

# Regularized Precision Matrix Estimation via ADMM

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## Abstract

ADMMsigma is an R package that estimates a penalized precision matrix via the alternating direction method of multipliers (ADMM) algorithm. This report will provide a brief overview of the algorithm and detail how it can be utilized to estimate precision matrices of jointly normal distributions. In addition, examples and simulation results will be provided.

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## 1 Introduction

Suppose we want to solve the following optimization problem:

$$\begin{aligned} & \text{minimize } f(x) + g(z) \\ & \text{subject to } Ax + Bz = c \end{aligned}$$

where  $x \in \mathbb{R}^n, z \in \mathbb{R}^m, A \in \mathbb{R}^{p \times n}, B \in \mathbb{R}^{p \times m}, c \in \mathbb{R}^p$  – though we will later consider cases where  $x$  and  $z$  are matrices. Further, we will assume  $f$  and  $g$  are convex. The *augmented lagrangian* is constructed as follows:

$$L_\rho(x, z, y) = f(x) + g(z) + y^T(Ax + Bz - c) + \frac{\rho}{2} \|Ax + Bz - c\|_2^2$$

where  $y \in \mathbb{R}^p$  is the lagrange multiplier. The optimal value is

$$p^* = \inf \{f(x) + g(z) | Ax + Bz = c\}$$

Clearly, the minimization under the augmented lagrangian is equivalent to that of the usual lagrangian since any feasible point  $(x, z)$  satisfies the constraint  $\rho \|Ax + Bz - c\|_2^2 / 2 = 0$ .

The alternating direct method of multipliers (ADMM) algorithm consists of the following repeated iterations:

$$x^{k+1} := \arg \min_x L_\rho(x, z^k, y^k) \quad (1)$$

$$z^{k+1} := \arg \min_z L_\rho(z^{k+1}, z, y^k) \quad (2)$$

$$y^{k+1} := y^k + \rho(Ax^{k+1} + Bz^{k+1} - c) \quad (3)$$

A more complete introduction to the algorithm – specifically how it arose out of *dual ascent* and *method of multipliers* – can be found in Boyd et al. (2011).

## 2 Regularized Precision Matrix Estimation

We now consider the case where  $X_1, \dots, X_n$  are iid  $N_p(\mu, \Sigma)$  and we are tasked with estimating the precision matrix, denoted  $\Omega \equiv \Sigma^{-1}$ . The maximum likelihood estimator for  $\Omega$  is

$$\hat{\Omega} = \arg \min_{\Omega \in S_+^p} \{Tr(S\Omega) - \log \det(\Omega)\}$$

where  $S = \sum_{i=1}^n (X_i - \bar{X})(X_i - \bar{X})^T / n$ . It is straight forward to show that when the solution exists,  $\hat{\Omega} = S^{-1}$ .

We can construct a *penalized* likelihood estimator by adding a penalty term,  $P(\Omega)$ , to the likelihood:

$$\hat{\Omega}_\lambda = \arg \min_{\Omega \in S_+^p} \{Tr(S\Omega) - \log \det(\Omega) + P(\Omega)\}$$

Throughout the rest of this document we will take  $P(\Omega)$  to be  $P(\Omega) = \lambda \left[ \frac{1-\alpha}{2} \|\Omega\|_F^2 + \alpha \|\Omega\|_1 \right]$  so that the full penalized likelihood is as follows:

$$\hat{\Omega}_\lambda = \arg \min_{\Omega \in S_+^p} \left\{ Tr(S\Omega) - \log \det(\Omega) + \lambda \left[ \frac{1-\alpha}{2} \|\Omega\|_F^2 + \alpha \|\Omega\|_1 \right] \right\}$$

where  $0 \leq \alpha \leq 1$ ,  $\lambda > 0$ ,  $\|\cdot\|_F^2$  is the Frobenius norm and we define  $\|A\|_1 = \sum_{i,j} |A_{ij}|$ . This *elastic-net* penalty was explored by Hui Zou and Trevor Hastie (Zou and Hastie 2005) and is identical to the penalty used in the popular penalized regression package **glmnet**. Clearly, when  $\alpha = 0$  the elastic-net reduces to a ridge-type penalty and when  $\alpha = 1$  this reduces to a lasso-type penalty.

