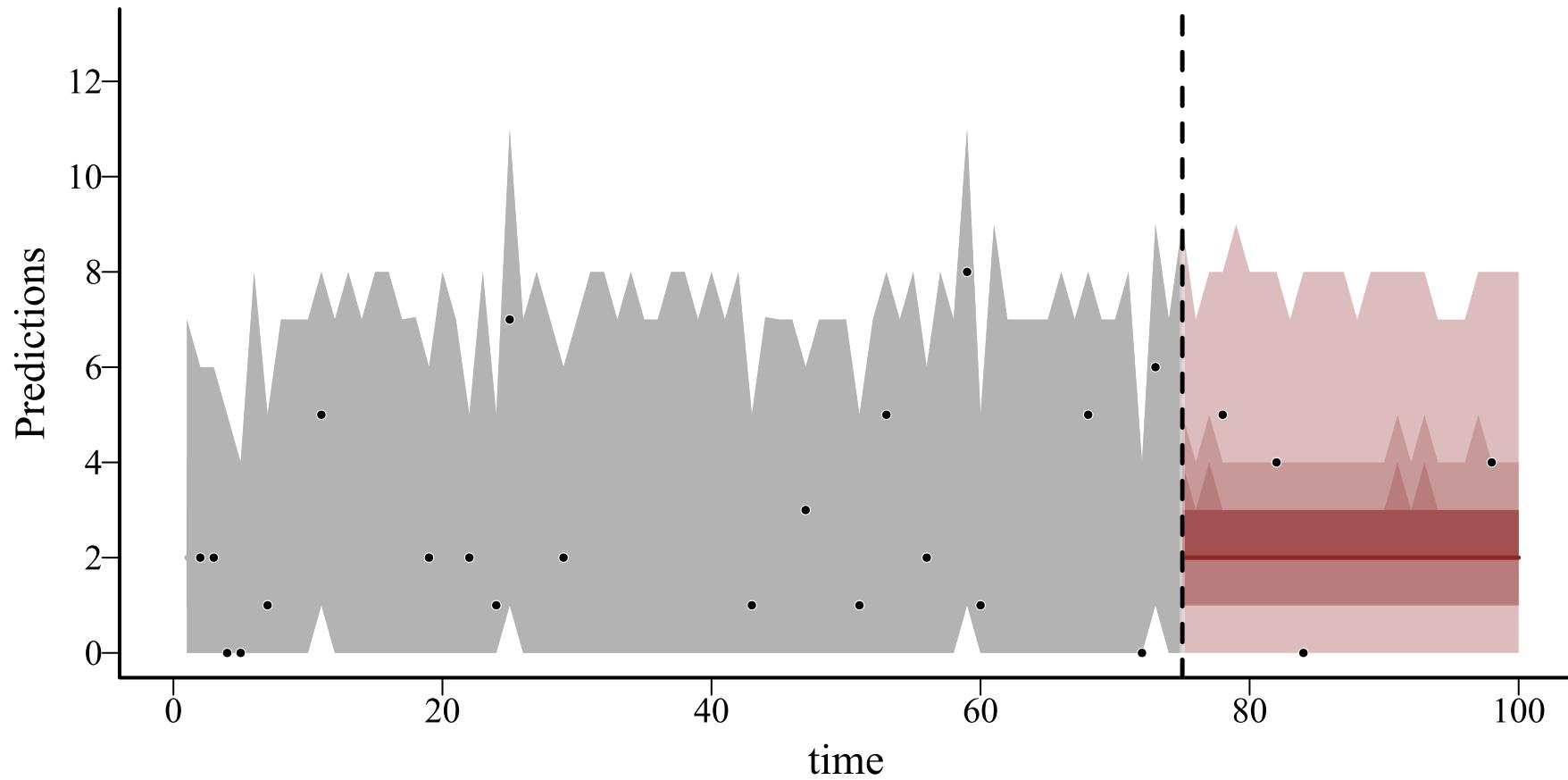
# Same model

```
mod_example2 ← mvgam(y ~ 1,  
                      trend_model = 'AR1',  
                      data = sim_data2$data_train,  
                      newdata = sim_data2$data_test,  
                      family = poisson())
```

# The latent trend



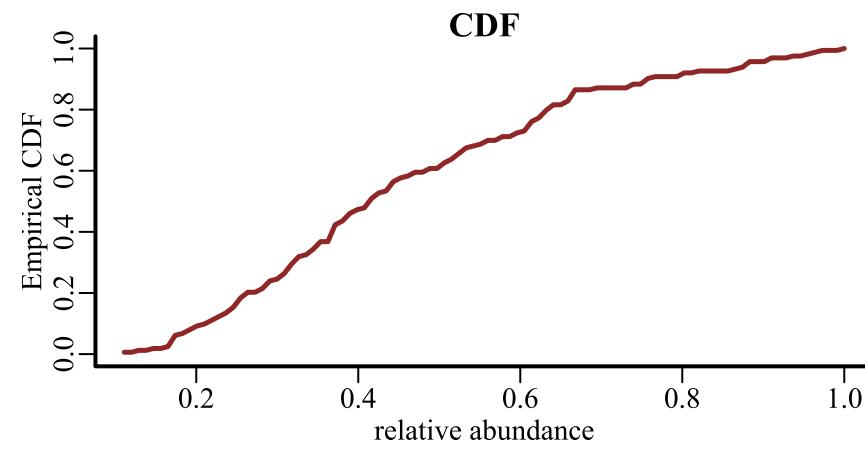
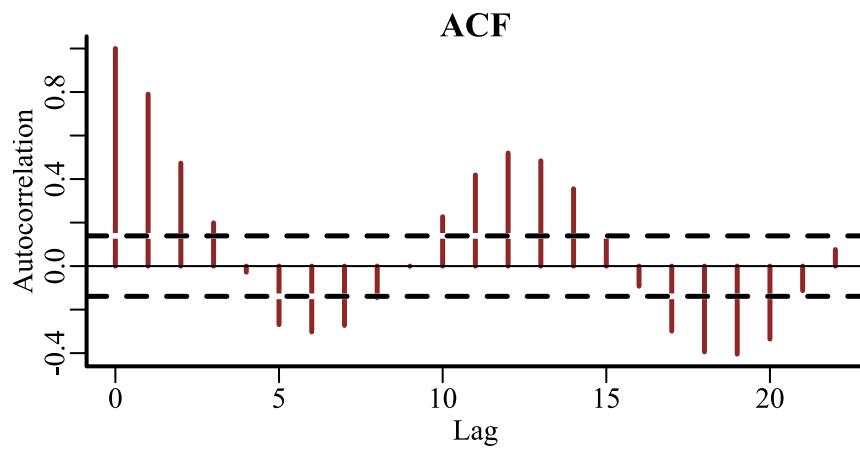
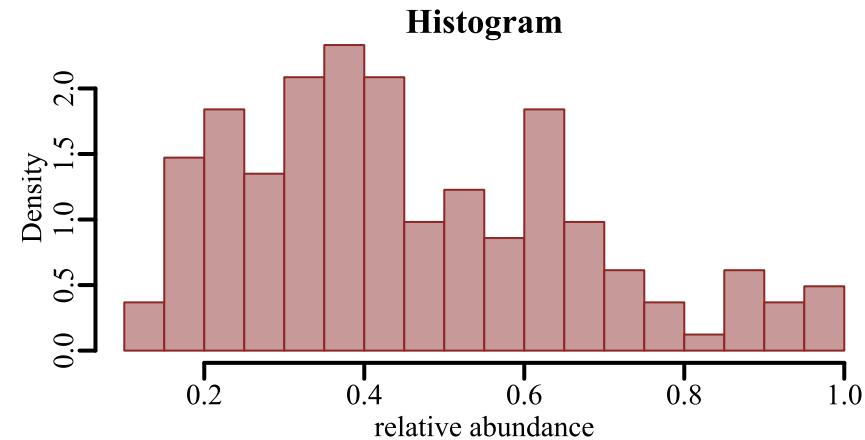
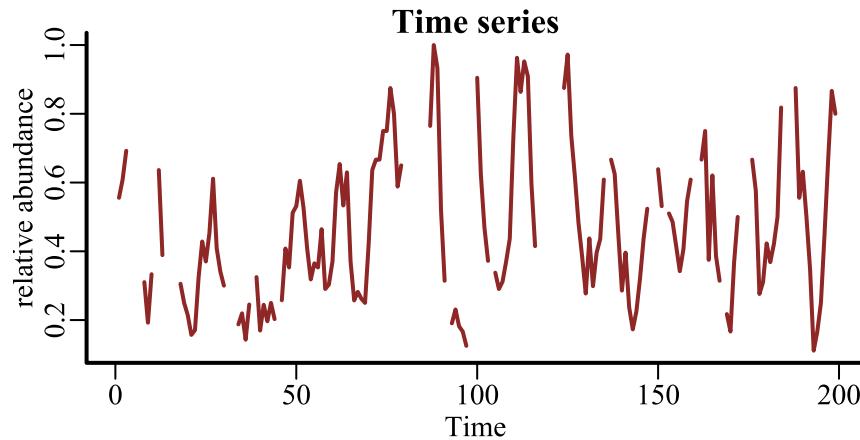
# Forecasts



*Some packages exist to model count-valued time series using autoregressive terms*

**But you must not have missing data or measurement error, and you cannot handle multiple series at once**

**Fine for some situations. But what if your data look like this?**



Properties of Merriam's kangaroo rat relative abundance time series from a long-term monitoring study in  
Portal, Arizona, USA

# Dynamic Beta GAM

```
mod_beta ← mgcv::mgcv(mgcv::relabund ~  
    mgcv::te(mintemp, ndvi),  
    mgcv::trend_model = 'AR3',  
    mgcv::family = betar(),  
    mgcv::data = dm_data)
```

Beta regression using the `mgcv`  's `betar` family

# Dynamic Beta GAM

```
mod_beta ← mgcv::mgcv(mvgam(relabund ~  
    te(mintemp, ndvi),  
    trend_model = 'AR3',  
    family = betar(),  
    data = dm_data)
```

Beta regression using the `mgcv`  's `betar` family

AR3 dynamic trend model

# Dynamic Beta GAM

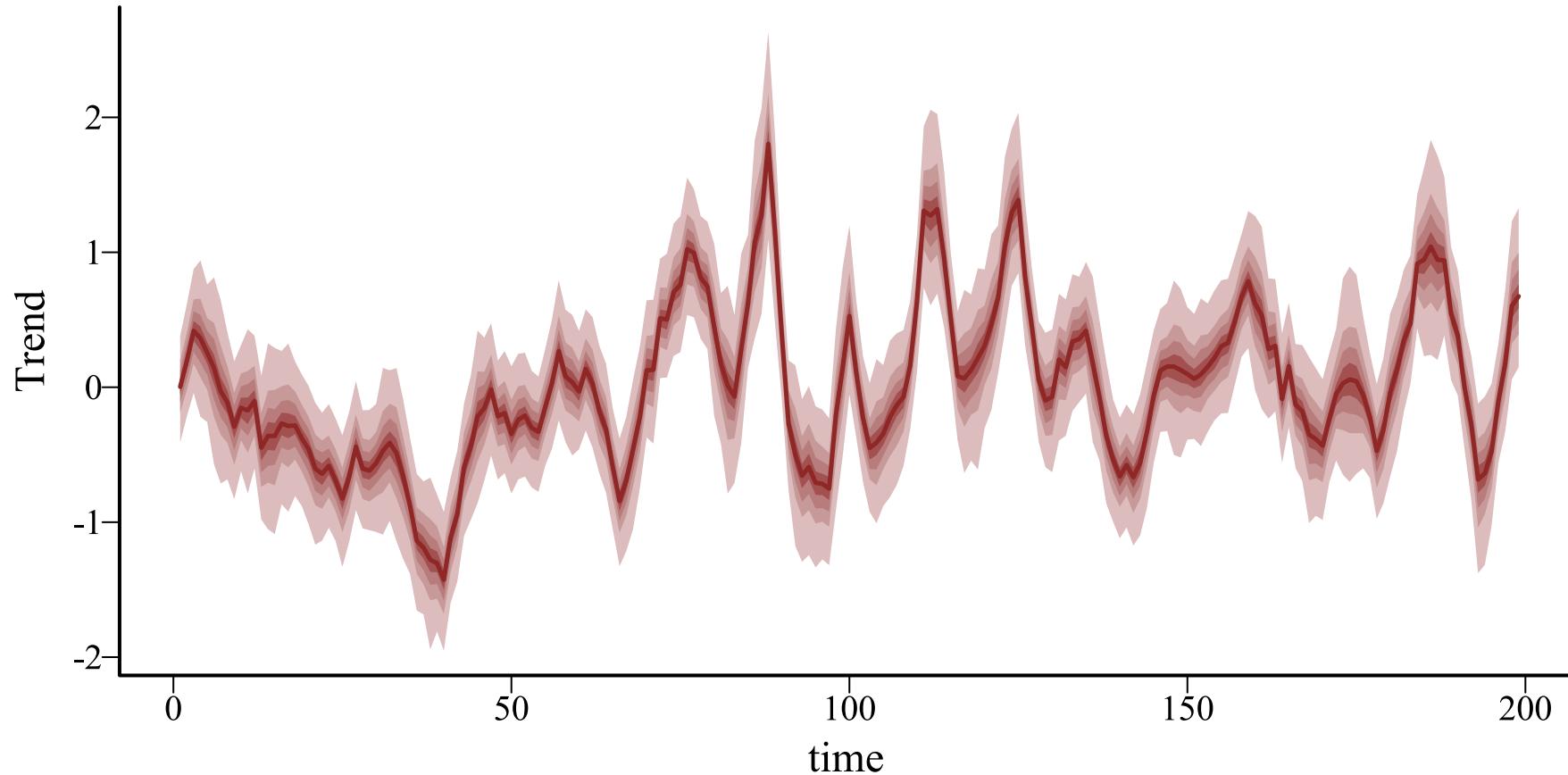
```
mod_beta ← mgcv::mgcv(mgcv::relabund ~  
    mgcv::te(mintemp, ndvi),  
    trend_model = 'AR3',  
    family = betar(),  
    data = dm_data)
```

