# Example codes for "Backward importance sampling for online estimation of state space models"

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## 1 About the codes

- All the codes presented below are R codes that should work for a R version > 3.4.
- The proposed algorithm is mainly embedded in two R packages, one for the first section (that also contains code for the state of the art *Grand PaRIS* algorithm) and one for the second. To read the raw code of the proposed algorithm, the reader should navigate in the **src** directory of the packages after having downloaded them.

## 2 Comparison for the Sine Model

## 2.1 Installation of the required package

The required codes are packages into the GrandParisPackage package for the R software, available on github. To install this package in R, simply run:

## 2.2 Code for the comparison

Then, the full comparison is done in the following code:

```
# Cleaning ------
rm(list = ls())
# Packages -----
library(GrandParisPackage) # For the main algorithm
library(parallel) # For parallel computing (do not work on windows)
library(tidyverse) # For data processing
library(tictoc) # For time comparisons
library(gridExtra) # For multiple plots
# Simulating data -----
my_seed <- 333 # For all experiments, the random seed
set.seed(my_seed) # For reproducibility
trueTheta <- pi / 4; trueSigma2 <- 1;</pre>
n \leftarrow 11; times \leftarrow seq(from = 0, to = 5, length = n);
SINEprocess <- SINE_simulate(theta = trueTheta, sigma2 = trueSigma2,
                         x0 = 10, times = times)
observations <- SINEprocess[, "observations"]</pre>
# Experiment_function -----
# A function that computes the approximated estimate of E[X_0 \mid Y_0:n]
# for different tilde(N)
# It does it 400 times
compare_on_ntilde <- function(n_particle = 100, n_tilde = NULL){</pre>
 if(is.null(n_tilde)){
   stop("You must provide ntilde!")
 n_rep <- 400
 AR <- do.call(rbind,
              mclapply(1:n_rep, function(i){
               tic()
                res <- GrandParisPackage:::E_track(observations = observations,</pre>
                                               ind_tracked = 0,
                                               observationTimes = times,
                                               thetaStart = trueTheta,
                                               particleSize = n particle,
                                               backwardSampleSize = n_tilde,
                                               sigma2Start = trueSigma2,
```

```
nIterations = 1)
                 my_tictoc <- toc()</pre>
                 dur <- my_tictoc$toc - my_tictoc$tic</pre>
                 names(dur) <- NULL
                 data.frame(N = n_particle,
                           N_tilde = n_tilde, Xhat = res, Time = dur,
                           Method = factor("AR",
                                           levels = c("AR", "IS")))
               },
               mc.cores = detectCores() - 1)) # Number of cores for parallelization
  IS <- do.call(rbind,</pre>
               mclapply(1:n_rep, function(i){
                 tic()
                 res <- GrandParisPackage:::E_track_IS(observations = observations,</pre>
                                                   ind_tracked = 0,
                                                   observationTimes = times,
                                                   thetaStart = trueTheta,
                                                   particleSize = n_particle,
                                                   backwardSampleSize = n_tilde,
                                                   sigma2Start = trueSigma2,
                                                   nIterations = 1)
                 my_tictoc <- toc()</pre>
                 dur <- my_tictoc$toc - my_tictoc$tic</pre>
                 names(dur) <- NULL</pre>
                 data.frame(N = n_particle,
                           N_tilde = n_tilde, Xhat = res, Time = dur,
                           Method = factor("IS",
                                           levels = c("AR", "IS")))
               },
               mc.cores = detectCores() - 1)) # Number of cores for parallelization
 return(rbind.data.frame(AR, IS))
# Obtaining estimates -----
set.seed(my_seed) # For reproducibility
my_ntildes \leftarrow c(2, 5, 10, 20, 30)
# Run only once (experiment) -----
n_tilde_comparison_results <- lapply(my_ntildes, function(n_t)</pre>
  compare_on_ntilde(n_particle = 100, n_tilde = n_t)) %>%
 bind_rows()
# Plotting results ------
main_plot <- ggplot(n_tilde_comparison_results,</pre>
                   aes(x = factor(N_tilde), fill = Method)) +
  labs(x = expression(tilde(N))) +
  scale_fill_viridis_d(option = "G") +
  theme(legend.position = c(.9, .125))
```

```
p1 <- main_plot +
