```
function [uk,ctrlstate] = PIDControl2(yk, U, ctrlPar, ctrlState)
%
% PIDControl()
%
% DESCRIPTION:
% This function implements a discretized proportional-integral-derivative
% (PID) controller for controlling the insulin flow rate. This differs from
% PIDControl by additionally having the insulin vector as input
% INPUT:
% yk
             - Current blood glucose concentration
% U
             - Insulin vector of both basal and bolus insulin
%
% ctrlPar
             - vector of the following:
%
                    * Ts

    Sampling time, 5 min

%
                    * Kp

    Proportional gain

%
                    * Ki

    Integrator gain

%
                                       - Derivative gain
                    * Kd
                                       - The traget glucose concentration, y=108
%
                    * vbar
%
                                       - Nominal insulin flow rate
                    * ubar
%
                    * Ti

    Tuned parameters

                    * Td

    Tuned parameters

%
%
% crtlState - vector of the following:
%
                    * Ik

    The integral term (Ik)

%
                     * vkm1
                                       - Previous glucose concentration, yk-1
%
% OUTPUT:
%

    a vector of manipulated inputs

    crtlstate - the updated controller state
%
%
% PROJECT:
% Fagprojekt 2022
% A diabetes case study - Meal detection
% GENEREL:
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                            : The Technical University of Denmark (DTU)
                            : Applied Mathematics and Computer Science
% Department
% AUTHORS:
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% Unpack control parameters
Ts
       = ctrlPar(1); %
                                   Sampling time
        = ctrlPar(2); %
Κp
                                   Proportional gain
       = ctrlPar(3); %
%Ki
                                   Integrator gain (not used since we calculate it in the ✓
function line 68)
%Kd
     = ctrlPar(4); %
                                   Derivative gain (not used since we calculate it in the ✓
```

```
function line 69)
     = ctrlPar(5); %
                                 Target blood glucose concentration
ybar
       = ctrlPar(6); %
ubar
                                 Nominal insulin flow rate
Τi
      = ctrlPar(7); %
                                 Tuned parameters
Td
      = ctrlPar(8); %
                                 Tuned parameters
% Unpack control state
Ik = ctrlState(1); %
                                Value of integral at previous time step
ykm1 = ctrlState(2); %
                                Previous observed glucose concentration
% Computing
ek = yk-ybar;
                          % Setpoint error
Ki = Kp * Ts/Ti;
                          % Helps controlling the steady state
Kd = Kp * Td/Ts;
                          % The top
Pk = Kp * ek;
                          % Proportional term. Controls how fast the error change
Ikp1 = Ik + Ki * ek;
                          % Integral term. The area of the error
Dk = Kd * (yk-ykm1);
                         % Derivative term. The top of the curve
uba = ubar + Pk + Ik + Dk; % Basal insulin flow rate
ubo = U;
                          % Bolus insulin flow rate
% OUTPUT
% The controlled manipulated inputs at time step
uk = [uba,ubo];
% Controller state OUTPUT
ctrlstate = [Ikp1; yk];
end
```

```
function [T,X] = OpenLoopSimulation(x0, tspan, U, D, p, simModel, simMethod, NK)
% OpenLoopSimulation()
% DESCRIPTION:
% Function peforms an open-loop simulation for given initial condition of
% the state vector, time, intervals, disturbance variables, parameters, and
% simulation model and methods. The open-loop simulation uses the MVPmodel
% and ExplicitEuler to compute both the subcutaneous glucose concentration,
% Gsc(t), the blood glucose concentration G(t) and the statevector x(t) for
% each time step.
%
%
% INPUT:
                 - initial state vector
                                                                  (dimension: 7)
% x0

    time interval to integrate over

                                                                  (dimension N+1)
% tspan

    bolus and basal insulin (manipulated input)

                                                                  (dimension nu \times N)
% U
                 - meal rate (disturbance)
                                                                  (dimension nd \times N)
% D
                 parameter values
                                                                  (dimension np)
% p

    simulation model, MVPmodel

                                                                  (function handle)
% simModel

    simulation method, ExplicitEuler

                                                                  (function handle)
% simMethod

    Number of timesteps in each time interval

% NK
% OUTPUT:
                                                                             (dimension: ∠
% T - The control state of time for each step
% X - The statevector x(t) for each time step stored in a matrix
                                                                             (dimension: nx

✓
\times N+1)
%
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% Number of control steps
N = numel(tspan) - 1;
% Number of states
nx = numel(x0);
% Number of time steps in each control interval
Nk=NK;
```

```
% Allocate memory
T = zeros(1, N+1);
X = zeros(nx, N+1);
% Initial condition in each control interval
xk = x0;
% Store solution
T(1) = tspan(1);
X(:,1) = x0;
% Loop for each time step. Computes Gsc(t), G(t), x(t) from k = 0 to N.