By letting  $f$  be equal to the non-penalized likelihood and  $g$  equal to  $P(\Omega)$ , our goal is to minimize the full augmented lagrangian where the constraint is that  $\Omega - Z$  is equal to zero:

$$L_\rho(\Omega, Z, \Lambda) = f(\Omega) + g(Z) + Tr[\Lambda(\Omega - Z)] + \frac{\rho}{2} \|\Omega - Z\|_F^2$$

The ADMM algorithm for regularized precision matrix estimation is

$$\Omega^{k+1} = \arg \min_{\Omega} \left\{ Tr(\Omega) - \log \det(\Omega) + Tr[\Lambda^k(\Omega - Z^k)] + \frac{\rho}{2} \|\Omega - Z^k\|_F^2 \right\} \quad (4)$$

$$Z^{k+1} = \arg \min_Z \left\{ \lambda \left[ \frac{1-\alpha}{2} \|Z\|_F^2 + \alpha \|Z\|_1 \right] + Tr[\Lambda^k(\Omega^{k+1} - Z)] + \frac{\rho}{2} \|\Omega^{k+1} - Z\|_F^2 \right\} \quad (5)$$

$$\Lambda^{k+1} = \Lambda^k + \rho(\Omega^{k+1} - Z^{k+1}) \quad (6)$$

## 2.1 Condensed-Form ADMM

An alternate form of the ADMM algorithm can be constructed by scaling the dual variable. Let us define  $R^k = \Omega - Z^k$  and  $U^k = \Lambda^k / \rho$ . Then

$$\begin{aligned} \text{Tr} [\Lambda^k (\Omega - Z^k)] + \frac{\rho}{2} \|\Omega - Z^k\|_F^2 &= \text{Tr} [\Lambda^k R^k] + \frac{\rho}{2} \|R^k\|_F^2 \\ &= \frac{\rho}{2} \|R^k + \Lambda^k / \rho\|_F^2 - \frac{\rho}{2} \|\Lambda^k / \rho\|_F^2 \\ &= \frac{\rho}{2} \|R^k + U^k\|_F^2 - \frac{\rho}{2} \|U^k\|_F^2 \end{aligned}$$

The condensed-form can now be written as follows:

$$\Omega^{k+1} = \arg \min_{\Omega} \left\{ \text{Tr}(\Omega) - \log \det(\Omega) + \frac{\rho}{2} \|\Omega - Z^k + U^k\|_F^2 \right\} \quad (7)$$

$$Z^{k+1} = \arg \min_Z \left\{ \lambda \left[ \frac{1-\alpha}{2} \|Z\|_F^2 + \alpha \|Z\|_1 \right] + \frac{\rho}{2} \|\Omega^{k+1} - Z + U^k\|_F^2 \right\} \quad (8)$$

$$U^{k+1} = U^k + \Omega^{k+1} - Z^{k+1} \quad (9)$$

More generally (in vector form),

$$x^{k+1} := \arg \min_x \left\{ f(x) + \frac{\rho}{2} \|Ax + Bz^k - c + u^k\|_2^2 \right\} \quad (10)$$

$$z^{k+1} := \arg \min_z \left\{ g(z) + \frac{\rho}{2} \|Ax^{k+1} + Bz - c + u^k\|_2^2 \right\} \quad (11)$$

$$u^{k+1} := u^k + Ax^{k+1} + Bz^{k+1} - c \quad (12)$$

Note that there are limitations to using this method. For instance, because the dual variable is scaled by  $\rho$  (the step size), this form limits one to using a constant step size (without making further adjustments to  $U^k$ ) – a limitation that could prolong the convergence rate. Because of this, we will only consider the non-condensed form for the remainder of this report.

