MATLAB CODES

We implemented Algorithms 2.1, 2.2, NHZ1 and NHZ2 for constrained nonlinear monotone equations with the following Matlab codes:

(1) Code for Algorithm 2.1

end

```
function DLclustering(fnum,dimnum,xnum,xlrange,tol,maxit,
  maxfev)
% A modied Dai-Liao method for constrained systems and image
  de-blurring via clustering of eigenvalues
% Mohammed Yusuf Waziri, Kabiru Ahmed, Abubakar Sani Halilu,
  Salisu Murtala, Habibu Abdullahi,
% and Ya'u Balarabe Musa 2024
% Global convergence method
% call: dlcs(f,x0,tol,maxit)
% Input: dimnum= dimension
%
         fnum = function number
%
        xnum= initial iterate number
%
        tol= stoping tolerance
%
        maxit= maximum number of iteration
tic;
\%\%\% default maxit, fev and tol, constant input
  if nargin <7</pre>
   maxfev = 2000;
end
if nargin < 6
   maxit=1000; % default max. iter
end
if nargin < 5
    tol=10^(-10); % default tolerance
%%%%%%%%%%% variable input
  if nargin<4</pre>
   xlrange=[]; % excel range
end
if nargin < 3
   xnum=1; % default initial point
end
if nargin<2</pre>
    dimnum=1; % default problem
end
if nargin<1</pre>
    fnum=1; % default dimension
```

```
%%%%%%%%%%%%%%%% defining dimension
  switch dimnum
   case 1
      dim=5000;
   case 2
      dim=10000;
   case 3
      dim=50000;
   otherwise
      end
%%%%%%%%%%%%%% defining problems
  problem=fnum;
switch problem
     case 1
    1=0; u=+inf;
      f = 'cp0';
      proj = 'Pj';
     case 2
     1=0; u=+inf;
      f='cp5';
      proj = 'Pj';
     case 3
     1=0; u=+inf;
      f = 'k26';
      proj='Pj';
     case 4
     1=0; u=+inf;
      f='kab2';
      proj = 'Pj';
     case 5
     1=0; u=+inf;
      f='kab4';
      proj = 'Pj';
    case 6
     1=0; u=+inf;
      f='kab5';
      proj='Pj';
   otherwise
      f='fnum'; %for any other problem
end
%%%%%%%%%%%%%%%%% defining initial points
  guess=xnum;
```

```
switch guess
    case 1
        x0 = (((1:dim)')/dim);
        x0=1-(((1:dim)')/dim);
    case 3
        x0=2-(((1:dim)')/dim);
    case 4
        x0=((dim-(1:dim))/dim)';
    case 5
        x0 = (-1*((-1).^(1:dim)-2)')/4;
    case 6
        x0=(-2*((-1).^(1:dim)-2)')/2;
    otherwise
        x0=xnum; %for any other initial point
end
%Step 0 Initialization
ITER=0; %iteration
FEV=0; % function evaluation
bck=0; % backtracking counter
% line search parameters
% Step 1 stopping rule
F0=feval(f,x0); % evaluating F(x0);
FEV = FEV + 1;
norm_F0 = sqrt(sum(F0.^2)); % norm of F(x0)
d0=-F0; % initial direction
while(ITER<=maxit && norm_F0>tol)
    ro=0.49; m=0; sig=0.0001; %w=0.6;
F0=feval(f,x0);
   bita=1;
   % Step 3: line search
   while (-((feval(f,x0+bita*(ro)^m*d0))'*d0) <sig*bita*(ro)^</pre>
      m*(norm(d0))^2 && m<=10)
              m=m+1;
        FEV = FEV + 1;
    end
    if FEV>=maxfev
            disp('maximum number of function evalution reached
               ')
            return;
    end
   % backtracking counter
```

```
if m
        bck=bck+1;
    end
    alph=bita*(ro)^(m);
    z=x0+alph*d0;
    Fz = feval(f,z); % computing f(z)
    FEV = FEV + 1;
    if (feval(proj,(z),l,u)==z \& norm(Fz)<tol)
        x0=z;
        F0=Fz:
        norm_F0= norm(F0);
        disp('zk is in the convex set and its the solution at
           iteration number')
        disp((num2str(ITER)))
        break
    else
   zetak=Fz'*(x0-z)/(Fz'*Fz); % computing zetak
        P=feval(proj,(x0-1.8*zetak*Fz),l,u); % projection on
           convex set
    x=P;
    F1=feval(f,x);
    s=z-x0;
    y=Fz-F0;
    r = 0.01;
    gam=4;
    wk = y + r * s;
    %tk=2*gam*(s'*wk)/(s'*s)-(wk'*wk)/(s'*wk);
    tk = (1/gam)*(wk'*wk)/(s'*wk)+(1/gam)*(s'*wk)/(s'*s);
    bk1 = (F1'*wk)/(d0'*wk);
    bk2=(F1'*s)/(d0'*wk);
    d1 = -(1/gam)*F1+(1/gam)*bk1*d0-tk*bk2*d0;
    end
    x0=x;
    F0=F1;
    d0=d1;
    norm_F0=sqrt(sum(F0.^2));
    ITER = ITER + 1;
end
x0;
disp([num2str(ITER) ' / ' num2str(FEV) ' / ' num2str(bck)
   ' / ' num2str(toc) ' / ' num2str(norm_F0) ])
disp((num2str(f)))
disp((num2str(dim)))
table1='ClusteringDK.xlsx';
T={ITER,FEV,toc,norm_F0};
 sheet=fnum;
```

```
xclRange=xlrange;
 xlswrite(table1,T,sheet,xclRange);
  table1='dlcs.xlsx';
% T={num2str(ITER),num2str(FEV),num2str(toc),num2str(norm_F0)
  };
% sheet=fnum;
% xlRange=xlrange;
% xlswrite(table1,T,sheet,xlRange);
%winopen(table1)
toc;
(2) Code for Algorithm 2.2
function DKclustering (fnum, dimnum, xnum, xlrange, tol, maxit,
  maxfev)
% Image de-blurring with a Dai-Kou-type method via clustering
   of eigenvalues
% Mohammed Yusuf Waziri, \textbf{Kabiru Ahmed}, Abubakar Sani
  Halilu, Salisu Murtala, Habibu Abdullahi,
% and Ya'u Balarabe Musa 2024
% Global convergence method
% call: dlcs(f,x0,tol,maxit)
\% Input: dimnum= dimension
%
         fnum = function number
%
        xnum= initial iterate number
%
        tol= stoping tolerance
%
        maxit = maximum number of iteration
tic;
%%%%% default maxit, fev and tol, constant input
   if nargin<7</pre>
    maxfev = 2000;
end
if nargin < 6
    maxit=1000; % default max. iter
end
if nargin <5</pre>
    tol=10^(-10); % default tolerance
%%%%%%%%%%% variable input
   if nargin <4
    xlrange=[]; % excel range
end
if nargin < 3
    xnum=1; % default initial point
end
```

```
if nargin<2</pre>
   dimnum=1; % default problem
end
if nargin<1
   fnum=1; % default dimension
end
%%%%%%%%%%%%%%%%% defining dimension
  switch dimnum
   case 1
      dim = 5000;
   case 2
       dim=10000;
   case 3
       dim=50000;
   otherwise
       end
%%%%%%%%%%%%%%% defining problems
  problem=fnum;
switch problem
    case 1
     1=0; u=+inf;
      f='kab4';
      proj = 'Pj';
      case 2
     1=0; u=+inf;
      f='cp5';
      proj='Pj';
      case 3
     1=0; u=+inf;
      f='k26';
      proj='Pj';
      case 4
      1=0; u=+inf;
      f='kab2';
      proj = 'Pj';
     case 5
     1=0; u=+inf;
      f='kab4';
       proj='Pj';
     case 6
     1=0; u=+inf;
      f='kab5';
      proj='Pj';
```

```
otherwise
       f='fnum'; %for any other problem
end
%%%%%%%%%%%%%%%% defining initial points
  guess=xnum;
switch guess
   case 1
       x0 = (((1:dim)')/dim);
   case 2
       x0=1-(((1:dim)')/dim);
   case 3
       x0=2-(((1:dim)')/dim);
    case 4
       x0 = ((dim - (1:dim))/dim)';
   case 5
       x0 = (-1*((-1).^(1:dim)-2)')/4;
   case 6
       x0 = (-2*((-1).^(1:dim)-2)')/2;
    otherwise
       x0=xnum; %for any other initial point
end
%Step 0 Initialization
ITER=0; %iteration
FEV=0; % function evaluation
bck=0; % backtracking counter
% line search parameters
% Step 1 stopping rule
F0=feval(f,x0); % evaluating F(x0);
FEV = FEV + 1;
norm_F0 = sqrt(sum(F0.^2)); % norm of F(x0)
d0=-F0; % initial direction
while(ITER<=maxit && norm_F0>tol)
   ro=0.48; m=0; sig=0.0001;
F0=feval(f,x0);
   bita=1;
%
     tau=1;
%
     e=1;
     nm=(norm(((feval(f,x0+bita*(ro)^m*d0)))));
   % Step 3: line search
   while (-((feval(f,x0+bita*(ro)^m*d0))'*d0) <sig*bita*(ro)^</pre>
      m*(norm(d0))^2 && m<=10)
```

```
m=m+1;
                  FEV = FEV + 1;
    end
    if FEV>=maxfev
                                 disp('maximum number of function evalution reached
                                            ')
                                return;
    end
   % backtracking counter
                  bck=bck+1;
    end
    alph=bita*(ro)^(m);
   z=x0+alph*d0;
   Fz= feval(f,z); % computing f(z)
   FEV = FEV + 1;
    if (feval(proj,(z),l,u)==z \& norm(Fz)<tol)
                  x0=z;
                  F0=Fz;
                  norm_F0= norm(F0);
                  disp('zk is in the convex set and its the solution at
                             iteration number')
                  disp((num2str(ITER)))
                  break
   else
zetak=Fz'*(x0-z)/(Fz'*Fz); % computing zetak
                  P=feval(proj,(x0-1.8*zetak*Fz),1,u); % projection on
                             convex set
   x = P;
   F1=feval(f,x);
   s=z-x0;
   y=Fz-F0;
   r = 0.01;
   gam = 0.25;
   wk = y + r * s;
   %tk=2*gam*(s'*wk)/(s'*s)-(wk'*wk)/(s'*wk);
   %tauk=2*gam*(s'*wk)/(s'*s);
   bk2 = (gam*(F1'*wk)/(d0'*wk) - (gam*(wk'*wk)/(s'*wk) + gam*(s'*wk) + g
              wk)/(s'*s))*(F1'*s)/(d0'*wk));
   d1 = -gam*F1+bk2*d0;
    end
   x0=x;
   FO = F1;
   d0=d1;
   norm_F0=sqrt(sum(F0.^2));
```

```
ITER = ITER + 1;
end
x0;
disp([num2str(ITER) ' / ' num2str(FEV) ' / ' num2str(bck)
   ' / ' num2str(toc) ' / ' num2str(norm_F0) ])
 disp((num2str(f)))
disp((num2str(dim)))
table1='ClusteringDK.xlsx';
 T={ITER,FEV,toc,norm_F0};
 sheet=fnum;
 xclRange=xlrange;
 xlswrite(table1,T,sheet,xclRange);
 table1='dlcs.xlsx';
% T={num2str(ITER),num2str(FEV),num2str(toc),num2str(norm_F0)
% sheet=fnum;
% xlRange=xlrange;
% xlswrite(table1,T,sheet,xlRange);
%winopen(table1)
toc;
(3) Code for NHZ1
function NHZM1(fnum,dimnum,xnum,xlrange,tol,maxit,maxfev)
% Sparse signal reconstruction via Hager-Zhang-type schemes
   for constrained system of nonlinear equations,
% Mohammed Yusuf Waziri, Kabiru Ahmed, Abubakar Sani Halilu,
   and Salisu Murtala.
