Speeding up parameter inference with compiled models

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Differential equation (DE) models in R are easy to implement and allow simple interactive development using readable code and access to R's many high-level functions. This was part of our motivation to develop deBInfer as an R based inference package. However, numerically solving DE models specified as R functions is also relatively slow. The deSolve package also allows the evaluation of DE models that are defined in lower-level languages such as C and FORTRAN. These compiled models have the benefit of increased simulation speed. As the DE model is evaluated many times during the MCMC procedure, even moderate speed-ups from using compiled models can result in large absolute time savings.

We demonstrate the speed-up using an example from the deSolve documentation (Soetaert et al. 2010). Full details on the model specification can be found in (Soetaert et al. 2009) which can be displayed with the command vignette("compiledCode"). Further details on the set up of the deBInfer inference procedure are described in (Boersch-Supan & Johnson 2016) and annotated examples are available in the vignettes vignette("logistic_ode_example") and vignette("vignette_chytrid_dede_example").

1 Specifying the ODE model

Following Soetaert et al. (2009), ee use the following simple ODE:

$$\begin{aligned} \frac{dy_1}{dt} &= -k_1 \cdot y_1 + k_2 \cdot y_2 \cdot y_3 \\ \frac{dy_2}{dt} &= k_1 \cdot y_1 - k_2 \cdot y_2 \cdot y_3 - k_3 \cdot y_2 \cdot y_2 \\ \frac{dy_3}{dt} &= k_3 \cdot y_2 \cdot y_2 \end{aligned}$$

where y_1 , y_2 and y_3 are state variables, and k_1 , k_2 and k_3 are parameters.

1.1 ODE model implementation in R

We implement this model as an R function:

```
modelR <- function(t, Y, parameters) {
    with (as.list(parameters), {

        dy1 = -k1*Y[1] + k2*Y[2]*Y[3]
        dy3 = k3*Y[2]*Y[2]
        dy2 = -dy1 - dy3

        list(c(dy1, dy2, dy3))
    })
}</pre>
```

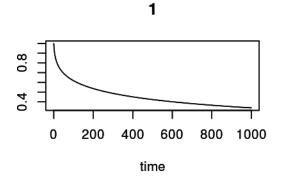
And also specify the Jacobian $(\frac{\partial y'}{\partial y})$ for it:

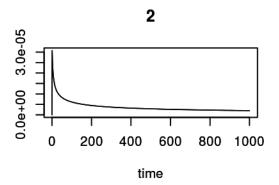
```
jacR <- function (t, Y, parameters) {
    with (as.list(parameters),{

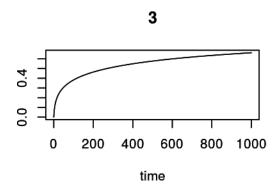
        PD[1,1] <- -k1
        PD[1,2] <- k2*Y[3]
        PD[1,3] <- k2*Y[2]
        PD[2,1] <- k1
        PD[2,3] <- -PD[1,3]
        PD[3,2] <- k3*Y[2]
        PD[2,2] <- -PD[1,2] - PD[3,2]

    return(PD)
    })
}</pre>
```

This model can then be run as follows:







1.2 ODE model implementation in C

To create compiled models (dynamic-link libraries (.dll) on Windows or shared objects (.so) on other systems) a recent version of the GNU compiler suite is required. This is usually the case for Linux systems. Windows users can install the required toolchain by following the instructions on [https://cran.r-project.org/bin/windows/Rtools/]. OSX users need to download and install an appropriate version of Xcode from the Apple developer website [https://developer.apple.com/] or the OSX App Store.