for k = 1:N
    % Times
         = tspan(k);
    tk
    tkp1 = tspan(k+1);
    % Manipulated inputs and disturbance variables
    uk = U(:,k);
    dk = D(:,k);
    % Time interval
    tspank = linspace(tk, tkp1, Nk+1);
    % Solve initial value problem
    [Tk, Xk] = simMethod(simModel, tspank, xk, uk, dk, p);
    % Update initial condition
    xk = Xk(end, :)';
    % Store solution
    T(k+1) = Tk(end)';
    X(:, k+1) = Xk(end, :)';
end
```

end

```
function [ Gfm_vec , filt_prev , flag, zero_one ] = GRID_func( ...
         delta_G , G_vec , tau, tspan , filt_prev , Gmin, Gfm_vec , t_vec, flag)
% GRID_func()
%
% DESCRIPTION:
% The function is a part of the GRID algortihm. This is part of the
% dectection logic, the last part of the algoriihm.
% The function takes in the glucose measurements, calls the two filter
% functions and finds the derivatives. After this it is able to detect
% weather or not there a meal has been detected. Lastly, it counts down
% such that a meal will not be detected twice within two hours
% INPUT:
                         - The maximum ROC (rate of change)
% delta_G
%
% G_vec

    Vector consisting of the glucose value and the

%
                          two previous glucose measurements.
%
                           As follows: [Gm-2, Gm-1, Gm].
                        - The interval step given as a number
% tspan
%
% filt_prev

    Vector of previous filteret glucose measurements

%
                          As follows: [G_{F,NS}(k-1), G_{F}(k-2)].
                          For equation (1)&(3).
%
%
% Gmin
                         - Vector of minumum glucose measurements
%
                          As follows: [G_{\min,1},G_{\min,2},G_{\min,3}].
%
                          For equation (4).
%
% Gfm_vec
                        - Vector of previous derivatives
                          As follows: [G'_{F}(k-2), G'_{F}(k-2)].
%
%
                          For equation (4).
%
% t_vec
                        - Vector of sampling time respectively for G
%
% flag
                         - For counting the time from last detected meal
%
% OUTPUT:
% Gfm vec
                         - The stored new vector of the previous filtered
%
                           glucose measurements.
%
                          As follows: [G'_{F}(k-1), G'_{F}(k)].
%
% G_prev
                        - The stored new vector of previous glucose
%
                          measurements.
%
                          As follows: [G_{F,NS}(k-2), G_{F}(k-2)].
%
% zero_one
                        - 1 or 0 for detected meal.
%
                         - For counting the time from last detected meal
% flag
%
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% REFERENCE:
% MANGLER FRA ARTIKEL
% Inisializing all values
                         % The second previous derivative used in eugation 4
Gfm m2 = Gfm vec(1);
Gfm_m1 = Gfm_vec(2);
                         % The previous derivative used in equation 4
Gfnsm2 prev = filt prev(1); % The second previous noise-spike filtered value
                         % used in equation 1
Gfm2 prev
            = filt_prev(2); % The second previous low filterd value used in
                         % equation 2
% Minimum values used in equation 4
Gmin1 = Gmin(1);
Gmin2 = Gmin(2);
Gmin3 = Gmin(3);
% The two previous measured gluscose values and the one at control state
        = G_vec(1);
                                % The second previous glucose value
Gm2
                                 % The previous glucose value
        = G_{vec}(2);
Gm1
                                 % The glucose value at control state
G
        = G_{\text{vec}}(3);
% COMPUTING
% The noise-spike filter at the 3 sampling times
Gfnsm2 = spikefilt_func(Gm2,Gfnsm2_prev,delta_G);
Gfnsm1 = spikefilt_func(Gm1,Gfnsm2,delta_G);
       = spikefilt_func(G,Gfnsm1,delta_G);
% The low filter at the 3 sampling times
Gfm2 = lowfilt func(tau,tspan,Gfnsm2,Gfm2 prev);
Gfm1 = lowfilt_func(tau,tspan,Gfnsm1,Gfm2);
     = lowfilt_func(tau,tspan,Gfns,Gfm1);
% Inisializing input for estimate_lagrange
Gf_vec = [ Gfm2 , Gfm1 , Gf ];
% Computing the first derivative using lagrange
       = estimate_lagrange(t_vec,Gf_vec); % Returns the derivative
% The detection part from equation 4
if (Gf > Gmin1) && ...