Beta regression using the `mgcv`  's `betar` family

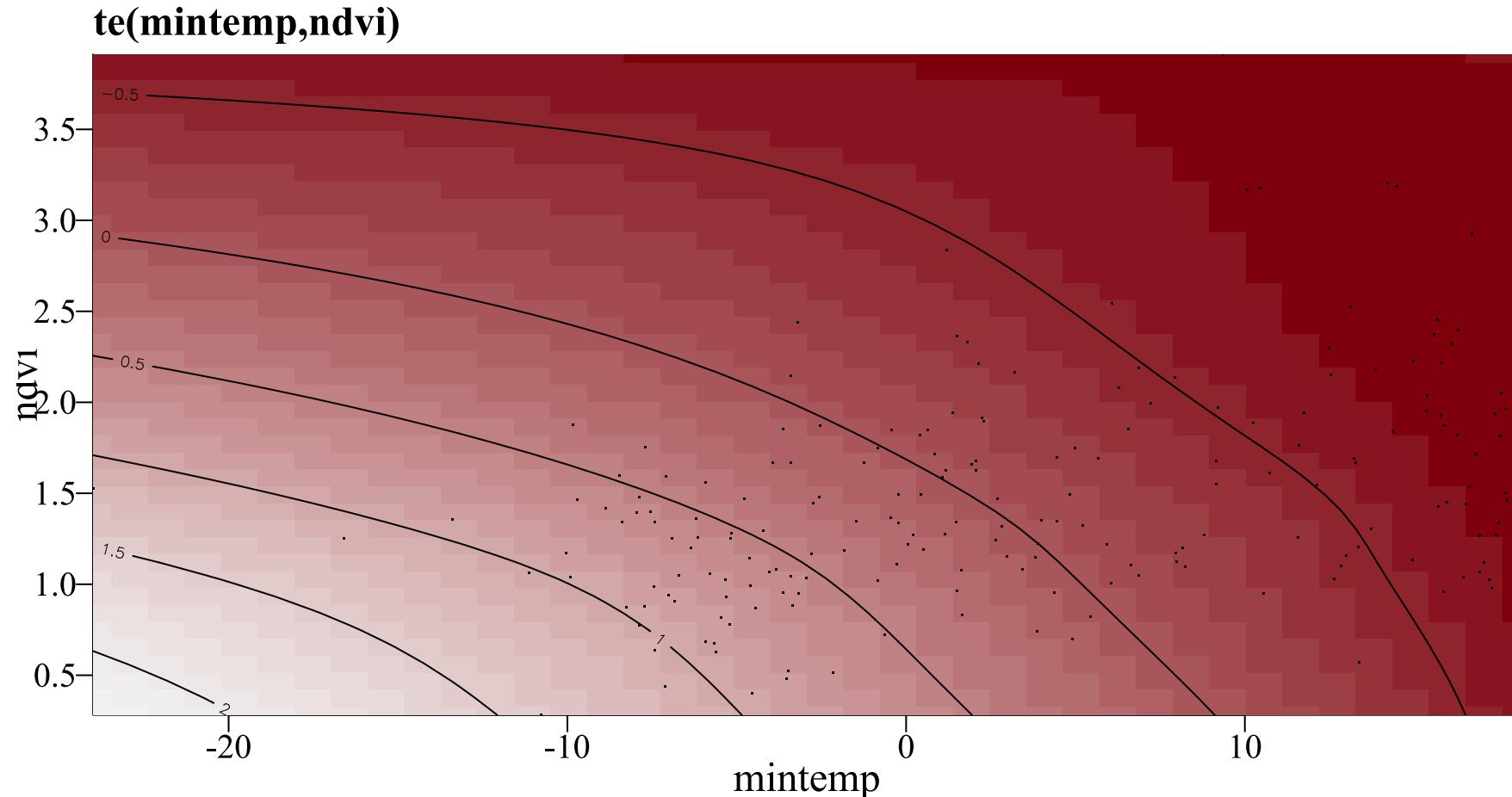
AR3 dynamic trend model

Multidimensional tensor product smooth function for nonlinear covariate interactions (using `te`).

# The latent trend



# Multidimensional smooth



# Huh?



# marginaleffects for clarity

---

Code    Plot

---

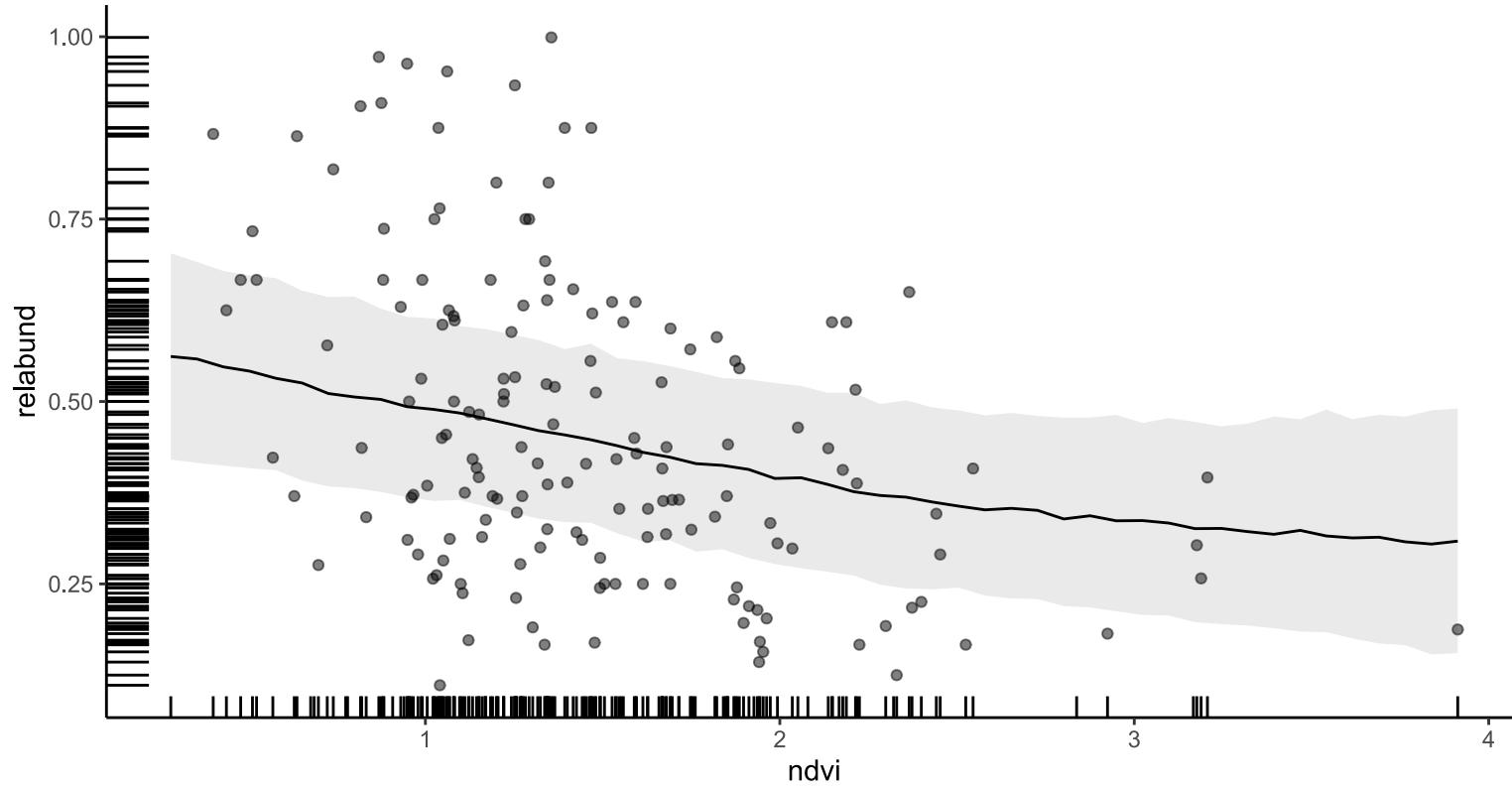
```
# plot conditional effect of NDVI on the outcome scale
plot_predictions(mod_beta, condition = 'ndvi',
                  points = 0.5, conf_level = 0.8, rug = TRUE) +
  theme_classic()
```

# marginaleffects for clarity

---

Code Plot

---



# marginaleffects for clarity

---

Code    Plot

---

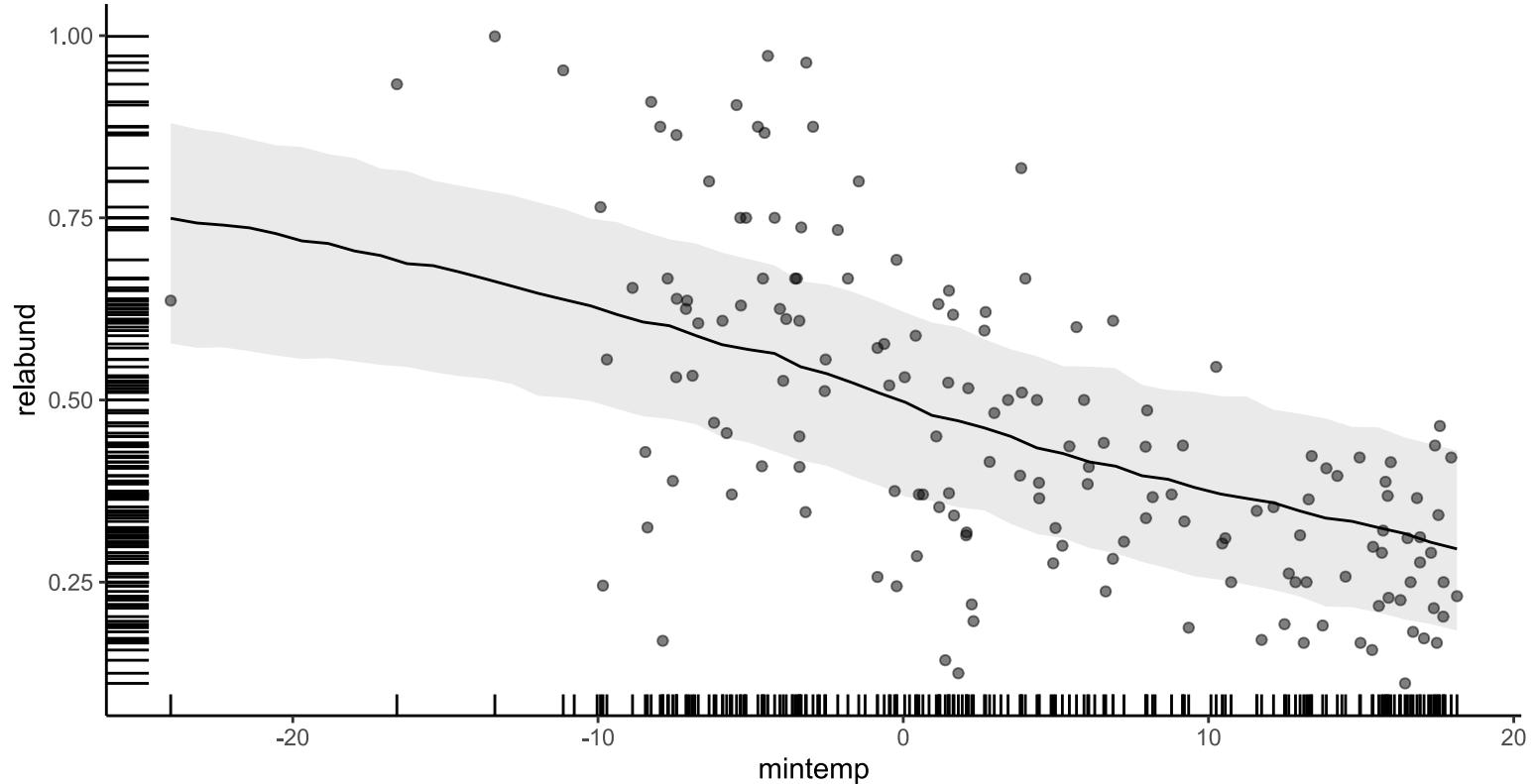
```
# plot conditional effect of Min Temp on the outcome scale
plot_predictions(mod_beta, condition = 'mintemp',
                  points = 0.5, conf_level = 0.8, rug = TRUE) +
  theme_classic()
```

