

Generator (dependent case - s0.6)

June 14, 2023

```
[1]: import warnings
warnings.filterwarnings('ignore')
```

```
[2]: ##### Importing packages
      <-#####

import numpy as np          # to handle arrays and matrices
import pickle

from scipy.linalg import toeplitz # to generate toeplitz matrix
from scipy.stats import chi2      # to have chi2 quantiles
from scipy.special import chdtri

import matplotlib.pyplot as plt  # to plot histograms ...
import pandas as pd             # to handle and create dataframes

from scipy.linalg import toeplitz # to generate toeplitz matrix
from scipy.stats import chi2      # to have chi2 quantiles
from scipy.special import chdtri

import time
import concurrent.futures
import random
import os

from itertools import product

##### For printing with colors #####
class color:
    PURPLE = '\033[95m'
    BLACK = '\033[1;90m'
    CYAN = '\033[96m'
    DARKCYAN = '\033[36m'
    BLUE = '\033[94m'
    GREEN = '\033[1;92m'
    YELLOW = '\033[93m'
```

```

RED = '\033[1;91m'
BOLD = '\033[1m'
UNDERLINE = '\033[4m'
END = '\033[0m'
BCKGRND = '\033[0;100m'
RBCKGRND = '\033[0;101m'

print(color.BLUE + color.BOLD + '***** Starting the program !\n
↳*****' + color.END )

```

***** Starting the program ! *****

0.1 2.1 Toeplitz Matrix

Creating a function that gets two parameters s and q such that : - s (dependency threshold or intraparameter of dispersion) - and q (the dimension)

then returns a symmetric toeplitz matrix S^2 that can be used as a covariance matrix for generating normal vectors

```

[3]: def cov_toep(s, q):
      """A function that takes the dependency thresholds $s$
      and the dimension $q$ and returns a $q \times q$-toeplitz matrix.
      """
      row = np.array([])
      for k in range(q):
          row = np.append(row, float(s**k))
      return toeplitz(row, row)

```

```

[4]: print(cov_toep(0.1, 4))
      print()
      print(cov_toep(0.99, 4))

```

```

[[1.    0.1  0.01 0.001]
 [0.1   1.   0.1  0.01 ]
 [0.01  0.1  1.   0.1  ]
 [0.001 0.01 0.1  1.   ]]

```

```

[[1.    0.99  0.9801 0.970299]
 [0.99   1.   0.99  0.9801  ]
 [0.9801 0.99  1.   0.99  ]
 [0.970299 0.9801 0.99  1.   ]]

```

```

[5]: # importing q_list, n_list, S2 diagonals, for different sigma (sigma means here
      ↳the std of underlying normal
      # distribution that generates lognormal vectors), having the same alpha
      q_list = list(pickle.load(open("q_list", "rb")))

```

```
n_list = list(pickle.load(open("n_list", "rb")))
print(q_list)
```

[50, 100, 150, 200, 250, 300, 350, 400]

```
[6]: def scalar(A,B):
      """Takes two symmetric matrices A and B of sizes q
      and returns the modified frobenius scalar of A and B
      """
      return(np.trace(A.dot(np.transpose(B)))/A.shape[0])

def norm(A):
      """Takes a symmetric matrix A of sizes q
      and returns the norm of A
      This norm is associated to the modified frobenius scalar
      """
      return np.sqrt(scalar(A,A))

def alphaaa1(S_2):
      q = len(S_2)
      I_q = np.diag(np.ones(q))
      sigma2 = scalar(S_2,I_q)
      alpha2 = norm(S_2 - sigma2*I_q)**2
      return alpha2

def alphaaa2(vec):
      q = len(vec)
      I_q = np.diag(np.ones(q))
      S_2 = np.diag(vec)
      sigma2 = scalar(S_2,I_q)
      alpha2 = norm(S_2 - sigma2*I_q)**2
      return alpha2
```

```
[7]: ##### Preparing true covariance matrices for different
      ↪ values of q #####

valeur_propre_collection = [cov_toep(0.6,q) for q in q_list]
print(*valeur_propre_collection[q_list.index(50)][0][:10])
print([alphaaa1(cov_toep(0.6,q)) for q in q_list])
```

```
1.0 0.6 0.36 0.21599999999999997 0.1296 0.07775999999999998 0.04665599999999999
0.027993599999999993 0.016796159999999994 0.010077695999999997
[1.0898437500000002, 1.107421875, 1.1132812499999998, 1.1162109374999998,
1.1179687499999995, 1.119140625, 1.1199776785714288, 1.1206054687499998]
```

```
[8]: # list(product(n_list, range(K)))[12]
```

```

[9]: ##### Choosing dimension
      ↪#####

# K = int(input("Number of Monte Carlo iterations is : "))
print()
print("list of q values : ", q_list, "\n")
print("list of n values : ", n_list, "\n")

##### Generator function #####

def generator(n, q):
    """
    Creating a function that gets into paramaters :
        the number of observations
        the dimension
        dependency threshold
        2 covariance parameter
    and returns a matrix of n observations (n rows), where each row represents
    a q-vector normally distributed with a mean 0 and covariance matrix  $S^2$ 
    ↪defined
    by a toeplitz matrix with a threshold s
    """

    # defining eigenvalues, lognormal variables
    valeur_propre = valeur_propre_collection[q_list.index(q)]

    # defining a mean vector and a covariance matrix
    mean = np.zeros(q) ; cov = valeur_propre

    # z_intermediate = q rows and n columns we still need to transpose
    z_int = np.random.multivariate_normal(mean, cov, n).T

    # z sample matrix, having n rows and q columns
    z = np.transpose(z_int)

    # Returning the matrix of n-observations of dimension q
    return z

##### Statistics functions #####

def sn2(z):
    return np.cov(np.transpose(z))

def zbar(z):
    """takes a n x q sample matrix and return the vector mean of each column
  