The call to the derivative and Jacobian function is more complex for compiled code compared to R-code, because it has to comply with the interface needed by the integrator. The requirements for this are detailed in Soetaert et al. (2009). A C implementation of the example model is found below.

```
/* file mymod.c */
#include <R.h>
static double parms[3];
#define k1 parms[0]
#define k2 parms[1]
#define k3 parms[2]
/* initializer */
void initmod(void (* odeparms)(int *, double *))
{
    int N=3;
    odeparms(&N, parms);
}
/* Derivatives and 1 output variable */
void derivs (int *neq, double *t, double *y, double *ydot,
             double *yout, int *ip)
{
    if (ip[0] <1) error("nout should be at least 1");</pre>
    ydot[0] = -k1*y[0] + k2*y[1]*y[2];
    ydot[2] = k3 * y[1]*y[1];
    ydot[1] = -ydot[0]-ydot[2];
    yout[0] = y[0]+y[1]+y[2];
}
/* The Jacobian matrix */
void jac(int *neq, double *t, double *y, int *ml, int *mu,
           double *pd, int *nrowpd, double *yout, int *ip)
 pd[0]
                      = -k1;
 pd[1]
                      = k1;
 pd[2]
                      = 0.0;
  pd[(*nrowpd)]
                      = k2*y[2];
  pd[(*nrowpd) + 1]
                      = -k2*y[2] - 2*k3*y[1];
  pd[(*nrowpd) + 2]
                      = 2*k3*y[1];
 pd[(*nrowpd)*2]
                      = k2*y[1];
 pd[2*(*nrowpd) + 1] = -k2 * y[1];
  pd[2*(*nrowpd) + 2] = 0.0;
/* END file mymod.c */
```

2 Running ODE models implemented in compiled code

To run the C implementation of the model it must first be compiled. This can be done using the R system statement

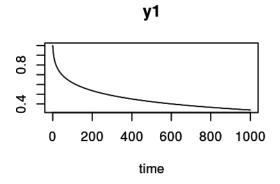
```
system("R CMD SHLIB mymod.c")
```

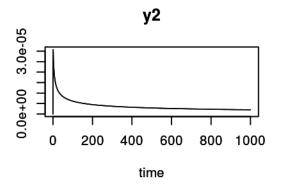
which will create the file mymod.dll on Windows, or mymod.so on other platforms.

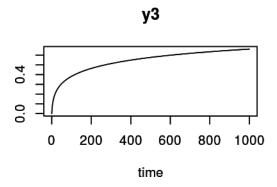
We can then load the compiled model with

```
dyn.load(paste("mymod", .Platform$dynlib.ext, sep = ""))
```

The model can now be run as follows:







We can determine the speed improvement as a function of desired output time points using the microbenchmark package

```
timepoints <- c(10,100, 500, 1000,2000, 5000, 10000)
solve_time <- lapply(timepoints, function(x){</pre>
                                            times \leftarrow seq(0, 0.1*10<sup>4</sup>, length.out = x)
                                            comp <- microbenchmark::microbenchmark(</pre>
                                                 C = ode(Y, times, func = "derivs", parms = parms,
                                                                         jacfunc = "jac", dllname = "mymod",
                                                                         initfunc = "initmod", nout = 1, outnames = "Sum"),
                                                 R = ode(Y, times, modelR, parms = parms, jacfunc = jacR), times=10
                                                 return(comp)
                                                 }
                                            )
#saveRDS(solve_time, "examples/Soetaert-r-vs-c-solve_time.RDS")
solve_speedup <- sapply(solve_time, function(x)mean(x$time[x$expr=="R"]/x$time[x$expr=="C"]))</pre>
solve_R <- sapply(solve_time, function(x)mean(x$time[x$expr=="R"]/1e9))</pre>
solve_C <- sapply(solve_time, function(x)mean(x$time[x$expr=="C"]/1e9))</pre>
#pdf("examples/c-vs-r-solve.pdf")
par(mfrow=c(1,2))
plot(timepoints, solve_R, type="p", ylab = "Execution time (s)", xlab = "Time points")
points(timepoints, solve_C, col="red", type="p", pch=c(16))
legend("topleft", legend = c("R ODE", "C ODE"), pch = c(1,16), col= c("black", "red"))
plot(rep(timepoints, each=1), solve_speedup, type="p", ylab = "Relative execution time R/C", xlab = "Time" = "Relative execution time R/C", xlab = "Time" = 
                                                                                                    0
                                                                                                                                                                                           0
                                                                                                                                                                                                                              0
                                 R ODE
                                                                                                                         Relative execution time R/C
                                        C ODE
                                                                                                                                                                      0
                                                                                                                                         35
Execution time (s)
               9
               o.
                                                                                                                                                               0
                                                                                                                                         30
                                                                 0
              0.4
              0.2
                                                                                                                                         25
                                            0
                                00
                              0
                                                     4000
                                                                                8000
                                                                                                                                                       0
                                                                                                                                                                              4000
                                                                                                                                                                                                          8000
                                                  Time points
                                                                                                                                                                            Time points
```

