



Stellenbosch

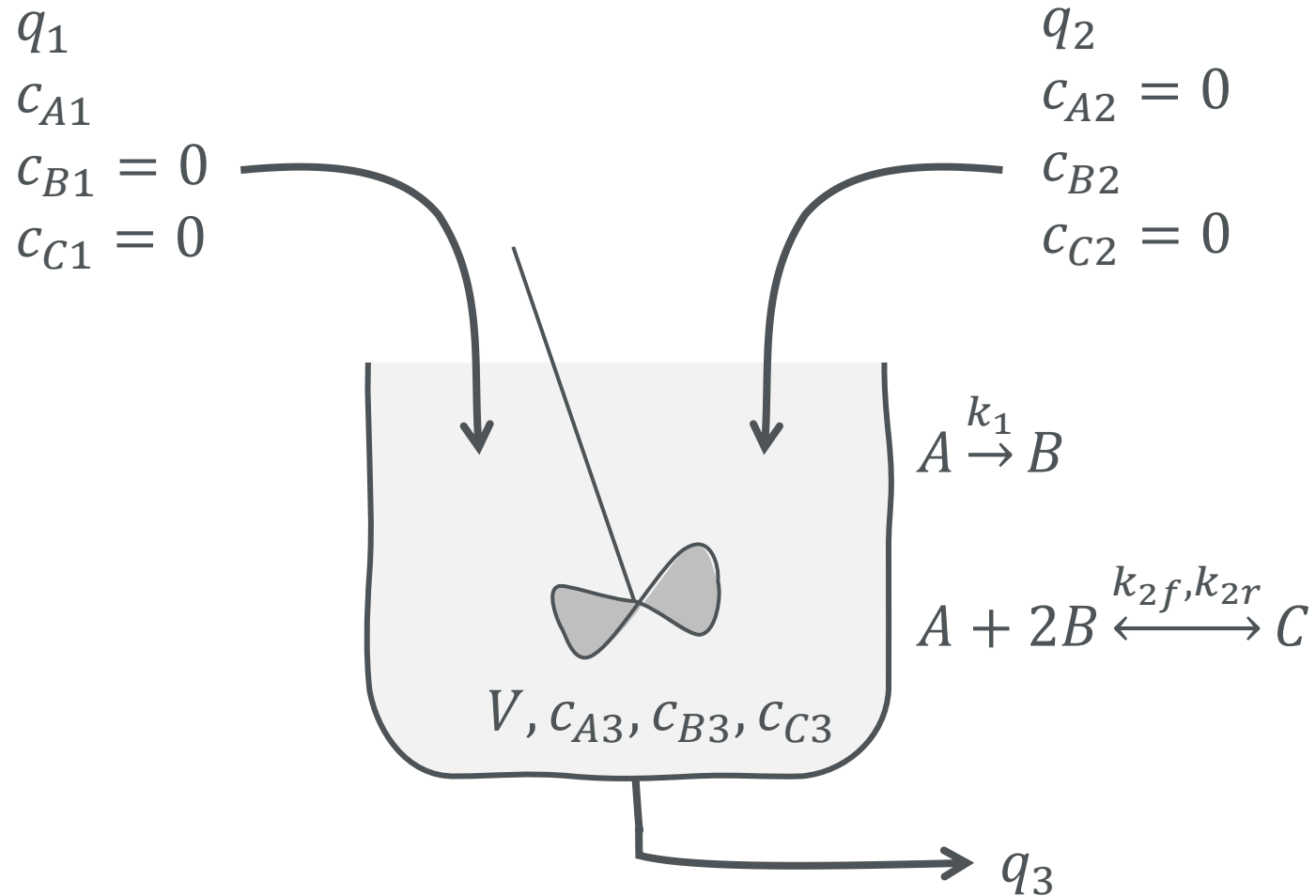
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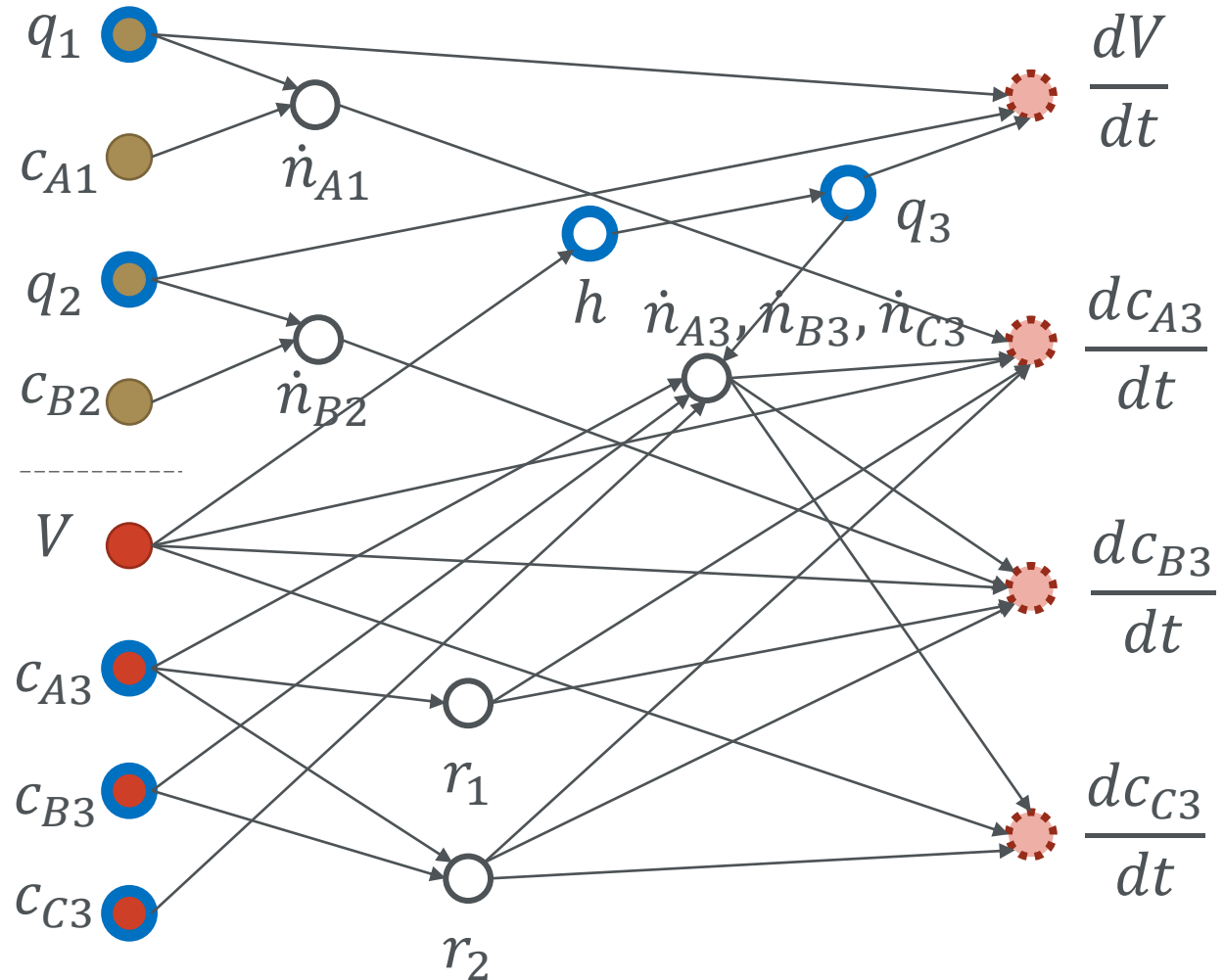
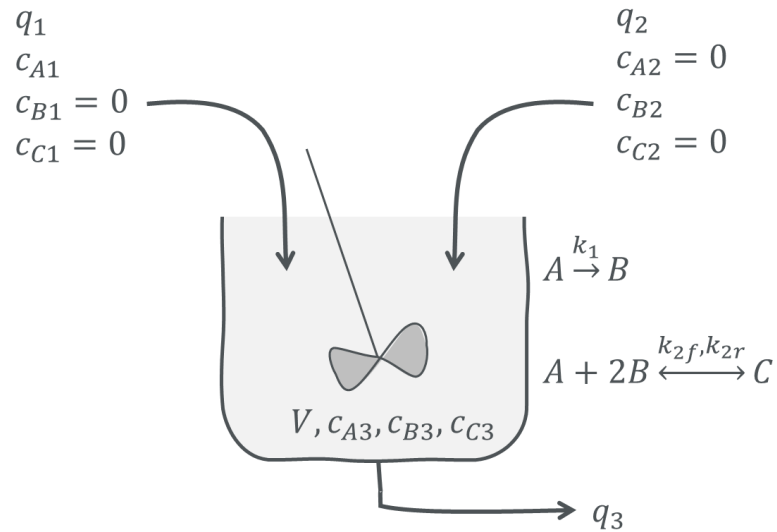
Developing and implementing systems of ODEs

