MSc Project IIS1 | BWFLnet_Water Analysis

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```
clc
clear
close all
```

Model Setup and Initialisation

Initialisation of EPANET Matlab toolkit and loading of .net file.

```
epanet_path = 'E:\Program Files\MATLAB\R2021b\toolbox\epanet\EPANET-Matlab-Toolkit-master';
net_id = 'BWFLnet_MSc_2022_calibrated';
run( [epanet_path, '\start_toolkit'])

EPANET-MATLAB Toolkit Paths Loaded.

net = epanet([net_id,'.inp']);

EPANET version {20200} loaded (EMT version {v2.2.0}).
Loading File "BWFLnet_MSc_2022_calibrated.inp"...
Input File "BWFLnet_MSc_2022_calibrated.inp" loaded sucessfuly.

load("wq_data.mat");
```

Network Data Loading

Load general network data, noting that this network has already been hydraulically calibrated, thus we can directly perform the quality analysis.

```
% Save element count variables
nn = net.NodeJunctionCount;
n0 = net.NodeReservoirCount;
np = net.LinkCount; % count of pipes and valves
D = net.getLinkDiameter';
R = net.getLinkRoughnessCoeff';
% Save node and link index vectors
Reservoir_Idx = net.getNodeReservoirIndex;
Junction_Idx = net.getNodeJunctionIndex;
Link Idx = double(net.getLinkIndex);
% Specify hydraulic and quality time steps (in seconds)
net.setTimeQualityStep(15*60); % 15 minute quality time steps
net.setTimeHydraulicStep(15*60); % 15 minute hydraulic time steps
nt = net.getTimeSimulationDuration./3600; % number of time steps in hours
days = 7; % simulation duration in days
net.setTimeSimulationDuration(nt*3600*days/7); % Set the simulation duration by changing days
nt = net.getTimeSimulationDuration./3600; % get the new nt
```

Graph Creation Preparation

Use graph theory to plot network connectivity and spatial coordinates.

```
% Create A12 and A10 incidence matrices
LinkNodesList = net.getLinkNodesIndex;
A = zeros(np,nn+n0);
for k=1:np
    i = LinkNodesList(k,1);
    j = LinkNodesList(k,2);
    A(k,i) = -1;
    A(k,j) = 1;
end
A12 = A(:,Junction_Idx);
A10 = A(:,Reservoir_Idx);
A12 = sparse(A12);
A10 = sparse(A10);
% Obstain node XY and elevation information
XY = zeros(nn+n0,2);
XY(:,1) = net.getNodeCoordinates{1};
XY(:,2) = net.getNodeCoordinates{2};
elev = double(net.getNodeElevations(Junction Idx))';
% Create adjacency matrix
A = [A12,A10];
AdjA = sparse(size(A,2),size(A,2));
for k = 1:size(A,1)
    node_in = find(A(k,:) == -1);
    node_out = find(A(k,:) == 1);
    AdjA(node in, node out) = 1;
    AdjA(node_out,node_in) = 1;
end
gr = graph(AdjA);
```

Chlorine Simulation

Setup general chemical simulation parameters.

```
% Initialise EPANET simulation type
net.setQualityType('Chemical','mg/L');
net.setNodeSourceQuality(1:nn+n0,zeros(nn+n0,1)); % set all node source quality to zero
% Load the water quality data
Junction_SelectIdx = net.getNodeIndex(wq_data.node_ids);
Junction_SelectIdx = Junction_SelectIdx([1,4,7,9]);
cext = wq_data.chlorine([3,6],:);
```

Assign chlorine decay parameters.

```
% Link bulk reaction coefficients
```

```
lambda = 0.5*ones(np,1); % units of days^-1
net.setLinkBulkReactionCoeff(1:np, -lambda);
net.setOptionsPipeBulkReactionOrder(1);
% Link wall decay reaction coefficients... apply by pipe size for S&G?
% theta = ones(np,1); % units of m/day
% for k = 1:np
%
      if D(k) <= 75
%
          theta(k) = 0.25;
%
      elseif D(k) > 75 \&\& D(k) <= 150
%
          theta(k) = 0.175;
%
      elseif D(k) > 150 \&\& D(k) <= 250
%
          theta(k) = 0.1;
%
      else
%
          theta(k) = 0.05;
%
      end
% end
%
% net.setLinkWallReactionCoeff(1:np, -theta);
% net.setOptionsPipeWallReactionOrder(1);
% Initial concentrations at nodes (mg/L)
c0 = zeros(nn+n0,nt);
net.setNodeInitialQuality(net.NodeIndex,c0);
% % Assign source chlorine with time varying pattern
% cext = ones(n0,nt);
% % Constant source concentration for n0 reservoirs
% % cext = cext*0.25;
%
% % Varying source concentration by source element
% source_c = [1.00, 0.85];
% for j = 1:n0
      cext(j,:) = source_c(j)*cext(j,:);
% end
% Varying source concentration by source element with
% concentration spike (pulse) at t = 12
% t = 144;
% for i = 1:4*days*nt
%
      for j = 1:n0
%
          if i == t
%
              cext(j,i) = 1*cext(j,i); % Pulse of 1 mg/L
%
          else
%
              cext(j,i) = source_c(j)*cext(j,i);
%
          end
%
      end
% end
base_cext = ones(n0,1);
pattern_cext = cext./(base_cext*ones(1,size(cext,2))); % extend vector over nt columns
```

