# **EPANET WQ Tests | BWFLnet**

Prepared by: Bradley W Jenks

Date: 01 November 2021

```
clc
clear
close all
```

## **Model Setup and Initialisation**

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

```
epanet_path = 'C:\Users\bradw\OneDrive - Imperial College London\9_Software\EPANET-Matlab-Tool\
run([epanet_path,'\start_toolkit']);

EPANET-MATLAB Toolkit Paths Loaded.

net = epanet('BWFLnet.inp');

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

Load network data and setup general simulation parameters.

```
% Save element count variables
nt = net.getTimeSimulationDuration./3600; % number of time steps in hours
nn = net.NodeJunctionCount;
n0 = net.NodeReservoirCount;
ntank = net.NodeTankCount;
np = net.LinkCount; % count of pipes and valves
D = net.getLinkDiameter';
% Save node index vectors
Reservoir Idx = net.getNodeReservoirIndex;
Junction_Idx = net.getNodeJunctionIndex;
Tank_Idx = net.getNodeTankIndex;
% Initialise EPANET simulation type
net.setQualityType('Chlorine','mg/L');
net.setNodeSourceQuality(1:nn+n0+ntank,zeros(nn+n0+ntank,1)); % set all node source quality to
% Specify hydraulic and quality time steps (in seconds)
net.setTimeQualityStep(5*60); % 5 minute quality time steps
net.setTimeHydraulicStep(60*60); % 1 hour hydraulic time steps
days = 7; % simulation duration in days; change this variable to change overall simualtion time
net.setTimeSimulationDuration(days*(nt*3600)); % load modified simulation duration from initial
```

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 = 0*ones(nn+n0+ntank,nt);
net.setNodeInitialQuality(net.NodeIndex,c0);
% Assign source chlorine with time varying pattern
cext = ones(n0,days*nt);
% Constant source concentration for n0 reservoirs
% cext = cext*0.25;
% Varying source concentration by source element
source_c = [0.25, 0.35];
% for j = 1:n0
     cext(j,:) = source c(j)*cext(j,:);
% end
% Varying source concentration by source element with
% concentration spike (pulse) at t = 72 (Day 3)
t = 72;
for i = 1: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:nt*days,1:nn).';
q = 1e-3*hydraulic_res.Flow(1:nt*days,:).';

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

## **Graph Creation**

Use graph theory to plot network connectivity and spatial coordinates.

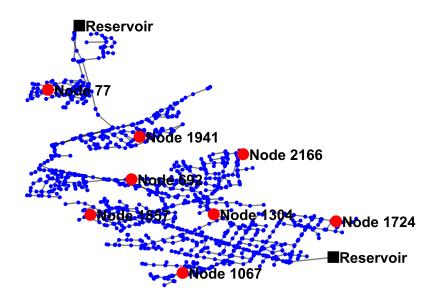
```
% Create A12 and A10 incidence matrices
LinkNodesList = net.getLinkNodesIndex;
A = zeros(np,nn+n0+ntank);
for k=1:np
    i = LinkNodesList(k,1);
    j = LinkNodesList(k,2);
    A(k,i) = -1;
    A(k,j) = 1;
end
A12 = A(:,Junction_Idx);
A10 = A(:,Reservoir_Idx);
A12 = sparse(A12);
A10 = sparse(A10);
% Obstain node XY and elevation information
XY = zeros(nn+n0+ntank,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);
```

## **Results Plotting**

Asssign results to graph nodes.

```
% Index vector for time series plot of selected junctions
Junc select = [77, 1941, 692, 2166, 1857, 1724, 1067, 1304];
for i = 1:length(Junc_select)
    Junction_Name{i} = sprintf('Node %d', Junc_select(i));
end
% Plot of network with selected junctions for visualisation
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,Reservoir_Idx,'NodeColor','k','Marker','s','MarkerSize',10);
highlight(p1,Junc_select,'NodeColor','r','MarkerSize',8);
labelnode(p1,Junc_select,Junction_Name);
labelnode(p1,Reservoir_Idx,{'Reservoir','Reservoir'});
p1.NodeFontSize = 11;
p1.NodeLabelColor = 'k';
p1.NodeFontWeight = 'bold';
axis('off')
```



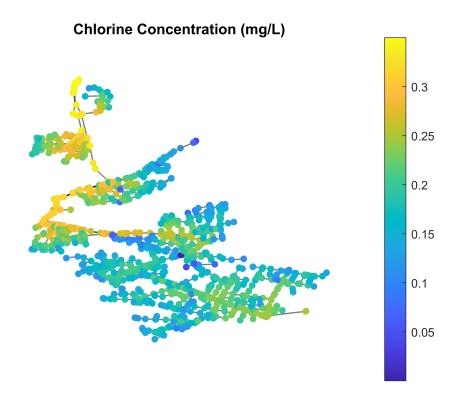
```
% Plot of Cl2 concentration distributed across network

% Select period to show results
i = 168;

figure
p2 = plot(gr);
p2.XData = XY(:,1);
p2.YData = XY(:,2);
p2.LineWidth = 1;
p2.EdgeColor = 'k';
p2.MarkerSize = 4;
p2.NodeLabel = '';

p2.NodeCData = c_nodes(:,i);

colorbar
axis('off')
title('Chlorine Concentration (mg/L)')
```



Plot time series of chlorine concentration for select junctions.

