

TR_1D_model1_SS\stack_state

TR_1D_model1_SS\stack_state.m

```

% TR_1D_model1_SS\stack_state.m
%
% function [x_state,iflag] = stack_state(State, ...
%   num_species,num_pts);
%
% This procedure stacks the concentration and temperature
% profiles into a single master array for use with the
% general solver.
%
% INPUT :
% =====
% State          see TR_1D_model1_SS.m for details
%                data structure containing concentration
%                and temperature profiles
% num_species    INT
%                the number of species in the system
% num_pts        INT
%                the number of grid points
%
% OUTPUT :
% =====
% x_state        REAL(num_DOF= (num_species+1)*num_pts)
%                this is the 1-D master array of state
%                variables
%
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% Department of Chemical Engineering
% 7/2/2001
%
% Version as of 7/25/2001

```

```

function [x_state,iflag] = stack_state(State, ...
    num_species,num_pts);

```

```

iflag = 0;

```

```

func_name = 'stack_state';

```

```

% This flag controls what to do in case of an
% assertion failure. See the assertion routines
% for further details.

```

```

i_error = 2;

```

```
% check for errors in input parameters
```

```
% check dimensions
```

```
assert_scalar(1,num_species,'num_species', ...  
    func_name,1,1,1);  
assert_scalar(1,num_pts,'num_pts',func_name,1,1,1);
```

```
% check if State is proper structure type
```

```
StateType.num_fields = 2;  
% .conc  
ifield = 1;  
FieldType.name = 'conc';  
FieldType.is_numeric = 1;  
FieldType.num_rows = num_pts;  
FieldType.num_columns = num_species;  
FieldType.check_real = 1;  
FieldType.check_sign = 2;  
FieldType.check_int = 0;  
StateType.field(ifield) = FieldType;  
% .Temp  
ifield = 2;  
FieldType.name = 'Temp';  
FieldType.is_numeric = 1;  
FieldType.num_rows = num_pts;  
FieldType.num_columns = 1;  
FieldType.check_real = 1;  
FieldType.check_sign = 1;  
FieldType.check_int = 0;  
StateType.field(ifield) = FieldType;  
% call assertion routine for structure  
assert_structure(i_error,State,'State', ...  
    func_name,StateType);
```

```
% allocate x_state column vector and
```

```
% initialize to zeros
```

```
num_DOF = (num_species+1)*num_pts;  
x_state = linspace(0,0,num_DOF);
```

```
%PDL> First, we stack the concentration profiles
```

```
%PDL> Set pos_counter to zero
```

```
pos_counter = 0;
```

```
%PDL> FOR ispecies FROM 1 TO ProbDim.num_species
```

```
for ispecies = 1:num_species
```

```
% PDL> x_state(pos_counter+1:pos_counter+num_pts) =  
%   state_data:conc(:,ispecies)
```

```
   x_state(pos_counter+1:pos_counter+num_pts) = ...  
     State.conc(:,ispecies);
```

```
% PDL> Increment pos_counter by num_pts
```

```
   pos_counter = pos_counter + num_pts;
```

```
%PDL> ENDFOR
```

```
end
```

```
%PDL> Next, we stack the temperature profile
```

```
%PDL> x_state(pos_counter+1:pos_counter+num_pts) =  
%   state_data:Temp
```

```
   x_state(pos_counter+1:pos_counter+num_pts) = ...  
     State.Temp;
```

```
iflag = 1;
```

```
return;
```