Mathematical modelling and machine learning, 2nd of March 2022

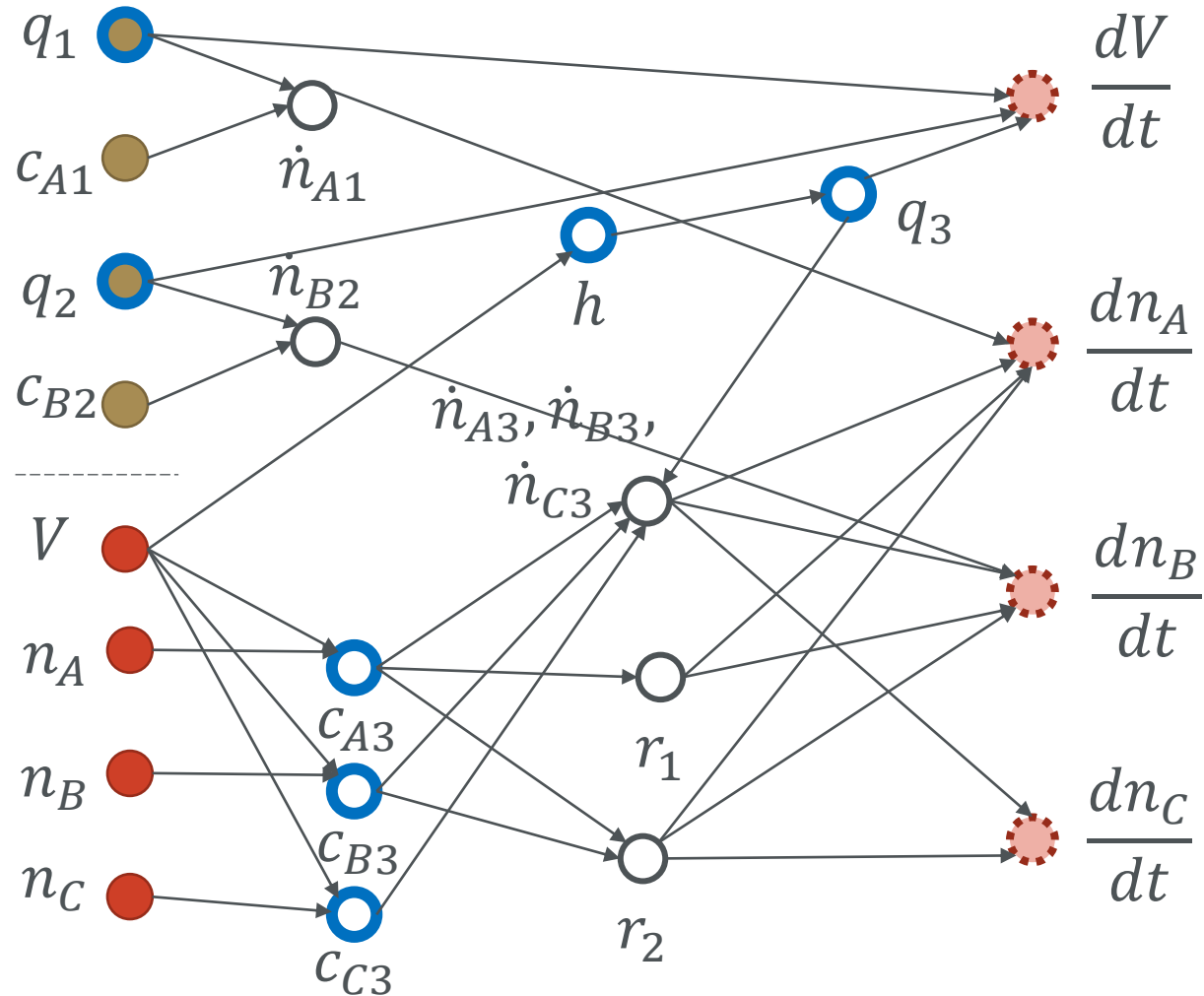
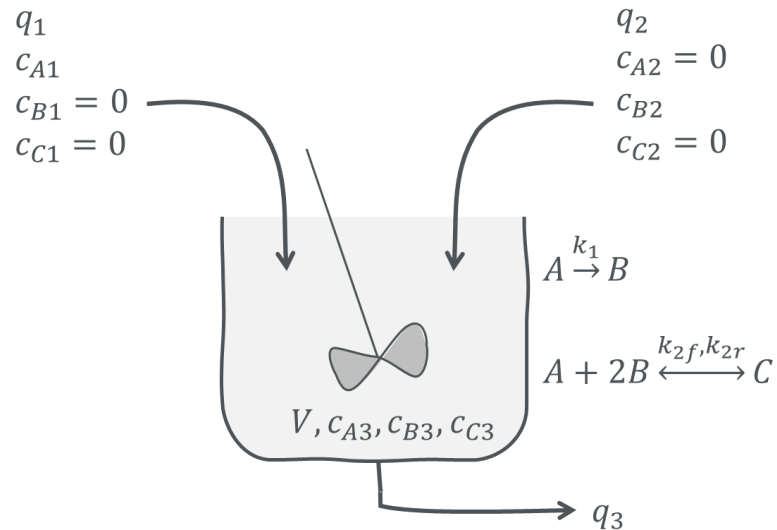
Consider the following system



