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# The Dominant Pole Design Toolbox

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<i>Title and subtitle</i> The Dominant Pole Design Toolbox		
<i>Abstract</i> <p>This report describes the Matlab routines that were used in the thesis Persson, P.: "Towards Autonomous PID Control." The routines are used to compute PID controllers based on placement of a few dominant poles. The user must specify a transfer function of the plant to be controlled and the desired degree of stability of the closed loop system. The stability is expressed as the maximum of the sensitivity function of the plant. From these specifications are the PID controller parameters computed, by minimizing the cost functional <math>IE = \int_0^{\infty} e(t) dt</math>.</p> <p>The routines can also be used for computing frequency responses, amplitude and phase margins, sensitivity functions, etc.</p> <p>A method of connecting Matlab and Simnon is also briefly described. By this connection, Simnon can be used as a computation engine for Matlab. Simnon commands can be given from Matlab, parameters can be set, values in Simnon can be retrieved to Matlab, and simulation results can be taken from Simnon into Matlab.</p>		
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# 1. Introduction

Ever since the publication of the famous articles [Ziegler and Nichols, 1942] and [Ziegler and Nichols, 1943] there has been an interest in systematic tuning methods for PID controllers. This interest has increased recently due to research on automatic tuning of controllers. See, e.g., the article [Åström and Hägglund, 1984] and the book [Åström and Hägglund, 1988]. Several controllers with automatic tuning are commercially available.

The PhD thesis [Persson, 1992b] describes methods for the tuning of PID controllers based on pole placement. During the work with the thesis a number of routines for computation of PID controllers were written. This report presents an improved version of these routines as a Matlab toolbox. For readings on Matlab, see [MathWorks, 1990]. The toolbox will be called the DPD Toolbox (the Dominant Pole Design Toolbox). This report is intended to be a manual for the DPD Toolbox, but also to further illustrate the computations carried out in [Persson, 1992b].

The methods require that the transfer function of the process is known. The toolbox contains design routines and also some utility functions of general interest. Apart from using the routines in the toolbox for computing controllers there are also routines for computing frequency responses, amplitude and phase margins, sensitivity functions, etc.

This report also describes a way to connect Matlab and Simnon. By doing this Simnon can be used as a computing engine for Matlab. There are routines for giving Simnon commands from Matlab, to set Simnon parameters from matlab, to retrieve Simnon values to Matlab, and to get time signals from Simnon to Matlab.

## 2. Design Methods

This chapter will give a brief review of the results and methods presented in the thesis [Persson, 1992b]. We also give references to the implementation of the different methods.

### Design of PID controllers

All methods that are presented here are based on the placement of a few (two or three) of the dominant poles of the closed loop system. We will use the parallel form of the PID controller

$$U(s) = k(\beta Y_r(s) - Y(s) + \frac{1}{sT_i}(Y_r(s) - Y(s)) - \frac{sT_d}{1 + s\frac{T_d}{N}}Y(s)), \quad (2.1)$$

where  $u(t)$  is the controller output,  $y(t)$  is the plant output,  $y_r(t)$  is the reference signal,  $\beta$  is the set point weighting factor,  $N$  is a filter constant, and  $k$ ,  $T_i$ ,  $T_d$  are the controller parameters. The parameter  $\beta$  will affect a zero of the closed loop system, but does not affect the stability of the system. Notice that the derivative only acts on the output from the process,  $y(t)$ . The parameters  $\beta$  and  $N$  are not determined by pole placement. For computation of the poles of the system the transfer function

$$G_c(s) = k(1 + \frac{1}{sT_i} + sT_d) = k + \frac{k_i}{s} + sk_d \quad (2.2)$$

will be used. The transfer function of the plant will be denoted  $G_p(s)$ , and is assumed to be known.

We will require that the characteristic equation,  $1 + G_c(s)G_p(s) = 0$ , has roots in specified locations. For controllers with two parameters we specify poles in the locations  $p_{1,2} = \omega_0(-\zeta_0 \pm i\sqrt{1 - \zeta_0^2})$  and for controllers with three parameters we specify poles in the locations  $p_{1,2} = \omega_0(-\zeta_0 \pm i\sqrt{1 - \zeta_0^2})$  and  $p_3 = -\alpha_0\omega_0$ . The equations  $1 + G_c(p_i)G_p(p_i) = 0$  are then solved with respect to the controller parameters. The controller parameters will be functions of  $\omega_0$ ,  $\zeta_0$ , and  $\alpha_0$ . The design routines will accept  $\zeta_0 > 1$  which corresponds to two poles on the negative real axis.

The control error is defined as  $e(t) = y_r(t) - y(t)$ . It is easy to show that

$$\text{IE} = \int_0^\infty e(\tau) d\tau = \frac{1}{k_i}, \quad (2.3)$$

when the input to the closed system is a step on the input of the plant. The integrated error, IE, can be used as an optimization criterion. This is advantageous because we have a simple analytical expression of the cost functional. Since the distance of the poles from the origin is closely coupled to the performance of the system we have chosen to determine  $\omega_0$  by maximizing  $k_i$ .

For a well damped closed system  $\text{IE} \approx \text{IAE} = \int_0^\infty |e(\tau)| d\tau$ , which is the conventional optimization criterion.

To determine a suitable value of  $\zeta_0$  we will specify a value of the maximum of the sensitivity function  $M_s$ , defined as

$$M_s = \max_{\omega > 0} \frac{1}{|1 + G_c(i\omega)G_p(i\omega)|}$$

This way of determining  $\zeta_0$  gives good control of the shape of the closed loop step responses. A small value of  $M_s$  will give a well damped system and a large value will give an oscillatory system. Normally a value in the interval 1.5...2.0 is recommended. It is not recommended to choose a value of  $\zeta_0$  directly, since the result of design with  $\zeta_0$  directly will vary much depending on the process.

The parameter  $\alpha_0$  must be chosen for the design of PID controllers. Normally  $\alpha_0 = 1$  is a good choice. For resonant processes it is necessary to choose  $\alpha_0$  smaller,  $\alpha_0 \approx 0.2 \dots 0.5$ .

**Design of PI controllers** In case of a PI controller the formulas for  $k$  and  $k_i$  become

$$k = -\frac{\sqrt{1 - \zeta_0^2}A + \zeta_0 B}{\sqrt{1 - \zeta_0^2}(A^2 + B^2)} \quad (2.4)$$

$$k_i = -\frac{\omega_0 B}{\sqrt{1 - \zeta_0^2}(A^2 + B^2)}, \quad (2.5)$$

where  $A = \text{Re } G_p(\omega_0(-\zeta_0 + i\sqrt{1 - \zeta_0^2}))$  and  $B = \text{Im } G_p(\omega_0(-\zeta_0 + i\sqrt{1 - \zeta_0^2}))$ . The expressions for  $\zeta_0 > 1$  look similar, but require that  $G_p(s)$  is evaluated twice. The functions  $k(\omega_0, \zeta_0)$  and  $k_i(\omega_0, \zeta_0)$  are implemented in the function 'dppi'. The recommended design method is to choose a  $\zeta_0$ , then maximize  $k_i(\omega_0, \zeta_0)$  with respect to  $\omega_0$ , giving  $k^*(\zeta_0)$  and  $k_i^*(\zeta_0)$ . The parameter  $\zeta_0$  should be chosen such that the system gets a prescribed  $M_s$ -value. This means solving a nonlinear equation with respect to  $\zeta_0$ . The equation becomes

$$M_s = \max_{\omega > 0} \frac{1}{|1 + (k^*(\zeta_0) + \frac{k_i^*(\zeta_0)}{i\omega})G_p(i\omega)|}. \quad (2.6)$$

This is an approximate, but often sufficiently accurate computation method. This is implemented by the function 'pidesms'.

Another possibility is to determine a function  $\zeta_0 = \zeta_0(\omega_0)$  from the equation

$$M_s = \max_{\omega > 0} \frac{1}{|1 + (k(\omega_0, \zeta_0(\omega_0)) + \frac{k_i(\omega_0, \zeta_0(\omega_0))}{i\omega})G_p(i\omega)|} \quad (2.7)$$

and then maximize  $k_i = k_i(\omega_0, \zeta_0(\omega_0))$  with respect to  $\omega_0$ . This is implemented by the function 'pides2'. This is computationally *very* demanding, but gives the controller with maximal  $k_i$  for a specified  $M_s$ . The approximate method usually gives results very close to the true optimum. For systems with poorly damped poles it is necessary to use the exact design method.