```

```

    """
    return z.mean(axis=0)

def max_p(M):
    """Largest eigenvalue of a given matrix M"""
    val_p = np.linalg.eigvals(M)
    return max(val_p.real)

def min_p(M):
    """Smallest eigenvalue of a given matrix M that is not null"""
    val_p = np.linalg.eigvals(M)
    val_p = val_p[val_p>=10**-6]
    return min(val_p.real)

def product_vect(z, i):
    return np.matmul(z[i].reshape(z[i].shape[0], 1), np.transpose(z[i].
    ↪reshape(z[i].shape[0], 1)))

##### Algebra functions
↪#####

def scalar(A,B):
    """Takes two symmetric matrices A and B of sizes q
    and returns the modified frobenius scalar of A and B
    """
    return(np.trace(A.dot(np.transpose(B)))/A.shape[0])

def norm(A):
    """Takes a symmetric matrix A of sizes q
    and returns the norm of A
    This norm is associated to the modified frobenius scalar
    """
    return np.sqrt(scalar(A,A))

```

list of q values : [50, 100, 150, 200, 250, 300, 350, 400]

list of n values : [50, 75, 100, 125, 150, 175, 200]

1 Time comparison

```
[10]: def monte_carlo(k):
    # random.seed(k)
    np.random.seed(int(os.getpid() * time.time()) % 123456789)
    z = generator(n,q)
    sn_2 = sn2(z)
    # beta = (norm(sn_2 - s_2))**2
    # calculate empirical mean
    z_bar = zbar(z)
    #empirical covariance matrix
    sn_2 = sn2(z)

    # Calculating empirical eigenvalues and eigenvectors (of  $Sn^2$ )
    emp_val_p, emp_vec_p = np.linalg.eigh(sn_2)

    # Calculating true (theoretical) eigenvalues and eigenvectors (of  $S^2$ )
    # val_p, vec_p = np.linalg.eigh(s_2)

    # Identity matrix of size q
    I_q = np.diag(np.ones(q))

    # sigma_n ( ^ 2 )
    sigma_n = scalar(sn_2, I_q)

    # delta_n ( ^ 2 )
    delta_n = norm(sn_2 - sigma_n*I_q)**2

    # intermediate beta_n
    beta_bar_n = (1/n**2)*0
    for i in range(n):
        beta_bar_n += (1/n**2)*norm(product_vect(z, i) - sn_2)**2

    # beta_n ( ^ 2 )
    beta_n = min(beta_bar_n, delta_n)

    # alpha_n ( ^ 2 )
    alpha_n = delta_n - beta_n

    # rho_n ( * 2 )
    rho_n = (beta_n/alpha_n)*sigma_n

    rho_1_n = (beta_n/delta_n)*sigma_n

    rho_2_n = alpha_n/delta_n

    Sigma_n_hat_ast = sn_2 + rho_n*I_q
```

```

Sigma_n_hat = rho_1_n*I_q + rho_2_n*sn_2

self_norm_sum = n*z_bar.dot(np.linalg.inv(Sigma_n_hat).dot(np.
↳transpose(z_bar)))
self_norm_sum_ast = n*z_bar.dot(np.linalg.inv(Sigma_n_hat_ast).dot(np.
↳transpose(z_bar)))
return np.array([self_norm_sum, self_norm_sum_ast, beta_n, sigma_n,
↳alpha_n, rho_n, max_p(sn_2),
min_p(sn_2)], dtype=np.int)

t1 = time.perf_counter()
dico = {}
for n in n_list[:3]:
    for q in q_list[:3]:
        if n <= q :
            dico[(n,q)] = np.stack(list(map(monte_carlo, range(10))))

print(dico[list(dico.keys())[0]])

t2 = time.perf_counter()

print(f'Finished in {t2-t1} seconds')