% Optimization, Vol 73, Issue 6, pp. 1949 - 1980, 2023.
% call: dlcs(f,x0,tol,maxit)
% Input: dimnum= dimension
%
          fnum= function number
%
        xnum= initial iterate number
%
        tol= stoping tolerance
%
         maxit = maximum number of iteration
%%%% default maxit, fev and tol, constant input
   if nargin <7
    maxfev = 2000;
end
if nargin < 6
    maxit=1000; % default max. iter
end
if nargin<5</pre>
    tol=10^(-10); % default tolerance
end
```

```
%%%%%%%%%%% variable input
  if nargin<4</pre>
   xlrange=[]; % excel range
end
if nargin<3</pre>
   xnum=1; % default initial point
end
if nargin<2</pre>
   dimnum=1; % default problem
end
if nargin<1</pre>
   fnum=1; % default dimension
\%\%\%\%\%\%\%\%\%\%\%\%\%\%\% defining dimension
  switch dimnum
   case 1
       dim = 5000;
   case 2
       dim=10000;
   case 3
      dim = 50000;
   otherwise
       end
%%%%%%%%%%%%%%% defining problems
  problem=fnum;
switch problem
     case 1
     1=0; u=+inf;
      f = 'cp0';
      proj='Pj';
      case 2
     1=0; u=+inf;
      f='cp5';
      proj = 'Pj';
      case 3
     1=0; u=+inf;
      f = 'k26';
      proj='Pj';
      case 4
      1=0; u=+inf;
      f='kab2';
      proj='Pj';
```

```
case 5
     1=0; u=+inf;
       f='kab4';
       proj = 'Pj';
     case 6
     1=0; u=+inf;
       f='kab5';
       proj='Pj';
    otherwise
       f='fnum'; %for any other problem
end
%%%%%%%%%%%%%%%%% defining initial points
  guess=xnum;
switch guess
   case 1
       x0 = (((1:dim)')/dim);
   case 2
       x0=1-(((1:dim)')/dim);
   case 3
       x0=2-(((1:dim)')/dim);
   case 4
       x0 = ((dim - (1:dim))/dim)';
   case 5
       x0 = (-1*((-1).^(1:dim)-2)')/4;
   case 6
       x0 = (-2*((-1).^(1:dim)-2)')/2;
    otherwise
       x0=xnum; %for any other initial point
end
%Step 0 Initialization
ITER=0; %iteration
FEV=0; % function evaluation
bck=0; % backtracking counter
% line search parameters
% Step 1 stopping rule
F0=feval(f,x0); % evaluating F(x0);
FEV = FEV + 1;
norm_F0 = sqrt(sum(F0.^2)); % norm of F(x0)
d0=-F0; % initial direction
while(ITER<=maxit && norm_F0>tol)
   ro=0.51; sig=0.0001; m=0;
F0=feval(f,x0);
```

```
bita=1;
   % Step 3: line search
   while (-((feval(f,x0+bita*(ro)^m*d0))'*d0) < sig*bita*(ro)
             )^m*(norm(d0))^2 && m <= 10)
                m=m+1;
                FEV = FEV + 1;
   end
   if FEV>=maxfev
                               disp('maximum number of function evalution reached
                                         ')
                              return;
   end
   % backtracking counter
   if m
                 bck=bck+1;
    end
   alph=bita*(ro)^(m);
   z=x0+alph*d0;
   Fz = feval(f,z); % computing f(z)
   FEV = FEV + 1;
   if (feval(proj,(z),1,u)==z & norm(Fz)<tol)</pre>
                 x0=z;
                 FO=Fz;
                 norm_F0=norm(F0);
                 disp('zk is in the convex set and its the solution at
                           iteration number')
                 disp((num2str(ITER)))
             break
   else
zetak=Fz'*(x0-z)/(Fz'*Fz); % computing zetak
                 P=feval(proj,(x0-1.8*zetak*Fz),1,u); % projection on
                           convex se
   x=P;
   F1=feval(f,x);
   s=z-x0;
   y=Fz-F0;
   c = 0.01;
   wk = y + c * s;
   gam=4;
   e1=(2)/(4*gam);
   \texttt{thet1=(1/gam)*sqrt((s'*wk)/(norm(s)*norm(wk)))^3;}
   thet2=max(thet1,e1);
   % modified y using line search
   betak = (F1 * wk)/(d0 * wk) - (gam*thet2*(wk * wk)*(F1 * d0))/((d0 * wk) + (gam*thet2*(wk * wk) * (gam*thet2*(wk) * (gam*thet2*(wk * wk) * (gam*thet2*(wk * wk) * (gam*thet2*(wk * wk
              '*wk)^2);
```

```
d1= -F1+betak*d0; % spectral Dai-Liao direction
    end
    x0=x;
    F0=F1;
    d0=d1;
    norm_F0=sqrt(sum(F0.^2));
    ITER = ITER + 1;
end
x0;
disp([num2str(ITER) ' / ' num2str(FEV) ' / ' num2str(bck)
   ' / ' num2str(toc) ' / ' num2str(norm_F0) ])
 disp((num2str(f)))
disp((num2str(dim)))
table1='MHZnew.xlsx';
 T={ITER,FEV,toc,norm_F0};
 sheet=fnum;
 xclRange=xlrange;
 xlswrite(table1,T,sheet,xclRange);
  table1='dlcs.xlsx';
% T={num2str(ITER),num2str(FEV),num2str(toc),num2str(norm_F0)
   };
% sheet=fnum;
% xlRange=xlrange;
% xlswrite(table1,T,sheet,xlRange);
%winopen(table1)
toc;
(4) Code for NHZ2
function NHZM2(fnum,dimnum,xnum,xlrange,tol,maxit,maxfev)
% Sparse signal reconstruction via Hager-Zhang-type schemes
   for constrained system of nonlinear equations,
% Mohammed Yusuf Waziri, Kabiru Ahmed, Abubakar Sani Halilu,
   and Salisu Murtala.
% Optimization, Vol 73, Issue 6, pp. 1949 - 1980, 2023.
% call: dlcs(f,x0,tol,maxit)
% Input: dimnum= dimension
%
         fnum = function number
%
         xnum= initial iterate number
%
        tol= stoping tolerance
         maxit = maximum number of iteration
%
%%%%% default maxit, fev and tol, constant input
   if nargin<7</pre>
    maxfev = 2000;
end
```

```
if nargin<6</pre>
   maxit=1000; % default max. iter
end
if nargin <5
   tol=10^(-10); % default tolerance
end
%%%%%%%%%%%% variable input
  if nargin<4</pre>
   xlrange=[]; % excel range
end
if nargin<3</pre>
   xnum=1; % default initial point
end
if nargin < 2
   dimnum=1; % default problem
end
if nargin<1
   fnum=1; % default dimension
%%%%%%%%%%%%%%%% defining dimension
  switch dimnum
   case 1
       dim = 5000;
   case 2
       dim=10000;
   case 3
       dim=50000;
   otherwise
                  % for any other dimension
       dim=dimnum;
end
\%\%\%\%\%\%\%\%\%\%\%\%\%\%\% defining problems
  problem=fnum;
switch problem
     case 1
     1=0; u=+inf;
       f = 'cp0';
       proj = 'Pj';
      case 2
     1=0; u=+inf;
       f = 'cp5';
       proj = 'Pj';
      case 3
     1=0; u=+inf;
```

```
f='k26';
        proj = 'Pj';
       case 4
       1=0; u=+inf;
        f='kab2';
        proj = 'Pj';
      case 5
      1=0; u=+inf;
        f='kab4';
        proj = 'Pj';
      case 6
      1=0; u=+inf;
        f='kab5';
        proj = 'Pj';
    otherwise
        f='fnum'; %for any other problem
end
%%%%%%%%%%%%%%%% defining initial points
   guess=xnum;
switch guess
    case 1
        x0 = (((1:dim)')/dim);
    case 2
        x0=1-(((1:dim)')/dim);
    case 3
        x0=2-(((1:dim)')/dim);
    case 4
        x0=((dim-(1:dim))/dim)';
    case 5
        x0 = (-1*((-1).^(1:dim)-2)')/4;
    case 6
        x0=(-2*((-1).^(1:dim)-2)')/2;
    otherwise
        x0=xnum; %for any other initial point
end
%Step 0 Initialization
ITER=0; %iteration
FEV=0; % function evaluation
bck=0; % backtracking counter
% line search parameters
% Step 1 stopping rule
F0=feval(f,x0); % evaluating F(x0);
FEV = FEV + 1;
```

```
norm_F0 = sqrt(sum(F0.^2)); % norm of F(x0)
d0=-F0; % initial direction
while(ITER<=maxit && norm_F0>tol)
    ro=0.50; sig=0.01; m=0;
F0=feval(f,x0);
    bita=1;
    % Step 3: line search
    while (-((feval(f,x0+bita*(ro)^m*d0))'*d0) < sig*bita*(ro</pre>
       ) m*(norm(d0)) 2 && m<=10)
        m=m+1;
        FEV = FEV + 1;
    end
    if FEV>=maxfev
            disp('maximum number of function evalution reached
               ')
            return;
    % backtracking counter
    if m
        bck=bck+1;
    end
    alph=bita*(ro)^(m);
    z=x0+alph*d0;
    Fz= feval(f,z); % computing f(z)
    FEV = FEV + 1:
    if (feval(proj,(z),l,u)==z \& norm(Fz)<tol)
        x0=z;
        F0=Fz;
        norm_F0=norm(F0);
        disp('zk is in the convex set and its the solution at
           iteration number')
        disp((num2str(ITER)))
       break
    else
   zetak=Fz'*(x0-z)/(Fz'*Fz); % computing zetak
        P=feval(proj,(x0-1.8*zetak*Fz),1,u); % projection on
           convex set
    x = P:
    F1=feval(f,x);
    s=z-x0;
    y = Fz - F0;
    c = 0.1;
    wk = y + c * s;
    gam=4;
```

```
e1=(2)/(4*gam);
    thet1=(1/gam)*(((s'*wk)^2)/((s'*s)*(wk'*wk)));
    thet2=max(thet1,e1);
    betak=((F1'*wk)/(d0'*wk))-(gam*thet2*(wk'*wk)*(F1'*d0))/((
       d0'*wk)^2);
    d1 = -F1 + betak * d0;
    end
    x0=x;
    F0=F1;
    d0=d1;
    norm_F0=sqrt(sum(F0.^2));
    ITER = ITER + 1;
end
x0;
disp([num2str(ITER) ' / ' num2str(FEV) ' / ' num2str(bck)
   ' / ' num2str(toc) ' / ' num2str(norm_F0) ])
 disp((num2str(f)))
disp((num2str(dim)))
table1='MHZnew.xlsx';
T={ITER,FEV,toc,norm_F0};
 sheet=fnum;
xclRange=xlrange;
xlswrite(table1,T,sheet,xclRange);
toc;
```

We implemented Algorithms 2.1, 2.2, NHZ1 and NHZ2 for image de-blurring and signal recovery with the following Matlab codes:

(1) Code for Algorithm 2.1 Image de-blurring

```
function [x,x_debias,objective,times,debias_start,mses,taus]=
...