Routines for design of PI controllers with  $\zeta_0$  chosen to get specified values of  $A_m$  and  $\varphi_m$  are also available, 'pidesam' and 'pidespm'. Their implementation is very similar to the one of 'pidesms',  $\zeta_0$  is chosen to get the prescribed value of  $A_m$  or  $\varphi_m$ .

**Design of PD controllers** In case of a PD controller the formulas for  $k$  and  $k_d$  become

$$k = \frac{-\sqrt{1 - \zeta_0^2}A + \zeta_0 B}{\sqrt{1 - \zeta_0^2}(A^2 + B^2)} \quad (2.8)$$

$$k_d = \frac{B}{\omega_0 \sqrt{1 - \zeta_0^2}(A^2 + B^2)}, \quad (2.9)$$

where  $A = \operatorname{Re} G_p(\omega_0(-\zeta_0 + i\sqrt{1 - \zeta_0^2}))$  and  $B = \operatorname{Im} G_p(\omega_0(-\zeta_0 + i\sqrt{1 - \zeta_0^2}))$ . The expressions for  $\zeta_0 > 1$  look similar, but require that  $G_p(s)$  is evaluated twice. The functions  $k(\omega_0, \zeta_0)$  and  $k_d(\omega_0, \zeta_0)$  are implemented in the function 'dppd'.

In the PD case it is natural to maximize the controller constant  $k$ , to get minimum offset to a load disturbance. Design with specified value of  $M_s$  is implemented in the function 'pddesms' in the same way as in 'pidesms'.

**Design of PID controllers** In the case of a PID controller the formulas for  $k$ ,  $k_i$ , and  $k_d$  become

$$k = -\frac{\sqrt{1 - \zeta_0^2}(-2\alpha_0\zeta_0(A^2 + B^2) + (1 + \alpha_0^2)AC) + (\alpha_0^2 - 1)\zeta_0 BC}{(1 - 2\alpha_0\zeta_0 + \alpha_0^2)\sqrt{1 - \zeta_0^2}(A^2 + B^2)C} \quad (2.10)$$

$$k_i = -\alpha_0\omega_0 \frac{(\alpha_0 - \zeta_0)BC + \sqrt{1 - \zeta_0^2}(AC - A^2 - B^2)}{(1 - 2\alpha_0\zeta_0 + \alpha_0^2)\sqrt{1 - \zeta_0^2}(A^2 + B^2)C} \quad (2.11)$$

$$k_d = -\frac{(\alpha_0\zeta_0 - 1)BC + \alpha_0\sqrt{1 - \zeta_0^2}(AC - A^2 - B^2)}{\omega_0(1 - 2\alpha_0\zeta_0 + \alpha_0^2)\sqrt{1 - \zeta_0^2}(A^2 + B^2)C}, \quad (2.12)$$

where  $A = \operatorname{Re} G_p(\omega_0(-\zeta_0 + i\sqrt{1 - \zeta_0^2}))$ ,  $B = \operatorname{Im} G_p(\omega_0(-\zeta_0 + i\sqrt{1 - \zeta_0^2}))$ , and  $C = G_p(-\alpha\omega_0)$ . The expressions for  $\zeta_0 > 1$  look similar, but require that  $G_p(s)$  is evaluated three times. The functions  $k(\omega_0, \zeta_0, \alpha_0)$ ,  $k_i(\omega_0, \zeta_0, \alpha_0)$ , and  $k_d(\omega_0, \zeta_0, \alpha_0)$  are implemented in the function 'dppid'.

The recommended design method is very similar to the one for PI controllers. The only difference is that we must also have a term  $i\omega k_d^*(\zeta_0)$  in the equation for  $M_s$ . The recommended design method is implemented in the function 'piddesms'.

A second method for PID controller design is described in [Persson, 1992b] on pages 97 – 101. The method consists of modifying a well tuned PI controller by increasing the derivative gain  $k_d$ , until the  $M_s$ -function reaches a specified value. This design method is implemented in the function 'piddesms2'. In general this method is inferior to the one implemented in 'piddesms'.

## 3. Implementation

In this chapter the Matlab implementation of the design routines will be discussed. The data structures will be described, and the structure of the design and utility functions will be presented.

### 3.1 Hardware and Software

The implementation which is described here works under Matlab 4.0 on a SPARC. Versions of the routines for Matlab 3.5j on SPARC, PC 386, and PC 486 are also available, but are not described here. All function names have been limited to eight characters to make it possible to use the functions unchanged on PC systems.

### 3.2 Data Structures

The only data structure available in Matlab is the matrix. Everything in Matlab must be expressed with matrices. In this section the conventions used in the DPD Toolbox will be described.

**Transfer functions** Transfer functions are described as strings, where the 's' is the Laplace transform variable. All Matlab math operators and functions are allowed in the transfer functions strings.

The design routines require that these strings can be evaluated when 's' assumes the value of a vector, hence the multiplicative operators '\*', '^', '/', and '\' must be replaced by the operators '.\*', '.^', './', and '\.'. The function 'convert' does this transformation.

#### EXAMPLE 3.1

The transfer function

$$G(s) = \frac{e^{-s}}{(s+1)^2} \quad (3.1)$$

can be described as a string as

```
pstr = 'exp(-s)./(s+1)./(s+1)'
```

or

```
pstr = 'exp(-s)./(s+1).^2'
```

or

```
pstr = 'exp(-s).*(s+1).^(-2)'. □
```

#### EXAMPLE 3.2

The transfer function

$$G(s) = \frac{1}{(s+1)^8} \quad (3.2)$$

is described as

```
pstr = '(s+1).^(-8)' or pstr = '1./(s+1).^8'. □
```



In Matlab Version 3, it is possible to use global variables in the definition of transfer function strings. This is not possible in Matlab Version 4, without changes in the code. To use global variables in the transfer function strings, the variables should be declared 'global' in the routine 'evals'.

**The Controller Data Structure** The design routines return a row vector describing the controller. The vector has the content

$$c = [\omega_0 \quad \zeta_0 \quad \alpha_0 \quad k \quad k_i \quad T_i \quad k_d \quad T_d \quad N_r]. \quad (3.3)$$

The parameter  $\omega_0$  is the distance of the dominant poles from the origin,  $\zeta_0$  is their relative damping. In case that the design routine uses a third pole  $-\alpha_0\omega_0$ , the value of  $\alpha_0$  is returned, in other cases NaN is returned. The controller parameters are returned as  $k$ ,  $k_i$ ,  $T_i$ ,  $k_d$ , and  $T_d$ . The parameter  $N_r$  tells which of the fundamental pole placement routines has computed the controller. This parameter is never used. This data structure should be accessed by the access functions 'getw0', 'getz0', 'geta0', 'getk', 'getki', 'getti', 'getkd', and 'gettd'. The routine 'con2str' converts the controller data structure to the controller transfer function expressed as a string.

**Frequency responses** Frequency responses generated with 'sfrcol' are stored as

$$\begin{bmatrix} \omega_1 & G_1(\omega_1) & \dots & G_m(\omega_1) \\ \vdots & \vdots & & \vdots \\ \omega_n & G_1(\omega_n) & \dots & G_m(\omega_n) \end{bmatrix}. \quad (3.4)$$

This data structure is also used in the toolboxes described in [Gustafsson *et al.*, 1990a] and [Gustafsson *et al.*, 1990b], which are recommended for plotting.