```
% For loop to assign new patterns to source contrations at reservoirs and
% source type
for i=1:n0
    patternId = sprintf('Res_C_%d',i);
    net.addPattern(patternId,pattern_cext(i,:));
    net.setNodeSourcePatternIndex(net.NodeReservoirIndex(i),net.getPatternIndex(patternId));
    net.setNodeSourceQuality(net.NodeReservoirIndex(i),base_cext(i));
    net.setNodeSourceType(net.NodeReservoirIndex(i),'CONCEN');
end
```

Results from EPANET Simulation

Simulate hydraulic and quality analyses using EPANET's solvers.

```
hydraulic_res = net.getComputedHydraulicTimeSeries;
quality_res = net.getComputedQualityTimeSeries;

% Assign hydraulic results to network elements
h = hydraulic_res.Head(1:1+4*nt,1:nn).';
q = 1e-3*hydraulic_res.Flow(1:1+4*nt,:).';

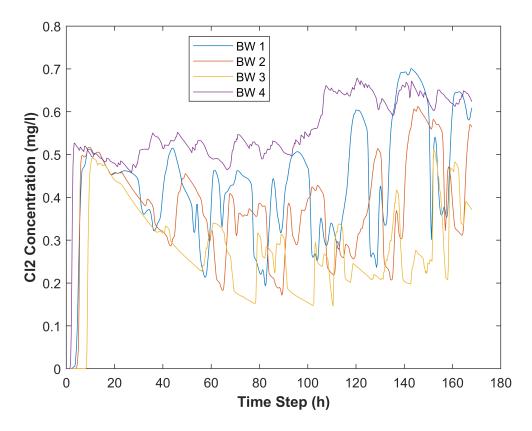
% Assign quality results to network elements
c_nodes = quality_res.NodeQuality';
c_pipes = quality_res.LinkQuality';
```

Genetic Algorithm Preparation

Implement genetic algorithm to solve the parameter optimization problem.

```
% Get the simulated and observed data
c_nodes_Sim = c_nodes(Junction_SelectIdx,:);
c_nodes_Obs = wq_data.chlorine([1,4,7,9],:);
```

Plot the simulated chlorine concentration for visualisation.



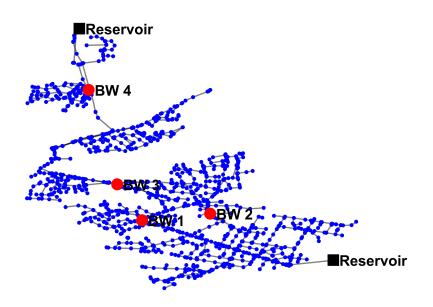
According to the water age analysis, the longest water age before periodic behaviour is approximately 40 hours, thus we split the data into three parts, the first part is a 2-day stabilisation period, the second part is a 1-day train data, the last part is the 4-day validation time.

```
% Data splitting
c_nodes_Sim = c_nodes_Sim(:,2:end);
Sim_train = c_nodes_Sim(:,2*96+1:3*96);
Sim\ val = c\ nodes\ Sim(:,3*96+1:end);
Obs_train = c_nodes_Obs(:,2*96+1:3*96);
Obs val = c nodes Obs(:,3*96+1:end);
mse function = \emptyset(sim,obs) (1/size(obs,1)/size(obs,2))*sum(sum((sim - obs).^2));
mse_before = mse_function(Sim_val,Obs_val);
% Plot the sensor location on the network map
figure
p1 = plot(gr);
p1.XData = XY(:,1);
p1.YData = XY(:,2);
p1.LineWidth = 1;
p1.EdgeColor = 'k';
p1.MarkerSize = 2;
p1.NodeColor = 'b';
p1.NodeLabel = '';
highlight(p1,Junction SelectIdx,'NodeColor','r','MarkerSize',8);
labelnode(p1,Junction SelectIdx,Junction Name);
highlight(p1,Reservoir_Idx,'NodeColor','k','Marker','s','MarkerSize',10);
labelnode(p1,Reservoir_Idx,{'Reservoir','Reservoir'});
```

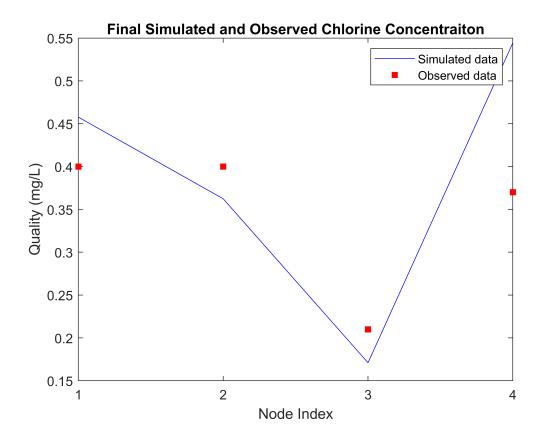
```
p1.NodeFontSize = 11;
p1.NodeLabelColor = 'k';
p1.NodeFontWeight = 'bold';
p1.EdgeFontSize = 11;
p1.EdgeLabelColor = 'k';
p1.EdgeFontWeight = 'bold';