# marginaleffects for clarity

---

Code Plot

---



# marginaleffects for clarity

---

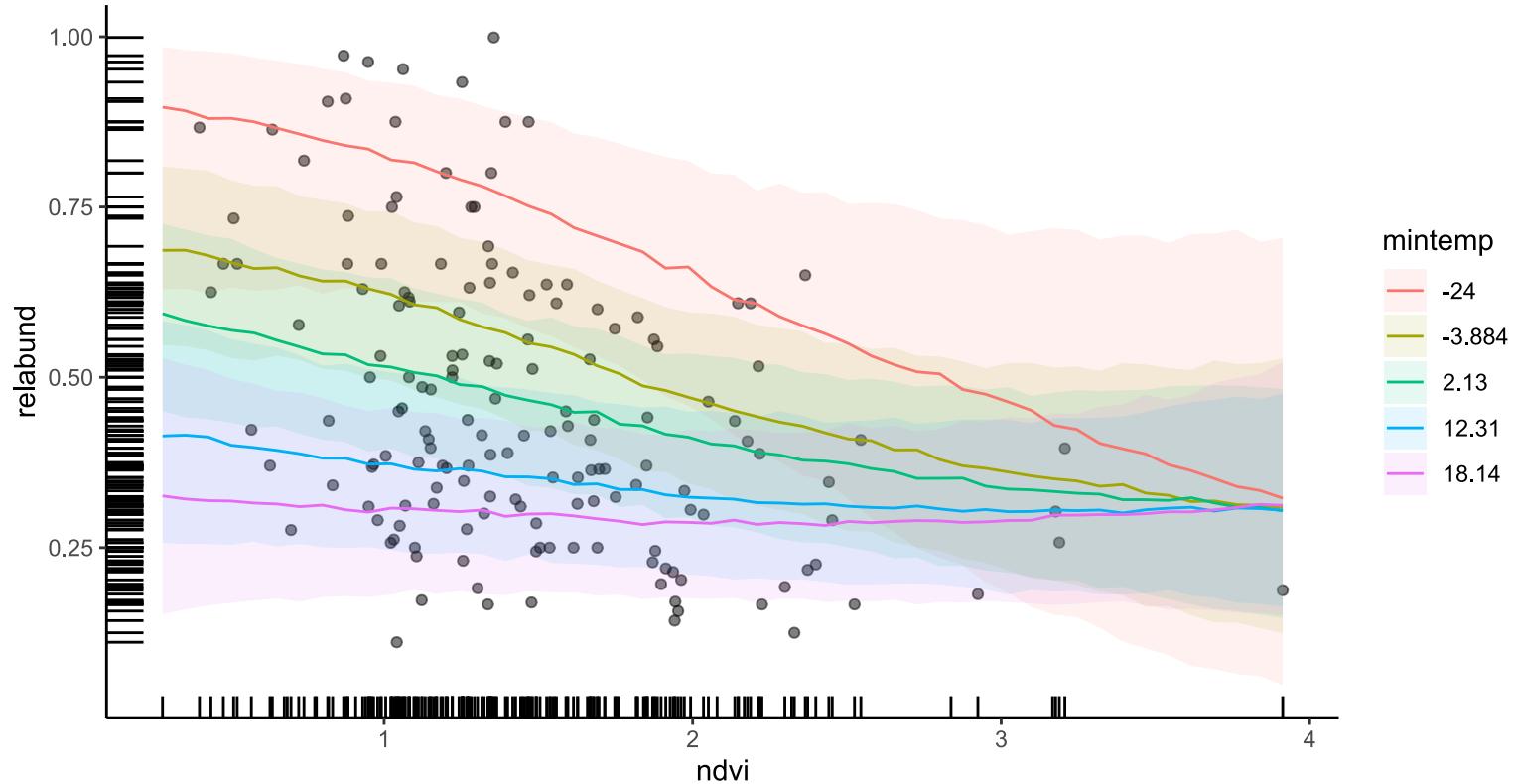
Code    Plot

---

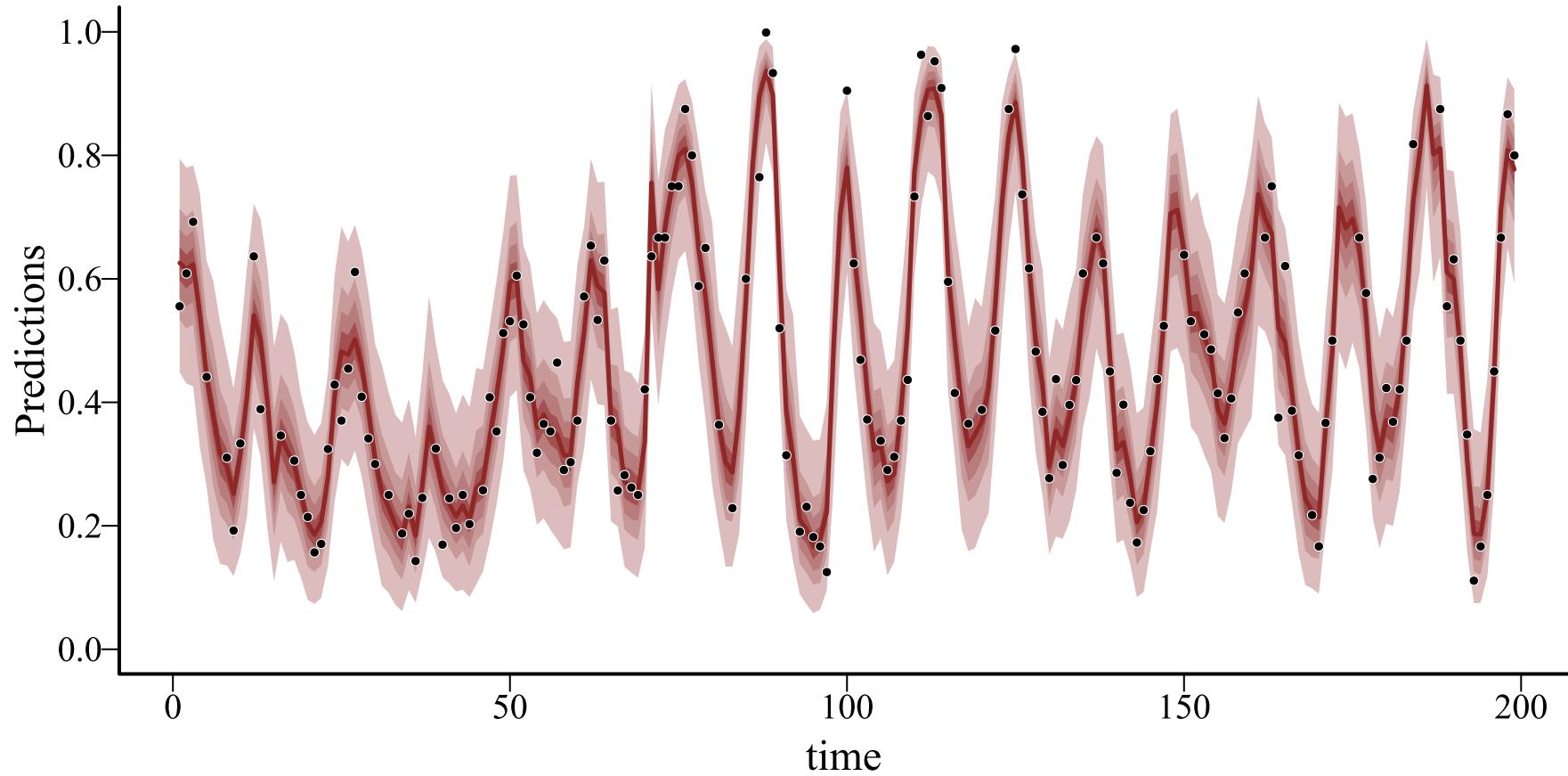
```
# plot conditional effect of BOTH covariates on the outcome scale
plot_predictions(mod_beta, condition = c('ndvi', 'mintemp'),
                  points = 0.5, conf_level = 0.8, rug = TRUE) +
  theme_classic()
```

# marginaleffects for clarity

Code Plot



# Hindcasts



We can estimate latent dynamic residuals for *many* types of GLMs / GAMs, thanks to the link function

We do not need to regress the outcome on its own past values

Very advantageous for ecological time series. But what kinds of dynamic processes are available in the `mvgam` and `brms`  's?

# Random walks

Simple stochastic processes that can fit a wide range of data

$$z_t \sim \text{Normal}(\alpha + z_{t-1}, \sigma)$$

Where:

$\sigma$  determines the spread (or flexibility) of the process

$\alpha$  is an optional intercept or *drift* parameter

Process at time  $t$  is centred around it's own value at time  $t - 1$ , with spread determined by probabilistic error

# A Random Walk

---

Code    Plot

---

```
# set seed and number of timepoints
set.seed(1111); T <- 100

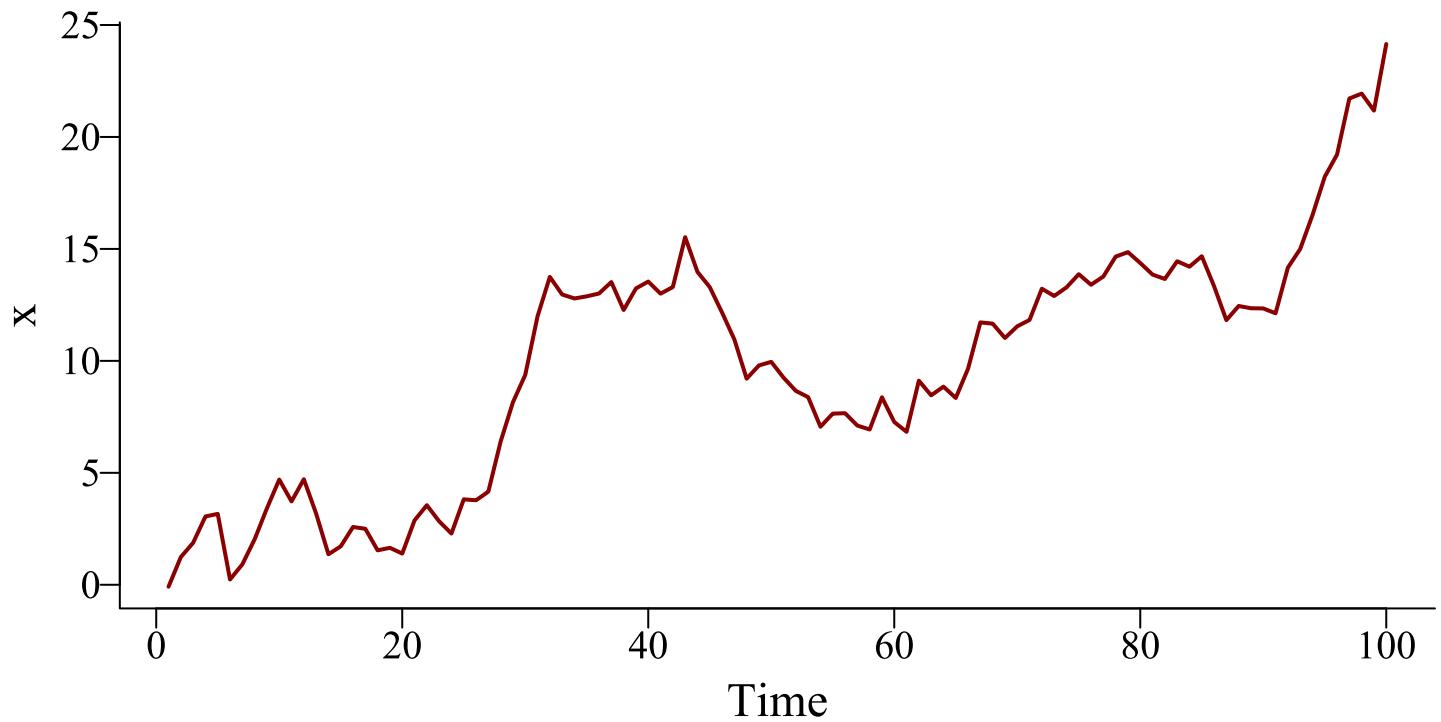
# initialize first value
series <- vector(length = T); series[1] <- rnorm(n = 1, mean = 0, sd = 1)

# compute values 2 through T
for (t in 2:T) {
  series[t] <- rnorm(n = 1, mean = series[t - 1], sd = 1)
}

# plot the time series as a line
plot(series, type = 'l', bty = 'l', lwd = 2,
      col = 'darkred', ylab = 'x', xlab = 'Time')
```

# A Random Walk

[Code](#)   [Plot](#)



# AR1

Similar to a Random Walk and can fit a wide range of data

$$z_t \sim \text{Normal}(\alpha + \phi * z_{t-1}, \sigma)$$

Where:

$\sigma$  determines the spread (or flexibility) of the process

$\alpha$  is an optional intercept or *drift* parameter

$\phi$  is a coefficient estimating correlation between  $z_t$  and  $z_{t-1}$

Process at time  $t$  is a *function* of it's own value at time  $t - 1$

# AR2 and AR3

As with AR1, but with additional autoregressive terms

$$z_t \sim \text{Normal}(\alpha + \phi_1 * z_{t-1} + \phi_2 * z_{t-2} + \phi_3 * z_{t-3}, \sigma)$$

# An AR1

---

Code    Plot

---

```
# set seed and number of timepoints
set.seed(1111); T ← 100

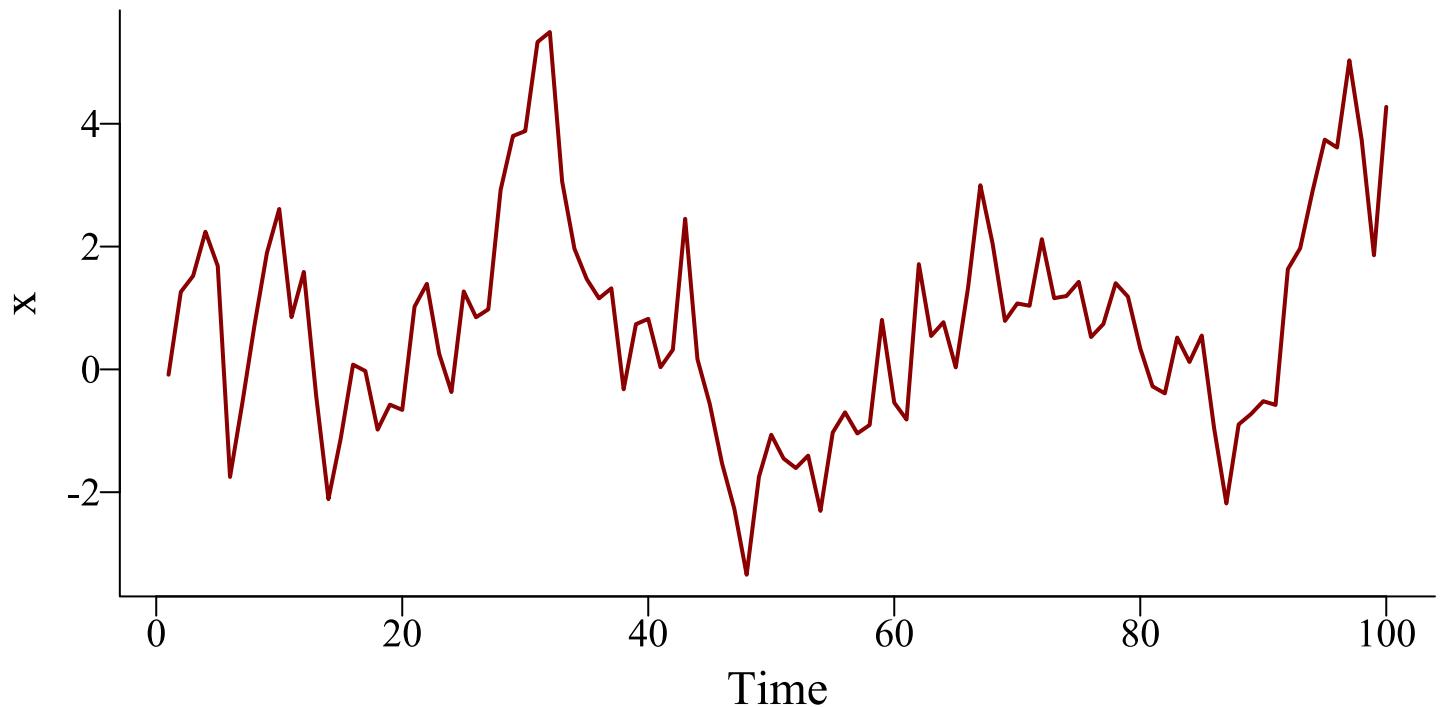
# initialize first value
series ← vector(length = T); series[1] ← rnorm(n = 1, mean = 0, sd = 1)

# compute values 2 through T, with phi = 0.7
for (t in 2:T) {
    series[t] ← rnorm(n = 1, mean = 0.7 * series[t - 1], sd = 1)
}

# plot the time series as a line
plot(series, type = 'l', bty = 'l', lwd = 2,
      col = 'darkred', ylab = 'x', xlab = 'Time')
```

# An AR1

Code Plot



# Alternative AR1 simulation

---

[Code](#)    [Plot](#)

---

```
# set seed for reproducibility
set.seed(1111)

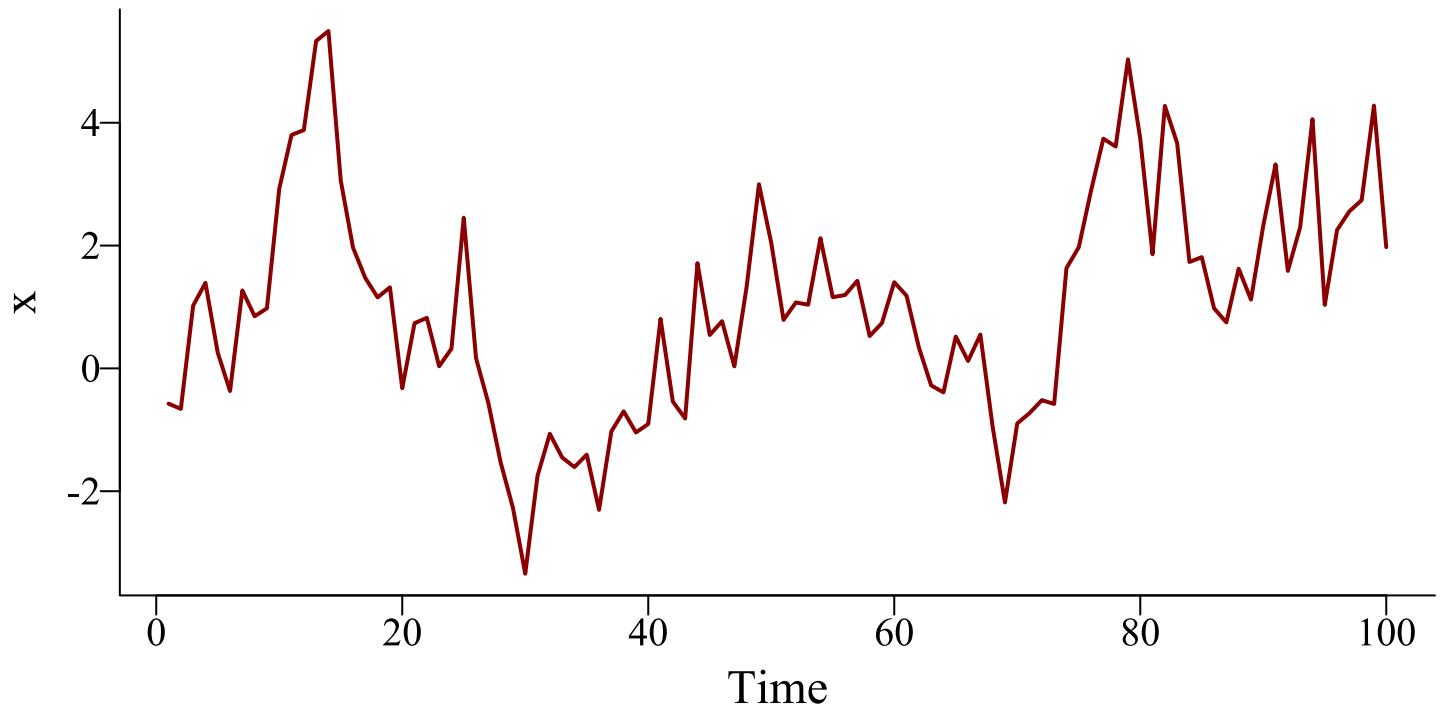
# number of timepoints
T ← 100

# use arima.sim to simulate from a AR1 model (phi = 0.7)
series ← arima.sim(model = list(ar = 0.7), n = T, sd = 1)

# plot the time series as a line
plot(series, type = 'l', bty = 'l', lwd = 2,
      col = 'darkred', ylab = 'x', xlab = 'Time')
```

