

Short Tutorial on Matlab

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Part 5. Using S-function blocks in Simulink®

I. Motivation: With the complexity of medium-size to large-size nonlinear models, it may be more efficient to use a set of differential equations written in an m-file. These m-files will be accessed by Simulink through the S-function block. Thus, this method mixes the advantages of an m-file which can be run directly by solvers such as **ode45**, with the graphical links to other Simulink blocks.

II. Example System:

Suppose we want to model the nonisothermal CSTR,

$$\frac{dC_a}{dt} = \left(\frac{F}{V}\right) \cdot (C_{af} - C_a) - k_0 \cdot \exp\left[-\frac{E_a}{R \cdot (T + 460)}\right] \cdot C_a$$

$$\frac{dT}{dt} = \left(\frac{F}{V}\right) \cdot (T_f - T) - \frac{\Delta H}{\rho \cdot C_p} \cdot \left[k_0 \cdot \exp\left[-\frac{E_a}{R \cdot (T + 460)}\right] \cdot C_a \right] - \left(\frac{U \cdot A}{\rho \cdot C_p \cdot V}\right) \cdot (T - T_j)$$

We want to model this system in which we will treat the jacket temperature, T_j , as the input (i.e. manipulated variable). We will also want to monitor concentration and temperature of the liquid in the CSTR as our outputs.

III. Write the m-file.

Recall that we could model the process by writing an m-file to be used by Matlab solvers such as **ode45**. One such file, which we will name as **reactor.m**, is shown in Figure 1.

Test the model to make sure it works. For instance, with $T_j=55$:

```
>> [t,x]=ode45(@reactor,[0 10],[0.1;40],[ ],55);
```

Note/Recall: The command-line specifies: a simulation-time span of **[0 10]**, an initial-value column vector: **[0.1;40]**, a null placeholder, **[]**, for default options, and setting T_j with a value equal to 55.

```

function    dx = reactor(t,x,Tj)
%
%    model for reactor
%
Ca = x(1)          ;    % lbmol/ft^3
T  = x(2)          ;    % oF

Ea = 32400         ;    % BTU/lbmol
k0 = 15e12         ;    % hr^-1
dH = -45000        ;    % BTU/lbmol

U  = 75            ;    % BTU/hr-ft^2-oF
rhocp = 53.25      ;    % BTU/ft^3
R  = 1.987         ;    % BTU/lbmol-oF
V  = 750           ;    % ft^3
F  = 3000          ;    % ft^3/hr
Caf = 0.132        ;    % lbmol/ft^3
Tf = 60            ;    % oF

A = 1221           ;    % ft^2

ra = k0*exp(-Ea/(R*(T+460)))*Ca;
dCa = (F/V)*(Caf-Ca)-ra;
dT  = (F/V)*(Tf-T)-(dH)/(rhocp)*ra...
      -(U*A)/(rhocp*V)*(T-Tj);

dx = [dCa;dT];

```

Figure 1. File saved as **reactor.m**

Remarks:

1. We treat T_j as an argument/parameter. This is in anticipation that we will be varying T_j later as an input/manipulated variable.
2. The arguments \mathbf{x} and \mathbf{dx} are column vectors for state and derivative, respectively.
3. Writing a model first for direct ODE45 implementation is advisable, specially for complex processes. This way, one can check the validity of the model, prior to its incorporation to a Simulink model.

IV. Write an S-function file.

This file will also be saved as an m-file. It contains the protocol in which Simulink can access information from Matlab.

For our example, we show one such S-function file in Figure 2. We will save this file as **reactor_sfcn.m**.

```

function [sys,x0,str,ts]=...
    reactor_sfcn(t,x,u,flag,Cinit,Tinit)

```

```

switch flag

    case 0 % initialize

        str=[] ;
        ts = [0 0] ;

        s = simsizes ;

            s.NumContStates = 2 ;
            s.NumDiscStates = 0 ;
            s.NumOutputs = 2 ;
            s.NumInputs = 1 ;
            s.DirFeedthrough = 0 ;
            s.NumSampleTimes = 1 ;

        sys = simsizes(s) ;

        x0 = [Cinit, Tinit] ;

    case 1 % derivatives

        Tj = u ;
        sys = reactor(t,x,Tj) ;

    case 3 % output

        sys = x;

    case {2 4 9} % 2:discrete
                % 4:calcTimeHit
                % 9:termination

        sys =[];

    otherwise

        error(['unhandled flag =',num2str(flag)]) ;

end

```

Figure 2. File saved as **reactor_sfcn.m**.

Let us deconstruct the S-function file given in Figure 2 to understand what the file needs to contain.

1. The first line specifies the input and output arguments.

```

function [sys,x0,str,ts]=...
        reactor_sfcn(t,x,u,flag,Cinit,Tinit)

```

As it is with any Matlab functions, the variable names themselves are not as crucial as the positions of the variables in the list.

a) input arguments

- (1) **t** - the time variable
- (2) **x** - the column-vector of state variables
- (3) **u** - the column-vector of input variables (whose value will come from other Simulink blocks)
- (4) **flag** - indicator of which group of information and/or calculations is being requested by Simulink.