### 3.3 Utility Routines

The transfer functions represented as strings are evaluated with the function 'evals',

**EXAMPLE 3.3**—The use of evals

To evaluate the transfer function  $G(s) = e^{-s}/(s+1)^2$ , use the following commands.

```
>> pstr = 'exp(-s)./(s+1)./(s+1)';
>> w = [0.1 0.2 0.3];
>> evals(pstr, i*w)
```

```
ans =
```

```
0.9461 - 0.2920i    0.7964 - 0.5388i    0.5825 - 0.7088i
```

□

There are routines for computing the maximum of the sensitivity function, 'mscl', amplitude margin, 'amarg', and phase margin, 'pmarg'. The routines for computing amplitude and gain margins are based on the general routines 'asolveol' and 'psolveol' which solve the equations

$$a_1 = |G_p(i\omega)G_c(i\omega)| \quad (3.5)$$

$$a_2 = \arg G_p(i\omega)G_c(i\omega) \quad (3.6)$$

with respect to  $\omega$  for given  $a_1$  and  $a_2$ .

All these routines require the process and the controller expressed as strings. There is a conversion routine that converts the controller data structure to a string, 'con2str'.

The function 'makep' returns one of a number of standard processes with parameters given as arguments to the function. The following processes are provided

$$G_1 = \frac{p_1 e^{-sp_2}}{sp_3 + 1} \quad (3.7)$$

$$G_2 = \frac{1}{(s + 1)^{p_1}} \quad (3.8)$$

$$G_3 = \frac{(1 - sp_1)}{(s + 1)^3} \quad (3.9)$$

$$G_4 = \frac{e^{-sp_1}}{(sp_2 + 1)(sp_3 + 1)} \quad (3.10)$$

$$G_5 = \frac{p_2^2 e^{-sp_1}}{(s^2 + 2sp_2p_3 + p_2^2)} \quad (3.11)$$

$$G_6 = \frac{p_2^3 p_4 e^{-sp_1}}{(s + p_2 p_4)(s^2 + 2sp_2p_3 + p_2^2)} \quad (3.12)$$

$$G_7 = \frac{1}{(s + 1)(sp_1 + 1)(sp_1^2 + 1)} \quad (3.13)$$

$$G_8 = \frac{1}{(s + 1)(sp_1 + 1)(sp_1^2 + 1)(sp_1^3 + 1)} \quad (3.14)$$

$$G_9 = \frac{(1 - sp_1/2)}{(1 + sp_1/2)(1 + sp_2)} \quad (3.15)$$

$$G_{10} = \frac{e^{-sp_1}}{(1 + sp_2)(1 + sp_3)(1 + sp_4)} \quad (3.16)$$

$$G_{11} = \frac{e^{-sp_1}}{s(sp_2 + 1)} \quad (3.17)$$

$$G_{12} = \frac{e^{-sp_1}}{s(sp_2 + 1)(sp_3 + 1)} \quad (3.18)$$

Many of the routines require a tolerance for the solution of an equation or an optimization. When a tolerance is required in a function it can be given as a parameter to the function. If the tolerance parameter is omitted the tolerance is taken from the global variable 'GTOL', and if 'GTOL' is not defined the tolerance is given from the function 'deftol'. Normally the tolerance is given from 'deftol' and is set to  $10^{-4}$  as default.

The optimization is carried out with the functions 'opt' and 'optg'. These functions implement a golden ratio optimization algorithm.

The solution of an equation is carried out by the functions 'solve' and 'solveb'. In these functions a simple bisection method is implemented. The function 'solveb' is specially implemented for solving phase equations to get the correct phase value across  $n\pi$  borders.

### 3.4 Design Routines

The design routines are implemented in a number of layers, see Figure 3.1.

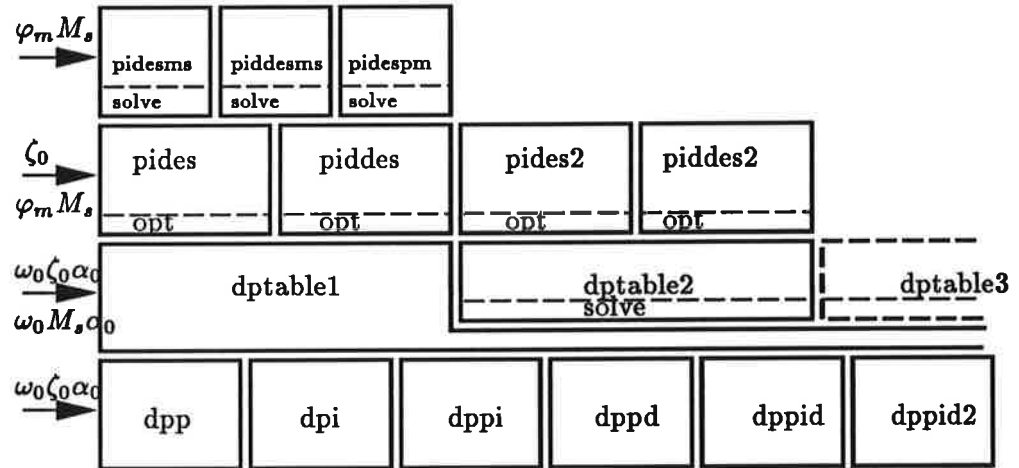


Figure 3.1 The structure of the design routines.

The lowest level consists of the basic pole placement routines, i.e., routines that compute controller parameters given  $\omega_0$ ,  $\zeta_0$ , and  $\alpha_0$ . The next level consists of interface routines to the basic primitives. There may be several of these interface routines. In this version four interface routines are available, 'dptable1', 'dptable2', 'dptable3', and 'dptable4'. In 'dptable1' controllers are computed given  $\omega_0$ ,  $\zeta_0$ , and  $\alpha_0$ ,  $\omega_0$  may be a vector. In 'dptable2' controllers are computed given  $\omega_0$ ,  $M_s$ , and  $\alpha_0$ . This involves solving the  $M_s$ -equation with respect to  $\zeta_0$ . The routine 'dptable3' takes the input parameters  $\omega_0$ ,  $A_m$ ,  $\alpha_0$  and 'dptable4' takes the input parameters  $\omega_0$ ,  $\varphi_m$ ,  $\alpha_0$ .

The third level optimizes  $k_i$ , given a  $\zeta_0$  or  $M_s$ . Finally, the fourth level solves equations with respect to  $\zeta_0$  such that some stability requirements are fulfilled. These requirements may be  $M_s$ ,  $A_m$ , or  $\varphi_m$ . These are the approximate, but recommended design functions.

## 4. Examples

This chapter will present a number of examples of the use of the design routines, and some discussions of the examples. All examples in this chapter have been executed in Matlab Version 4.0 on an SPARC ELC, exactly as they are written here.

### 4.1 Hints for the use of the routines

When design is carried out with  $M_s$  as design parameter the default search interval of  $\zeta_0$  is 'linspace(0.01, 2, 10)', this can handle most cases, but requires a lot of computation. If the plant is well damped and we require a moderate  $M_s$ , 1.6 to 2, say, then a smaller search interval can be supplied, e.g.,  $\zeta_0 = [0.1 \ 0.5 \ 0.9]$

The  $\omega_0$  should be chosen such that we are sure to detect a maximum. The computations are relatively cheap, don't use a too sparse discretization.

Normally a tolerance of  $10^{-4}$  in the computations is sufficient. A larger tolerance may cause errors in the routines.

Problems may also appear if the process transfer function is evaluated at a process pole. To avoid this choose the initial guesses of  $\omega_0$  and  $\alpha_0$  'irregularly.' For example, if the routine 'piddesms' is used for  $G_p(s) = e^{-s}/(s+1)$  choose  $w0s = [0.11:0.1:3.0]$  rather than  $w0s = [0.1:0.1:3.0]$  or  $\alpha = 1.01$  rather than  $\alpha = 1.0$ .

Observe that the optimization is meaningless for processes of too low order. For a second order process we can place the poles anywhere with a PID controller, and get the closed system arbitrarily fast.

In Example 4.1 the function 'polyadd' is used. The function is defined as:

```
function p = polyadd(p1, p2)
dn = length(p1) - length(p2);
p = [zeros([1 -dn]) p1] + [zeros([1 dn]) p2];
```

### 4.2 Examples

In the following a number of examples of the use of the routines will be presented. The reader may need to consult Chapter 5 for a brief description of the function.