   DLcluster(y,A,tau,varargin)

%   HTTCGP_CS 1.0, Nov. 29, 2019

%   This function solves the convex problem
% arg min_x = 0.5*|| y - A x ||_2^2 + tau || x ||_1
% using the algorithm modified three-term conjugate gradient method, described in the following paper
%
%   This code is to use the well-known code CG_DESCENT to solve \ell_1 norm
% regularization least square problems.
%
%
```

```
% Copyright (2019): Jianghua Yin
   ______
%
%
% The first version of this code by Jianghua Yin, Nov. 29,
  2019
% test for number of required parametres
if (nargin-length(varargin)) ~= 3
  error('Wrong number of required parameters');
end
% flag for initial x (can take any values except 0,1,2)
Initial_X_supplied = 3333;
% Set the defaults for the optional parameters
stopCriterion = 3;
tolA = 0.01;
tolD = 0.0001;
debias = 0;
maxiter = 10000;
maxiter_debias = 500;
miniter = 5;
miniter_debias = 5;
init = 0;
compute_mse = 0;
AT = 0;
verbose = 1;
continuation = 0;
cont_steps = -1;
firstTauFactorGiven = 0;
% Set the defaults for outputs that may not be computed
debias_start = 0;
x_{debias} = [];
mses = [];
% Read the optional parameters
if (rem(length(varargin),2)==1)
  error('Optional parameters should always go by pairs');
else
  for i=1:2:(length(varargin)-1)
    switch upper(varargin{i})
    case 'STOPCRITERION'
```

```
stopCriterion = varargin{i+1};
case 'TOLERANCEA'
 tolA = varargin{i+1};
case 'TOLERANCED'
 tolD = varargin{i+1};
case 'DEBIAS'
 debias = varargin{i+1};
case 'MAXITERA'
 maxiter = varargin{i+1};
case 'MAXITERD'
 maxiter_debias = varargin{i+1};
case 'MINITERA'
 miniter = varargin{i+1};
case 'MINITERD'
 miniter_debias = varargin{i+1};
case 'INITIALIZATION'
 as array
        used below
        x = varargin{i+1};
 else
        init = varargin{i+1};
 end
case 'MONOTONE'
  enforceMonotone = varargin{i+1};
case 'CONTINUATION'
  continuation = varargin{i+1};
case 'CONTINUATIONSTEPS'
 cont_steps = varargin{i+1};
case 'FIRSTTAUFACTOR'
 firstTauFactor = varargin{i+1};
 firstTauFactorGiven = 1;
case 'TRUE_X'
 compute_mse = 1;
 true = varargin{i+1};
case 'ALPHAMIN'
 alphamin = varargin{i+1};
case 'ALPHAMAX'
 alphamax = varargin{i+1};
case 'AT'
 AT = varargin{i+1};
case 'VERBOSE'
 verbose = varargin{i+1};
otherwise
% Hmmm, something wrong with the parameter string
```

```
error(['Unrecognized option: ''' varargin{i} '''']);
    end;
  end;
end
%%%%%%%%%%%%%%%%%%%
if (sum(stopCriterion == [0 1 2 3 4 5]) == 0)
  error(['Unknown stopping criterion']);
end
% if A is a function handle, we have to check presence of AT,
if isa(A, 'function_handle') & ~isa(AT, 'function_handle')
  error(['The function handle for transpose of A is missing'])
end
% if A is a matrix, we find out dimensions of y and x,
% and create function handles for multiplication by A and A',
% so that the code below doesn't have to distinguish between
% the handle/not-handle cases
if ~isa(A, 'function_handle')
 AT = @(x) (x'*A)'; %A'*x;
 A = 0(x) A*x;
end
\% from this point down, A and AT are always function handles.
% Precompute A'*y since it'll be used a lot
Aty = AT(y);
% Initialization
switch init
             % initialize at zero, using AT to find the size
    case 0
       of x
       x = AT(zeros(size(y)));
    case 1 % initialize randomly, using AT to find the size
       x = randn(size(AT(zeros(size(y)))));
    case 2
             % initialize x0 = A'*y
       x = Aty;
    case Initial_X_supplied % initial x was given by user
       % initial x was given as a function argument; just
          check size
       if size(A(x)) ~= size(y)
          error(['Size of initial x is not compatible with A'
             ]);
       end
```

```
otherwise
       error(['Unknown ''Initialization'' option']);
end
% now check if tau is an array; if it is, it has to
\% have the same size as x
if prod(size(tau)) > 1
   try,
      dummy = x.*tau;
   catch,
      error(['Parameter tau has wrong dimensions; it should be
          scalar or size(x)']),
   end
end
% if the true x was given, check its size
if compute_mse & (size(true) ~= size(x))
   error(['Initial x has incompatible size']);
end
% if tau is scalar, we check its value; if it's large enough,
\% the optimal solution is the zero vector
if prod(size(tau)) == 1
   aux = AT(y);
   max_tau = max(abs(aux(:)));
   if tau >= max_tau
      x = zeros(size(aux));
      if debias
         x_debias = x;
      objective(1) = 0.5*(y(:)'*y(:));
      times(1) = 0;
      if compute_mse
          mses(1) = sum(true(:).^2);
      end
      return
   end
                          %
\verb"end"
\% initialize u and v
u = x.*(x >= 0);
v = -x.*(x < 0);
\% define the indicator vector or matrix of nonzeros in x
nz_x = (x = 0.0);
```

```
num_nz_x = sum(nz_x(:));
% start the clock
t0 = cputime;
% store given tau, because we're going to change it in the
% continuation procedure
final_tau = tau;
% store given stopping criterion and threshold, because we're
% to change them in the continuation procedure
final_stopCriterion = stopCriterion;
final_tolA = tolA;
% set continuation factors
if continuation&&(cont_steps > 1)
   % If tau is scalar, first check top see if the first factor
   % too large (i.e., large enough to make the first
   % solution all zeros). If so, make it a little smaller than
      that.
   % Also set to that value as default
   if prod(size(tau)) == 1
      if (firstTauFactorGiven == 0)|(firstTauFactor*tau >=
         firstTauFactor = 0.5*max_tau / tau;
         if verbose
             fprintf(1,'\n setting parameter FirstTauFactor\n'
                )
         end
      end
   end
   cont_factors = 10.^[log10(firstTauFactor):...
                    log10(1/firstTauFactor)/(cont_steps-1):0];
end
if ~continuation
  cont_factors = 1;
  cont_steps = 1;
end
iter = 1;
if compute_mse
       mses(iter) = sum((x(:)-true(:)).^2);
end
```

```
keep_continuation = 1;
cont_loop = 1;
iter = 1;
taus = [];
sigma = 0.0001;
% loop for continuation
while keep_continuation
    % Compute and store initial value of the objective
       function
    resid = y - A(x);
    if cont_steps == -1
        gradq = AT(resid);
        tau = max(final_tau, 0.2*max(abs(gradq)));
        if tau == final_tau
            stopCriterion = final_stopCriterion;
            tolA = final_tolA;
                                                      % stop
            keep_continuation = 0;
               continuation
        else
            stopCriterion = 1;
            tolA = 1e-5;
        end
    else
        tau = final_tau * cont_factors(cont_loop); %
        if cont_loop == cont_steps
            stopCriterion = final_stopCriterion;
            tolA = final_tolA;
                                                     %
            keep_continuation = 0;
        else
            stopCriterion = 1;
            tolA = 1e-5;
        end
    end
    taus = [taus tau];
    if verbose
        fprintf(1, '\nSetting tau = \%0.5g\n', tau)
    end
    % if in first continuation iteration, compute and store
    % initial value of the objective function
    if cont_loop == 1
        alpha = 1.0;
        f = 0.5*(resid(:)'*resid(:)) + ...
             sum(tau(:).*u(:)) + sum(tau(:).*v(:));
```

```
objective(1) = f;
    if compute_mse
        mses(1) = (x(:)-true(:))'*(x(:)-true(:));
    end
    if verbose
        fprintf(1, 'Initial obj=%10.6e, alpha=%6.2e,
           nonzeros=%7d\n',...
