

Variable selection with the strong heredity constraint and its oracle property

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Trying to code the Strong Heredity Interaction Model (SHIM) in [Choi et al. \(2010\)](#).

1 Simulate the data

```
set.seed(123456)
p = 10 # number of predictors
n = 200 # number of test subjects
m = 200 # number of validation subjects
rho = 0.5
signal_to_noise_ratio = 4

main_effect_names <- paste0("x", 1:p) # names of the main effects
# names of the active set
true_var_names <- c("x1", "x2", "x3", "x4", "x1:x2", "x1:x3",
  "x1:x4", "x2:x3", "x2:x4", "x3:x4")
beta1 <- c(7, 2, 1, 1, 0, 0, 0, 0, 0, 0) %>% magrittr::set_names(true_var_names)
beta2 <- c(7, 2, 1, 1, 1, 0, 0, 0.5, 0.4, 0.1) %>% magrittr::set_names(true_var_names)
beta3 <- c(7, 2, 1, 1, 7, 7, 7, 2, 2, 1) %>% magrittr::set_names(true_var_names)
beta4 <- c(7, 2, 1, 1, 14, 14, 14, 4, 4, 2) %>% magrittr::set_names(true_var_names)
beta5 <- c(0, 0, 0, 0, 7, 7, 7, 2, 2, 1) %>% magrittr::set_names(true_var_names)

# simulate Toeplitz like structure
H <- abs(outer(1:p, 1:p, "-"))
cor <- rho^H

DT <- MASS::mvrnorm(n = n, mu = rep(0, p), Sigma = cor) %>% magrittr::set_colnames(paste0("x",
```

```

1:p)) %>% set_rownames(paste0("Subject", 1:n))
head(DT)

##           x1      x2      x3      x4      x5      x6      x7      x8
## Subject1 -0.16 -0.92 -1.54 -0.79 -1.04 -0.148  0.257 -0.28
## Subject2  0.37  0.28  0.36 -0.69  0.60 -0.151  0.011  0.73
## Subject3  0.71  1.10  1.26 -0.19 -0.83  0.323 -0.223  1.16
## Subject4 -0.44  0.66  0.44  0.50  0.33  0.025 -1.229 -0.99
## Subject5  1.30  0.15 -0.82 -1.04 -1.32 -1.796 -0.782 -1.47
## Subject6 -0.82 -1.71  0.93  0.87 -0.09 -0.321 -0.616 -1.68
##           x9      x10
## Subject1 -0.267  1.28
## Subject2 -0.098  0.32
## Subject3 -1.157  0.17
## Subject4  0.597 -0.48
## Subject5 -3.419 -2.32
## Subject6 -0.901 -1.24

# create X matrix which contains all main effects and
# interactions each column is standardized to mean 0 and sd 1
X <- model.matrix(as.formula(paste0("~(", paste0(main_effect_names,
  collapse = "+"), ")^2-1")), DT %>% as.data.frame()) %>% scale
crossprod(X) %>% diag

##      x1      x2      x3      x4      x5      x6      x7      x8
## 199    199    199    199    199    199    199    199
##      x9      x10  x1:x2  x1:x3  x1:x4  x1:x5  x1:x6  x1:x7
## 199    199    199    199    199    199    199    199
## x1:x8  x1:x9  x1:x10  x2:x3  x2:x4  x2:x5  x2:x6  x2:x7
## 199    199    199    199    199    199    199    199
## x2:x8  x2:x9  x2:x10  x3:x4  x3:x5  x3:x6  x3:x7  x3:x8
## 199    199    199    199    199    199    199    199
## x3:x9  x3:x10  x4:x5  x4:x6  x4:x7  x4:x8  x4:x9  x4:x10
## 199    199    199    199    199    199    199    199
## x5:x6  x5:x7  x5:x8  x5:x9  x5:x10  x6:x7  x6:x8  x6:x9
## 199    199    199    199    199    199    199    199
## x6:x10  x7:x8  x7:x9  x7:x10  x8:x9  x8:x10  x9:x10
## 199    199    199    199    199    199    199

apply(X, 2, mean) %>% sum

## [1] 1.6e-16