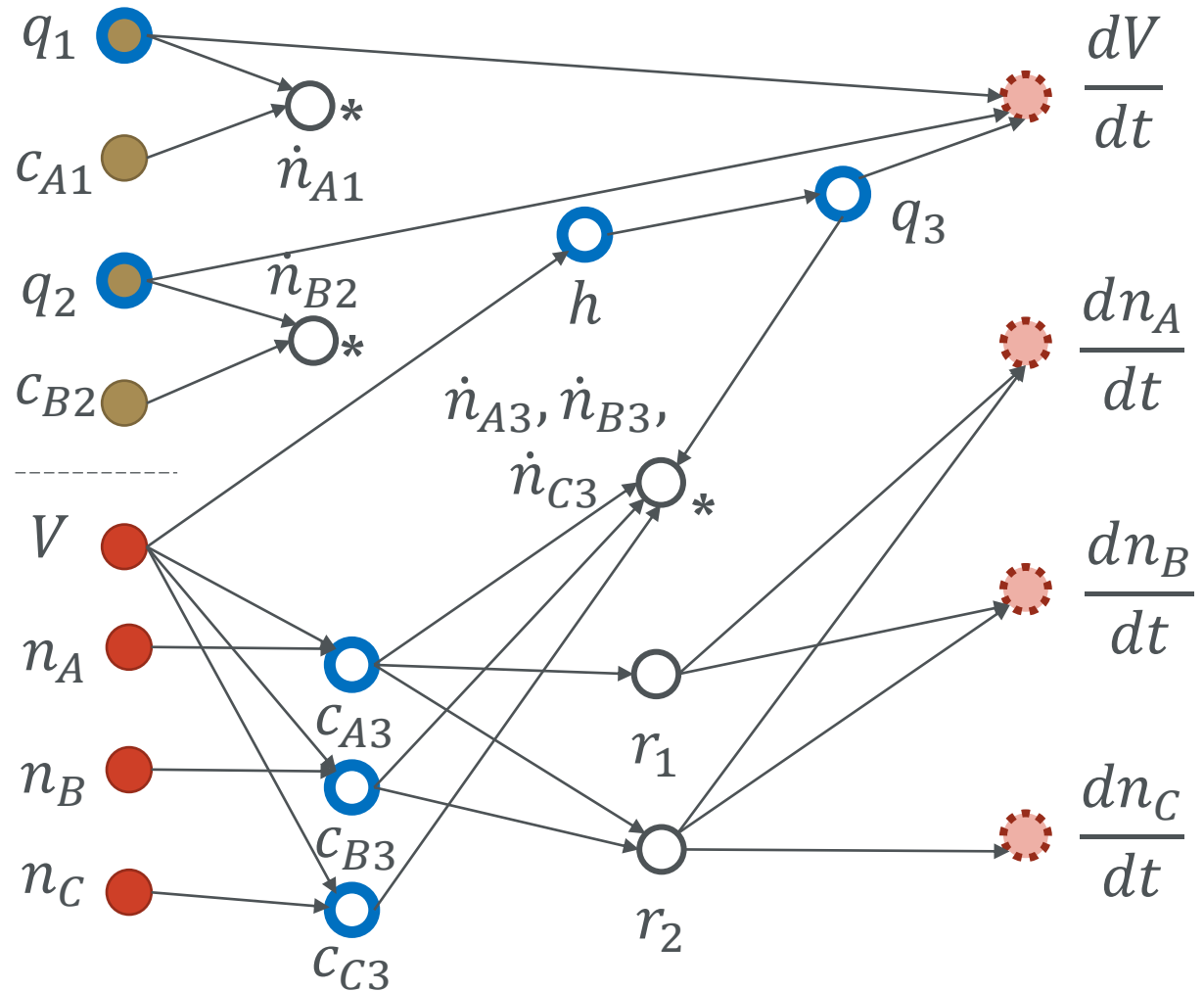
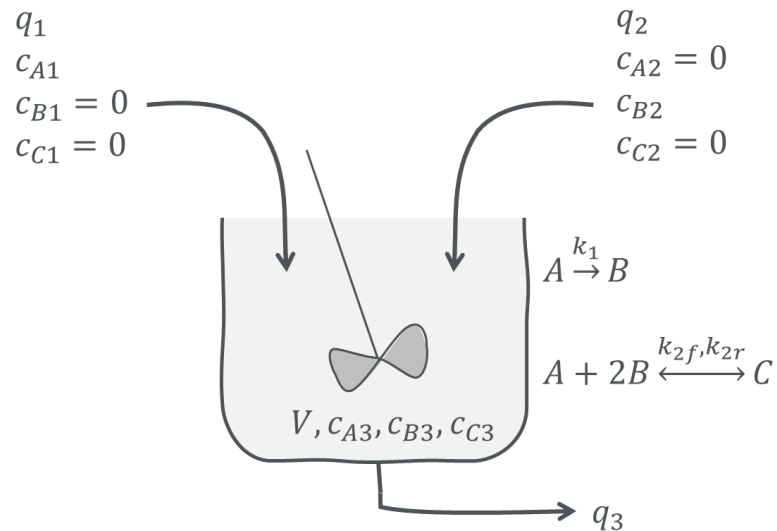
Consider the following system



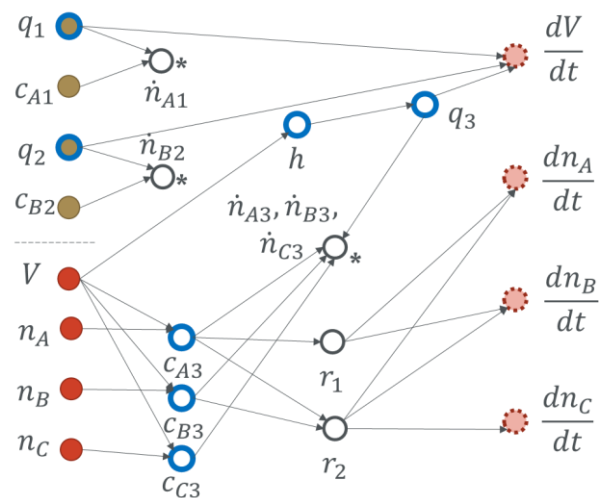
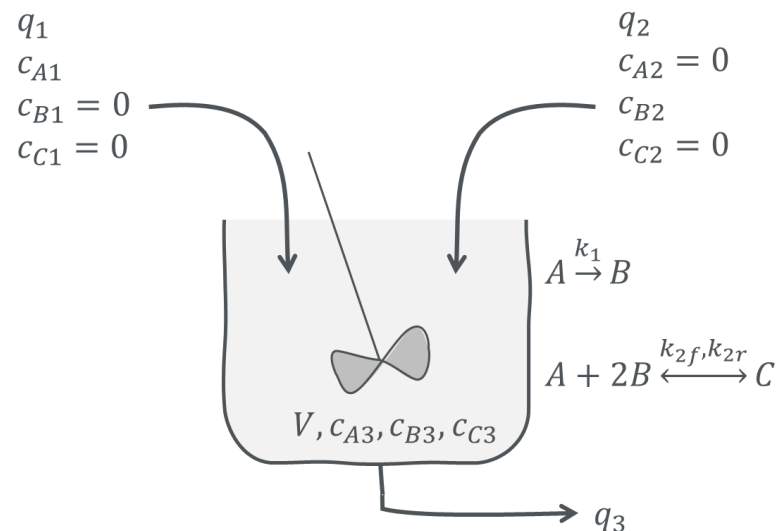
Consider the following system