            f,alpha,num_nz_x);
    end
end
% Compute the initial gradient and the useful
% quantity resid_base
resid_base = y - resid;
% control variable for the outer loop and iteration
   counter
keep_going = 1;
if verbose
    fprintf(1,'\nInitial obj=%10.6e, nonzeros=%7d\n',f,
       num_nz_x);
end
temp = AT(resid_base);
term = temp - Aty;
gradu = term + tau; % Hz+c w.r.t. u
gradv = -term + tau; % Hz+c w.r.t. v
Lu = min(u, gradu); % F(z) = min(z, Hz+c) w.r.t. u, z = [u
    v]'
Lv = min(v, gradv); % F w.r.t. v
du = - Lu;
dv = - Lv;
while keep_going
    % compute dx
    dx = du - dv;
    auv = A(dx);
    Bdu = AT(auv);
    Bdv = -Bdu;
    % initial steplength
    betas = 1;
                % betas = 10 for paper;
    old_Lu = Lu;
    old_Lv = Lv;
    %NormF = Lu(:)'*Lu(:) + Lv(:)'*Lv(:);
    %NormFs = sqrt(NormF);
    Lu = min(u+betas*du, gradu+betas*Bdu); % F(z+betas*d)
       w.r.t. u where d=[du;dv];
    Lv = min(v+betas*dv, gradv+betas*Bdv); % F(z+betas*d)
       w.r.t. v;
```

```
Luvduv = Lu(:)'*du(:) + Lv(:)'*dv(:); % F(z+betas*d)
   '*d
dudv = du(:) '*du(:) + dv(:) '*dv(:); % ||d||^2
normFz = sqrt(Lu(:)'*Lu(:)+Lv(:)'*Lv(:));
% - Luvduv < sigma*betas*max(0.001,min(0.8,NormFz))*
  dudv
while - Luvduv < sigma*betas*normFz*dudv</pre>
   % -Luvduv < sigma*betas*dudv
   betas = 0.5*betas;
                      % betas = 0.5*betas;
   Lu = min(u+betas*du,gradu+betas*Bdu);
   Lv = min(v+betas*dv,gradv+betas*Bdv);
   normFz = sqrt(Lu(:)'*Lu(:) + Lv(:)'*Lv(:));
   Luvduv = Lu(:)'*du(:) + Lv(:)'*dv(:);
end
lambda = -1.8*Luvduv*betas/normFz^2;  % lambda = -
  1.6*Luvduv*betas/(Lu(:)'*Lu(:)+Lv(:)'*Lv(:));
old_u = u;
old_v = v;
u = old_u - lambda * Lu;
v = old_v - lambda * Lv;
uvmin = 0; % min(u,v);
u = u - uvmin;
v = v - uvmin;
x = u - v;
% calculate nonzero pattern and number of nonzeros (do
   this *always*)
nz_x_prev = nz_x;
nz_x = (x^-=0.0);
num_nz_x = sum(nz_x(:));
% update residual and function
ALuv = A(Lu-Lv);
prev_f = f;
f = 0.5*(resid(:)'*resid(:)) + sum(tau(:).*u(:)) +
 sum(tau(:).*v(:));
% compute new alpha
dd = Lu(:)'*Lu(:) + Lv(:)'*Lv(:);
% print out stuff
if verbose
   fprintf(1, 'It=%4d, obj=%9.5e, alpha=%6.2e, nz=%8d
    iter, f, alpha, num_nz_x);
end
% update iteration counts, store results and times
```

```
iter = iter + 1;
    objective(iter) = f;
    times(iter) = cputime-t0;
    % compute the next direction
    temp = AT(resid_base);
    term = temp - Aty;
    gradu = term + tau; % Hz+c w.r.t. u
    gradv = -term + tau; % Hz+c w.r.t. v
    %Lu = min(u, gradu); %F(x)
    Lv = min(v, gradv);
    sku = betas*du;
    skv = betas*dv;
    %norms=sqrt(sksk);
    r = 0.01;
    yku = (Lu-old_Lu)+r*sku;
    ykv = (Lv-old_Lv)+r*skv;
    %с
    gam=4;
     skyk=sku(:)'*yku(:)+skv(:)'*ykv(:);
     sksk=sku(:)'*sku(:)+skv(:)'*skv(:);
     Fksk=old_Lu(:)'*sku(:)+old_Lv(:)'*skv(:);
     Fkyk=old_Lu(:)'*yku(:)+old_Lv(:)'*ykv(:);
     dkyk = du(:)'*yku(:)+dv(:)'*ykv(:);
     ykyk=yku(:)'*yku(:)+ykv(:)'*ykv(:);
     bk1 = (Fkyk)/(dkyk);
     bk2=(Fksk)/(dkyk);
     bk3 = (ykyk)/(gam*(skyk));
     bk4=(skyk)/(gam*(sksk));
     tk=bk4+bk3;
        if ((iter > 1))
        du = -1/gam*min(u,gradu)+1/gam*bk1*du-tk*bk2*du;
        dv = -1/gam*min(v,gradv)+1/gam*bk1*dv-tk*bk2*dv;
        end
    %end
%end
    if compute_mse
        err = true - x;
        mses(iter) = (err(:)'*err(:));
    end
    switch stopCriterion
```

```
case 0,
    % compute the stopping criterion based on the
   % of the number of non-zero components of the
       estimate
   num_changes_active = (sum(nz_x(:)~=nz_x_prev
       (:)));
    if num_nz_x >= 1
        criterionActiveSet = num_changes_active;
    else
        criterionActiveSet = tolA / 2;
    end
    keep_going = (criterionActiveSet > tolA);
    if verbose
        fprintf(1, 'Delta n-zeros = %d (target = %e
           )\n',...
          criterionActiveSet , tolA)
    end
case 1,
   % compute the stopping criterion based on the
      relative
   % variation of the objective function.
    criterionObjective = abs(f-prev_f)/(prev_f);
    keep_going = (criterionObjective > tolA);
    if verbose
        fprintf(1,'Delta obj. = %e (target = %e)\n
          criterionObjective , tolA)
    end
case 2,
    % stopping criterion based on relative norm of
        step taken
    delta_x_criterion = norm(Lu(:)-Lv(:))/norm(x
    keep_going = (delta_x_criterion > tolA);
    if verbose
        fprintf(1,'Norm(delta x)/norm(x) = %e (
           target = %e)\n',...
          delta_x_criterion,tolA)
    end
case 3,
   % compute the "LCP" stopping criterion - again
        based on the previous
   \% iterate. Make it "relative" to the norm of x
```

```
w = [ min(gradu(:), old_u(:)); min(gradv(:),
               old_v(:)) ];
            criterionLCP = norm(w(:), inf);
            criterionLCP = criterionLCP / ...
              max([1.0e-6, norm(old_u(:),inf), norm(old_v
                  (:),inf)]);
            keep_going = (criterionLCP > tolA);
            if verbose
                fprintf(1, 'LCP = \%e (target = \%e) \n',
                   criterionLCP, tolA)
            end
        case 4.
            % continue if not yeat reached target value
               tolA
            keep_going = (f > tolA);
            if verbose
                 fprintf(1,'Objective = %e (target = %e)\n'
                    ,f,tolA)
            end
        case 5,
            % stopping criterion based on relative norm of
                step taken
            delta_x_criterion = sqrt(dd)/sqrt(x(:)'*x(:));
            keep_going = (delta_x_criterion > tolA);
            if verbose
                 fprintf(1,'Norm(delta x)/norm(x) = %e (
                   target = \%e) \n',...
                   delta_x_criterion, tolA)
            end
        otherwise,
            error(['Unknown stopping criterion']);
    end % end of the stopping criteria switch
    % take no less than miniter...
    if iter<=miniter</pre>
        keep_going = 1;
    elseif iter > maxiter %and no more than maxiter
       iterations
            keep_going = 0;
    end
end % end of the main loop of keep_going
% increment continuation loop counter
cont_loop = cont_loop+1;
```

```
end % end of the continuation loop
  %
% Print results
%
  if verbose
  fprintf(1,'\nFinished the main algorithm!\nResults:\n')
  fprintf(1, '||A x - y ||_2^2 = %10.3e\n', resid(:)'*resid(:))
  fprintf(1, ||x||_1 = %10.3e\n', sum(abs(x(:))))
  fprintf(1,'Objective function = %10.3e\n',f);
  nz_x = (x^=0.0); num_nz_x = sum(nz_x(:));
  fprintf(1,'Number of non-zero components = %d\n',num_nz_x);
  fprintf(1, 'CPU time so far = %10.3e\n', times(iter));
  fprintf(1, ' n');
end
% If the 'Debias' option is set to 1, we try to remove the
  bias from the 11
% penalty, by applying CG to the least-squares problem
  obtained by omitting
% the l1 term and fixing the zero coefficients at zero.
% do this only if the reduced linear least-squares problem is
% overdetermined, otherwise we are certainly applying CG to a
  problem with a
% singular Hessian
if (debias & (sum(x(:)^-=0)^-=0))
 if (num_nz_x > length(y(:)))
   if verbose
     fprintf(1, '\n')
     fprintf(1, Debiasing requested, but not performed \n');
     fprintf(1, 'There are too many nonzeros in x\n\n');
     fprintf(1, 'nonzeros in x: %8d, length of y: %8d\n',...
         num_nz_x, length(y(:));
   end
 elseif (num_nz_x==0)
   if verbose
     fprintf(1, '\n')
     fprintf(1, 'Debiasing requested, but not performed\n');
     fprintf(1,'x has no nonzeros\n\n');
```

```
end
else
  if verbose
    fprintf(1, '\n')
    fprintf(1, 'Starting the debiasing phase...\n\n')
  end
 x_{debias} = x;
 zeroind = (x_debias~=0);
  cont_debias_cg = 1;
 debias_start = iter;
 % calculate initial residual
 resid = A(x_debias);
 resid = resid-y;
 resid_prev = eps*ones(size(resid));
 rvec = AT(resid);
 % mask out the zeros
 rvec = rvec .* zeroind;
 rTr_cg = rvec(:)'*rvec(:);
 \% set convergence threshold for the residual |\ |\ RW
     x_{debias} - y ||_2
 tol_debias = tolD * (rvec(:)'*rvec(:));
 % initialize pvec
 pvec = -rvec;
 % main loop
 while cont_debias_cg
    % calculate A*p = Wt * Rt * R * W * pvec
    RWpvec = A(pvec);
    Apvec = AT(RWpvec);
    % mask out the zero terms
    Apvec = Apvec .* zeroind;
    % calculate alpha for CG
    alpha_cg = rTr_cg / (pvec(:)'* Apvec(:));
    % take the step
    x_debias = x_debias + alpha_cg * pvec;
    resid = resid + alpha_cg * RWpvec;
```

```
rvec = rvec + alpha_cg * Apvec;
    rTr_cg_plus = rvec(:)'*rvec(:);
    beta_cg = rTr_cg_plus / rTr_cg;
    pvec = -rvec + beta_cg * pvec;
    rTr_cg = rTr_cg_plus;
    iter = iter+1;
    objective(iter) = 0.5*(resid(:)'*resid(:)) + ...
        sum(tau(:).*abs(x_debias(:)));
    times(iter) = cputime - t0;
    if compute_mse
      err = true - x_debias;
      mses(iter) = (err(:)'*err(:));
    end
    \% in the debiasing CG phase, always use convergence
       criterion
    % based on the residual (this is standard for CG)
    if verbose
      fprintf(1,' Iter = %5d, debias resid = %13.8e,
         convergence = %8.3e\n', ...
          iter, resid(:)'*resid(:), rTr_cg / tol_debias);
    end
    cont_debias_cg = ...