## 2.2 Algorithm

$$\begin{aligned} \Omega^{k+1} &= \arg \min_{\Omega} \left\{ \text{Tr}(\Omega) - \log \det(\Omega) + \text{Tr} [\Lambda^k (\Omega - Z^k)] + \frac{\rho}{2} \|\Omega - Z^k\|_F^2 \right\} \\ Z^{k+1} &= \arg \min_Z \left\{ \lambda \left[ \frac{1-\alpha}{2} \|Z\|_F^2 + \alpha \|Z\|_1 \right] + \text{Tr} [\Lambda^k (\Omega^{k+1} - Z)] + \frac{\rho}{2} \|\Omega^{k+1} - Z\|_F^2 \right\} \\ \Lambda^{k+1} &= \Lambda^k + \rho (\Omega^{k+1} - Z^{k+1}) \end{aligned}$$

1. Decompose  $S + \Lambda^k - \rho Z^k = VQV^T$ .

$$\Omega^{k+1} = \frac{1}{2\rho} V \left[ -Q + (Q^2 + 4\rho I_p)^{1/2} \right] V^T$$

2. Elementwise soft-thresholding for all  $i = 1, \dots, p$  and  $j = 1, \dots, p$ .

$$\begin{aligned} Z_{ij}^{k+1} &= \frac{1}{\lambda(1-\alpha) + \rho} \text{sign}(\rho\Omega_{ij}^{k+1} + \Lambda_{ij}^k) (|\rho\Omega_{ij}^{k+1} + \Lambda_{ij}^k| - \lambda\alpha)_+ \\ &= \frac{1}{\lambda(1-\alpha) + \rho} \text{Soft}((\rho\Omega_{ij}^{k+1} + \Lambda_{ij}^k), \lambda\alpha) \end{aligned}$$

3. Update  $\Lambda$ .

$$\Lambda^{k+1} = \Lambda^k + \rho (\Omega^{k+1} - Z^{k+1})$$

### 2.2.1 Proof of (1):

$$\Omega^{k+1} = \arg \min_{\Omega} \left\{ \text{Tr}(\Omega) - \log \det(\Omega) + \text{Tr}[\Lambda^k (\Omega - Z^k)] + \frac{\rho}{2} \|\Omega - Z^k\|_F^2 \right\}$$

#### Code snippet:

Note this is not the actual code. The real code is written in c++.

```
# ridge penalized precision matrix
# function
RIDGEsigma = function(S, lam) {

  # dimensions
  p = dim(S)[1]

  # gather eigen values of S (spectral
# decomposition)
  e.out = eigen(S, symmetric = TRUE)

  # augment eigen values for omega hat
  new.evs = (-e.out$val + sqrt(e.out$val^2 +
    4 * lam))/(2 * lam)

  # compute omega hat for lambda (zero
# gradient equation)
  omega = tcrossprod(e.out$vec * rep(new.evs,
    each = p), e.out$vec)

  # compute gradient
  grad = S - qr.solve(omega) + lam * omega

  return(list(omega = omega, gradient = grad))
}
```

### 2.2.2 Proof of (2)

$$Z^{k+1} = \arg \min_Z \left\{ \lambda \left[ \frac{1-\alpha}{2} \|Z\|_F^2 + \alpha \|Z\|_1 \right] + \text{Tr} [\Lambda^k (\Omega^{k+1} - Z)] + \frac{\rho}{2} \|\Omega^{k+1} - Z\|_F^2 \right\}$$

Code snippet:

Note this is not the actual code. The real code is written in c++.

```
# ADMMsigma function
ADMMsigma = function(X = NULL, S = NULL,
  lam, alpha = 1, rho = 2, mu = 10, tau1 = 2,
  tau2 = 2, tol1 = 1e-04, tol2 = 1e-04,
  maxit = 1000) {

  # compute sample covariance matrix, if
  # necessary
  if (is.null(S)) {

    # covariance matrix
    n = dim(X)[1]
    S = (n - 1)/n * cov(X)

  }

  # allocate memory
  p = dim(S)[1]
  criterion = TRUE
  iter = lik = s = r = eps1 = eps2 = 0
  new.Z = Y = Omega = matrix(0, nrow = p,
    ncol = p)

  # loop until convergence
  while (criterion && (iter <= maxit)) {

    # ridge equation (1) gather eigen values
    # (spectral decomposition)
    Z = new.Z
    Omega = sigma_ridge(S + Y - rho *
      Z, lam = rho)$omega

    # penalty equation (2) soft-thresholding
    new.Z = soft(Y + rho * Omega, lam *
      alpha)/(lam * (1 - alpha) + rho)

    # update U (3)
    Y = Y + rho * (Omega - new.Z)

    # calculate new rho
    s = sqrt(sum((rho * (new.Z - Z))^2))
    r = sqrt(sum((Omega - new.Z)^2))
  }
}
```

```

rho = rho * (tau1 * (r > mu * s) +
  (s > mu * r)/tau2 + (s/mu <=
    r & r <= mu * s))
iter = iter + 1

# stopping criterion
eps1 = p * tol1 + tol2 * max(sqrt(sum(Omega^2)),
  sqrt(sum(new.Z^2)))
eps2 = p * tol1 + tol2 * sqrt(sum(Y^2))
criterion = (r >= eps1 || s >= eps2)

}
return(list(Iterations = iter, Omega = Omega))
}

```

## 3 R Package

### 3.1 Installation

```

# The easiest way to install is from CRAN
install.packages("ADMMsigma")

# You can also install the development
# version from GitHub:
# install.packages('devtools')
devtools::install_github("MGallow/ADMMsigma")

```

If there are any issues/bugs, please let me know: [github](#). You can also contact me via my website. Pull requests are welcome!

A (possibly incomplete) list of functions contained in the package can be found below:

- `ADMMsigma()` computes the estimated precision matrix (ridge, lasso, and elastic-net type regularization optional)
- `RIDGESigma()` computes the estimated ridge penalized precision matrix via closed-form solution
- `plot.ADMMsigma()` produces a heat map for cross validation errors
- `plot.RIDGESigma()` produces a heat map for cross validation errors

## 3.2 Usage

```
library(ADMMsigma)

# generate data from a dense matrix first
# compute covariance matrix
S = matrix(0.9, nrow = 5, ncol = 5)
diag(S) = 1

# generate 100 x 5 matrix with rows drawn
# from iid  $N_p(0, S)$ 
Z = matrix(rnorm(100 * 5), nrow = 100, ncol = 5)
out = eigen(S, symmetric = TRUE)
S.sqrt = out$vectors %*% diag(out$values^0.5) %*%
         t(out$vectors)
X = Z %*% S.sqrt

# elastic-net type penalty (use CV for
# optimal lambda and alpha)
ADMMsigma(X)

##
## Iterations:
## [1] 35
##
## Tuning parameters:
##      log10(lam)  alpha
## [1,]        -2.5    0.6
##
## Omega:
##      [,1]    [,2]    [,3]    [,4]    [,5]
## [1,]  7.50103 -1.71813 -2.17055 -1.21055 -1.95217
## [2,] -1.71813  6.88957 -2.51084 -1.42759 -1.14920
## [3,] -2.17055 -2.51084  8.17508 -1.79449 -1.38145
## [4,] -1.21055 -1.42759 -1.79449  6.25581 -1.97081
## [5,] -1.95217 -1.14920 -1.38145 -1.97081  6.53472

# ridge penalty (use CV for optimal
# lambda)
ADMMsigma(X, alpha = 0)

##
## Iterations:
## [1] 39
##
## Tuning parameters:
##      log10(lam)  alpha
## [1,]         -3      0
##
## Omega:
##      [,1]    [,2]    [,3]    [,4]    [,5]
## [1,]  7.92134 -1.80920 -2.32991 -1.24122 -2.07704
```

```
## [2,] -1.80920  7.24953 -2.70397 -1.48210 -1.17634
## [3,] -2.32991 -2.70397  8.69320 -1.89807 -1.43455
## [4,] -1.24122 -1.48210 -1.89807  6.54511 -2.08445
## [5,] -2.07704 -1.17634 -1.43455 -2.08445  6.84977
```

```
# lasso penalty (lam = 0.1)
ADMMsigma(X, lam = 0.1, alpha = 1)
```

```
##
## Iterations:
## [1] 10
##
## Tuning parameters:
##      log10(lam)  alpha
## [1,]         -1      1
##
## Omega:
##      [,1]      [,2]      [,3]      [,4]      [,5]
## [1,]  2.80422 -0.62667 -0.67987 -0.55554 -0.66056
## [2,] -0.62667  2.66766 -0.75422 -0.60328 -0.54369
## [3,] -0.67987 -0.75422  2.88706 -0.65217 -0.57631
## [4,] -0.55554 -0.60328 -0.65217  2.53979 -0.69921
## [5,] -0.66056 -0.54369 -0.57631 -0.69921  2.60963
```