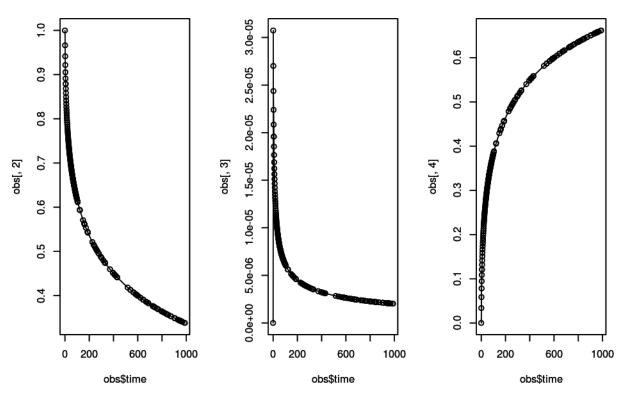
which shows us that the C version of the models is solved 21-39 times faster than the R version.

#dev.off()

3 Conducting inference with compiled models

We can now simulate a dataset to demonstrate the inference procedure:

```
set.seed(143)
#force include the first time-point (t=0)
obs <- as.data.frame(outR[c(1:100,runif(100, 0, nrow(outR))),])
obs <- obs[order(obs$time),]
par(mfrow=c(1,3))
plot(obs$time, obs[,2], type='o')
plot(obs$time, obs[,3], type='o')
plot(obs$time, obs[,4], type='o')</pre>
```



4 Defining an observation model and parameters for inference

For simplicity we assume a normal log-likelihood for these data

$$\ell(\mathcal{Y}|\boldsymbol{\theta}) = \sum_{i} \sum_{t} \ln \left(\frac{1}{\sigma_{obs} \sqrt{2\pi}} \exp\left(-\frac{(\tilde{y}_{t,i} - (y_{t,i})^2}{2\sigma_{obs}^2}\right) \right)$$
 (1)

where $\tilde{y}_{t,i}$ are the observations, and $y_{t,i}$ are the predictions of the DE model given the current MCMC sample of the parameters $\boldsymbol{\theta}$.

```
# the observation model
obs_model <- function(data, sim.data, samp){

llik.y1 <- sum(dnorm(obs[,2], mean = sim.data[,2], sd = samp[['sd.y1']], log = TRUE))</pre>
```

```
llik.y2 <- sum(dnorm(obs[,3], mean = sim.data[,3], sd = samp[['sd.y2']], log = TRUE))
llik.y3 <- sum(dnorm(obs[,4], mean = sim.data[,4], sd = samp[['sd.y3']], log = TRUE))
return(llik.y1 + llik.y2 + llik.y3)
}</pre>
```