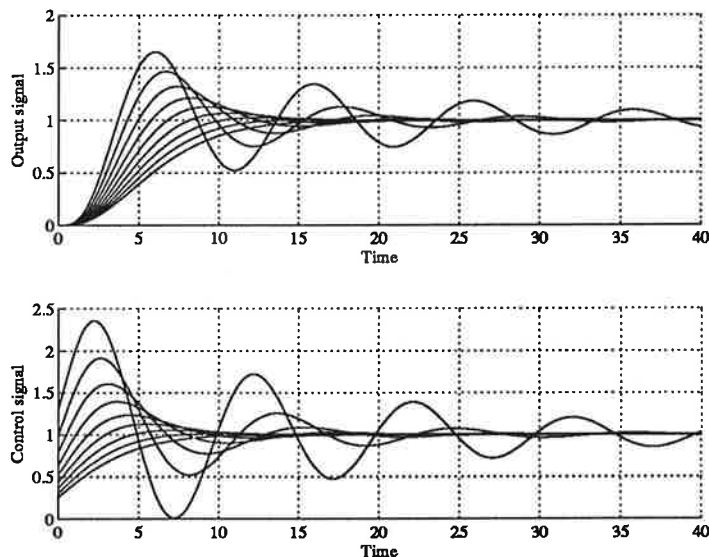
**EXAMPLE 4.1—Controllers with different  $\zeta_0$**

In the following example a number of PI controllers are designed with different values of  $\zeta_0$  for the process

$$G(s) = \frac{1}{(s+1)^4}.$$

The corresponding family of step responses and control signals are simulated and plotted.

```
pstr = '1./(s+1).^4';
b = 1; a = conv([1 2 1], [1 2 1]);
tv = [0:0.1:40]; w0s = [0.1:0.1:2]; z0s = [0.1:0.1:0.9];
for ix = z0s,
    con = pides(pstr, w0s, ix);
    r = [1 0]; s = [getk(con) getki(con)]; t = s;
    ysp = step(conv(b, t), polyadd(conv(a, r), conv(b, s)), tv);
    usp = step(conv(a, t), polyadd(conv(a, r), conv(b, s)), tv);
    subplot(2, 1, 1); hold on; plot(tv, ysp); drawnow;
    subplot(2, 1, 2); hold on; plot(tv, usp); drawnow;
end;
subplot(2, 1, 1);
grid; xlabel('Time'); ylabel('Output signal');
subplot(2, 1, 2);
grid; xlabel('Time'); ylabel('Control signal');
```



**Figure 4.1** Step responses of  $G(s) = 1/(s+1)^4$  controlled by PI controllers designed with different  $\zeta_0$ .

Designing controllers with  $\zeta_0$  as the design parameter may produce very different looking step responses with the same  $\zeta_0$  for different processes. For this reason it is better to design controllers with specifications on  $M_s$ .

In this case the simulation is carried out by computing the transfer functions and using the standard function 'step'.  $\square$

EXAMPLE 4.2—Computation of  $A_m$  and  $\varphi_m$   
 In this example we will design PI controllers for

$$G(s) = \frac{e^{-sL}}{s+1}$$

with different values of the design parameter  $M_s$ . Phase and amplitude margins are then computed for the final systems as functions of  $M_s$ .

```
dels = [0.1 0.5 1];
mss = [1.8:0.2:3.0]; ws = [0.1:0.1:20]; res = [];
for jx = dels,
    tmp = [];
    pstr = makep(1, 1, jx, 1);
    for ix = mss,
        con = pidesms(pstr, ws, ix);
        tmp = [tmp;
            amarg(pstr, con2str(con), ws) ...
            pmarg(pstr, con2str(con), ws)];
    end;
    res = [res tmp];
end;
c1 = 1:2:cols(res); c2 = 2:2:cols(res);
subplot(2, 1, 1); plot(mss, res(:, c1));
xlabel('Ms'); ylabel('Am');
text(mss(2), res(2, c1(1)), 'L=0.1')
text(mss(2), res(2, c1(2)), 'L=0.5')
text(mss(2), res(2, c1(3)), 'L=1.0')
subplot(2, 1, 2); plot(mss, res(:, c2));
xlabel('Ms'); ylabel('Pm');
text(mss(2), res(2, c2(1)), 'L=0.1')
text(mss(2), res(2, c2(2)), 'L=0.5')
text(mss(2), res(2, c2(3)), 'L=1.0')
```

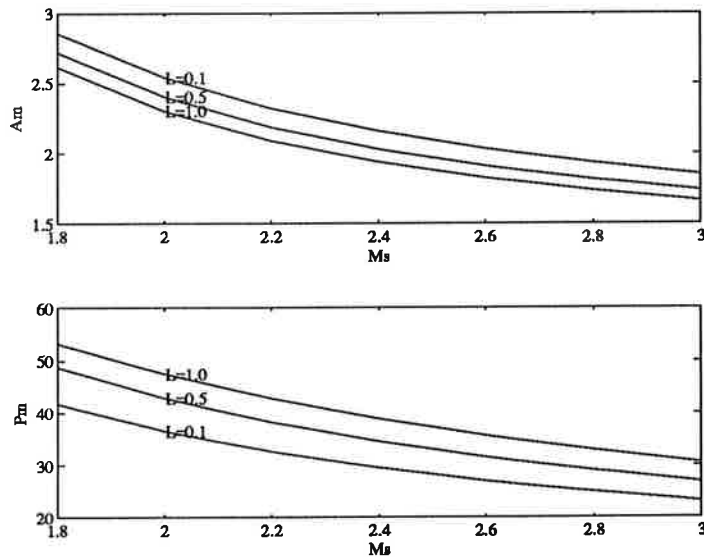


Figure 4.2 Amplitude and phase margins for systems designed with different  $M_s$ .

This example illustrates that a constant  $A_m$  or  $\varphi_m$  as design parameter may give very different  $M_s$  in the closed system for different processes. □

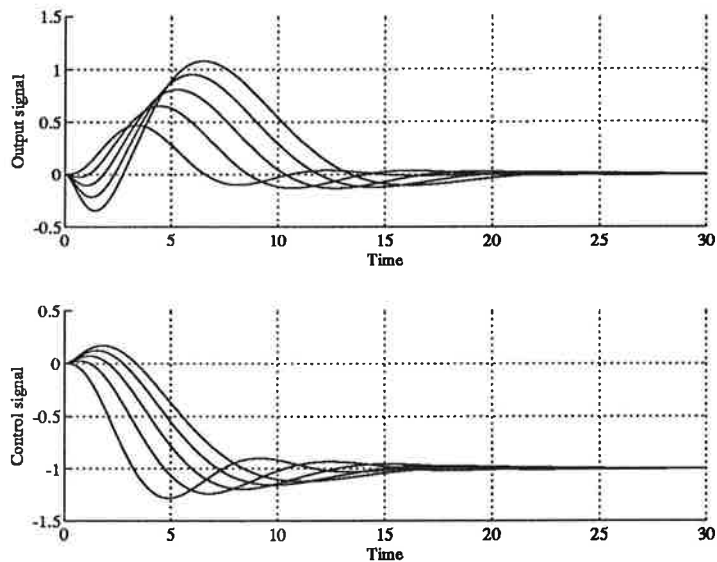
**EXAMPLE 4.3—A process with different parameters**  
 In this example the process

$$G_p(s) = \frac{1 - \alpha s}{(s + 1)^3}$$

is controlled by a PI controller designed with  $M_s = 2$  for different values of the parameter  $\alpha$ . Load responses and the corresponding control signals are simulated and plotted.

```

clg;
tv = [0:0.1:30]; alphas = [0:0.5:2]; w0s = [0.1:0.1:2];
for ix = alphas,
    pstr = makep(3, ix);
    b = [-ix 1];
    a = [1 3 3 1];
    con = pidesms(pstr, w0s, 2);
    r = [1 0]; s = [getk(con) getki(con)]; t = s;
    ysp = step(conv(b, r), polyadd(conv(a, r), conv(b, s)), tv);
    usp = step(-conv(b, s), polyadd(conv(a, r), conv(b, s)), tv);
    subplot(2, 1, 1); hold on; plot(tv, ysp); drawnow;
    subplot(2, 1, 2); hold on; plot(tv, usp); drawnow;
end;
subplot(2, 1, 1); grid; xlabel('Time');
ylabel('Output signal');
subplot(2, 1, 2); grid; xlabel('Time');
ylabel('Control signal');
  
```



**Figure 4.3** Load responses for  $G_p(s) = (1 - s\alpha)/(s + 1)^3$  for different values of  $\alpha$  with PI controllers designed with  $M_s = 2$ .