```

```

apply(X, 2, sd) %>% sum
## [1] 55

# # alternative, according to Bien et al 2013, they first
# # standardize the main effects, # then take all pairwise
# # products for interactions, then center those products #
# # first scale the main effects X_temp <- DT %>% scale
# # apply(X_temp, 2, mean) %>% sum apply(X_temp, 2, sd) %>% sum
# # then create all pairwise interactions X_temp_2 <-
# # model.matrix(as.formula(paste0('~(',paste0(main_effect_names,
# # collapse = '+'),')^2-1')), X %>% as.data.frame())
# # apply(X_temp_2[, (p+1):ncol(X_temp_2)], 2, sd) %>% hist
# # apply(X_temp_2[, (p+1):ncol(X_temp_2)], 2, mean) %>% hist #
# # then center the product terms and combine with standardized
# # main effects X <- cbind(X_temp_2[,1:p],
# # scale(X_temp_2[, (p+1):ncol(X_temp_2)], center = T, scale =
# # F)) dim(X) str(X) X %>% crossprod() %>% diag

y.star <- X[, names(beta4)] %*% beta4
error <- rnorm(n)
k <- sqrt(var(y.star)/(signal_to_noise_ratio * var(error)))
Y <- (y.star + k * error) %>% scale(center = TRUE, scale = FALSE)
colnames(Y) <- "Y"

(b0 <- mean(y.star + k * error))

## [1] 0.27

# names of interaction variables assuming interaction terms
# contain a ':'
interaction_names <- colnames(X) %>% grep(":", ., value = T)

```

References

Nam Hee Choi, William Li, and Ji Zhu. Variable selection with the strong heredity constraint and its oracle property. *Journal of the American Statistical Association*, 105(489):354–364, 2010. [1](#)

A R Code

```
sessionInfo()

getPkg <- function(pkg) install.packages(pkg, repos = "http://cran.r-project.org")

pkg = try(require(knitr))
if (!pkg) {
  cat("Installing 'knitr' from CRAN\n")
  getPkg("knitr")
  require(knitr)
}

pkg = try(require(data.table))
if (!pkg) {
  cat("Installing 'data.table' from CRAN\n")
  getPkg("data.table")
  require(data.table)
}

pkg = try(require(magrittr))
if (!pkg) {
  cat("Installing 'magrittr' from CRAN\n")
  getPkg("magrittr")
  require(magrittr)
}

pkg = try(require(glmnet))
if (!pkg) {
  cat("Installing 'glmnet' from CRAN\n")
  getPkg("glmnet")
  require(glmnet)
}

pkg = try(require(stringr))
if (!pkg) {
  cat("Installing 'stringr' from CRAN\n")
  getPkg("stringr")
  require(stringr)
}
```

```

pckg = try(require(plyr))
if (!pckg) {
  cat("Installing 'plyr' from CRAN\n")
  getPkg("plyr")
  require(plyr)
}
set.seed(123456)
p = 10 # number of predictors
n = 200 # number of test subjects
m = 200 # number of validation subjects
rho = 0.5
signal_to_noise_ratio = 4

main_effect_names <- paste0("x", 1:p) # names of the main effects
# names of the active set
true_var_names <- c("x1", "x2", "x3", "x4", "x1:x2", "x1:x3",
  "x1:x4", "x2:x3", "x2:x4", "x3:x4")
beta1 <- c(7, 2, 1, 1, 0, 0, 0, 0, 0, 0) %>% magrittr::set_names(true_var_names)
beta2 <- c(7, 2, 1, 1, 1, 0, 0, 0.5, 0.4, 0.1) %>% magrittr::set_names(true_var_names)
beta3 <- c(7, 2, 1, 1, 7, 7, 7, 2, 2, 1) %>% magrittr::set_names(true_var_names)
beta4 <- c(7, 2, 1, 1, 14, 14, 14, 4, 4, 2) %>% magrittr::set_names(true_var_names)
beta5 <- c(0, 0, 0, 0, 7, 7, 7, 2, 2, 1) %>% magrittr::set_names(true_var_names)

# simulate Toeplitz like structure
H <- abs(outer(1:p, 1:p, "-"))
cor <- rho^H

DT <- MASS::mvrnorm(n = n, mu = rep(0, p), Sigma = cor) %>% magrittr::set_colnames(paste0("x",
  1:p)) %>% set_rownames(paste0("Subject", 1:n))
head(DT)

# create X matrix which contains all main effects and
# interactions each column is standardized to mean 0 and sd 1
X <- model.matrix(as.formula(paste0("~(", paste0(main_effect_names,
  collapse = "+"), ")^2-1")), DT %>% as.data.frame()) %>% scale
crossprod(X) %>% diag

apply(X, 2, mean) %>% sum