Consider the following system



Consider the following system



Equation

$$\frac{dV}{dt} = q_1 + q_2 - q_3$$

$$\frac{dn_A}{dt} = \dot{n}_{A1} - \dot{n}_{A3} + v_{A1}r_1V + v_{A2}r_2V$$

$$\frac{dn_B}{dt} = \dot{n}_{B2} - \dot{n}_{B3} + v_{B1}r_1V + v_{B2}r_2V$$

$$\frac{dn_C}{dt} = -\dot{n}_{C3} + v_{C1}r_1V + v_{C2}r_2V$$

$$q_3 = c_V\sqrt{h}$$

$$h = V/A$$

$$\dot{n}_{A1} = c_{A1}q_1$$

$$\dot{n}_{B2} = c_{B2}q_2$$

$$\dot{n}_{j3} = c_{j3}q_3, \quad j = A, B, C$$

$$c_{j3} = n_{j3}/V, \quad j = A, B, C$$

$$r_1 = k_1c_{A3}$$

$$r_2 = k_{2f}c_Ac_B^2 - k_{2r}c_C$$

x

V

n_A

n_B

n_C

q_3

h

\dot{n}_{A1}

\dot{n}_{B2}

\dot{n}_{j3}

c_{j3}

r_1

r_2

u

q_1, q_2

c_{A1}

c_{B2}

p

v_{A1}, v_{A2}

v_{B1}, v_{B2}

v_{C1}, v_{C2}

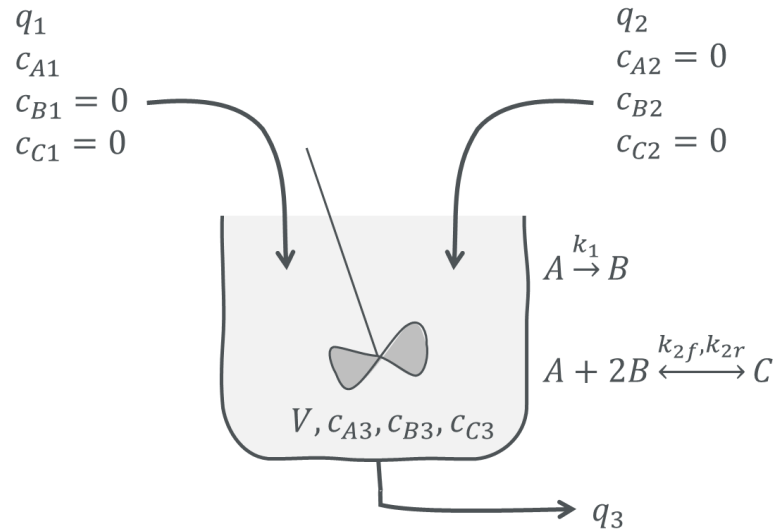
c_V

A

k_1

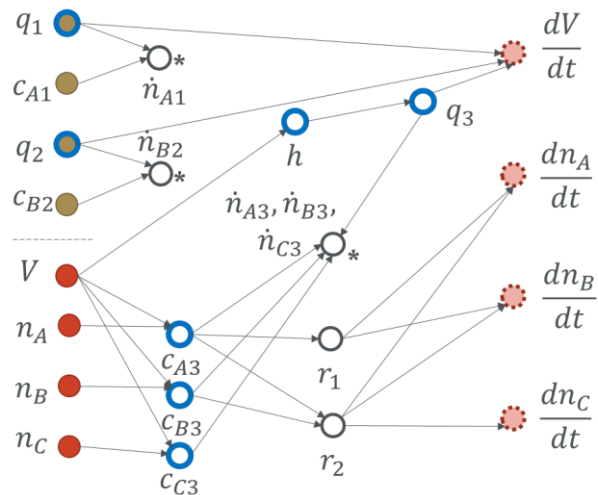
k_{2f}, k_{2r}

Consider the following system

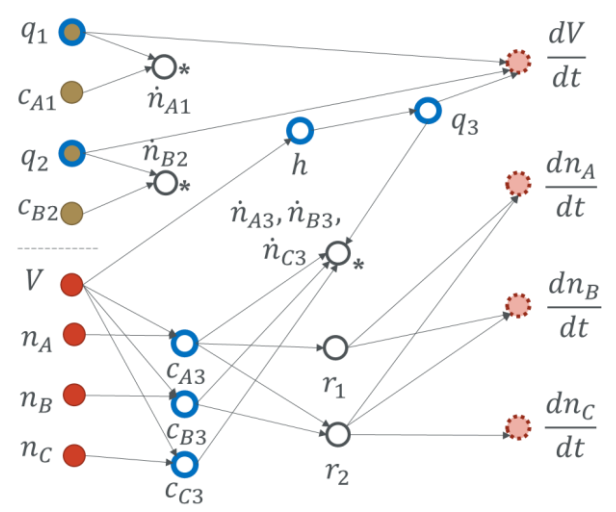
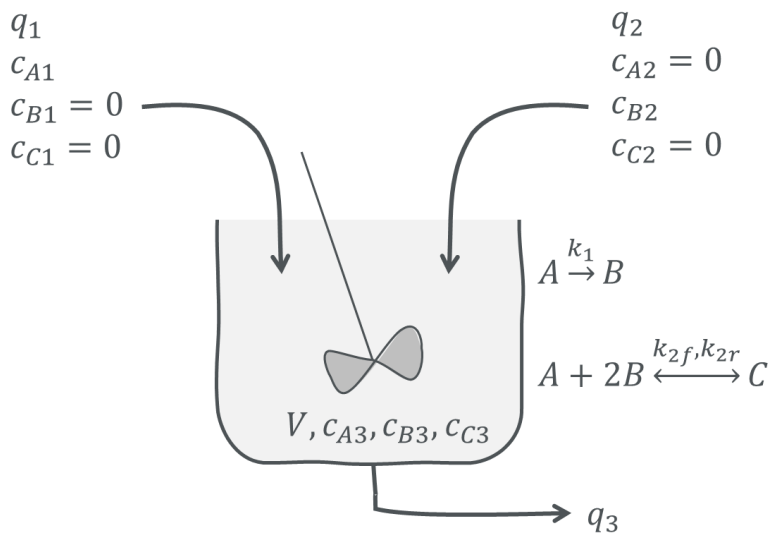


Reactions can be written as follows:

$$\begin{aligned}
 \dot{\mathbf{s}} &= \begin{bmatrix} \text{Generation|consumption of } A \\ \text{Generation|consumption of } B \\ \text{Generation|consumption of } C \end{bmatrix} \\
 &= \begin{bmatrix} \nu_{A1}r_1 + \nu_{A2}r_2 \\ \nu_{B1}r_1 + \nu_{B2}r_2 \\ \nu_{C1}r_1 + \nu_{C2}r_2 \end{bmatrix} \\
 &= \begin{bmatrix} \nu_{A1} & \nu_{B1} & \nu_{C1} \\ \nu_{A2} & \nu_{B2} & \nu_{C2} \end{bmatrix} \begin{bmatrix} r_1 \\ r_2 \end{bmatrix} \\
 &= \mathbf{N}\mathbf{r}
 \end{aligned}$$

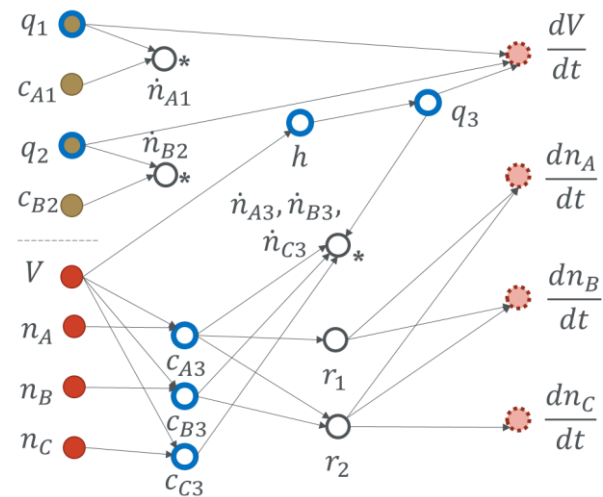
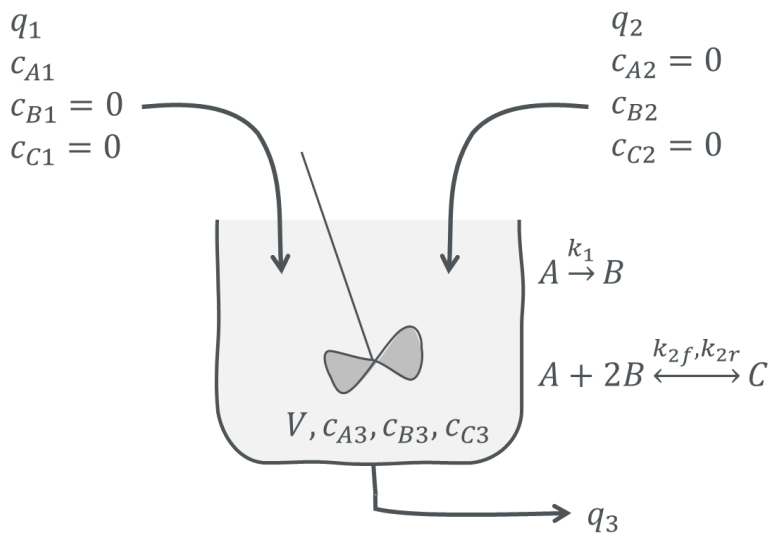


