

A brief introduction to the QPADM-slack algorithm

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Distributed big data

- The QPADM-slack algorithm is designed for solving estimation problems in distributed big data. In these estimation problems, the parameter vector to be estimated is $\beta \in \mathbb{R}^p$, i.e., the variable beta in our `paraQPADMsLackCcpp()` function in the code.
- In distributed environments, the full data (X, \mathbf{y}) is split into K partitions, with each partition stored on a local machine. In our code, $X \in \mathbb{R}^{n \times p}$ and $\mathbf{y} \in \mathbb{R}^n$ are respectively denoted as `x` and `y`. They are input variables of `paraQPADMsLackCcpp()`. The k th ($k = 1, 2, \dots, K$) partition of them are respectively denoted as `xk` and `yk` with

```
xk = x.rows(k*nk, k*nk+nk-1)
yk = y.subvec(nk*k, nk*k+nk-1)
```

Here `.rows()` and `.subvec()` are Rcpp functions.

`x.rows(k*nk, k*nk+nk-1)` is the sub-matrix of `x` from the $(k*nk+1)$ th row of `x` to the $(nk*k+nk)$ th row, and `y.subvec(nk*k, nk*k+nk-1)` is the sub-vector of `y` from the $(k*nk+1)$ th element of `y` to the $(nk*k+nk)$ th element. Note that, in Rcpp, the subscript starts from 0.

QPADM-slack is an iterative algorithm, and in each iteration, it needs to update $(1+5K)$ variables, i.e., the central variable β^{s+1} and the local variables $\xi_k^{s+1}, \eta_k^{s+1}, \beta_k^{s+1}, u_k^{s+1}$ and v_k^{s+1} ($k = 1, 2, \dots, K$), where s is the number of iterations that the algorithm has already executed (denoted by iteration in the code). In the following, we give the corresponding forms of these variables used in our code.

- β^{s+1} : beta;
- ξ_k^{s+1} : in the code, we use xi to represent the long vector composed of $\xi_1^{s+1}, \xi_2^{s+1}, \dots, \xi_K^{s+1}$, and its k th subvector, i.e., `xi.subvec(k* n_k , k* n_k + n_k -1)`, is then ξ_k^{s+1} ;
- η_k^{s+1} : similarly, we use eta to represent the long vector composed of $\eta_1^{s+1}, \eta_2^{s+1}, \dots, \eta_K^{s+1}$, and its k th subvector `eta.subvec(k* n_k , k* n_k + n_k -1)`, is then η_k^{s+1} . Besides, the long vector composed of $\eta_1^s, \eta_2^s, \dots, \eta_K^s$ is denoted by `etaini`. That is, `etaini` is the value of eta in the previous iteration.

- β_k^{s+1} : in the code, we use a matrix z to represent $\beta_1^{s+1}, \beta_2^{s+1}, \dots, \beta_K^{s+1}$, and the k th column of z , i.e., $z.col(k)$ is β_k^{s+1} . Correspondingly, $zini$ is the value of z in the previous iteration. Besides, $zmean$ is the mean of $zini$ calculated by rows.
- u_k^{s+1} : similarly, we use u to represent the matrix composed of $u_1^{s+1}, u_2^{s+1}, \dots, u_K^{s+1}$, and the k th column of u , i.e., $u.col(k)$ is u_k^{s+1} . Correspondingly, $uini$ is the value of u in the previous iteration, and $umean$ is the mean of $uini$.
- v_k^{s+1} : we use v to represent the long vector composed of $v_1^{s+1}, v_2^{s+1}, \dots, v_K^{s+1}$, and its k th subvector $v.subvec(k*nk, k*nk+nk-1)$ is v_k^{s+1} . $vini$ is then the value of v in the previous iteration. Besides, for simplicity, we use $vinik$ to represent $vini.subvec(nk*k, nk*k+nk-1)$ in the code.

Next, we give the update rules for these variables, and their specific implementation in the code.

β^{s+1} (a p -dimensional vector) is updated by element. Denote the j th element of β^{s+1} as β_j^{s+1} . In our algorithm, β_j^{s+1} is updated as follows.