# Alternative AR1 simulation

Code Plot



# Properties of an AR1

$\phi = 0$  and  $\alpha = 0$ , process is white noise

$\phi = 1$  and  $\alpha = 0$ , process is a Random Walk

$\phi = 1$  and  $\alpha \neq 0$ , process is a Random Walk with drift

$|\phi| < 1$ , process oscillates around  $\alpha$  and is ***stationary***

# Stationarity

*"A stationary time series is one whose statistical properties do not depend on the time at which the series is observed"* ([Hyndman and Athanasopoulos, Forecasting Principles and Practice](#))

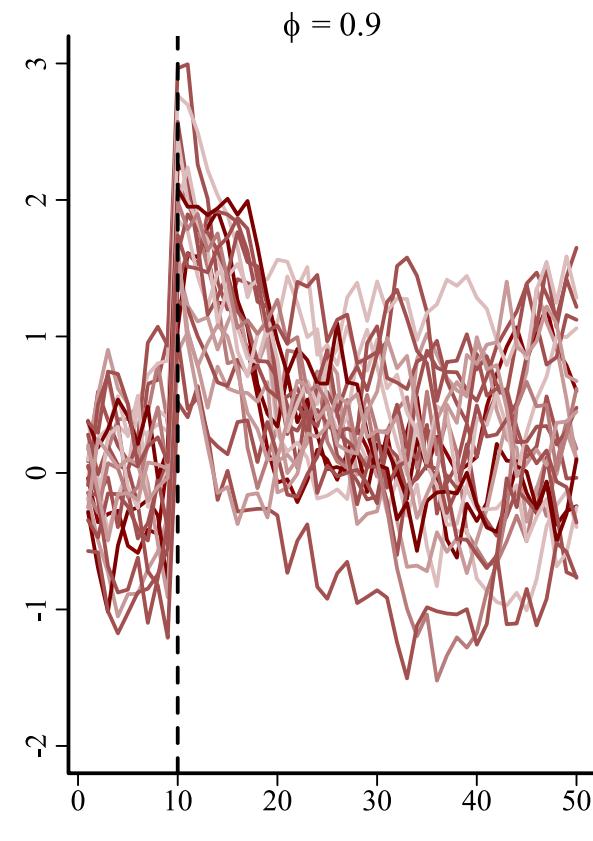
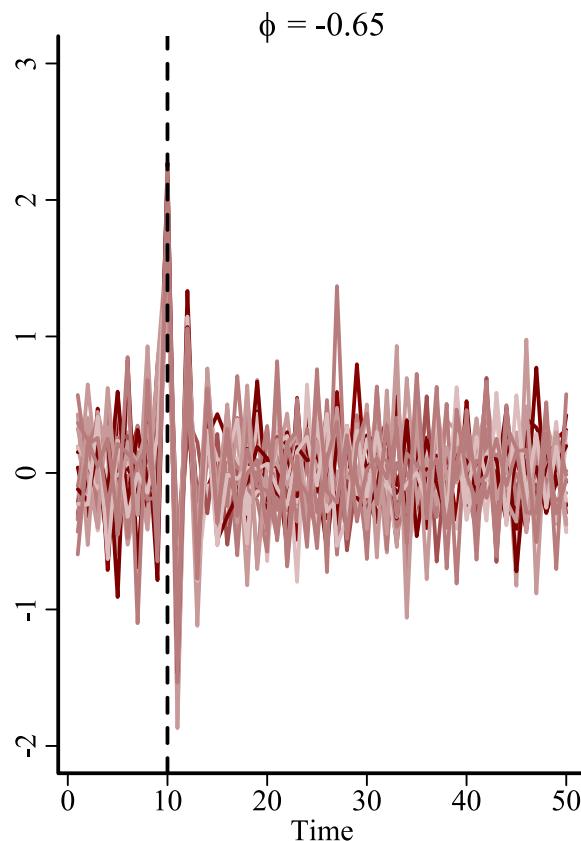
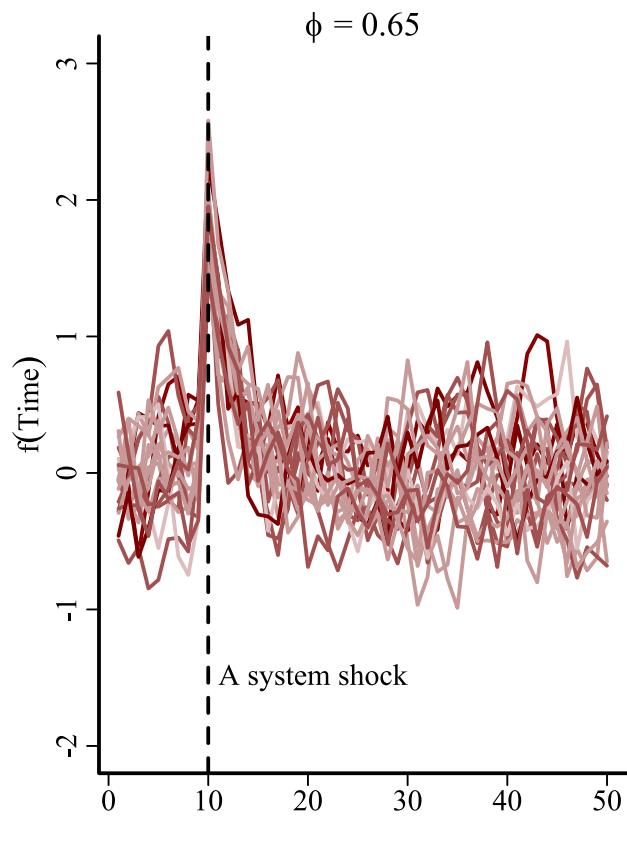
Non-stationary series are more difficult to predict

Either mean, variance, and/or autocorrelation structure can change over time

Random Walk is nonstationary because it has no long-term mean

Stationary time series are useful for inferring properties of **stability**

# Stationarity $\Rightarrow$ stability



# Enforcing AR stationarity

For AR processes up to order 3 (AR1, AR2 or AR3), process remains stationary as long as:

$$|\phi| < 1 \text{ for all } \phi$$

# Enforcing AR stationarity

Enforcing stationarity (by restricting  $|\phi| < 1$ ) is the default behaviour in brms 

```
library(brms)
brm(y ~ x + ... +
     ar(p = 1, time = time, cov = TRUE),
     family = poisson(),
     data = data)
```

# Enforcing AR stationarity

This ***is not*** the default behaviour in `mvgam` .

Instead,  $|\phi| < 1.5$  to allow nonstationary residual processes if they are supported by the data

```
mvgam(y ~ x + ... ,  
       trend_model = 'AR1' ,  
       data = data)
```

# Enforcing AR stationarity

This can be modified by changing upper and lower boundaries in the prior

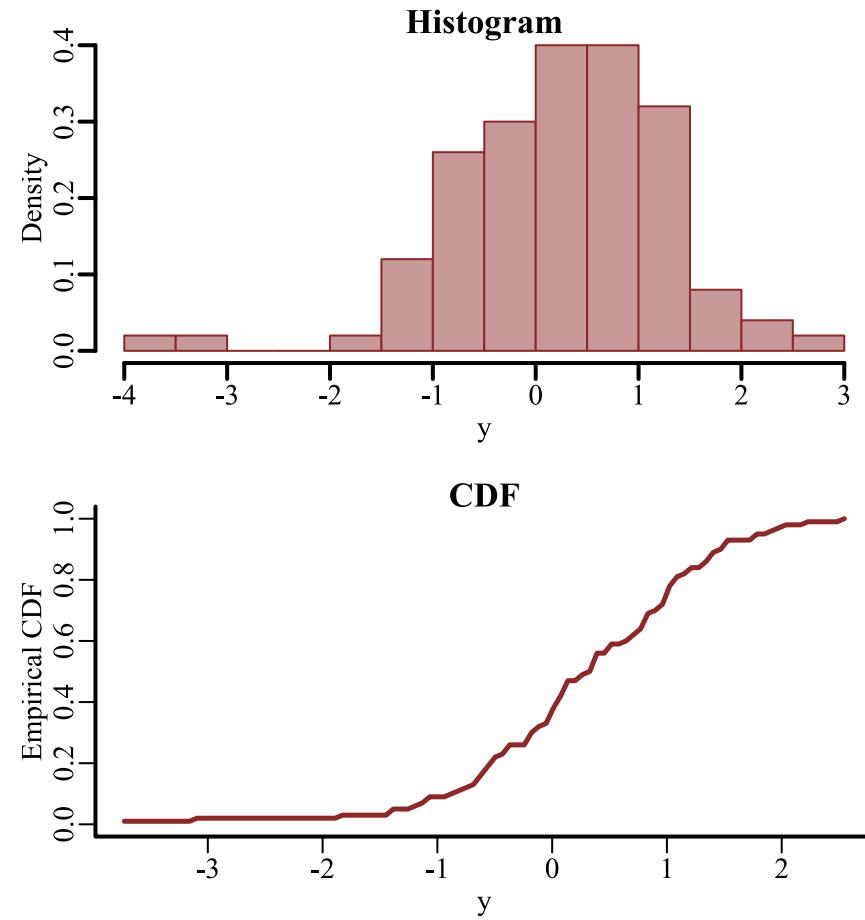
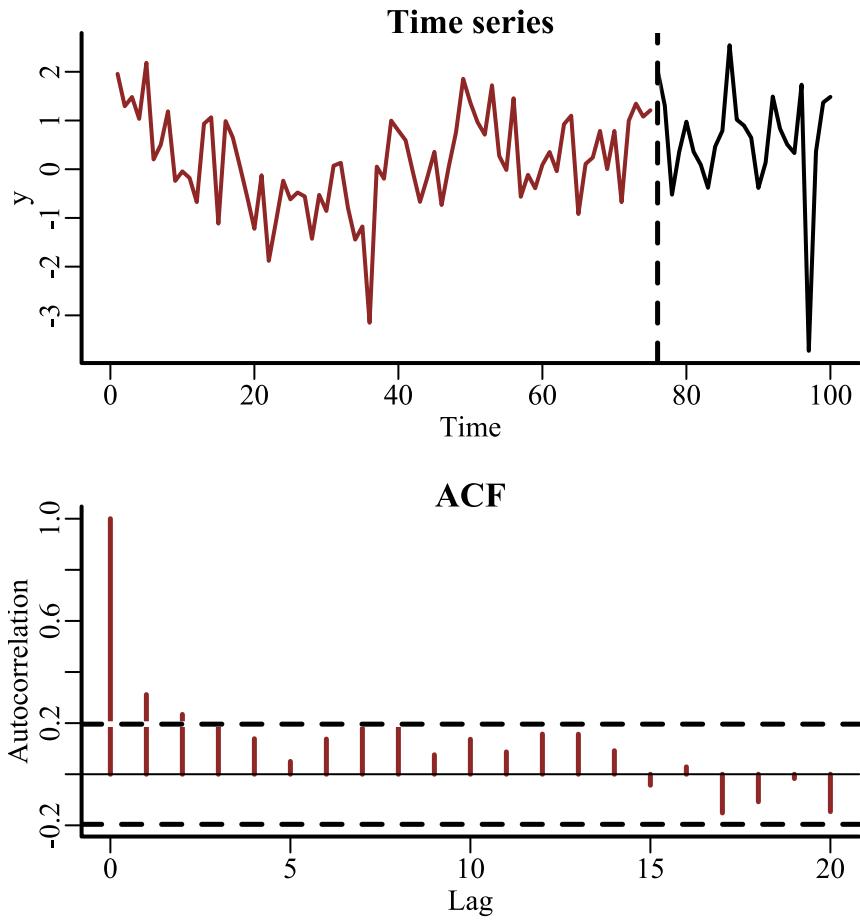
```
priors ← get_mvgam_priors(y ~ x + ... ,  
                           trend_model = 'AR1',  
                           data = data)  
priors$new_lowerbound[1] ← 1  
priors$new_upperbound[1] ← 1  
  
mvgam(y ~ x + ... ,  
       trend_model = 'AR1',  
       priors = priors,  
       data = data)
```

# An example

Simulate a time series with heavy-tailed Student-T observations over a latent Random Walk process that also has seasonal variation

```
set.seed(555)
t_data ← sim_mvgam(T = 100, n_series = 1,
                     family = student_t(),
                     trend_model = 'RW',
                     trend_rel = 0.6)
```

# Simulated data



# Fit RW and AR1 models

```
t_mod_rw ← mgvam(  
    y ~ s(season, bs = 'cc', k = 6),  
    knots = list(season = c(0.5, 12.5)),  
    trend_model = 'RW', family = student_t(),  
    data = t_data$data_train, newdata = t_data$data_test)  
  
t_mod_ar1 ← mgvam(  
    y ~ s(season, bs = 'cc', k = 6),  
    knots = list(season = c(0.5, 12.5)),  
    trend_model = 'AR1', family = student_t(),  
    data = t_data$data_train, newdata = t_data$data_test)
```

In both models, a cyclic smooth captures repeated periodic variation by forcing the smooth to join at the boundaries

# RW MCMC diagnostics ☺

```
## n_eff / iter looks reasonable for all parameters
## Rhat looks reasonable for all parameters
## 0 of 2000 iterations ended with a divergence (0%)
## 0 of 2000 iterations saturated the maximum tree depth of 10 (0%)
## Chain 1: E-FMI = 0.1385
## Chain 2: E-FMI = 0.1088
## Chain 3: E-FMI = 0.0873
## *E-FMI below 0.2 indicates you may need to reparameterize your model
```