   ((Gfm > Gmin3) && (Gfm_m1 > Gmin3) && (Gfm_m2 > Gmin3) ...
```

```
|| (Gfm > Gmin2) && (Gfm_m2 > Gmin2) )
    zero_one = 1; % A meal has been detected
else
    zero_one = 0; % No meal has been detected
end
% Output
filt_prev = [Gfnsm2,Gfm2]; % Outputting the updated filtered values
Gfm_vec = [Gfm_m1,Gfm]; % Outputting the updated derivative values
% COUNTING PART
if flag > 0
% flag larger than 0 means that a meal has been detected within 120 min
% implying that a meal should not be detected again already. So zero_one
% is set to 0. Therefore, flag is subtracted by -1, such that it will
% count down so a meal can be detetected again after 120 min.
    flag = flag-1;
    zero_one = 0;
elseif flag == 0 && zero_one == 1
 % flag equal to zero means a meal may be detected again
 % but only is if zero_one equals 1. When this happen flag start over
 % counting down 120 min.
    flag = 120/tspan;
end
end
```

```
function [truepositive, falsepositive] = evaluationfunction(stride, D, D_detected, Ts, N)
%
% detectionrates()
%
% DESCRIPTION:
% This function computes the TP, FP, TN, FN based on a true vector with
% detected meals D, and a computed vector with detected meals D_detected.
% It uses a stride since the meal will be detected a time period after the
% meals was given.
% It uses the indices for the true meals and the detected meals to compare
% if in the stride a meals should be detected or not and vise versa.
% INPUT:
                    - The maximal time period it takes from the meal to be
% stride
% detected
% D
                    - The true vector of real meals.
                    - The estimated vecor of 0 or 1. 1 meaning meals is
% D_detected
% detected
                    - Bolus insulin
% U
%
% Ts
                    - The time between control steps
                    - The number of observations
% N
%
% OUTPUT:
% Two outputs being TP, FP
% PROJECT:
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% Initializing
falsenegative
                = 0;
falsepositive
                = 0;
                = 0;
truepositive
truemeals
                = zeros(1,N);
mealdetec
                = zeros(1,N);
% Changing datatype of D to binary
for i = 1:N
    if D(1,i) >= 50/Ts \% Not considering the snackmeals
        D(1,i) = 1;
```

```
else
        D(1,i) = 0;
    end
end
% Finding the indices for the truemeals
for i = 1 : N
    if D(1,i) == 1
        truemeals(i) = i;
    end
end
% Finding the indices for the detected meals
for i = 1 : N
    if D_detected(i) == 1
        mealdetec(i) = i;
    end
end
% Removing all the zeros so there is only the indices left
idxdetecmeals = nonzeros(mealdetec)';
idxtruemeals = nonzeros(truemeals)';
% Examine if there are no true meals where there are detected meals
for i = 1:length(idxdetecmeals)
    % The idx value when meal has been detected
   k = idxdetecmeals(i);
    if (k-stride) < 1</pre>
        j = k-1;
        if sum(D(1,k-j:k)) == 0
        falsepositive = falsepositive + 1;
        end
    elseif sum(D(1,k-stride:k)) == 0
        falsepositive = falsepositive + 1;
    end
end
% Examine if there are true meals where there are detected meals
for i = 1:length(idxdetecmeals)
    % The idx value when meal has been detected
   k = idxdetecmeals(i);
    if (k-stride) < 1</pre>
        j = k-1;
        if sum(D(1,k-j:k)) == 0
        falsepositive = falsepositive + 1;
        end
    elseif sum(D(1,k-stride:k)) == 1
        truepositive = truepositive + 1;
```

end end

end

```
function [T, X, Y, U, ctrlState] = ClosedLoopSimulation_withnoise2(tspan,x0,D,U,p, ...
    ctrlAlgorithm, simMethod, simModel, observationMethod, ctrlPar,ctrlState0,NK,∠
intensity)
%
% ClosedLoopSimulation()
%
% DESCRIPTION:
% Performs a closed-loop simulation of a model-based control algorithm for
% given time range, initial condition, disturbance variables, insulin
% levels, parameters, control algorithm, simulation model, observation
% model, control parameters, control state, control intervals and intensity
% level.