  geom_boxplot(aes(y = Time)) +
  scale_y_continuous(trans = "log10") +
  labs(y = "Comput. time (seconds)", fill = "")
p2 <- main_plot +
  geom_boxplot(aes(y = Xhat)) +
  labs(y = expression(hat("\U1d53c")~"["~X[0]~"|"~Y[0:n]~"]"),
       fill = "")
my_plot <- gridExtra::grid.arrange(p1, p2, nrow = 1)</pre>
# 12 * 4 inches
# Controlling N --
compare_on_n <- function(n_particle = 100, # Number of particles</pre>
                          alphas = NULL, # Power of N to obtain Ntilde
                          \# 0.5 and 0.6 in the article
                          frac = NULL){ # Fraction of N to obtain Ntilde (0.1 in the article)
  n_tilde_AR <- 2
  levels_IS <- paste0("IS_a_", alphas, "_p_", frac)</pre>
  n rep <- 50
  AR <- do.call(rbind,
                mclapply(1:n_rep, function(i){
                   res <- GrandParisPackage:::E_track(observations = observations,</pre>
                                                        ind tracked = 0,
                                                        observationTimes = times,
                                                        thetaStart = trueTheta,
                                                        particleSize = n_particle,
                                                        backwardSampleSize = n_tilde_AR,
                                                        sigma2Start = trueSigma2,
                                                       nIterations = 1)
                   my_tictoc <- toc()</pre>
                   dur <- my_tictoc$toc - my_tictoc$tic</pre>
                   data.frame(N = n_particle,
                              N_tilde = n_tilde_AR, Xhat = res, Time = dur,
                              Method = factor("AR",
                                               levels = c("AR", levels_IS)))
                 }, mc.cores = detectCores() - 2))
  IS <- do.call(rbind.data.frame,</pre>
                 mapply(function(my_alpha, my_prop){
                   n_tilde_IS = ceiling(my_prop * n_particle^my_alpha)
                   do.call(rbind.data.frame,
                           mclapply(1:n_rep, function(i){
                     res <- GrandParisPackage:::E_track_IS(observations = observations,</pre>
                                                             ind_tracked = 0,
                                                             observationTimes = times,
                                                             thetaStart = trueTheta,
                                                             particleSize = n_particle,
                                                             backwardSampleSize = n_tilde_IS,
                                                             sigma2Start = trueSigma2,
                                                             nIterations = 1)
                     my_tictoc <- toc()</pre>
```

```
dur <- my_tictoc$toc - my_tictoc$tic</pre>
                    names(dur) = NULL
                    data.frame(N = n_particle,
                                N_tilde = n_tilde_IS, Xhat = res, Time = dur,
                                Method = factor(pasteO("IS_a_", my_alpha, "_p_", my_prop),
                                                levels = c("AR", levels_IS)))
                  }, mc.cores = detectCores() - 2))
                }, alphas, frac, SIMPLIFY = F))
 return(rbind.data.frame(AR, IS))
}
my_n_particles <- c(50, 100, 200, 500, 1000, 2000)
set.seed(my_seed)
res_npart <- lapply(my_n_particles,</pre>
                 function(my_n) compare_on_n(n_particle = my_n,
                                              alphas = c(0.5, 0.6, 1),
                                              frac = c(1, 1, 0.1))
res_df_npart <- do.call(rbind.data.frame, res_npart)</pre>
# save(res_df_npart, file = "res_df_npart.RData")
# Plotting results
my_labels <- expression("AR,"~tilde(N)==2, "IS,"~tilde(N)== N^0.5,
                        "IS, "~tilde(N) == N^0.6, "IS, "~tilde(N) == N/10)
main_plot <- res_df_npart %>%
  mutate(Method = factor(Method, labels = my_labels)) %>%
  ggplot(aes(x = factor(N), fill = Method)) +
  labs(x = expression(N), fill = "") +
  scale_fill_viridis_d(labels = my_labels, option = "G") +
  theme(legend.text.align = 0,
        legend.position = c(.75, .2),
        legend.background = element_blank())
p1 <- main_plot +
  geom_boxplot(aes(y = Time)) +
  scale_y_continuous(trans = "log10") +
  labs(y = "Comput. time (seconds)")
p2 <- main_plot +
  geom_boxplot(aes(y = Xhat)) +
  labs(y = expression(hat("\U1d53c")~"["~X[0]~"|"~Y[0:n]~"]"))
my_plot <- gridExtra::grid.arrange(p1, p2, nrow = 1)</pre>
ggplot(res_df_npart, aes(x = N, y = Time)) +
  geom_point() + geom_smooth() +
  facet_wrap(~Method, scales = "free_y") +
  labs(x = expression(tilde(N)),
      y = "Comput. time (seconds)")
```