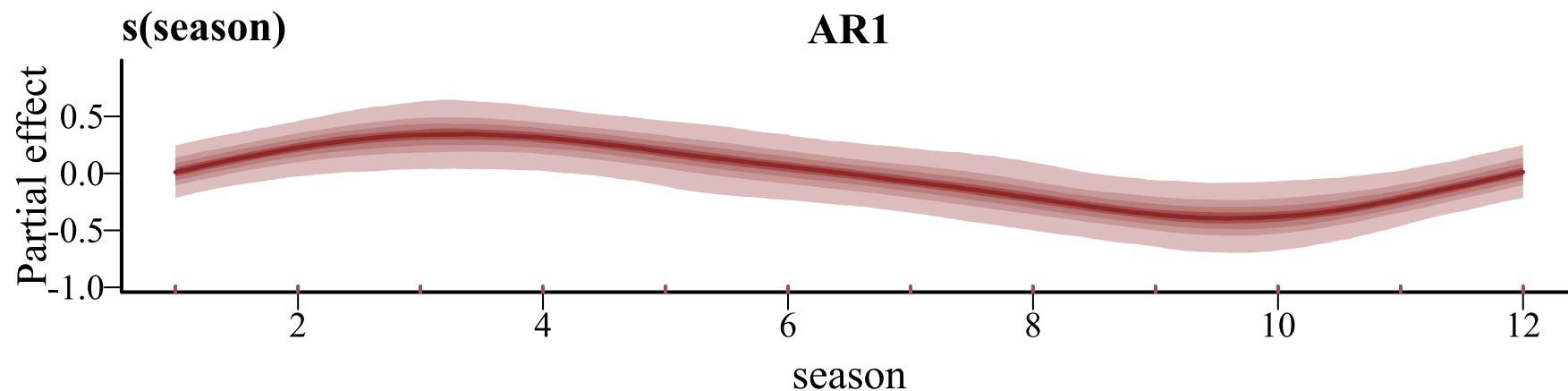
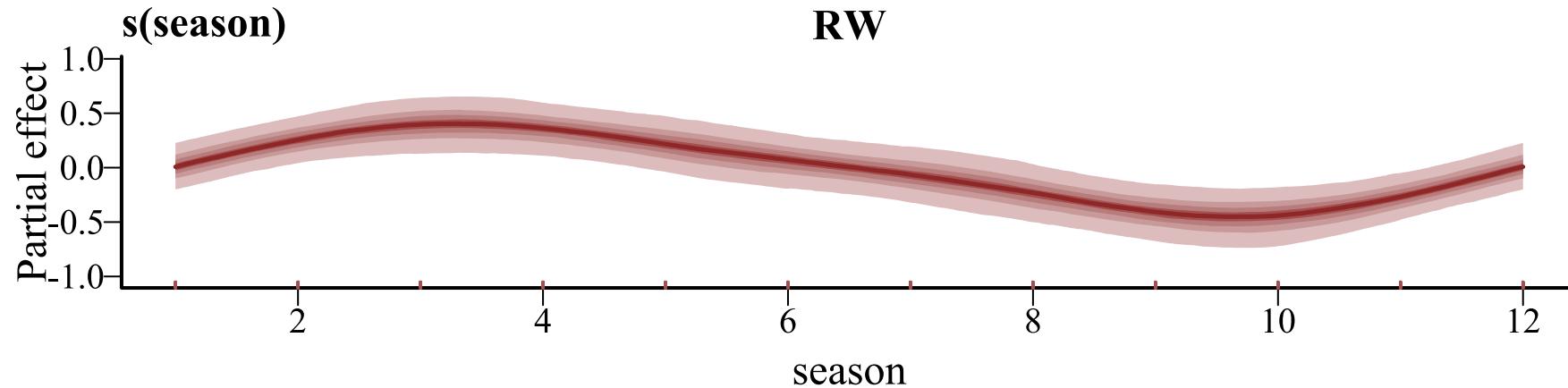
Good convergence, no divergences encountered

# AR1 MCMC diagnostics 😓

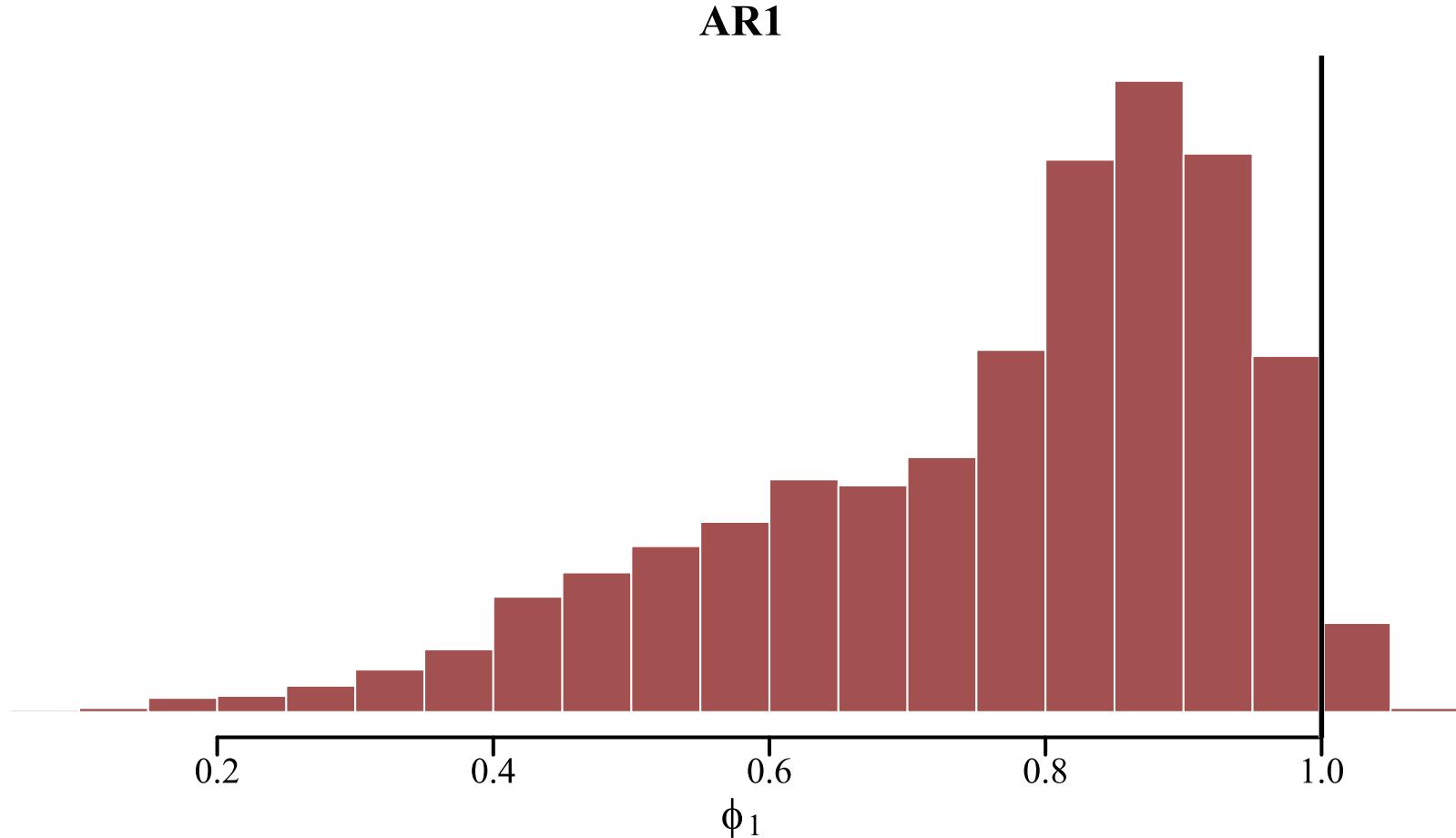
```
## n_eff / iter looks reasonable for all parameters
## Rhats above 1.05 found for 45 parameters
## *Diagnose further to investigate why the chains have not mixed
## 0 of 2000 iterations ended with a divergence (0%)
## 0 of 2000 iterations saturated the maximum tree depth of 10 (0%)
## Chain 1: E-FMI = 0.1215
## Chain 2: E-FMI = 0.1225
## Chain 3: E-FMI = 0.1438
## Chain 4: E-FMI = 0.0758
## *E-FMI below 0.2 indicates you may need to reparameterize your model
```

No divergences encountered, but bad convergence

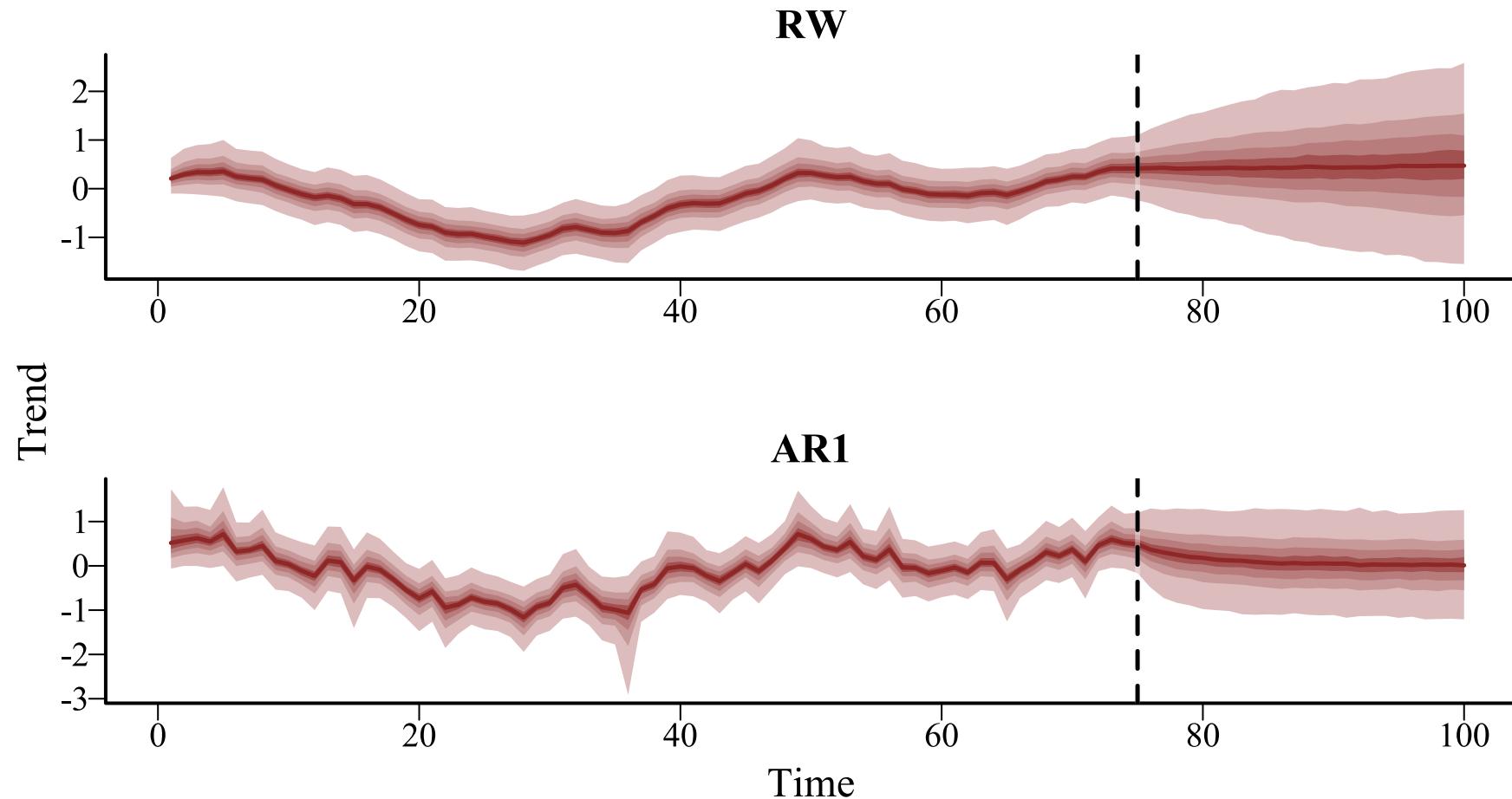
# Similar smooth inferences



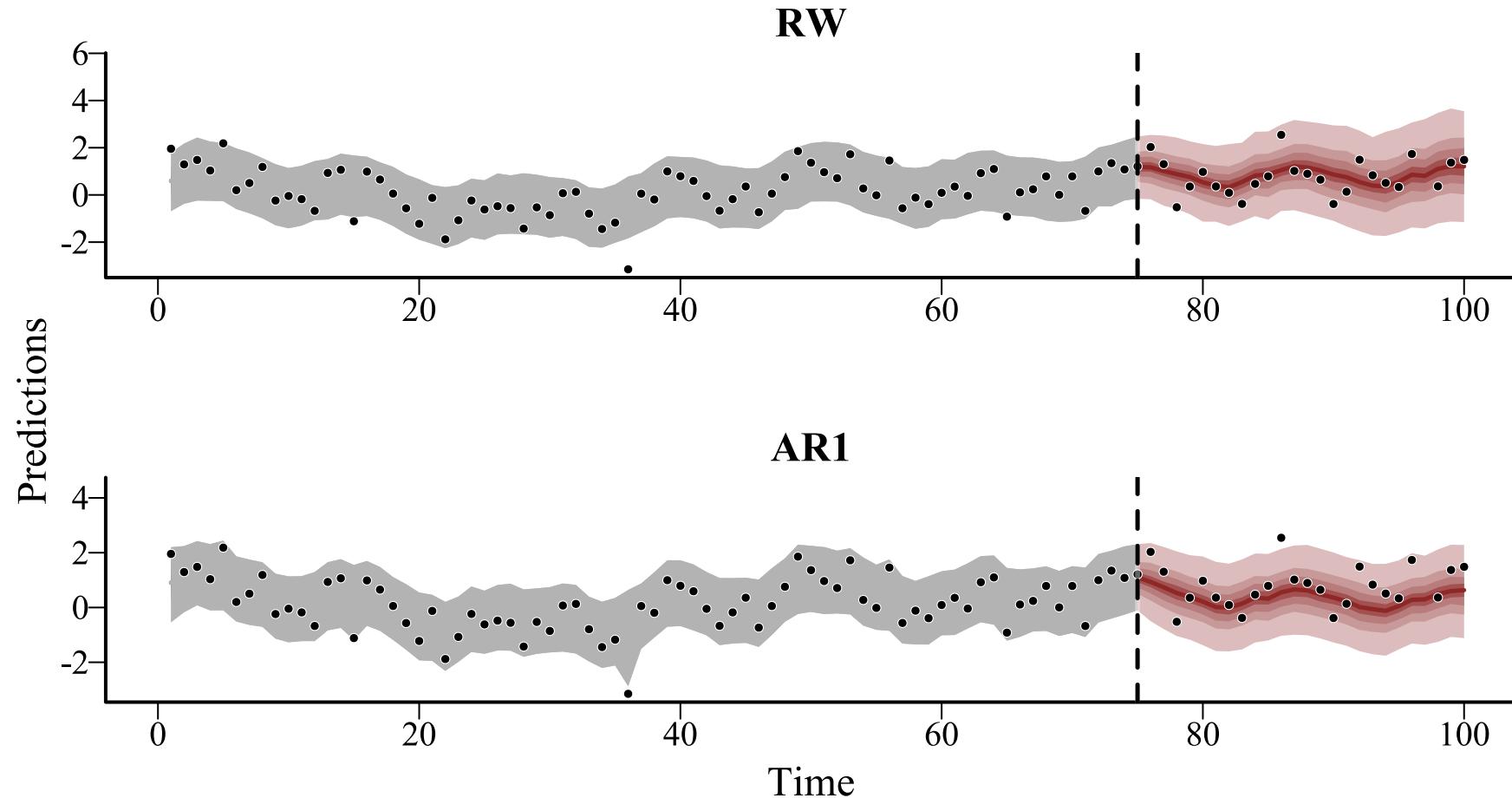
# AR1 mostly stationary ...