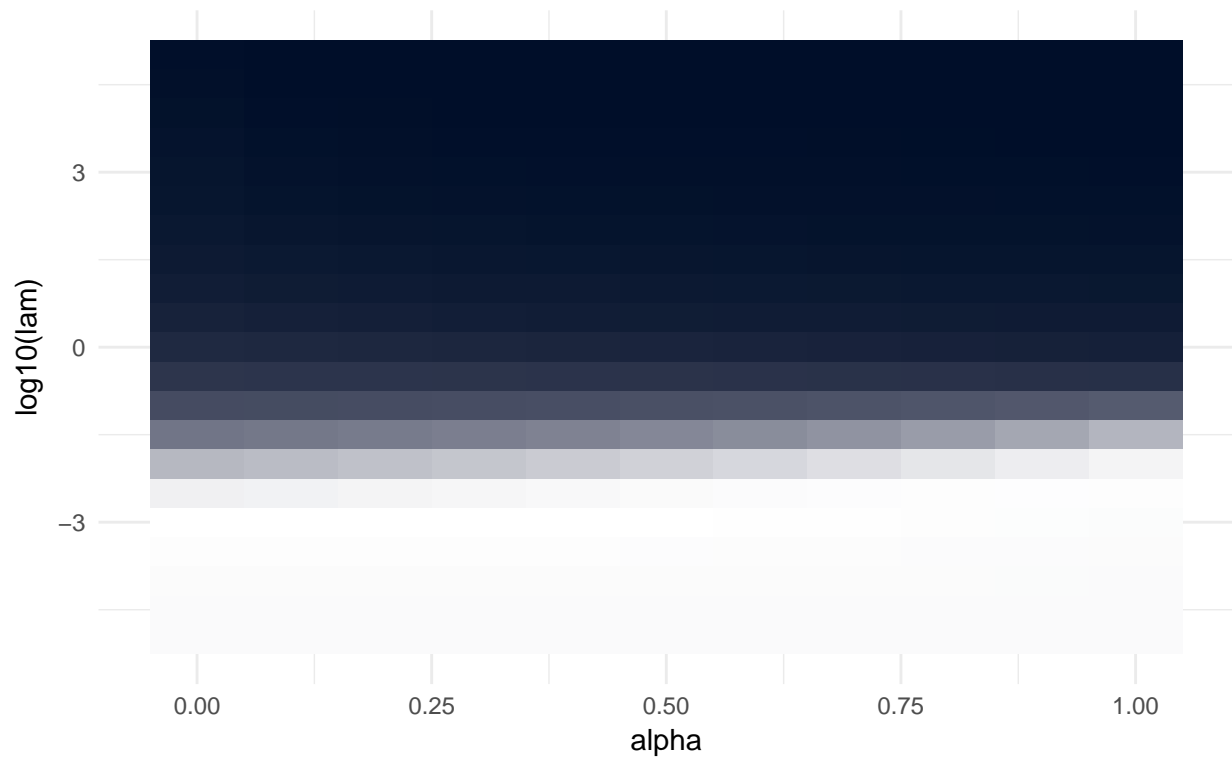
```
# ridge penalty no ADMM
RIDGEsigma(X, lam = 10^seq(-8, 8, 0.01))
```

```
##
## Tuning parameter:
##      lam log10(lam)
## [1,]  0.001      -3.1
##
## Omega:
##      [,1]      [,2]      [,3]      [,4]      [,5]
## [1,]  8.08365 -1.83697 -2.40822 -1.24051 -2.12903
## [2,] -1.83697  7.38256 -2.79786 -1.49474 -1.17783
## [3,] -2.40822 -2.79786  8.92397 -1.94165 -1.44583
## [4,] -1.24051 -1.49474 -1.94165  6.63854 -2.12649
## [5,] -2.12903 -1.17783 -1.44583 -2.12649  6.95522
```

```
# produce CV heat map for ADMMsigma
ADMMsigma(X) %>% plot
```