## 3 Online estimation in the Sine model

In this section, the online estimation is performed

```
rm(list = ls())
library(GrandParisPackage)
library(tidyverse)
library(parallel)
library(gridExtra)
```

## 3.1 Simulating data

#### 3.1.1 Simulation parameters

#### 3.1.2 Simulation

```
# The following code creates 500 trajectories in a directory "simulated_data"
# that can be created with the following code
# dir.create("simulated_data")
n_traj <- 500
# For windows user, replace by lapply and remove the mc.cores argument below
if(!dir.exists("simulated_data/")){
  dir.create("simulated_data") # Create a folder to stock simulated data
}
# Simulated 500 trajectories and write it in the folder "simulated_data"
# Simulation is done exactly using exact algorithms
mclapply(1:n_traj,
         function(i){
           seed <- 100 + i
           set.seed(seed)
           simulated_POD <- SINE_simulate(theta = trueTheta,</pre>
                                           sigma = trueSigma2, x0 = 0,
                                           times = times)
           write.table(simulated_POD, paste0("simulated_data/simul_data_seed", seed, ".txt"),
                       col.names = T, row.names = F, sep = ";")
         },
         mc.cores= detectCores() - 1)
```

## 3.2 Estimation

Code for estimation is only shown for one trajectory. This gives the result of the first part of section 5.3 (the Figure 4)

```
# Get the appropriate trajectory
seed <- 122
set.seed(seed) # For reproducibility
simulated POD <- read.table(paste0("simulated data/simul data seed", seed, ".txt"),
                             sep = ";", header = TRUE)
observations <- simulated_POD[, "observations"]</pre>
# Estimation parameters -----
## One set of observations, several starting points for the algorithm
get_estimation_one_obs_several_start <- function(){</pre>
  gradientSteps <- get_grad_steps(0.6, cst = 8, n = n)</pre>
  n_start_points <- 50</pre>
  n_particles <- 100</pre>
  N_tilde <- n_particles / 10
  allRes <- mclapply(1:n_start_points, function(seed){</pre>
    set.seed(seed)
    thetaStart <- runif(1, 0, 2 * pi)</pre>
    fastTangOR(observations, times, particleSize = n_particles,
               thetaModel = thetaStart, sigma2 = trueSigma2,
               updateOrders = rep(TRUE, n),
               gradientSteps = gradientSteps,
               all = FALSE, estimateTheta = TRUE, estimateSigma2 = FALSE,
               randomWalkParam = 1, backwardSampleSize = N_tilde, IS = TRUE)
  },
  mc.cores = detectCores() - 1)
  thetaEst <- sapply(allRes, function(x) x$Estimates[,1]) %% (2*pi)</pre>
  return(thetaEst)
res_one_obs_several_start <- get_estimation_one_obs_several_start()</pre>
res_several_obs_one_start <- mclapply(1:length(dir("simulated_data/")), function(i){</pre>
  seed <- 100 + i
  set.seed(seed)
  simulated_POD <- read.table(paste0("simulated_data/simul_data_seed", seed, ".txt"),</pre>
                               sep = ";", header = T)
  observations <- simulated_POD[, "observations"]</pre>
  thetaStart <- trueTheta
  Res <- fastTangOR(observations, times,</pre>
                     thetaModel = thetaStart, sigma2 = trueSigma2, particleSize = 500,
                     updateOrders = rep(TRUE, length(observations)),
                     gradientSteps = get_grad_steps(0.6, cst = 8,
                                                     n = length(observations)),
                     all = TRUE, estimateSigma2 = FALSE, randomWalkParam = 1)
  Res
```

```
},
mc.cores = detectCores() - 1)
```