axis('off')
title('Sensor Location in the Network')
```

Sensor Location in the Network

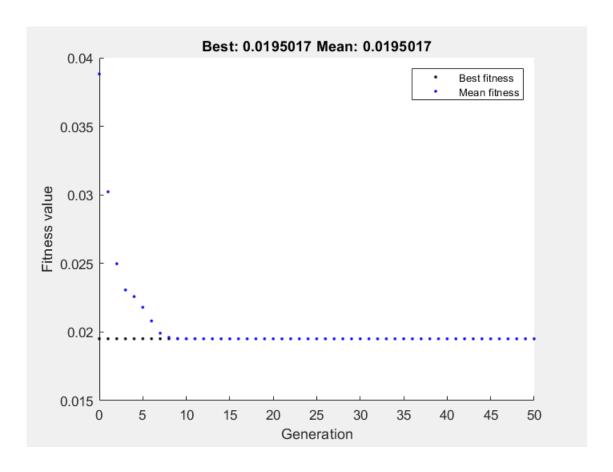


```
% Plot the calibrated quality-time figure
figure,
plot(1:size(c_nodes_Obs,1),Sim_train(:,96),'-b')
hold on
scatter(1:size(c_nodes_Obs,1),Obs_train(:,96),'red','filled','square')
legend({'Simulated data','Observed data'})
xlabel('Node Index')
ylabel('Quality (mg/L)')
xticks(1:1:size(c_nodes_Obs,1));
title('Final Simulated and Observed Chlorine Concentraiton')
```



No Group Genetic Algorithm Implementation

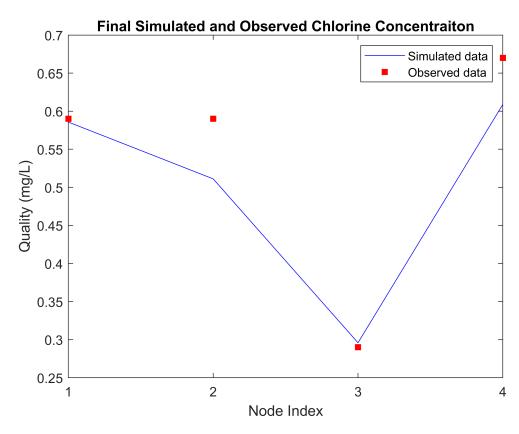
Implement genetic algorithm to solve the parameter optimization problem based on one group.



```
% Test the ga result
theta_NoGroup = ones(2281,1); % units of m/day
for i = 1:2281
         theta_NoGroup(i) = x_NoGroup(1);
end
net.setLinkBulkReactionCoeff(1:np, theta_NoGroup);
net.setOptionsPipeBulkReactionOrder(1);
quality_res_NoGroup = net.getComputedQualityTimeSeries;
```

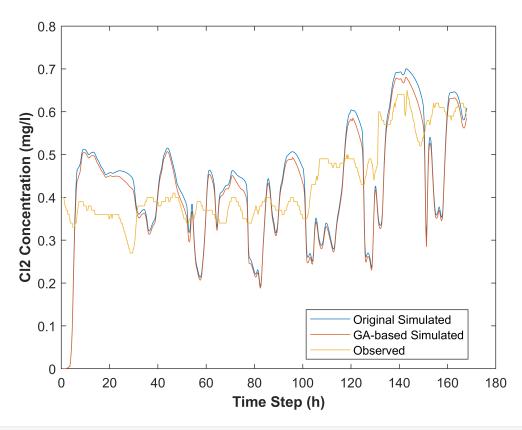
```
net.getLinkBulkReactionCoeff;
c_nodes_NoGroup = quality_res_NoGroup.NodeQuality';
c_nodes_Sim_NoGroup = c_nodes_NoGroup(Junction_SelectIdx,2:end);
Sim_NoGroup_val = c_nodes_Sim_NoGroup(:,3*96+1:end);
mse_after_NoGroup = mse_function(Sim_NoGroup_val,Obs_val);