```

```

apply(X, 2, sd) %>% sum

# # alternative, according to Bien et al 2013, they first
# standardize the main effects, # then take all pairwise
# products for interactions, then center those products #
# first scale the main effects X_temp <- DT %>% scale
# apply(X_temp, 2, mean) %>% sum apply(X_temp, 2, sd) %>% sum
# # then create all pairwise interactions X_temp_2 <-
# model.matrix(as.formula(paste0('~(', paste0(main_effect_names,
# collapse = '+'), ')^2-1')), X %>% as.data.frame())
# apply(X_temp_2[, (p+1):ncol(X_temp_2)], 2, sd) %>% hist
# apply(X_temp_2[, (p+1):ncol(X_temp_2)], 2, mean) %>% hist #
# then center the product terms and combine with standardized
# main effects X <- cbind(X_temp_2[, 1:p],
# scale(X_temp_2[, (p+1):ncol(X_temp_2)], center = T, scale =
# F)) dim(X) str(X) X %>% crossprod() %>% diag

y.star <- X[, names(beta4)] %*% beta4
error <- rnorm(n)
k <- sqrt(var(y.star)/(signal_to_noise_ratio * var(error)))
Y <- (y.star + k * error) %>% scale(center = TRUE, scale = FALSE)
colnames(Y) <- "Y"

(b0 <- mean(y.star + k * error))

# names of interaction variables assuming interaction terms
# contain a ':'
interaction_names <- colnames(X) %>% grep(":", ., value = T)
#' variables: character vector of variable names for which you want the univariate regression est
#' must be contained in the column names of x
#' x: matrix that includes all data corresponding to variables with corresponding column names
#' y: response (matrix form)
#' returns OLS coefficients as a p x 1 data.frame

uni_fun <- function(variables, x, y) {

  res <- plyr::ldply(variables, function(i) {

```

```

    fit <- lm.fit(x = x[, i, drop = F], y = y) # dont need to add intercept because y has be
    fit$coefficients[1]
  }) %>% magrittr::set_rownames(variables) %>% magrittr::set_colnames("univariate_beta") %>%
    as.matrix
  return(res)
}

# function that takes a vector of betas (which are the main
# effects) and alphas which are the interaction effects, and
# outputs the main effects, and converts the alphas to gammas
# note that  $\alpha_{ij} = \gamma_{ij} * \beta_i * \beta_j$ ,  $i < j$ 
convert <- function(betas.and.alphas, main.effect.names, interaction.names,
  epsilon = 1e-04) {

  # betas.and.alphas is the result from uni_fun create output
  # matrix
  betas_and_gammas <- matrix(nrow = nrow(betas.and.alphas)) %>%
    magrittr::set_rownames(rownames(betas.and.alphas))

  for (k in interaction.names) {
    # get names of main effects corresponding to interaction
    (main <- betas.and.alphas[k, , drop = F] %>% rownames %>%
      stringr::str_split(":") %>% unlist)
    # convert alpha to gamma BUT NEED TO CHECK IF BETAS ARE 0!!!!
    betas_and_gammas[k, ] <- if (any(abs(betas.and.alphas[main,
      ]) < epsilon))
      0 else betas.and.alphas[k, ]/prod(betas.and.alphas[main,
      ])
  }

  # add back the main effects which dont need to be transformed
  for (j in main.effect.names) {
    betas_and_gammas[j, ] <- betas.and.alphas[j, ]
  }

  return(betas_and_gammas)
}