        (iter-debias_start <= miniter_debias ) | ...
        ((rTr_cg > tol_debias) & ...
        (iter-debias_start <= maxiter_debias));</pre>
  end
  if verbose
    fprintf(1,'\nFinished the debiasing phase!\nResults:\n')
    fprintf(1,'||A x - y ||_2^2 = 10.3e\n',resid(:)'*resid
       (:))
    fprintf(1, ||x||_1 = %10.3e\n', sum(abs(x(:))))
    fprintf(1,'Objective function = %10.3e\n',f);
    nz = (x_debias^{=0.0});
    fprintf(1, 'Number of non-zero components = %d\n', sum(nz
    fprintf(1, 'CPU time so far = %10.3e\n', times(iter));
    fprintf(1, '\n');
  end
end
```

```
mses = mses/length(true(:));
  end
end
(2) Code for Algorithm 2.2 Image de-blurring
function [x,x_debias,objective,times,debias_start,mses,taus]=
   DKcluster(y, A, tau, varargin)
%
% HTTCGP_CS 1.0, Nov. 29, 2019
% This function solves the convex problem
% arg min_x = 0.5*|| y - A x ||_2^2 + tau || x ||_1
% using the algorithm modified three-term conjugate gradient
  method, described in the following paper
%
\% This code is to use the well-known code CG_DESCENT to solve
  \ell_1 norm
% regularization least square problems.
%
    ._____
% Copyright (2019): Jianghua Yin
%
%
% The first version of this code by Jianghua Yin, Nov. 29,
  2019
% test for number of required parametres
if (nargin-length(varargin)) ~= 3
  error('Wrong number of required parameters');
end
% flag for initial x (can take any values except 0,1,2)
Initial_X_supplied = 3333;
% Set the defaults for the optional parameters
stopCriterion = 3;
tolA = 0.01;
tolD = 0.0001;
```

if compute_mse

```
debias = 0;
maxiter = 10000;
maxiter_debias = 500;
miniter = 5;
miniter_debias = 5;
init = 0;
compute_mse = 0;
AT = 0;
verbose = 1;
continuation = 0;
cont_steps = -1;
firstTauFactorGiven = 0;
% Set the defaults for outputs that may not be computed
debias_start = 0;
x_{debias} = [];
mses = [];
% Read the optional parameters
if (rem(length(varargin),2)==1)
 error('Optional parameters should always go by pairs');
else
 for i=1:2:(length(varargin)-1)
   switch upper(varargin{i})
    case 'STOPCRITERION'
      stopCriterion = varargin{i+1};
    case 'TOLERANCEA'
      tolA = varargin{i+1};
    case 'TOLERANCED'
      tolD = varargin{i+1};
    case 'DEBIAS'
      debias = varargin{i+1};
    case 'MAXITERA'
      maxiter = varargin{i+1};
    case 'MAXITERD'
      maxiter_debias = varargin{i+1};
    case 'MINITERA'
      miniter = varargin{i+1};
    case 'MINITERD'
      miniter_debias = varargin{i+1};
    case 'INITIALIZATION'
      as array
             used below
             x = varargin{i+1};
```

```
else
              init = varargin{i+1};
       end
     case 'MONOTONE'
       enforceMonotone = varargin{i+1};
     case 'CONTINUATION'
       continuation = varargin{i+1};
     case 'CONTINUATIONSTEPS'
       cont_steps = varargin{i+1};
     case 'FIRSTTAUFACTOR'
       firstTauFactor = varargin{i+1};
       firstTauFactorGiven = 1;
     case 'TRUE_X'
       compute_mse = 1;
       true = varargin{i+1};
     case 'ALPHAMIN'
       alphamin = varargin{i+1};
     case 'ALPHAMAX'
       alphamax = varargin{i+1};
     case 'AT'
       AT = varargin{i+1};
     case 'VERBOSE'
       verbose = varargin{i+1};
     otherwise
      % Hmmm, something wrong with the parameter string
      error(['Unrecognized option: ''' varargin{i} '''']);
    end;
  end;
end
%%%%%%%%%%%%%%%%%%
if (sum(stopCriterion == [0 1 2 3 4 5]) == 0)
  error(['Unknown stopping criterion']);
end
% if A is a function handle, we have to check presence of AT,
if isa(A, 'function_handle') & ~isa(AT, 'function_handle')
  error(['The function handle for transpose of A is missing'])
end
% if A is a matrix, we find out dimensions of y and x,
\% and create function handles for multiplication by A and A',
% so that the code below doesn't have to distinguish between
% the handle/not-handle cases
if ~isa(A, 'function_handle')
```

```
AT = Q(x) (x'*A)'; %A'*x;
 A = O(x) A*x;
\% from this point down, A and AT are always function handles.
% Precompute A'*y since it'll be used a lot
Aty = AT(y);
% Initialization
switch init
    case 0 % initialize at zero, using AT to find the size
       of x
       x = AT(zeros(size(y)));
    case 1 % initialize randomly, using AT to find the size
       of x
       x = randn(size(AT(zeros(size(y)))));
    case 2
             % initialize x0 = A'*y
       x = Aty;
    case Initial_X_supplied % initial x was given by user
       % initial x was given as a function argument; just
          check size
       if size(A(x)) ~= size(y)
          error(['Size of initial x is not compatible with A'
             ]);
       end
    otherwise
       error(['Unknown ''Initialization'' option']);
end
% now check if tau is an array; if it is, it has to
\% have the same size as x
if prod(size(tau)) > 1
   try,
      dummy = x.*tau;
   catch,
      error(['Parameter tau has wrong dimensions; it should be
          scalar or size(x)']),
   end
end
% if the true x was given, check its size
if compute_mse & (size(true) ~= size(x))
   error(['Initial x has incompatible size']);
end
```

```
% if tau is scalar, we check its value; if it's large enough,
\% the optimal solution is the zero vector
if prod(size(tau)) == 1
   aux = AT(y);
   max_tau = max(abs(aux(:)));
   if tau >= max_tau
      x = zeros(size(aux));
      if debias
         x_debias = x;
      end
      objective(1) = 0.5*(y(:)'*y(:));
      times(1) = 0;
      if compute_mse
          mses(1) = sum(true(:).^2);
      end
      return
                         %
   end
end
% initialize u and v
u = x.*(x >= 0);
v = -x.*(x < 0);
\% define the indicator vector or matrix of nonzeros in x
nz_x = (x = 0.0);
num_nz_x = sum(nz_x(:));
% start the clock
t0 = cputime;
% store given tau, because we're going to change it in the
% continuation procedure
final_tau = tau;
% store given stopping criterion and threshold, because we're
   going
% to change them in the continuation procedure
final_stopCriterion = stopCriterion;
final_tolA = tolA;
% set continuation factors
if continuation&&(cont_steps > 1)
   % If tau is scalar, first check top see if the first factor
       is
   % too large (i.e., large enough to make the first
```

```
% solution all zeros). If so, make it a little smaller than
       that.
   \% Also set to that value as default
   if prod(size(tau)) == 1
      if (firstTauFactorGiven == 0)|(firstTauFactor*tau >=
         max_tau)
         firstTauFactor = 0.5*max_tau / tau;
         if verbose
             fprintf(1,'\n setting parameter FirstTauFactor\n'
                )
         end
      end
   end
   cont_factors = 10.^[log10(firstTauFactor):...
                    log10(1/firstTauFactor)/(cont_steps-1):0];
end
if ~continuation
  cont_factors = 1;
  cont_steps = 1;
end
iter = 1;
if compute_mse
       mses(iter) = sum((x(:)-true(:)).^2);
end
keep_continuation = 1;
cont_loop = 1;
iter = 1;
taus = [];
sigma = 0.001;
% loop for continuation
while keep_continuation
    % Compute and store initial value of the objective
       function
    resid = y - A(x);
    if cont_steps == -1
        gradq = AT(resid);
        tau = max(final_tau, 0.2*max(abs(gradq)));
        if tau == final_tau
            stopCriterion = final_stopCriterion;
            tolA = final_tolA;
            keep_continuation = 0;
                                                      % stop
               continuation
```

```
else
        stopCriterion = 1;
        tolA = 1e-5;
    end
else
    tau = final_tau * cont_factors(cont_loop); %
    if cont_loop == cont_steps
        stopCriterion = final_stopCriterion;
        tolA = final_tolA;
                                                 %
        keep_continuation = 0;
    else
        stopCriterion = 1;
        tolA = 1e-5;
    end
end
taus = [taus tau];
if verbose
    fprintf(1, '\nSetting tau = %0.5g\n', tau)
\% if in first continuation iteration, compute and store
% initial value of the objective function
if cont_loop == 1
    alpha = 1.0;
    f = 0.5*(resid(:)'*resid(:)) + ...
         sum(tau(:).*u(:)) + sum(tau(:).*v(:));
    objective(1) = f;
    if compute_mse
        mses(1) = (x(:)-true(:))'*(x(:)-true(:));
    end
    if verbose
        fprintf(1,'Initial obj=%10.6e, alpha=%6.2e,
           nonzeros=%7d\n',...