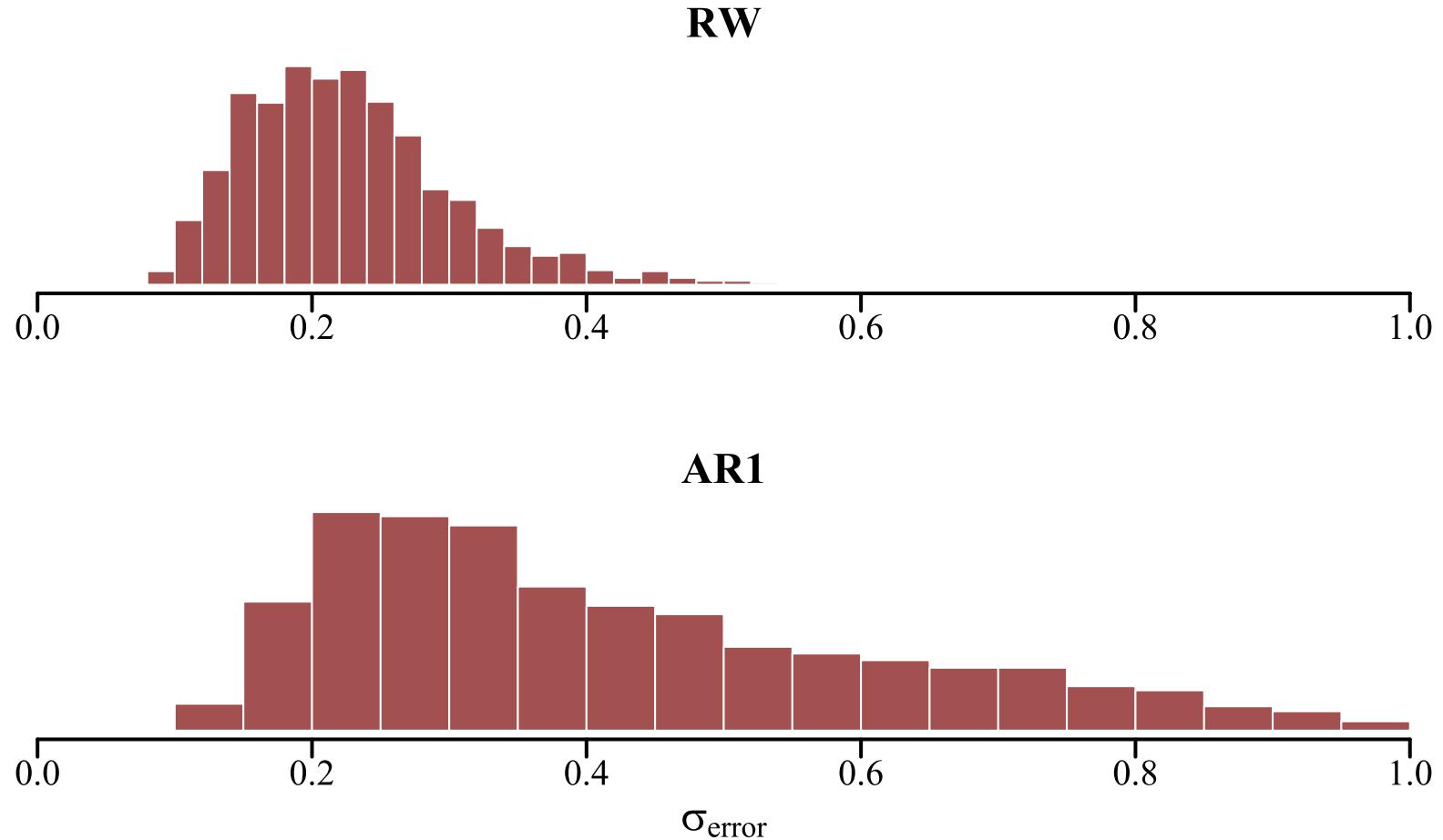
# ... so trends differ ...



# ... as do forecasts



# Process error must compromise



**It is straightforward to fit latent dynamic models with RW or AR models up to order 3 in `mvgam`. Bayesian regularization helps shrink un-needed AR coefficients toward 0**

**In `brms`, only AR1 can be fit for non-Gaussian observations (though can also handle ARMA(1,1)) models. However, implementation is different and much slower**

**But what if we think the latent dynamic process is *smooth*?**

# Gaussian Processes

# Gaussian Processes

*"A Gaussian Process defines a probability distribution over functions; in other words every sample from a Gaussian Process is an entire function from the covariate space  $X$  to the real-valued output space."* (Betancourt; [Robust Gaussian Process Modeling](#))

$$z \sim \text{MVNormal}(0, \Sigma)$$

$$\Sigma_{t_i, t_j} = \alpha^2 * \exp(-0.5 * ((|t_i - t_j|/\rho))^2)$$

Where:

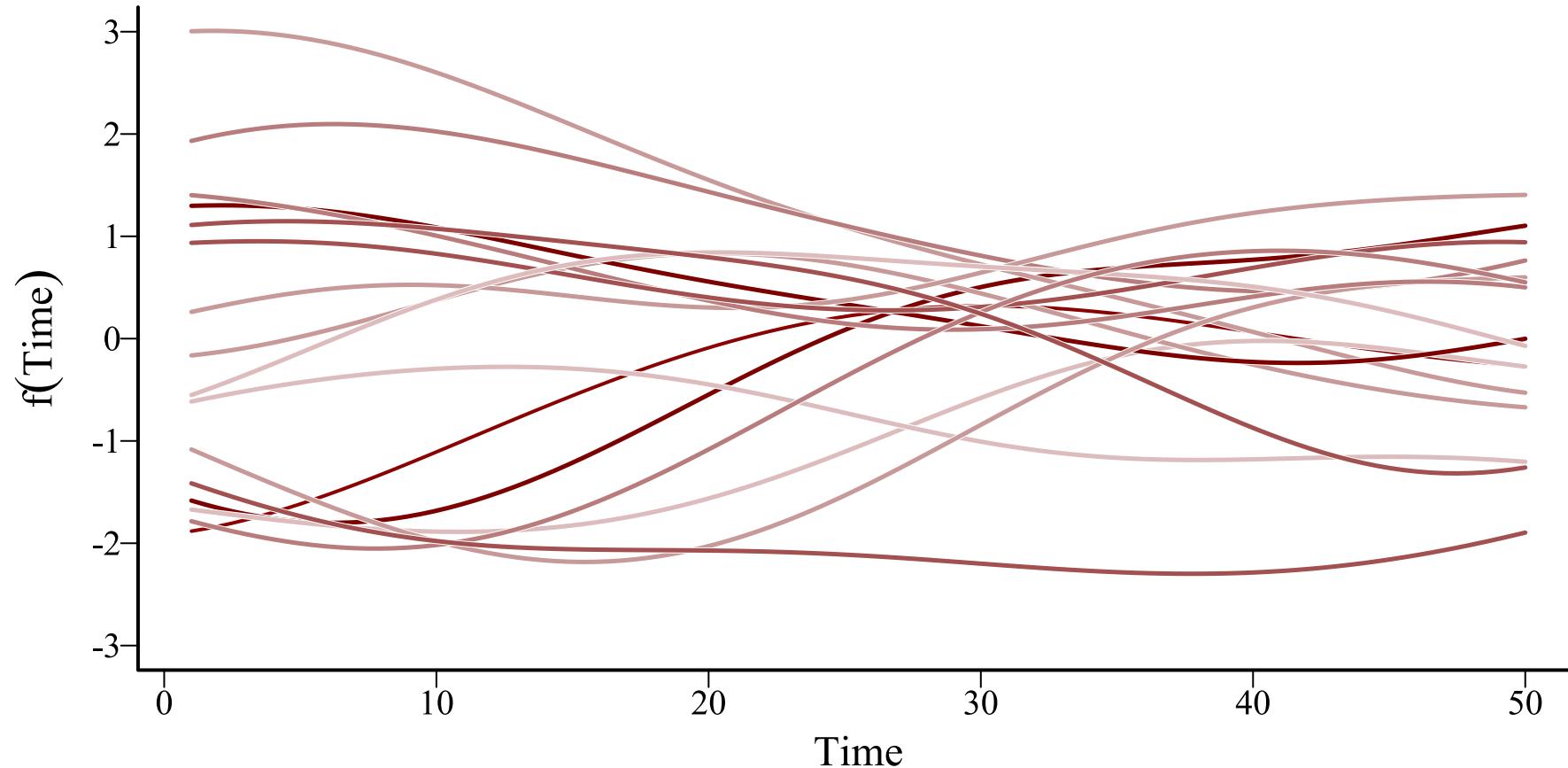
$\alpha$  controls the marginal variability (magnitude) of the function