The non-minimum phase character of the responses can clearly be seen for large values of  $\alpha$ . □

**EXAMPLE 4.4—Nyquist curves from PI design on different processes**  
 In this example we design PI controllers for a number of different processes. The processes are described in the array 'a', which contains the process number and the parameters of the process.

PI controllers are designed with constant  $\zeta_0$  and with constant  $M_s$ . As can be seen from the plotted Nyquist curves we get much more similarity of different loop transfer functions with controllers designed with  $M_s$  than with constant  $\zeta_0$ .

```
w = logspace(-1, 2, 200)'; resms = [w]; resz = [w];
a = [1 1 1 1
     2 4 0 0
     3 1 0 0
     4 1 2 2];
for ix = a'
    pstr = makep(ix');
    cms = pidesms(pstr, [0.1:0.1:10], 2);
    cz = pides(pstr, [0.1:0.1:10], 0.5);
    f = sfrcol(pstr, con2str(cms), w);
    resms = [resms f(:, 2)];
    f = sfrcol(pstr, con2str(cz), w);
    resz = [resz f(:, 2)];
end;
nypl(resms); nygrid;
plotc(-1, 1/2, -pi/2, pi/2);
axis([-1 0.5 -1 0.5])
nypl(resz); nygrid;
plotc(-1, 1/2, -pi/2, pi/2);
axis([-1 0.5 -1 0.5])
```

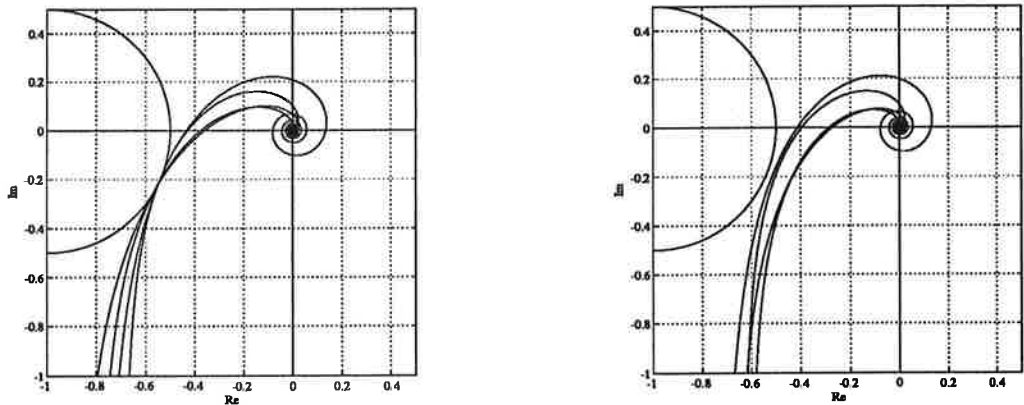


Figure 4.4 To the left: Nyquist plots for processes designed with constant  $M_s$ , to the right: Nyquist plots for processes designed with constant  $\zeta_0$ .



**EXAMPLE 4.5—Timing**

In this example controllers are computed with different tolerance in the optimization and equation solving routines, and the time it takes to compute the different controllers are recorded.

```
pstr = 'exp(-s)/(s + 1)'; w0s = [0.1:0.1:10]; res = [];  
for ix = 2:8  
    tic;  
    c = pidesms(pstr, w0s, 2, [], 10^(-ix));  
    t1 = toc;  
    tic;  
    c = pidesms(pstr, w0s, 2, [0.2:0.2:0.6], 10^(-ix));  
    t2 = toc;  
    res = [res; t1 t2];  
end;  
res
```

```
res =  
  
    11.0203     9.1207  
    18.1734    15.8076  
    28.4095    25.4336  
    39.2039    35.6356  
    52.6931    48.4410  
    69.8643    65.0888  
    93.5680    83.4929
```

Some time can be saved by guessing a correct interval for  $\zeta_0$  instead of relying on the default values from the algorithm. □

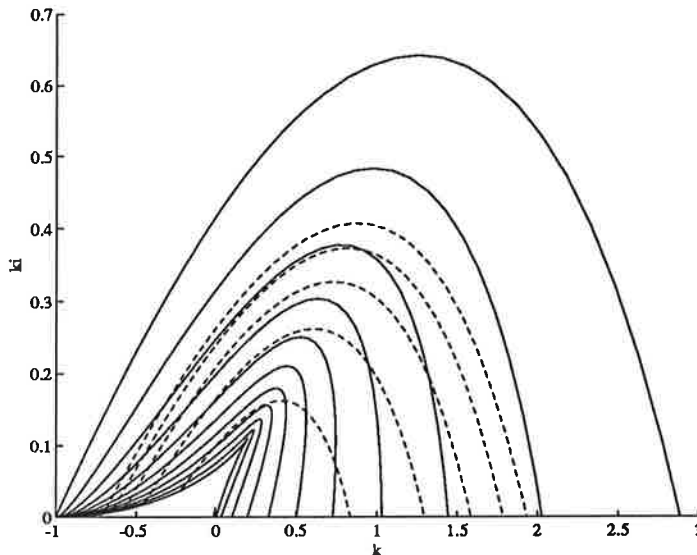
**EXAMPLE 4.6—Computation of  $k_i(k)$  curves**

This is an example where we need to use the low level routines directly. We compute the  $k$  and  $k_i$  from Equations 2.4 and 2.5 as functions of  $\omega_0$  for

$$G_p(s) = \frac{1}{(s+1)^5}. \quad (4.1)$$

In the first case we choose constant  $\zeta_0$ . The curves  $k_i(k)$  are shown with solid lines in Figure 4.5. In the second case we choose  $\zeta_0$  such that we get a prescribed  $M_s$  value for all  $\omega_0$ . The  $k_i(k)$  are shown with dashed lines.  $\zeta_0$  is chosen in the interval 0.0:0.1:1, and  $M_s$  in the interval 1.5:0.5:3.5. As can be seen a maximum of a dotted line corresponds reasonably well to a maximum of a dashed line. This Matlab computation takes rather long time.

```
clg;
pstr = '1./(s+1).^5'; w0s = [0.01:0.02:1];
z0s = [0:0.1:0.9 0.999]; mss = [1.5:0.5:3.5];
for ix = z0s,
    c = dptable1(pstr, 'pi', w0s, ix);
    hold on; plot(getk(c), getki(c));
end;
axis([-1 3 0 0.7])
for ix = mss,
    c = dptable2(pstr, 'pi', w0s, ix);
    hold on; plot(getk(c), getki(c), '--');
end;
xlabel('k'); ylabel('ki');
```



**Figure 4.5** Plots of  $k_i(k)$  for constant  $\zeta_0$  (solid lines) and constant  $M_s$  (dashed lines) for varying  $\omega_0$ .

These kinds of curves are very useful to have to get a feel for how a system react for choices of different  $\zeta_0$  and  $\omega_0$ . □

#### EXAMPLE 4.7—Interface to Simnon

In some cases it may be desirable to use Simnon for simulation, for example when we want to handle time delays or non-linearities. This example demonstrates how easily it is done with the Simnon interface routines.

The routines 'makep' and 'setsimnon' are designed to accept the same parameters describing a transfer function. This requires that all the necessary systems are implemented in Simnon, or that Matlab can generate the necessary Simnon code. In these examples the Simnon code is written by hand.