% Plot the ga-modified quality-time figure
figure,
plot(1:size(c_nodes_Obs,1),Sim_NoGroup_val(:,end),'-b')
hold on
scatter(1:size(c_nodes_Obs,1),Obs_val(:,end),'red','filled','square')
legend({'Simulated data','Observed data'})
xlabel('Node Index')
ylabel('Quality (mg/L)')
xticks(1:1:size(c_nodes_Obs,1));
title('Final Simulated and Observed Chlorine Concentraiton')
```



```
% BW1 as an example
figure
plot(0.25:0.25:days*24,c_nodes_Sim(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Sim_NoGroup(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Obs(1,:),'-')

legend({'Original Simulated','GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 Concentration (mg/l)','fontweight','bold')
```



fprintf('The mse before calibration is %.4f, after calibration it is %.4f', mse_before, mse_after

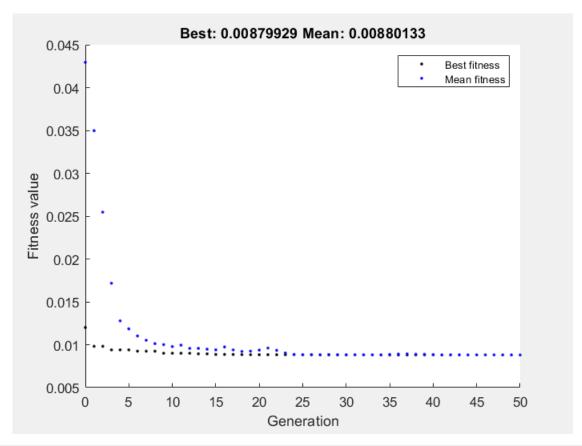
Diameter based Group Genetic Algorithm Implementation

Implement genetic algorithm to solve the parameter optimization problem based on diameter groups.

```
% % genetic algorithm implementation
rng default
options = optimoptions('ga','Display','none','PopulationType','DoubleVector','PopulationSize',!
    'Generations',50,"ConstraintTolerance",1e-6,"PlotFcn","gaplotbestf");

lb = [-10,-10,-10,-10];
ub = [0,0,0,0];

Mse_function = @(x) Mse_D(x,Obs_train,Junction_SelectIdx,D);
[x_D, fval_D] = ga(Mse_function,4,[],[],[],[],lb,ub,[],options);
```

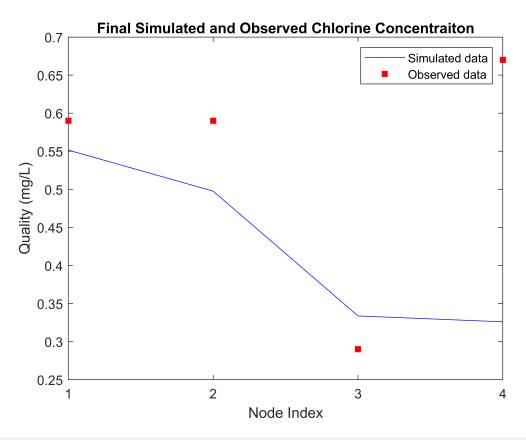


```
% Test the ga result
theta_D = ones(2281,1); % units of m/day
for i = 1:2281
    if D(i) <= 75
        theta_D(i) = x_D(1);
    elseif D(i) > 75 \&\& D(i) <= 150
        theta D(i) = x D(2);
    elseif D(i) > 150 \&\& D(i) <= 250
        theta_D(i) = x_D(3);
    else
        theta_D(i) = x_D(4);
    end
end
net.setLinkBulkReactionCoeff(1:np, theta_D);
net.setOptionsPipeBulkReactionOrder(1);
quality_res_D = net.getComputedQualityTimeSeries;
c_nodes_D = quality_res_D.NodeQuality';
c nodes Sim D = c nodes D(Junction SelectIdx,2:end);
```

```
Sim_D_val = c_nodes_Sim_D(:,3*96+1:end);
mse_after_D = mse_function(Sim_D_val,Obs_val);

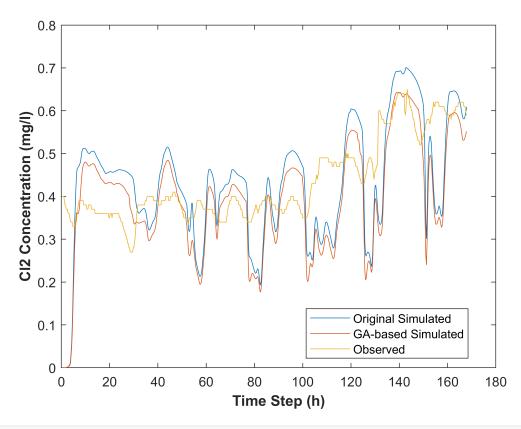
% Plot the ga-modified quality-time figure
figure,
plot(1:size(c_nodes_Obs,1),Sim_D_val(:,end),'-b')
```

