es3

2024-01-13

Load libraries

```
shhh = function(lib_name){ # It's a library, so shhh!
  suppressWarnings(suppressMessages(require(lib_name, character.only = TRUE)))
shhh("tidyverse")
shhh("ACutils")
shhh("mvtnorm")
shhh("salso")
shhh("FGM")
shhh("gmp")
shhh("mcclust")
shhh("mcclust.ext")
shhh("logr")
shhh("tidygraph")
shhh("ggraph")
shhh("igraph")
shhh("Rcpp")
shhh("RcppArmadillo")
shhh("RcppEigen")
## Load custom functions
source("functions/utility_functions.R");
source("functions/bulky_functions.R");
source("functions/data_generation.R")
sourceCpp("functions/wade.cpp")
Rcpp::sourceCpp('functions/UpdateParamsGSL.cpp')
library('RcppGSL')
library(fda)
library(tidyverse)
library(coda)
library(lattice)
```

Goal

Stima della partizione e grafo con bulky functions 2024.

Simulated data

Simuliamo i dati in sim_data.R con p=40 basi e 3 cluster di dimensioni [13,13,14]. Grafo e matrice di precisione sono generati attraverso la funzione Generate BlockDiagonal.

```
# Import simulated data
BaseMat = as.matrix(read_csv('simulated_data/BaseMat.csv'))  # matrix Phi, dim = 200 x 40
y_hat_true = as.matrix(read_csv('simulated_data/y_hat_true.csv'))  # y true, dim = 300 x 200
beta_true = as.matrix(read_csv('simulated_data/beta_true.csv'))  # beta, dim = 300 x 40
mu_true = as.matrix(read_csv('simulated_data/mu_true.csv'))  # mu true, dim = 40 x 1

# object containing the true graph G, the true precision matrix K
simKG <- readRDS("simulated_data/simKG.rds")

n <- dim(y_hat_true)[1]  # n=300
r <- dim(y_hat_true)[2]  # r=200
p <- dim(BaseMat)[2]  # p=40</pre>
```

Initialization

```
# Compute quantities for function UpdateParamGSL
tbase base = t(BaseMat) % * % BaseMat
                                              # p x p (phi_t * phi)
tbase_data = t(BaseMat)%*%t(y_hat_true)
                                              \# p \times n \quad (phi_t + Y_t)
Sdata = sum(diag(y_hat_true%*%t(y_hat_true))) # initialize phi*beta = 0
set_UpdateParamsGSL_list = set_UpdateParamsGSL (
    tbase_base = tbase_base,
   tbase_data = tbase_data,
   Sdata
               = Sdata,
   a_tau_eps = 2000,
   b_{tau_eps} = 2,
   sigma_mu
                = 100,
                = r,
   Update Beta = FALSE,
   Update Mu = FALSE,
   Update_Tau = FALSE
# Set the number of iterations and burn-in
niter <- 50000
burn_in <- 1000
# Set the initialization values of the chains
initialization_values = set_initialization(
   Beta
                 = t(beta_true),
                  = mu_true,
   mu
                  = 100,
   tau_eps
                  = matrix(0,p,p),
   G
                  = matrix(0,p,p),
                  = rep(1,p),
   Z
   rho
                  = p,
```

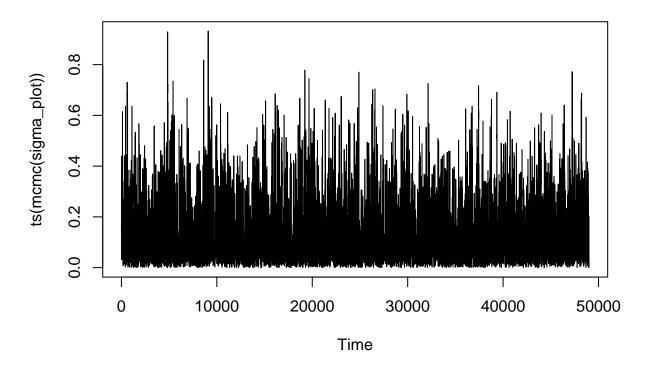
```
a_sigma
                 = 1,
   b_sigma
                 = 1,
   c_sigma
                 = 1,
   d_sigma
                 = 1
   c_theta
               = 100,
               = 10,
   d_theta
   sigma
                = 0.5,
   theta
               = 1,
   weights_a0 = rep(1,p-1),
weights_d0 = rep(1,p-1),
   total_weights = 0,
   total_K = matrix(0,p,p),
   total_graphs = matrix(0,p,p),
   graph_start = NULL,
   graph_density = sum(simKG$Graph) / (p*(p-1)),
   beta_sig2 = 0.2,
)
```