In this example we design a PID controller for the process

$$G(s) = \frac{e^{-2s}}{s+1}$$

and simulate the step and load responses for different values of the set point weighting factor  $\beta$ . The plots show a family of step responses when  $\beta$  assumes the values 0:0.2:1.

```
clg; w0s = [0.1:0.1:10]; alpha0 = 1.01; ms = 2.0;
betas = 0:0.2:1; ts = 30;
pstr = makep(1, 1, 2, 1);
con = piddesms(pstr, w0s, ms, alpha0);
setsimnon(1, 1, 2, 1);
par('t1', ts/2); % t1 is the time when the load starts acting
setpid(con);
for ix = betas,
    par('b', ix);
    y = simu(0, ts);
    subplot(2, 1, 1); hold on; plot(y(:, 1), y(:, 2)); drawnow;
    subplot(2, 1, 2); hold on; plot(y(:, 1), y(:, 3)); drawnow;
end;
subplot(2, 1, 1); grid; xlabel('Time'); ylabel('Output signal');
subplot(2, 1, 2); grid; xlabel('Time'); ylabel('Control signal');
```

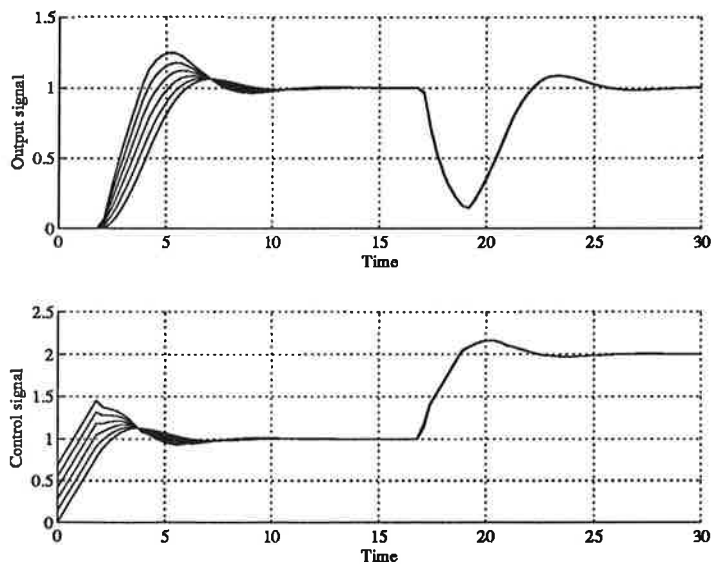


Figure 4.6 Step and load responses for different  $\beta$  in the interval [0 1].

□

## 5. The Routines

Only a brief description of the routines will be given in this report. The complete source code listing is given in [Persson, 1992a]. The help text of the functions will also explain the meaning of the parameters.

### Brief Description of the Routines

Design Routines	
<b>ides</b>	I controller with maximal $k_i$ .
<b>pid</b>	PI controller with specified $\zeta_0$ and maximal $k_i$ .
<b>pid2</b>	PI controller with specified $M_s$ and maximal $k_i$ .
<b>pidesm</b>	PI controller with maximal $k_i$ and $\zeta_0$ chosen to get specified $M_s$ .
<b>pidesam</b>	PI controller with maximal $k_i$ and $\zeta_0$ chosen to get specified $A_m$ .
<b>pidespm</b>	PI controller with maximal $k_i$ and $\zeta_0$ chosen to get specified $\varphi_m$ .
<b>betades</b>	Design of the set point weighting factor.
<b>pidesk</b>	PID controller with specified $\zeta_0$ and $k_d$ and maximal $k_i$ .
<b>pid</b>	PID controller with specified $\zeta_0$ and $\alpha_0$ and maximal $k_i$ .
<b>pid2</b>	PID controller with specified $M_s$ and $\alpha_0$ and maximal $k_i$ .
<b>pidesm</b>	PID controller with maximal $k_i$ and $\zeta_0$ chosen to get specified $M_s$ .
<b>pidesam</b>	PID controller with maximal $k_i$ and $\zeta_0$ chosen to get specified $A_m$ .
<b>pidespm</b>	PID controller with maximal $k_i$ and $\zeta_0$ chosen to get specified $\varphi_m$ .
<b>pidesm2</b>	PID controller with maximal $k_i$ and specified $M_{s1}$ and $M_{s2}$ .
<b>pd</b>	PD controller with specified $\zeta_0$ and maximal $k$ .
<b>pdesm</b>	PD controller with maximal $k$ and $\zeta_0$ chosen to get specified $M_s$ .
<b>pdesam</b>	PD controller with maximal $k$ and $\zeta_0$ chosen to get specified $A_m$ .
<b>pdespm</b>	PD controller with maximal $k$ and $\zeta_0$ chosen to get specified $\varphi_m$ .
<b>zn1pi</b>	PI according to Ziegler-Nichols oscillation method.
<b>zn1pid</b>	PID according to Ziegler-Nichols oscillation method.
<b>zn2pi</b>	PI according to Ziegler-Nichols step response method.
<b>zn2pid</b>	PID according to Ziegler-Nichols step response method.

Computation Routines	
<b>amarg</b>	Computes the amplitude margin.
<b>pmarg</b>	Computes the phase margin.
<b>asolve1</b>	Solves $ L(i\omega)/(1 + L(i\omega))  = a$ , with $L(i\omega) = G_c(i\omega)G_p(i\omega)$
<b>asolve01</b>	Solves $ G_c(i\omega)G_p(i\omega)  = a$ .
<b>psolve1</b>	Solves $\arg L(i\omega)/(1 + L(i\omega)) = \alpha$ , with $L(i\omega) = G_c(i\omega)G_p(i\omega)$
<b>psolve01</b>	Solves $\arg G_c(i\omega)G_p(i\omega) = \alpha$ .
<b>mpbeta</b>	Finds the $M_p$ value of the closed system as function of $\beta$ .
<b>mscl</b>	Finds the $M_s$ value.
<b>mshelp</b>	Help routine used in 'pidesm2'.
<b>sfrcol</b>	Finds the frequency responses. For Bode and Nyquist plots.

<b>Basic Pole Placement Routines</b>	
<b>dptable1</b>	Interface to the pole placement routines. Argument $\omega_0 \zeta_0 \alpha_0$ .
<b>dptable2</b>	Interface to the pole placement routines. Argument $\omega_0 M_s \alpha_0$ .
<b>dptable3</b>	Interface to the pole placement routines. Argument $\omega_0 A_m \alpha_0$ .
<b>dptable4</b>	Interface to the pole placement routines. Argument $\omega_0 \varphi_m \alpha_0$ .
<b>dpi</b>	Pole placement of an I controller.
<b>dpp</b>	Pole placement of an P controller.
<b>dppd</b>	Pole placement of an PD controller.
<b>dppi</b>	Pole placement of an PI controller.
<b>dppid</b>	Pole placement of an PID controller.
<b>dppid2</b>	Pole placement of an PID controller with specified $k_d$ .

<b>Help Routines</b>	
<b>closeit</b>	Computes $x/(1+x)$ .
<b>cols</b>	Returns the number of columns of a matrix.
<b>evals</b>	Evaluates a string for a given value of 's'.
<b>evalx</b>	Evaluates a string for a given value of 'x'.
<b>kdguess</b>	A guess of the maximal $k_d$ . Used in 'pidesm2'.
<b>locmax</b>	Finds a local maximum in an array.
<b>phase</b>	Computes the phase of a vector of complex numbers.
<b>plotc</b>	Plots a circle.
<b>rows</b>	Returns the number of rows of a matrix.

<b>Default Handling Routines</b>	
<b>defprint</b>	Controls the printout form some of the routines.
<b>deftol</b>	Gives the tolerance for the optimization and solving routines.

<b>Optimization and Solving Routines</b>	
<b>opt</b>	One parameter optimizer.
<b>optg</b>	Computes the global optimum on an interval. Used in 'mscl'.
<b>solve</b>	One parameter equation solver.
<b>solveb</b>	Solving routine used in 'psolve*', to get the phase right. Inefficient.

<b>Conversion Routines</b>	
<b>con2str</b>	Convert the controller matrix to the PID controller string
<b>convert</b>	Converts a transfer function string to accept vector arguments.
<b>geta0</b>	Access function to get $\alpha_0$ .
<b>getk</b>	Access function to get $k$ .
<b>getkd</b>	Access function to get $k_d$ .
<b>getki</b>	Access function to get $k_i$ .
<b>gettd</b>	Access function to get $T_d$ .
<b>getti</b>	Access function to get $T_i$ .
<b>getw0</b>	Access function to get $\omega_0$ .
<b>getz0</b>	Access function to get $\zeta_0$ .
<b>makep</b>	Returns a number of standard systems as strings.
<b>numtostr</b>	Converts a numerical value to a string.
<b>par2con</b>	Converts PID parameter to the standard controller data structure.
<b>qstring</b>	Returns a string with quotes.
<b>sfun</b>	Implements the functions of 'makep'.
<b>tf2str</b>	Converts transfer function to a string.