```
hold on
scatter(1:size(c_nodes_Obs,1),Obs_val(:,end),'red','filled','square')
legend({'Simulated data','Observed data'})
xlabel('Node Index')
ylabel('Quality (mg/L)')
xticks(1:1:size(c_nodes_Obs,1));
title('Final Simulated and Observed Chlorine Concentraiton')
```



```
% Node 753 as an example
figure
plot(0.25:0.25:days*24,c_nodes_Sim(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Sim_D(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Obs(1,:),'-')

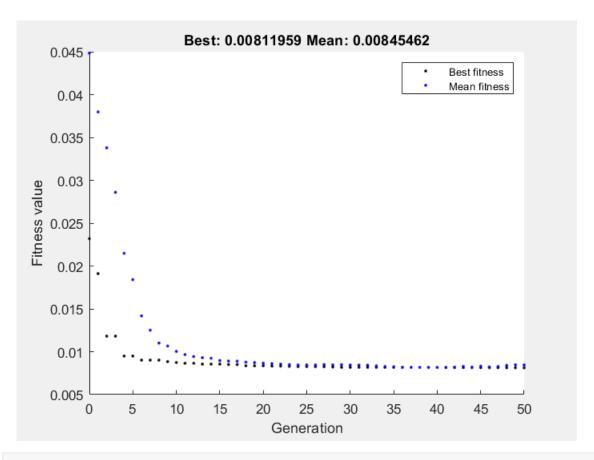
legend({'Original Simulated','GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 Concentration (mg/l)','fontweight','bold')
```



fprintf('The mse before calibration is %.4f, after calibration it is %.4f', mse_before, mse_after

Roughness based Group Genetic Algorithm Implementation

Implement genetic algorithm to solve the parameter optimization problem based on roughness groups.

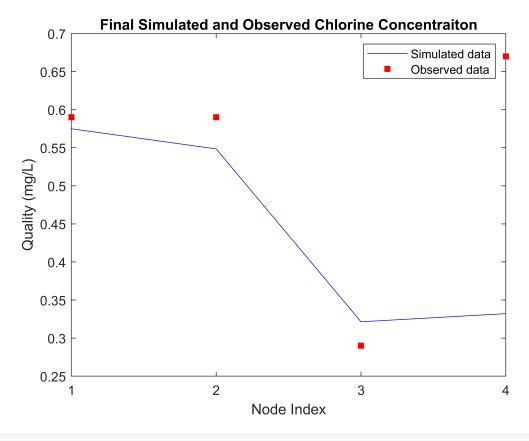


```
% Test the ga result
theta_R = ones(2281,1); % units of m/day
for i = 1:2281
    if abs(R(i) - 0) <= 0.01
        theta_R(i) = x_R(1);
    elseif abs(R(i) - 0.1000) <= 0.01
        theta_R(i) = x_R(2);
    elseif abs(R(i) - 0.1399) <= 0.01
        theta_R(i) = x_R(3);
    elseif abs(R(i) - 1.5267) <= 0.01
        theta_R(i) = x_R(4);
    elseif abs(R(i) - 2.1576) <= 0.01
        theta_R(i) = x_R(5);
    elseif abs(R(i) - 9.0114) <= 0.01
        theta_R(i) = x_R(6);
    elseif abs(R(i) - 10.3677) <= 0.01</pre>
        theta_R(i) = x_R(7);
    elseif abs(R(i) - 11.6809) <= 0.01
        theta_R(i) = x_R(8);
    elseif abs(R(i) - 14.8687) <= 0.01
        theta R(i) = x R(9);
    elseif abs(R(i) - 16.2377) <= 0.01
        theta_R(i) = x_R(10);
    else
        theta_R(i) = x_R(11);
    end
end
```

```
net.setLinkBulkReactionCoeff(1:np, theta_R);
net.setOptionsPipeBulkReactionOrder(1);
quality_res_R = net.getComputedQualityTimeSeries;
c_nodes_R = quality_res_R.NodeQuality';
c_nodes_Sim_R = c_nodes_R(Junction_SelectIdx,2:end);
```

```
Sim_R_val = c_nodes_Sim_R(:,3*96+1:end);
mse_after_R = mse_function(Sim_R_val,Obs_val);

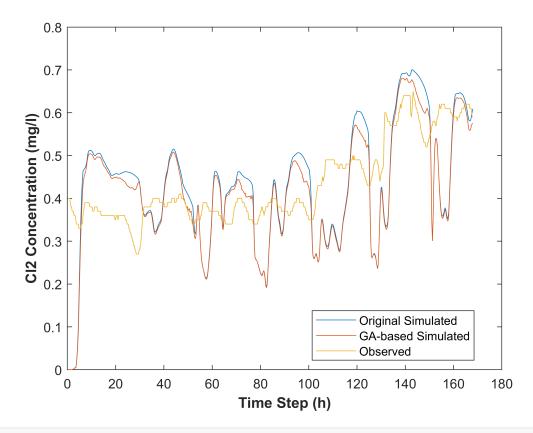
% Plot the ga-modified quality-time figure
figure,
plot(1:size(c_nodes_Obs,1),Sim_R_val(:,end),'-b')
hold on
scatter(1:size(c_nodes_Obs,1),Obs_val(:,end),'red','filled','square')
legend({'Simulated data','Observed data'})
xlabel('Node Index')
ylabel('Quality (mg/L)')
xticks(1:1:size(c_nodes_Obs,1));
title('Final Simulated and Observed Chlorine Concentraiton')
```