            f,alpha,num_nz_x);
    end
end
% Compute the initial gradient and the useful
% quantity resid_base
resid_base = y - resid;
% control variable for the outer loop and iteration
   counter
keep_going = 1;
if verbose
    fprintf(1,'\nInitial obj=%10.6e, nonzeros=%7d\n',f,
       num_nz_x);
end
```

```
temp = AT(resid_base);
term = temp - Aty;
gradu = term + tau; % Hz+c w.r.t. u
gradv = -term + tau; % Hz+c w.r.t. v
Lu = min(u, gradu); % F(z) = min(z, Hz+c) w.r.t. u, z = [u
    יןע
Lv = min(v,gradv); % F w.r.t. v
du = - Lu;
dv = - Lv;
while keep_going
    % compute dx
    dx = du - dv;
    auv = A(dx);
    Bdu = AT(auv);
    Bdv = -Bdu;
    % initial steplength
    betas = 1;  % betas = 10 for paper;
    old_Lu = Lu;
    old_Lv = Lv;
    %NormF = Lu(:)'*Lu(:) + Lv(:)'*Lv(:);
    %NormFs = sqrt(NormF);
    Lu = min(u+betas*du, gradu+betas*Bdu); % F(z+betas*d)
       w.r.t. u where d=[du;dv];
    Lv = min(v+betas*dv, gradv+betas*Bdv); % F(z+betas*d)
       w.r.t. v;
    Luvduv = Lu(:)'*du(:) + Lv(:)'*dv(:); % F(z+betas*d)
       '*d
    dudv = du(:)'*du(:)+dv(:)'*dv(:);
                                            % ||d||^2
    normFz = sqrt(Lu(:)'*Lu(:)+Lv(:)'*Lv(:));
    % - Luvduv < sigma*betas*max(0.001,min(0.8,NormFz))*
       dudv
    while - Luvduv < sigma*betas*normFz*dudv</pre>
        % -Luvduv < sigma*betas*dudv
        betas = 0.9*betas;
                            \% betas = 0.5*betas;
        Lu = min(u+betas*du,gradu+betas*Bdu);
        Lv = min(v+betas*dv,gradv+betas*Bdv);
        normFz = sqrt(Lu(:)'*Lu(:) + Lv(:)'*Lv(:));
        Luvduv = Lu(:)'*du(:) + Lv(:)'*dv(:);
    end
    lambda = -1.8*Luvduv*betas/normFz^2;  % lambda = -
       1.6*Luvduv*betas/(Lu(:)'*Lu(:)+Lv(:)'*Lv(:));
    old_u = u;
    old_v = v;
    u = old_u - lambda * Lu;
    v = old_v - lambda * Lv;
    uvmin = 0; % min(u,v);
```

```
u = u - uvmin;
v = v - uvmin;
x = u - v;
\% calculate nonzero pattern and number of nonzeros (do
   this *always*)
nz_x_prev = nz_x;
nz_x = (x^-=0.0);
num_nz_x = sum(nz_x(:));
% update residual and function
ALuv = A(Lu-Lv);
prev_f = f;
f = 0.5*(resid(:)'*resid(:)) + sum(tau(:).*u(:)) +
 sum(tau(:).*v(:));
% compute new alpha
dd = Lu(:)'*Lu(:) + Lv(:)'*Lv(:);
% print out stuff
if verbose
   fprintf(1, 'It=%4d, obj=%9.5e, alpha=%6.2e, nz=%8d
       ٠,...
    iter, f, alpha, num_nz_x);
end
% update iteration counts, store results and times
iter = iter + 1;
objective(iter) = f;
times(iter) = cputime-t0;
% compute the next direction
temp = AT(resid_base);
term = temp - Aty;
gradu = term + tau; % Hz+c w.r.t. u
gradv = -term + tau; % Hz+c w.r.t. v
%Lu = min(u, gradu); % F(x)
%Lv = min(v,gradv);
sku = betas*du;
skv = betas*dv;
%
r = 0.01;
yku = (Lu-old_Lu)+r*sku;
ykv = (Lv-old_Lv)+r*skv;
%с
%
gam = 0.25;
skyk=sku(:)'*yku(:)+skv(:)'*ykv(:);
sksk=sku(:)'*sku(:)+skv(:)'*skv(:);
```

```
Fksk=old_Lu(:)'*sku(:)+old_Lv(:)'*skv(:);
     Fkyk=old_Lu(:)'*yku(:)+old_Lv(:)'*ykv(:);
     dkyk=du(:)'*yku(:)+dv(:)'*ykv(:);
     ykyk=yku(:)'*yku(:)+ykv(:)'*ykv(:);
     bk1 = (Fkyk)/(dkyk);
     bk2=(Fksk)/(dkyk);
     bk3=gam*(ykyk)/((skyk));
     bk4=gam*(skyk)/((sksk));
     tauk=bk4+bk3;
        if ((iter > 1))
         du = -gam*min(u,gradu)+gam*bk1*du-tauk*(bk2)*du;
         dv = -gam*min(v,gradv)+gam*bk1*dv-tauk*(bk2)*dv;
        end
    %end
%end
    if compute_mse
        err = true - x;
        mses(iter) = (err(:)'*err(:));
    end
    switch stopCriterion
        case 0,
            % compute the stopping criterion based on the
               change
            % of the number of non-zero components of the
               estimate
            num_changes_active = (sum(nz_x(:)~=nz_x_prev
               (:)));
            if num_nz_x >= 1
                criterionActiveSet = num_changes_active;
            else
                criterionActiveSet = tolA / 2;
            end
            keep_going = (criterionActiveSet > tolA);
            if verbose
                fprintf(1,'Delta n-zeros = %d (target = %e
                   )\n',...
                  criterionActiveSet , tolA)
            end
        case 1,
            % compute the stopping criterion based on the
               relative
            % variation of the objective function.
```

```
criterionObjective = abs(f-prev_f)/(prev_f);
    keep_going = (criterionObjective > tolA);
    if verbose
        fprintf(1,'Delta obj. = %e (target = %e)\n
          criterionObjective, tolA)
    end
case 2,
   % stopping criterion based on relative norm of
        step taken
    delta_x_criterion = norm(Lu(:)-Lv(:))/norm(x
    keep_going = (delta_x_criterion > tolA);
    if verbose
        fprintf(1, 'Norm(delta x)/norm(x) = %e (
           target = %e) \n',...
          delta_x_criterion, tolA)
    end
case 3,
   % compute the "LCP" stopping criterion - again
        based on the previous
   \% iterate. Make it "relative" to the norm of x
   w = [ min(gradu(:), old_u(:)); min(gradv(:),
       old_v(:)) ];
    criterionLCP = norm(w(:), inf);
    criterionLCP = criterionLCP / ...
      max([1.0e-6, norm(old_u(:),inf), norm(old_v
         (:),inf)]);
    keep_going = (criterionLCP > tolA);
    if verbose
        fprintf(1, 'LCP = \%e (target = \%e) \n',
           criterionLCP, tolA)
    end
case 4,
   % continue if not year reached target value
       tolA
   keep_going = (f > tolA);
    if verbose
        fprintf(1,'Objective = %e (target = %e)\n'
           ,f,tolA)
    end
case 5,
    % stopping criterion based on relative norm of
        step taken
    delta_x_criterion = sqrt(dd)/sqrt(x(:)'*x(:));
```

```
keep_going = (delta_x_criterion > tolA);
              if verbose
                  fprintf(1,'Norm(delta x)/norm(x) = %e (
                    target = %e)\n',...
                    delta_x_criterion, tolA)
              end
          otherwise,
              error(['Unknown stopping criterion']);
       end % end of the stopping criteria switch
       % take no less than miniter...
       if iter<=miniter</pre>
          keep_going = 1;
       elseif iter > maxiter %and no more than maxiter
          iterations
              keep_going = 0;
       end
   end % end of the main loop of keep_going
   % increment continuation loop counter
   cont_loop = cont_loop+1;
end % end of the continuation loop
%
  % Print results
  if verbose
  fprintf(1,'\nFinished the main algorithm!\nResults:\n')
  fprintf(1, | | A x - y | |_2^2 = %10.3e |_r, resid(:) | *resid(:) |
  fprintf(1, ||x||_1 = %10.3e\n', sum(abs(x(:))))
  fprintf(1,'Objective function = %10.3e\n',f);
  nz_x = (x^-=0.0); num_nz_x = sum(nz_x(:));
  fprintf(1, 'Number of non-zero components = %d\n', num_nz_x);
  fprintf(1, 'CPU time so far = %10.3e\n', times(iter));
  fprintf(1, '\n');
end
% If the 'Debias' option is set to 1, we try to remove the
  bias from the 11
```

```
% penalty, by applying CG to the least-squares problem
   obtained by omitting
% the l1 term and fixing the zero coefficients at zero.
% do this only if the reduced linear least-squares problem is
% overdetermined, otherwise we are certainly applying CG to a
   problem with a
% singular Hessian
if (debias & (sum(x(:)^-=0)^-=0))
  if (num_nz_x > length(y(:)))
    if verbose
      fprintf(1, ' n')
      fprintf(1, 'Debiasing requested, but not performed\n');
      fprintf(1, 'There are too many nonzeros in x \in x);
      fprintf(1, 'nonzeros in x: %8d, length of y: %8d\n',...
          num_nz_x, length(y(:)));
  elseif (num_nz_x==0)
    if verbose
      fprintf(1, '\n')
      fprintf(1, 'Debiasing requested, but not performed\n');
      fprintf(1,'x has no nonzeros\n\n');
    end
  else
    if verbose
      fprintf(1, '\n')
      fprintf(1,'Starting the debiasing phase...\n\n')
    end
    x_{debias} = x;
    zeroind = (x_debias~=0);
    cont_debias_cg = 1;
    debias_start = iter;
    % calculate initial residual
    resid = A(x_debias);
    resid = resid-y;
    resid_prev = eps*ones(size(resid));
    rvec = AT(resid);
    % mask out the zeros
    rvec = rvec .* zeroind;
    rTr_cg = rvec(:)'*rvec(:);
```

```
% set convergence threshold for the residual || RW
   x_{debias} - y ||_2
tol_debias = tolD * (rvec(:)'*rvec(:));
% initialize pvec
pvec = -rvec;
% main loop
while cont_debias_cg
  % calculate A*p = Wt * Rt * R * W * pvec
  RWpvec = A(pvec);
  Apvec = AT(RWpvec);
  % mask out the zero terms
  Apvec = Apvec .* zeroind;
  % calculate alpha for CG
  alpha_cg = rTr_cg / (pvec(:)'* Apvec(:));
  % take the step
  x_debias = x_debias + alpha_cg * pvec;
  resid = resid + alpha_cg * RWpvec;
  rvec = rvec + alpha_cg * Apvec;
  rTr_cg_plus = rvec(:)'*rvec(:);
  beta_cg = rTr_cg_plus / rTr_cg;
  pvec = -rvec + beta_cg * pvec;
  rTr_cg = rTr_cg_plus;
  iter = iter+1;
  objective(iter) = 0.5*(resid(:)'*resid(:)) + ...
      sum(tau(:).*abs(x_debias(:)));
  times(iter) = cputime - t0;
  if compute_mse
    err = true - x_debias;
    mses(iter) = (err(:)'*err(:));
  end
  % in the debiasing CG phase, always use convergence
     criterion
  % based on the residual (this is standard for CG)
```

```
if verbose
        fprintf(1,' Iter = %5d, debias resid = %13.8e,
           convergence = \%8.3e\n', ...
            iter, resid(:)'*resid(:), rTr_cg / tol_debias);
      end
      cont_debias_cg = ...
          (iter-debias_start <= miniter_debias ) | ...
          ((rTr_cg > tol_debias) & ...