Consider the fol



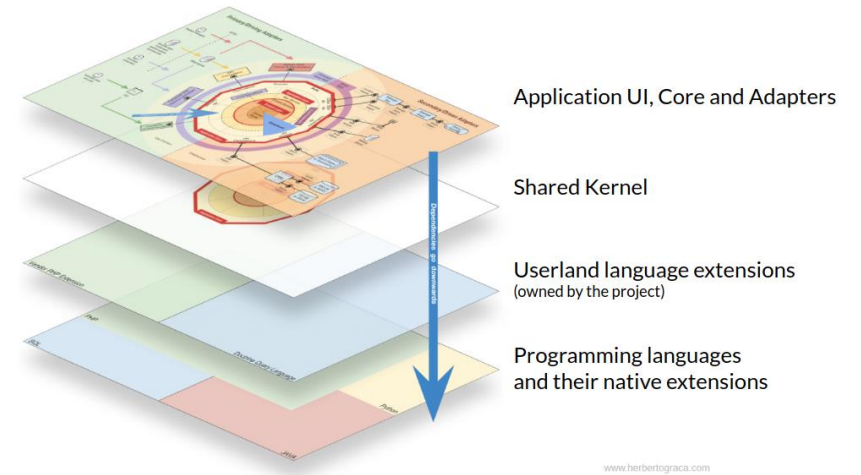
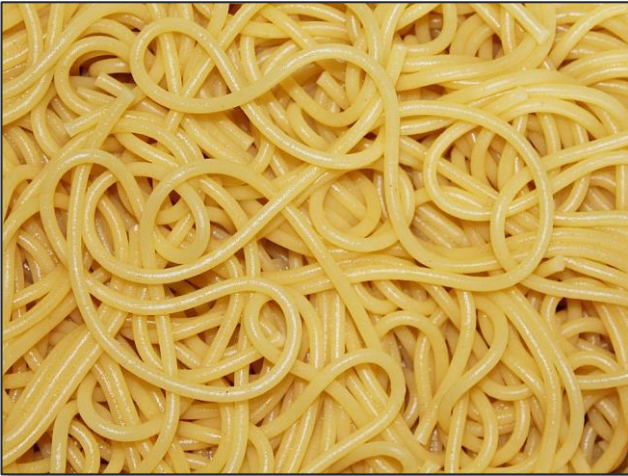
Equation	x	u	p
$\frac{dV}{dt} = q_1 + q_2 - q_3 + \dot{s}_1$	V	q_1, q_2	
$\frac{dn_A}{dt} = \dot{n}_{A1} - \dot{n}_{A3} + \dot{s}_2$	n_A		
$\frac{dn_B}{dt} = \dot{n}_{B2} - \dot{n}_{B3} + \dot{s}_3$	n_B		
$\frac{dn_C}{dt} = -\dot{n}_{C3} + \dot{s}_4$	n_C		
$q_3 = c_V \sqrt{h}$	q_3		c_V
$h = V/A$	h		A
$\dot{n}_{A1} = c_{A1} q_1$	\dot{n}_{A1}	c_{A1}	
$\dot{n}_{B2} = c_{B2} q_2$	\dot{n}_{B2}	c_{B2}	
$\dot{n}_{j3} = c_{j3} q_3, \quad j = A, B, C$	\dot{n}_{j3}		
$c_{j3} = n_{j3}/V, \quad j = A, B, C$	c_{j3}		
$r_1 = k_1 c_{A3}$	r_1		k_1
$r_2 = k_{2f} c_A c_B^2 - k_{2r} c_C$	r_2		k_{2f}, k_{2r}
$\dot{s} = N r$	\dot{s}		N

Consider the fol



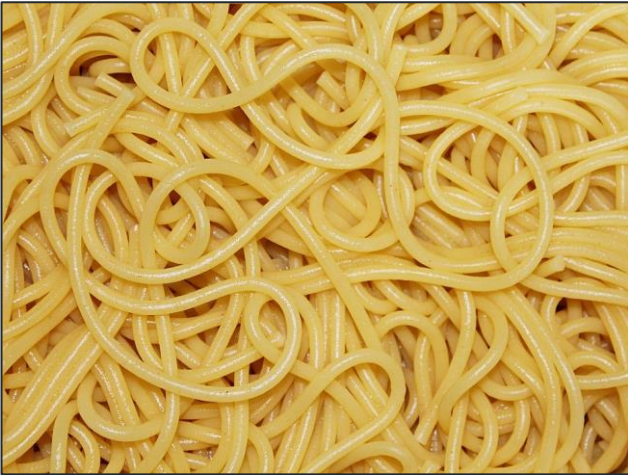
Equation	x	u	p
$\frac{dV}{dt} = q_1 + q_2 - q_3 + \dot{s}_1$	V	q_1, q_2	Measured
$\frac{dn_A}{dt} = \dot{n}_{A1} - \dot{n}_{A3} + \dot{s}_2$	n_A		
$\frac{dn_B}{dt} = \dot{n}_{B2} - \dot{n}_{B3} + \dot{s}_3$	n_B		
$\frac{dn_C}{dt} = -\dot{n}_{C3} + \dot{s}_4$	n_C		
$q_3 = c_V \sqrt{h}$	q_3	Measured	c_V
$h = V/A$	h		A
$\dot{n}_{A1} = c_{A1} q_1$	\dot{n}_{A1}	c_{A1}	
$\dot{n}_{B2} = c_{B2} q_2$	\dot{n}_{B2}	c_{B2}	
$\dot{n}_{j3} = c_{j3} q_3, \quad j = A, B, C$	\dot{n}_{j3}		
$c_{j3} = n_{j3}/V, \quad j = A, B, C$	Measured	c_{j3}	
$r_1 = k_1 c_{A3}$	r_1	Unknown	k_1
$r_2 = k_{2f} c_A c_B^2 - k_{2r} c_C$	r_2		k_{2f}, k_{2r}
$\dot{s} = \mathbf{N} \mathbf{r}$	\dot{s}		\mathbf{N}