## 4 Experiments for the Lotka Volterra Model

## 4.1 Installation of the required package

The required codes are packages into the LotkaVoltR package for the R software, available on github. To install this package in R, simply run:

```
devtools::install_github("papayoun/LotkaVoltR")
```

## 4.2 Smoothing on synthetic data

## 4.2.1 Simulating data

```
# Cleaning
rm(list = ls()) # Cleaning environment
# Librairies
library(LotkaVoltR) # Dedicated library
library(tidyverse) # For data processing
# Dynamics parameters
a1 <- c(12, 0.05, 1) # Prey parameters
a2 <- c(2, 0.2, 0.1) # Predator parameters
Gamma \leftarrow matrix(c(0.5, 0.1, 0.1, 0.2), nrow = 2) \# Diffusion parameters
mu0 <- c(50,20) # Mean of the inital distribution
Sigma0 <- diag(1, 2) # Variance of the initial distribution
# Observation process parameters
Sigma_obs <- matrix(c(0.01, 0.005, 0.005, 0.01), ncol = 2) # Observation noise
q_{values} \leftarrow c(0.2, 0.3) \# Known \ q \ values
# Model creation
# Creation of a Partially Observed Lotka Volterra model
POLV_model <- POLV_create(a1 = a1, a2 = a2, gam = Gamma, mu0 = mu0,
                           sigma0 = Sigma0, cov = Sigma_obs, qs = q_values)
# Synthetic data simulation
# Simulation times for the Euler scheme
simulation_times <- seq(from = 0, by = 1e-6, # Simulation time step, small!!
```

#### 4.2.2 Performing smoothing

## 4.2.3 Plotting results

```
Y2 = Y2 / q_values[2]) %>%
  rename(X1 = Y1, X2 = Y2) %>%
  mutate(method = my_levels["obs"])
true_values <- simulated_process %>%
  as.data.frame() %>%
  select(X1, X2) %>%
 mutate(method = my_levels["true"])
concatened_results <- bind_rows(observed_values,</pre>
                                true_values,
                                smoothing_mean) %>%
 mutate(method = factor(method, levels = my_levels)) # Pour l'ordre
ggplot(concatened_results) +
  aes(x = X1, y = X2, color = method, linetype = method) +
  geom_point() +
  geom_path() +
  theme(legend.position = "none") +
  facet_wrap(~method, labeller = label_parsed) +
  labs(x = "Number of preys", y = "Number of predators") +
  theme(strip.text = element_text(size = 24))
```

## 4.3 Experiments on Lynx data

#### 4.3.1 Data set

## 4.3.2 Performing EM for estimation

## Initial parameter

```
gam = Gamma)
# Model creation
```

EM functions: A generalized EM is performed. At each step, candidates are generated using a home made (not optimized!) evolution strategy, by sampling around the current parameter through a Gaussian distribution. Each parameter lives in a constrained space. Each new offspring is generated in the unconstrained  $\mathbb{R}$  space before being put back to the constrained space. This is mainly done with the logistic function.