## The function calls

An alphabetic list of the function with input and output arguments is also given for reference. The following argument conventions have been used

<b>am</b>	= amplitude margin
<b>con</b>	= the controller data structure (see Section 3.2)
<b>contype</b>	= a string describing the controller type
<b>cstr</b>	= a controller expressed as a string
<b>k, ti, td</b>	= PID controller parameters
<b>ms</b>	= the $M_s$ parameter
<b>pm</b>	= phase margin
<b>pstr</b>	= a transfer function expressed as a string
<b>str</b>	= a string
<b>tol</b>	= the tolerance for equation solving and optimization
<b>ws, w0, w0s</b>	= an array of frequencies
<b>z0</b>	= the $\zeta_0$ parameter
<b>zguess</b>	= changes default value of $\zeta_0$ search interval

The design routines return the controller parameters  $k$ ,  $T_i$ , and  $T_d$ . If only one output argument is present in a call of the design functions, the complete controller structure is returned. For a complete description of the functions, use the help texts of the functions. Parameters which have default values are written in *italics*.

```
[am, wx] = amarg(cstr, pstr, ws, tol)
  wx = asolvecl(cstr, pstr, y, ws, tol)
  wx = asolveol(cstr, pstr, y, ws, tol)
  b = betades(pstr, con, mp, tol)
res = closeit(f)
res = cols(matrix)
str = con2str(k, ti, td, n)
str = convert(pstr)
  dp = defprint
  dt = deftol
con = dpi(pstr, w0)
con = dpp(pstr, w0)
con = dppd(pstr, w0, z0)
con = dppi(pstr, w0, z0)
con = dppid(pstr, w0, z0, alpha0)
con = dppid2(pstr, w0, z0, kd0)
con = dptable1(pstr, contype, w0s, z0, p1)
con = dptable2(pstr, contype, w0s, ms, p1, tol)
con = dptable3(pstr, contype, w0s, pm, p1, tol)
con = dptable4(pstr, contype, w0s, am, p1, tol)
  r = evals(str, s)
  r = evalx(str, x)
a0 = geta0(con)
k = getk(con)
kd = getkd(con)
ki = getki(con)
```

```

        td = gettd(con)
        ti = getti(con)
        w0 = getw0(con)
        z0 = getz0(con)
        con = ides(pstr, w0s, tol)
        kd = kdguess(pstr, w0s, z0)
        res = locmax(array)
        str = makep(sysnr, p1, p2, p3, p4)
        mpr = mpbeta(pstr, k, ti, td, betas, tol)
    [ms, ws] = mscl(cstr, pstr, wx, tol)
        str = numtostr(num, n)
    [xsol, fx] = opt(str, x0s, tol)
    [xsol, fx] = optg(str, x0s, tol)
        con = par2con(k, ti, td)
    [k, td] = pddes(pstr, w0s, z0, tol)
    [k, td] = pddesam(pstr, w0s, am, zguess, tol)
    [k, td] = pddesms(pstr, w0s, ms, zguess, tol)
    [k, td] = pddespm(pstr, w0s, pm, zguess, tol)
        phi = phase(g)
    [k, ti, td] = pidde(pstr, w0s, z0, alpha0, tol)
    [k, ti, td] = pidde2(pstr, w0s, ms, alpha0, tol)
    [k, ti, td] = pidesam(pstr, w0s, am, alpha0, zguess, tol)
    [k, ti, td] = piddekd(pstr, w0s, z0, kd0, tol)
        cons = pidde2m(pstr, w0s, ms1, ms2, kdr, zguess, tol)
    [k, ti, td] = pidesms(pstr, w0s, ms, alpha0, zguess, tol)
    [k, ti, td] = pidespm(pstr, w0s, pm, alpha0, zguess, tol)
    [k, ti] = pides(pstr, w0s, z0, tol)
    [k, ti] = pides2(pstr, w0s, ms, tol)
    [k, ti] = pidesam(pstr, w0s, am, zguess, tol)
    [k, ti] = pidesms(pstr, w0s, ms, zguess, tol)
    [k, ti] = pidespm(pstr, w0s, pm, zguess, tol)
        plotc(z, rad, phi1, phi2, lt, ddeg)
    [pm, wx] = pmarg(cstr, pstr, ws, tol)
        wx = psolvecl(cstr, pstr, y, ws, tol)
        wx = psolveol(cstr, pstr, y, ws, tol)
        str = qstring(str)
        res = rows(matrix)
        fr = sfrcol(cstr, pstr, w1, w2, n)
        xx = solve(str, y, x0, fol, tol)
        xx = solveb(str, y, x0, fol, tol)
        str = tf2str(num, den)
    [k, ti] = zn1pi(pstr, w0s, tol)
    [k, ti, td] = zn1pid(pstr, w0s, tol)
    [k, ti] = zn2pi(kp, l, t)
    [k, ti, td] = zn2pid(kp, l, t)

```

### Interface to Simnon

Matlab is excellent for many kinds of numerical computations, but there exists better tools for simulation of dynamical systems. When the systems contain non-linearities or time delays Matlab is normally not adequate. The simulation

program Simnon is then a better suited program. Simnon is described in [SSPA, 1990].

A facility has been written to run simulations in Simnon from Matlab, with Simnon working as a 'computation engine.' With this software it is possible to issue Simnon commands from Matlab, to set Simnon parameters from Matlab, and to get numerical results back from Simnon to Matlab.

It is now described how to use the interface functions. Matlab and Simnon communicate via named Unix-pipes. In the following we assume that Matlab will be run from the directory /home/nisse/matlabdir and that Simnon will be run from the directory /home/nisse/simnondir. The file pipsimu is located in the directory /home/nisse/lib.

First we have to make two named pipes in Unix, the pipes will be called ipipe and opipe, and will appear to the user as two files in the directory /home/nisse/simnondir. The named pipes are created with the following commands

```
% /usr/etc/mknod /home/nisse/simnondir/ipipe p
% /usr/etc/mknod /home/nisse/simnondir/opipe p
```

Open a new window, and give the following commands

```
% cd /home/nisse/simnondir
% /home/nisse/lib/pipsimu ipipe
```

This will start Simnon in the window and Simnon will now accept commands from the named pipe ipipe.

The program pipsimu is a Perl script, which starts Simnon, reads commands from ipipe and sends them along to Simnon. This seems unnecessary, but Simnon refuses to read commands directly from a named pipe. For readings on Perl, see [Wall and Schwartz, 1991]. The Perl script is quite short and simple:

```
#!/usr/local/bin/perl
$pipename = shift;
open(SIMNON, "|simnon") || die "cannot pipe\n";
select(SIMNON);
$|=1;
print "x\n";
#print "algor rkf45\n";
while(open(PIPE, $pipename)){
    while(<PIPE>){
        if (eof(PIPE)) {close(PIPE);}
        print;
        if ($_ eq "stop\n") {exit;};
    }
}
```

Now open a window where Matlab will be run, and give the following commands

```
% cd /home/nisse/matlabdir
% matlab
>> global inpipe1 outpipe1 simdir
>> inpipe1 = '/home/nisse/simnondir/ipipe';
```



```
>> outpipe1 = '/home/nisse/simnondir/opipe';
>> simdir   = '/home/nisse/simnondir/';
```

Two Matlab variables 'inpipe1' and 'outpipe1' are now defined and have values of the named pipes. The variables must have the full path names. Simnon will now accept commands from Matlab. A typical command sequence will be

```
>> gp = makep(2, 8);           % Define a transfer function
>> cp = pidesms(gp, [0.1:0.1:5], 2); % Design a controller
>> setsimnon(2, 8);          % Compile the correct model
>> setpid(cp);               % Set the PID parameters
>> r = simu(0, 100);         % Simulate
>> plot(r(:,1), r(:, 2));    % Plot the results
```

The functions 'makep' and 'setsimnon' must be written such that the parameters give Simnon commands to compile the correct model files. The function 'setsimnon' typically contains a number of Simnon commands as 'syst' and 'par'.