```

```

[[49 26 0 0 1 0 6 0]
 [62 32 1 1 1 0 5 0]
 [74 41 1 1 1 0 7 0]
 [43 23 0 0 1 0 6 0]
 [62 31 1 1 1 0 5 0]
 [58 30 1 1 1 0 6 0]
 [40 21 1 1 1 0 6 0]
 [48 26 0 0 1 0 7 0]
 [48 29 1 1 1 0 8 0]
 [43 22 0 0 0 0 5 0]]
Finished in 2.470908208997571 seconds

```

```
[11]: len(dico.keys())
```

```
[11]: 7
```

```
[12]: t1 = time.perf_counter()
data = {}
for q in q_list[:3]:
    for n in n_list[:3]:
        if n <= q :
            with concurrent.futures.ProcessPoolExecutor() as executor:

```

```

        f1 = np.stack(list(executor.map(monte_carlo, range(10))))
        #f2 = f1.result()
        data[(n,q)] = f1
        # if q/n%1==0 : print(str(i)+" "+str(j)+"",
                                #color.BLUE + color.BOLD + f"for n = %d \t",
↪and q = %d"%(n,q) + color.END, "\n",
                                #f1, "\n\n")

print(data[list(data.keys())[0]])

t2 = time.perf_counter()

print(f'Finished in {t2-t1} seconds')

```

```

[[48 26  1  1  1  0  6  0]
 [37 18  0  0  0  1  5  0]
 [74 37  1  0  1  0  6  0]
 [45 22  0  0  0  1  5  0]
 [63 31  0  0  0  0  6  0]
 [61 33  1  1  1  0  6  0]
 [62 34  1  1  1  0  7  0]
 [43 22  0  0  0  0  6  0]
 [67 35  1  1  1  0  7  0]
 [45 22  1  0  1  0  6  0]]
Finished in 3.6140988240003935 seconds

```

```
[13]: data[list(data.keys())[1]]
```

```
[13]: array([[102, 38,  2,  0,  1,  1,  9,  0],
           [ 88, 30,  1,  0,  1,  1,  7,  0],
           [ 74, 29,  2,  1,  1,  1,  8,  0],
           [101, 37,  1,  0,  1,  1,  8,  0],
           [ 99, 36,  2,  1,  1,  1,  8,  0],
           [115, 41,  2,  1,  1,  1,  8,  0],
           [104, 43,  2,  1,  1,  1,  9,  0],
           [ 90, 35,  2,  1,  1,  1,  8,  0],
           [117, 39,  2,  1,  1,  1,  7,  0],
           [ 96, 37,  1,  0,  1,  1,  8,  0]])
```

2 Generating step

```
[14]: K = int(input("Number of Monte Carlo iterations is : "))
```

```
Number of Monte Carlo iterations is : 999
```



```
[15]: t_init = time.perf_counter()
dico = {}
for n in n_list:
    for q in q_list:
        t1 = time.perf_counter()
        if n <= q :
            dico[(n,q)] = np.stack(list(map(monte_carlo, range(10))))
            t2 = time.perf_counter()
            if q==n or q==2*n :
                print(f"q = {q} and n = {n} --> ",f'Finished in {round(t2-t1,4)} seconds')

print(dico[list(dico.keys())[0]][0][:10])

t_fin = time.perf_counter()
print()
print(f'All loops finished in {round(t_fin-t_init, 3)} seconds')
```

```
q = 50 and n = 50 --> Finished in 0.077 seconds
q = 100 and n = 50 --> Finished in 0.1482 seconds
q = 150 and n = 75 --> Finished in 0.5359 seconds
q = 100 and n = 100 --> Finished in 0.2312 seconds
q = 200 and n = 100 --> Finished in 1.1857 seconds
q = 250 and n = 125 --> Finished in 1.9085 seconds
q = 150 and n = 150 --> Finished in 0.6856 seconds
q = 300 and n = 150 --> Finished in 3.45 seconds
q = 350 and n = 175 --> Finished in 6.3775 seconds
q = 200 and n = 200 --> Finished in 2.2651 seconds
q = 400 and n = 200 --> Finished in 10.6307 seconds
```

```
[[63 35 0 0 1 0 6 0]
 [63 33 0 0 1 0 6 0]
 [34 19 1 1 1 0 6 0]
 [45 22 1 1 1 1 5 0]
 [59 32 1 1 1 0 6 0]
 [32 18 1 1 1 0 7 0]
 [43 24 0 0 1 0 6 0]
 [29 16 1 1 1 0 7 0]
 [45 23 1 1 1 1 6 0]
 [45 23 1 1 1 0 6 0]]
```