```



```

# function that converts gammas to alphas
convert2 <- function(betas.and.gammas, main.effect.names, interaction.names) {

  # betas.and.gammas = rbind2(beta_hat_next,
  # gamma_hat_previous) betas.and.gammas is the result from
  # shim function create output matrix
  betas.and.alphas <- matrix(nrow = nrow(betas.and.gammas)) %>%
    magrittr::set_rownames(rownames(betas.and.gammas))

  for (k in interaction.names) {
    # k='x1:x10' get names of main effects corresponding to
    # interaction
    (main <- betas.and.gammas[k, , drop = F] %>% rownames %>%
      stringr::str_split(":") %>% unlist)
    # convert alpha to gamma
    betas.and.alphas[k, ] <- betas.and.gammas[k, ] * prod(betas.and.gammas[main,
      ])
  }

  # add back the main effects which dont need to be transformed
  for (j in main.effect.names) {
    betas.and.alphas[j, ] <- betas.and.gammas[j, ]
  }

  return(betas.and.alphas)
}

# function to calculate xtilde in step 3 of algorithm
xtilde <- function(interaction.names, data.main.effects, beta.main.effects) {

  #' interaction.names character vector of interaction names
  #' data.main.effects data frame or matrix containing the main effects data
  #' beta.main.effects data frame or matrix containing the coefficients of main effects
  #' Only the main effects are used in this step,
  #' however you can provide this function, either the betas.and.alphas or
  #' betas.and.gammas because only the betas (main effect) paramters
  #' are used in the calculation of xtilde

```

```

# create output matrix
xtildas <- matrix(ncol = length(interaction.names), nrow = nrow(data.main.effects)) %>%
  magrittr::set_colnames(interaction.names)

for (k in interaction.names) {
  # get names of main effects corresponding to interaction
  (main <- k %>% stringr::str_split(":") %>% unlist)

  # step 3 to calculate x tilda
  xtildas[, k] <- prod(beta.main.effects[main, ]) * data.main.effects[,
    main[1], drop = F] * data.main.effects[, main[2],
    drop = F]
}

return(xtildas)
}

# adaptive weights used in fitting algorithm, intercept set
# to 0 because Y is centered
ridge_weights <- function(x, y, main.effect.names, interaction.names) {
  #' interaction.names character vector of interaction names
  #' main.effect.names character vector of main effect names
  #' must be contained in the column names of x
  #' x: matrix that includes all data corresponding to variables with corresponding column names
  #' all columns have mean 0 and variance 1
  #' y: centered response (matrix form)
  #' returns ridge weights as a p x 1 data.frame
  #' STILL NEED TO DECIDE IF I SHOULD MULTIPLY WEIGHTS BY LOG(N)/N AS DONE IN SECTION 3.1 AND IN
  #' PAPER BY WANG LI AND TSAI 2007 JRSSB

  n <- length(y)
  # fit the ridge to get betas and alphas
  (fit <- glmnet::cv.glmnet(x = x, y = y, alpha = 0, standardize = F,
    intercept = F))
  # remove intercept
  (betas.and.alphas <- coef(fit, s = "lambda.1se") %>% as.matrix() %>%
    magrittr::extract(-1, , drop = F))
}

```

```

# (fit <- lm.fit(x = x, y = y)) # remove intercept
# (betas.and.alphas <- coef(fit) %>% as.matrix())

# create output matrix
weights <- matrix(nrow = nrow(betas.and.alphas)) %>% magrittr::set_rownames(rownames(betas.and.alphas))

# main effects weights
for (j in main.effect.names) {
  weights[j, ] <- abs(1/betas.and.alphas[j, ]) ## log(n)/n
}

for (k in interaction.names) {
  # get names of main effects corresponding to interaction
  (main <- betas.and.alphas[k, , drop = F] %>% rownames %>%
    stringr::str_split(":") %>% unlist)
  weights[k, ] <- abs(prod(betas.and.alphas[main, ])/betas.and.alphas[k,
    ]) ## log(n)/n
}

return(weights)
}

soft <- function(x, y, beta, lambda, weight) {
  # user must supply x AND y, or beta.. but not both i set it
  # up this way because to get the sequence of lambdas, I use
  # the beta argument so that I only compute this once. I use
  # the x, y argument for the CV folds lambda can be a vector
  # and this functions will return each thresholded beta for
  # each lambda e.g. soft(0.25, lambda =
  # seq(0.001,0.65,length.out = 100), 1.5)

  if (missing(x) & missing(y) & missing(beta))
    stop("user must supply x AND y, or beta but not both")
  if (missing(x) & missing(y))
    return(list(beta = sign(beta) * pmax(0, abs(beta) - lambda *
      weight)))
  if (missing(beta)) {
    # (beta <- lm.fit(x = cbind2(rep(1, length(y)),x), y = y) %>%