Gibbs sampler

Useful plots

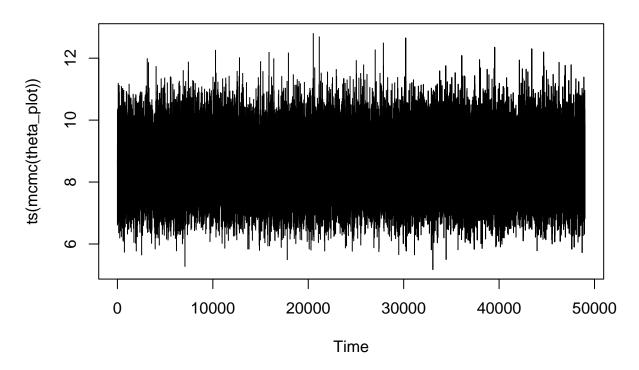
```
sigma_plot <- as.vector(chains$sigma)
sigma_plot <- sigma_plot[(burn_in+1):niter]
plot(ts(mcmc(sigma_plot)), main='Trace plot sigma')</pre>
```

Trace plot sigma



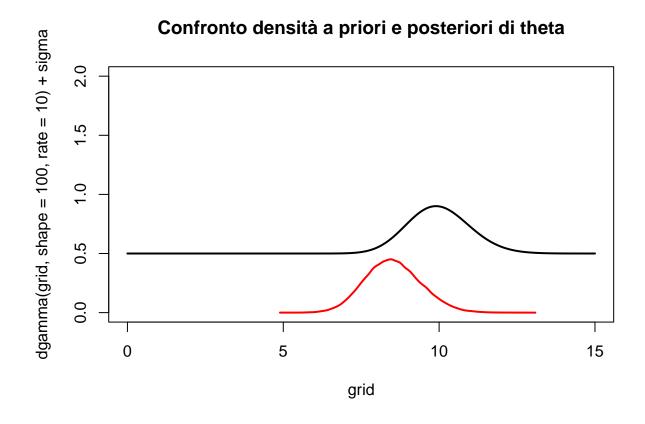
```
theta_plot <- as.vector(chains$theta)
theta_plot <- theta_plot[(burn_in+1):niter]
plot(ts(mcmc(theta_plot)), main='Trace plot theta')</pre>
```

Trace plot theta



Confronto densità di theta

```
# confronto densità a priori
min_val <- 0
max_val <- 15
sigma <- 0.5
grid <- seq(min_val, max_val, length.out=10000)
plot(grid, dgamma(grid,shape=100, rate=10)+sigma, type='l', lwd=2, ylim=c(0,2), main='Confronto densità
lines(density(mcmc(theta_plot))$x, density(mcmc(theta_plot))$y, type='l', lwd=2, col='red') #posterior</pre>
```



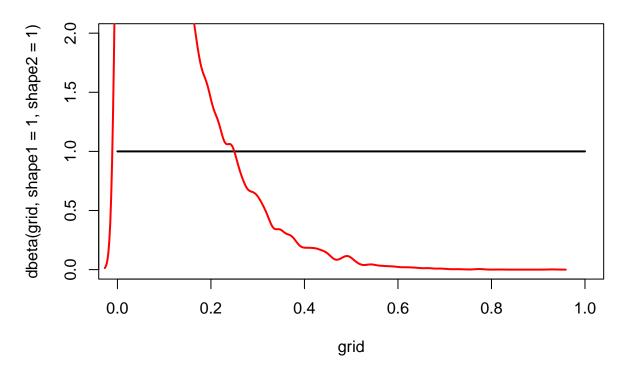
```
# numero atteso di cluster, formula Martinez and Mena
theta_mean <- 1
n_mean_cluster <- lpochhammer(theta_mean + sigma, p, log=FALSE)/(sigma * lpochhammer(theta_mean + 1, p-1)
n_mean_cluster</pre>
```