          (iter-debias_start <= maxiter_debias));</pre>
    end
    if verbose
      fprintf(1,'\nFinished the debiasing phase!\nResults:\n')
      fprintf(1,'||A x - y ||_2^2 = 10.3e\n',resid(:)'*resid
         (:))
      fprintf(1, ||x||_1 = 10.3e\n', sum(abs(x(:)))
      fprintf(1,'Objective function = %10.3e\n',f);
      nz = (x_debias^{=0.0});
      fprintf(1,'Number of non-zero components = %d\n',sum(nz
      fprintf(1, 'CPU time so far = %10.3e\n', times(iter));
      fprintf(1, '\n');
    end
  end
  if compute_mse
    mses = mses/length(true(:));
  end
end
(3) Code for NHZ1 signal reconstruction
function [x,x_debias,objective,times,debias_start,mses,taus]=
     NHZ1(y,A,tau,varargin)
%
\% CGD_CS version 1.0, December 3, 2009
% This function solves the convex problem
\% arg min_x = 0.5*|| y - A x ||_2^2 + tau || x ||_1
% using the algorithm modified PRP conjugate gradient method,
   described in the following paper
% This code is to use the well-known code CG_DESCENT to solve
  \ell_1 norm
% regularization least square problems.
```

```
%
%
% Copyright (2010): Yunhai Xiao and Hong Zhu
%
\% The first version of this code by Yunhai Xiao, Oct. 15. 2010
% test for number of required parametres
if (nargin-length(varargin)) ~= 3
  error('Wrong number of required parameters');
end
% flag for initial x (can take any values except 0,1,2)
Initial_X_supplied = 3333;
% Set the defaults for the optional parameters
stopCriterion = 3;
tolA = 0.01;
tolD = 0.0001;
debias = 0;
maxiter = 10000;
maxiter_debias = 500;
miniter = 5;
miniter_debias = 5;
init = 0;
compute_mse = 0;
AT = 0;
verbose = 1;
continuation = 0;
cont_steps = -1;
firstTauFactorGiven = 0;
% Set the defaults for outputs that may not be computed
debias_start = 0;
x_{debias} = [];
mses = [];
% Read the optional parameters
if (rem(length(varargin),2)==1)
  error('Optional parameters should always go by pairs');
else
```

```
for i=1:2:(length(varargin)-1)
  switch upper(varargin{i})
  case 'STOPCRITERION'
     stopCriterion = varargin{i+1};
  case 'TOLERANCEA'
    tolA = varargin{i+1};
  case 'TOLERANCED'
    tolD = varargin{i+1};
  case 'DEBIAS'
     debias = varargin{i+1};
  case 'MAXITERA'
    maxiter = varargin{i+1};
  case 'MAXITERD'
     maxiter_debias = varargin{i+1};
  case 'MINITERA'
    miniter = varargin{i+1};
  case 'MINITERD'
     miniter_debias = varargin{i+1};
  case 'INITIALIZATION'
     if prod(size(varargin{i+1})) > 1  % initial x supplied
        as array
            used below
           x = varargin{i+1};
     else
           init = varargin{i+1};
     end
  case 'MONOTONE'
     enforceMonotone = varargin{i+1};
  case 'CONTINUATION'
     continuation = varargin{i+1};
  case 'CONTINUATIONSTEPS'
     cont_steps = varargin{i+1};
  case 'FIRSTTAUFACTOR'
     firstTauFactor = varargin{i+1};
    firstTauFactorGiven = 1;
  case 'TRUE_X'
     compute_mse = 1;
    true = varargin{i+1};
  case 'ALPHAMIN'
     alphamin = varargin{i+1};
  case 'ALPHAMAX'
     alphamax = varargin{i+1};
  case 'AT'
    AT = varargin{i+1};
  case 'VERBOSE'
```

```
verbose = varargin{i+1};
     otherwise
      \% Hmmm, something wrong with the parameter string
      error(['Unrecognized option: ''' varargin{i} '''']);
    end;
  end;
end
if (sum(stopCriterion == [0 1 2 3 4 5]) == 0)
  error(['Unknown stopping criterion']);
end
% if A is a function handle, we have to check presence of AT,
if isa(A, 'function_handle') & ~isa(AT, 'function_handle')
  error(['The function handle for transpose of A is missing'])
end
% if A is a matrix, we find out dimensions of y and x,
% and create function handles for multiplication by A and A',
% so that the code below doesn't have to distinguish between
% the handle/not-handle cases
if ~isa(A, 'function_handle')
 AT = Q(x) (x'*A)'; %A'*x;
 A = 0(x) A*x;
\% from this point down, A and AT are always function handles.
% Precompute A'*y since it'll be used a lot
Aty = AT(y);
% Initialization
switch init
    case 0 % initialize at zero, using AT to find the size
      x = AT(zeros(size(y)));
    case 1 % initialize randomly, using AT to find the size
      x = randn(size(AT(zeros(size(y)))));
    case 2
            % initialize x0 = A'*y
       x = Aty;
    case Initial_X_supplied % initial x was given by user
      % initial x was given as a function argument; just
          check size
       if size(A(x)) ~= size(y)
```

```
error(['Size of initial x is not compatible with A'
             ]);
       end
    otherwise
       error(['Unknown ''Initialization'' option']);
end
% now check if tau is an array; if it is, it has to
\% have the same size as x
if prod(size(tau)) > 1
   try,
      dummy = x.*tau;
   catch,
      error(['Parameter tau has wrong dimensions; it should be
          scalar or size(x)']),
   end
end
% if the true x was given, check its size
if compute_mse & (size(true) ~= size(x))
   error(['Initial x has incompatible size']);
end
% if tau is scalar, we check its value; if it's large enough,
\% the optimal solution is the zero vector
if prod(size(tau)) == 1
   aux = AT(y);
   max_tau = max(abs(aux(:)));
   if tau >= max_tau
      x = zeros(size(aux));
      if debias
         x_{debias} = x;
      objective(1) = 0.5*(y(:)'*y(:));
      times(1) = 0;
      if compute_mse
          mses(1) = sum(true(:).^2);
      end
      return
   end
end
\% initialize u and v
u = x.*(x >= 0);
v = -x.*(x < 0);
```

```
\% define the indicator vector or matrix of nonzeros in x
nz_x = (x = 0.0);
num_nz_x = sum(nz_x(:));
% start the clock
t0 = cputime;
% store given tau, because we're going to change it in the
% continuation procedure
final_tau = tau;
% store given stopping criterion and threshold, because we're
% to change them in the continuation procedure
final_stopCriterion = stopCriterion;
final_tolA = tolA;
% set continuation factors
if continuation&&(cont_steps > 1)
   % If tau is scalar, first check top see if the first factor
   % too large (i.e., large enough to make the first
   % solution all zeros). If so, make it a little smaller than
      that.
   % Also set to that value as default
   if prod(size(tau)) == 1
      if (firstTauFactorGiven == 0)|(firstTauFactor*tau >=
         max_tau)
         firstTauFactor = 0.5*max_tau / tau;
         if verbose
             fprintf(1,'\n setting parameter FirstTauFactor\n'
                )
         end
      end
   cont_factors = 10.^[log10(firstTauFactor):...
                    log10(1/firstTauFactor)/(cont_steps-1):0];
end
if ~continuation
  cont_factors = 1;
  cont_steps = 1;
end
iter = 1;
```

```
if compute_mse
       mses(iter) = sum((x(:)-true(:)).^2);
end
keep_continuation = 1;
cont_loop = 1;
iter = 1;
taus = [];
% loop for continuation
while keep_continuation
    % Compute and store initial value of the objective
       function
    resid = y - A(x);
    if cont_steps == -1
        gradq = AT(resid);
        tau = max(final_tau, 0.2*max(abs(gradq)));
        if tau == final_tau
            stopCriterion = final_stopCriterion;
            tolA = final_tolA;
                                                      % stop
            keep_continuation = 0;
               continuation
        else
            stopCriterion = 1;
            tolA = 1e-5;
        end
    else
        tau = final_tau * cont_factors(cont_loop); %
        if cont_loop == cont_steps
            stopCriterion = final_stopCriterion;
            tolA = final_tolA;
                                                     %
            keep_continuation = 0;
        else
            stopCriterion = 1;
            tolA = 1e-5;
        end
    end
    taus = [taus tau];
    if verbose
        fprintf(1, '\nSetting tau = \%0.5g\n', tau)
    end
    \% if in first continuation iteration, compute and store
    % initial value of the objective function
```

```
if cont_loop == 1
    alpha = 1.0;
    f = 0.5*(resid(:)'*resid(:)) + ...
         sum(tau(:).*u(:)) + sum(tau(:).*v(:));
    objective(1) = f;
    if compute_mse
        mses(1) = (x(:)-true(:))'*(x(:)-true(:));
    end
    if verbose
        fprintf(1, 'Initial obj=%10.6e, alpha=%6.2e,
           nonzeros=%7d\n',...