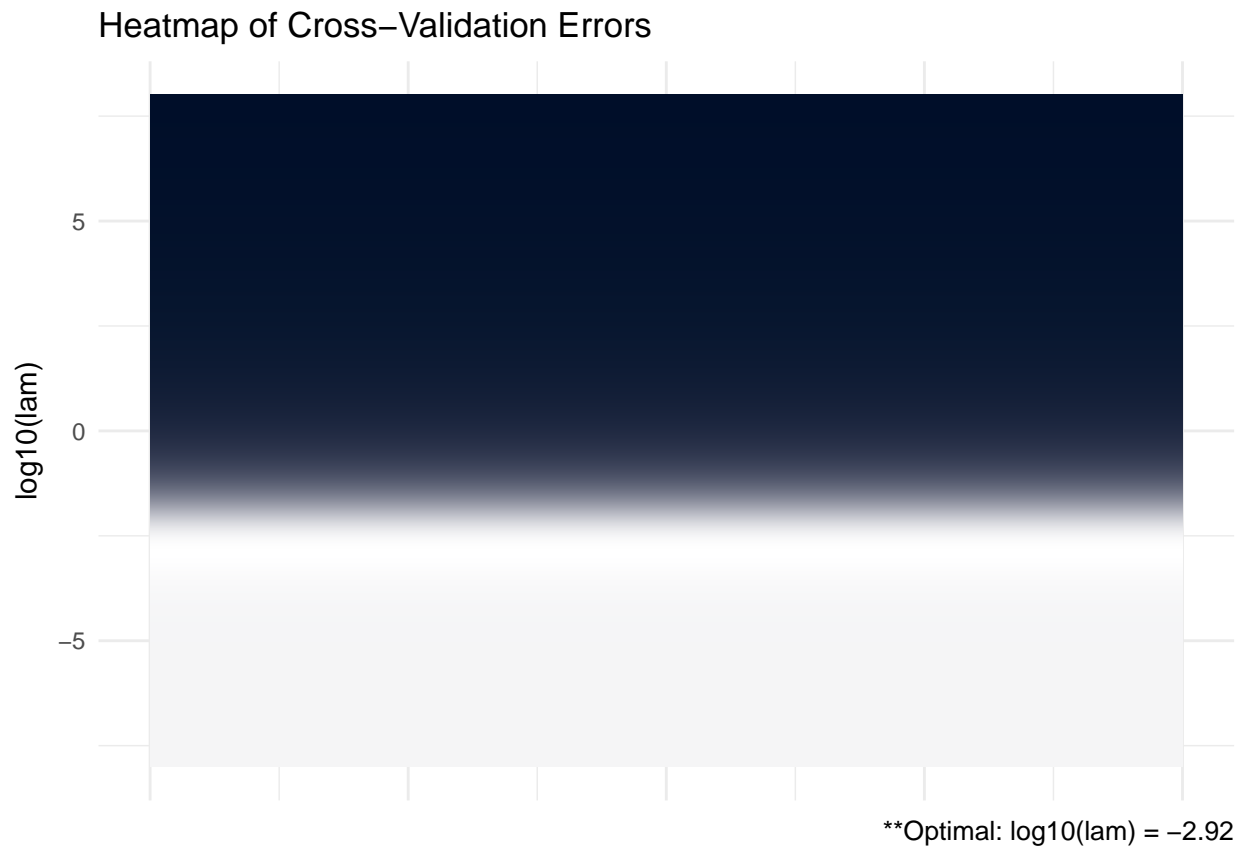


Heatmap of Cross-Validation Errors



**\*\*Optimal:  $\log_{10}(\text{lam}) = -3$ ,  $\alpha = 0.2$**

```
# produce CV heat map for RIDGESigma
RIDGESigma(X, lam = 10^seq(-8, 8, 0.01)) %>%
  plot
```