Correct level of code complexity



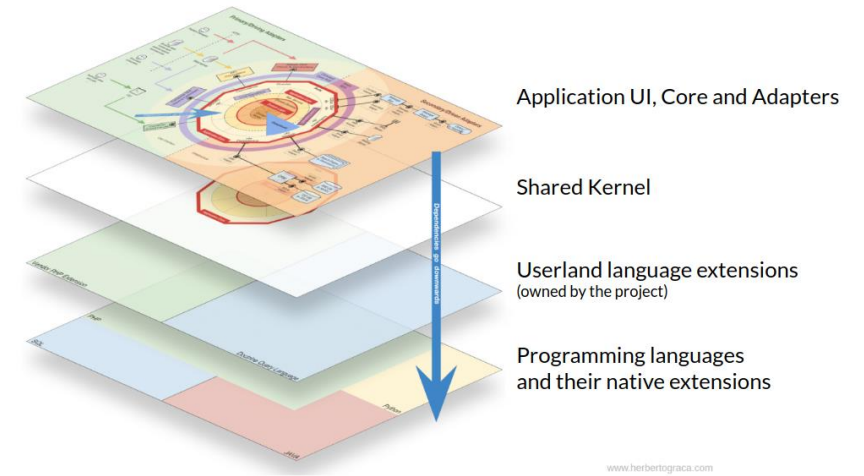
<https://herbertograca.com/2019/06/05/reflecting-architecture-and-domain-in-code/>

Correct level of code complexity



Code must be

- Easy to read
- Easy to reuse and modify



<https://herbertograca.com/2019/06/05/reflecting-architecture-and-domain-in-code/>

Implementing the model

- Create file “MAIN_System_of_ODEs.m”

Implementing the model

- Create file “MAIN_System_of_ODEs.m”
- Initialize
- Provide time vector (from 0 to 1200 s)
- Provide parameter values as structure

Implementing the model

```
%% System of ODEs: CSTR with two reactions
```

```
% Tobi Louw, 2022-03-02
```

```
% This code is used to illustrate the implementation of ODEs in MATLAB
```

```
% The example system is a CSTR facilitating two reactions, described in
```

```
% the accompanying PDF
```

```
clc
```

```
clear
```

```
clf
```

```
%% Define time region of interest
```

```
t = linspace(0, 1200); % s, Time over which to perform integration
```

```
%% Define parameters
```

```
p.cV = 0.045; % m2.5/s, Outlet flowrate coefficient
```

```
p.A =
```

Implementing the model

```
% System of ODEs: CSTR with two reactions
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t = linspace(0, 1200); % s, Time over which to perform integration
```

```
% Define parameters
```

```
p.cV = 0.045; % m2.5/s, Outlet flowrate coefficient
```

```
p.A =
```

$$\begin{aligned}c_V &= 0.045 \text{ m}^{2.5} \cdot \text{s}^{-1} \\ A &= 2 \text{ m}^2 \\ k_1 &= 0.05 \text{ s}^{-1} \\ k_{2f} &= 2.5 \text{ m}^6 \cdot \text{mol}^{-2} \cdot \text{s}^{-1} \\ k_{2r} &= 0.05 \text{ s}^{-1}\end{aligned}$$

Implementing the model

% Define parameters

```
p.cV = 0.045;    % m2.5/s, Outlet flowrate coefficient  
p.A = 2;         % m2, Cross-sectional area of reactor  
p.k1 = 5.0e-2;   % 1/s, Reaction 1 rate constant  
p.k2f = 2.5e+0;  % m6/mol2.s, Reaction 2 forward rate constant  
p.k2r = 5.0e-2;  % 1/s, Reaction 2 reverse rate constant
```

% Matrix of stoichiometric coefficients for each state variable

```
p.Nu.V  = [ 0 0];  
p.Nu.nA = [-1 -1];  
p.Nu.nB = [+1 -2];  
p.Nu.nC = [ 0 +1];
```


Implementing the model

- Create file “MAIN_System_of_ODEs.m”
- Initialize
- Provide time vector (from 0 to 1200 s)
- Provide parameter values as structure
- **Provide exogeneous inputs as functions**

Implementing the model

% Define exogeneous inputs

```
u.q1 = @(t) 0.02 + 0*t;           % m3/s, Inlet flowrate 1 (constant)
u.q2 = @(t) 0.01 + 0.05*(t > 400); % m3/s, Inlet flowrate 2 (step-change)
```

$c_{A1} = 1.50 \text{ mol.m}^{-3}$
with a step increase of 0.50 mol.m^{-3} at $t = 800 \text{ s}$

Implementing the model

% Define exogeneous inputs

```
u.q1 = @(t) 0.02 + 0*t;           % m3/s, Inlet flowrate 1 (constant)
u.q2 = @(t) 0.01 + 0.05*(t > 400); % m3/s, Inlet flowrate 2 (step-change)
u.cA1 = @(t) 1.50 + 0.50*(t > 800); % mol/m3, Inlet concentration A (step-change)
```

Implementing the model

```
% Define exogeneous inputs
```

```
u.q1 = @(t) 0.02 + 0*t;           % m3/s, Inlet flowrate 1 (constant)
u.q2 = @(t) 0.01 + 0.05*(t > 400); % m3/s, Inlet flowrate 2 (step-change)
u.cA1 = @(t) 1.50 + 0.50*(t > 800); % mol/m3, Inlet concentration A (step-change)
```

```
% Concentration cB2 will vary stochastically
```

```
y(1) = 0;
for i = 2:length(t)
    y(i) = 0.8*y(i-1) + 0.05*randn;
end
y = y + 2.0;
u.cB2 = griddedInterpolant(t, y);
```

Implementing the model

- Create file “MAIN_System_of_ODEs.m”
- Initialize
- Provide time vector (from 0 to 1200 s)
- Provide parameter values as structure
- Provide exogeneous inputs as functions
- **Define state structure and provide initial conditions**

Implementing the model

```
% Define state structure and initial conditions
```

```
p.state_fields = {'V', 'nA', 'nB', 'nC'}; % Field names for each state  
x0.V = 0.9; % m3, initial tank volume  
x0.nA = 0.15; % mol/m3, initial concentration of A  
x0.nB = 0.25; % mol/m3, initial concentration of A  
x0.nC = 0.30; % mol/m3, initial concentration of A  
x0_vec = xS2xV(x0, p.state_fields);
```

```
function xV = xS2xV(xS, fields)
```

```
% Map all elements in structure "xS" and indexed by "fields"  
% to the corresponding element in the vector "xV"  
xV(1) = xS.V;  
xV(2) = xS.nA;  
xV(3) = xS.nB;  
xV(4) = xS.nC;
```

Implementing the model

```
% Define state structure and initial conditions
```

```
p.state_fields = {'V', 'nA', 'nB', 'nC'}; % Field names for each state
x0.V = 0.9;    % m3, initial tank volume
x0.nA = 0.15;  % mol/m3, initial concentration of A
x0.nB = 0.25;  % mol/m3, initial concentration of A
x0.nC = 0.30;  % mol/m3, initial concentration of A
x0_vec = xS2xV(x0, p.state_fields);
```

```
function xV = xS2xV(xS, fields)
```

```
% Map all elements in structure "xS" and indexed by "fields"
% to the corresponding element in the vector "xV"
```

```
n = length(fields);
xV = zeros(n, length(xS.(fields{1})));
for i = 1:n
    xV(i,:) = xS.(fields{i});
end
```

Implementing the model

```
function xV = xS2xV(xS, fields)
% Map all elements in structure "xS" and indexed by "fields"
% to the corresponding element in the vector "xV"
n = length(fields);
xV = zeros(n, length(xS.(fields{1})));
for i = 1:n
    xV(i,:) = xS.(fields{i});
end
```

```
function xS = xV2xS(xV, fields)
% Maps all elements in vector "xV" to the structure "xS",
% using the elements in "fields" as field names.
n = length(fields);
for i = 1:n
    xS.(fields{i}) = xV(i,:);
end
```