[1] 12.40632

Confronto densità di theta

```
min_val <- 0
max_val <- 1
sigma <- 0.5
grid <- seq(min_val, max_val, length.out=10000)
plot(grid, dbeta(grid, shape1 = 1, shape2 = 1), type='l', lwd=2, ylim=c(0,2), main='Confronto densità a
lines(density(mcmc(sigma_plot))$x, density(mcmc(sigma_plot))$y, type='l', lwd=2, col='red') #posterior</pre>
```

Confronto densità a priori e posteriori di sigma



Posterior analysis

```
## Recomputing the partition in other forms and the number of groups

rho_true = c(13,13,14)
r_true = rho_to_r(rho_true)
z_true = rho_to_z(rho_true)
p = length(z_true)
num_clusters_true = length(rho_true)
rho <- chains*rho
r = do.call(rbind, lapply(chains*rho, rho_to_r))

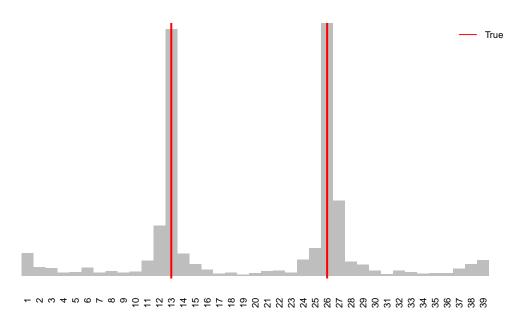
## Warning in (function (..., deparse.level = 1) : number of columns of result is
## not a multiple of vector length (arg 2)

z = do.call(rbind, lapply(chains*rho, rho_to_z))
num_clusters = do.call(rbind, lapply(chains*rho, length))
num_clusters = as.vector(num_clusters)

### Barplot of changepoints</pre>
```

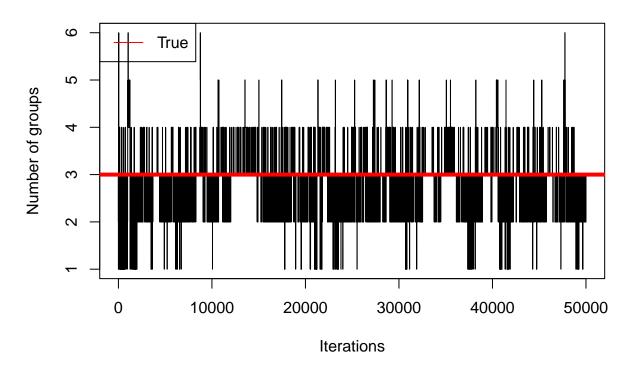
```
bar_heights = colSums(r)
cp_true = which(r_true==1)
color <- ifelse(seq_along(bar_heights) %in% c(cp_true), "red", "gray")</pre>
barplot(
  bar_heights[1:(p-1)],
  names = seq_along(bar_heights[1:(p-1)]),
  border = "NA",
  space = 0,
  yaxt = "n",
  main="Changepoint frequency distribution",
  \#col = color,
  cex.names=.6,
 las=2
abline(v=cp_true-0.5, col="red", lwd=2)
legend("topright", legend=c("True"), col=c("red"),
       bty = "n",
       lty = 1,
       cex = 0.6)
```

Changepoint frequency distribution



```
### Evolution of the number of clusters
plot(
    x = seq_along(num_clusters),
```

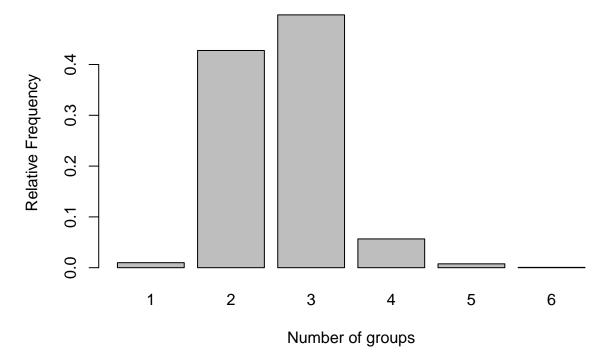
Number of groups - Traceplot



```
barplot(
  prop.table(table(num_clusters)),
  xlab = "Number of groups",
  ylab = "Relative Frequency",
  main = paste(
    "Number of groups - Relative Frequency\n",
    "Last:",
  tail(num_clusters, n = 1),
    "- Mean:",
  round(mean(num_clusters), 2),
    "- True:",
```