% Closed-loop is used when the input depends on the
% output; this function is part of the PID-controller.
%
% INPUT:

    boundaries of the control intervals

                                                                                (dimension: ∠
%
    tspan
N+1
      )
                                                                                (dimension: ∠
%
    x0

    initial state

nx
%

    disturbance variables for each control interval

    D
                                                                                (dimension: ∠
nd \times N)

    Insulin levels of both bolus and basal

    U
                                                                                (dimension nu∠
%
\times N)
                                                                                (dimension: ∠
%
                         parameters
    р
    )
np
%
    simModel
                         - simulation model
                                                                                (function ∠
handle)
                                                                                (function ∠
%
    ctrlAlgorithm

    control algorithm

handle)
%
                         controller parameters
    ctrlPar
                         - initial controller state
    ctrlState0
                                                                                (dimension: ∠
%
nc)

    simulation method

                                                                                (function ∠
%
    simMethod
handle)

    Number of steps in each control interval

%
                                                                                (scalar)
                         - The intensity value used for Euler Maruyama
%
    intensity
                                                                                (scalar)
%
% OUTPUT:
   T - boundaries of control intervals (=tspan)
                                                      (dimension:
                                                      (dimension: nx \times N+1)
   X – the states in the simulation model
   X - the observed variables
                                                      (dimension: ny \times N+1)
    U - the computed manipulated inputs
                                                      (dimension: nu \times N)
%
    ctrlState - matrix of controller states
                                                      (dimension: nc \times N+1)
%
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% Initial time
t0 = tspan(1);
% Observed variables of glucose at steady state
y0 = observationMethod(x0,p);
% manipulated inputs calculated
uDummy = ctrlAlgorithm(y0,U(1,1), ctrlPar, ctrlState0);
% Number of each variable
nx = numel(x0);
                        % states
                        % glucose concentration
ny = numel(y0);
nu = numel(uDummy);
                       % manipulated inputs
nc = numel(ctrlState0); % glucose concentration and integral term
% Number of control intervals
N = numel(tspan)-1;
% Number of time steps in each control interval
Nk = NK;
% Initialising Output
T = zeros(1, N+1);
X = zeros(nx, N+1);
Y = zeros(ny, N+1);
ctrlState = zeros(nc, N+1);
% Storing solution
T(1) = t0;
X(:,1) = x0;
Y(:,1) = y0;
ctrlState(:, 1) = ctrlState0;
% Copying initial condition to another name
tk = t0;
xk = x0;
yk = y0;
for k = 1:N
    %%% Initializing the loop
    % Time
    tkp1 = tspan(k+1);
    % Time interval
    tspank = linspace(tk, tkp1, Nk+1);
    % Controller state
    ctrlStatek = ctrlState(:, k);
```

```
% Disturbance variables
dk = D(:, k);
%%% Start computing
% Compute manipulated inputs after the PID control
[uk, ctrlStatekp1] = ctrlAlgorithm(yk, U(2 ,k), ctrlPar, ctrlStatek);
% Solving the differential equation with euler maruyama
[Tk, Xk] = simMethod(simModel, tspank, xk, uk, dk, p,intensity);
%% Overwriting for the next loop
% States at the next time step
tkp1 = Tk(end);
xkp1 = Xk(end,:)';
% Observed variables at the next time step
ykp1 = xkp1(7);
% Update initial condition
tk = tkp1;
xk = xkp1;
yk = ykp1;
%%% Storing solution
% Store solution and updating conditions
T(k+1) = tkp1;
X(:, k+1) = xkp1;
Y(:, k+1) = ykp1;
U(:, k) = uk;
ctrlState(:, k+1) = ctrlStatekp1;
```

end