Implementing the model

- Create file “MAIN_System_of_ODEs.m”
- Initialize
- Provide time vector (from 0 to 1200 s)
- Provide parameter values as structure
- Provide exogeneous inputs as functions
- Define state structure and provide initial conditions
- Create a function that calculates all intermediate variables, given time, state variables and exogeneous inputs

```

function v = CalculateIntermediates(t, x, u, p)
% Calculate intermediate process variables
% (i.e. all variables which are neither exogeneous inputs nor state variables)

% Calculate concentrations in CSTR
v.cA3 = x.nA ./ x.V;           % mol/m3, concentration of A

% Calculate all flowrates into / out of CSTR
v.h = x.V/p.A;                 % m, liquid level in CSTR
v.q3 = p.cV*sqrt(v.h);         % m3/s, flowrate out of CSTR
v.nA1 = u.q1(t).*u.cA1(t);     % mol/s, molar flowrate of A into CSTR

% Calculate reaction rates and source terms for each state variable
r(1,:) = p.k1*v.cA3;           % mol/m3.s, reaction rate 1
r(2,:) = p.k2f*v.cA3.*v.cB3.^2 - p.k2r*v.cC3; % mol/m3.s, reaction rate 2

Nu = xS2xV(p.Nu, p.state_fields); % Convert structured coefficients to vector
S_vec = Nu*r;                  % Vector of source terms

% Package vector of source terms as a structure
v.S = xV2xS(S_vec, p.state_fields);

```

```

function v = CalculateIntermediates(t, x, u, p)
% Calculate concentrations in CSTR
v.cA3 = x.nA ./ x.V; % mol/m3, concentration of A
v.cB3 = x.nB ./ x.V; % mol/m3, concentration of B
v.cC3 = x.nC ./ x.V; % mol/m3, concentration of C

% Calculate all flowrates into / out of CSTR
v.h = x.V/p.A; % m, liquid level in CSTR
v.q3 = p.cV*sqrt(v.h); % m3/s, flowrate out of CSTR
v.nA1 = u.q1(t).*u.cA1(t); % mol/s, molar flowrate of A into CSTR
v.nB2 = u.q2(t).*u.cB2(t); % mol/s, molar flowrate of B into CSTR
v.nA3 = v.q3.*v.cA3; % mol/s, molar flowrate of A out of CSTR
v.nB3 = v.q3.*v.cB3; % mol/s, molar flowrate of B out of CSTR
v.nC3 = v.q3.*v.cC3; % mol/s, molar flowrate of C out of CSTR

% Calculate reaction rates and source terms for each state variable
r(1,:) = p.k1*v.cA3; % mol/m3.s, reaction rate 1
r(2,:) = p.k2f*v.cA3.*v.cB3.^2 - p.k2r*v.cC3; % mol/m3.s, reaction rate 2

Nu = xS2xV(p.Nu, p.state_fields); % Convert structured coefficients to vector
S_vec = Nu*r; % Vector of source terms
v.S = xV2xS(S_vec, p.state_fields);

```

Implementing the model

- Create file “MAIN_System_of_ODEs.m”
- Initialize
- Provide time vector (from 0 to 1200 s)
- Provide parameter values as structure
- Provide exogeneous inputs as functions
- Define state structure and provide initial conditions
- Create a function that calculates all intermediate variables, given time, state variables and exogeneous inputs
- Create a function that calculates the derivative of the state variables, given time, state variables and exogeneous inputs

Implementing the model

```
function dxdt = SystemODEs(t, x_vec, u, p)
% Calculate the time-derivative of all state variables

% Map state vector to structure and calculate intermediate variables
x = xV2xS(x_vec, p.state_fields);
v = CalculateIntermediates(t, x, u, p);

% Calculate state derivatives as structure
ddt.V = u.q1(t) + u.q2(t) - v.q3 + v.S.V;
ddt.nA = v.nA1 - v.nA3 + v.S.nA;
ddt.nB = v.nB2 - v.nB3 + v.S.nB;
ddt.nC = - v.nC3 + v.S.nC;

% Map state derivative structure to vector
dxdt = xS2xV(ddt, p.state_fields);
```

Implementing the model

- Create file “MAIN_System_of_ODEs.m”
- Initialize
- Provide time vector (from 0 to 1200 s)
- Provide parameter values as structure
- Provide exogeneous inputs as functions
- Define state structure and provide initial conditions
- Create a function that calculates all intermediate variables, given time, state variables and exogeneous inputs
- Create a function that calculates the derivative of the state variables, given time, state variables and exogeneous inputs
- **Simulate system and plot results**

Implementing the model

```
<in MAIN_System_of_ODEs>

%% Simulate system of ODEs
[~, x_vec] = ode45(@(t, x) SystemODEs(t, x, u, p), t, x0_vec);
x = xV2xS(x_vec', p.state_fields);
v = CalculateIntermediates(t, x, u, p);

%% Plot simulation results
tiledlayout flow
ax_height = nexttile;
plot(t, v.h, 'LineWidth', 2);
legend('h', 'Location', 'northwest');

ax_concentration = nexttile;
plot(t, v.cA3, t, v.cB3, t, v.cC3, 'LineWidth', 2)
legend('c_A', 'c_B', 'c_C', 'Location', 'northwest')
```

Adding measurements

- Running simulation = running plant
- **Simulate measurement instrumentation**

Adding measurements

```
<in MAIN_System_of_ODEs>

%% Define measurement noise, frequency and delay
% Create measurement structures:
% fields: names of measurements
% var: assume gaussian noise with variance "var"
% T: measurement period T = 1/frequency
% D: measurement delay  $y \sim y(t - D)$ 
meas.fields = {'h', 'cC3'};

meas.h = struct('func', @(t, x, u, v, p) v.h, 'var', 0.1, 'T', 5, 'D', 2);
meas.cC3 = struct('func', @(t, x, u, v, p) v.cC3, 'var', 0.02, 'T', 60, 'D', 60);
```

Adding measurements

```
<in MAIN_System_of_ODEs>
```

```
%% Define measurement noise, frequency and delay
```

```
% Create measurement structures:
```

```
% fields: names of measurements
```

```
% var: assume gaussian noise with variance "var"
```

```
% T: measurement period  $T = 1/\text{frequency}$ 
```

```
% D: measurement delay  $y \sim y(t - D)$ 
```

```
meas.fields = {'h', 'cC3'};
```

```
function y = Measurements(t, x, u, v, p, meas)
```

```
mea % Calculate measurement values for each field in "meas"
```

```
mea for i = 1:length(meas.fields)
```

```
    current = meas.(meas.fields{i});
```

```
    values = current.func(t, x, u, v) + current.var*randn(size(t));
```

```
    times = 0 : current.T : t(end);
```

```
    interp_values = interp1(t, values, times);
```

```
    y.(meas.fields{i}) = timeseries(interp_values, times + current.D);
```

```
end
```