```

```

# coef %>% magrittr::extract(2))
(beta <- lm.fit(x = x[, 1, drop = F], y = y) %>% coef %>%
  magrittr::extract(1))

# lm.fit(x = cbind2(rep(1, length(y_tilde_2)), x_tilde_2), y =
# y_tilde_2) %>% coef %>% magrittr::extract(2)
b_lasso <- sign(beta) * pmax(0, abs(beta) - lambda *
  weight)
# return(list('beta' = b_lasso, 'df' = nonzero(b_lasso)))
return(b_lasso)
}

}

cv_lasspath <- function(outlist, lambda, x, y, foldid) {
  # typenames <- c(misclass = 'Misclassification Error', loss =
  # 'Margin Based Loss') if (pred.loss == 'default') pred.loss
  # <- 'loss' if (!match(pred.loss, c('loss'), FALSE)) {
  # warning('Only 'loss' available for least squares
  # regression; 'loss' used') pred.loss <- 'loss' }
  y <- as.double(y)
  nfolds <- max(foldid)
  predmat <- matrix(NA, length(y), length(lambda))
  nlams <- double(nfolds)
  for (i in seq(nfolds)) {
    # i=1
    which <- foldid == i
    # this gives be the fitted object for each CV fold
    fitobj <- outlist[[i]]

    # this gives the predicted responses for the subjects in the
    # held-out fold for each lambda so if each fold has 20
    # subjects, and there are 100 lambdas, then this will return
    # a 20 x 100 matrix
    preds <- x[which, , drop = FALSE] %*% t(fitobj$beta)
    # preds <- predict(fitobj, x[which, , drop = FALSE], type =
    # 'link')
    nlami <- length(fitobj$lambda)
    predmat[which, seq(nlami)] <- preds
  }
}

```

```

      nlams[i] <- nlami
    }
    cvraw <- (y - predmat)^2
    cvob <- cvcompute(cvraw, foldid, nlams)
    cvraw <- cvob$cvraw
    N <- cvob$N
    cvm <- apply(cvraw, 2, mean, na.rm = TRUE)
    cvsd <- sqrt(apply(scale(cvraw, cvm, FALSE)^2, 2, mean, na.rm = TRUE)/(N -
      1))
    list(cvm = cvm, cvsd = cvsd, name = "MSE")
  }

cvcompute <- function(mat, foldid, nlams) {
  nfolds <- max(foldid)
  outmat <- matrix(NA, nfolds, ncol(mat))
  good <- matrix(0, nfolds, ncol(mat))
  mat[is.infinite(mat)] <- NA
  for (i in seq(nfolds)) {
    mati <- mat[foldid == i, ]
    outmat[i, ] <- apply(mati, 2, mean, na.rm = TRUE)
    good[i, seq(nlams[i])] <- 1
  }
  N <- apply(good, 2, sum)
  list(cvraw = outmat, N = N)
}

lamfix <- function(lam) {
  llam <- log(lam)
  lam[1] <- exp(2 * llam[2] - llam[3])
  lam
}

nonzero <- function(beta, bystep = FALSE) {
  beta <- as.matrix(beta)
  nr = nrow(beta)
  if (nr == 1) {
    if (bystep)
      apply(beta, 2, function(x) if (abs(x) > 0)
        1 else NULL) else {