$\rho$  controls how correlations decay as a function of time lag

$\Sigma$  is the kernel, in this case a squared exponential kernel

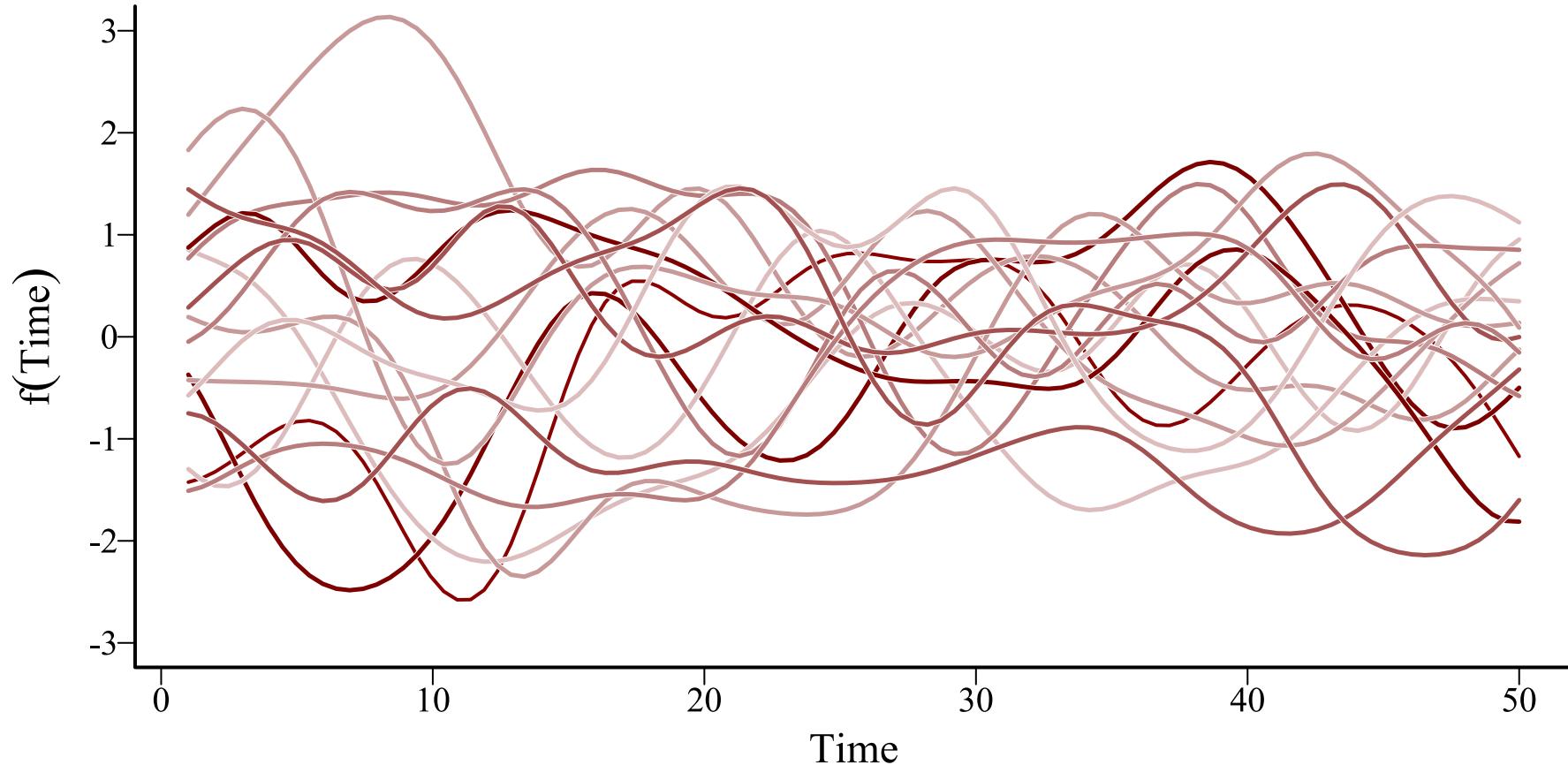
# Random functions

$\alpha = 1; \rho = 16$



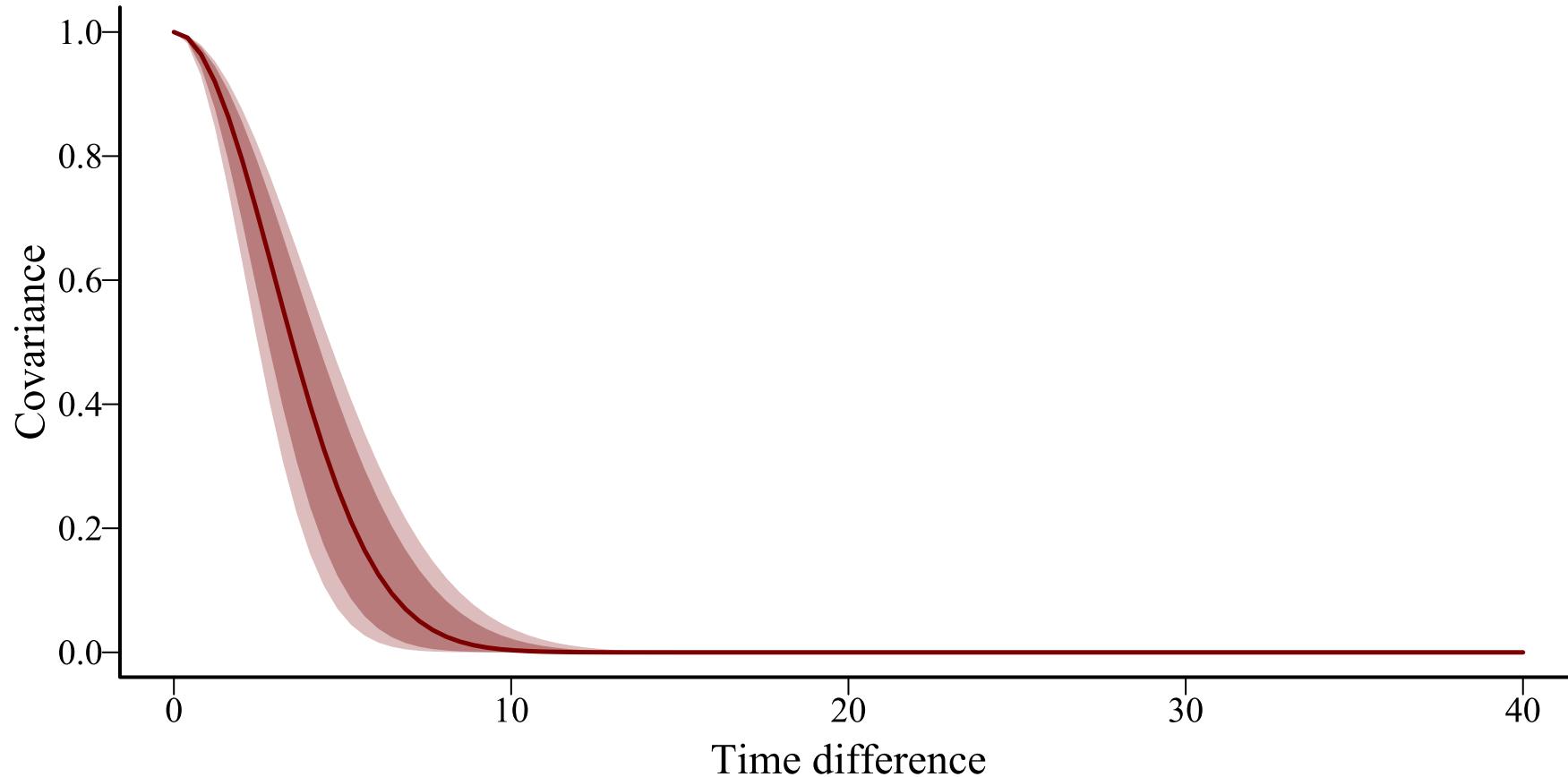
# Length scale $\Rightarrow$ memory

$$\alpha = 1; \rho = 4$$

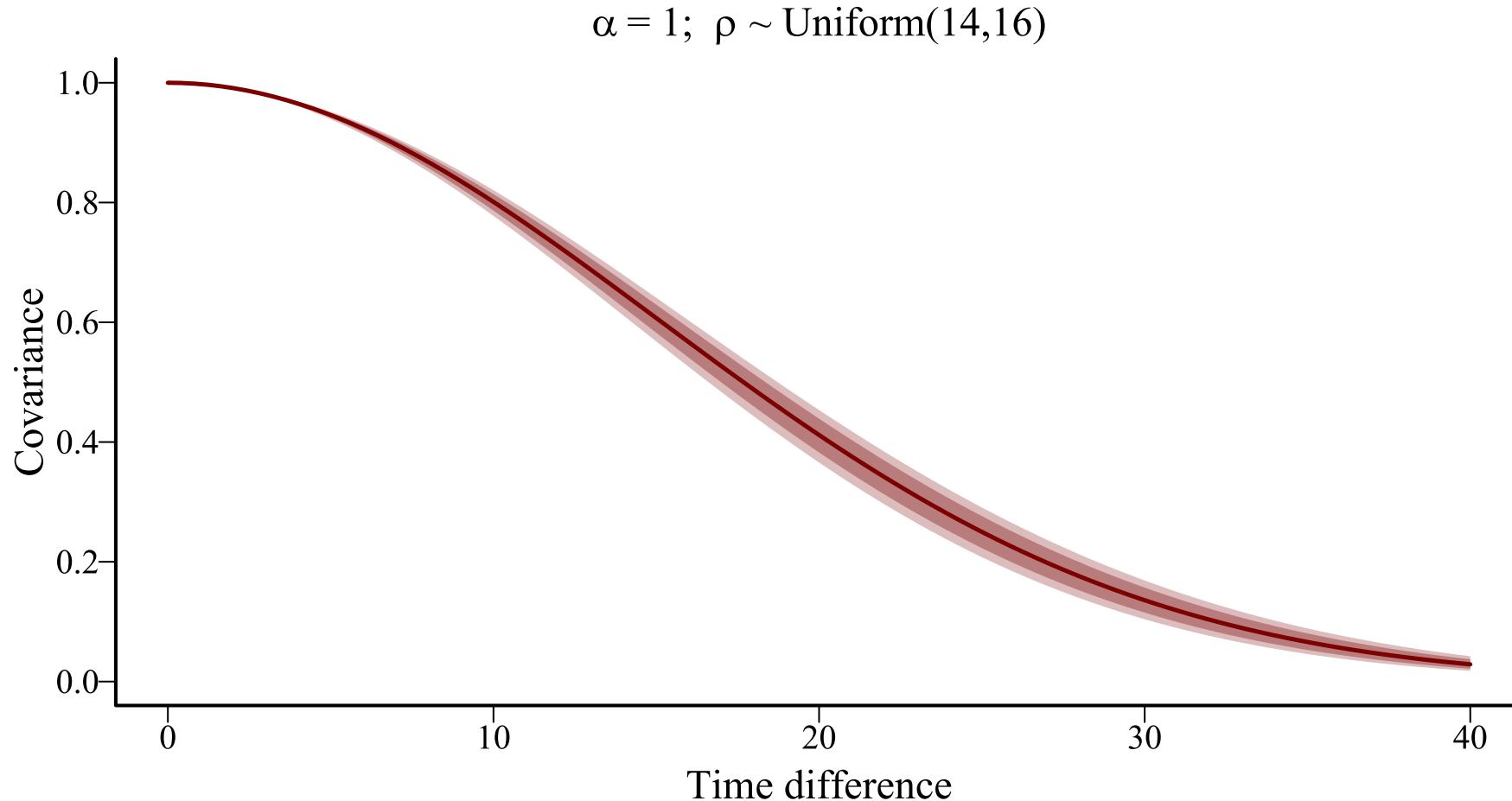


# Kernel $\Rightarrow$ covariance decay

$\alpha = 1; \rho \sim \text{Uniform}(2,4)$

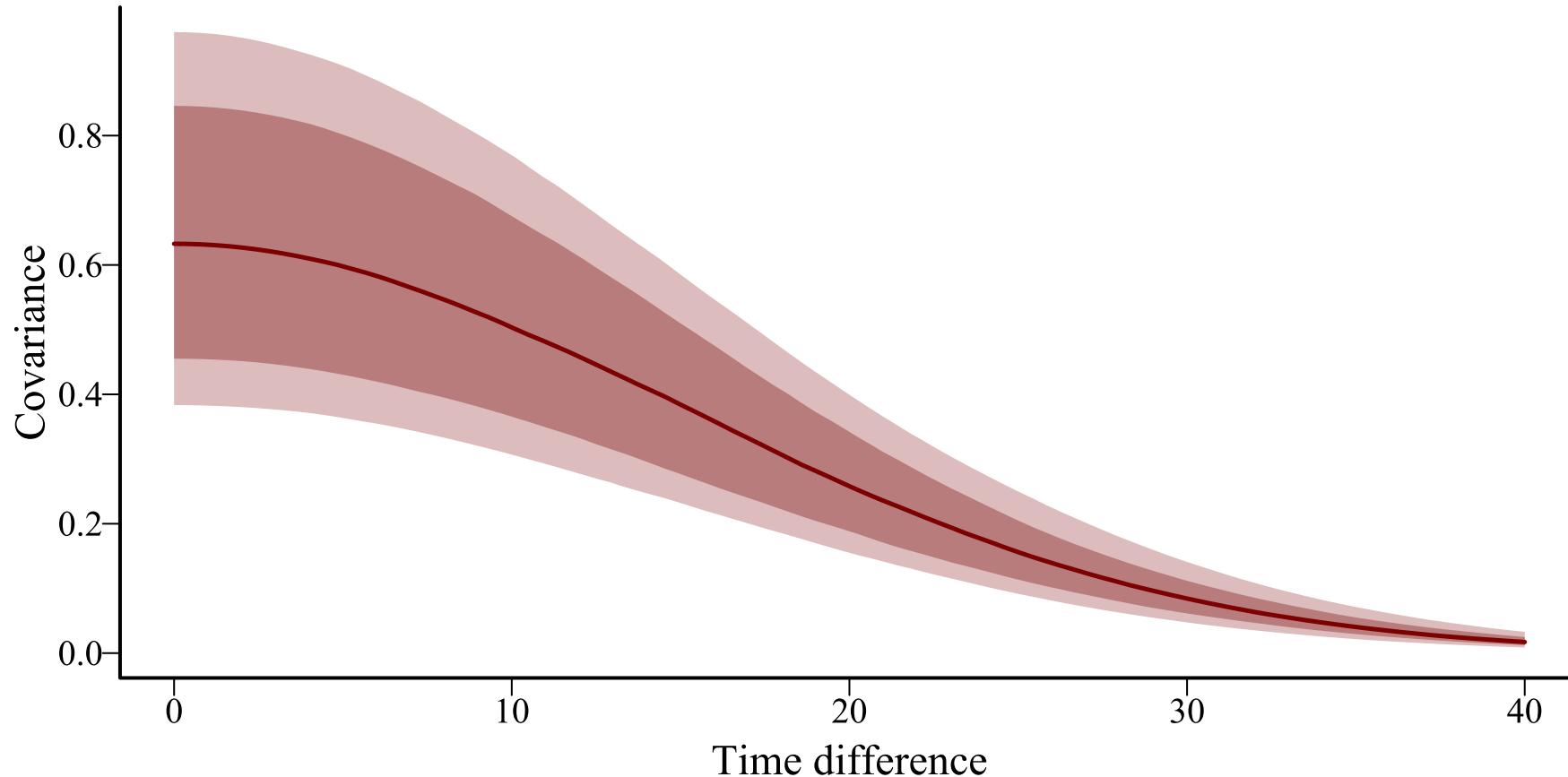


# Kernel $\Rightarrow$ covariance decay

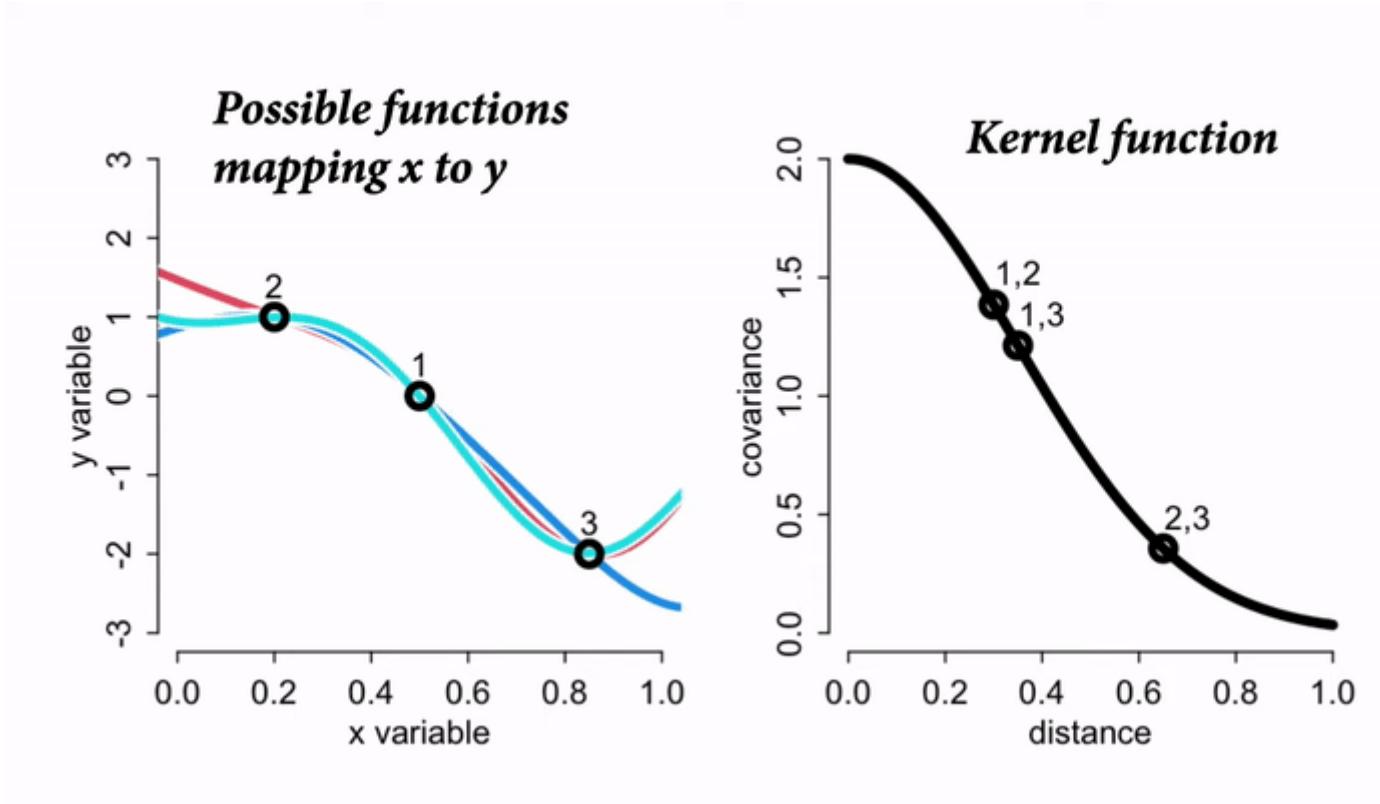


# Kernel $\Rightarrow$ covariance decay

$\alpha = \text{Uniform}(0.6,1); \rho \sim \text{Uniform}(14,16)$



# Kernel smoothing in action

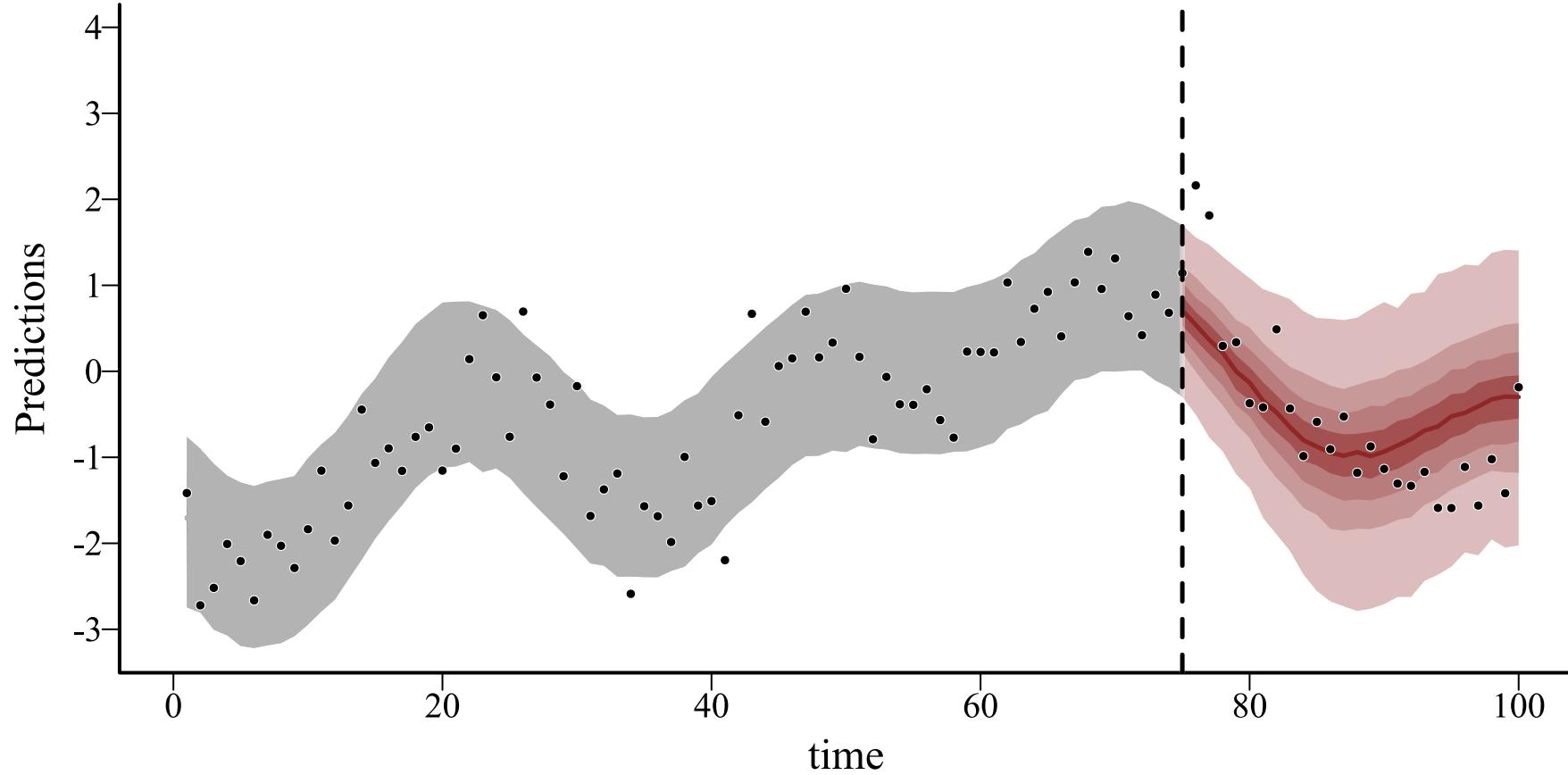


**A latent GP allows prediction for *any* time point because all we need is the distance to each training time point**

**The cross-covariance for prediction vs training time points provides the kernel used to extend functions forward in time**

**Allows GPs to make much better predictions than splines, but at a high computational cost**

# Global knowledge ✓



# Approximating GPs

A quick note that both the `mvgam` and `brms` 's can employ an approximation method to improve computational efficiency for estimating Gaussian Process parameters

Relies on basis expansions to reduce dimensionality of the problem

Details not focus of this lecture, but can be found in this reference

Riutort-Mayol et al 2023; Practical Hilbert space approximate Bayesian Gaussian processes for probabilistic programming

Both packages use automatic, informative priors for length scales  $\rho$ , but these can be changed (more on this in Tutorial 2)