```
% Node 753 as an example
figure
plot(0.25:0.25:days*24,c_nodes_Sim(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Sim_R(1,:),'-')
```

```
hold on
plot(0.25:0.25:days*24,c_nodes_Obs(1,:),'-')

legend({'Original Simulated','GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 Concentration (mg/l)','fontweight','bold')
```



fprintf('The mse before calibration is %.4f, after calibration it is %.4f', mse_before, mse_after

The mse before calibration is 0.0371, after calibration it is 0.0247

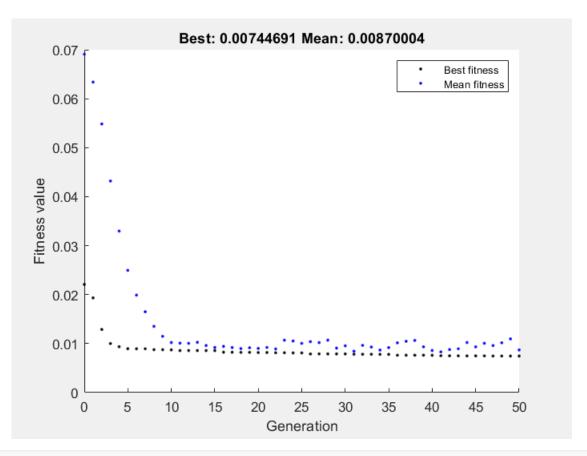
Roughness and Diameter based Group Genetic Algorithm Implementation1

Implement genetic algorithm to solve the parameter optimization problem based on roughness and diameter groups, in this we consider the roughness groups affect the bulk decay and diameter groups affect the wall decay.

```
% genetic algorithm implementation
rng default
options = optimoptions('ga','Display','none','PopulationType','DoubleVector','PopulationSize',!
    'Generations',50,"ConstraintTolerance",1e-6,"PlotFcn","gaplotbestf");

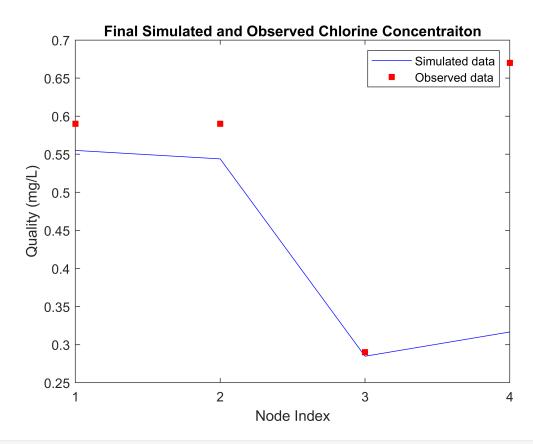
lb = -10*ones(15,1);
ub = 0*ones(15,1);

Mse_function = @(x) Mse_DR(x,Obs_train,Junction_SelectIdx,D,R);
[x_DR, fval_DR] = ga(Mse_function,15,[],[],[],[],lb,ub,[],options);
```



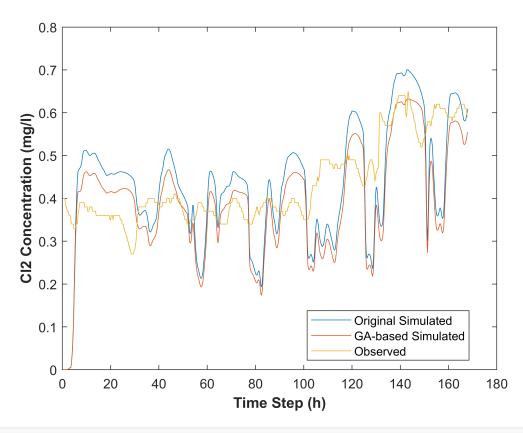
```
% Test the ga result
thetaR = ones(2281,1); % units of m/day
for i = 1:2281
    if abs(R(i) - 0) <= 0.01
        thetaR(i) = x_DR(1);
    elseif abs(R(i) - 0.1000) <= 0.01
        thetaR(i) = x DR(2);
    elseif abs(R(i) - 0.1399) <= 0.01
        thetaR(i) = x_DR(3);
    elseif abs(R(i) - 1.5267) <= 0.01
        thetaR(i) = x_DR(4);
    elseif abs(R(i) - 2.1576) <= 0.01
        thetaR(i) = x_DR(5);
    elseif abs(R(i) - 9.0114) <= 0.01
        thetaR(i) = x DR(6);
    elseif abs(R(i) - 10.3677) <= 0.01
        thetaR(i) = x_DR(7);
    elseif abs(R(i) - 11.6809) <= 0.01
        thetaR(i) = x_DR(8);
    elseif abs(R(i) - 14.8687) <= 0.01
        thetaR(i) = x DR(9);
    elseif abs(R(i) - 16.2377) <= 0.01
        thetaR(i) = x_DR(10);
    else
        thetaR(i) = x_DR(11);
    end
end
```