```

```

        if (any(abs(beta) > 0))
          1 else NULL
      }
    } else {
      beta = abs(beta) > 0
      which = seq(nr)
      ones = rep(1, ncol(beta))
      nz = as.vector((beta %*% ones) > 0)
      which = which[nz]
      if (bystep) {
        if (length(which) > 0) {
          beta = as.matrix(beta[which, , drop = FALSE])
          nz1 = function(x, which) if (any(x))
            which[x] else NULL
          which = apply(beta, 2, nz1, which)
          if (!is.list(which))
            which = data.frame(which)
          which
        } else {
          dn = dimnames(beta)[[2]]
          which = vector("list", length(dn))
          names(which) = dn
          which
        }
      } else which
    }
  }
}

getmin <- function(lambda, cvm, cvsd) {
  cvmin <- min(cvm)
  idmin <- cvm <= cvmin
  lambda.min <- max(lambda[idmin])
  idmin <- match(lambda.min, lambda)
  semin <- (cvm + cvsd)[idmin]
  idmin <- cvm <= semin
  lambda.1se <- max(lambda[idmin])
  list(lambda.min = lambda.min, lambda.1se = lambda.1se)
}

```

```

lambda.interp <- function(lambda, s) {
  if (length(lambda) == 1) {
    nums <- length(s)
    left <- rep(1, nums)
    right <- left
    sfrac <- rep(1, nums)
  } else {
    s[s > max(lambda)] <- max(lambda)
    s[s < min(lambda)] <- min(lambda)
    k <- length(lambda)
    sfrac <- (lambda[1] - s)/(lambda[1] - lambda[k])
    lambda <- (lambda[1] - lambda)/(lambda[1] - lambda[k])
    coord <- approx(lambda, seq(lambda), sfrac)$y
    left <- floor(coord)
    right <- ceiling(coord)
    sfrac <- (sfrac - lambda[right])/(lambda[left] - lambda[right])
    sfrac[left == right] <- 1
  }
  list(left = left, right = right, frac = sfrac)
}

plot.cv.uninet <- function(x, sign.lambda = 1, ...) {
  cvobj <- x
  xlab <- "log(Lambda)"
  if (sign.lambda < 0)
    xlab <- paste("-", xlab, sep = "")
  plot.args <- list(x = sign.lambda * log(cvobj$lambda), y = cvobj$cvu,
    ylim = range(cvobj$cvupper, cvobj$cvlo), xlab = xlab,
    ylab = cvobj$name, type = "n")
  new.args <- list(...)
  if (length(new.args))
    plot.args[names(new.args)] <- new.args
  do.call("plot", plot.args)
  error.bars(sign.lambda * log(cvobj$lambda), cvobj$cvupper,
    cvobj$cvlo, width = 0.01, col = "darkgrey")
  points(sign.lambda * log(cvobj$lambda), cvobj$cvu, pch = 20,
    col = "red")
  axis(side = 3, at = sign.lambda * log(cvobj$lambda), labels = paste(cvobj$nz),
    tick = FALSE, line = 0)
}

```

```

    abline(v = sign.lambda * log(cvobj$lambda.min), lty = 3)
    abline(v = sign.lambda * log(cvobj$lambda.1se), lty = 3)
    invisible()
  }

coef.cv.uninet <- function(object, s = c("lambda.1se", "lambda.min"),
  ...) {
  if (is.numeric(s))
    lambda <- s else if (is.character(s)) {
    s <- match.arg(s)
    lambda <- object[[s]]
  } else stop("Invalid form for s")
  coef.uninet(object$uninet.fit, s = lambda, ...)
}

coef.uninet <- function(object, s = NULL, type = c("coefficients",
  "nonzero"), ...) {
  type <- match.arg(type)
  # b0 <- t(as.matrix(object$b0)) rownames(b0) <- '(Intercept)'
  nbeta <- rbind2(object$beta)
  if (!is.null(s)) {
    vnames <- dimnames(nbeta)[[1]]
    # dimnames(nbeta) <- list(NULL, NULL)
    lambda <- object$lambda
    lamlist <- lambda.interp(lambda, s)
    nbeta <- nbeta[, lamlist$left, drop = FALSE] * lamlist$frac +
      nbeta[, lamlist$right, drop = FALSE] * (1 - lamlist$frac)
    # dimnames(nbeta) <- list(vnames, paste(seq(along = s)))
  }
  if (type == "coefficients")
    return(nbeta)
  if (type == "nonzero")
    return(nonzero(nbeta[1, , drop = FALSE], bystep = TRUE))
}

# fits weighted lasso with single predictor and no intercept
uninet <- function(x, y, nlambdas = 100, method = c("ls"), lambda.factor = ifelse(nobs <
  nvars, 0.01, 1e-04), lambda = NULL, penalty.factor = rep(1,
  nvars)) {