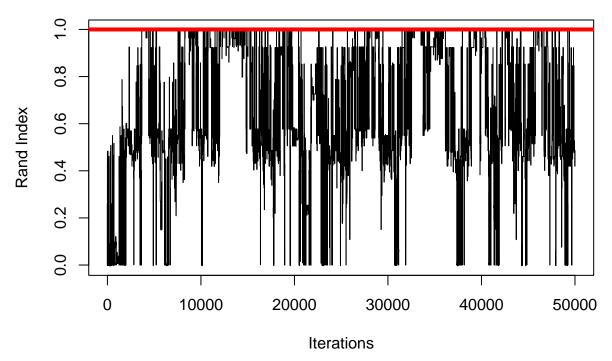
```
num_clusters_true
)
)
```

Number of groups – Relative Frequency Last: 2 – Mean: 2.63 – True: 3



```
### Evolution of the Rand Index
# computing rand index for each iteration
rand_index = apply(z, 1, mcclust::arandi, z_true)
# plotting the traceplot of the index
plot(
  x = seq_along(rand_index),
  y = rand_index,
  type = "n",
  xlab = "Iterations",
  ylab = "Rand Index",
  main = paste(
    "Rand Index - Traceplot\n",
    "Last:",
    round(tail(rand_index, n=1), 3),
    "- Mean:",
    round(mean(rand_index), 2)
  )
)
lines(x = seq_along(rand_index), y = rand_index)
```

Rand Index - Traceplot Last: 0.486 - Mean: 0.66



```
### Retrieving best partition using VI on visited ones (order is guaranteed here)
# compute VI
sim_matrix <- salso::psm(z)
dists <- VI_LB(z, psm_mat = sim_matrix)

# select best partition (among the visited ones)
best_partition_index = which.min(dists)
rho_est = rho[[best_partition_index]]
z_est = z[best_partition_index,]

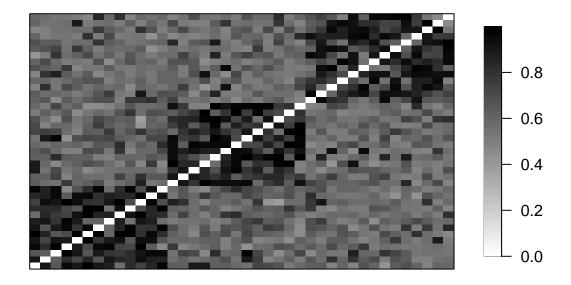
# VI loss
dists[best_partition_index]</pre>
```

[1] 0.6645961

```
# select best partition
unname(z_est)
```

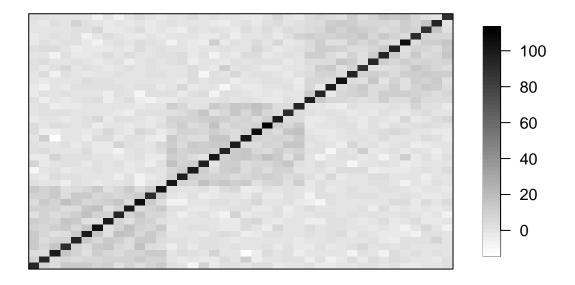
```
# compute Rand Index
mcclust::arandi(z_est, z_true)
## [1] 1
## Graph
# Extract last plinks
last_plinks = tail(chains$G, n=1)[[1]]
# Criterion 1 to select the threshold (should not work very well) and assign final graph
threshold = 0.5
G_est <- matrix(0,p,p)</pre>
G_est[which(last_plinks>threshold)] = 1
#Criterion 2 to select the threshold
bfdr_select = BFDR_selection(last_plinks, tol = seq(0.1, 1, by = 0.001))
# Inspect the threshold and assign final graph
bfdr_select$best_treshold
## [1] 0.904
G_est = bfdr_select$best_truncated_graph
### Standardized Hamming distance
SHD = sum(abs(simKG$Graph - G_est)) / (p^2 - p)
SHD
## [1] 0.225641
### Plot estimated matrices
ACutils::ACheatmap(
 last_plinks,
  use_x11_device = F,
 horizontal = F,
 main = "Estimated plinks matrix",
 center_value = NULL,
  col.upper = "black",
  col.center = "grey50",
  col.lower = "white"
```

