

Assume we have iid covariates  $X \in \mathbb{R}^{n \times p}$  and  $Y \in \mathbb{R}^p$  response. Also, we have  $\Omega \in \mathbb{R}^{k \times p}$  random projection matrix which follows  $\mathbb{E}[\Omega^T \Omega] = kI_p$  and has Rademacher random variable and iid elements. Let

$$\beta_n^\Omega = \Omega^T \underset{\beta \in \mathbb{R}^k}{\operatorname{argmin}} \mathbb{P}_n \{Y(\Omega X)^T \beta\}^2$$

Now assume we have a distribution for the dimension parameter  $k$ , i.e.  $k_1, \dots, k_B \sim Q$ , thus giving us  $B$  values of  $\Omega$  (call each individual realization  $\Omega(k(b))$  for an integer  $k \in [1, B]$  and a potential estimator ]

$$\hat{\beta}_n^{\text{ave}} = \frac{1}{B} \sum_{b=1}^B \beta_n^{\Omega(k_b)}$$

I implemented this streaming approach with  $k$  following a distribution by starting with a specific case and then using that to create a generalized function. The following gives some context for the R code.

We assume  $k$  follows a uniform distribution centered at an integer value near  $\log(p)$  in order to guarantee that  $p \gg k$ . Using this  $k$ ,  $B$  omega matrices are generated, each unique. For my algorithm, I made each entry of a given omega matrix a standard normal draw. The rest of the information needed to complete the problem must be inputted by the user. In order to generate a general algorithm for all cases of  $p, X, Y, B$ , I first started with a specific case (i.e. I chose  $p, X, Y, B$ ). The first thing I did was use an optimization function from class. The first part of the below code is the gradient/tuning function. The second part is a general case, which is a function that returns the desired  $\hat{\beta}_n^{\text{ave}}$ . The final part is my specific case, where I used  $p = 100$ ,  $X$  follows a uniform distribution between 5 and 15,  $Y = 10$ , and  $B = 1000$ . For my specific case, the general optimization solution turns out to be equivalent to OLS.

This general case function utilizes the **readline()** function in R, which asks the user to input values for  $p, Y$ , and  $B$ . For this reason, do not simply ctrl+A to run the code. If you would like to scroll to the general case, it's the first instance of the line with multiple **#** in a row (to signify a break), the second being the general case. The **genav** (general) function requires an input of  $X$  and returns  $\hat{\beta}_n^{\text{ave}}$  (assuming the user inputted  $p, y$ , and  $B$ ).

```
# gradient function , takes beta_hat , and sigma matrix as input
compute_min <- function (beta_hat , sig , maxit=500, tol=1e-8){
  fn = function (b){
    return (as.numeric(
      t(b-beta_hat) %*% sig %*% (b-beta_hat)))
  }
  gr = function (b){
    return(as.numeric(2*sig%*%(b-beta_hat)))
  }
  beta_cur <- beta_hat
  direction_cur <- gr (beta_cur)

  # Line search objective
  ls_obj = function (alpha){
    return (fn(beta_cur - alpha*direction_cur))
  }

  alpha_opt <- optimize (ls_obj , interval=c(0,1))$minimum
  beta_next <- beta_cur - alpha_opt*gr(beta_cur)
  for (k in 1:maxit){
    direction_next <- gr (beta_next)
    dx = beta_next - beta_cur
    if (max(abs(dx)) < tol){
      break
    }
  }
}
```

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    }
    dg <- direction_next - direction_cur
    alpha_opt <- sum(dg*dx)/sum(dg^2)
    beta_cur <- beta_next
    direction_cur <- direction_next
    beta_next <- beta_next - alpha_opt*direction_next
  }
  return (beta_next)

}

#####
# this is the general case
# this assumes x is defined correctly
# we assume a specific structure of omega
# however each omega will be different
p <- as.numeric(readline("enter value for p "))
y <- as.numeric(readline("enter scalar for y "))
B <- as.numeric(readline("enter large integer for B "))
genav <- function(x){
  bomg <- 0
  for (i in 1:B){
    # we draw K in a uniform distribution centered at log(N)
    k <- runif(1,1,2*ceiling(log(p)))
    omg <- matrix(0,nrow = k,ncol = p)
    for (c in 1:p){
      for (d in 1:k){
        omg[d,c] <- rnorm(1,0,1)
      }
    }
    q <- omg%%x
    bt <- t(solve(t(q)%%q)%%t(q)*y)
    sig <- q%%t(q)
    bomg1 <- t(omg)%%compute_min(bt,sig)
    bomg <- (bomg*(i-1))/i + bomg1/i
    rm(omg)
  }
  return(bomg)
}
far <- matrix(0,nrow=2,ncol = 2)
dim(far)
dim

#####

#example with a specific x, omega, y, B, p
# B from the problem
B <- 1000
# bomg is running average, see code from #2
bomg <- 0
# p is fixed far greater than what k could be
p <- 100

```

```

# y is a scalar
y <- 10
for (i in 1:B){
  # we draw K in a uniform distribution centered at log(N), per Dr. Laber
  k <- runif(1,1,10)
  x <- as.matrix(runif(p,5,15))
  omg <- matrix(0,nrow = k,ncol = p)
  for (c in 1:p){
    for (d in 1:k){
      omg[d,c] <- rnorm(1,0,1)
    }
  }
  q <- omg%*%x
  bt <- t(solve(t(q)%*%q)%*%t(q)*y)
  sig <- q%*%t(q)
  bomg1 <- t(omg)%*%compute_min(bt,sig)
  bomg <- (bomg*(i-1))/i + bomg1/i
  rm(omg)
}
# the next line returns the target
bomg

```