We declare the DE model parameter r, assign a prior $r \sim \mathcal{N}(0,1)$ and a random walk sampler with a Normal kernel (samp.type="rw") and proposal variance of 0.005. Similarly, we declare $K \sim \ln \mathcal{N}(1,1)$ and $\ln(\sigma_{obs}^2) \sim \mathcal{N}(0,1)$. Note that we are using the asymmetric uniform proposal distribution $\mathcal{U}(\frac{a}{b}\theta^{(k)}, \frac{b}{a}\theta^{(k)})$ for the variance parameter (samp.type="rw-unif"), as this ensures strictly positive proposals.

```
library(deBInfer)
k1 <- debinfer_par(name = "k1", var.type = "de", fixed = FALSE,</pre>
                value = 0.05, prior = "lnorm", hypers = list(meanlog = 0, sdlog = 1),
                 prop.var = c(998,1000), samp.type = "rw-unif")
k2 <- debinfer_par(name = "k2", var.type = "de", fixed = FALSE,</pre>
                value = 8000, prior = "norm", hypers = list(mean = 5000, sd = 1000),
                prop.var = 500, samp.type="rw")
k3 <- debinfer_par(name = "k3", var.type = "de", fixed = FALSE,
                value = 3e7, prior = "norm", hypers = list(mean = 1e7, sd = 5e6),
                prop.var = 5e7, samp.type="rw")
sd.y1 <- debinfer_par(name = "sd.y1", var.type = "obs", fixed = FALSE,</pre>
                value = 0.05, prior = "lnorm", hypers = list(meanlog = 0, sdlog = 1),
                prop.var = c(3,4), samp.type = "rw-unif")
sd.y2 <- debinfer_par(name = "sd.y2", var.type = "obs", fixed = FALSE,
                value = 0.05, prior = "lnorm", hypers = list(meanlog = 0, sdlog = 1),
                prop.var = c(3,4), samp.type = "rw-unif")
sd.y3 <- debinfer_par(name = "sd.y3", var.type = "obs", fixed = FALSE,</pre>
                value = 0.05, prior = "lnorm", hypers = list(meanlog = 0, sdlog = 1),
                prop.var = c(3,4), samp.type = "rw-unif")
#Lastly, we provide an initial value N_0=0.1 for the DE:
y1 <- debinfer_par(name = "y1", var.type = "init", fixed = TRUE, value = 1)
y2 <- debinfer_par(name = "y2", var.type = "init", fixed = TRUE, value = 0)
y3 <- debinfer_par(name = "y3", var.type = "init", fixed = TRUE, value = 0)
#The declared parameters are then collated using the `setup_debinfer` function. Note that for models wi
mcmc.pars <- setup_debinfer(k1, k2, k3, sd.y1, sd.y2, sd.y3, y1, y2, y3)
```