Then, a usual EM approach is performed. To ensure good fit, the EM is performed from three different starting points.

```
library(LotkaVoltR) # Dedicated library
# Function generation candidat -----
generate_candidate_list <- function(par_list, # Current parameter</pre>
                                    standard dev list, # Standard Deviation for
                                     # the future offspring generation
                                    up_lim_list){ # Limits of the constrained space
  out <- par_list
  # Generating a1, in [0, 1] * [0, 0.05] * [0, 0.05]
  sigmoid <- function(x){</pre>
   1 / (1 + \exp(-x))
  logit <- function(x){</pre>
   log(x / (1 - x))
  out$a1 <- (par_list$a1 / up_lim_list$a1) %>% # Back to [0, 1]
   logit() %>% # Back to the real world
   rnorm(n = 3, sd = standard_dev_list$a1) %>% # Moving a bit
   sigmoid() %>% # Retour dans [0,1]
   {. * up_lim_list$a1} # Back to the constrained space
  out$a2 <- (par_list$a2 / up_lim_list$a2) %>% # Back to [0, 1]
    logit() %>% # Back to the real world
   rnorm(n = 3, sd = standard_dev_list$a2) %>% # Moving a bit
    sigmoid() %>% # Retour dans [0,1]
    {. * up_lim_list$a2} # Back to [0, 1] * [0, 0.1] * [0, 0.1]
  diag(out$gam) <- (diag(par_list$gam) / up_lim_list$gam) %% # Back to [0, 1]
   logit() %>% # Back to the real world
   rnorm(n = 2, sd = standard_dev_list$gam) %>% # Moving a bit
    sigmoid() %>% # Back to [0,1]
    {. * up_lim_list$gam}
  diag(out$cov) <- (diag(par_list$cov) / up_lim_list$cov) %>% # Back to [0, 1]
   logit() %>% # Back to the real world
   rnorm(n = 2, sd = standard_dev_list$cov) %>% # Moving a bit
    sigmoid() %>% # Back to [0,1]
    {. * up_lim_list$cov}
  return(out)
}
EM_function <- function(obs, obs_times, initial_param, initial_sd, up_lims,
                        n cands, n iter,
```

```
seed,
                         name = NULL){
  param_0 <- initial_param</pre>
  out <- list(param_0)</pre>
  set.seed(seed)
  final_E_steps <- matrix(NA, nrow = n_iter, ncol = n_cands + 1)</pre>
  for(i in 1:n_iter){
    print(paste("Iteration", i))
    par_list <- c(purrr::rerun(n_cands,</pre>
                                 generate_candidate_list(param_0,
                                                           initial_sd,
                                                           up_lims)),
                   list(param_0))
    E_step_evals <- get_E_step(obs_ = t(obs),</pre>
                                 obsTimes_ = obs_times,
                                 myParams = param_0,
                                 testedParams = par_list,
                                 n_part = 200, n_dens_samp = 100)
    initial_sd <- purrr::map(initial_sd,</pre>
                               function(x) 0.9 * x) # Reducing the exploration
    param_0 <- par_list[[which.max(E_step_evals)]]</pre>
    out <- c(out, list(param_0))</pre>
    final_E_steps[i, ] <- E_step_evals</pre>
  return(list(out, final_E_steps))
}
observations <- as.matrix(hares_lynx_data[, c("Hares", "Lynx")])
observation_times <- hares_lynx_data$Year</pre>
initial_param <- list(a1 = a1, a2 = a2, mu0 = mu0,
                       sigma0 = Sigma0, RWC = diag(0.005, 2),
                       qs = q_values, cov = Sigma_obs,
                       wD = 1, w0 = 1,
                       gam = Gamma)
sd_list \leftarrow list(a1 = c(1, 1, 1),
                 a2 = c(1, 1, 1),
                 cov = 0.5,
                 gam = 0.5)
upper_lims_list <- list(a1 = c(1, 0.01, 0.05),
                         a2 = c(1, 0.05, 0.01),
                         cov = 0.5,
                         gam = 0.5)
library(parallel) # For parallel computing
results <- mclapply(1:30, function(my_seed){
  EM_function(obs = observations,
              obs_times = observation_times,
              initial_param = initial_param,
              initial_sd = sd_list,
              up_lims = upper_lims_list,
              n_{cands} = 10,
              n_iter = 30, seed = my_seed,
              name = "lynx")},
```

```
mc.cores = detectCores() - 1)
```

The results are then used to perform the smoothing as in previous section