```



```

# x=x_tilde_2; y=y_tilde_2; nlambdas = 100; method = c('ls');
# lambda.factor = ifelse(nobs < nvars, 0.01, 1e-04); lambda =
# NULL; penalty.factor = rep(1, nvars); standardize = TRUE;
# eps = 1e-08; maxit = 1e+06

method <- match.arg(method)
this.call <- match.call()
y <- drop(y)
x <- as.matrix(x)
np <- dim(x)
nobs <- as.integer(np[1])
nvars <- as.integer(np[2])
vnames <- colnames(x)
N = nrow(x)

if (is.null(vnames))
  vnames <- paste("V", seq(nvars), sep = "")
if (length(y) != nobs)
  stop("x and y have different number of observations")
if (length(penalty.factor) != nvars)
  stop("The size of L1 penalty factor must be same as the number of input variables")
# if (lambda2 < 0) stop('lambda2 must be non-negative') maxit
# <- as.integer(maxit) lam2 <- as.double(lambda2)
penalty.factor <- as.double(penalty.factor)
# pf2 <- as.double(pf2) isd <- as.integer(standardize) eps <-
# as.double(eps) dfmax <- as.integer(dfmax) pmax <-
# as.integer(pmax) if (!missing(exclude)) { jd <-
# match(exclude, seq(nvars), 0) if (!all(jd > 0)) stop('Some
# excluded variables out of range') jd <-
# as.integer(c(length(jd), jd)) } else jd <- as.integer(0)

if (is.null(lambda)) {
  if (lambda.factor >= 1)
    stop("lambda.factor should be less than 1")
  # flmin <- as.double(lambda.factor)
  max_lam <- double(1)
  # loop to figure out max lambda its the minimum lambda that
  # gives a beta of 0
  beta_tilde <- lm.fit(x = x, y = y) %>% coef

```

```

    # ulam = 0
    b = 1
    while (b != 0) {
      max_lam <- max_lam + 0.001
      b <- soft(beta = beta_tilde, lambda = max_lam, weight = penalty.factor)
    }

    lambda <- seq(lambda.factor * max_lam, max_lam, length.out = 100)
  } else {
    flmin <- as.double(1)
    if (any(lambda < 0))
      stop("lambdas should be non-negative")
    ulam <- as.double(rev(sort(lambda)))
    nlam <- as.integer(length(lambda))
  }

  fit <- switch(method, ls = soft(x = x, y = y, lambda = lambda,
    weight = penalty.factor))

  fit$lambda <- if (is.null(lambda))
    lamfix(lambda) else lambda
  fit$call <- this.call
  # fit$ldf <- nonzero(fit$beta)
  class(fit) <- c("uninet", class(fit))
  fit
}

cv.uninet <- function(x, y, lambda = NULL, nfolds = 10, foldid,
  ...) {

  # penalty.factor = adaptive.weights[colnames(x),] lambda =
  # NULL x = x_tilde_2 ; y = y_tilde_2
  np <- dim(x)
  nobs <- as.integer(np[1])
  nvars <- as.integer(np[2])
  # lambda.factor = ifelse(nobs < nvars, 0.01, 1e-04) nfolds =
  # 10
  N = nrow(x)