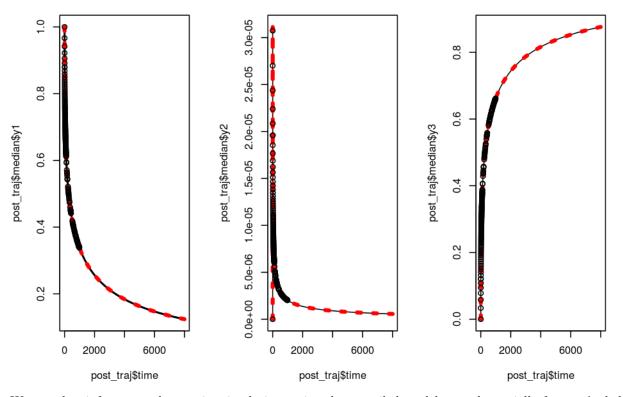
5 Conduct inference

Finally we use deBInfer to estimate the parameters of the original model

```
# do inference with deBInfer
# MCMC iterations
iter <- 1000
# inference call</pre>
```

```
Rt <- system.time(mcmc_samples <- de_mcmc(N = iter, data = obs, de.model = modelR,
                          obs.model = obs_model, all.params = mcmc.pars,
                          Tmax = max(obs$time), data.times = obs$time, cnt = 500,
                          plot = FALSE, solver = "ode", verbose.mcmc = FALSE))
## Order of initial conditions is y1, y2, y3
## initial posterior probability = 568.058099497889
Ct <- system.time(mcmc_samplesC <- de_mcmc(N = iter, data = obs, de.model = "derivs",
           jacfunc = "jac", dllname = "mymod",
           initfunc = "initmod", nout = 1, outnames = "Sum",
                          obs.model = obs_model, all.params = mcmc.pars,
                          Tmax = max(obs$time), data.times = obs$time, cnt = 1250,
                          plot = FALSE, solver = "ode", verbose.mcmc = FALSE))
## Order of initial conditions is y1, y2, y3
## initial posterior probability = 568.058099497889
print(Rt)
      user system elapsed
           0.028 94.157
## 94.050
print(Ct)
##
      user system elapsed
## 13.429
           0.004 13.444
#coda::rejectionRate(mcmc_samplesC$samples)
#plot(mcmc_samplesC)
, and simulate posterior trajectories.
system.time(post_traj <- post_sim(mcmc_samples, n=100, times=0:8000, burnin=100, output = 'all', prob =
##
      user system elapsed
## 80.080
           0.104 80.226
system.time(post_trajC <- post_sim(mcmc_samplesC, n=100, times=0:8000, burnin=100, output = 'all', prob</pre>
      user system elapsed
## 11.697 0.108 11.812
```

```
par(mfrow=c(1,3))
plot(post_traj$time, post_traj$median$y1, type='l',lwd=2)
lines(post_trajC$time, post_trajC$median$y1, col="red", lty=3, lwd=4)
points(obs$time, obs[,2])
plot(post_traj$time, post_traj$median$y2, type='l')
lines(post_trajC$time, post_trajC$median$y2, col="red", lty=3, lwd=4)
points(obs$time, obs[,3])
plot(post_traj$time, post_traj$median$y3, type='l')
lines(post_trajC$time, post_trajC$median$y3, col="red", lty=3, lwd=4)
points(obs$time, obs[,4])
```



We can that inference and posterior simulations using the compiled model are substantially faster. And this speed up is consistent across a range of MCMC iterations:

```
}
#saveRDS(timings, "examples/Soetaert-r-vs-c-timings.RDS")
iter_reps <- rep(c(100,1000,2000,5000,10000),3)
if(!exists("timings")) timings <- readRDS("Soetaert-r-vs-c-timings.RDS")</pre>
#pdf("examples/c-vs-r-iters.pdf")
par(mfrow=c(1,2))
plot(iter_reps, timings[,1], ylim=c(0, max(timings[,1])), type="p", ylab = "Execution time (s)", xlab =
points(iter_reps, timings[,2], col="red", type="p", pch=c(1,16))
legend("topleft", legend = c("R ODE", "C ODE"), pch = c(1,16), col= c("black", "red"))
plot(iter_reps, timings[,1]/timings[,2], ylim=c(0, max(timings[,1]/timings[,2])), type="p", ylab = "Rel
     9
                                       0
                                                             0 0
             R ODE
                                                                         0
                                                                                       0
     500
                                               Relative execution time R/C
               C ODE
                                                     9
Execution time (s)
     400
                                                     2
                                                      4
     300
                         0
                                                     က
     200
                                                     S
     100
                 0
     0
                                                     0
           0
               2000
                          6000
                                    10000
                                                           0
                                                              2000
                                                                         6000
                                                                                    10000
                 MCMC iterations
                                                                MCMC iterations
```

#dev.off()

Boersch-Supan, P. & Johnson, L. (2016). deBinfer: Bayesian inference for dynamical models of biological systems. arXiv, 1605.00021.

Soetaert, K., Petzoldt, T. & Setzer, R.W. (2009). R package desolve, writing code in compiled languages.

Soetaert, K., Petzoldt, T. & Setzer, R.W. (2010). Solving differential equations in R: Package deSolve. *Journal of Statistical Software*, **33**, 1–25.