Under SCAD penalty:

$$\beta_j^{s+1} = \arg \min_{x \in \{x_1, x_2, x_3, 0\}} \frac{1}{\rho K} p_{\lambda}^{\text{SCAD}}(x) + \frac{1}{2} (x - \psi_j^s)^2,$$

where

$$p_{\lambda}^{\text{SCAD}}(x) = \lambda |x| \mathbb{I}(0 \leq |x| \leq \lambda) + \frac{-x^2 + 2a\lambda|x| - \lambda^2}{2(a-1)} \mathbb{I}(\lambda < |x| \leq a\lambda) + \frac{(a+1)\lambda^2}{2} \mathbb{I}(|x| > a\lambda)$$

is the SCAD penalty function (implemented by `gpenalty()` in our code), and $\psi_j^s = z_j^s + u_j^s/\rho$, where z_j^s and u_j^s respectively are the j th element of `zmean` and `umean`. More specifically, β_j^{s+1} is one of the four values x_1 , x_2 , x_3 and 0, such that the function $\frac{1}{\rho K} p_{\lambda}^{\text{SCAD}}(x) + \frac{1}{2} (x - \psi_j^s)^2$ reaches its minimum. The definitions of x_1 , x_2 and x_3 are given in the next page.

$$\begin{aligned}x_1 &= \text{sign}(\psi_j^k) \min\left(\lambda, \max\left(0, |\psi_j^k| - \lambda/(\rho K)\right)\right), \\x_2 &= \text{sign}(\psi_j^k) \min\left(a\lambda, \max\left(\lambda, \frac{\rho K |\psi_j^k| (a-1) - a\lambda}{\rho K (a-1) - 1}\right)\right), \\x_3 &= \text{sign}(\psi_j^k) \max(a\lambda, |\psi_j^k|).\end{aligned}$$

Code:

```

zmean = mean(zini, 1);
umean = mean(uini, 1);
for(int j = 0; j < p; j++){
    phi = zmean(j)+umean(j)/pho;
    //calculate x1
    xscad(0) = sign(phi)*min(lambda, max(0.0, abs(phi)-lambda/(pho*K)));
    //calculate x2
    xscad(1) = sign(phi)*min(a*lambda, max(lambda, (pho*K*(a-1)*abs(phi)-a*lambda)/(pho*K*(a-1)-1)));
    //calculate x3
    xscad(2) = sign(phi)*max(a*lambda, abs(phi));
    arma::vec hscad(4, fill::zeros);
    //calculate the function value under x1, x2, x3 and 0
    for(int i = 0; i < 4; i++){
        hscad(i) = 0.5*(xscad(i)-phi)*(xscad(i)-phi)+gpenalty(xscad(i), a, lambda, penalty)/(pho*K);
    }
    beta(j) = xscad(hscad.indexmin());
}

```

Note: if the model contains an intercept, i.e., the input logic variable `intercept == TRUE`, we should update the first element of β^{s+1} by

```
beta(0) = zmean(0)+umean(0)/pho;
```

where `pho` is ρ and `phi` is ψ_j^s .

Under MCP penalty:

$$\beta_j^{s+1} = \arg \min_{x \in \{x_1, x_2, 0\}} \frac{1}{\rho K} p_\lambda^{\text{MCP}}(x) + \frac{1}{2} (x - \psi_j^k)^2,$$

where

$$p_\lambda^{\text{MCP}}(x) = \left(\lambda |x| - \frac{x^2}{2a} \right) \mathbb{I}(0 \leq |x| \leq a\lambda) + \frac{a\lambda^2}{2} \mathbb{I}(|x| > a\lambda)$$

is the MCP penalty function (implemented by `gpenalty()` in our code), and the definition of ψ_j^s is the same as that under the SCAD penalty. Thus, β_j^{s+1} is one of the three values x_1 , x_2 and 0, such that the function $\frac{1}{\rho K} p_\lambda^{\text{MCP}}(x) + \frac{1}{2} (x - \psi_j^k)^2$ reaches its minimum. Here,

$$x_1 = \text{sign}(\psi_j^k) \min \left(a\lambda, \max \left(0, \frac{a(\rho K |\psi_j^k| - \lambda)}{a\rho K - 1} \right) \right),$$

$$x_2 = \text{sign}(\psi_j^k) \max(a\lambda, |\psi_j^k|).$$

Code:

```
for(int j = 0; j < p; j++){
    phi = zmean(j)+umean(j)/pho;
    //calculate x1
    xmcj(0) = sign(phi)*min(a*lambda, max(0.0, a*(pho*K*abs(phi)-lambda)/(pho*K*a-1)));
    //calculate x2
    xmcj(1) = sign(phi)*max(a*lambda, abs(phi));
    //calculate the function value under x1, x2 and 0
    for(int i = 0; i < 3; i++){
        hmcj(i) = 0.5*(xmcj(i)-phi)*(xmcj(i)-phi)+gpenalty(xmcj(i), a, lambda, penalty)/(pho*K);
    }
    beta(j) = xmcj(hmcj.indexmin());
}
```