There are six types of request that Simulink performs, each of which is designated by an integer number:

flag value	Job/Data Request
0	<u>Initialization:</u> a) Setup of input/output vector sizes and other setup modes b) Specification/calculation of initial conditions for the state variables.
1	<u>Derivative Equation Updating:</u> a) Calculations involving input vectors b) Calculation of the derivatives
2	<u>Discrete Equation Updating</u> (will not be used for our example)
3	<u>Output Calculations:</u> Evaluating output variables as a function of the elements of the state vector (and in some case, also the elements of the input vector)
4	<u>Get Time of Next Variable Hit</u> (will not be used for our example)
9	<u>Termination:</u> Additional routines/calculations at the end of the simulation run. (will not be used for our example)

- (5) **Cinit, Tinit** - additional supplied parameters.

In our case, these are the initial conditions for concentration and temperature.

Note: We do not specify what the values of the input arguments are. Their values will be specified by Simulink during a simulation run.

b) output arguments

- (1) **sys** - the main vector of results requested by Simulink. Depending on the **flag** sent by Simulink, this vector will hold different information.

If flag = 0	sys = [a,b,c,d,e,f,g] where, a = number of continuous time states b = number of discrete time states c = number of outputs (<i>Note: this is not necessarily the number of states</i>) d = number of inputs e = 0 (<i>required to be 0, not currently used</i>) f = 0(no) or 1(yes) for direct algebraic feed through of input to output. (<i>this is relevant only if during flag=3, the output variables depend algebraically on the input variables.</i>) g = number of sample times. (<i>for continuous process, we set this equal to 1</i>)
If flag = 1	sys = a column vector of the derivatives of the state variables
If flag = 3	sys = a column vector of the output variables
If flag = 2,4,9	since these flags are not used in our example, they can just send out a null vector: sys =[]

The next set of 3 output arguments are needed by Simulink only when **flag** = 0, otherwise they are ignored:

- (2) **x0** - column vector of initial conditions.
- (3) **str** - need to be set to null. This is reserved for use in future versions of Simulink.
- (4) **ts** - an array of two columns to specify sampling time and time offsets. Since our example will deal only with continuous systems, this will be set to [0 0] during initiation phase.

2. After the first line, the S-function file is split into the different cases determined by **flag**. As shown in Figure 3, we show the bare structure of the “containers” for the different cases. We have left out the details for case 1, 2 and 3. For case 2, 4, and 9, we simply set **sys=[]**. The last two lines to catch an exceptional case where a bug occurs during the Simulink run.

```

switch flag
    case 0                % initialize
        %...
    case 1                % derivatives
        %...
    case 3                % output
        %...
    case {2 4 9}         % 2:discrete
                        % 4:calcTimeHit
                        % 9:termination
        sys =[];
    otherwise
        error(['unhandled flag =',num2str(flag)]) ;
end

```

Figure 3.

Now, let us fill the details.

For case 0 (initialization),

- a) define **str**, **ts** and **x0**

```

str=[] ;
ts = [0 0] ;
x0 = [Cinit, Tinit] ;

```

- b) create a row vector which specifies the number of inputs and outputs, etc. To aid in this, we invoke the **simsizes** command. Without arguments, **simsizes** will creates a strucure variable which we can then fill with the required values:

```

s = simsizes ;

```

```

s.NumContStates = 2 ;
s.NumDiscStates = 0 ;
s.NumOutputs    = 2 ;
s.NumInputs     = 1 ;
s.DirFeedthrough = 0 ;
s.NumSampleTimes = 1 ;

```

Using the command **simsizes** again with the structure variable as the argument actually translates the values in the structure, **s**, into a row vector which gets sent to Simulink via **sys**:

```

sys = simsizes(s) ;

```

For case 1 (derivative calculations)

We set the input **u** to **Tj** and then apply it to the m-file we wrote earlier, i.e. **reactor.m**:

```

case 1 % derivatives
Tj = u ;
sys = reactor(t,x,Tj) ;

```

For case 3 (output calculations)

```

case 3 % output
sys = x;

```

V. Insert the S-Function block into the Simulink.

In the Simulink Library browser, go to the **[User-Define Functions]** subdirectory. Then drag-drop the **S-Function** block (see Figure 4).

Double-click on the S-function block and fill in the parameters. Change the S-function name to **reactor_sfcn**. Also, fill in the parameters. In our case, we input **0.1,40** (which is the value for **Cinit** and **Tinit**) as shown in Figure 5.