```
thetaD = ones(2281,1); % units of m/day
for i = 1:2281
    if D(i) <= 75
        thetaD(i) = x_DR(12);
    elseif D(i) > 75 \&\& D(i) <= 150
        thetaD(i) = x_DR(13);
    elseif D(i) > 150 \&\& D(i) <= 250
        thetaD(i) = \times DR(14);
    else
        thetaD(i) = x_DR(15);
    end
end
net.setLinkBulkReactionCoeff(1:np, thetaR);
net.setOptionsPipeBulkReactionOrder(1);
net.setLinkWallReactionCoeff(1:np, thetaD);
quality_res_DR = net.getComputedQualityTimeSeries;
c_nodes_DR = quality_res_DR.NodeQuality';
c nodes Sim DR = c nodes DR(Junction SelectIdx,2:end);
Sim DR val = c nodes Sim DR(:,3*96+1:end);
mse_after_DR = mse_function(Sim_DR_val,Obs_val);
% Plot the ga-modified quality-time figure
figure,
plot(1:size(c_nodes_Obs,1),Sim_DR_val(:,end),'-b')
hold on
scatter(1:size(c nodes Obs,1),Obs val(:,end),'red','filled','square')
legend({'Simulated data','Observed data'})
xlabel('Node Index')
ylabel('Quality (mg/L)')
xticks(1:1:size(c nodes Obs,1));
title('Final Simulated and Observed Chlorine Concentraiton')
```



```
% Node 753 as an example
figure
plot(0.25:0.25:days*24,c_nodes_Sim(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Sim_DR(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Obs(1,:),'-')

legend({'Original Simulated','GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 Concentration (mg/l)','fontweight','bold')
```



fprintf('The mse before calibration is %.4f, after calibration it is %.4f', mse_before, mse_after

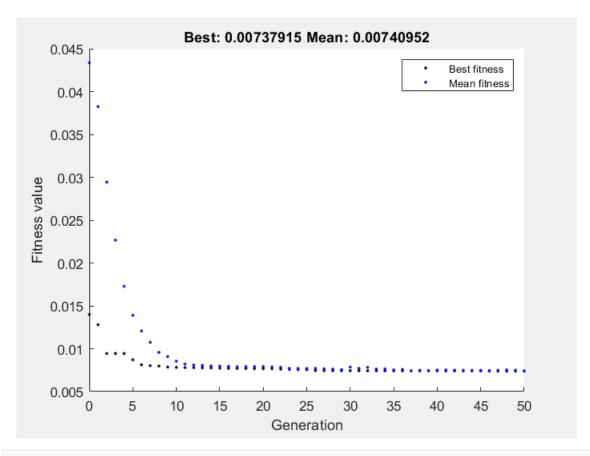
Roughness and Diameter based Group Genetic Algorithm Implementation2

Implement genetic algorithm to solve the parameter optimization problem based on roughness and diameter groups, in this we consider the roughness groups and diameter groups affect the bulk decay together.

```
% genetic algorithm implementation
rng default
options = optimoptions('ga','Display','none','PopulationType','DoubleVector','PopulationSize',!
    'Generations',50,"ConstraintTolerance",1e-6,"PlotFcn","gaplotbestf");

lb = -10*ones(22,1);
ub = 0*ones(22,1);

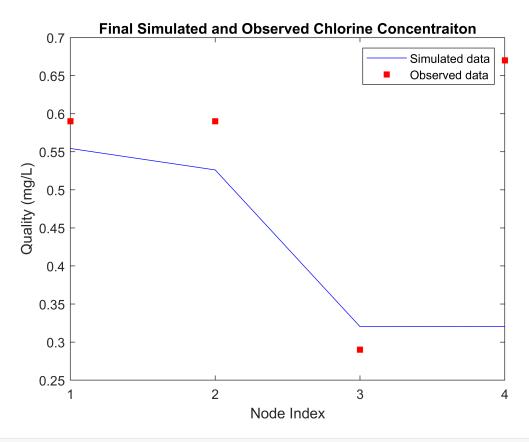
Mse_function = @(x) Mse_DandR(x,Obs_train,Junction_SelectIdx,D,R);
[x_DandR, fval_DandR] = ga(Mse_function,22,[],[],[],[],lb,ub,[],options);
```