Similarly, when the input logic variable `intercept == TRUE`, we update the first element of β^{s+1} by

```
beta(0) = zmean(0)+umean(0)/pho;
```

$$\xi_k^{s+1} = \max(\mathbf{0}, \mathbf{y}_k - X_k \boldsymbol{\beta}_k^s + \boldsymbol{\eta}_k^s + \mathbf{v}_k^s / \rho - \tau \mathbf{1} / \rho)$$

```
arma::vec vinik = vini.subvec(nk*k, nk*k+nk-1), etainik = etaini.subvec(k*nk, k*nk+nk-1);
yx.subvec(k*nk, k*nk+nk-1) = yk-xk*zini.col(k);
xi.subvec(k*nk, k*nk+nk-1) = yx.subvec(k*nk, k*nk+nk-1)+etainik+vinik/pho-tau*arma::ones(nk)/(n*pho);
for(int i = k*nk; i<k*nk+nk; i++){
    if(xi(i) < 0){
        xi(i) = 0;
    }
}
```

$$\eta_k^{s+1} = \max(\mathbf{0}, -\mathbf{y}_k + X_k \boldsymbol{\beta}_k^s + \xi_k^{s+1} - \mathbf{v}_k^s / \rho - (1 - \tau) \mathbf{1} / \rho)$$

```
eta.subvec(k*nk, k*nk+nk-1) = -yx.subvec(k*nk, k*nk+nk-1)+xi.subvec(k*nk, k*nk+nk-1)-vinik/pho-(1-tau)*arma::ones(nk)/(n*pho);
for(int i = k*nk; i < k*nk+nk; i++){
    if(eta(i) < 0){
        eta(i) = 0;
    }
}
```

$$\beta_k^{s+1} = (I_p + X_k^T X_k)^{-1} [\beta^{s+1} - \mathbf{u}_k^s / \rho + X_k^T (\mathbf{y}_k - \xi_k^{s+1} + \eta_k^{s+1} + \mathbf{v}_k^s / \rho)]$$

Step 1: calculate $(I_p + X_k^T X_k)^{-1}$ (this step is completed before the “while” loop as the matrix inversion does not need to be updated)

```
arma::mat tmp2, xk = x.rows(k*nk, k*nk+nk-1);
if(nk > p) tmp2 = inv(arma::eye(p, p) + xk.t() * xk);
else tmp2 = arma::eye(p, p) - xk.t() * inv(arma::eye(nk, nk) + xk * xk.t()) * xk;
tmp.slice(k) = tmp2;
```

Step 2: calculate β_k^{s+1}

```
z.col(k) = tmp.slice(k) * (beta - uini.col(k) / pho + xk.t() * (yk - xi.subvec(k*nk, k*nk+nk-1) + eta.subvec(k*nk, k*nk+nk-1) + vi
```

$$\mathbf{u}_k^{s+1} = \mathbf{u}_k^s + \rho (\boldsymbol{\beta}_k^{s+1} - \boldsymbol{\beta}^{s+1})$$

```
u.col(k) = uini.col(k)+pho*(z.col(k)-beta);
```

$$\mathbf{v}_k^{s+1} = \mathbf{v}_k^s + \rho (\mathbf{y}_k - X_k \boldsymbol{\beta}_k^{s+1} - \boldsymbol{\xi}_k^{s+1} + \boldsymbol{\eta}_k^{s+1})$$

```
yx.subvec(k*nk,k*nk+nk-1) = yk-xk*z.col(k);
v.subvec(k*nk,k*nk+nk-1) = vinik+pho*(yx.subvec(k*nk,k*nk+nk-1)-xi.subvec(k*nk,k*nk+nk-1)+eta.subvec(k*nk,k*nk+nk-1));
```