Adding measurements

```
<in MAIN_System_of_ODEs>

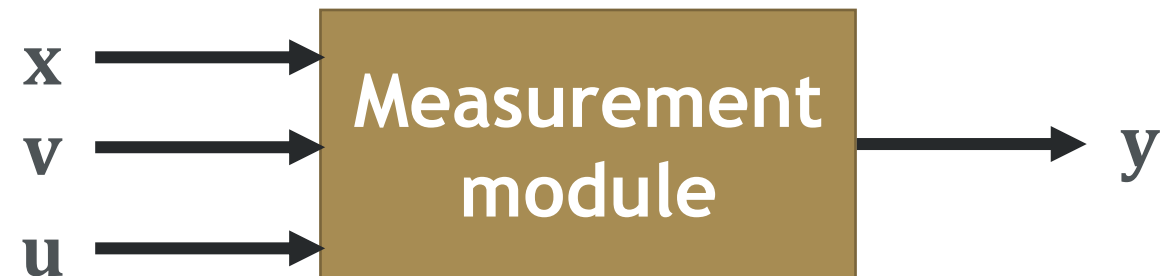
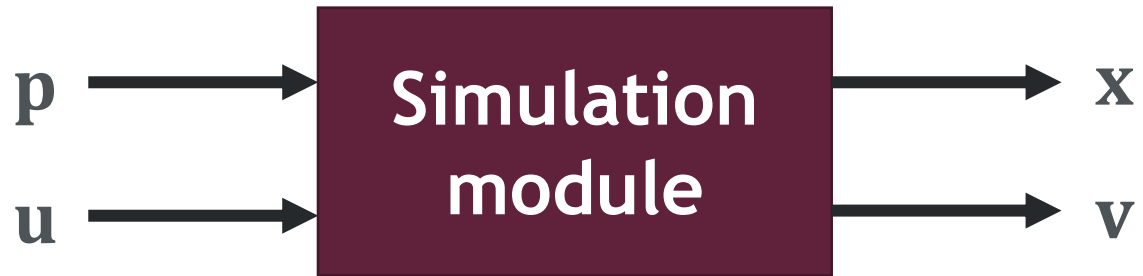
% Record measurements
y = Measurements(t, x, u, v, p, meas);

%% Plot measurements
axes(ax_height)
hold on
plot(y.h, 'k.', 'MarkerSize', 8)
legend('h', 'Location', 'northwest')

axes(ax_concentration)
hold on
plot(y.cC3, 'k.', 'MarkerSize', 20)
legend('c_C', 'Location', 'northwest')
```

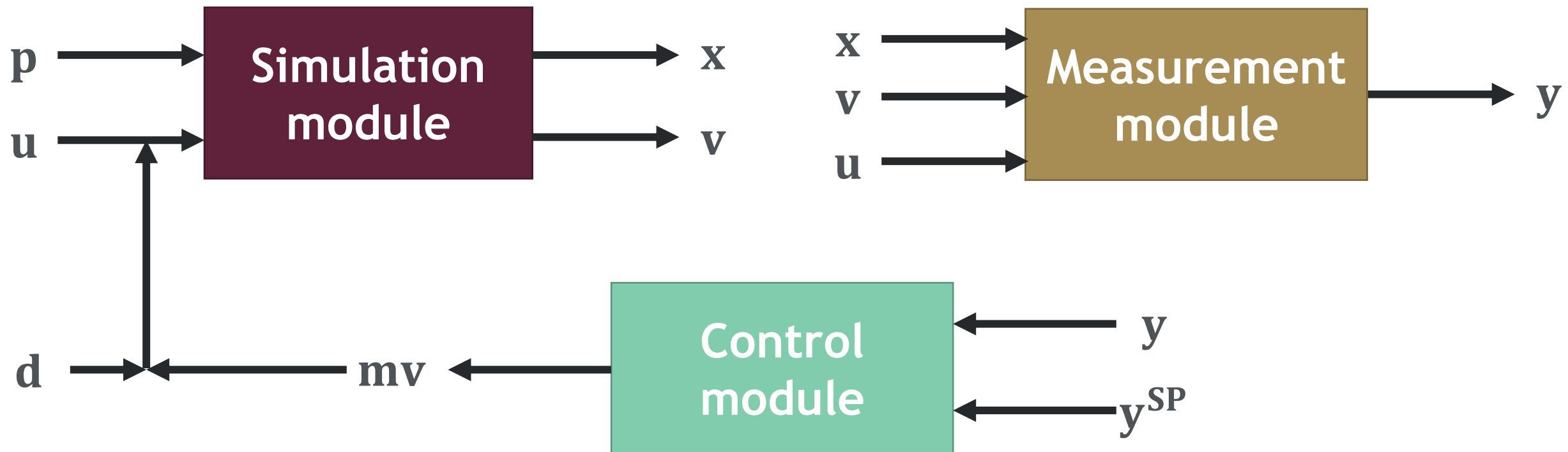
Adding measurements

- Running simulation = running plant
- Simulate measurement instrumentation



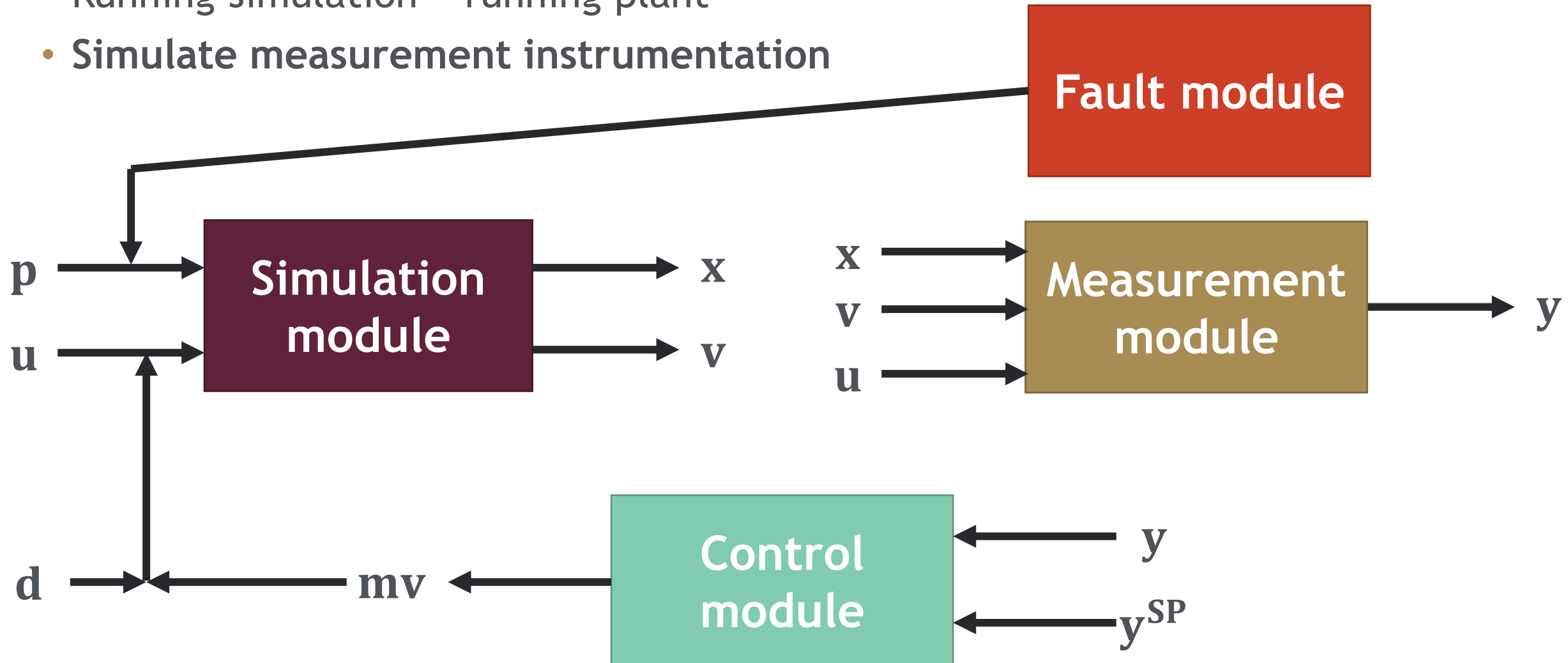
Adding measurements and control

- Running simulation = running plant
- Simulate measurement instrumentation



Adding measurements, control and faults...

- Running simulation = running plant
- Simulate measurement instrumentation



Repository

- <https://github.com/tmlouw/Introduction-to-ODEs>