## 3.3 Benchmark

### 3.3.1 Computer Specs:

- MacBook Pro (Late 2016)
- Processor: 2.9 GHz Intel Core i5
- Memory: 8GB 2133 MHz
- Graphics: Intel Iris Graphics 550

```
# generate data from tri-diagonal  
# (sparse) matrix compute covariance  
# matrix (can confirm inverse is  
# tri-diagonal)  
S = matrix(0, nrow = 100, ncol = 100)  
  
for (i in 1:100) {  
  for (j in 1:100) {  
    S[i, j] = 0.7^(abs(i - j))  
  }  
}  
  
# generate 1000 x 100 matrix with rows  
# drawn from iid N_p(0, S)
```

```

Z = matrix(rnorm(1000 * 100), nrow = 1000,
           ncol = 100)
out = eigen(S, symmetric = TRUE)
S.sqrt = out$vectors %*% diag(out$values^0.5) %*%
         t(out$vectors)
X = Z %*% S.sqrt

# glasso (for comparison)
microbenchmark(glasso(s = S, rho = 0.1))

## Unit: milliseconds
##           expr      min       lq      mean    median      uq
##  glasso(s = S, rho = 0.1) 51.45916 53.67822 55.97162 55.03115 56.41237
##           max neval
## 101.6477    100

# benchmark ADMMsigma - default tolerance
microbenchmark(ADMMsigma(S = S, lam = 0.1,
                        alpha = 1, tol1 = 1e-04, tol2 = 1e-04))

## Unit: milliseconds
##           expr
##  ADMMsigma(S = S, lam = 0.1, alpha = 1, tol1 = 1e-04, tol2 = 1e-04)
##           min      lq      mean    median      uq      max neval
##  41.03643 42.0598 45.47058 42.81414 44.21306 203.617    100

# benchmark ADMMsigma - tolerance 1e-8
microbenchmark(ADMMsigma(S = S, lam = 0.1,
                        alpha = 1, tol1 = 1e-08, tol2 = 1e-08))

## Unit: milliseconds
##           expr
##  ADMMsigma(S = S, lam = 0.1, alpha = 1, tol1 = 1e-08, tol2 = 1e-08)
##           min      lq      mean    median      uq      max neval
## 189.5281 194.7297 204.5684 199.469 207.7964 301.8445    100

# benchmark ADMMsigma CV - default
# parameter grid
microbenchmark(ADMMsigma(X), times = 5)

## Unit: seconds
##           expr      min       lq      mean    median      uq      max neval
##  ADMMsigma(X) 48.25459 48.34069 49.79247 48.5219 48.78645 55.0587     5

# benchmark ADMMsigma parallel CV
microbenchmark(ADMMsigma(X, cores = 3), times = 5)

## Unit: seconds
##           expr      min       lq      mean    median      uq
##  ADMMsigma(X, cores = 3) 29.46813 30.01207 30.07205 30.21053 30.31711
##           max neval
##  30.35241     5

# benchmark ADMMsigma CV - likelihood
# convergence criteria
microbenchmark(ADMMsigma(X, crit = "loglik"),

```

```
times = 5)
```

```
## Unit: seconds
```

```
##          expr      min      lq      mean  median
## ADMMsigma(X, crit = "loglik") 76.10179 76.32729 76.84281 77.04886
##          uq      max neval
## 77.23379 77.5023      5
```

## References

- Boyd, Stephen, Neal Parikh, Eric Chu, Borja Peleato, Jonathan Eckstein, and others. 2011. “Distributed Optimization and Statistical Learning via the Alternating Direction Method of Multipliers.” *Foundations and Trends in Machine Learning* 3 (1). Now Publishers, Inc.: 1–122.
- Zou, Hui, and Trevor Hastie. 2005. “Regularization and Variable Selection via the Elastic Net.” *Journal of the Royal Statistical Society: Series B (Statistical Methodology)* 67 (2). Wiley Online Library: 301–20.