```
% Test the ga result
thetaDandR = ones(2281,1); % units of m/day
for i = 1:2281
    if abs(R(i) - 0) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(1);
        else
            thetaDandR(i) = x_DandR(2);
        end
    elseif abs(R(i) - 0.1000) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(3);
        else
            thetaDandR(i) = x_DandR(4);
        end
    elseif abs(R(i) - 0.1399) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(5);
        else
            thetaDandR(i) = x_DandR(6);
        end
    elseif abs(R(i) - 1.5267) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(7);
        else
            thetaDandR(i) = x_DandR(8);
        end
```

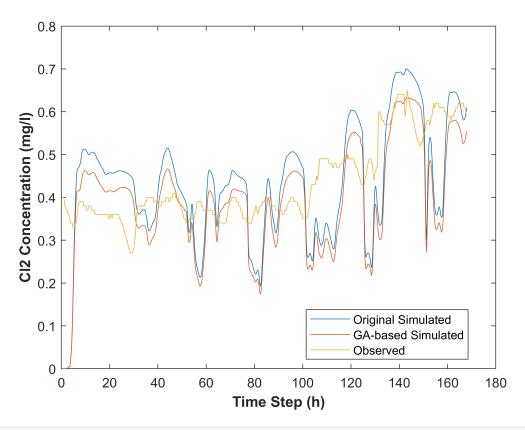
```
elseif abs(R(i) - 2.1576) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(9);
        else
            thetaDandR(i) = x_DandR(10);
        end
    elseif abs(R(i) - 9.0114) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(11);
        else
            thetaDandR(i) = x_DandR(12);
        end
    elseif abs(R(i) - 10.3677) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x DandR(13);
        else
            thetaDandR(i) = x_DandR(14);
        end
    elseif abs(R(i) - 11.6809) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(15);
        else
            thetaDandR(i) = x_DandR(16);
        end
    elseif abs(R(i) - 14.8687) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(17);
        else
            thetaDandR(i) = x_DandR(18);
        end
    elseif abs(R(i) - 16.2377) <= 0.01
        if D(i) <= 150
            thetaDandR(i) = x_DandR(19);
        else
            thetaDandR(i) = x_DandR(20);
        end
    else
        if D(i) <= 150
            thetaDandR(i) = x_DandR(21);
        else
            thetaDandR(i) = x_DandR(22);
        end
    end
end
net.setLinkBulkReactionCoeff(1:np, thetaDandR);
net.setOptionsPipeBulkReactionOrder(1);
quality_res_DandR = net.getComputedQualityTimeSeries;
c_nodes_DandR = quality_res_DandR.NodeQuality';
c_nodes_Sim_DandR = c_nodes_DandR(Junction_SelectIdx,2:end);
Sim_DandR_val = c_nodes_Sim_DandR(:,3*96+1:end);
mse after DandR = mse function(Sim DandR val,Obs val);
```

```
% Plot the ga-modified quality-time figure
figure,
plot(1:size(c_nodes_Obs,1),Sim_DandR_val(:,end),'-b')
hold on
scatter(1:size(c_nodes_Obs,1),Obs_val(:,end),'red','filled','square')
legend({'Simulated data','Observed data'})
xlabel('Node Index')
ylabel('Quality (mg/L)')
xticks(1:1:size(c_nodes_Obs,1));
title('Final Simulated and Observed Chlorine Concentraiton')
```



```
% Node 753 as an example
figure
plot(0.25:0.25:days*24,c_nodes_Sim(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Sim_DandR(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_Obs(1,:),'-')

legend({'Original Simulated','GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 Concentration (mg/l)','fontweight','bold')
```



fprintf('The mse before calibration is %.4f, after calibration it is %.4f', mse_before, mse_after

Test the hydrants

Find out the regulations in the node with hydrants.

```
% Here we based
c_nodes_hydrantObs = wq_data.chlorine([2,5,8],:);
Junction_HydrantIdx = net.getNodeIndex(wq_data.node_ids([2,5,8]));
c_nodes_hydrantSim = c_nodes_DR(Junction_HydrantIdx,2:end);
% Plot figure for visualization
figure
subplot(3,1,1)
plot(0.25:0.25:days*24,c_nodes_hydrantSim(1,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_hydrant0bs(1,:),'-')
legend({'GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 (mg/l)','fontweight','bold')
subplot(3,1,2)
plot(0.25:0.25:days*24,c_nodes_hydrantSim(2,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_hydrant0bs(2,:),'-')
legend({'GA-based Simulated','Observed'},'location','best');
```

```
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 (mg/l)','fontweight','bold')

subplot(3,1,3)
plot(0.25:0.25:days*24,c_nodes_hydrantSim(3,:),'-')
hold on
plot(0.25:0.25:days*24,c_nodes_hydrantObs(3,:),'-')
legend({'GA-based Simulated','Observed'},'location','best');
xlabel('Time Step (h)','fontweight','bold')
ylabel('Cl2 (mg/l)','fontweight','bold')
```