All loops finished in 133.603 seconds

```
[16]: def monte_carlo(k):
    # random.seed(k)
    np.random.seed(int(os.getpid() * time.time()) % 123456789)
    z = generator(n,q)
    sn_2 = sn2(z)
```

```

s_2 = valeur_propre_collection[q_list.index(q)]
beta = (norm(sn_2 - s_2))**2
# calculate empirical mean
z_bar = zbar(z)
#empirical covariance matrix
sn_2 = sn2(z)

# Calculating empirical eigenvalues and eigenvectors (of  $S_n^2$ )
emp_val_p, emp_vec_p = np.linalg.eigh(sn_2)

# Calculating true (theoretical) eigenvalues and eigenvectors (of  $S^2$ )
# val_p, vec_p = np.linalg.eigh(s_2)

# Identity matrix of size q
I_q = np.diag(np.ones(q))

# sigma_n ( ^2 )
sigma_n = scalar(sn_2, I_q)

# delta_n ( ^2 )
delta_n = norm(sn_2 - sigma_n*I_q)**2

# intermediate beta_n
beta_bar_n = (1/n**2)*0
for i in range(n):
    beta_bar_n += (1/n**2)*norm(product_vect(z, i) - sn_2)**2

# beta_n ( ^2 )
beta_n = min(beta_bar_n, delta_n)

# alpha_n ( ^2 )
alpha_n = delta_n - beta_n

# rho_n ( *^2 )
rho_n = (beta_n/alpha_n)*sigma_n

rho_1_n = (beta_n/delta_n)*sigma_n

rho_2_n = alpha_n/delta_n

Sigma_n_hat_ast = sn_2 + rho_n*I_q
Sigma_n_hat = rho_1_n*I_q + rho_2_n*sn_2

self_norm_sum = n*z_bar.dot(np.linalg.inv(Sigma_n_hat).dot(np.
↪transpose(z_bar)))
self_norm_sum_ast = n*z_bar.dot(np.linalg.inv(Sigma_n_hat_ast).dot(np.
↪transpose(z_bar)))

```

```

    return np.array([self_norm_sum, self_norm_sum_ast, beta_n, sigma_n,
↳alpha_n, rho_n, max_p(sn_2),
                    min_p(sn_2), beta])

```

```

[17]: print(color.BLUE + color.BOLD + '***** Starting the extraction !'
↳*****' + color.END )

K = int(input("Number of Monte Carlo iterations is : "))
t_init = time.perf_counter()
dico = {}
for n in n_list:
    for q in q_list:
        t1 = time.perf_counter()
        if n <= q :
            dico[(n,q)] = np.stack(list(map(monte_carlo, range(K))))
            t2 = time.perf_counter()
            if q==n or q==2*n :
                print(f"q = {q} and n = {n} --> ",f'Finished in {round(t2-t1,
↳4)} seconds')

print(dico[list(dico.keys())[0]][0][:10])

t_fin = time.perf_counter()
print()
print(f'All loops finished in {round(t_fin-t_init, 3)} seconds')