# Estimation in brms

In `brms` 📦, use the gp function with `time` as the covariate

```
brm(y ~ x + ... +
      gp(time, c = 5/4, k = 20, scale = FALSE),
      family = poisson(),
      data = data)
```

Requires arguments to determine behaviour of the approximation (`c` and `k`). Good defaults are `5/4` and `20`, but depends on number of timepoints and expected smoothness

# Estimation in mvgam

In `mvgam` 📦, use `trend_model = 'GP'`

```
mvgam(y ~ x + ... +
       trend_model = "GP",
       family = poisson(),
       data = data)
```

Doesn't require arguments, but the number of basis functions `k` can be altered by changing the priors

No examples here as we will go deeper into GPs in the tutorial

But if you want extra detail, watch this lecture: - Statistical  
Rethinking 2023 - 16 - Gaussian Processes

# Dynamic coefficient models

# Dynamic coefficients

Major advantage of flexible interfaces such as `brms`, `mgcv` and `mvgam`'s is ability to handle many types of nonlinear effects

These can include smooth functions of covariates, as we have been using so far

But they can also include other types of nonlinearities

Spatial autocorrelation functions

Distributed lag functions

***Time-varying effects***

# Smooth time-varying effects

If a covariate effect changes over time, we'd usually expect this change to be *smooth*

Splines and Gaussian Processes provide useful tools to estimate these effects

But as we've seen previously, splines will often give poor predictions about how effects will change in the future

# In mvgam

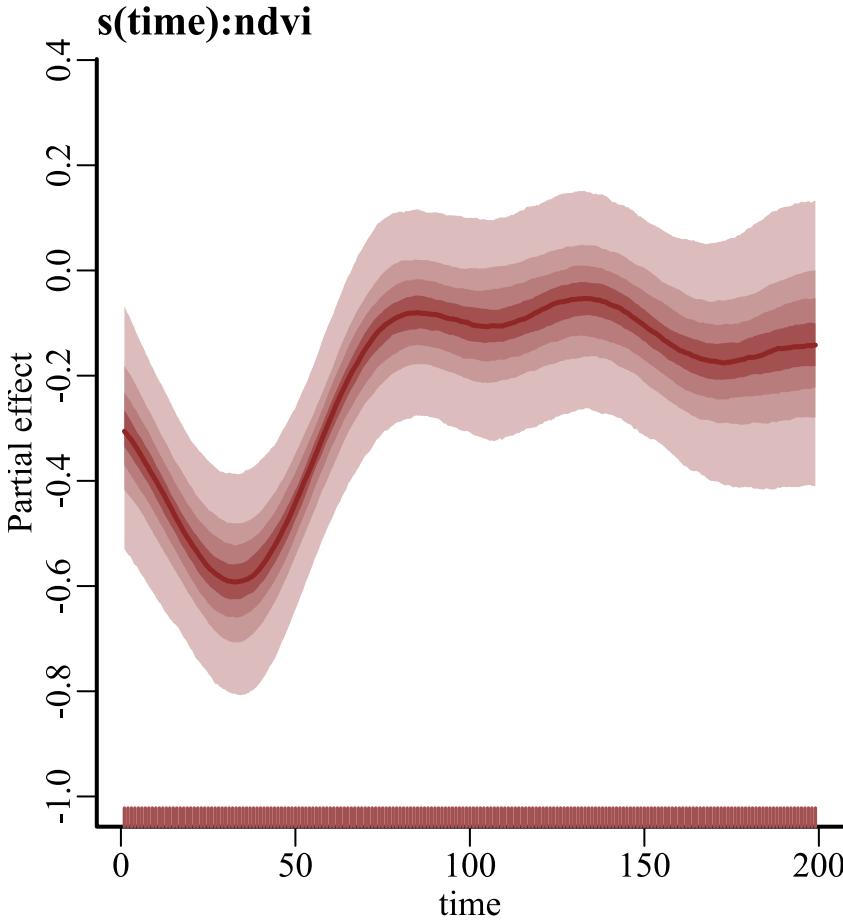
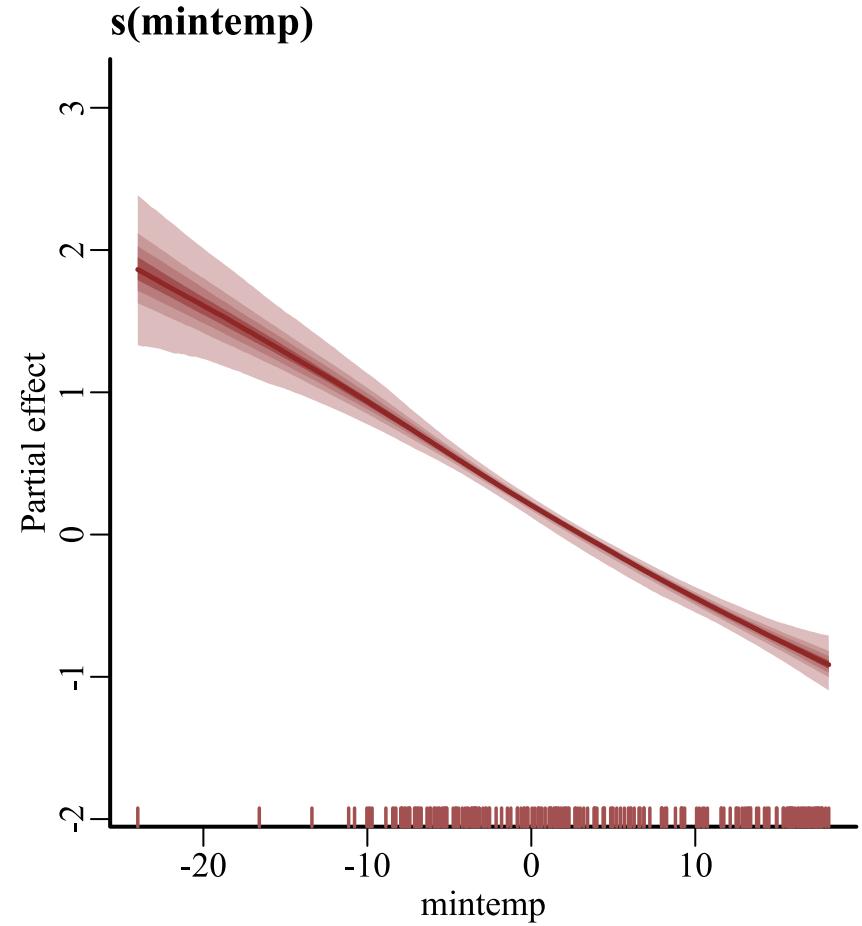
In `mvgam` 📦, use `dynamic` to set up time-varying effects

```
mod_beta_dyn ← mvgam(relabund ~ s(mintemp, k = 6) +  
                      dynamic(ndvi, rho = 28),  
                      family = betar(),  
                      data = dm_data)
```

Requires user to set the length scale  $\rho$ , as the function is approximated using a low-rank GP smooth from the `mgcv` 📦

No uncertainty in the GP parameters will be estimated, but function will resemble a squared exponential GP

# Estimated smooths



# Predicted effects

---

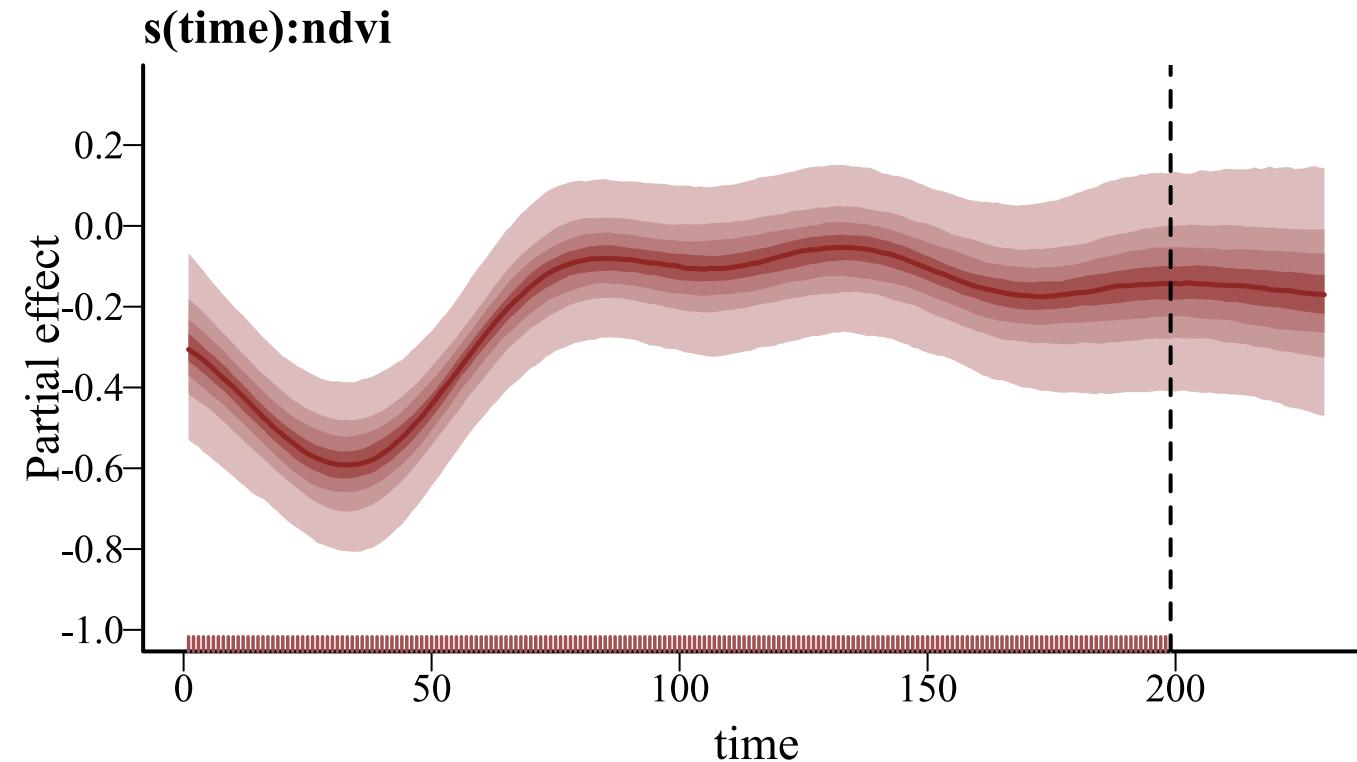
Code    Plot

---

```
# use mvgam's plot_mvgam_smooth to view predicted effects
plot_mvgam_smooth(mod_beta_dyn, smooth = 2,
                    # datagrid from marginaleffects is useful
                    # to set up prediction scenarios
                    newdata = datagrid(time = 1:230,
                                        model = mod_beta_dyn))
abline(v = max(dm_data$time), lwd = 2, lty = 'dashed')
```

# Predicted effects

Code Plot



# In brms

In `brms` 📦, use `gp` with the `by` argument

```
brm_beta_dyn ← brm(relabund ~ s(mintemp, k = 6) +  
                     gp(time, by = ndvi, c = 5/4, k = 20),  
                     family = Beta(),  
                     data = dm_data,  
                     chains = 4,  
                     cores = 4,  
                     backend = 'cmdstanr')
```

A GP specifying time-varying effects of `ndvi`. This estimates a full GP, so is more flexible than the `mvgam` / `mgcv` version

# Time-vaying effect

---

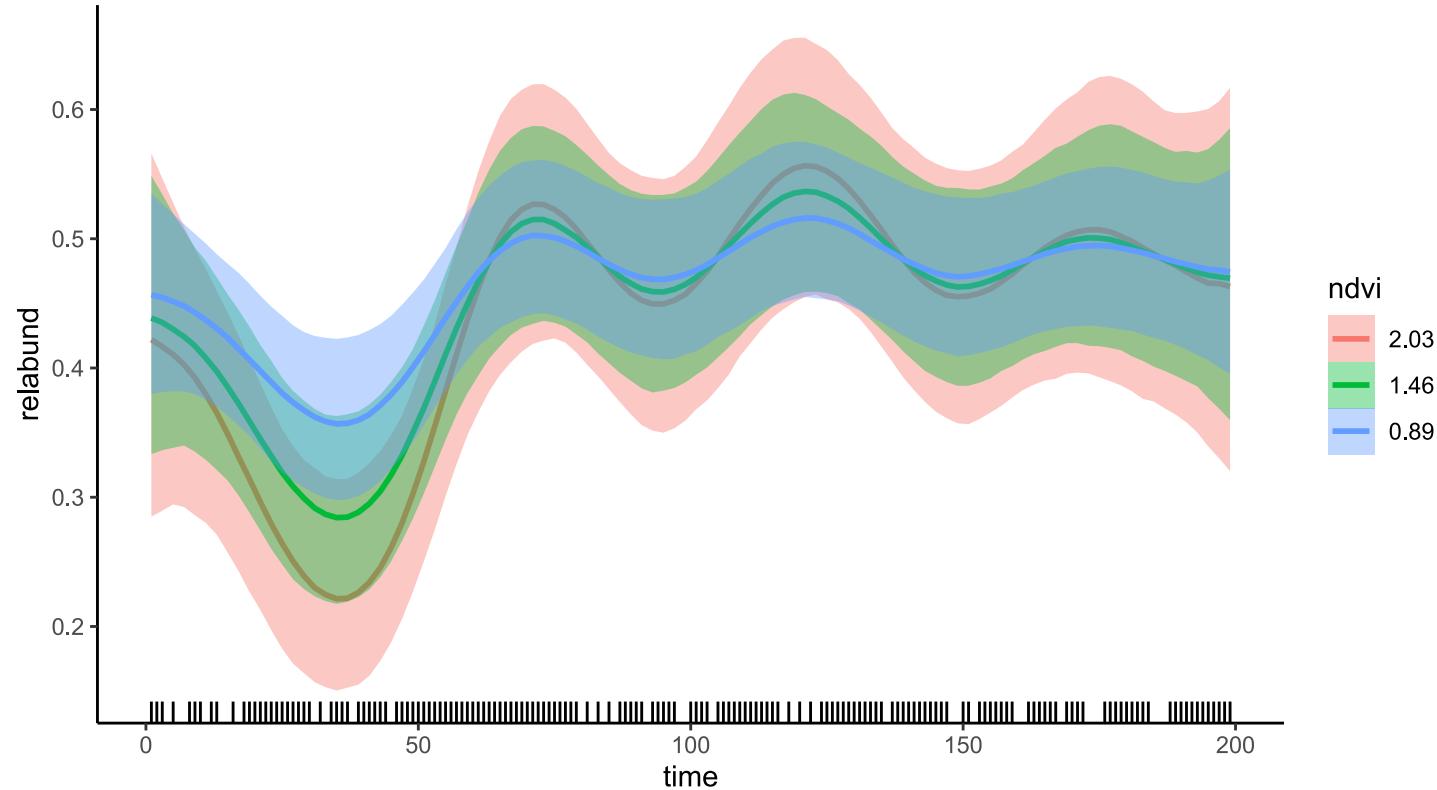
Code    Plot

---

```
# use brms' conditional_effects to view predictions
plot(conditional_effects(brm_beta_dyn, effects = c('time:ndvi')),
      theme = theme_classic(),
      mean = FALSE,
      rug = TRUE)
```

# Time-vaying effect

Code Plot



**We have seen many ways to handle dynamic components in Bayesian regression models**

**These flexible processes can capture time-varying effects and give realistic forecasts, while also allowing us to respect the properties of the observations**

**But how do we evaluate and compare dynamic GAMs / GLMs?**

# In the next lecture, we will cover

Forecasting from dynamic models

Bayesian posterior predictive checks

Point-based forecast evaluation

Probabilistic forecast evaluation