Estimated plinks matrix



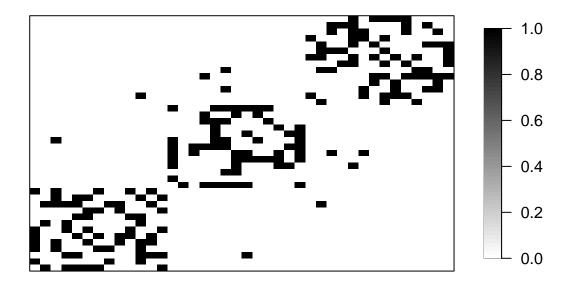
```
ACutils::ACheatmap(
  tail(chains$K,n=1)[[1]],
  use_x11_device = F,
  horizontal = F,
  main = "Estimated Precision matrix",
  center_value = NULL,
  col.upper = "black",
  col.center = "grey50",
  col.lower = "white"
)
```

Estimated Precision matrix



```
ACutils::ACheatmap(
   G_est,
   use_x11_device = F,
   horizontal = F,
   main = "Estimated Graph",
   center_value = NULL,
   col.upper = "black",
   col.center = "grey50",
   col.lower = "white"
)
```

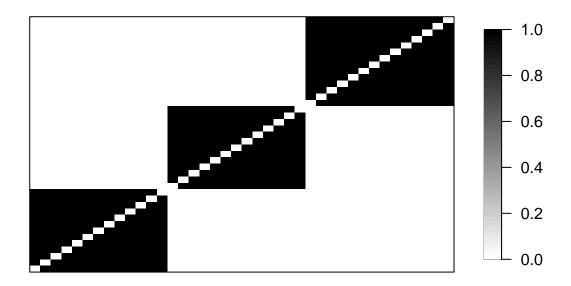
Estimated Graph



```
# Plot the true graph
# remove self loops
simKG$Graph[col(simKG$Graph)==row(simKG$Graph)] = 0

ACutils::ACheatmap(
    simKG$Graph,
    use_x11_device = F,
    horizontal = F,
    main = "Simulated Graph",
    center_value = NULL,
    col.upper = "black",
    col.center = "grey50",
    col.lower = "white"
)
```

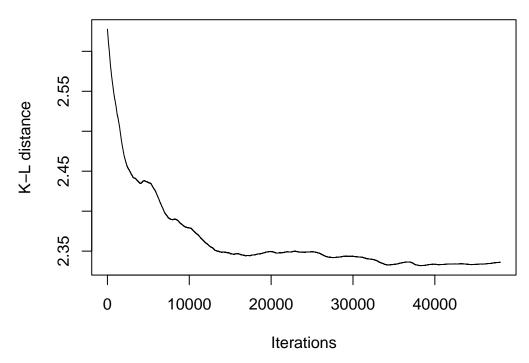
Simulated Graph



```
### Evolution of the Kullback-Leibler
kl_dist = do.call(rbind, lapply(chains$K, function(k) {
    ACutils::KL_dist(simKG$Prec, k)
}))

last = round(tail(kl_dist, n=1), 3)
plot(
    x = seq_along(kl_dist[2000:length(kl_dist)]),
    y = kl_dist[2000:length(kl_dist)],
    type = "n",
    xlab = "Iterations",
    ylab = "K-L distance",
    main = paste("Kullback-Leibler distance\nLast value:", last)
)
lines(x = seq_along(kl_dist[2000:length(kl_dist)]), y = kl_dist[2000:length(kl_dist)])
```

Kullback-Leibler distance Last value: 2.336



non c'è burn in !!!!