### Listing of functions for interfacing Simnon

In this section a few of the commands for the Matlab-Simnon interface will be listed. Some of these functions must be modified to suit the the Simnon models with which will be used.

The basic command is 'remcommand' which sends text strings to a named pipe. The function 'simc' calls 'remcommand' but uses 'inpipe1' as default for the named pipe, 'simc' also accepts several input parameters.

---

```
function remcommand(com, inpipe)

%REMMCOMMAND Sends a command string to be executed in a remote
%              process. The command is sent via the named pipe inpipe.

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Tue Dec  8 13:41:16 1992

eval(['!/usr/5bin/echo ' ' ' com ' ' ' ' >> ' inpipe]);
```

---

```
function simc(p1, p2, p3, p4, p5, p6, p7, p8, p9, p10)

%SIMC A number of text strings are sent to a remote process (Simnon).
%      The command is sent via a named pipe which name is in the
%      global variable inpipe1.
%
%      SIMC(p1)
%      .
%      .
%      .
%      SIMC(p1, p2, p3, p4, p5, p6, p7, p8, p9, p10)

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Tue Dec  8 13:41:16 1992
```

```

global inpipe1
for ix=1:nargin,
    remcommand(eval(sprintf('p%g', ix)), inpipe1);
end;

```

---

The commands 'getsimval' and 'getsimvec' transfer Simnon variables, parameters and simulation results back to Matlab. 'getsimval' gets the value of a variable or parameter, and 'getsimvec' gets the simulation result stored in the file 'store.d' or any other file the user specifies. The Matlab user must know what has been stored in 'store.d' and in which order the signals are stored.

The communication goes via the named pipe 'outpipe1' to get synchronization of Matlab and Simnon. This way Matlab does not go on computing until Simnon has provided the required value.

---

```

function res = getsimval(val, outpipe)

%GETSIMVAL Gets a value of a parameter or variable form Simnon.
%
%           res = getsimval(val)
%           val = the name of the parameter or variable as a string
%
%           The value is sent back via a named pipe, outpipe, to
%           synchronize the two processes.

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Sat Dec 5 16:29:50 1992

if nargin==1
    global outpipe1
    outpipe = outpipe1;
end;

simc(['disp ' val '/ztmp']);
simc('write (dk zzfile FF)');
simc('write (dk zzfile) ztmp. ');
simc(['$cat zzfile.t >>' outpipe]);
delete('zzfile.txt');
eval(['!cat < ' outpipe ' > zzfile.txt']);
load('zzfile.txt');
res = zzfile;

```

---

```

function res = getsimvec(file)

%GETSIMVEC Gets the simulation values from a file stored during a
%           Simnon simulation.
%
%           res = getsimvec(file)
%           file = the store file where the values are stored.
%                (default: 'store')

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Thu Jan 7 11:42:29 1993

```

```

if nargin==0, file = 'store'; end;
global outpipe1

simc('write (dk zzfile FF)');
simc(['export zzfile <' file ' /0']);
simc(['$cat zzfile.t >> ' outpipe1]);
delete('zzfile.txt');
eval(['!cat < ' outpipe1 ' > zzfile.txt']);
load('zzfile.txt');
res = zzfile;

```

---

The following two commands implement the Simnon commands 'simu' and 'par' in Matlab. If 'simu' has an output argument then a 'getsimvec' is automatically given to get the simulation result back to Matlab.

---

```

function res = simu(t0, t1, dt)

%SIMU Makes the simu command in simnon. The simulation result is
%      returned if there is an out parameter.
%
%      y = SIMU(t0, t1, dt)

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Sat Dec 5 16:29:50 1992

if nargin==2,
    par('dt[logger]', (t1-t0)/100);
    simc(['simu ' num2str(t0) ' ' num2str(t1)]);
else
    par('dt[logger]', dt);
    simc(['simu ' num2str(t0) ' ' num2str(t1) ' ' num2str(dt)]);
end;
if nargout==1, res = getsimvec; end;

```

---

```

function par(p1, v1, p2, v2, p3, v3, p4, v4, p5, v5)

%PAR Set parameters in Simnon. The parameter name is given as a
%      string and its value as a number.
%
%      PAR(p1, v1)
%      .
%      .
%      .
%      PAR(p1, v1, p2, v2, p3, v3, p4, v4, p5, v5)
%      pi = the parameter name expressed as a string
%      vi = the numerical value of the parameter

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Tue Dec 8 13:39:30 1992

```

```

for ix=1:(nargin/2),
    tmp1 = eval(sprintf('p%g', ix));
    tmp2 = numtostr(eval(sprintf('v%g', ix)));
    simc(['par ' tmp1 ' : ' tmp2]);

```

```
end;
```

---

Finally, a command for transferring the standard controller data structure to the PID controller parameters in Simnon. This function is only given as an example, in general a function like this is very dependent on how the controller is implemented in Simnon.

---

```
function setpid(k, ti, td, p4, p5, p6, p7, p8, p9, p10, p11)

%SETPID Set the PID controller parameters in the controller in
%       Simnon.
%
%       SETPID(con)
%       SETPID(k, ti)
%       SETPID(k, ti, td)

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%Lund Institute of Technology, Lund, Sweden
%
%LastEditDate : Sat Dec 5 16:30:53 1992

global simdir
if rows(k)>1,
    disp('More than one row in the gain vector, using the last.');
```

```
    k = k(rows(k), :);
end;
if nargin==1 & (cols(k) > 1),
    tmp = k;
    k = getk(tmp);
    ki = getki(tmp);
    ti = getti(tmp);
    td = gettd(tmp);
elseif nargin==2,
    td = 0;
end;

mn = -1e4; mx = 1e4;
k = max(mn, min(k, mx));

fn = [simdir 'pargen.t'];
delete(fn);
fd = fopen(fn, 'w');
fprintf(fd, 'macro pargen\n');
```

```
kcp = 1;
if ti==Inf | ki == 0, ti = 1; kci = 0; else kci = 1; end;
if td==0, kcd = 0; td = 1; else kcd = 1; end;

fn = [simdir 'pargen.t'];
delete(fn);
fd = fopen(fn, 'w');
fprintf(fd, 'macro pargen\n');
fprintf(fd, 'par k: %g\n', k);
fprintf(fd, 'par kcp: %g\n', kcp);
fprintf(fd, 'par ti: %g\n', ti);
fprintf(fd, 'par kci: %g\n', kci);
fprintf(fd, 'par kcd: %g\n', kcd);
fprintf(fd, 'par td: %g\n', td);
```

```
for ix=4:2:nargin,  
    fprintf(fd, ['par ' eval(sprintf('p%g', ix)) ' : ' ...  
               num2str(eval(sprintf('p%g', ix+1))) '\n']);  
end;  
fprintf(fd, 'end\n');  
fclose(fd);  
simc('pargen');
```

---

## 6. Conclusions

The Matlab toolbox presented in this report implements a simple and systematic tuning procedure for PID controllers, with a good design variable. The design routines are modular and has proven to be easy to extend to implement new design ideas.

The design is built on a string representation of transfer functions, and can in principle handle any transfer function. All optimization and equation solution can be carried out with an arbitrary precision.

The design method computes a controller which is a compromise between robustness (the condition on  $M_s$ ) and performance (the maximization of  $k_i$ ). The method can handle a broad spectrum of dynamics uniformly in the sense that the time and frequency responses of the controlled systems will look very similar. It is, however, difficult to investigate the closed loop systems obtained with the method analytically, except in the simplest cases.

### Systems with poorly damped poles

Some care must be taken if the design method presented is used for systems with poorly damped poles. The principle to maximize  $k_i$  for a given value of  $M_s$  is still a fruitful one, but the routines 'pidesms' and 'piddesms' cannot be used for systems with poorly damped poles. It may work for PID controllers, but certainly not for PI controllers. The reason is that the approximations for finding  $\zeta_0$  are not longer valid. Instead we must use the general routines in 'dptable2', but this gives rise to several computational problems. It is also important to choose the design parameter  $\alpha_0$  properly for oscillating systems,  $\alpha_0 < 0.5$ . These problems are currently (December 1992) topics for further research.

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