This is a note of paper "A Newton-Type Method for Positive-Semidefinite Linear Complementarity Problems", which provides a newton method of solving LCP problem which we meet in Ben. Moll (2017). The file "LCP.m" utilizes a slightly modified newton method, which could accelerate the algorithm, while the main idea is not changed.

LCP problem:

$$Mx + q = y, y^Tx = 0, y, x \ge 0. M$$
 is a $n \times n$ matrix, x, y is $n \times 1$ vector (1)

Notation: $|\cdot|$, norm;

Define a mapping:

$$\mathbf{T}(\mathbf{w}) \coloneqq [(\mathbf{M}\mathbf{x} + \mathbf{q} - \mathbf{y})_{\mathbf{n} \times \mathbf{1}}, \boldsymbol{\phi}(\mathbf{w}_1), \boldsymbol{\phi}(\mathbf{w}_2), \dots, \boldsymbol{\phi}(\mathbf{w}_{\mathbf{n}})]^T$$

Where: $\mathbf{w_i} := (x_i, y_i)$ and the function: $\phi(\mathbf{a}, \mathbf{b}) := \sqrt{a^2 + b^2} - a - b$. We have to note that ϕ equals to zero if and only if $\mathbf{ab} = 0, \mathbf{a}, \mathbf{b} \geq 0$.

 $\label{eq:General idea: w = (w_1; w_2; ...) is the solution of LCP if and only if $|T(w)| = 0$ $\to T(w)_i = 0$. Since, $Mx + q = y \ (n \times 1)$, and $y^Tx = 0 \Leftrightarrow \varphi(a,b) = 0$.}$

As some elements will be zero and $\nabla T(x, y)$ may be singular/non-exist. Therefore, we introduce G(w):

$$\mathbf{G}(\mathbf{w}) = \begin{bmatrix} M & diag(-1) \\ diag(\mu(w_i)) & diag(\nu(w_i)) \end{bmatrix}, \tag{2}$$

where
$$\mu(w_i) = -1 + \frac{x_i}{\sqrt{x_i^2 + y_i^2}}, \nu(w_i) = -1 + \frac{y_i}{\sqrt{x_i^2 + y_i^2}}$$
 (3)

Algorithm:

Step 0. Choose w^0 and $\lambda \in (0,1)$. Set k = 0.

Step 1. If |T(w)| = 0, stop.

Step 2. Determine $\widehat{\mathbf{w}}^k$ as solution of:

$$G(\overline{\mathbf{w}}^{\mathbf{k}})(\mathbf{w} - \overline{\mathbf{w}}^{\mathbf{k}}) + T(\overline{\mathbf{w}}^{\mathbf{k}}) = 0$$

Step 3. Update w.

Algorithm A.

Step 0. Choose $w^0 \in \mathbb{R}^{2n}$ and $\lambda \in (0, 1)$. Set k := 0.

Step 1. If $||T(\bar{w}^k)|| = 0$, then stop.

Step 2. Determine $\hat{w}^k \in R^{2n}$ as solution of $G(\bar{w}^k)(w - \bar{w}^k) + T(\bar{w}^k) = 0$.

Step 3. Find the largest number $t_k \in \{(1/2)^j | j \in N\}$ such that $||T(\bar{w}^k + t_k(\hat{w}^k - \bar{w}^k))|| \le (1 - \lambda t_k) ||T(\bar{w}^k)||.$

Step 4. Set $w^{k+1} := \bar{w}^k + t_k(\hat{w}^k - \bar{w}^k)$ and k := k+1. Go to Step 1.

Reference:

- [1] A. Fischer, A Newton-Type Method for Positive-Semidefinite Linear Complementarity Problems, *Journal of Optimization Theory and Applications*: Vol. 86, No. 3, pp. 585-608, 1995.
- [2] Roos, C. "Linear optimization: Theorems of the alternative." *Ecyclopedia of optimization*. Vol. III. Kluwer, 2001. 181-184.

```
function x = LCP(M,q,l,u,x0,display)
%LCP Solve the Linear Complementarity Problem.
%
% USAGE
%
   x = LCP(M,q) solves the LCP
%
%
           x >= 0
%
      Mx + q >= 0
%
  x'(Mx + q) = 0
%
%
   x = LCP(M,q,l,u) solves the generalized LCP (a.k.a MCP)
%
%
   1 < x < u =>
                   Mx + q = 0
%
       x = u
               =>
                   Mx + q < 0
%
   1 = x
               =>
                   Mx + q > 0
%
%
   x = LCP(M,q,l,u,x0,display) allows the optional initial value 'x0' and
%
   a binary flag 'display' which controls the display of iteration data.
%
%
   Parameters:
%
                Termination criterion. Exit when 0.5*phi(x)'*phi(x) < tol.
   tol
%
                Initial value of Levenberg-Marquardt mu coefficient.
   mu
%
   mu step
            - Coefficient by which mu is multiplied / divided.
%
             - Value below which mu is set to zero (pure Gauss-Newton).
   mu_min
%
   max iter - Maximum number of (succesful) Levenberg-Marquardt steps.
%
             - Tolerance of degenerate complementarity: Dimensions where
   b tol
%
                \max(\min(abs(x-1),abs(u-x)), abs(phi(x))) < b_tol
%
                are clamped to the nearest constraint and removed from
%
                the linear system.
%
% ALGORITHM
   This function implements the semismooth algorithm as described in [1],
%
   with a least-squares minimization of the Fischer-Burmeister function using
   a Levenberg-Marquardt trust-region scheme with mu-control as in [2].
%
%
   [1] A. Fischer, A Newton-Type Method for Positive-Semidefinite Linear
%
   Complementarity Problems, Journal of Optimization Theory and
   Applications: Vol. 86, No. 3, pp. 585-608, 1995.
%
%
   [2] M. S. Bazarraa, H. D. Sherali, and C. M. Shetty, Nonlinear
%
%
   Programming: Theory and Algorithms. John Wiley and Sons, 1993.
%
%
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%
```

```
mu_min
               = 1e-5;
max_iter
               = 20;
b_tol
               = 1e-6;
              = size(M,1);
n
if nargin < 3 || isempty(1)</pre>
   l = zeros(n,1);
   if nargin < 4 || isempty(u)</pre>
     u = inf(n,1);
     if nargin < 5 || isempty(x0)</pre>
                                                output
        x0 = min(max(ones(n,1),1),u);
         if nargin < 6
           display = false;
        end
     end
   end
end
lu
              = [1 u];
              = x0;
Х
[psi,phi,J]
               = FB(x,q,M,1,u);
new_x
              = true;
warning off MATLAB:nearlySingularMatrix
for iter = 1:max iter
   if new_x
     [mlu,ilu]
                   = min([abs(x-1),abs(u-x)],[],2);
                    = max(abs(phi),mlu) < b_tol;</pre>
     bad
     psi
                    = psi - 0.5*phi(bad)'*phi(bad);
                    = J(~bad,~bad);
     J
     phi
                    = phi(~bad);
                    = false;
     new x
     nx
                    = x;
                    = lu(find(bad)+(ilu(bad)-1)*n);
     nx(bad)
                 = J'*J + mu*speye(sum(~bad));
   Н
   Jphi
                 = J'*phi;
   d
                 = -H\Jphi;
```

tol

mu

mu_step

= 1.0e-12;

= 1e-3; = 5;

Initialize the output

Clear the singular part and consider the optimization problem in a subspace

Intuitively, we should follow [1]'s newton method, so we need to find out the Jacobian. [2] develops a new method with drift mu, so H is kind of new Jacobian.

Newton Method

```
Newton Update
           = x(\sim bad) + d;
  nx(~bad)
  [npsi,nphi,nJ] = FB(nx,q,M,l,u);
              = (psi - npsi) / -(Jphi'*d + 0.5*d'*H*d);
                                                         % actual reduction
/ expected reduction
  if r < 0.3
                     % small reduction, increase mu
     mu = max(mu*mu_step,mu_min);
  end
  if r > 0
                     % some reduction, accept nx
     x = nx;
                                                      In addition, I guess it is
     psi = npsi;
                                                      common to use adjustable mu
     phi = nphi;
                                                      to accelerate the algorithm
     J = nJ;
                                                      and this criteria can be
     new x = true;
                                                      found in some convex
     if r > 0.8 % large reduction, decrease mu
                                                      optimization book.
       mu = mu/mu_step * (mu > mu_min);
     end
  end
  if display
     disp(sprintf('iter = %2d,
                                    psi = %3.0e, r =
                                                                %3.1f,
= %3.0e',iter,psi,r,mu));
  end
  if psi < tol
     break;
  end
warning on MATLAB:nearlySingularMatrix
x = min(max(x,1),u);
function [psi,phi,J] = FB(x,q,M,l,u)
    = length(x);
Z1
  = l >-inf & u==inf;
Zu = l = -\inf \& u < \inf;
Zlu = 1 >-inf & u <inf;</pre>
                             Four possible boundaries
Zf
     = l==-inf & u==inf;
     = x;
     = M*x+q;
b
a(Z1) = x(Z1)-1(Z1);
a(Zu) = u(Zu)-x(Zu);
b(Zu) = -b(Zu);
```

More details from

```
note.
if any(Zlu)
  nt
      = sum(Zlu);
  at = u(Zlu)-x(Zlu);
  bt = -b(Zlu);
       = sqrt(at.^2 + bt.^2);
  st
  a(Zlu) = x(Zlu)-l(Zlu);
  b(Zlu) = st -at -bt;
end
      = sqrt(a.^2 + b.^2);
S
      = s - a - b;
phi
                                                    G-Matrix depends
phi(Zu) = -phi(Zu);
                                                    on J
phi(Zf) = -b(Zf);
       = 0.5*phi'*phi;
psi
if nargout == 3
  if any(Zlu)
     M(Zlu,:)
                      -sparse(1:nt,find(Zlu),at./st-ones(nt,1),nt,n)
sparse(1:nt,1:nt,bt./st-ones(nt,1))*M(Zlu,:);
  end
  da
         = a./s-ones(n,1);
  db
         = b./s-ones(n,1);
  da(Zf) = 0;
  db(Zf) = -1;
         = sparse(1:n,1:n,da) + sparse(1:n,1:n,db)*M;
  J
```

end