            f,alpha,num_nz_x);
    end
end
% Compute the initial gradient and the useful
% quantity resid_base
resid_base = y - resid;
% control variable for the outer loop and iteration
   counter
keep_going = 1;
if verbose
  fprintf(1,'\nInitial obj=%10.6e, nonzeros=%7d\n',f,
     num_nz_x);
end
while keep_going
  % compute gradient
  temp = AT(resid_base);
  term = temp - Aty;
  gradu = term + tau;
  gradv = -term + tau;
  %
  Lu = min(u,gradu); %1A
  Lv = min(v,gradv); %1A
  %NormF = Lu(:)'*Lu(:) + Lv(:)'*Lv(:);
  %NormF2 = sqrt(NormF);
  if (iter > 1)
     %(esku, eskv)^T=x_k-x_{k-1}=s_{k-1}
     su = u - old_u;
     sv = v - old_v;
     %Fkdk = F_k^d = \{k-1\}
```

```
Fkdk= Lu(:) '*old_du(:)+Lv(:) '*old_dv(:);
   %sksk = ||s_{k-1}||^2
   sksk = su(:)'*su(:)+sv(:)'*sv(:);
   Normsk=sqrt(sksk);
   m = 0.01;
   %(rku, rkv)^T = F_k - F_{k-1} = y_{k-1}
   rku = min(u,gradu)-min(old_u,old_gradu);
   rkv = min(v,gradv)-min(old_v,old_gradv);
   wk1=rku+m*su;
   wk2=rkv+m*sv;
   %skyb=s_{k-1}^Tyb
   %skyb = su(:)'*wk1(:)+sv(:)'*wk2(:);
   %tk=1+max(0,-(skyb/sksk));
   %wk = yb + tk * sk
   %wk1=yb1+tk*su;
   %wk2=yb2+tk*sv;
   %skwk=s_{k-1}^Tw_{k-1}
   skwk=su(:)'*wk1(:)+sv(:)'*wk2(:);
   dkwk = d_{k-1}^Tw_{k-1}
   dkwk=old_du(:)'*wk1(:)+old_dv(:)'*wk2(:);
   %Fkwk=F_{k-1}^Tw_{k-1}
   Fkwk=Lu(:)'*wk1(:)+Lv(:)'*wk2(:);
   % | | wk | |^2 = (yb + tk * sk)^2
   wkwk=wk1(:)'*wk1(:)+wk2(:)'*wk2(:);
   % | | wk | | = sqrt (wkwk)
   Normwk=sqrt(wkwk);
   gamma=4;
   ts=(1/gamma)*(sqrt(skwk/(Normsk*Normwk)))^3;
   L=(2/(4*gamma));
   tsb=max(ts,L);
   %Betak
   Betak=(Fkwk)/(dkwk)-gamma*tsb*((wkwk)*(Fkdk)/(dkwk)
      ^2);
end
old_gradu = gradu;
old_gradv = gradv;
% computation of search direction vector
du = - min(u, gradu);
dv = - min(v, gradv);
if (iter > 1)
   du = -min(u, gradu) +Betak*old_du;
   dv = -min(v, gradv) +Betak*old_dv;
end
dx = du - dv;
```

```
old_u = u;
old_v = v;
old_du = du;
old_dv = dv;
%Old_NormF = NormF;
% calculate useful matrix-vector product involving dx
auv = A(dx);
Bdu = AT(auv);
Bdv = -Bdu;
% preparetion for line search
sigma = 0.0001;
betas = 1;
Lu = min(u+betas*du, gradu+betas*Bdu);
Lv = min(v+betas*dv, gradv+betas*Bdv);
Luvduv = Lu(:)'*du(:) + Lv(:)'*dv(:);
dudv = du(:)'*du(:)+dv(:)'*dv(:);
% line search process
while - Luvduv < sigma*betas*dudv</pre>
    betas = 0.51*betas;
    Lu = min(u+betas*du,gradu+betas*Bdu);
    Lv = min(v+betas*dv,gradv+betas*Bdv);
    Luvduv = Lu(:)'*du(:) + Lv(:)'*dv(:);
end
% compute the projection steplength
lambda = - 1.8*Luvduv*betas/(Lu(:)'*Lu(:)+Lv(:)'*Lv(:));
    % lamda_k=-(alpha*d_k^T*F(z_k))/(norm(z_k))^2
%
u = old_u - lambda * Lu;
v = old_v - lambda * Lv;
uvmin = 0; % min(u,v);
u = u - uvmin;
v = v - uvmin;
x = u - v;
% calculate nonzero pattern and number of nonzeros (do
  this *always*)
nz_x_prev = nz_x;
nz_x = (x^-=0.0);
num_nz_x = sum(nz_x(:));
% update residual and function
ALuv = A(Lu-Lv);
resid = y - resid_base + lambda*ALuv;
prev_f = f;
f = 0.5*(resid(:)'*resid(:)) + sum(tau(:).*u(:)) + ...
    sum(tau(:).*v(:));
% compute new alpha
dd = Lu(:)'*Lu(:) + Lv(:)'*Lv(:);
```

```
%
resid_base = resid_base - lambda*ALuv;
% print out stuff
if verbose
   fprintf(1, 'It=%4d, obj=%9.5e, alpha=%6.2e, nz=%8d ',
       iter, f, alpha, num_nz_x);
end
% update iteration counts, store results and times
iter = iter + 1;
objective(iter) = f;
times(iter) = cputime-t0;
if compute_mse
  err = true - x;
 mses(iter) = (err(:)'*err(:));
end
switch stopCriterion
    case 0,
        % compute the stopping criterion based on the
           change
        \% of the number of non-zero components of the
           estimate
        num_changes_active = (sum(nz_x(:)~=nz_x_prev(:))
        if num_nz_x >= 1
            criterionActiveSet = num_changes_active;
        else
            criterionActiveSet = tolA / 2;
        end
        keep_going = (criterionActiveSet > tolA);
        if verbose
            fprintf(1, 'Delta n-zeros = %d (target = %e)\
               n',...
                criterionActiveSet , tolA)
        end
    case 1,
        % compute the stopping criterion based on the
           relative
        % variation of the objective function.
        criterionObjective = abs(f-prev_f)/(prev_f);
        keep_going = (criterionObjective > tolA);
        if verbose
            fprintf(1, 'Delta obj. = %e (target = %e)\n',
```

```
criterionObjective , tolA)
    end
case 2,
    % stopping criterion based on relative norm of
       step taken
    delta_x_criterion = norm(Lu(:)-Lv(:))/norm(x(:))
    keep_going = (delta_x_criterion > tolA);
    if verbose
        fprintf(1,'Norm(delta x)/norm(x) = %e (
           target = %e) \n',...
            delta_x_criterion, tolA)
    end
case 3,
    % compute the "LCP" stopping criterion - again
       based on the previous
    \% iterate. Make it "relative" to the norm of x.
    w = [\min(gradu(:), old_u(:)); \min(gradv(:),
       old_v(:)) ];
    criterionLCP = norm(w(:), inf);
    criterionLCP = criterionLCP / ...
        max([1.0e-6, norm(old_u(:),inf), norm(old_v
           (:),inf)]);
    keep_going = (criterionLCP > tolA);
    if verbose
        fprintf(1, 'LCP = \%e (target = \%e) \n',
           criterionLCP,tolA)
    end
case 4,
    % continue if not year reached target value tolA
    keep_going = (f > tolA);
    if verbose
        fprintf(1,'Objective = %e (target = %e)\n',f
           ,tolA)
    end
case 5,
 % stopping criterion based on relative norm of
     step taken
  delta_x_criterion = sqrt(dd)/sqrt(x(:)'*x(:));
  keep_going = (delta_x_criterion > tolA);
  if verbose
      fprintf(1,'Norm(delta x)/norm(x) = %e (target
         = \%e) \n',...
          delta_x_criterion, tolA)
  end
otherwise,
```

```
error(['Unknown stopping criterion']);
     end % end of the stopping criteria switch
     % take no less than miniter...
     if iter<=miniter</pre>
            keep_going = 1;
     elseif iter > maxiter %and no more than maxiter
        iterations
            keep_going = 0;
     end
  end % end of the main loop of the BB-QP algorithm
   % increment continuation loop counter
   cont_loop = cont_loop+1;
end % end of the continuation loop
%
  %
% Print results
  if verbose
  fprintf(1,'\nFinished the main algorithm!\nResults:\n')
  fprintf(1, '|A x - y ||_2^2 = %10.3e\n', resid(:)'*resid(:))
  fprintf(1, ||x||_1 = %10.3e\n', sum(abs(x(:))))
  fprintf(1, 'Objective function = %10.3e\n',f);
  nz_x = (x^-=0.0); num_nz_x = sum(nz_x(:));
  fprintf(1,'Number of non-zero components = %d\n',num_nz_x);
  fprintf(1, 'CPU time so far = %10.3e\n', times(iter));
  fprintf(1, '\n');
end
% If the 'Debias' option is set to 1, we try to remove the
  bias from the 11
% penalty, by applying CG to the least-squares problem
  obtained by omitting
\% the l1 term and fixing the zero coefficients at zero.
\% do this only if the reduced linear least-squares problem is
% overdetermined, otherwise we are certainly applying CG to a
  problem with a
% singular Hessian
```

```
if (debias & (sum(x(:)~=0)~=0))
  if (num_nz_x > length(y(:)))
    if verbose
      fprintf(1, ' n')
      fprintf(1, 'Debiasing requested, but not performed\n');
      fprintf(1, 'There are too many nonzeros in x\n\n');
      fprintf(1, 'nonzeros in x: %8d, length of y: %8d\n',...
          num_nz_x, length(y(:));
  elseif (num_nz_x==0)
    if verbose
      fprintf(1, ' n')
      fprintf(1, 'Debiasing requested, but not performed\n');
      fprintf(1,'x has no nonzeros\n\n');
    end
  else
    if verbose
      fprintf(1, '\n')
      fprintf(1, 'Starting the debiasing phase...\n\n')
    end
    x_{debias} = x;
    zeroind = (x_debias~=0);
    cont_debias_cg = 1;
    debias_start = iter;
    % calculate initial residual
    resid = A(x_debias);
    resid = resid-y;
    resid_prev = eps*ones(size(resid));
    rvec = AT(resid);
    % mask out the zeros
    rvec = rvec .* zeroind;
    rTr_cg = rvec(:)'*rvec(:);
    \% set convergence threshold for the residual || RW
       x_{debias} - y ||_2
    tol_debias = tolD * (rvec(:)'*rvec(:));
    % initialize pvec
    pvec = -rvec;
```

```
% main loop
while cont_debias_cg
  % calculate A*p = Wt * Rt * R * W * pvec
  RWpvec = A(pvec);
  Apvec = AT(RWpvec);
  % mask out the zero terms
  Apvec = Apvec .* zeroind;
  % calculate alpha for CG
  alpha_cg = rTr_cg / (pvec(:)'* Apvec(:));
  % take the step
  x_debias = x_debias + alpha_cg * pvec;
  resid = resid + alpha_cg * RWpvec;
  rvec = rvec + alpha_cg * Apvec;
  rTr_cg_plus = rvec(:)'*rvec(:);
  beta_cg = rTr_cg_plus / rTr_cg;
  pvec = -rvec + beta_cg * pvec;
  rTr_cg = rTr_cg_plus;
  iter = iter+1;
  objective(iter) = 0.5*(resid(:)'*resid(:)) + ...
      sum(tau(:).*abs(x_debias(:)));
  times(iter) = cputime - t0;
  if compute_mse
    err = true - x_debias;
    mses(iter) = (err(:)'*err(:));
  end
  % in the debiasing CG phase, always use convergence
     criterion
  % based on the residual (this is standard for CG)
  if verbose
    fprintf(1,' Iter = %5d, debias resid = %13.8e,
       convergence = \%8.3e\n', ...
        iter, resid(:)'*resid(:), rTr_cg / tol_debias);
  end
  cont_debias_cg = ...
      (iter-debias_start <= miniter_debias ) | ...
      ((rTr_cg > tol_debias) & ...
```

```
(iter-debias_start <= maxiter_debias));</pre>
    end
    if verbose
      fprintf(1,'\nFinished the debiasing phase!\nResults:\n')
      fprintf(1,'||A x - y ||_2^2 = 10.3e\n',resid(:)'*resid
         (:))
      fprintf(1, '||x||_1 = %10.3e\n', sum(abs(x(:))))
      fprintf(1,'Objective function = %10.3e\n',f);
      nz = (x_debias^{=0.0});
      fprintf(1,'Number of non-zero components = %d\n',sum(nz
         (:)));
      fprintf(1, 'CPU time so far = %10.3e\n', times(iter));
      fprintf(1, '\n');
    end
  end
  if compute_mse
    mses = mses/length(true(:));
  end
end
```