```

```

y <- drop(y)

# if (is.null(lambda)) { if (lambda.factor >= 1)
# stop('lambda.factor should be less than 1') #flmin <-
# as.double(lambda.factor) ulam <- double(1) } # loop to
# figure out max lambda # its the minimum lambda that gives a
# beta of 0 beta_tilde <- lm.fit(x = x, y = y) %>% coef #ulam
# = 0 b = 1 while (b != 0) { ulam <- ulam + 0.001 b <-
# soft(beta = beta_tilde, lambda = ulam, weight =
# penalty.factor) } lambda <- seq(lambda.factor*ulam, ulam,
# length.out = 100)

# betas for each lambda sapply(lambda, function(i) soft(beta
# = beta_tilde, lambda = i, weight = penalty.factor))

uninet.object <- uninet(x, y, lambda = lambda, ...)

# uninet.object <- uninet(x, y, lambda = lambda,
# penalty.factor = adaptive.weights[colnames(x_tilde_2),] )

lambda <- uninet.object$lambda
nz <- sapply(coef(uninet.object, type = "nonzero"), length)

if (missing(foldid))
  foldid = sample(rep(seq(nfolds), length = N)) else nfolds = max(foldid)
if (nfolds < 3)
  stop("nfolds must be bigger than 3; nfolds=10 recommended")
outlist = as.list(seq(nfolds))

for (i in seq(nfolds)) {
  which = foldid == i
  if (is.matrix(y))
    y_sub = y[!which, ] else y_sub = y[!which]
  outlist[[i]] = uninet(x = x[!which, ], drop = FALSE,
    y = y_sub, lambda = lambda, ...)
}

cvstuff <- do.call(cv_lspath, list(outlist, lambda, x, y,

```

```

    foldid))
  cvm <- cvstuff$cvm
  cvsd <- cvstuff$cvsd
  cvname <- cvstuff$name
  out <- list(lambda = lambda, cvm = cvm, cvsd = cvsd, cvupper = cvm +
    cvsd, cvlo = cvm - cvsd, name = cvname, nzero = nz, name = cvname,
    uninet.fit = uninet.object)
  lamin <- getmin(lambda, cvm, cvsd)
  obj <- c(out, as.list(lamin))
  class(obj) <- "cv.uninet"
  obj
}

Q_theta <- function(x, y, beta, gamma, weights, lambda.beta,
  lambda.gamma, main.effect.names, interaction.names) {

  betas.and.alphas <- convert2(betas.and.gammas = rbind2(beta,
    gamma), main.effect.names = main.effect.names, interaction.names = interaction.names)
  crossprod(y - x %*% betas.and.alphas - b0) + lambda.beta *
    (crossprod(weights[main.effect.names, ], abs(beta))) +
    lambda.gamma * (crossprod(weights[interaction.names,
      ], abs(gamma)))
}

```

B Session Information

```

sessionInfo()

## R version 3.2.2 (2015-08-14)
## Platform: x86_64-pc-linux-gnu (64-bit)
## Running under: Ubuntu 14.04 LTS
##
## locale:
##  [1] LC_CTYPE=en_CA.UTF-8      LC_NUMERIC=C
##  [3] LC_TIME=en_CA.UTF-8      LC_COLLATE=en_CA.UTF-8
##  [5] LC_MONETARY=en_CA.UTF-8  LC_MESSAGES=en_CA.UTF-8
##  [7] LC_PAPER=en_CA.UTF-8     LC_NAME=C
##  [9] LC_ADDRESS=C             LC_TELEPHONE=C
## [11] LC_MEASUREMENT=en_CA.UTF-8 LC_IDENTIFICATION=C
##
## attached base packages:
## [1] stats      graphics  grDevices  utils      datasets
## [6] base
##
## other attached packages:
## [1] plyr_1.8.3      stringr_1.0.0   glmnet_2.0-2
## [4] foreach_1.4.3   Matrix_1.2-2    magrittr_1.5
## [7] data.table_1.9.4 knitr_1.11
##
## loaded via a namespace (and not attached):
##  [1] Rcpp_0.12.2      lattice_0.20-33  codetools_0.2-14
##  [4] MASS_7.3-44      chron_2.3-45     grid_3.2.2
##  [7] formatR_1.2.1    evaluate_0.8     highr_0.5.1
## [10] stringi_1.0-1    reshape2_1.4.1   iterators_1.0.8
## [13] tools_3.2.2      methods_3.2.2

```