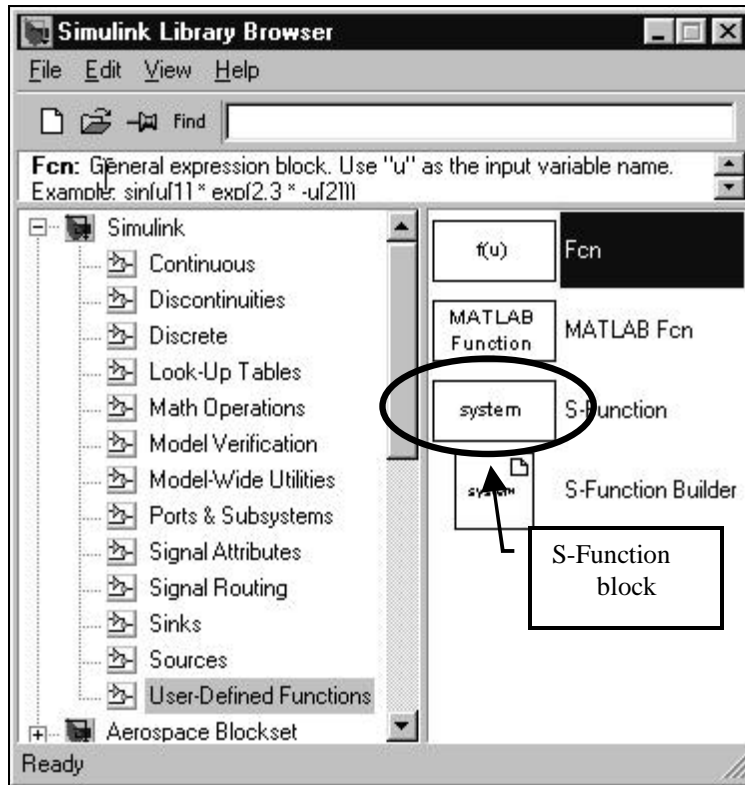


Figure 4.

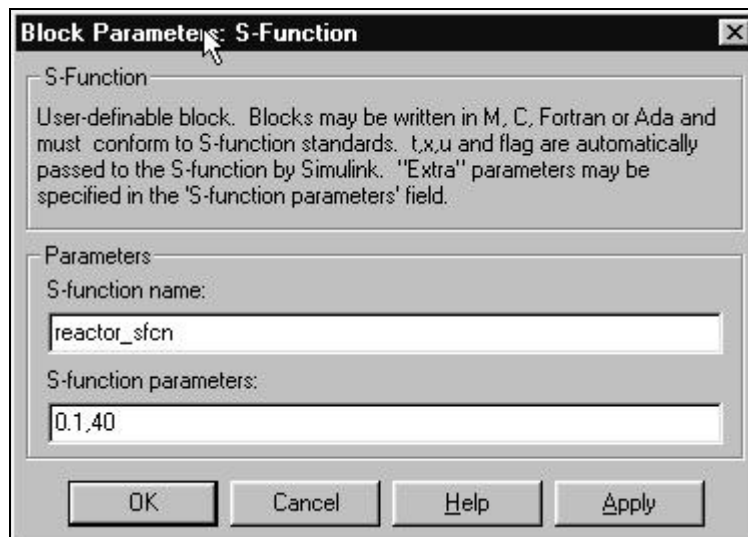


Figure 5.

VI. Add other Simulink blocks and simulate.

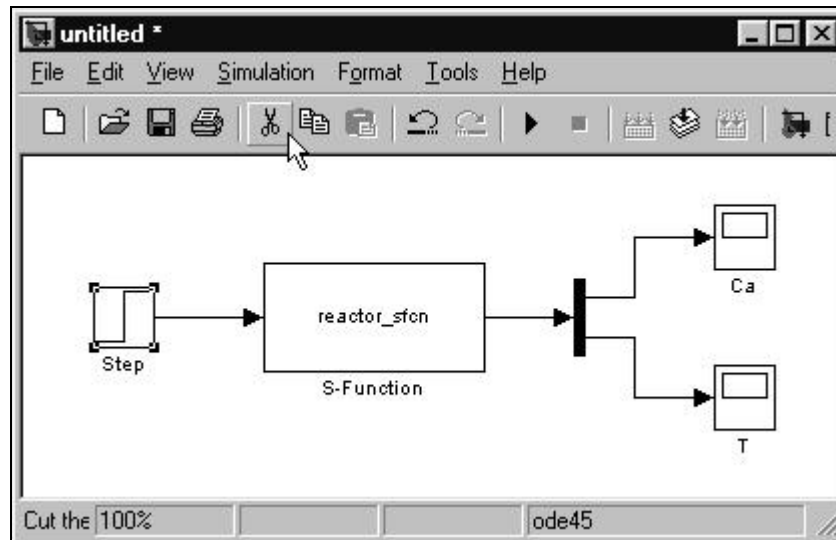


Figure 6.

Remark: In figure 6, we include a **demux** block (which stands for demultiplexer) to split the output vector to the 2 elements. In other applications where the input vectors has more than one element, we need a **mux** block (which stands for multiplexer). Both **mux** and **demux** blocks reside in the **Signal Routing** subdirectory of the Simulink Library browser.