```

***** Starting the extraction ! *****

```

Number of Monte Carlo iterations is : 999
q = 50 and n = 50 --> Finished in 6.9363 seconds
q = 100 and n = 50 --> Finished in 20.505 seconds
q = 150 and n = 75 --> Finished in 53.5549 seconds
q = 100 and n = 100 --> Finished in 22.8282 seconds
q = 200 and n = 100 --> Finished in 108.3973 seconds
q = 250 and n = 125 --> Finished in 212.258 seconds
q = 150 and n = 150 --> Finished in 74.594 seconds
q = 300 and n = 150 --> Finished in 387.8014 seconds
q = 350 and n = 175 --> Finished in 664.2058 seconds
q = 200 and n = 200 --> Finished in 90.4089 seconds
q = 400 and n = 200 --> Finished in 539.5729 seconds
[[4.36163387e+01 2.29947615e+01 1.01058375e+00 9.94405566e-01
 1.12688434e+00 8.91777509e-01 6.76539144e+00 2.92378122e-04
 1.05684057e+00]
 [3.90776396e+01 2.43712288e+01 1.19192070e+00 1.07589405e+00
 1.97523192e+00 6.49230289e-01 9.01775325e+00 9.11361451e-04
 1.53953565e+00]
 [3.77803838e+01 1.97392666e+01 8.70796720e-01 9.31511522e-01
 9.52761872e-01 8.51374517e-01 5.92505744e+00 1.12011788e-03

```

```

9.65588366e-01]
[4.76576721e+01 2.49843607e+01 9.04682944e-01 9.46184550e-01
9.96895627e-01 8.58662634e-01 6.09685173e+00 1.85657343e-04
9.67960342e-01]
[5.28018508e+01 3.10856052e+01 1.07685739e+00 1.03343395e+00
1.54146183e+00 7.21951696e-01 7.40632959e+00 4.79023767e-04
1.32548358e+00]
[6.35447898e+01 3.37727223e+01 1.13486920e+00 1.02885141e+00
1.28736851e+00 9.06975560e-01 7.12971728e+00 4.34679267e-04
1.26854087e+00]
[4.08157855e+01 2.27036168e+01 1.09457021e+00 1.03256946e+00
1.37204456e+00 8.23748584e-01 7.81942168e+00 9.61351707e-04
1.29813534e+00]
[6.00682046e+01 3.04724475e+01 9.66956786e-01 9.69569352e-01
9.95600139e-01 9.41674903e-01 5.70593205e+00 9.60919817e-04
9.34965244e-01]
[3.09277276e+01 1.64594267e+01 9.82699628e-01 9.66607448e-01
1.11793863e+00 8.49675248e-01 6.23098512e+00 9.45408080e-05
1.01211637e+00]
[5.44032587e+01 2.83334117e+01 1.00524975e+00 9.93486050e-01
1.09253249e+00 9.14116156e-01 5.80753461e+00 1.61521465e-03
1.15697731e+00]]

```

All loops finished in 12750.635 seconds

```
[18]: pickle.dump(dico, open("data_d_s0.6", "wb"))
```

```
[19]: teest = pickle.load(open("data_d_s0.6", "rb"))
len(teest[list(teest.keys())[0]])
```

```
[19]: 999
```

```
[20]: teest[list(teest.keys())[0]][:100,0]
```

```
[20]: array([43.61633868, 39.07763963, 37.78038384, 47.65767212, 52.80185077,
63.54478979, 40.81578545, 60.06820461, 30.9277276 , 54.40325868,
53.92542133, 40.7806492 , 61.91762462, 45.53339006, 46.01145098,
57.9904726 , 45.8014757 , 51.65765489, 51.99714112, 42.08810106,
48.90455907, 48.35879879, 66.13698697, 34.69292988, 36.92586344,
67.40620951, 40.74739076, 61.96337079, 37.28401014, 68.90944204,
63.65948672, 44.47166404, 36.93888047, 48.2314378 , 51.49386381,
51.70780367, 48.0522432 , 57.07099258, 54.70883515, 43.95208731,
41.61446553, 43.5558393 , 36.47337442, 64.30745563, 38.39860469,
41.46017687, 36.47325054, 52.44642022, 51.68395228, 34.84017003,
45.35290214, 30.11200769, 50.83476017, 56.97882203, 83.37324586,
35.64845542, 55.78649682, 44.74674965, 67.7756861 , 62.00105905,
49.7587842 , 57.20680248, 33.8051873 , 77.41768783, 44.98385818,
```

56.61212951, 56.45081119, 46.01350391, 33.94984957, 33.58601856,
52.43267599, 59.94202846, 42.56985293, 43.92285685, 42.10736153,
40.94954339, 31.93847436, 49.96644167, 68.7511026 , 42.08953012,
39.00069594, 65.85563002, 20.24246909, 55.51626525, 48.55358617,
65.19782297, 49.9842238 , 23.73902728, 61.15717356, 44.24313986,
38.03946275, 41.01256628, 55.3776742 , 74.00487526, 66.20808614,
26.72228924, 79.64053046, 39.51449616, 80.02408484, 59.61516822])

[]: