Introduction to Matlab for Engineers

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Preface

These lecture notes are developed for the introductory course in Matlab for Master of Science students at Telemark University College. The lecture notes consists of three parts, and the presentation of each part is scheduled to take 8 hours and to be presented within a single week. Thus, a total of three weeks spread over the opening of the fall semester are used to present the content of the course, with a total of ca. 24 hours of arranged presentations (lectures + exercises). The presentations are integrated with exercises in the courses Numerical Methods, and Modeling of Dynamic Systems, but Matlab is also used in other courses to reinforce the learning.

The choice of course content is the result of a couple of meetings between Are Mjaavatten, Bernt Lie, and Randi Holta in the early summer of 2003. Our background experiences for composing the course are:

- Continuous use of Matlab for some 10 years,
- The arrangement of a NIF³ course in the use of Matlab (BL), where the course content was highly inspired by talks with student representatives Marte S. Lerdal and Bjørn Erik Thorsteinsen in June 1999,
- Experience with using Matlab in courses in Numerical Methods (AM) and Modeling of Dynamic Systems (BL).

Compared to the lecture notes for 2004, we have moved some of the material which is not of immediate use to appendices. We have also marked material which is more advanced with an asterisk (*). We encourage readers to give feedback on the course content and the presentation style.

Finally, it should be mentioned that there exist a useful freeware version of Matlab, named Octave. A brief introduction to the use of Octave and how it deviates from using MathWorks' Matlab version, see www.techteach.no/octave/index.htm. For more information on Octave, see www.octave.org.

Are Mjaavatten, Bernt Lie Porsgrunn, August 27 2005

 $^{^{3}}$ NIF = Norske Sivilingeniørers Forening, which has been transformed into Tekna.

Part I

Analysis of technical data with Matlab

Chapter 1

The Matlab workplace

The Matlab graphical user interface is programmed in Java, and thus looks practically the same on all supported computer platforms. We assume that the reader knows how to start up a program on her/his respective computer. When Matlab is started, it looks more or less as in fig. 1.1 (shows Matlab v. 6.5).

The main window in fig. 1.1 is the *Command Window*. This is where commands are written, and we will show excerpts of such command list in these notes. As an introductory example, assume that we type the command:

>> sin(0.7)

Here, the symbol >> is the Matlab command prompt, and the user does not type this symbol. We simply typed sin(0.7). Matlab responds with:

ans =

0.6442

>>

Here, 0.6442 is Matlab's evaluation of sin(0.7), and the symbol >> signifies that Matlab's Command Window is ready to accept another command. In the Matlab window, we now have the following situation, fig. 1.2.

Note that by double-clicking on the item sin(0.7) in the Command History window, this command is copied into the Command Window, and is immediately (re-) executed. Also, by double-clicking on the item ans in the Workspace, the so-called Array Editor is opened, containing the data of variable ans. The Array Editor is a spreadsheet-like window which makes it possible to interactively modify the content of the chosen variable.

Workspace: table of user defined variables and data structures Click here to	o change current directory
A MATLAB	
Eile Edit View Web Window Help	
🗋 🗅 😅 🐰 🛅 💼 🗠 🖓 👸 🥐 Current Directory: 🕬	MATLAB6p5\work
Workspace R	Command Window
🚰 🛛 🗄 Stack: Base 🗸	Using Toolbox Path Cache. Type "help toolbox_path_cache" f
Name Size Bytes glass	To get started, select "MATLAB Help" from the Help menu.
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Vvorkspace Current Directory	
Command History	
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bench	
% 6/23/03 2:28 PM%	
bench	
% 7/12/03 12:57 PM%	
bench	
x=linspace(0,2*pi);	
plot(x,sin(x))	
% 7/12/03 1:46 PM%	
A Start	
/	
/	Command window: type commands here
Command history: previous commands	

Figure 1.1: The opening window in Matlab 6.5. It looks almost the same in Matlab 7.0.

Variable is named "an is listed in Wo		Command and re Command	
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		ans =	
		0.6442	
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Command History	N		
bench	~		
% 6/23/03 2:28 PM	*		
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\$ 7/12/03 12:57 PM	4		
bench			
x=linspace(0,2*pi);			
plot(x,sin(x)) % 7/12/03 1:46 PM			
sin(0.7)	`		
	×	<	
Start			
The latest command is	listed in		
Command History			

Figure 1.2: The Matlab window after entering a command.

Chapter 2

Arrays: Matlab's basic data structure

2.1 Overview of learning goals

After having completed this chapter, you should have a clear understanding of what an array is, what distinguishes an array from vectors and matrices, and how you can operate on arrays. In particular, you should master:

- How to define variables by assigning a value to a name, and how to save, load, and import data,
- The creation of arrays, how to pick out subarrays and how to build superarrays,
- How to operate on variables using functions and operators,
- How to adjust the presentation format of numbers,
- Basic housekeeping of variables, commands, and files,
- How to use Matlab's help system.

2.2 Arrays and data structures

By array, we will mean a *table of numbers* similar to a spreadsheet, where we require that every element in the table contains a number. Thus, there can be no vacant elements, and there can be no text strings or other objects in any of the elements. In principle, an array can hold a single element. In that case, we will use the term *scalar* to denote the array. If the array consists of a single row of numbers, we will denote it a *row array*. If the array consists of a single column of numbers, we will denote it a *column array*. A scalar is a special case of both a row array and of a column array. If the array contains several rows (or columns) of numbers where each row (column) does not have to be a scalar, then we will simply call it a (two dimensional) *array*. Obviously, both row and column arrays are arrays.

Even more complicated arrays could be imagined. We can consider arrays consisting of several layers of two dimensional arrays, much like a spreadsheet can contain several sheets of spreadsheets. In Matlab parlance, we say that the array consists of several *pages* of (two dimensional) arrays. Together these pages could be considered a volume. We can also consider arrays with several *volumes*. Together, several volumes could constitute a library. Then we can imagine arrays consisting of several libraries. And so on. Finally, we can also imagine arrays which contains zero elements; these we will denote *empty arrays*.

Originally, Matlab only supported two dimensional arrays (and character strings). From version 5 of of Matlab, more general arrays are supported, as well as other *data structures* — these other data structures can be considered various types of tables. Often, row and column arrays are denoted row and column *vectors* in Matlab literature. Likewise, two dimensional arrays are often denoted *matrices* (Matlab was originally short for MATrix LABoratory). In these notes, we prefer to use the word array. The reason is that vectors are mathematical objects with very specific properties; vectors are not necessarily arrays. Likewise, matrices are mathematical operators with specific properties (matrices are arrays, however). Thus, we will use the term *array* when we do not imply such specific properties.

Exercise 2.1 Write down examples of various types of arrays.

Exercise 2.2 Write down examples of tables which are not arrays.

2.3 Assignment: Naming data structures

During computations, it is very important to be able to give names to data structures, e.g. to arrays. In practice, some data structures have constant value during these computations, while others vary/are changed. In this section, we will look at how we can give names (or: *assign* names) to data structures in Matlab. Some computer languages distinguish between how constants and variables are named — in Matlab, there is no such distinction

We have already seen that Matlab automatically gives a name to a result from a computation, see e.g. fig. 1.2 where the result of a computation was given the name **ans**. In practice, we need to be able to decide our own names for the data structures.

Various computer languages differ in how data structures are assigned a name. Symbolically, we could describe the assignment as <data structure> -> <name>, meaning "put the content of the data structure <data structure> into a computer memory location which bears the name <name>". In most computer languages, this assignment is written the reverse way as <name> <- <data structure>, and other symbols may be used for the assignment. In e.g. the computer algebra system Maple, the assignment is written <name> := <data structure>.

In *Matlab*, the assignment symbol is identical to the symbol which is normally used for writing equality in mathematics: <name> = <data structure>:

Here¹, **a** and **b** are the names we want to use for the data structures, while 0.7 and sin(a) (where **a** is a replacement for 0.7) are the data structures (they are arrays, or more precisely: scalars). In the assignment it is important both that the name we want to use is a valid name (see below), and that the data structure has a value (e.g. 0.7, and **a** as a replacement for 0.7). If we switch the sequence of the above commands, an error occurs:

```
>> clear all
>> b = sin(a)
??? Undefined function or variable 'a'.
>> a = 0.7
a =
0.7000
```

At first, name a does not have a value, hence sin(a) does not have a value when we try to assign sin(a) to b. We'll come back to the effect of command clear all later.

The chosen assignment symbol in Matlab (and many other languages) may be confusing at first. Consider a common type of assignment:

>> i = i+1;

¹To get a compact response in the Command Window, the command format compact has first been issued.

This assignment looks strange: how can i be equal to i+1? Well, it can not! The point is that i=i+1 does not mean that i is equal to i+1. Instead, it means: add 1 to the value of data structure i (assuming that a i already has a value), and put the result into data structure i. This is somewhat clearer if we write the assignment as $i+1 \rightarrow i$, or as in Maple: i := i+1. Still, as long as we accept that = means assignment and not equality in Matlab, the statement i = i+1 makes sense:

>> i = 1; >> i = i+1 i = 2

Valid names for data structures in Matlab can consist of *any number of characters* as long as they *start with a letter*, and the *subsequent characters* are either letters in the English alphabet (a, ..., z, A, ..., Z), digits (0, ..., 9), or underscore (_). Note that Matlab distinguishes between lowercase and uppercase letters. See Section 2.4 below for some more technical details about variable names.

Exercise 2.3 (Assigning and changing the value of a variable) Open Matlab and make sure you see the Command Window, the Workspace Window and the Command History Window. In the command window, write x = 3. Then press the "Enter" key and watch what happens in the workspace window. Next write:

x = x + 1

x = x + 1

Make sure you understand what is going on! Now write:

x = 3 * x - 4

Note that Matlab does not interpret this as an equation; interpreted as an equation, the answer would be x = 2. Instead, Matlab multiplies 3 with the current value of x (which is 5) and subtracts 4. The result (11) is then assigned to x. The interpretation of the = sign is thus: take the result of the instructions to the right and put the result into the variable at the left.

2.4 Naming data structures: some technicalities*

Although names of data structures can consist of any number of characters, Matlab only uses the N first letters to distinguish between names. The number N may vary from computer platform to computer platform, and the number is found by issuing the command namelengthmax in Matlab's Command Window. Here is the response from the author's computer when issuing this command:

```
>> namelengthmax
ans =
63
```

Thus, if you name one data structure by starting with 63 letters **a**, followed by digit **1**, and another data structure by starting with 63 letters **a** followed by digit **2**, Matlab will accept these names as valid names on the data structures, but Matlab will not be able to distinguish between the names. In practice, much shorter names are used.

There is a built in command isvarname in Matlab that can be used to check whether a name is valid. The following excerpt from the Matlab Command Window illustrates hot to use the command in order to check the potential names sin, namelengthmax, and 5ball:

ans = 0

The response from Matlab is 1 ("true") for legal names, and 0 ("false") for illegal names.

Although both sin and namelengthmax are legal names of data structures, it is probably not a good idea to use these names: using the names may confuse Matlab as to how the names should be interpreted. Matlab also has a number of other built in command names which should be avoided. In order to check that a potential name is not used by Matlab for other purposes (e.g. a built-in command, etc.), issue the command which -all <name>, where <name> is replaced by the potential name. Examples:

```
>> which -all sin
sin is a built-in function.
C:\MATLAB6p5\toolbox\matlab\elfun\sin.m % Shadowed
>> which -all namelengthmax
namelengthmax is a built-in function.
C:\MATLAB6p5\toolbox\matlab\lang\namelengthmax.m % Shadowed
>> which -all 5ball
5ball not found.
```

Note that **5ball** is not found; still **5ball** is an illegal name, and can not be used. It is very important to plan the names for data structures well, and to avoid names that are already used by Matlab. Consider e.g. that we have a data structure that contains the feed temperature of some process stream. We would like to denote the variable by **Tf**. Is this a good choice? Let us check:

```
>> which -all Tf
C:\MATLAB6p5\toolbox\control\control\@tf\tf.m % tf method
C:\MATLAB6p5\toolbox\control\control\@zpk\tf.m % zpk method
C:\MATLAB6p5\toolbox\control\control\@ss\tf.m % ss method
C:\MATLAB6p5\toolbox\control\control\@frd\tf.m % frd method
C:\MATLAB6p5\toolbox\ident\ident\@idmodel\tf.m % idmodel method
C:\MATLAB6p5\toolbox\ident\ident\@idfrd\tf.m % idfrd method
```

No, it is not a good choice. What about using T_f as name of the data structure instead?

>> which -all T_f T_f not found

Hence, T_f is both unused and a legal variable name, and can thus be used as the name for the data structure of feed temperatures.

Note that the built-in functions may vary from Matlab setup to setup: Depending on which *Toolboxes*² are installed on your computer, you may or may not get the same response as above on your computer.

Although we have used scalars as values for data structures in examples above, the right hand side (rhs) in the assignment (<name> = <data structure>) can be any data structure with value.

2.5 Built-in constants in Matlab

Matlab has some built-in names for constants. We have already seen the constant namelengthmax (in reality, this is not an assigned scalar, but a so-called function — more about functions later). Usually, it is a bad idea to "overwrite" the names of built-in constants, i.e. to use names for our own variables/constants which coincide with built-in Matlab names. Some built-in Matlab constants/names are shown in Table 2.1.

 $^{^{2}}$ A Toolbox is a collection of Matlab functions (or: classes) with a somewhat specialized common topic. Most Toolboxes are for sale at a cost that comes in addition to the cost of basic Matlab. Examples of Toolboxes are the Control Toolbox, the Statistics Toolbox, etc.

Name	Math name	Value	Description
i	i	$\sqrt{-1}$	used in Mathematics
j	j	$\sqrt{-1}$	used in Electrical Engineering
pi	π	—	
eps		2.2402×10^{-16}	Smallest number in Matlab > 0
inf	∞		e.g., 1/0
NaN			e.g., $0/0$, Not a Number

Table 2.1: Some built-in Matlab constants

The following excerpt from Matlab's Command Window illustrates how these constants are used.

```
>> i
ans =
        0 + 1.0000i
>> j
ans =
        0 + 1.0000i
>> pi
ans =
    3.1416
>> eps
ans =
  2.2204e-016
>> NaN
ans =
   NaN
>> inf
ans =
   Inf
>> 1/0
Warning: Divide by zero.
ans =
   Inf
>> 0/0
Warning: Divide by zero.
ans =
  NaN
```

Although it is possible to re-assign these constant names, we strongly recommend that this is avoided. Whether a name is a built in constant or not, can also be checked by the statement which -all <name>, as illustrated earlier.

2.6 Basic creation of arrays

2.6.1 Dimension of arrays

When organizing arrays, one could consider a scalar to be a *zero-dimensional* array, row arrays and column arrays to be *one-dimensional* arrays, and two dimensional arrays in general to be *two-dimensional* arrays. What is the significance of such a notion of dimension?

Consider the two dimensional array in Table 2.2.

Why is this array two-dimensional? Because we (normally) need two numbers to address each element. In Matlab, rows and columns are numbered with integers starting from 1. Furthermore, when addressing the elements, the row number is given first, and then the column number, e.g. the content of element (1,1) is 1, while the content of element (2,2) is -2×10^{-3} .

However, if we consider only the first column in Table 2.2, this column is one-dimensional because we only need a single number (the row number) to address each element, the content of element (2) of

	Column 1	Column 2
$\mathbf{Row} \ 1$	1	2
$\mathbf{Row} \ 2$	3.7	-2×10^{-3}
Row 3	8	1.3×10^5

Table 2.2: Example array which we want to name A.

the first column is 3.7. Similarly, a row array is one dimensional. If we consider the scalar in the element of the third row and the second column (value: 1.3×10^5), we don't need any number to address this element.

Assuming that we want to name the two dimensional array of Table 2.2 by A, how do we assign value to A? There are (at least) two ways to do this. We can give the value of each element in separate command lines — it doesn't really matter much which element we start with. Or we can assign the whole data structure in one batch.

2.6.2 Element-wise assignment

Let us start by assigning element no. (3, 1):

```
>> A(3,1) = 8
A =
0
8
```

Notice that Matlab immediately realized that since the row and column numbers must start with digit 1, A must at least be a 3×1 array. Since we, the user, so far has only specified the value of element (3, 1), Matlab is designed to assume that all the other elements have value 0.

Next, let us give the value of element (2,2):

```
>> A(2,2) = -2e-3

A = 0 0 0

0 -0.0020

8.0000 0
```

Notice that -2e-3 is Matlab syntax for -2×10^{-3} .³ This time, Matlab realizes that the element must have two columns. Since Matlab already knows that there are three columns, the array is augmented to become a 3×2 array, which is the final size in this case. Thus filling the remaining values will not change the size of the array:

```
>> A(1,1) = 1;
>> A(2,1) = 3.7;
>> A(1,2) = [2];
>> A(3,2) = 1.3e5
A =
1.0e+005 *
0.0000 0.0000
0.0000 -0.0000
0.0001 1.3000
```

Note:

 $^{^{3}-2 \}times 10^{-3}$ can also be written as -2*10⁽⁻³⁾ in Matlab, but the notation -2e-3 is clearly the preferred (and simplest!) notation.

- When we end a command line with a *semicolon* (";"), the result is not printed in the Command Window. If we do *not* end the command line with semicolon, the result is printed.
- Since a scalar is a zero-dimensional array, there is no problem in specifying the scalar as either 2 or [2].
- Even though it may seem like only two elements in the final result contain values different from zero, this is not so: here, all elements have values that differ from zero. The reason why 4 out of 6 elements appear to be zero, is the way Matlab by default formats the numbers. We will return to how these numbers can be formatted.

Suppose we want to change the value of one element, e.g. assume that element (2,2) should be $+2 \times 10^{-3}$ instead of the current value. We then simply rewrite the value:

```
>> A(2,2) = 2e-3
A =
    1.0e+005 *
    0.0000    0.0000
    0.0000    0.0000
    0.0001    1.3000
```

We can also check the value of a certain element, e.g. element (2,2), as follows:

>> A(2,2) ans = 0.0020

Finally, it should be commented that this element-wise way of giving values to A is somewhat inefficient in that we have changed the size of the array twice: the first time when we gave the value of element (3, 1) and thus creating a 3×1 array, and the second time when we gave the value of element (2, 2) and thus augmented the array to a 3×2 array.

2.6.3 Assigning row arrays

If we consider the two dimensional array of Table 2.2, we can define three row arrays:

Here we notice that arrays are indicated by including the numbers in square braces [and]. Furthermore, we can separate the various elements of each row either by comma, or by a space. Since each of the row arrays are one-dimensional, it suffices to specify one number to indicate which element we want to check, e.g. $A_r2(2)$.

2.6.4 Assigning column arrays

Column arrays are constructed similar to the way row arrays are constructed. However, for column arrays, either semicolon or Carriage Return (CR) is used to separate the elements:

```
>> A_c1 = [1; 3.7; 8]
A_c1 =
    1.0000
    3.7000
    8.0000
>> A_c2 = [2
-2e-3
     1.3e5]
A_c2 =
  1.0e+005 *
    0.0000
   -0.0000
    1.3000
>> A_c2(2)
ans =
   -0.0020
```

Alternatively, we can pretend that the column array is a column matrix, and use the operation of *transposing* the matrix to turn the row array into a column array. The matrix transpose of a matrix M, M^T , is performed using the notation M' in Matlab⁴. Thus, we would have defined e.g. A_c2 by first creating it as a row array, and then transposing the result:

```
>> A_c2 = [2, -2e-3, 1.3e5]'
A_c2 =
    1.0e+005 *
    0.0000
    -0.0000
    1.3000
```

The techniques of either separating elements by semicolon, or by transposing a row array, are normally preferred since fewer lines in the Command Window are needed. However, separating elements by CR makes the Command Window more readable.

2.6.5 Assigning two-dimensional arrays

We now consider the two dimensional array in Table 2.2 as a number of rows below each other. The following notation is used to assign the value of the array to the name of the array:

```
>> A = [1, 2; 3.7 -2e-3;
8, 1.3e5]
A =
    1.0e+005 *
    0.0000    0.0000
    0.0000    -0.0000
    0.0001    1.3000
```

Notice that we list the first row first (either separated by comma or space), then separate the first row from the second row (either by semicolon or CR), and so on. Often, the two dimensional array is easier to read if CR is used to separate rows.

 $^{^{4}}$ Actually, operator "," takes both the transpose, and the conjugate of matrix M. We will return with a more precise description.

Exercise 2.4 (Chaos) Write:

x = 0.7

x = 2 * x * (1 - x)

Repeat the last instruction several times. (Use the up arrow!) Do you see any relationship between the result and the equation?

We can make a side-by-side comparison of the iteration sequence from two nearby initial values for x (e.g. 0.7 and 0.7001) by

 $x = [0.7 \ 0.7001]$

x = 2 * x . * (1 - x)

Note the .* operator, signifying element by element multiplication rather than the standard matrix multiplication! Again, repeat the last command several times to see how x changes.

Now, try with a bigger coefficient in the right-hand side:

 $x = [0.7 \ 0.7001]$

x = 4*x.*(1-x)

Note that when we repeat the last command, the values do not tend towards a constant, but rather flutter wildly about. This is an example of *chaotic behaviour*.

The results from the two close starting points diverge rapidly. In a chaotic system, a small change in the initial value may yield very different long-time results. This is called "the butterfly effect": if the weather is a chaotic phenomenon, then the very small change in the motion of the air caused by the flapping of a butterfly's wings in Brazil might in theory trigger a storm in the Indian Ocean some weeks or months later. This effect may indeed limit the feasibility of long-range weather forecasts.

2.6.6 Line continuation

For large arrays, it may be difficult to fit one row on one command line. We may then split the row across several command lines by using the line continuation operator ...:

```
>> A = [1, 2; 3.7 ...
-2e-3; 8, 1.3e5]
A =
1.0e+005 *
0.0000 0.0000
0.0000 -0.0000
0.0001 1.3000
```

Here, Matlab will consider the assignment as having taken place on a single command line.

2.6.7 One-dimensional addressing of array elements*

For any array, we can address the element element by a single number. For a scalar, the number is 1. For row or column arrays, the number is the location of the element in the array. For two-dimensional arrays, we imagine that the columns are stacked on top of each other, with the first column on the top, and the last column on the bottom.

```
>> A
A =
    1.0e+005 *
    0.0000    0.0000
    0.0001    1.3000
>> A(6)
ans =
        130000
>> A(5)
ans =
        -0.0020
```

Usually, it is more complicated to figure out the element address using a single number address.

2.7 Exporting and importing data from files

It is important to be able to save data from the workspace/a Matlab session to a file, so that we can easily import it again at a later time. It may also be important to be able to export the data to some file format that can be imported into other programs, or to import data from other programs such as Excel, etc. Here, we will only discuss saving data to and loading data from basic Matlab formats, and import of data from Excel.⁵

2.7.1 Saving and loading data

Suppose we have two arrays, A and B:

$$A = \begin{pmatrix} 1 & 2 \\ 3 & 4 \end{pmatrix}, \qquad B = \begin{pmatrix} 824 & -65 & -814 & -741 \\ -979 & -764 & 216 & 663 \\ 880 & 916 & 617 & -535 \end{pmatrix}.$$

We create these in Matlab as follows:

Suppose these two data structures are the only ones in the workspace of Matlab at the time being. First, we want to save all data structures of the workspace to a file. Before doing so, we need to determine:

- 1. The location of the file on the computer. The default directory where the file will be placed, is the *Current Directory*, see fig. 1.1 on p. 4. Clicking the down-arrow (\vee) for the Current Directory brings up a list of previously used directories, which can be selected. Clicking the ellipsis icon (...)to the right of the Current Directory list, brings up a file browser that can be used to select the directory.
- 2. What file type to save the workspace in. The default file type is Matlab's own .mat type. The alternative is to save the workspace as an .ascii file. Note that the .mat type works both faster, and preserves more information.

The basic syntax for saving files is save <filename> {-<filetype>}, where <filetype> can be either mat or ascii (mat is the default), and {} indicates that what is between these braces is optional. Here is an example of saving the workspace to both file types:

```
>> save myworkspacem1 -mat
>> save myworkspacea -ascii
>> save myworkspacem2
```

Note that saving to myworkspacem2 also leads to a file of type .mat, i.e. myworkspacem2.mat. The reason is that .mat is the default file type.

Next, imagine that we clear the workspace so that there is nothing there. This will typically be the case when we start up Matlab, but we can also enforce an empty workspace (more about that later). So, the workspace is empty. Let us then load the file containing the workspace we just saved. The syntax is load $\{-<filetype>\}$ $<filename>^6$. We first see what happens when we load the .mat files.

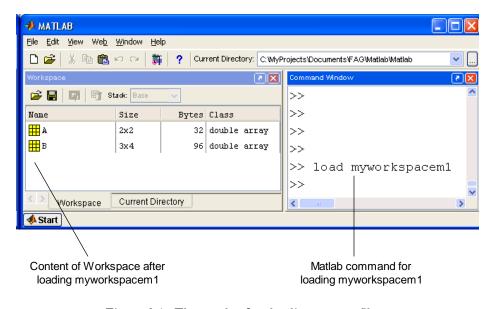


Figure 2.1: The result after loading a .mat file.

The result is shown in fig. 2.1. We see that the original variable names are preserved. What then if we load a .ascii file? It turns out that Matlab can not load the file myworkspacea. The reason is that the arrays have a different number of columns:

```
>> load myworkspacea
??? Error using ==> load
Number of columns on line 3 of ASCII file
C:\MyProjects\Documents\FAG\Matlab\Matlab\myworkspaceamust be the same as
previous lines.
```

See FILEFORMATS for a list of known file types and the functions used to read them.

If we open the file in a text editor (e.g. Notepad, or the Matlab editor), the ASCII file looks as in fig. 2.2. This indicates that the ASCII format is not useful for storing more than one data structure, and perhaps it is not suitable at all for some data structures. Note also that there is no information about the array names in the file.

It is possible to use more details when saving and loading the workspace. We can choose to only save part of the workspace. The syntax is: save <filename> {<variablename1> ... <variablename1> ... <variablenameN>} {-<filetype>}, where {<variablename1> ... <variablenameN>} is meant to indicate that we can list any number of variable names, and that they should be separated by *space* (*not* comma!). Similarly, the syntax for loading files is load {-<filetype>} <filename> {<variablename1> ... <variablename1> ... <variablena

```
>> save myworkspacem
>> save myworkspacemA A
>> save myworkspacemB B
>> save myworkspaceaA A -ascii
>> save myworkspaceaB B -ascii
```

⁵It is possible to import data using commands such as fileread, fread, etc.

 $^{^{6}}$ Note that for saving files, the file type is given after the filename, while for loading files, the file type is given before the filename.

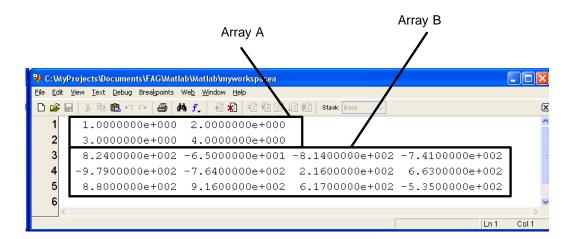


Figure 2.2: The content of the ASCII file. Note that the first two rows come from the A array, while the three next rows come from the B array.

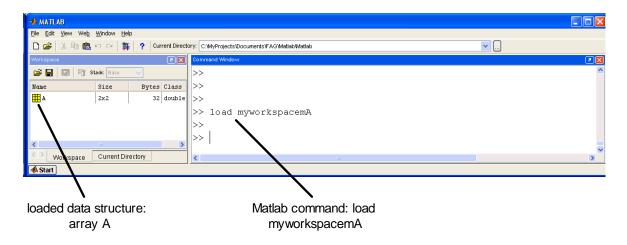


Figure 2.3: Situation after loading file myworkspacemA (a .mat file), which only contains array A.

This leads to the following files:

19.07.2003	16:01	66	myworkspaceaA
19.07.2003	16:01	195	myworkspaceaB
19.07.2003	16:01	264	myworkspacem.mat
19.07.2003	16:01	184	myworkspacemA.mat
19.07.2003	16:01	208	myworkspacemB.mat
	5 File(s)		917 bytes

What happens if we start with an empty workspace and load these files? The following examples illustrates what takes place.

>> load myworkspacemA

This operation leads to the situation in fig. 2.3. Similarly:

>> load myworkspaceaB

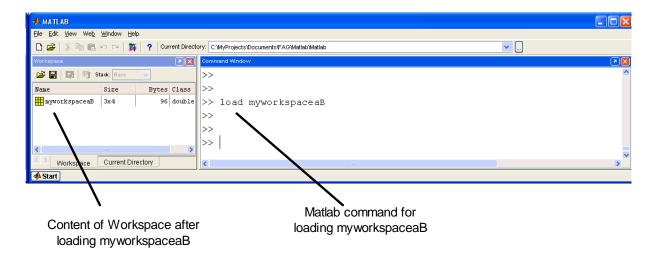


Figure 2.4: Situation after loading file myworkspaceaB (an .ascii file), which only contains array B.

This operation leads to the situation in fig. 2.4. Note in particular that since the ASCII file does not contain information about the name of the stored array, the loaded variable takes its name from the loaded file. Thus, what used to be array B now is named myworkspaceaB.

2.7.2 File import

The simplest way of importing data from Excel is via the Matlab File/Import Data... command. This command can also be used to import data saved using the save command of Matlab (see previous subsection). This command simply opens a file browser which shows files that Matlab recognizes. Suppose we have an Excel spreadsheet as in fig. 2.5. In the Matlab window, we choose the command File/Data Import..., which opens a file browser.⁷ We choose the relevant file (MatlabImportData.xls, see fig. 2.5), and click Open. The Import Wizard is opened, and leads to the result in fig. 2.6. Clicking Finish in the Matlab Import Wizard populates the Matlab workspace with the variables data, textdata, and colheaders. These data look as follows:

```
>> data
data =
             78.0000
    1.0000
                         0.3000
    2.0000
             79.0000
                         0.4000
    3.0000
             82.0000
                         0.5500
             85.0000
    4.0000
                         0.5700
                         0.5900
    5.0000
             92.0000
>> colheaders
colheaders =
    'Sample no.'
                     'Temperature'
                                       'Mole fraction'
>> textdata
textdata =
    'Sample no.'
                     'Temperature'
                                       'Mole fraction'
```

So far, we have learned how to *build* the data structure (array) of variable data (but not how to build the data structures colheaders and textdata — which are cell arrays). In order to analyze the data contained in variable data, we need to be able to deconstruct the variable, and pick out elements of the array. We'll get back to that problem soon.

 $^{^{7}}$ It is required that Excel is installed on your computer for this to work: Matlab communicates with Excel in order to import the data.

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2							
	A1	▼ fs	Sample no.				
	А	В	С	~			
1	Sample no.	Temperature	Mole fraction				
2	1	78.00	0.30				
3	2	79.00	0.40				
4	3	82.00	0.55				
5	4 85.00 0.57		0.57				
6	5	92.00	0.59	~			
K (H + H Sheet1 Sheet: < >						
	Sum=433.41						

Figure 2.5: Excel spreadsheet with data that we want to import into Matlab.

📣 Import Wizard									
Select value	Select variables to import using checkboxes								
 Create 	● Create variables matching preview.								
O <u>C</u> reat	O <u>C</u> reate vectors from each column using column names.								
○ C <u>r</u> eate	e vectors fro	m each rov	v using row	names.					
Variables	C Variables in C:\Documents and Settings\berntl\My Documents\ArchiveDocuments\Courses\MSc\Matlab\D;								
Import	Name 🔺	Size	Bytes	Class		1	2	3	
	Colhe	1x3	248	cell	1	1.0	78.0	0.3	
	🗄 data	5x3	120	double	2	2.0	79.0	0.4	
	🖸 textd	1x3	248	cell	3	3.0	82.0	0.55	
					4	4.0	85.0	0.57	
					5	5.0	92.0	0.59	
<		Ш		>					-
Help						< Back	Next >	Finish	

Figure 2.6: The result of the Matlab Import Wizard.

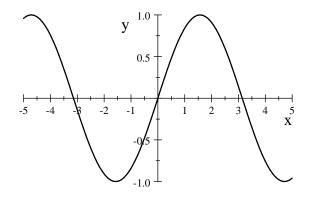


Figure 2.7: The function $y = \sin x$, where x is the input argument, and y is the output argument. This is an example of a relationship between x and y, which is also a function.

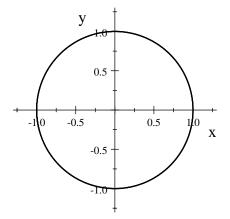


Figure 2.8: Implicit plot of $x^2 + y^2 = 1$. This is an example of a relationship between x and y which is not a function.

2.8 Functions and operations

The basic idea of *functions* is known from high school mathematics: A function is an operation that maps its input argument(s) uniquely into its output argument(s). The basic property is the uniqueness: with a specified (input) argument, the answer (output argument) is uniquely determined. There is no room for doubt or randomness. An example of a function is $y = \sin x$, which is shown in fig. 2.7. When we fix the input argument (the value of x), the value for the output argument (y) is given uniquely.

An example of a curve that does not satisfy the property of a function, is the implicit plot of $x^2 + y^2 = 1$, fig. 2.8. This is not a function simply because if we fix the input argument (e.g. x), there is not a unique output argument (y when x is the input argument). This non-uniqueness causes the mapping between x and y to not be a function.

Matlab has many built-in functions. The typical structure of calling these functions are: output_argument = function_name(input_argument). The function name must satisfy the requirements for Matlab variable names. The input and output arguments can be any data structure, and can also be lists (comma separated) of variables/data structures. There can even be a variable number of input and output arguments, depending on the properties of the input arguments and how much details of the output arguments the user is interested in. The requirement for making these functions is still that when the input arguments are given, there is a unique answer (output argument).

Let us consider what takes place during the common mathematical operation of multiplication, e.g. consider

y = 3x.

Here, the input argument x is multiplied by the number 3 in order to obtain the output argument y.

There is a unique relationship between x and y, so this is a function. In order to compute the output argument in Matlab, we can use e.g. the built-in function c = mtimes(a,b):

```
>> x = 5;
>> y = mtimes(3,x)
y =
15
```

This is a simple example of using a function in Matlab. There are some restrictions on what the input arguments can be for this function to work, but let us not worry about this now.

In the particular case of function **mtimes** (and a few other functions), we know from mathematics that it is more convenient to introduce the *operator* multiplication. Thus, we can instead express $y = 3 \cdot x$ as follows:

>> x = 5; >> y = 3*x y = 15

Every operation in Matlab has an equivalent function. But there are many functions that do not have an equivalent operator. Although we in the examples above have used scalar arguments, many functions accept or need other data structures as arguments (e.g. arrays). We will get back to this later.

Exercise 2.5 Consider the mathematical expression $z = 3x - 2y + \frac{(x-y)^2}{x+y}$. Using the operator +, -, *, /, and ^, this can be written as:

 $z = 3*x - 2*y + ((x-y)^2)/(x+y)$

in Matlab. Check this, by first assigning values to x and y.

Next, try to repeat this using the corresponding functions plus, -, mtimes, mrdivide, and mpower: z = plus(mtimes(3,x), plus(mtimes(-2,y), mrdivide(mpower(plus(x,-y),2), plus(x,y))));Do you get the same result? Which formulation is simplest?

2.9 Numeric format and accuracy

All commands in Matlab are carried out in double precision⁸. 8 bytes are used to store each element in an array, according to some IEEE standard⁹. In practice, this means that there is ca. 16 decimal digits accuracy in the computations, and a range of roughly $[10^{-308}, 10^{+308}]$. Many times, the accuracy in observed data that are used in computations is much lower, but still, Matlab always works with double precision, and the full accuracy is always stored to files. We can, however, influence how the data are presented to the user in the Matlab Command Window.

Table 2.3 shows the available commands for manipulating the format of the data as they are presented to the user.

The command format compact is useful for saving space when inserting Matlab commands/responses into a word processor. As an example, consider our familiar array:

```
>> format compact
>> a = [3, 2; 3.7, -2e-3; 8, 1.3e5]
a =
    1.0e+005 *
    0.0000    0.0000
    0.0000    -0.0000
    0.0001    1.3000
```

 $^{^{8}}$ Since ca. version 6.5 of Matlab, there is also support for other data types such as single precision, integer, boolean, etc.

 $^{{}^{9}}$ IEEE = The Institute of Electrical and Electronics Engineers, Incorporated, is an engineering organization with over 365 000 members in more than 105 countries; about 40% of the members are from outside of the USA.

Table 2.3: Available commands to format data structures in Matlab.				
Command	Comment			
format	Default. Same as short (see below).			
format short	Scaled fixed point format with 5 digits.	default		
format long	Scaled fixed point format with 15 digits.			
format short e	Floating point format with 5 digits.			
format long e	Floating point format with 15 digits.			
format short g	Best of fixed or floating point format with 5 digits.			
format long g	Best of fixed or floating point format with 15 digits.			
format hex	Hexadecimal format.			
format +	The symbols +, - and blank are printed for positive, negative and			
	zero elements. Imaginary parts are ignored.			
format bank	Fixed format for dollars and cents.			
format rat	Approximation by ratio of small integers.			
format compact	Suppress extra line-feeds.			
format loose	Puts the extra line-feeds back in.	default		

This (default) formatting is with 5 digits accuracy. Let's change this to 15 digits accuracy:

```
>> format long
>> a
a =
    1.0e+005 *
    0.0000300000000    0.000020000000
    0.0000370000000    -0.0000002000000
    0.0000800000000    1.3000000000000
```

Exercise 2.6 Experiment with each of the other **format** statements using the above array **a**.

See Appendix C for a more detailed example on how Hexadecimal numbers work.

2.10 Functions for creating arrays

Table 2.4 shows some functions for creating matrices. Let us illustrate the basic use of some of these functions.

2.10.1 Creating arrays

The colon operator The colon operator is often used with integers, but can also be used with floating point numbers:

 Table 2.4: Some commands for creating and manipulating arrays in Matlab.

 Notation
 Description

Command	Notation	Description
<pre>m:k:n, colon(m,k,n)</pre>	m:k:n	Colon operator for creating row array where the first element has
		value m , the last element has value n , and the elements in between
		are evenly spaced with distance k .
linspace(x,y,n)		Creates row array with n elements, where the first element has
		value x , the last element has value y , and the elements in between
		are evenly (linearly) spaced.
logspace(x,y,n)		Creates row array with n elements, where the first element has
		value 10^x , the last element has value 10^y , and the elements in
	0	between are logarithmically spaced.
zeros(m,n)	$0_{m,n}$	$m \times n$ -array filled with 0.
ones(m,n)	$1_{m,n}$	$m \times n$ -array filled with 1.
eye(m,n)	$I_{m,n}$	$m \times n$ -array with 1 on the main diagonal, and 0 elsewhere.
diag(v)	$\operatorname{diag}\left(A,B\right)$	If v is an array of length n: creates an $n \times n$ -array with the elements
1()		of v on the main diagonal.
rand(m,n)		$m \times n$ -array with uniformly distributed random numbers in the
		intervall $(0, 1)$.
randn(m,n)		$m \times n$ -array with normally distributed random numbers $\sim \mathcal{N}(0, 1)$. Produces a row vector where the first element is the number of
<pre>size(A), size(A,n)</pre>		Froduces a row vector where the first element is the number of rows of array A , and the second number is the number of columns
		of A.
length(A)		The maximum of the two numbers in size(A).
reshape(A,k,1)		$m \times n$ -array A is reshaped into a $k \times \ell$ -array; this is only possible
100110000000000000000000000000000000000		if $m \cdot n = k \cdot \ell$.
rot90(A,k)		Array A is "rotated" $k \times 90^{\circ}$ counterclock-wise.
blkdiag(A,B)	$\operatorname{diag}(A,B)$	Block-diagonal matrix with matrices A and B on the main diagonal
0.0	0())	blocks.
<pre>toeplitz(c,r)</pre>		Produces a Toeplitz array with c as the first column, and r as the
-		first row. With a single argument, c and r are assumed equal.
hankel(c,r)		Produces a Hankel array with c as the first column, and r as the <i>last</i>
		row. With a single argument, c is given, and the array is square.
repmat(A,m,n)		Replicates a matrix: matrix A is stacked together into a super-
		matrix with m block rows of copies of A , and n block columns of
		copies of A .
kron(A,B)	$A\otimes B$	Produces the Kronecker-product of A and $B, A \otimes B$.

```
>> 1:2:6
ans =
     1
           3
                  5
>> 6:-3:-6
ans =
     6
           3
                  0
                       -3
                              -6
>> pi:pi/10:2*pi
ans =
  Columns 1 through 5
                                                 4.0841
       3.1416
                     3.4558
                                   3.7699
                                                               4.3982
  Columns 6 through 10
       4.7124
                     5.0265
                                   5.3407
                                                 5.6549
                                                                5.969
  Column 11
       6.2832
```

The colon operator is used extensively in operation on arrays, as we will see soon. Note that the end point (n of m : k : n) is only included if it can be written as $n = m + k \cdot i$ where *i* is some integer number. Also notice that when using the colon operator, we do not explicitly specify the number of elements in the row array.

The *function* for doing the same as the colon *operator*, is colon(m,k,n):

>> colon(1,2,6) ans = 1 3 5

Linspace and logspace With the linspace command, we do not specify the distance between the numbers, but we do specify the number of elements in the row array:

```
>> linspace(0,2*pi,5)
ans =
0 1.5708 3.1416 4.7124 6.2832
```

If the number of elements is left out (e.g., we write linspace(x,y)), then the default number of elements is 50.

When using the logspace command, the result is as follows:

Zeros and ones Using function zeros is exemplified by the following Matlab session:

Notice that when we only use one argument in the function, the resulting array is square. Function **ones** works similarly to function **zeros**:

Eye and diag Function eye also works similarly:

0

0

Function eye is typically used to generate the *identity matrix* in linear algebra. Function diag works as follows:

Note that diag can be used to create square "eye" arrays:

Note also that function diag can have a second argument: diag(v,k). The second argument is then an integer, and specifies whether the diagonal should be placed above the main diagonal (positive k) or below the main diagonal (negative k).

liag(1:2,-1)		
=			
0	0	0	
1	0	0	
0	2	0	
liag(1:2,2)		
=			
0	0	1	0
0	0	0	2
0	0	0	0
0	0	0	0
	= 0 1 0 diag(= 0 0	1 0 0 2 diag(1:2,2) = 0 0 0 0	= 0 0 0 1 0 0 0 2 0 diag(1:2,2) = 0 0 1 0 0 0

Finally, function diag can also be used to construct an array by *extracting* the diagonal from a matrix — the result is a column array:

```
>> A = [1, 2; 3.7 -2e-3; 8, 1.3e5]
A =
                          2
            1
          3.7
                     -0.002
            8
                   1.3e+005
>> diag(A)
ans =
             1
       -0.002
>> diag(A,-1)
ans =
          3.7
     1.3e+005
```

Random numbers Uniformly distributed random numbers are created using function rand:

0.60684	
0.48598	
0.45647	
0.018504	
0.61543	0.92181
0.79194	0.73821
	0.48598 0.45647 0.018504 0.61543

Here, we notice that in the first two calls to the function (calling rand(2) twice), the result is different each time. The reason is that the numbers are supposed to be random, and thus they should not be the same. For comparisons, we often want to enforce the results of two drawings to be the same. We can achieve this by *resetting* the algorithm with command rand('state', 0):

The random generator for *normally distributed* numbers works in the same way, e.g.:

```
>> randn(2)
ans =
-0.43256 0.12533
-1.6656 0.28768
```

2.10.2 Size of arrays

Size Sometimes we don't know the size of an array, and it is useful to find the size. This can be done by using function **size**. Suppose we have an array A with an unknown size $m \times n$. We find the size as follows:

Here, we can argue that we already knew the size: we had just created matrix A to be of size 20×30 . But we can easily get into a situation where we do not really know the size, or where it is too complicated to figure it out; thus it may be easier to use command **size** to find the size. Let us expand slightly on the example:

We see that we can easily pick out the number of rows and the number of columns.

If we only care about a specific dimension (e.g. the number of rows, or the number of columns), than we can pick out this as follows:

Length Sometimes — typically for one-dimensional arrays, we want to find the length of the array. For this case, it is simpler to use the **length** function:

```
>> v = 1:3:20;
>> length(v)
ans =
7
>> size(v)
ans =
1 7
```

Application of size Let us finally consider a simple application of the size command: suppose we have an array of unit elements (1), A, and that we want to create an array B of the same size, but with uniformly distributed random numbers. This can be achieved as follows:

```
>> A = ones(2,3)
A =
     1
            1
                  1
     1
            1
                  1
>> B = rand(size(A))
B =
     0.012863
                    0.68312
                                 0.035338
      0.38397
                   0.092842
                                    0.6124
```

2.10.3 Rearranging arrays

Reshape Sometimes, it is useful to rearrange the shape of an array. Suppose we have an array A of dimension $m \times n$, and that we want to turn this array into an array B of dimension $j \times k$. As long as $m \cdot n = j \cdot k$, arrays A and B will contain the same number of elements. When this dimension requirement $m \cdot n = j \cdot k$ is fulfilled, Matlab function reshape(A, j, k) can be used to change the shape of A from being a 3×2 array, into being a 2×3 array:

```
>> A = rand(3,2)
A =
                    0.20277
      0.81317
    0.0098613
                    0.19872
      0.13889
                    0.60379
>> B = reshape(A,2,3)
B =
      0.81317
                    0.13889
                                 0.19872
    0.0098613
                    0.20277
                                 0.60379
```

Conceptually, this command may be carried out as follows:

- 1. First, matrix A is made into a column array with $m \cdot n$ rows (and one column).
- 2. Then, the first j elements of the column array are placed in the first column of array B, the next j elements of the column array are placed in the next column of B, and so on, until B has k columns.

>> A = rand(2,3)A = 0.46599 0.84622 0.20265 0.52515 0.67214 0.41865 >> rot90(A) ans = 0.20265 0.67214 0.84622 0.52515 0.46599 0.41865 >> rot90(ans)ans = 0.52515 0.41865 0.67214 0.20265 0.84622 0.46599 >> rot90(ans)ans = 0.46599 0.41865 0.52515 0.84622 0.20265 0.67214 >> rot90(ans) ans = 0.46599 0.84622 0.20265 0.41865 0.52515 0.67214

Rotate When considering the geometric layout of an array, it is sometimes of interest to rotate this layout. Function **rot90** rotates an array counterclockwise by 90°:

We see that after 4 rotations, we are back where we started. Sometimes, it may be useful to combine the rot90 function with the operation of transposing an array.

2.11 Subarrays and superarrays

2.11.1 Picking subarrays

Indexing the array In this section, we only give a basic presentation of how to index arrays. See Appendix B.1 p. 159 for a more complete discussion.

Suppose array A is given. We often wish to pick out element (i, j) from array A. Let us denote element (i, j) of array A by $A_{i,j}$, which in Matlab is given as A(i,j).

For array A, it can also be useful to be able to pick a subarray which has indices over the rectangle given by row i_1 to row i_2 , and column j_1 to j_2 . This submatrix can be denoted as $A_{i_1:i_2,j_1:j_2}$. In Matlab, this subarray is specified as $A(i1:i_2, j_1:j_2)$:

```
>> rand('state',0)
>> A = rand(3, 4)
A =
                    0.48598
      0.95013
                                  0.45647
                                                 0.4447
      0.23114
                     0.8913
                                 0.018504
                                                0.61543
      0.60684
                     0.7621
                                  0.82141
                                               0.79194
>> A(1,3)
ans =
      0.45647
>> A(2:3,2:3)
ans =
                   0.018504
       0.8913
       0.7621
                    0.82141
```

Properties of elements Often, it is desirable to find elements of arrays with certain properties. We can use the function find to find the indices of such elements. Let us find the elements of array A with elements which satisfy A(i, j) > 0.5. This can be done as follows:

>> A				
A =				
	0.95013	0.48598	0.45647	0.4447
	0.23114	0.8913	0.018504	0.61543
	0.60684	0.7621	0.82141	0.79194
>> in	d = find(A >	0.5)		
ind =	:			
	1			
	3			
	5			
	6			
	9			
1	1			
1	2			

Clearly, here the index must be interpreted as a one-dimensional addressing of the elements of A, i.e. the row number in col (A). We can now find the value of these elements by using the following syntax:

>> A(ind) ans = 0.95013 0.60684 0.8913 0.7621 0.82141 0.61543 0.79194

We can also use the found index to "clip" the values of A such that the maximal value is 0.5:

>> A(ind) = 0.5 A =			
0.5	0.48598	0.45647	0.4447
0.23114	0.5	0.018504	0.5
0.5	0.5	0.5	0.5

2.11.2 Building superarrays

Concatenation It is possible to concatenate several arrays into superarrays. The concatenation operator is [], and this operator works just as if the arrays we concatenate, are scalars:

```
>> A = rand(2,3)
A =
      0.41027
                   0.057891
                                  0.81317
      0.89365
                    0.35287
                               0.0098613
>> B = [A, A; A, A]
B =
                   0.057891
      0.41027
                                 0.81317
                                               0.41027
                                                            0.057891
                                                                           0.81317
      0.89365
                    0.35287
                               0.0098613
                                               0.89365
                                                             0.35287
                                                                         0.0098613
      0.41027
                   0.057891
                                 0.81317
                                               0.41027
                                                            0.057891
                                                                           0.81317
```

0.89365 0.35287 0.0098613 0.89365 0.35287 0.0098613

There are some obvious restrictions on how to concatenate arrays.

- Every element of each block row must have the same number of rows.
- The total number of columns in each block row must be the same.

>> C = [[1,2;2,4], A; A, [4,7;2,2]]					
C =					
	1	2	0.41027	0.057891	0.81317
	2	4	0.89365	0.35287	0.0098613
(0.41027	0.057891	0.81317	4	7
(0.89365	0.35287	0.0098613	2	2
>> D =	[[1;2],	A, [4;2];			
	A, [4,7	; 2,2]]			
D =					
	1	0.41027	0.057891	0.81317	4
	2	0.89365	0.35287	0.0098613	2
(0.41027	0.057891	0.81317	4	7
(0.89365	0.35287	0.0098613	2	2

Note here that the first array [1,2;2,4] of block row 1 of C has the same number of rows as array A. Likewise, array A has the same number of rows as array [4,7; 2,2]. This fulfills the requirement for the rows. Finally, the number of columns for [1,2; 2,4] and A are the same as the number of columns for A and [4,7; 2,2]. This fulfills the requirement for the columns. Similarly, the requirements are fulfilled for array D.

Replication of arrays Sometimes, we need to replicate an array many times. Function repmat (replicate matrix) makes a copy of the array in question: E = repmat(A,m,n) produces $m \cdot n$ copies of array A, with m copies of A below each other, and then n copies of this stacked array beside each other:

```
>> A
A =
      0.41027
                  0.057891
                                 0.81317
                   0.35287
      0.89365
                               0.0098613
>> repmat(A,3,2)
ans =
      0.41027
                  0.057891
                                 0.81317
                                               0.41027
                                                           0.057891
                                                                          0.81317
      0.89365
                   0.35287
                               0.0098613
                                               0.89365
                                                             0.35287
                                                                        0.0098613
                  0.057891
                                                           0.057891
      0.41027
                                 0.81317
                                               0.41027
                                                                          0.81317
      0.89365
                   0.35287
                               0.0098613
                                                             0.35287
                                                                        0.0098613
                                               0.89365
      0.41027
                   0.057891
                                 0.81317
                                               0.41027
                                                           0.057891
                                                                          0.81317
      0.89365
                   0.35287
                               0.0098613
                                               0.89365
                                                             0.35287
                                                                        0.0098613
```

2.12 Housekeeping

It is important to be able to administer both variables and commands in a computational program like Matlab. The functions in Table 2.5 are useful for housekeeping in Matlab.

The use of these commands is discussed in more detail in the following sections.

30

	Table 2.5: Some commands for basic housekeeping in Matlab.
Command	Description
who	List the variables defined in Matlab's workspace.
whos	List variables and type/size in Matlab's workspace.
clear	Delete all variables from Matlab's workspace; clear a b clears a and b.
↑,↓	Use arrows to navigate in Matlab's list of previous commands.
\longrightarrow , \longleftarrow	Use arrows to navigate in the current command line for editing the command.
cd	Change directory.
delete	Delete a file.
dir	Show the files in the directory.
what	Show the <i>Matlab</i> files in the directory.
mkdir	Create a directory.

2.12.1 Variables

We have already seen that the variables that have been introduced, are listen in the *Workspace window* — see fig. 1.1. We have also indicated how we can use the Array Editor to edit the variables. Let us now see how we can address the variables from the command line.

In order to get a list of variables, we can use the commands who and whos. Command who basically lists the variables, while command whos also gives some details about the variables:

<pre>>> a = linspace(1,10,20); >> b = zeros(1,20); >> c = ones(1,20); >> d = eye(1,20); >> rand(1,20); >> who</pre>					
Your variabl	les are:				
a ans b	c d				
>> whos					
Name	Size	Bytes	Class		
a	1x20	160	double	array	
ans	1x20	160	double	array	
b	1x20	160	double	array	
с	1x20	160	double	array	
d	1x20	160	double	array	

Grand total is 100 elements using 800 bytes

Remember that each element in an array occupies 8 bytes of memory. Since we have $5 \cdot (1 \times 20) = 100$ elements, this means $100 \cdot 8 = 800$ bytes.

The variable with name **ans** always contains the latest result that has been computed.

If we want to delete a variable from memory, this is done using command clear. By simply writing the command >>clear, every variable in the workspace is deleted. If we want to specify a specific variable to delete, or a set of variables to delete, this is done as follows:

>> clear ans >> clear a b >> who

Your variables are:

c d

Notice that when listing variables that are to be cleared, these must be separated by *space* and not by comma. If we write >>clear a, b, this is interpreted as clear a and then display b in the command window.

We can even use regular expressions to clear variables:

```
>> A1 = rand(2,30);
>> A2 = rand(2,40);
>> who
Your variables are:
A1 A2 c d
>> clear A*
>> who
Your variables are:
c d
```

2.12.2 Matlab commands

As we have seen, by double-clicking on a command in Matlab's Command History window, the command is copied to the Command Window and executed.

Another way to use previous commands, is to step through the previous commands starting with the most recent command, by using arrow-up (\uparrow) or arrow-down (\downarrow) on the keyboard. The previous commands are stored in a history list, which is seen in the Command History window. When the desired command has been found by stepping up and down in the list using arrow-up or arrow-down, the command is executed by hitting the carriage return (CR) key. It is also possible to edit the commands that are called back to the command window using the arrows: use the mouse to mark possible characters to delete using either the Delete or the Backspace keys. If you want to add new text in the command, either use the mouse to position the cursor, or use arrow-left (\leftarrow) or arrow-right (\rightarrow) to position the cursor. Then type in the missing characters.

Quite often, the history list is lengthy, and one may have to hit the arrow-keys many times to find the correct command. If so, the following possibility is useful: type the first characters in the command that you want to re-use, and hit the arrow-up key (\uparrow) . This way, the most recent command of the history list starting with exactly those characters, will be found.

2.12.3 OS commands

Sometimes, it is useful to issue operating systems (OS) commands from within Matlab. This used to be achieved by preceding the command with the character !. In the latest version of Matlab, preceding a command by ! and ending the command by &, opens a console window where OS operation can be performed on a command line. However, in the latest version of Matlab, there are built in Matlab commands for issuing the most common OS commands, such as cd (Change Directory), delete (DELETE file), dir (list content of DIRectory), mkdir (ReMove DIRectory), etc.

>> dir

	eulersim.m	reaksim1.m
	losandregrad.m	reaksim2.m
MatlabImportData.xls	matlab.mat	reaksim3.m
dampfjaer.m	myworkspaceaA	slett

```
data.m
                      myworkspaceaB
                                             slett.mat
dfsim1.m
                      myworkspacem.mat
                                             temp
dfsim2.m
                      myworkspacemA.mat
                                             temp.mat
etwadata.m
                      myworkspacemB.mat
etwaeqsol.m
                      odexeu.m
                      reakdata.m
etweq.m
>> cd ..
>> dir
                   Intro-kurs.ind
                                       Matlab-course.dmp
                   Intro-kurs.lof
                                       Matlab-course.dvi
. .
FHQOYRO0.bmp
                   Intro-kurs.log
                                       Matlab-course.idx
Figurer
                   Intro-kurs.lot
                                       Matlab-course.ilg
HJ3FG000.wmf
                   Intro-kurs.tex
                                       Matlab-course.ind
HJ3FG001.wmf
                   Intro-kurs.toc
                                       Matlab-course.lof
Intro-kurs.aaa
                   IntroFig
                                       Matlab-course.log
Intro-kurs.aab
                   Matlab
                                       Matlab-course.lot
Intro-kurs.aux
                   Matlab-course.aaa
                                       Matlab-course.tex
Intro-kurs.bak
                   Matlab-course.aab
                                       Matlab-course.toc
Intro-kurs.bbl
                   Matlab-course.aut
                                       figs
Intro-kurs.blg
                                       sw0000
                   Matlab-course.aux
Intro-kurs.dvi
                   Matlab-course.bak
Intro-kurs.idx
                   Matlab-course.bbl
Intro-kurs.ilg
                   Matlab-course.blg
>> cd Matlab
>> dir
                      eulersim.m
                                             reaksim1.m
•
                      losandregrad.m
                                             reaksim2.m
. .
MatlabImportData.xls
                      matlab.mat
                                             reaksim3.m
dampfjaer.m
                      myworkspaceaA
                                             slett
data.m
                      myworkspaceaB
                                             slett.mat
dfsim1.m
                      myworkspacem.mat
                                             temp
dfsim2.m
                      myworkspacemA.mat
                                             temp.mat
etwadata.m
                      myworkspacemB.mat
etwaeqsol.m
                      odexeu.m
                      reakdata.m
etweq.m
```

2.13 Basic functions

2.13.1 Overview of operations

We have already used the built-in Matlab commands eye, zeros, ones, linspace, colon, etc. There are a number of mathematical functions in Matlab, e.g.sin, cosh, tan, exp, etc., as well as other functions.

Typically, operations/functions are either unary or binary. Unary functions operate on a single object, while binary functions operate on two objects. The very fundamental binary operations for arrays are array addition, array subtraction, array multiplication, and array division: in Table 2.6, it is assumed that both A and B are arrays of the same size.

Let us next take a look at the mathematical (unary) functions. If we in Matlab let an array be the argument of a mathematical function, how is the result to be interpreted? A simple example illustrates this the best. If we consider applying the sinus function sin on the row array (x_1, x_2, \ldots, x_n) , this is interpreted as the following row array:

$$\sin(x_1, x_2, \dots, x_n) = (\sin x_1, \sin x_2, \dots, \sin x_n):$$

Matlab operatorMatlab functionDescriptionA+Bplus(A,B)Element by element addition, $A_{ij} + B_{ij}$.A-Bminus(A,B)Element by element subtraction, $A_{ij} - B_{ij}$.A.*Btimes(A,B)Element by element multiplication, $A_{ij} \cdot B_{ij}$.A./Brdivide(A,B)Element by element division, A_{ij}/B_{ij} .

Table 2.6: Basic binary array operations in Matlab. It is assumed that array A and B have the same shape.

>> x=0.3

x =

0.3000

>> sin(x)

ans =

0.2955

```
>> x=[0.3,0.7,0.18]
```

```
х =
```

0.3000 0.7000 0.1800

>> sin(x)

ans =

```
0.2955 0.6442 0.1790
>> x=[0.3;0.7;0.18];
>> sin(x)
```

ans =

0.2955 0.6442 0.1790

In general, let fun (x) denote a general mathematical function of a scalar x. If we use an $m \times n$ array A as argument to the Matlab function, we (normally) get:

$$\operatorname{fun}(A) = \begin{pmatrix} \operatorname{fun}(A_{1,1}) & \operatorname{fun}(A_{1,2}) & \cdots & \operatorname{fun}(A_{1,n}) \\ \operatorname{fun}(A_{2,1}) & \operatorname{fun}(A_{2,2}) & \cdots & \operatorname{fun}(A_{2,n}) \\ \vdots & \vdots & \ddots & \vdots \\ \operatorname{fun}(A_{m,1}) & \operatorname{fun}(A_{m,2}) & \cdots & \operatorname{fun}(A_{m,n}) \end{pmatrix}.$$

In other words: the answer is found by copying the input argument array, and then mapping the function to every element of the copied input argument. This is valid for most mathematical functions, e.g. when fun is sin, cos, tan, sinh, cosh, sec, exp, etc. As an example:

1 2 3 2 3 5 >> sinh(A)

ans =

1.1752	3.6269	10.0179
3.6269	10.0179	74.2032

where we have used the hyperbolic sine function sinh.

One more example: suppose we want to compute $\sin x$ for a number of values x in the interval $[0, 2\pi]$. This can be achieved as follows:

```
>> x = linspace(0,2*pi,10)
х =
  Columns 1 through 6
         0
              0.6981
                         1.3963
                                   2.0944
                                              2.7925
                                                        3.4907
  Columns 7 through 10
    4.1888
              4.8869
                         5.5851
                                   6.2832
>> y = sin(x)
у =
  Columns 1 through 6
         0
              0.6428
                         0.9848
                                   0.8660
                                              0.3420
                                                        -0.3420
  Columns 7 through 10
   -0.8660
             -0.9848
                        -0.6428
                                  -0.0000
```

In the sequel, some basic Matlab functions are discussed. More functions can be found e.g. in (?) or in the Matlab documentation.

2.13.2 Trigonometric functions

A partial list of available trigonometric functions is shown in Table 2.7. Examples of the use of these trigonometric functions are shown below:

Matlab function	Math function	Description
cos(x)	$\cos x$	x is given in radians.
cosh(x)	$\cosh x = \left(e^x + e^{-x}\right)/2$	
sin(x)	$\sin x$	"
sinh(x)	$\sinh x = \left(e^x - e^{-x}\right)/2$	"
tan(x)	$\tan x = \sin x / \cos x$	"
tanh(x)	$\tanh x = \sinh x / \cosh x$	"
acos(x)	$\cos^{-1}x$	Result is given in radians.
acosh(x)	$\cosh^{-1}x$	
asin(x)	$\sin^{-1}x$	"
asinh(x)	$\sinh^{-1}x$	"
atan(x)	$\tan^{-1}x$	"
atanh(x)	$\tanh^{-1} x$	"

Table 2.7: Selected trigonometric functions in Matlab.

```
ans =
```

0.4671 >> sin(A)ans = 0.8135 0.5703 0.7779 0.2291 0.4671 0.6904 >> asin(ans) ans = 0.6068 0.8913 0.9501 0.2311 0.4860 0.7621

Notice that $\sin (A_{2,2}) = (\sin A)_{2,2}$. Also notice that when we first take the sinus of a number, and next take the inverse sinus (arcsine), we will not necessarily get back the original result — the answer may be in a different quadrant:

2.13.3 Exponential functions

A partial list of available exponential functions is shown in Table 2.8. Some examples of the use of these functions:

Matlab function	Math function	Description
x.^p	x^p	Array exponentiation.
exp(x)	$\exp x = e^x$	Exponentiation of e .
log(x)	$\ln x = \log_e x = \exp^{-1} x$	Natural logarithm.
log10(x)	$\log_{10} x = \ln x / \ln 10$	Briggs'/logarithm with base 10.
log2(x)	$\log_2 x = \ln x / \ln 2$	Logarithm with base 2.
pow2(x)	$\log_2^{-1} x$	Inverse of $\log_2 x$.
sqrt(x)	$\sqrt{x} = x^{1/2}$	Square root of x .

Table 2.8: Selected exponential functions in Matlab

Table 2.9:	Selected	complex	functions	in	Matlab.
------------	----------	---------	-----------	----	---------

Matlab function	Math function	Description
abs(z)		Module/absolute value/magnitude of z .
angle(z)	$\angle z$	Phase angle in radians of z .
conj(z)	\overline{z}	Complex conjugate of z .
<pre>imag(z)</pre>	$\Im z$	Imaginary part of z .
real(z)	$\Re z$	Real part of z .
isreal(z)	—	True for real variables, false otherwise.
cplxpair(z)	—	Sort array into complex conjugate pairs.
<pre>complex(x,y)</pre>	$x + i \cdot y$	Form complex number $z = x + i \cdot y$.
z', ctranspose(z)	z^*	Complex conjugate and transpose of array z .

>> A(2,2)^2

ans =

0.2362

>> A.^2

ans =

0.9027	0.3683	0.7944
0.0534	0.2362	0.5808

Notice that for scalars s, we may write s^2 . This doesn't work for arrays though, or rather for array A, writing A^2 has a different meaning than A^2 . Operation A^2 makes sense when the array is a matrix with certain properties, but not for general arrays. Operation A^2 on the other hand, makes sense for any array, and produces the result discussed in the introduction.

2.13.4 Complex functions

A partial list of available complex functions is shown in Table 2.9. Some example of the use of these functions:

```
>> B = A-eye(size(A))
```

B =

-0.0499	0.6068	0.8913
0.2311	-0.5140	0.7621

>> C = sqrt(B)

C = 0 + 0.2233i 0.7790 0.9441 0.4808 0 + 0.7170i 0.8730 >> abs(C)ans = 0.2233 0.7790 0.9441 0.4808 0.7170 0.8730 >> angle(C) ans = 1.5708 0 0 0 1.5708 0 >> conj(C) ans = 0 - 0.2233i 0.7790 0.9441 0.4808 0 - 0.7170i 0.8730 >> imag(C) ans = 0 0 0 0.2233 0.7170 0 >> real(C) ans = 0.7790 0.9441 0 0 0.8730 0.4808 >> isreal(C) ans = 0 >> C' ans = 0 - 0.2233i 0.4808 0.7790 0 - 0.7170i 0.8730 0.9441

2.13.5 Rounding and remainder functions

A list of rounding and remainder functions is shown in Table 2.10.

Table 2.10:	Selected	rounding	and	$\operatorname{remainder}$	functions	in Matlab.	

Matlab function	Math function	Description
fix(x)	$\lfloor x \rfloor \cdot \operatorname{sign} x$	Round toward zero.
floor(x)	$\lfloor x \rfloor$	Round toward $-\infty$.
ceil(x)	$\lceil x \rceil$	Round toward $+\infty$.
round(x)	[x]	Round toward nearest integer.
mod(x,y)	$mod(x, y) \cdot sign y$	Modulus or signed remainder.
rem(x,y)	$mod(x, y) \cdot sign x$	Remainder after division.
sign(x)	$\operatorname{sign} x$	Signum function.

Some examples of how to use these functions:

>> D = 10*B D = -0.4987 6.0684 8.9130 -5.1402 2.3114 7.6210 >> fix(D) ans = 0 6 8 2 -5 7 >> floor(D) ans = 6 -6 -1 8 -6 2 7 >> ceil(D) ans = 0 7 9 3 -5 8 >> round(D) ans = 0 6 9 2 -5 8 >> sign(D) ans = -1 1 1 1 -1 1

Here, function **round** is probably the one that is most useful. Modulus and remainders are computed as follows:

Matlab function	Description
factor(N)	Prime factors of integer N .
isprime	True for prime numbers, false otherwise.
primes(N)	Generate list of prime numbers $\leq N$.
gcd(M,N)	Greatest common divisor of integers M and N .
lcm(M,N)	Least common multiple of integers M and N .
rats(x)	Rational approximation of floating point number x .

Table 2.11: Selected number theoretic functions in Matlab.

>> mod(D,2)

ans =

1.5013	0.0684	0.9130
0.3114	0.8598	1.6210

>> rem(D,2)

ans =

-0.4987	0.0684	0.9130
0.3114	-1.1402	1.6210

2.13.6 Number theoretic functions*

A partial list of number theoretic functions is shown in Table 2.11. Prime factors are found as follows:

```
>> N = 133
```

N =

133

>> factor(N)

ans =

7 19

Notice that the answer is given as elements in a row array. Is e.g. 19 a prime number?

```
>> isprime(ans(2))
ans =
    1
>> isprime(N)
ans =
    0
```

What are the prime numbers smaller than 133?

```
>> primes(N)
```

ans =

```
Columns 1 through 11
```

З Columns 12 through 22 Columns 23 through 32

What is a good rational approximation of π ?

>> rats(pi)

ans =

355/113

Suppose we want to simplify 144/244? We compute the greatest common divisor:

>> gcd(144,244)

ans =

Thus, $\frac{144}{4}/\frac{244}{4} = 36/61$ is the simplest equivalent expression where both numerator and denominator are integers. Because 4 is the GCD of (144, 244), we know that 144/4 and 244/4 are integer numbers. Suppose we want to add 1/12 + 1/16. What is the least common multiple of 12 and 16?

>> lcm(12,16)

ans =

Thus, we have $1/12 + 1/16 = \frac{48}{12}/48 + \frac{48}{16}/48 = 4/48 + 3/48 = (4+3)/48 = 7/48$. Because 48 is the LCM of (12, 16), we know that 48/12 and 48/16 are integer numbers.

2.14 Help

It is a common problem that Matlab has so many functions, that it is impossible to keep track of how each of them work. And sometimes, we don't even know whether a certain function is available.

If we know the name of a certain function, and want to find out how to use it, we can either type >>help <functionname>, >>helpwin <functionname>, or >>doc <functionname> to get a description of how to use the function:

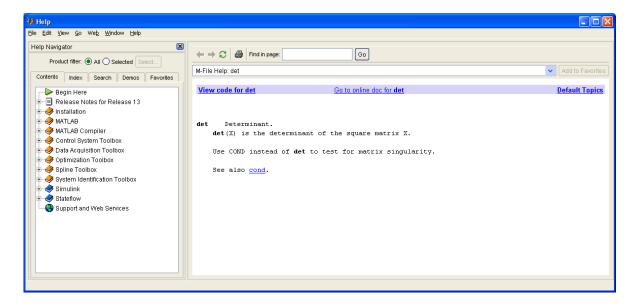


Figure 2.9: The help browser window which opens when using the doc command in Matlab's command window, or by choosing Help/MATLAB Help in the Matlab window.

Using doc instead of help (e.g. doc det) opens up a help browser in a separate window¹⁰, fig. 2.9. The help browser can also be used to search for keywords, etc., when one does not know the exact name of a function. The help browser can also be opened from the HELP menu of the Matlab window, by selecting MATLAB Help. From this help window, it is also possible to search the world wide web for information about Matlab functionality.

 $^{^{10}}$ As an alternative to command doc, command helpwin opens a help window. When the Matlab documentation files are installed, command helpwin appears to give the same response as command doc.

Chapter 3

Basic plotting in Matlab

3.1 Overview of learning goals

In order to study data from experiments, or results from computations, it is often useful to present the results graphically in plots. Matlab has very good support for plotting. Here, we will mainly go through the possibilities in two-dimensional plotting. Good references for plotting in Matlab are (Higham & Higham 2000) for a brief overview, and (Hanselman & Littlefield 2005) for more details.

After having completed this chapter, you should have a clear understanding of what a figure is, and how you can adjust properties of the figure interactively in Matlab. In particular, you should master:

- How to produce two-dimensional plots and how to modify plots,
- How to create arrays of plots,
- Presentation of experimental data, 3.5.

The description will indicate how you use the Matlab commands of Table 3.1.

3.2 Basic two dimensional plots

3.2.1 The plot function

The basic plot command is plot:

```
>> x=linspace(0,2*pi,20);
>> y = sin(x);
>> plot(y)
```

plot is a built-in Matlab function. The result of this command is shown in fig. 3.1, where the plot function has been used with a single argument. Then, the elements of the (row or column array) argument is displayed as a function of the element number: because array x has 20 elements, array y has 20 elements, too. When using the plot command, Matlab draws straight lines between the data points. Thus, the result is a graph with 19 straight line segments between 20 data points: $(1,0), (2, \pi/10), \ldots, (20, 2\pi)$, and the *abscissa* ("x-axis") ranges from 1 to 20.

Table	5.1. Dasic plotting commands.	
Matlab command	Description	Reference
plot	Plot 2D plots	p. 43
	Interactive modification of plots	p. 44
hold on / hold off	Hold/un-hold plots	p. 46
legend	Insert legend	p. 47
	IAT _E X typesetting in plots	p. 48
subplot	Create array plots	p. 51

Table 3.1: Basic plotting commands.

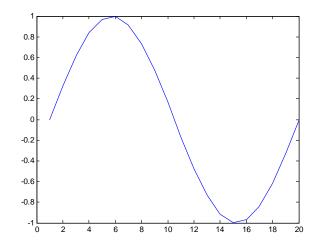


Figure 3.1: Function $y = \sin x$ plotted using command plot(y).

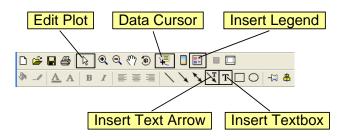


Figure 3.2: Figure Toolbar (upper row) and Plot Edit Toolbar (lower row). Select which toolbar to see from the View menu of the figure window.

3.2.2 Editing plots

In order to modify a plot, the Figure Toolbar and the Plot Edit Toolbar are useful, fig. 3.2. The Figure Toolbar is displayed by default, while the Plot Edit Toolbar is toggled on in the **View** menu of the figure window.

In order to edit a plotted graph, click on the Edit Plot icon (fig. 3.2), and then click on the curve to select the curve. Table 3.2 illustrates how we can modify some properties of the graph interactively.

Usually, we do not want to plot y as a function of the element number, but rather to plot y as a function of x:

>> plot(x,y)

The result is shown in fig. 3.3. Notice that the abscissa ranges from 0 to $6.28 \approx 2\pi$.

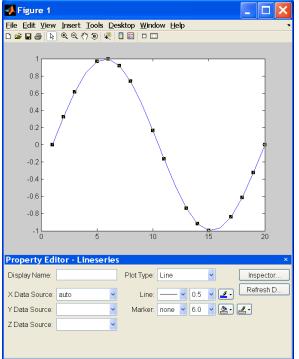
3.2.3 Multiple plots

We can plot several functions in the same plot window:

>> plot(x,y,x,cos(x))

The result is displayed in fig. 3.4.

Remark 3.1 In order to prepare a plot for printing, it is *advisable* to use black color on each graph, and to distinguish among several graphs by varying the line type (solid, dotted, dashed, etc.).



(a) Click Edit Tool (fig. 3.2), then click graph, right-click, and choose Properties...

📦 graph2d.lineseries		
- BeingDeleted	off	
- BusyAction	🔽 queue	
- ButtonDownFcn		
— Clipping	🔽 on	
🖅 Color	O 100 100 100 100 100 100 100 100 100 10	
— CreateFcn		
— DeleteFon		
— DisplayName	null	
— EraseMode	🔽 normal	
— HandleVisibility	💌 on	
— HitTest	🔽 on	
- Interruptible	🔽 on	
- LineStyle	.	
- LineWidth	0.5	
— Marker	🔽 none	
MarkerEdgeColor	🕒 auto	
MarkerFaceColor	🔕 none	
— MarkerSize	6.0	
 SelectionHighlight 	💌 on	
— Tag		
— UIContextMenu		
— UserData	🌐 (0x0 double array)	
— Visible	💌 on	
— XData	[1.0; 2.0; 3.0; 4.0; 5.0;	6.0; 7.0; 8.0; 9.0; 10.0
— XDataMode	💌 auto	
— XDataSource	null	
— YData	0.0; 0.324699469204	68346; 0.614212712
— YDataSource	null	
— ZData		
- ZDataSource	null	

Table 3.2: Mofidying properties of graphs.

(b) Click on Inspector... in Proper Editor (table element (a)).

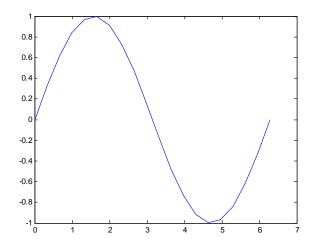


Figure 3.3: Function $y = \sin x$, plotted using command plot(x,y).

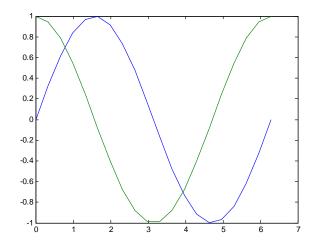


Figure 3.4: Function $y = \sin x$ and function $\cos x$ plotted using command plot(x,y,x,cos(x)).

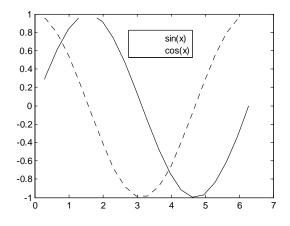


Figure 3.5: Example of edited line styles and line colors. To insert legend, click on the Insert Legend icon (fig. 3.2). To change the default legend text, double-click on the text and edit it.

By following the recipe in Table 3.2, the line styles and line colors of fig. 3.4 have been changed, and *legends* have been inserted, see fig. 3.5.

It is also possible to *add plots* to a figure window. By simply typing a new plot command:

>> plot(x,exp(-x))

the previous plot is deleted, and the new one is drawn. That is not what we want. To make sure that we preserve the original graphs in the figure window, we need to issue a command to **hold** the current graphs in the figure window:

```
>> plot(x,y,x,cos(x))
>> hold on
>> plot(x,exp(-x))
>> hold off
```

The result is shown in fig. 3.6.

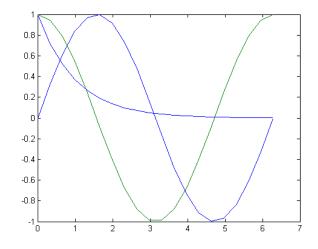


Figure 3.6: The graph of $\exp(-x)$ has been added to the graphs of $\sin x$ and $\cos x$.

Another way to plot several graphs in one figure, is to let the second argument be a two dimensional array. Either the number of rows, or the number of columns of the second argument, must be equal to the length of the first argument:

>> plot(x, [y; cos(x); exp(-x)]) >> plot(x', [y', cos(x)', exp(-x)'])

The command lines above give the same result; we could also have used the following command:

>> z = [y; cos(x); exp(-x)]; >> plot(x,z)

3.2.4 Editing axes properties

So far, we have seen how to change the properties of individual graphs. It is also possible to change the properties that are shared by all graphs. Table 3.3 shows how to select the so-called Axes Properties: Some of the Axes properties can only be used with three dimensional plots. Labels, etc. can be typeset using T_FX and I_TFX commands.

3.2.5 Command line plot editing

All the settings for Graph and Axes properties, can also be set on the command line. This is very useful if we need to create a script for producing many figures overnight.

In the simplest case, we can produce the plot in fig. 3.5 as follows:

```
>> x = linspace(0,2*pi,20);
>> plot(x,sin(x),'k-',x,cos(x),'k--')
>> legend('sin(x)','cos(x)')
```

The only necessary interactive modification, is to use the cursor to move the legend box. In the plot statement above, the first character of the text strings indicate black color. For other colors and line styles, as well as markers, see doc plot, and click on the LineSpec hyperlink.

For more advanced modifications, it is possible to take advantage of the Handle Graphics data structure, see Chapter 12. A number of Matlab books contain more information, e.g. (Hanselman & Littlefield 2005).

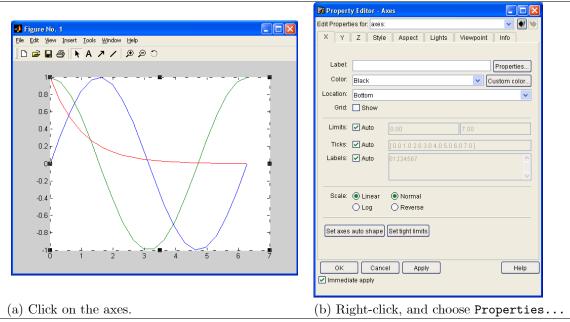


Table 3.3: Modifying properties of axes.

3.2.6 Fancy typesetting in plots*

The basic typesetting in Matlab plots is rather simple. To achieve a more professional typesetting, it is possible to take advantage of T_EX and $I\!\!\!A T_EX$ typesetting (which is used in this document). For an introduction to $I\!\!\!A T_EX$, see e.g. (Lamport 1986).

Table 3.4 illustrates some possibilities for $\[Mathbb{E}]X$ typesetting in Mathab plots. The legend text is modified as indicated in the two upper plots of Table 3.4. By selecting the legend box, right-clicking, and selecting Interpreter/latex, the result in the lower left plot of Table 3.4 is achieved.

In LAT_EX typesetting, mathematical expressions are bracketed by the \$ character, or by the \$\$ symbol: using \$ bracketing, indicates in-line math such as $\sin(x)$ and $\int_0^{2\pi} \sin(x) dx$, while \$\$ bracketing indicates displayed math such as

$$\sin(x) \\ \int_0^{2\pi} \sin(x) \, dx.$$

In Table 3.4, the lower left plot doesn't really illustrate the difference between in-line math and displayed math. To demonstrate the difference in a Matlab plot, we replace the $\cos(x)$ and the $\cos(x)$ and the $\cos(x)$ legend strings by $\int_0^{2\pi}.$ interpreter, the result is as shown in the lower right plot of Table 3.4.

Inserting typeset legends can also be done from the command line. For the legend in the lower right plot of Table 3.4, the following statement can be used:

3.2.7 Closing

For more information about LATEX typesetting, see the Help browser, e.g. (Hanselman & Littlefield 2005), or any LATEX book such as (Lamport 1986).

Exercise 3.1 Describe what the following functions can do: hold, axes, xlabel, ylabel, title, gtext, logspace, format. What are the arguments of these functions?

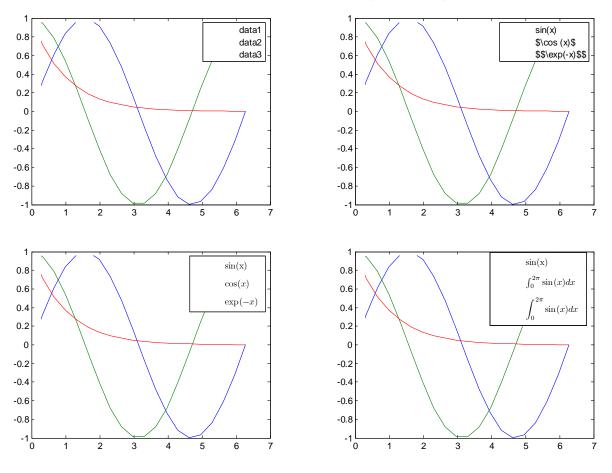


Table 3.4: Illustration of using TEX and $\ensuremath{\mathbb{P}}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspace{TEX}\xspac$

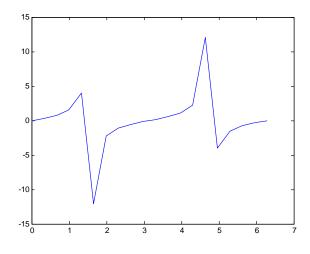


Figure 3.7: Function $y = \tan x$, plotted using command plot(x,tan(x)).

Exercise 3.2 Check out the help files for the following Matlab 2D functions: plot, loglog, semilogx, semilogy, plotyy, polar, fplot, fill, area, bar, barh, hist, pie, comet, errorbar, quiver, scatter.

3.3 Improving plots

In some cases, the plots become "ugly"/erroneous — this is typically the case where the ordinate value goes to $\pm \infty$ at some value of the abscissa, e.g. the function $y = \tan x$:

>> plot(x, tan(x))

The result is shown in fig. 3.7. The result is obviously incorrect: We know that there are asymptotes at $x = \frac{\pi}{2}i, i \in \{1, 3, 5, \ldots\}$, but the plotting algorithm draws straight lines between every point. We could wish to clip the function value e.g. at $y = \pm 3$, but in such a way that if $|\tan x| \ge 3$, nothing is plotted. We can achieve this as follows:

- 1. Find the indices for vector $\tan x$ where $|\tan x| \ge 3$.
- 2. Replace the functional value of $\tan x$ for the given indices $(|\tan x| \ge 3)$ with the value NaN (Not a Number). If the symbol NaN is found in an array that Matlab is asked to plot, no lines are drawn to this point.

The following Matlab sequence illustrates how this can be done:

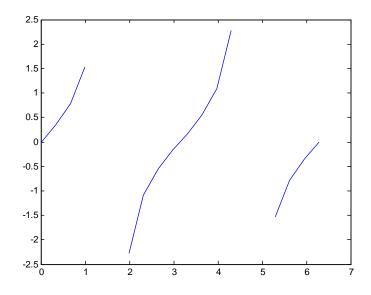


Figure 3.8: The function $y = \tan x$, plotted using command plot(x,y), but where the y values greater than 3 in absolute value have been replaced by NaN (Not a Number).

0	0.3433	0.7783	1.5306	NaN	NaN	-2.2798
Columns 8	through 14	4				
-1.0863	-0.5412	-0.1669	0.1669	0.5412	1.0863	2.2798
Columns 1	5 through 2	20				
NaN	NaN	-1.5306	-0.7783	-0.3433	-0.0000	
>> plot(x,y)					

The result is shown in fig. 3.8. In fig. 3.8 we can clearly see the indication of asymptotes at $x = \frac{\pi}{2}$ and at $x = \frac{3\pi}{2}$.

3.4 Array plots

In addition to plotting several graphs in a single axis system, it is possible to put graphs in an array of axes. The plots are considered as elements in a two dimensional array, and the command subplot is used to define which element the plot is put into. The the command has three integer arguments, subplot(m,n,p), and this command is followed by e.g. the plot command which actually inserts the plot. In the subplot command, the two first arguments (m, n) indicate that the plot is to be inserted into an $m \times n$ array. The third argument (p) specifies in which element of the array the plot will be placed — counting row-wise.

Suppose we want to produce an array with 1 column and 3 rows: the first element is specified as subplot(3,1,1), the second element as subplot(3,1,2), etc. The two first arguments define the size of the array, and the third argument specifies the location where the next plot should be put.

The following example illustrates the procedure:

>> subplot(1,3,1) >> plot(x,y)

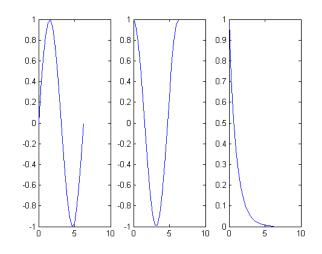


Figure 3.9: The use of the subplot command to produce an array of plots.

>> subplot(1,3,2)
>> plot(x,cos(x))
>> subplot(1,3,3);
>> plot(x,exp(-x))

This produces the plots in fig. 3.9.

It is possible to produce more complicated array plots if we "cheat" and mix the array division in the plot. The following example illustrates this:

>> subplot(2,2,1)
>> plot(x,y)
>> subplot(2,2,2)
>> plot(x,cos(x))
>> subplot(2,1,2)
>> plot(x,exp(-x))

Here, we have mixed freely between one plot array of size (2, 2), and another of size (2, 1). The result is displayed in fig. 3.10.

3.5 Presentation of experimental data

3.5.1 Case study: distillation of water and ethanol

For liquid mixtures of two components (e.g. water and ethanol), the unit operation of distillation can be used to separate the two components into their pure constituents. The distillation process takes advantage of the fact that the two components have different boiling points, i.e. by heating up the mixture, the component with the lowest boiling point will evaporate first. Thus, in principle, the vapor phase will contain the component with the lowest boiling point. In reality, part of the other component will also be found in the vapor phase.

In industrial practice, the mixture is continuously fed to a mixing vessel (feed stream F), where the mixture is heated to the boiling point, fig. 3.11. The resulting vapor leaves at the top of the mixing vessel (vapor stream V), while the (boiling) liquid leaves at the bottom of the vessel (liquid stream L). Usually, one assumes that the vapor and liquid are at thermodynamic equilibrium. It is customary to measure the content of "light" component (i.e., the one with the lowest boiling point, e.g. ethanol in a water-ethanol mixture) as the mole fraction of light component, i.e. the number of moles of light component divided by the total number of moles in the mixture. \dot{Q} indicates the added heat.

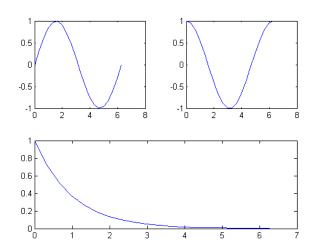


Figure 3.10: The use of the subplot command to produce an array of plots, where there is a mixture of array sizes.

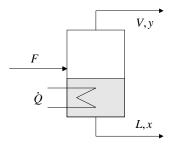


Figure 3.11: Sketch of a simple distillation unit, with feed stream (F), bottom liquid stream (L) with mole fraction x of light component, and top vapor stream (V) with mole fraction y of light component.

X 1	Aicro	soft Ex	ccel - Dist	illatio	m													
8	File	Edit	<u>V</u> iew In:	sert	F <u>o</u> rmat <u>T</u> o	ols <u>D</u> ata	<u>W</u> indow <u>H</u> elp	Acro <u>b</u> at								Туре а	question for h	elp 🔹 🗕 🗗 🗙
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1			Equil	ibri	um da	ata for	water/e	ethano	l: (x,y)	are mo	ole frac	tions c	of etha	nol in t	he liqu	id (x) a	and va	por (y) 🗌
2	х		0.00	000	0.0714	0.1429	0.2143	0.2857	0.3571	0.4286	0.5000	0.5714	0.6429	0.7143	0.7857	0.8571	0.9286	1.0000
3	у		0.00	000	0.3777	0.4913	0.5462	0.5814	0.6092	0.6353	0.6623	0.6921	0.7259	0.7648	0.8100	0.8631	0.9256	1.0000
- A 4 - 4	• • •	N \She	eet1/Sh	eet2 /	(Sheet3 /													
Read	ły																	

Figure 3.12: Equilibrium data for water and ethanol, as found from experiments: (x, y) are mole fractions of liquid and vapor at equilibrium.

Through experiments with continuous distillation of a mixture of water and ethanol, assume that we have found the following correlation between the ethanol content of the liquid (x) and the ethanol content of the vapor phase (y), fig. 3.12.

3.5.2 Plotting data

The equilibrium data are stored in an Excel file. First we import the data into Matlab, using File/Import Data... in Matlab. The import filter interprets the first row of the spreadsheet (see fig. 3.12) as NaN, and produces a variable data consisting of 3 rows and 15 columns:

>> data

data =

Columns 1	through 8						
NaN	NaN	NaN	NaN	NaN	NaN	NaN	NaN
0.0000	0.0714	0.1429	0.2143	0.2857	0.3571	0.4286	0.5000
0.0000	0.3777	0.4913	0.5462	0.5814	0.6092	0.6353	0.6623
Columns 9	through 15						
NaN	NaN	NaN	NaN	NaN	NaN	NaN	
0.5714	0.6429	0.7143	0.7857	0.8571	0.9286	1.0000	
0.6921	0.7259	0.7648	0.8100	0.8631	0.9256	1.0000	

Clearly, the second row contains the x-data, while the third row contains the y-data:

```
>> x = data(2,:);
>> y = data(3,:);
>> plot(x,y,'k-', x,x,'k:')
>> xlabel('x')
>> ylabel('y')
>> title('Equilibrium data for water-ethanol mixture')
```

The data are displayed in fig. 3.13. Clearly, the vapor phase is richer in ethanol than the liquid phase as long as y(x) > x.

There is a problem with the presentation in fig. 3.13: Matlab has drawn straight lines between each of the 15 data points. This is not considered a proper presentation of experimental data, since it may give the *false impression* that we actually know something about the value of y for other values of x than the experimental values. It is thus common to present experimental data as isolated data points:

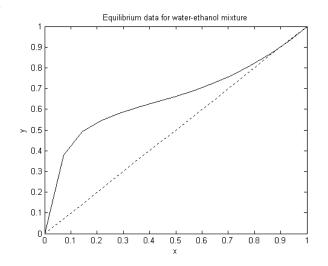


Figure 3.13: The equilibrium correlation for content of ethanol in the liquid phase (x) and in the vapor phase (y).

Matlab command	Description	Reference
meshgrid	Create array argument for 3D plot	p. 167
mesh / meshc	Basic 3D plotting of surfaces	p. 167
contour	Contour plot	p. 169
clf	Clear figure	p. 171
figure	open figure	p. 171
saveas	save plot	p. 171
open	open saved plot	p. 171
close	close figure window	p. 171

Table 3.5: Basic plotting commands.

```
>> plot(x,y,'kx', x,x,'k:')
>> xlabel('x')
>> ylabel('y')
>> title('Experimental equilibrium data for water-ethanol mixture')
>> legend('Experimental data: y(x)', 'y = x')
```

The result is shown in fig. 3.14.

In fig. 3.14, we do not make any assumptions about the values of y(x) outside of the experimental data.

3.6 Further study

Appendix D contains a discussion of more "advanced" topics such as

- How to produce three-dimensional plots, see accompagnying Appendix D,
- Basic housekeeping of plots, see accompagnying Appendix D,
- etc.

Some commands that are treated in Appendix D, are listed in Table 3.5.

Other useful plots not treated directly, are semilogx, semilogy, loglog, plot3, contourf, contour3, meshz, surf, surfc, waterfall, bar3, bar3h, pie, fill3, comet3, scatter3, stem3 — see the Matlab help system. Yet other useful commands are axis, title, xlabel, ylabel — see help system as well as Section 3.5.

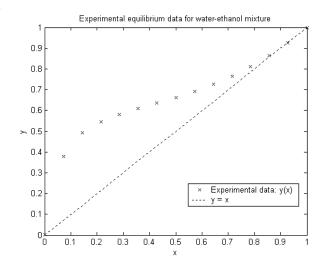


Figure 3.14: Experimentally found equilibrium correlation for content of ethanol in the liquid phase (x) and in the vapor phase (y).

Chapter 4

Simple data analysis*

4.1 Overview of learning goals

After having completed this chapter, you should have a clear understanding of what is meant by interpolation vs. data fitting, what are some advantages and disadvantages of standard interpolation models and fitted models, and basic use of these models. In particular, you should master:

- How to use functions for polynomial and other types of interpolation,
- How polynomials are used in Matlab, basic functions for operating on polynomials, and how these functions can be used to analyze models,
- How to build monovariable and multivariable interpolation models in Matlab.

In addition to using plotting functions that we studied in Chapter 3, see also Table 3.1 p. 43, we will use some additional functions to analyze the data, see Table 4.1.

In addition, Matlab has some very rudimentary functions for data analysis such as max, min, sum, mean, std, var. See the Matlab help system for the rudimentary data analysis tools.

4.2 Interpolation

4.2.1 Overview

In interpolation, we create a mathematical model that goes through every experimental data point. Interpolation is used to compute values of y(x) for x-values for which we do not have experimental values, i.e. in between the experimental values. The structure of such models is typically:

$$y_m\left(x\right) = \sum_{i=1}^n c_i \phi_i\left(x\right)$$

where $\phi_i(x)$ is a chosen function, and c_i is a constant, but the models can also be more complex. Typically in interpolation, an implicit assumption is that the experimental results are perfect.

Matlab command	Description	Reference
polyfit	Single-input polynomial model fit	p. 58
polyval	Evaluate polynomial model	p. 58
spline	Build spline single-input interpolation	p. 60
ppval	Evaluate interpolation models	p. 60
pchip	Build Hermitian cubic polynomials single-input interpolation	p. 60
interp1	Build various types of interpolation models	p. 61
	polynomial functions	Table 4.3 p. 64
interp2	Build various two-input interpolation models	Section 4.4.2

 Table 4.1: Table Caption

Various interpolation models have different sets of functions $\phi_i(x)$, which may be termed basis functions or trial functions. Depending on the chosen basis functions, the resulting interpolation models will have different properties (e.g. smoothness, etc.), and will predict different values $y_m(x)$ in between experimental data points.

The simple interpolation strategy of drawing straight lines between data points, is achieved by selecting $\phi_i(x)$ to be so-called first order spline basis functions. Another possibility is to use polynomial models, e.g.

$$y_m(x) = \sum_{i=1}^n c_i x^{i-1},$$

i.e. we have chosen $\phi_i(x) = x^{i-1}$. Sometimes, we want the interpolation formulae to be differentiable at every x; obviously, we can not achieve this if we use first order spline basis functions.

In the study, we use the experimental data from Section 3.5.

4.2.2 Polynomial interpolation

Let us experiment with polynomial interpolation. In order to find the interpolation formulae, we use the Matlab function polyfit(x,y,n), where x is the input variable, y is the output variable, and n is the order of the polynomial. Note that with N = 15 data points, it is necessary to use a model with 15 parameters c_i , i.e. to build a 14-th order polynomial in order to have true interpolation, i.e. that the model is $y_m(x) = \sum_{i=1}^{15} c_i x^{i-1}$. If we choose n < 15, then the model will not go through every data point. If we choose n > 15, then it is not possible to find unique values c_1 .

For now, we do not worry about the Matlab Warning.

If we want to plot the result, we do this by using command polyval(c,x):

>> ym = polyval(c,x);

Let us see how the interpolation function fits the data:

>> plot(x,y,'kx', x,x,'k:', x,ym,'k-')

The result is shown in fig. 4.1. At first sight, the result looks good. But wait — the model appears to be composed of straight lines between the experimental data points! That can not be correct for a polynomial model!

The problem is that we only computed model values for experimental values of x. Drawing a (solid) line through the computed vales (x, y_m) , Matlab simply uses a straight line between the computed data points (x, y_m) . To get a better view of the model, we need to compute the model value y_m for abscissa values in between the experimental points!, e.g. as follows:

```
>> xm = linspace(0,1,100);
>> ym = polyval(c,xm);
>> plot(x,y,'kx', x,x,'k:', xm,ym,'k-')
>> legend('Experimental data: y(x)', 'y = x', 'Model of data: ym(x)')
>> xlabel('x')
>> ylabel('y')
>> title('Experimental data and model data for water-ethanol equilibrium')
```

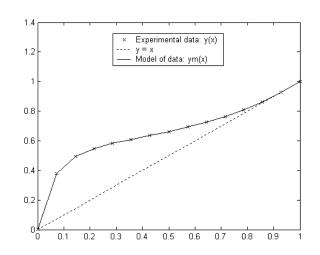


Figure 4.1: Experimental data $((y, x), \times)$ and model $((y_m(x), x), \text{ solid})$ for water-ethanol equilibrium. Notice flawed model representation.

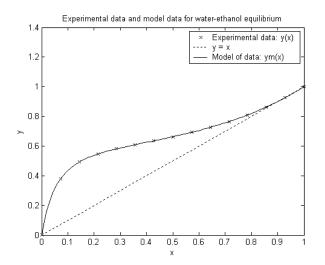


Figure 4.2: Experimental data $((y, x), \times)$ and model $((y_m(x), x), \text{ solid})$ for water-ethanol equilibrium.

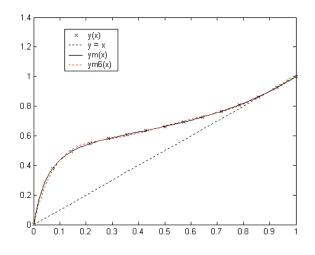


Figure 4.3: Experimental data $((y, x), \times)$, interpolation model $((y_m(x), x), \text{ solid})$, and low order model $(y_{m6}(x), x, \text{ dotted})$ for water-ethanol equilibrium.

The resulting plot is shown in fig. 4.2.

The Matlab warning is caused by the fact that in this case, we could have found almost as good a model by using fewer parameters c_i :

```
>> c6 = polyfit(x,y,6);
>> ym6 = polyval(c6,xm);
>> plot(x,y,'kx', x,x,'k:', xm,ym,'k-', xm,ym6,'r:')
>> legend('y(x)', 'y = x', 'ym(x)', 'ym6(x)')
```

The result is shown in fig. 4.3. Since model ym6 is not a true interpolation model (the model does not go through every experimental data point), we denote it a fitted model — in fact it is the model of 6-th order with the minimal squared error, i.e. the least squares model of 6-th order.

4.2.3 Other interpolation functions

Similar to function polyfit, function spline(x,y) calculates the data structure associated with cubic spline interpolation of the data set (x, y). Next, similar to function polyval, function ppval(c,xx) calculates the model value according to spline interpolation. Note that array xx must be different from array x if we want to compute the value $y_m(x)$ in between experimental data points x.

```
>> c = spline(x,y);
>> ym = ppval(c,xm);
>> plot(x,y,'kx', x,x,'k:', xm,ym,'k-')
>> legend('Experimental data y(x)', 'y = x', 'Spline model ym(x)')
>> xlabel('x')
>> ylabel('y')
>> ym = ppval(c,xm);
>> title('Experimental data and model data for water-ethanol equilibrium')
```

The result is shown in fig. 4.4.

Another interpolation function is pchip(x,y) which works similarly to polyfit and spline, but which is based on Hermitian cubic polynomials:

>> c = pchip(x,y);

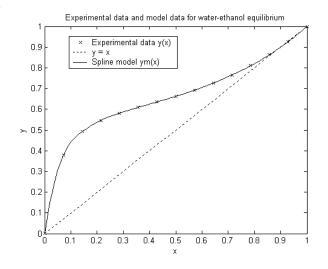


Figure 4.4: Experimental data $((y, x), \times)$ and spline model $((y_m(x), x), \text{ solid})$ for water-ethanol equilibrium.

Table 4.2: Data set for comparison of interpolation models.

```
i
                                    1
                                         2
                                              3
                                                   4
                                                         5
                                    0.8
                                         1.2
                                              2.5
                                                   5.1
                                                         10
                                x_i
                                         2.7
                                              1.4
                                                   0.85
                                                         0.55
                                    5.1
                               y_i
>> ym = ppval(c,xm);
>> plot(x,y,'kx', x,x,'k:', xm,ym,'k-')
>> xlabel('x')
>> ylabel('y')
>> title('Experimental data and model data for water-ethanol equilibrium')
>> legend('Experimental data y(x)', 'y = x', 'pchip model ym(x)')
```

The interpolation model is shown in fig. 4.5.

Finally, there is a generic interpolation function interp1(x,y,xx), which combines the functionality of spline and ppval in one function. An optional fourth argument specifies the type of interpolation: either 'nearest' for nearest neighbor interpolation, 'linear' for the default linear interpolation, 'spline' for cubic spline interpolation, and 'cubic' for cubic interpolation.

For the data set we have studied so far, there is not much difference to see in the different interpolation models, see figs. 4.2 - 4.5. In the next section, we will consider a new data set which better illustrates the properties of the various interpolation techniques.

4.2.4 Comparison of interpolation models

We consider the data set in Table 4.2.

We first create variables in Matlab to hold the data:

```
>> x = [0.8,1.2,2.5,5.1,10];
>> y = [5.1,2.7,1.4,0.85,0.55];
>> plot(x,y,'kx')
```

We want to interpolate the data in 100 data points between $\min x$ and $\max x$:

```
>> x_int = linspace(min(x),max(x),100);
```

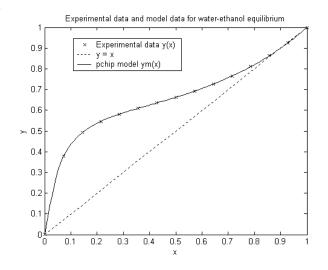


Figure 4.5: Experimental data $((y, x), \times)$ and pchip model $((y_m(x), x), \text{ solid})$ for water-ethanol equilibrium.

Next, we compute the interpolation values from some interpolation models:

```
>> cint_pol = polyfit(x,y,4);
>> ymint_pol = polyval(cint_pol,x_int);
>>
>> cint_spl = spline(x,y);
>> ymint_spl = ppval(cint_spl,x_int);
>>
>> cint_pch = pchip(x,y);
>>
  ymint_pch = ppval(cint_pch,x_int);
>>
>> ymint_i1n = interp1(x,y,x_int,'nearest');
>>
>> ymint_i1l = interp1(x,y,x_int);
>>
>> ymint_i1s = interp1(x,y,x_int,'spline');
>>
>> ymint_i1c = interp1(x,y,x_int,'cubic');
```

We also include a 3-order polynomial least squares model:

```
>> cint_pls = polyfit(x,y,3);
>> ymint_pls = polyval(cint_pls,x_int);
```

Next, we plot the various models:

```
>> plot(x_int,ymint_pol1,'k--', x_int,ymint_spl,'k:', x_int,ymint_pch,'k-', ...
x_int,ymint_i1n,'k-', x_int,ymint_i11,'k--', x_int,ymint_i1s,'k:', ...
x_int,ymint_i1c,'k-.', x_int,ymint_pls,'k:')
>> legend('polynom interp.', 'spline interp.', 'pchip interp.', ...
'nearest neigh. interp.', 'linear interp.', 'spline interp.', 'cubic interp.',...
'polynom least squares')
>> xlabel('x')
>> ylabel('y')
```

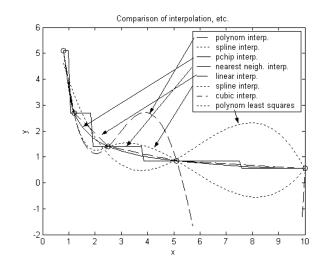


Figure 4.6: Comparison of various interpolation models.

```
>> title('Comparison of interpolation, etc.')
>> hold on;
>> plot(x,y,'ko')
```

The result is shown in fig. 4.6.

From the figure, it is somewhat difficult to compare the various models. We do, however see that the polynomial interpolation model (4-th order) probably is not very good at interpolating the data. Also the spline interpolation model appears to vary quite a lot between the data. The smoothest interpolation model appears to be the Hermitian cubic polynomial model (pchip). Although they are not differentiable, the nearest neighbor interpolation method and the linear interpolation method appear to give result similar to the Hermitian cubic polynomial model.

In sum, the various interpolation methods have different properties, and may give very different interpolation results, in particular if the number of available experimental data are few. We cannot really decide which interpolation method is "best", simply because we do not know the true values between the experimental data points. If we have some notion that the variation is relatively smooth, then we can say more, however. It should also be noted that the smoothness of various models may depend on the available data: in the distillation data, all methods give essentially the same result.

4.3 Analysis

4.3.1 Polynomials in Matlab

Polynomial models have certain advantages over other models when it comes to the ease at which we can analyze the resulting model. Let us see how polynomials are dealt with in Matlab.

Let us start by looking at a polynomial model as found by function polyfit. We consider the data given in fig. 3.12:

```
>> c = polyfit(x,y,4)
c =
    -6.3220 15.5972 -12.8875 4.5495 0.0432
>> length(c)
ans =
```

Table 4.3: Functions on polynomials.							
Function	Description						
conv(p,q)	Multiplication of polynomials, $p(x) \cdot q(x)$						
[q,r] = deconv(b,a)	Division of polynomials, $b(x) = a(x) \cdot q(x) + r(x)$						
polyder(p)	Differentiation of polynomials, $\frac{d}{dx}p(x)$						
<pre>polyint(p)</pre>	Integration of polynomials, $\int p(x) dx$						
polyval(p,x)	Evaluate polynomial, $y = p(x)$						
roots(p)	Roots of polynomials, $x : p(x) = 0$						
cplxpair(x)	Sort complex array \mathbf{x} into complex conjugate pairs						

5

MATLAB represents polynomials as row arrays containing coefficients ordered by descending powers. Thus, row array c represents the polynomial:

$$p(x) = -6.322x^4 + 15.5972x^3 - 12.8875x^2 + 4.5495x^1 + 0.0432.$$

4.3.2 Operations on polynomials

Matlab supports a number of functions for polynomials, see Table 4.3.

Suppose we want to multiply the polynomials $(1 + x - x^2)$ and $(2 + x^3)$. Using a CAS¹, we find:

$$(1+x-x^2) \cdot (2+x^3) = -x^5 + x^4 + x^3 - 2x^2 + 2x + 2.$$

Using Matlab, we find:

>> p = [-1, 1, 1];
>> q = [1,0,0,2];
>> conv(p,q)

ans =

-1 1 1 -2 2 2

We see that the result is identical in Matlab and the CAS.

We can also divide polynomials:

>> deconv(ans,p)

ans =

1 0 0 2

The answer is obviously correct: we have divided $p \cdot q$ by p, and the answer is q. We can also divide q by p to get:

```
> [a,b] = deconv(q,p)
```

a =

-1 -1

 $^{^1\}mathrm{CAS}$ = Computer Algebra System, here: MuPAD through Scientific WorkPlace.

b = 0 0 2 3

The result means that

 $\frac{2+x^3}{1+x-x^2} = -x - 1 + \frac{2x+3}{1+x-x^2}.$

It is trivial to show that this is the correct result. We can differentiate a polynomial, e.g. $q(x) = 2 + x^3$:

>> polyder(q)

ans =

3 0 0

The result is $q'(x) = 3x^2$, which is correct. We can integrate a polynomial, e.g. $p(x) = 1 + x - x^2$:

0

```
>> polyint(p)
ans =
    -0.3333    0.5000    1.0000
```

The result is $\int p(x) dx = -0.3333x^3 + 0.5x^2 + x$, which is correct.

We have previously used function polyval to evaluate polynomials. As an example, p(x = 0.5) can be found as follows:

```
>> polyval(p,0.5)
```

ans =

1.2500

Finally, we can find the roots polynomials, e.g. q(x) = 0:

>> roots(q)

ans =

-1.2599 0.6300 + 1.0911i 0.6300 - 1.0911i

where $\mathbf{i} = \sqrt{-1}$. It can be shown that the roots of polynomials with real number coefficients, must occur in complex conjugate pairs if the root is complex, e.g. if z = 0.6300 + 1.0911i is a root, then $\bar{z} = 0.6300 - 1.0911i$ must also be a root. It is possible to sort the roots in such a way that complex conjugate roots are grouped together using command cplxpair; in the example above, the **ans** array is already sorted.

4.3.3 Analysis of polynomial models

We consider the 6-th order least squares polynomial model for water-ethanol equilibrium:

```
>> c6 = polyfit(x,y,6)
c6 =
    -34.5134 119.0827 -161.7170 110.1621 -39.3068 7.2848 0.0049
```

In the original data, fig. 3.12, we see that the value of y(x) dips below y = x for some value x_1 , and we want to find an estimate of x_1 . We can do that as follows: find the roots of $y_m(x) - x = 0$:

What takes place here? First, we must find the polynomial description of $y_m(x) - x$. $y_m(x)$ is represented by the polynomial of array c6. We must find the array that represents polynomial x; this array must have the same length as array c6, or else we cannot subtract them.

We see that the complex conjugate roots are grouped already. The complex roots do not have a physical interpretation, and we are left with 2 real roots. Most likely, these roots, $x \in \{0.0008, 0.9915\}$ are related to the trivial roots x = 0 and x = 1. Thus, we can not find the location of x_1 from polynomial c6.

What if we choose a polynomial of odd order? In that case, there must be an odd number of real roots! We find:

```
>> c5 = polyfit(x,y,5)
c5 =
    15.5423 -45.1778 49.6511 -25.1126 6.0893 0.0157
>> roots(c5 - [0,0,0,0,1,0])
ans =
    0.9892
    0.8571
    0.5317 + 0.33061
    0.5317 - 0.33061
    -0.0030
```

Now, it is reasonable to associate the root 0.8571 with x_1 . From the data of fig. 3.12, it seems reasonable to bracket the root to lie in the interval $x_1 \in [0.8631, 1]$, and this doesn't really support the root we have found. Let us try with a higher dimensional polynomial of odd order:

```
>> c9 = polyfit(x,y,9);
>> x9 = zeros(size(c9));
>> x9(end-1) = 1
x9 =
     0
           0
                  0
                        0
                              0
                                     0
                                           0
                                                  0
                                                        1
                                                               0
>> roots(c9-x9)
ans =
   0.9047 + 0.1562i
   0.9047 - 0.1562i
   0.9999
   0.9000
   0.5249 + 0.3435i
   0.5249 - 0.3435i
   0.1139 + 0.2661i
   0.1139 - 0.2661i
  -0.0000
```

From the 9-th order model, it appears as if the root $x_1 \approx 0.9000$, which fits better into bracket of the solution. Still, we have too few data to really be sure about the value of x_1 .

We can also differentiate and integrate polynomials:

```
>> c6 = polyfit(x,y,6);
>> c6der = polyder(c6);
>> c6int = polyint(c6);
>> xx = linspace(0,1);
>> plot(xx, polyval(c6,xx),'k-', xx, polyval(c6der,xx),'k:', ...
xx, polyval(c6int,xx),'k-.')
>> legend('y_m(x)', 'd y_m(x)/dx', '\int y_m(x) dx')
>> xlabel('x')
```

The result is shown in fig. 4.7.

Although it should be possible to write similar functions as those of Table 4.3 for spline models, etc., such functions are not currently available in Matlab.

4.4 More on interpolation and prediction models

4.4.1 Case study: Saturated steam

The data in fig. 4.8 are taken from http://webbook.nist.gov.

It is of interest to build a model that correlates pressure and temperature, p(T). We start by importing the data in file SaturatedSteamVapor.xls into Matlab, and defining variables T and p:

```
>> T = data(3:end,1);
>> p = data(3:end,2);
>> plot(T,p,'kx')
>> xlabel('T [C^{\circ}]')
>> ylabel('p [atm]')
```

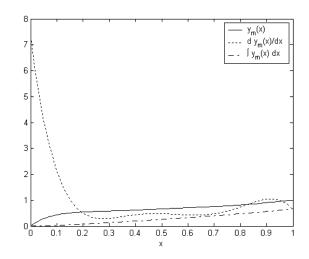


Figure 4.7: 6th order least squares polynomial model $y_m(x)$ of water-ethanol equilibrium, with the derivative $dy_m(x)/dx$ and the integral $\int y_m(x) dx$ of $y_m(x)$.

N 🛛	licrosoft Ex	cel - Satura	atedSteam\	/apor												
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	A	. ,,,	C	D	E	F	G	Н		J	К	L	М	N	0	TH
1 Saturated steam: Vapor Phase data. Taken from http://webbook.nist.gov																
2	Temperat ure (C)		Density (kg/m3)	Volume (m3/kg)	Internal Energy (kJ/kg)	Enthalpy (kJ/kg)	Entrany	Cv (J/g*K)	Ср (J/g*K)	Sound Spd. (m/s)	Joule- Thomson (F/atm)	Viscosity (uPa*s)	Therm. Cond. (W/m*K)	Phase		-
3	0.01	0.0060366	0.004855	205.99	2374.9	2500.9	9.1555	1.4184	1.8844	409	108.09	9.2163	0.017071	vapor	-	-
4	5.01	0.0086177	0.006807	146.91	2381.8	2510.1	9.0246	1.4226	1.8894	412.61	92.224	9.336	0.017339	vapor		
5	10.01	0.01213	0.009413	106.24	2388.7	2519.2	8.8995	1.4269	1.8947	416.17	78.898	9.4614	0.017622	vapor		
6	15.01	0.016846	0.012849	77.828	2395.5	2528.4	8.78	1.4314	1.9002	419.7	67.754]	9.5921	0.017918	vapor		
7	20.01	0.023102	0.017324	57.723	2402.3	2537.5	8.6658	1.4359	1.9059	423.19	58.461	9.7275	0.018228	vapor		
8	25.01	0.031303			2409.1	2546.5	8.5564	1.4405	1.9118	426.63	50.723	9.8671	0.018551	vapor		
9	30.01		0.030432		2415.9	2555.6	8.4518	1.4453	1.918	430.04	44.279	10.011	0.018887	vapor		-
l i i ∢ Read		et1 / Sheet	t2 / Sheet3						•							

Figure 4.8: Relationships between thermodynamic variables of vapor phase saturated steam.

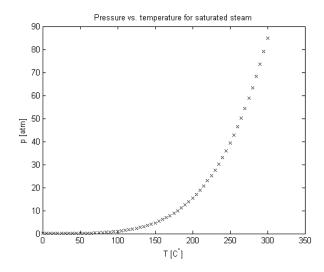


Figure 4.9: Pressure (p, atm) as a function of temparature $(T, ^{\circ}C)$ for saturated steam.

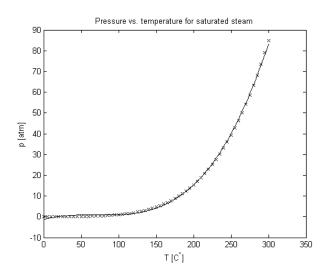


Figure 4.10: Experimental pressure data $(p, \text{ atm}, \times)$ as a function of temparature $(T, ^{\circ}C)$ for saturated steam, and 4th order least squares model fit (solid line).

>> title('Pressure vs. temperature for saturated steam')

The result is shown in fig. 4.9.

We choose to build a polynomial least squares model:

```
>> ppar = polyfit(T,p,3);
>> hold on;
>> TT = linspace(min(T),max(T));
>> plot(TT,polyval(ppar,TT), 'k-')
```

The result is not perfect, see fig. 4.10. In particular, the model is relatively poor at temperatures below ca. 80–90 °C.

Based on the model, what is the pressure at 100 $^\circ\mathrm{C?}$. Using the polynomial least squares model, Matlab suggests:

```
>> polyval(ppar,100)
```

ans =

0.9650

The results is somewhat surprising — we would have expected that $p^{\text{sat}}(100) = 1 \text{ atm}^2$ — the error is probably due to the relatively low quality of the polynomial model at low temperatures.

Let us turn the question around: at a pressure of 1 atm, what is the *temperature*? To answer this question, we must solve the equation $p^{\text{sat}}(T) = 1$. We must thus use the **roots** function for finding roots in Matlab:

```
>> length(ppar)
ans =
     4
>> fpol = ppar - [0,0,0,1]
fpol =
    0.0000
             -0.0015
                         0.1035
                                  -2.4529
>> roots(fpol)
ans =
  1.0e+002 *
   1.0223
   0.5597 + 0.1790i
   0.5597 - 0.1790i
>> format short g
>> ans
ans =
      102.23
      55.968 +
                   17.904i
      55.968 -
                   17.904i
```

First, we created polynomial fpol which is p(T) - 1. Next, we used the roots function to find the roots of p(T) - 1 = 0. The answer is two roots which are complex (a complex conjugate pair 55.968 ± 17.904*i*) and hence irrelevant, and the relevant real root $T^{\text{sat}}(1) = 102.23 \text{ °C}$. Again, we are somewhat puzzled — we know that the answer should have been 100 °C, but again, the inaccuracy is probably due to the poor quality of the model at "low" temperatures.

An interesting question is: could we have avoided the process of finding the roots of p(T) = 1? The answer is yes! We could simply have built another model, where we fit a polynomial p(T) to the data! However, with the current data, least squares polynomial models T(p) are very poor — see fig. 4.11. Let us check the "disastrous" prediction qualities of the 5-th order model:

>> Tpar = polyfit(p,T,5);

²The boiling point temperature for water at pressure p = 1 atm, is T = 100 °C.

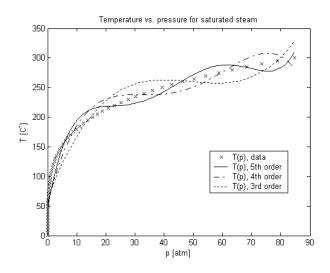


Figure 4.11: Experimental temparature $(T, ^{\circ}C, \times)$ as a function of pressure (p, atm) for saturated steam, and polynomial least squares models of degrees 3–5.

>> polyval(Tpar,1)

ans =

74.067

The correct answer should be $100 \,^{\circ}$ C.

Let us see if the situation improves by using the **spline** function:

99.974

Clearly, true interpolation using the spline function gives very good predictions.

In this section, we have tweaked around with mono variable functions, i.e. models where one variable is a function of another variable. It is often of interest to consider models where one variable is a function of more than one variable — so-called multivariable functions.

4.4.2 Multivariable interpolation

In this section, some of the plotting techniques from Appendix D are used.

To illustrate the ideas, we use superheated steam as a case study. Superheated steam is vapor which is "above" the saturation state of steam. Because of this, the thermodynamic properties are functions of two variables. Figure 4.12 shows the variables at p = 0.5 atm; data at other pressures can easily be created, e.g. for $p \in \{1, 2, 5, 10, 20, 50\}$ atm; these data are taken from http://webbook.nist.gov. Note

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D	🛩 🖪 🧉	3 🛕 🗠 🗸	· 🛍 🔣 1	.00% 🝷 😰) »	Arial	- 1	• B /	<u>u</u> ≣	= = =	§%,	◆.0 .00 • 00. €	F (F 🔛	- <u>ð</u> - <u>A</u>	- ₊
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	A	В	С	D	E	F	G	Н		J	K	L	M	N	-
1	Superh	neated s	steam: '	Taken fi	rom http	o://webb	ook.nis	st.gov							_
2	Pressure	e p = 0.5 a	atm												
	Temperat	Pressure	Density	Volume	Internal	Enthalpy	Entropy		Ср	Sound	Joule-	Viscosity	Therm.		
3	ure (C)	(atm)	(kg/m3)	(m3/kg)	Energy (kJ/kg)	(kJ/kg)	(J/g*K)	Cv (J/g*K)	(J/g*K)	Spd. (m/s)	Thomson (F/atm)	(uPa*s)	Cond. (W/m*K)	Phase	
3 4	ure (C) 100	(atm) 0.5	(kg/m3)	(m3/kg) 3.3736							(F/atm)			Phase vapor	
		0.5	(kg/m3)		(kJ/kg)	(kJ/kg)	(J/g*K)		(J/g*K)	(m/s)	(F/atm) 11.897	(uPa*s)	(W/m*K) 0.024483		
4	100	0.5	(kg/m3) 0.29642	3.3736	(kJ/kg) 2511.4	(kJ/kg) 2682.3	(J/g*K) 7.689	1.4856 1.4734	(J/g*K) 1.9755	(m/s) 474.89	(F/atm) 11.897 6.6802	(uPa*s) 12.311	(W/m*K) 0.024483	vapor	_
4 5 6 7	100 150 200 250	0.5	(kg/m3) 0.29642 0.26051	3.3736 3.8387	(kJ/kg) 2511.4 2585.7 2659.9 2735.1	(kJ/kg) 2682.3 2780.2 2877.7 2976.1	(J/g*K) 7.689 7.9351	1.4856 1.4734	(J/g*K) 1.9755 1.9483 1.9576	(m/s) 474.89 506.03 534.48 561.03	(F/atm) 11.897 6.6802 4.2898	(uPa*s) 12.311 14.208	(W/m*K) 0.024483 0.028526 0.03309	vapor vapor	
4 5	100 150 200 250 300	0.5 0.5 0.5 0.5 0.5	(kg/m3) 0.29642 0.26051 0.23261 0.2102 0.19176	3.3736 3.8387 4.299 4.7575 5.2149	(kJ/kg) 2511.4 2585.7 2659.9	(kJ/kg) 2682.3 2780.2 2877.7	(J/g*K) 7.689 7.9351 8.153 8.3507 8.5325	1.4856 1.4734 1.4886 1.513 1.5413	(J/g*K) 1.9755 1.9483 1.9576	(m/s) 474.89 506.03 534.48	(F/atm) 11.897 6.6802 4.2898	(uPa*s) 12.311 14.208 16.192 18.23 20.296	(W/m*K) 0.024483 0.028526 0.03309 0.038053	vapor vapor vapor	
4 5 6 7	100 150 200 250	0.5 0.5 0.5 0.5 0.5	(kg/m3) 0.29642 0.26051 0.23261 0.2102 0.19176	3.3736 3.8387 4.299 4.7575 5.2149 5.6717	(kJ/kg) 2511.4 2585.7 2659.9 2735.1 2811.5 2889.4	(kJ/kg) 2682.3 2780.2 2877.7 2976.1 3075.7 3176.8	(J/g*K) 7.689 7.9351 8.153 8.3507 8.5325 8.7015	1.4856 1.4734 1.4886 1.513 1.5413 1.5717	(J/g*K) 1.9755 1.9483 1.9576 1.9792 2.006 2.0355	(m/s) 474.89 506.03 534.48 561.03 586.04 609.76	(F/atm) 11.897 6.6802 4.2898 3.0028 2.2249 1.7147	(uPa*s) 12.311 14.208 16.192 18.23 20.296 22.374	(W/m*K) 0.024483 0.028526 0.03309 0.038053 0.043342 0.048908	vapor vapor vapor vapor	
4 5 6 7 8	100 150 200 250 300 350	0.5 0.5 0.5 0.5 0.5	(kg/m3) 0.29642 0.26051 0.23261 0.2102 0.19176 0.17631	3.3736 3.8387 4.299 4.7575 5.2149 5.6717	(kJ/kg) 2511.4 2585.7 2659.9 2735.1 2811.5	(kJ/kg) 2682.3 2780.2 2877.7 2976.1 3075.7	(J/g*K) 7.689 7.9351 8.153 8.3507 8.5325	1.4856 1.4734 1.4886 1.513 1.5413	(J/g*K) 1.9755 1.9483 1.9576 1.9792 2.006	(m/s) 474.89 506.03 534.48 561.03 586.04	(F/atm) 11.897 6.6802 4.2898 3.0028 2.2249 1.7147	(uPa*s) 12.311 14.208 16.192 18.23 20.296	(W/m*K) 0.024483 0.028526 0.03309 0.038053 0.043342 0.048908	vapor vapor vapor vapor vapor vapor	

Figure 4.12: Relationships between thermodynamic variables of superheated steam at p = 0.5 atm.

that for higher pressures, the fluid will liquidize for low temperatures. Since we are interested in vapor behavior, liquid results will be deleted in Excel, and will appear as NaN (not a number) when imported into Matlab.

Assume that we want to find a model for $\rho(T, p)$, i.e. how the vapor density ρ varies with temperature T and pressure p. We import the data, and associate the imported data to descriptive variable names — note that the data starts in the fourth row:

```
Import Wizard created variables in the current workspace.
>> T0_5 = data(4:end,1);
>> p0_5 = data(4:end,2);
>> rho0_5 = data(4:end,3);
Import Wizard created variables in the current workspace.
>> T1 = data(4:end,1);
>> p1 = data(4:end,2);
>> rho1 = data(4:end,3);
Import Wizard created variables in the current workspace.
>> T2 = data(4:end,1);
>> p2 = data(4:end,2);
>> rho2 = data(4:end,3);
Import Wizard created variables in the current workspace.
>> T5 = data(4:end,1);
>> p5 = data(4:end,2);
>> rho5 = data(4:end,3);
Import Wizard created variables in the current workspace.
>> T10 = data(4:end,1);
>> p10 = data(4:end,2);
>> rho10 = data(4:end,3);
Import Wizard created variables in the current workspace.
>> T20 = data(4:end,1);
>> p20 = data(4:end,2);
>> rho20 = data(4:end,3);
Import Wizard created variables in the current workspace.
>> T50 = data(4:end,1);
>> p50 = data(4:end,2);
>> rho50 = data(4:end,3);
```

```
Next, we construct the temperature vector:
```

```
>> T = T0_5';
>> p = [0.5,1,2,5,10,20,50];
>> rho = [rho0_5, rho1, rho2, rho5, rho10, rho20, rho50]
```

rho =

Columns 1 through 6

0.29642	0.59761	NaN	NaN	NaN	NaN
0.26051	0.52326	1.0559	NaN	NaN	NaN
0.23261	0.46645	0.93791	2.3849	4.9225	NaN
0.2102	0.42113	0.84523	2.1362	4.3557	9.099
0.19176	0.38399	0.76986	1.9391	3.9289	8.0794
0.17631	0.35294	0.70713	1.7774	3.5876	7.3144
0.16318	0.32658	0.65401	1.6416	3.3053	6.7032
0.15187	0.3039	0.60841	1.5256	3.0667	6.1973
0.14204	0.28418	0.5688	1.4253	2.8617	5.7687

Column 7

	NaN
	NaN
	NaN
	NaN
22.	405
19.	528
17.	537
16.	013
14.	782

Here, both T and p are row arrays, while **rho** is an array where each row holds the pressure variation for a fixed temperature. Finally, we can interpolate:

```
>> [TT,pp] = meshgrid(T,p);
>> size(TT)
ans =
     7
           9
>> size(rho)
ans =
           7
     9
>> Tvar = linspace(min(T),max(T));
>> pvar = linspace(min(p),max(p));
>> [TTvar, ppvar] = meshgrid(Tvar,pvar);
>> rhovar = interp2(TT,pp,rho',TTvar,ppvar);
>> meshc(TTvar,ppvar,rhovar)
>> xlabel('T [C^\circ]')
>> ylabel('p [atm]')
>> zlabel('\rho [kg/m^3]')
```

The resulting surface $\rho(T, p)$ is displayed in fig. 4.13.

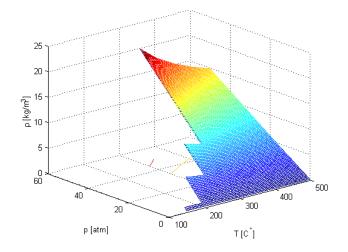


Figure 4.13: Interpolation model for $\rho(T, p)$.

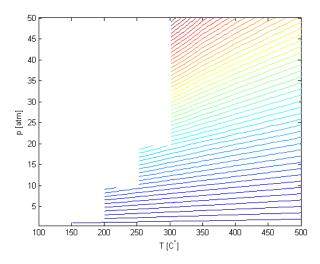


Figure 4.14: Contour plot of interpolation model for $\rho(T, p)$.

It is difficult to see from fig. 4.13 the valid region for (T, p) where ρ can be computed. A contour plot better illustrates the legal region:

```
>> contour(TTvar,ppvar,rhovar,50)
>> xlabel('T [C^\circ]')
>> ylabel('p [atm]')
```

Figure 4.14 illustrates the valid region for (T, p) for the interpolation model that we have developed: Suppose we want to find the vapor density of superheated steam at T = 375 °C and p = 3 atm — the point (T = 375 °C, p = 3 atm) is well within the valid region for (T, p), see fig. 4.14:

>> interp2(TT,pp,rho',375,3)

ans =

1.0235

The answer is: $\rho_{\rm 375\,^\circ C, 3\,atm} = 1.0235\,\rm kg/\,m^3.$

Chapter 5

Introduction to automation of tasks

5.1 Overview of learning goals

After having completed this chapter, you should have a clear understanding of why automation of tasks is useful, and how this can be achieved in Matlab. In particular, you should master:

- How to operate on strings,
- How to use the for-loop in Matlab,
- How to use the Matlab Editor, and restrictions on where you save files,
- What script files are, and know why they are useful.

5.2 Strings and basic string operations

It is useful to have some knowledge of strings in Matlab, since these are important for some types of task automation. A string is a sequence of characters. When defining the string, the sequence of characters must be surrounded by an apostrophe ':

```
>> clear
>> mystr1 = 'a';
>> mystr2 = 'abc';
>> whos
Name Size Bytes Class

mystr1 1x1 2 char array
mystr2 1x3 6 char array
```

Grand total is 4 elements using 8 bytes

We see that each character in a string takes up 2 bytes of memory. For our purpose, the most useful string operations are:

- Concatenation of strings s1 and s2: [s1, s2],
- Conversion of number n to a string s: s = num2str(n).

An example can be:

```
>> I = 2;
>> ['counter I has the value ', num2str(I)]
ans =
counter I has the value 2
```

A more thorough discussion of string operations is given in Appendix E.

5.3 Motivating examples

Suppose we want to find the sum

$$S_n = \sum_{k=1}^{n-1} \frac{1}{k^2}$$

for $n = 1, 2, 3, \cdots$. This is a quite laborious task if we are to do it manually. Although we can write the expression as

$$S_n = \sum_{k=1}^{n-1} \frac{1}{k^2} + \frac{1}{n^2} = S_{n-1} + \frac{1}{n^2},$$

this doesn't help much in saving work:

etc. $(S_n \text{ converges to } \frac{\pi^2}{6} \approx 1.6449).$

As another example, suppose we want to find the Fibonacci numbers f_k defined by:

$$\begin{array}{rcl}
f_1 &=& 1 \\
f_2 &=& 1 \\
f_{k+1} &=& f_k + f_{k-1} \text{ for } k \geq 2.
\end{array}$$

Straight ahead manual use of Matlab leads to:

etc.

In both of the examples above, we see that we repeat assignments that are almost the same. In the next section, we will introduce repetition statements which can be used to automate the computation of e.g. S_n and f_k .

Exercise 5.1 Before proceeding to the repetition statements, let us just note that we can compute S_n without using repetition loops. As an example, we can compute S_n as follows:

```
>> n = 20;

>> N = 1:n;

>> S = 1./N.^2;

>> cumsum(S)

ans =

Columns 1 through 5

1.0000 1.2500 1.3611 1.4236 1.4636

Columns 6 through 10

1.4914 1.5118 1.5274 1.5398 1.5498

Columns 11 through 15

1.5580 1.5650 1.5709 1.5760 1.5804

Columns 16 through 20

1.5843 1.5878 1.5909 1.5937 1.5962

Experiment with n (i.e. n) and see how large a value we must choose for n before the value \left|S_n - \frac{\pi^2}{6}\right| <
```

0.01.

In the Exercise above, we compute an array S which holds S_1 , S_2 , etc. See the Help browser for how Matlab function cumsum works.

5.4 The basic repetition statement

In order to repeat Matlab statements a specific number of times, the construct

```
for var = array
    statements
end
```

is used.¹ Here, for and end are keywords that should be used in verbatim², and array is an array. The columns of array are then assigned one at a time into var starting with the first column, and the statements are executed. Often, array is a row array, hence the "columns" are scalars. It should be remembered that e.g. 1:N creates a row array [1, 2, ..., N], thus the statement variable = 1:N is quite common.

Let us look at an example: suppose we want to print the numbers i and i^2 to the screen for $i \in \{1, \ldots, 10\}$: clearly we do not intend the meaning $i = \sqrt{-1}$. Since i by default denotes $\sqrt{-1}$, we use I as the *variable* name instead of i. The *statement* is $[I, I^2]$, which ensures that i and i^2 are displayed on the same line:

```
>> format compact
>> for I=1:10, [I, I^2], end
ans =
     1
            1
ans =
     2
            4
ans =
     3
            9
ans =
     4
           16
ans =
     5
           25
ans =
     6
           36
```

¹Alternatively, the statements can be written in a single line as for variable = array, statement 1, statement 2, ..., statement n, end.

²Verbatim = word for word, exactly as spoken or written.

ans = 7 49 ans = 8 64 ans = 9 81 ans = 10 100

Here, we have typed the command on a single line.

We can also type the commands on separate lines (*i* and \sqrt{i} , for a variation):

>> for I=1:10 [I, sqrt(I)] end ans = 1 1 ans = 2.0000 1.4142 ans = 3.0000 1.7321 ans = 4 2 ans = 5.0000 2.2361 ans = 6.0000 2.4495 ans = 7.0000 2.6458 ans = 8.0000 2.8284 ans = 9 3 ans = 10.0000 3.1623

In practice, we often use (much) more complex statements, and it is not practical to type the commands on the command line of Matlab. One reason is that it is easy to loose track of which commands we have carried out and which we have not carried out. Another reason is that it is easy to make mistakes: If we make mistakes in any line of the command sequence, we have to redo everything that depends on the erroneous command.

5.5 Scripts and the Matlab editor

A script is a file containing a sequence of Matlab statements. In order to run those statements, we simply type the name of the script file on Matlab's command line. Matlab's editor is convenient for typing and editing the statements in a script file, and we can start the editor by issuing the command >> edit on Matlab's command line. Figure 5.1 illustrates the opened Matlab Editor, containing three lines of Matlab commands.

Before running the script, it is necessary to save the file (File/Save). The script file must be stored either in a file in the *Current Directory* — see fig. 1.1 p. 4, or in Matlab's so-called *Path* — see help path. To run the script, either click on the *Save and run icon* of the Matlab Editor (see fig. 5.1), or type the file name in the Matlab Command Window. The file in fig. 5.1 has been saved with the file name RepetitionDemo1:

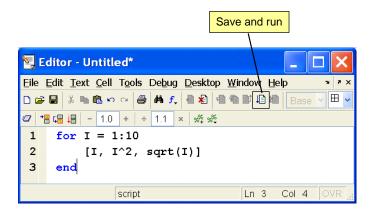


Figure 5.1: Matlab Editor window, containing a sequence of Matlab commands. Note: if the file has been saved, the "Save and run" icon changes to the "Run" icon.

>> I	Repetiti	ion	Demo1	
ans	=			
	1	1	1	
ans	=			
	2.0000		4.0000	1.4142
ans	=			
	3.0000		9.0000	1.7321
ans	=			
	4 1	16	2	
ans				
	5.0000		25.0000	2.2361
ans				
	6.0000		36.0000	2.4495
ans				
	7.0000		49.0000	2.6458
ans				
	8.0000		64.0000	2.8284
ans				
	9 8	31	3	
ans				
-	10.0000	1	00.0000	3.1623

It is considered good practice to *comment* script files in Matlab so that it is easier to understand what is taking place, when looking over the script at a later date. Characters appearing after the symbol % on a (command) line, are considered as comments — see fig. 5.2 which produces a multiplication table. The result of running the script (named RepetitionDemo2) is shown below:

>> Repet >> A A =	tiionD	emo2							
1	2	3	4	5	6	7	8	9	10
2	4	6	8	10	12	14	16	18	20
3	6	9	12	15	18	21	24	27	30
4	8	12	16	20	24	28	32	36	40
5	10	15	20	25	30	35	40	45	50
6	12	18	24	30	36	42	48	54	60
7	14	21	28	35	42	49	56	63	70
8	16	24	32	40	48	56	64	72	80
9	18	27	36	45	54	63	72	81	90
10	20	30	40	50	60	70	80	90	100

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	+∰ μ ∰ − 1.0 + ÷ 1.1 × ½ ½
1	8
2	% Computing the multiplication table
3	8
4	A = zeros(10,10); % Defines array A to store the result
5	
6	for $I = 1:10$
7	
8	······································
9	
10	- end
	script Ln 1 Col 1 OVR "

Figure 5.2: Matlab editor showing script file with comments and spaces.

In order to find the product 7×9 , find the intersection of the row with 7 in the first column, and the column with 9 in the first row. The result is 63.

5.6 Examples: Repetition and Sums

Suppose we want to use the **for** statement to find the sum $S_n = \sum_{k=1}^{n-1} \frac{1}{k^2} = S_{n-1} + \frac{1}{n^2}$ and the Fibonacci numbers f_k . First, we study how to find S_n . The following Matlab script will find S_n , fig. 5.3. The result is:

```
>> SumInvSquare
>> S
S =
             0
         0.25
      0.36111
      0.42361
      0.46361
      0.49139
       0.5118
      0.52742
      0.53977
      0.54977
      . . .
      0.63985
      0.63987
       0.6399
      0.63992
      0.63995
```

Similarly, we find the Fibonacci numbers as follows, fig. 5.4. The result is:

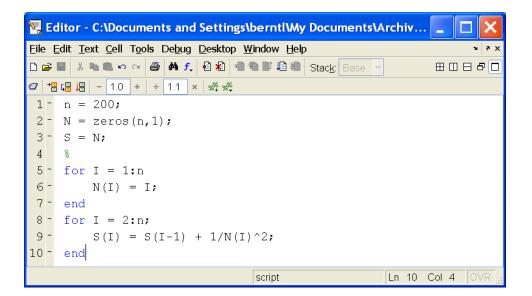


Figure 5.3: Script SumInvSquare for computing S_n .

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27 + ☐ ↓ ☐ ↓ ☐ 1.0 + ÷ 1.1 × ½ ½	
1 - k = 20;	
2 - f = zeros(k,1);	
3 - f(1) = 1;	
4 - f(2) = 1;	
5^{-} for I = 3:k	
6^{-} f(I) = f(I-1) + f(I-2);	
7 - end	
SumInvSquar× Fibonacci.m	
script Ln 7	Col 4 OVR

Figure 5.4: Script Fibonacci for computing f_k .

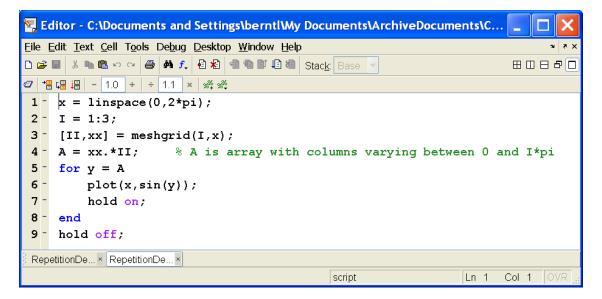


Figure 5.5: Matlab Editor with script where the repetition **array** has more than one row.

>> Fibonacci >> f
f =
1
1
2
3
5
8
13
21
34
55
89
144
233
377
610
987
1597
2584
4181
6765
0705

5.7 Example: Basic repetition over array*

This example uses some techniques introduced in Appendix D. Let us also study the case where variable = array, and array has more than one row. We consider the script of fig. 5.5.

The result of issuing the command

>> RepetitionDemo3

is fig. 5.6.

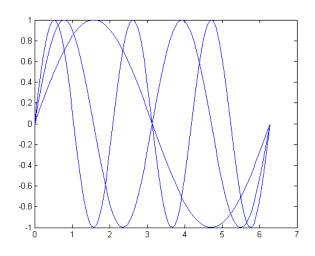


Figure 5.6: The result of running script RepetitionDemo3.

Chapter 6

Problems

Problem 6.1 We consider a *batch reactor*: a component A is converted to component R in a closed vessel through a chemical reaction. The concentration of component A, $c_A \pmod{1}$ can be described by the ordinary differential equation (ODE):

$$\frac{dc_A}{dt} = r_A,$$

where r_A is the so-called rate of production of component $A \pmod{(1 \cdot \min)}$. r_A is a function of c_A as given in fig. 6.1, taken from (Levenspiel 1972), p. 117.

- 1. Use the *Import Wizard* of Matlab to import the data from file ReactorDataLevenspiel.xls. (Hint: in Matlab, use command File/Import Data...).
- 2. Extract the values related to c_A and r_A from the **data** variable, and put them into variable names of your choice. Make sure that the chosen names are legal variable names in Matlab, and that the chosen variable names not already are used as function names.

(Note: the Excel spreadsheet gives $-r_A$, and not r_A . Since $-r_A$ is posivite, r_A is negative, which means that component A is *consumed*.)

Plot the experimental data to make sure that the data are correct. If you suspect that some data are outliers, then remove these faulty data. Replot the experimental data when possible outliers are removed.

- 3. Save the data in a .mat file. Clear the Matlab workspace, and check that the workspace is empty. Load the data from the file where you saved them.
- 4. Experiment with fitting models to the data for (c_A, r_A) . Use the following model types:
 - Polynomial least squares models of various orders. (Hint: polyfit, polyval.)
 - Spline interpolation models. (Hint: spline, ppval.)

N	Aicrosoft Excel - I	ReactorDataLevens	piel										×
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	A	В	С	D	E	F	G	Н		J	K	L	
1	Experimenta	al batch reacto	or data:	A> R									
2	c_A, mol/l	0.1000	0.2000	0.3000	0.4000	0.5000	0.6000	0.7000	0.8000	1.0000	1.3000	2.0000	
3	-r_A, mol/I min	0.1000	0.3000	0.5000	0.6000	0.5000	0.2500	0.1000	0.0600	0.0500	0.0045	0.0042	
4													
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Figure 6.1: Experimental batch reactor data for reaction with stoichiometry $A \rightarrow R$. Taken from (Levenspiel 1972), p. 117.

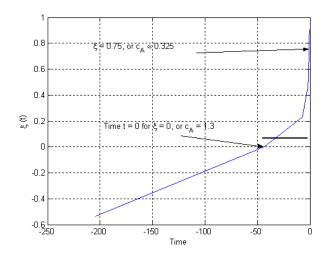


Figure 6.2: Extent of reaction $\xi(t)$ for reactor data based on trapezoidal integration.

• Hermitian cubic polynomials. (Hint: pchip, ppval.)

What model types are acceptable? (**Hint**: Is it acceptable that r_A becomes negative?)

5. * The *extent* ξ_A of reaction is defined as follows:

$$\xi_A(t) \triangleq \frac{c_A(0) - c_A(t)}{c_A(0)}.$$

Find $\xi_A(t)$, and plot $\xi_A(t)$ as a function of t when $c_A(0) = 1.3 \text{ mol/l}$. At $t = 0, \xi(0) = 0$. At what time does $\xi_A(t)$ rise above 0.75?

(**Hint**: The principle is based on finding $c_A(t)$ from the solution of the ODE

$$\frac{dc_A}{dt} = r_A \iff \frac{dc_A}{r_A(c_A)} = dt \iff \int_{c_A(0)}^{c_A(t)} \frac{dc_A}{r_A(c_A)} = \int_0^t dt = t.$$

This leads to $F(c_A(t), c_A(0)) = t$, which may be inverted to give $c_A(t) = G(t, c_A(0))$. Then we can find $\xi_A(t)$ from its definition. A possible procedure is as follows:

- Use array operations to compute the array containing $1/r_A$.
- Use Matlab function cumtrapz(x,y), which computes an array with a numeric approximation of $\left(\int_{x_1}^{x_2} y(x) dx, \int_{x_1}^{x_3} y(x) dx, \ldots, \int_{x_1}^{x_N} y(x) dx\right)$ where x_i is a discrete value of x. (Hint: In the current problem, x is c_A and y is $1/r_A$).
- Plot $c_A(t)$. (Hint: Since $\int_{c_A,1}^{c_{A,i}} \frac{dc_A}{r_A(c_A)} = F(c_A(t_i), c_{A,1}) = \int_0^t dt = t_i$, each pair $(F(c_A(t_i), c_{A,1}), c_{A,i}) \equiv (t_i, c_{A,i})$. We can use this fact to plot $(t, c_A(t))$. Verify that the result is similar to in fig. 6.2.)
- By zooming in at the points of fig. 6.2 where $\xi_A = 0$ (i.e. $c_A = 1.3$, which indicates time t = 0) and the time where $\xi_A = 0.75$ (the sought extent of reaction), find the time for achieving 75% extent of reaction.
- 6. * The calculations in the previous question are rather coarse due to relatively few data points. We can achieve smoother (but not necessarily more accurate) calculations as follows:
 - Use array operations to compute the array containing $1/r_A$.
 - Use your preferred interpolation model for $(c_A, 1/r_A)$.
 - Compute $1/r_A$ for an extended set of data points in the interval $(\min(c_A), \max(c_A))$ (e.g. using the linspace function of Matlab) with the help of the interpolation model.

- Use Matlab function cumtrapz(x,y) to compute the relevant array containing $\int_{c_{A,1}}^{c_{A,i}} \frac{dc_A}{r_A(c_A)} = F(c_A(t_i), c_{A,1}) = \int_0^t dt = t_i$, and plot the smoothed pair $(t_i, c_{A,i})$, as well as $(t_i, \xi_{A,i})$.
- Find the time necessary to achieve 75% extent of reaction. Is the answer different from the result in the previous question?

Part II

Exploiting the power of Matlab

Chapter 7

Revisiting Part I

(If you do not recall the following topics, you should re-check Part I or use Matlab's help facilities.)

The Matlab environment See Chapter 1.

- Command window
- Help window
- Workspace window
- Array editor

Basic operations

- The array data type (scalars, row and column arrays, 2 dimensional arrays), see Section 2.2.
- Assignment, Section 2.3.
- Variables, see Section 2.5
- Special variables: pi, i, j, etc., Table 2.1 p. 9.
- How to create arrays: zeros, ones, eye, diag, colon operator, rand, randn, linspace, logspace, size, length. See Sections 2.6.2 2.6.5, and 2.10.
- Selecting parts of arrays: A(2:3,5:8), A(1,3:end); building/concatenating arrays. See Section 2.11.
- Simple writing and reading of data from file: save, load, import wizard, see Section 2.7.
- Housekeeping of data: clear, who, whos, what, commands, and files. See Section 2.12.
- Basic binary array operations: +, -, .*, ./. See Table 2.6 p. 34.
- Basic unary array functions: sin, cos, tan, exp, log, log2, sqrt, abs, ceil, floor, transpose, etc. See Tables 2.7 p. 36, 2.8 p. 37, 2.9 p. 37, 2.10 p. 39, and 2.11 p. 40.
- Arrays vs. vectors and matrices, see Section A.

Plotting See Table 3.1 p. 43. In particular, note:

- Basic plotting: plot, semilogx, semilogy, loglog, linestyle, marker, color
- Tailoring your plot: axis, title, xlabel, ylabel, legend, hold, subplot
- figure, close

Simple data analysis* See Table 4.1 p. 57. In particular, you should check out rudimentary commands such as max, min, mean, sum, as well as model fitting techniques such as polyfit, polyval. See also how you can use the find command to spot outliers and missing data (NaN).

Introduction to automation of tasks

- Basic string operations. See Tab. E.1 p. 174, and Tab. E.2 p. 176 for a more detailed treatment.
- for-loops, see Section 5.4.
- Scripts, see Section 5.5.

More program flow control

8.1 Introduction

In Part I of these lecture notes, we started to look into Matlab programming by developing Matlab scripts using the Matlab editor, and automating repetitive tasks using the **for** loop. In this chapter, we will start by looking into program flow control by the **if** statement. To do this, we must first discuss the concepts of logical variables and relations, as well as logical expressions. Then we are ready for discussing the **if** statement. Thereafter, we will discuss the **switch** statement, and finally we will discuss an alternative repetition statement: the **while** loop.

8.2 Logical variables in Matlab

Writing scripts and functions, we often need to choose between alternative actions depending on the result of one or more tests. To do this we need relational and logical operators and logical (Boolean) variables.

Traditionally, Matlab only supported *floating point* numbers, but later versions of Matlab support various other data types similar to other programming languages, such as logical (Boolean) variables. Logical variables can only have two possible outcomes, **true** or **false**. Since the early Matlab versions did not support logical variables, such variables were "emulated" by interpreting the floating point number 0 to be **false**, and any floating point number $\neq 0$ to be **true**.

8.3 Relations: functions and operators

In programming languages, relations are statements which are either true or false, i.e. statements where the result is a logical quantity. Table 8.1 lists Matlab functions and operators for relational statements.

Matlab's parser translates the *operator expressions* into the function form. So if we write x > y, Matlab sees the function gt(x,y). The output of a relational expression is a *logical variable*, having the value true or false.

Examples Consider the following simple examples:

Table 8.1:	Matlab relational	functions and operators	. In the examples, i	t is assumed that $x = 3$	y = 4.
	Math operator	Matlab function	Matlab operator	Example	

Math operator	Matlab function	Matlab operator	Example
<	lt	<	x < y> true
\leq	le	<=	x<=y -> true
=	eq	==	x==y -> false
\neq	ne	~=	x~=y -> false
>	gt	>	x>y -> false
\geq	ge	>=	x>=y -> false

>> x = 2; >> y = 3; >> X = [1,2,3]; >> Y = [4,3,2]; >> x < y ans = 1 >> y < x ans = 0 >> x == y ans = 0

We can also give name to the result from a statement:

```
>> logic = x == y
logic =
0
>> logic = x < y
logic =
1</pre>
```

We can consider relations between arrays of equal size, where e.g. $[x, y, z] \leq [u, v, w]$ is interpreted as $[x \leq u, y \leq v, z \leq w]$:

>> X < Y
ans =
 1 1 0
>> X == Y
ans =
 0 0 0
>> log = X >= Y
log =
 0 0 1

8.4 Logical expressions

In the previous section, we considered statements about the *ranking* of floating point numbers. The result was logical quantities which assessed the truth of these statements.

It is also useful to work with logical expressions, i.e. combinations of *logical variables*. Table 8.2 gives an overview of Matlab operators for combining logical variables.

Any(a) is true if at least one element in a is true. All(a) is true if all elements in a are true.

Examples Consider the following simple examples:

```
>> 1 & 0
ans =
0
>> 1 & 1
ans =
1
```

Table 8.2: Matlab logical functions and operators.	In the examples, it is assumed that $a = [1, 0, 0]$,
b = [1, 1, 0].	

Logical operator	Matlab function	Matlab operator	Example
AND	and	&	a & b -> [1 0 0]
OR	or		a b> [1 1 0]
Exclusive OR	xor		xor(a,b)> [0 1 0
NOT	not	~	~a> [0 1 1]
Any	any		any(a)> 1
All	all		all(a)> 0

```
>> 1 | 0
ans =
    1
>> xor(1,0)
ans =
    1
>> any([1,0,1])
ans =
    1
>> all([1,0,1])
ans =
    0
>> all([1,1,1])
ans =
    1
```

Often, the logical expressions are combined with relations:

Finally, we can make the logical expressions simpler by introducing logical variables:

8.5 Making logical choices: the if statement

8.5.1 Motivating example

As a motivation, consider the formulae for the root $x \in \mathbb{C}$ of a second order polynomial $ax^2 + bx + c = 0$:

$$x = \begin{cases} \frac{-b \pm \sqrt{b^2 - 4ac}}{2a}, & a \neq 0\\ -\frac{c}{b}, & a = 0, b \neq 0\\ \emptyset & a = 0, b = 0, c \neq 0\\ \mathbb{C} & a = 0, b = 0, c = 0 \end{cases}.$$

We see that the answer depends on the values of coefficients (a, b, c). In order to write a Matlab script for the solution x with (a, b, c) is known, we need to be able to test which formulae is relevant.

8.5.2 Structure of if statements

The standard test statement in most programming languages is the **if** statement. In Matlab, the basic syntax is as follows:

```
if expression
statements
end
```

Here, **expression** is a logical variable/expression. If the logical variable is *true*, then the statements are executed. If not, they are not executed.

In many cases, we want to test for more than one logical variable/expression. We then use the following extended version of the **if** statement:

```
if expression1
   statements1
elseif expression2
   statements2
elseif expression3
   statements3
...
else
   statements
end
```

8.5.3 Example: use of if statement

The following script will find the root of the second order polynomial, fig. 8.1. Below are some examples of the use of this script:

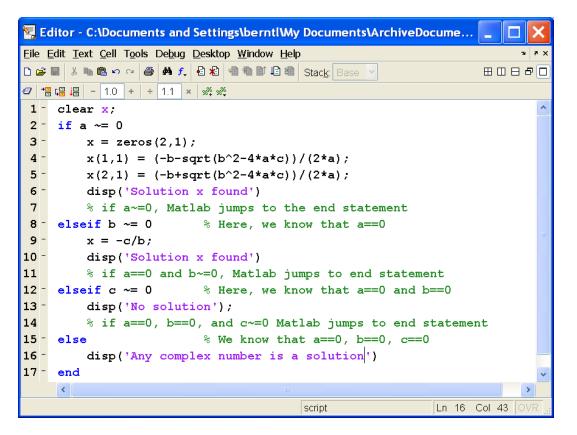


Figure 8.1: Script for finding root of $ax^2 + bx + c = 0$, where a, b, c are specified in the Matlab Command Window.

```
-0.2500 - 0.9682i

-0.2500 + 0.9682i

>> a=0;b=2;c=-3;

>> rootscript

Solution x found

>> x

x =

1.5000

>> a=0;b=0;c=2;

>> rootscript

No solution

>> a=0;b=0;c=0;

>> rootscript

Any complex number is a solution
```

8.6 Logical choices: the switch statement

When we need to test for a logical variable, or for a (complex) logical expression, we use the if statement as discussed in the previous section.

In many cases, the if statements are of the following type:

```
if variable == case_value1
   statements1
elseif variable == case_value2
   statements2
elseif variable == case_value3
```

```
statements3
...
else
statements
end
```

This particular kind of if statement structure can be expressed in the slightly more convenient form of a switch statement:

```
switch variable
   case case_value1
      statements1
   case case_value2
      statements2
   case case_value3
      statements3
...
   otherwise
      statements
end
```

8.7 Repetition loop: the while statement

8.7.1 Motivating example

We have already met the **for** statement:

```
for var = array
    statements
end
```

It is characteristic for the for statement that the variable var loops through the values of array a predetermined number of times.

Sometimes, we do not know in advance how many times we want to repeat an operation. As an example, suppose we want to find

$$S = \sum_{k=1}^{\infty} \frac{1}{k^2} = \frac{\pi^2}{6} \approx 1.6449.$$

to a certain accuracy. We thus use the approximation $S_n = \sum_{k=1}^n \frac{1}{k^2} = S_{n-1} + \frac{1}{n^2}$, where n is unknown but implicitly determined by the required accuracy $|S_n - S_{n-1}| < 10^{-6}$.

8.7.2 Structure of while statement

To find an approximation S_n , we can then use the while statement:

```
while expression
statements
end
```

Here, the statements are executed as long as expression is true.

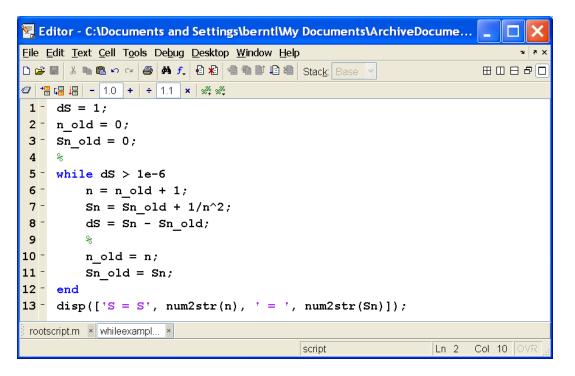


Figure 8.2: Script for finding $S_n = \sum_{k=1}^n S_{n-1} + \frac{1}{n^2}$ to a prespecified accuracy of $|S_n - S_{n-1}| < 10^{-6}$.

8.7.3 Example: use of while statement

The following example illustrates how we can compute S_n such that the accuracy requirement is fulfilled, fig. 8.2. Running this script gives the following answer:

>> whileexample S = S1000 = 1.6439

A possible problem with the use of the while statement, occurs if we use an erroneous test expression which never is fulfilled: for that case, the while loop will continue forever, and the only way to stop the script execution may be to hit the Ctrl+C key.¹

8.7.4 Breaking out from for loop

Finally, it should be noted that the same construct can be achieved by combining a for loop with an if statement, fig. ??. Executing this script leads to:

>> forwhileemulate
S = S1000 = 1.6439

The disadvantage of using the **for** statement here, is that we do not know in advance how large N should be. The advantage of using the **for** statement, is that the loop will not be repeated indefinitely even when we make a mistake in the program code.

 $^{^{1}}$ Ctrl+C is achieved by holding down the Ctrl key on the computer keyboard, and while your are doing this, press the C key. This is an emergency action only, which usually (but not always) works.

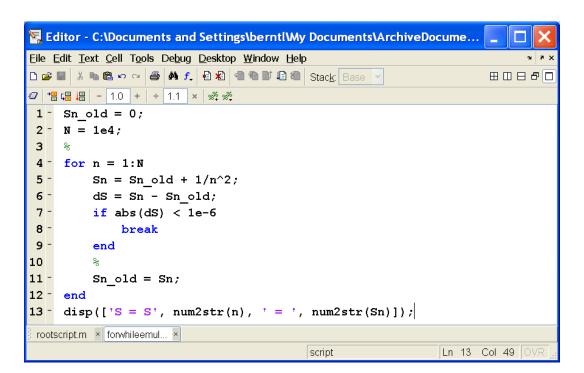


Figure 8.3: Script with for loop, where an if statement is used to beak out of the loop.

Writing your own functions

So far, we have used some of Matlab's built-in functions. In Matlab it is very easy to write your own functions if Matlab does not have one that suits your needs.

9.1 Example: Distance to horizon

9.1.1 Basic case

If you look out to the sea, the distance x_h to the horizon depends on how high above the sea level your eyes are, fig. 9.1. Approximating the earth with a perfect sphere and neglecting refraction, the

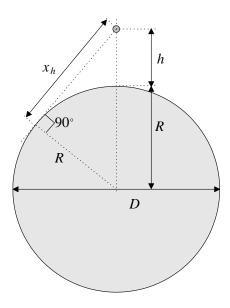


Figure 9.1: Distance x_h to visible horizon from height h. D and R are the diameter and radius, respectively, of the earth.

Pythagorean law leads to:

$$R^{2} + x_{h}^{2} = (R+h)^{2} \Leftrightarrow x_{h} = \sqrt{h(2R+h)} = \sqrt{h(D+h)},$$

where h is the elevation of your eyes above the sea level and D is the diameter of the earth. x_h is the length of the straight line between your eyes and the horizon. Knowing that the earth's diameter $D \approx 12000$ km, we can create a Matlab function for this. We open a new document in the editor by clicking the leftmost button on Matlab's toolbar and type:

function xh = horizon(h)

```
% Distance to the horizon
% Usage: xh = horizon(h)
d = 12e6; % Earth's diameter
xh = sqrt(h.*(h+d));
We save this file as horizon.m. Now we can test it:
>> horizon(2)
ans =
4899
```

We have thus chosen to name the function horizon; we thus also save the function file and give it the name horizon.m.

Note that we have used element by element operations $(.*, ./, .^)$ rather than matrix operations $(*, /, ^)$. This ensures the right answer if we want to calculate the function for an array of arguments at once, as we might want to do for plotting:

```
>> hp = 1:1000;
>> plot(hp,horizon(hp))
```

The first comment lines are echoed when we write help horizon:

```
>> help horizon
Distance to the horizon
Usage: xh = horizon(h)
```

Maybe we want to be able to find the distance to the horizon on the moon or some other planet. We can easily modify the function so that it takes the planet diameter as a second argument:

```
function xh = horizon2(h,d)
% Distance to the horizon
% Usage: xh = horizon(h,d)
% h : Observer distance above surface
% d : Planet diameter
% xh : Distance to horizon
xh = sqrt(h.*(h+d));
```

Note that Matlab searches for the *file* horizon2.m when we write:

```
>> horizon2(2,6e6)
ans =
3464.1
```

We should thus save the new function in a file horizon2. The function name itself (horizon2) is not important in itself, but it is considered good practice to let the function name be equal to the file name.

We can make a version that uses the earth's diameter as default if we give only h as the argument. As long as only one argument is used, the user will not note the difference from the original horizon.m, so we overwrite this file. Below, nargin knows the number of input arguments that the user has used:

```
function xh = horizon(h,d)
% Distance to the horizon
% Usage: xh = horizon(h) and xh = horizon(h,d)
% h : Observer distance from surface
% d : Planet diameter Default: 12e6 (Earth)
```

```
% xh : Distance to horizon
if nargin < 2
    d = 12e6; % Gives d a value if the user does not
end
xh = sqrt(h.*(h+d));</pre>
```

9.1.2 Returning more than one result

We may also return more than one value:

```
function [xh, angle] = horizon3(h,d)
% Distance to the horizon
% Usage: xh = horizon(h) and xh = horizon(h,d)
% h : Observer distance from surface
% d : Planet diameter Default: 12e6 (Earth)
% xh : Distance to horizon
% angle: angle between radius to horizon and to observer (degrees)
if nargin < 2
    d = 12e6; % Gives d a value if the user does not
end
xh = sqrt(h.*(h+d));
angle = atan(xh/(d/2))*180/pi;
Test:</pre>
```

From an airplane at 10 000 m above Oslo (60° N) you can see most of the way to Trondheim (63.5° N). We can ask for the second output value by specifying it in the calling statement (otherwise, it is discarded):

We can obtain the number of output arguments from within the function. This can save some computer time if some outputs require a lot of calculations. Below, **nargout** knows the number of output arguments that the user has asked for:

It is good practice to test that the arguments given by the user are valid:

```
function xh = horizon(h)
% Distance to the horizon
% Usage: xh = horizon(h) and xh = horizon(h,d)
if h < 0
        error('Argument must be positive')
end
xh = sqrt(h.*(h+d));</pre>
```

9.1.3 Exercises

Exercise 9.1 The Fibonacci numbers f_k are defined by:

$$f_1 = 1$$

$$f_2 = 1$$

$$f_{k+1} = f_k + f_{k-1}, \quad k > 2.$$

Thus the first 8 Fibonacci numbers are: 1, 1, 2, 3, 5, 8, 13, 21.

- Write a function fib(n) that returns Fibonacci number n.
- Let the function return the vector of numbers up to and including f_n in an optional second return variable. Also make the function able to take a two-element array as an optional second input variable. The two elements of the vector should be used as alternative starting values. E.g.: [fn,s]=fib(6,[8,3]) should return fn = 39 and s = [8,3,11,14,25,39].

Exercise 9.2 When calculating the distance to the horizon x_h from an elevation h, the distance from the "shore" to the horizon will also be approximately x_h when $h \ll D$.

- Write a function that computes h when x_h is known.
- Approximately, how high above the ground do you need to be to be able to see from one end of Lake Mjøsa¹ to the other end? Assume that Lake Mjøsa is 120 km long.
- Approximately, how high above the ground do you need to be to be able to see from the southern shore of Norway to the northern shore of Denmark? Assume that the distance is 150 km.
- A possible approximative expression for h is

$$h \approx \frac{1}{D} x_h^2.$$

Write another function for computing h based on this approximation. Compare the expression with the exact expression for a number of distances x_h , and suggest how large x_h can be before the simple formula $h \approx x_h^2/D$ becomes inaccurate.

¹Lake Mjøsa is the largest inland lake in Norway.

9.2 Workspaces and variables

Like most programming languages, Matlab uses several *workspaces* or *contexts*. As long as you issue commands in the *command window* or use *scripts*, you work in the *base workspace*. All variables that you have created are available to you until you clear them or until the session is terminated. Inside a function, you do not have access to the variables in the base workspace. This function will never work, even if you define the variable a before calling the function:

```
function s = testsum(x)
s = x + a;
```

We can try:

```
>> a = 4;
>> s = testsum(3)
??? Undefined function or variable 'a'.
Error in ==> C:\HiT\Matlabkurs\testsum.m
On line 2 ==> s = x + a;
```

Inside the function, Matlab does not "see" the **a** from the base workspace. Information to the function is passed only through the function's arguments. Likewise, variables that you create or change inside a function have no direct influence on the base workspace. Information can pass back only through the returned variables. Only the sequence of input and output variables is significant. Which variable names you use have no effect.

Example

Consider these examples:

```
function s = vartest(x,y)
a = 2;
s = x + a*y;
```

We test the function:

```
>> a = 4;
>> x = 1; y = 2;
>> z = vartest(x,y)
z =
5
>> a
a =
4
>> z = vartest(y,x)
z =
4
```

We see that the assignment a=2; in the function workspace does not affect the variable a in the base workspace. Furthermore, interchanging the sequence of x and y in the call gives two different results. In fact, we could have used some totally different variable names in the call:

>> arg1 = 1; arg2 = 2;z = vartest(arg1,arg2) z =

```
5
>> z=vartest(1,2)
z =
5
```

Assigning the answer to the variable \mathbf{s} in the function's workspace has no effect in the base workspace:

```
>> s
??? Undefined function or variable 's'.
```

Understanding this segregation between workspaces is essential to understanding how functions work.

9.3 Loops vs. matrix operations

Say we want to calculate the inner product of two vectors a and b. The inner product is defined as

$$a \cdot b = \sum_{i=1}^{n} a_i b_i,$$

where n is the number of elements of a and b. Say we represent both a and b as Matlab column vectors. Then we can calculate the inner product by using the formula above, but we can also exploit the fact that the inner product is the same as the matrix product between the transpose of a with b. The latter gives a much simpler expression:

```
% Inner product example
a = [3;6;-2;1];
b = [-3; 0; 4; 1];
% Loop formulation
s1 = 0;
for k = 1:length(a)
        s1 = s1+a(k)*b(k);
end
% Matrix/vector expression:
s2 = a'*b;
```

In older versions of Matlab, the matrix expression was much quicker than the loop form, but from Matlab v. 6.5, the loop runs almost as fast. You should use the formulation that results in the most easily readable code. Very often, this is the matrix form.

Exercise 9.3 Element C_{ij} in the product $C = A \cdot B$ is defined as $C_{ij} = \sum_{k=1}^{n} A_{ik} B_{kj}$.

• Write function C = matmult(A,B) that multiplies two square matrices without using Matlab's matrix operations. Test that the matrices are indeed square and of the same size, and abort using the error statement if this is not the case.

(Hint: The value of C(i,j) is the inner product of A(i,:) with B(:,j).

• Using two random matrices (A = rand(3); B= rand(3);), compare the output to A*B.

Advanced use of the Matlab editor

10.1 Errors and the debugger

I really hate this damn machine. I wish that they would sell it. It never does quite what I mean, but only what I tell it! — The programmer's lament

We inevitably make mistakes when we write functions and scripts in Matlab. Some are *syntax errors*: we write something that does not make sense to Matlab, no matter what the variables contain. Matlab spots these errors before it tries to execute your code. We may for instance forget the multiplication operator *****:

>> a = 2b
??? a = 2b
|
Error: Missing operator, comma, or semicolon.

Matlab tells you what it believes is wrong and marks the spot where it had to give up.

Syntax errors are usually relatively easy to correct. However, even when there are no syntax errors in the Matlab code, there may be logical errors or algorithmic errors which occur during run-time and lead to unintended results or even program crash — these are not always easy to find. Run-time errors depend on the contents of your variables. When the running of the Matlab code terminates, Matlab reverts to the base workspace, and we no longer have access to the variables in the function workspace. This makes it difficult to find out what went wrong.

One simple remedy is to have the function print intermediate results, either by using the **disp** statement, or simply by *omitting the semicolon* at the end of selected statements. However, these are not good strategies for finding errors, because when the code has been debugged and errors fixed, we need to remove the **disp** commands and insert semicolons; this may be a major job.

A more advanced solution is to use the *debugger*. The debugger is a powerful tool to help you understand what is going on inside your functions, fig. 10.1. In order to activate the debugger, a *breakpoint* must have been defined in the Matlab file, see fig. 10.1. With one or more breakpoints in the Matlab file, the debugger can be started as follows:

- if you want to debug a script file, either click on the (Saven and) Run icon in the editor, or type the name of the script file in the command window;
- if you want to debug a function file, either call the function (with appropriate arguments) from the command line, or run a script that calls the function.

The debugger allows you to stop execution inside a script or function. You can stop at a specified line, or when a warning or error is triggered. A breakpoint marks one or more lines of code where you

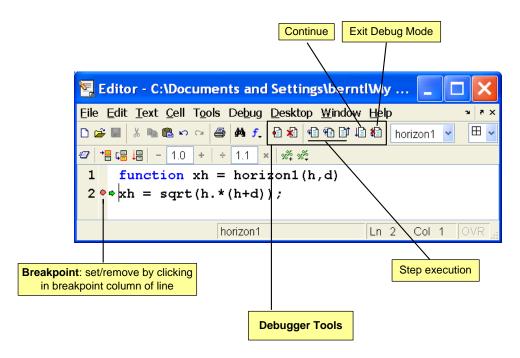


Figure 10.1: Some debugger tools.

want execution to pause. You define the breakpoints by clicking the button showing a red circle on the edge of the paper, or from the Breakpoints menu.

When the debugger halts execution, you will see a special prompt in the *command window*:

K>>

At the same time: in the *editor window*, a green arrow points to the line about to be executed, see fig. 10.1. You may now examine the contents of variables. This can be done by letting the cursor hover over a variable in the editor window — the value of the chosen variable will then appear in a yellow box over the editor window. Or you can browse the workspace browser and use the array editor. Or you can type commands in the command window:

You may *step* from line to line or into subfunctions, using one of the three Step execution icons, see fig. 10.1. All debugger buttons show yellow information balloons when you let the cursor hover over them, so finding your way around the debugger is quite intuitive.

To get out of the debugger, click the Exit Debug Mode icon (fig. 10.1, or enter the command

🔄 Editor - C:\Documents and Settings\berntl\My Documents\ArchiveDocuments\Courses\MSc\Matla	. 💶 🗖 🗙
Eile Edit Text <u>C</u> ell T <u>o</u> ols De <u>b</u> ug <u>D</u> esktop <u>W</u> indow <u>H</u> elp	X 5 K
🗅 🗃 🖩 🐇 ங 🎕 🕫 斗 🞒 👫 🗜 🖥 📽 🖷 🎕 🕼 🎜 Kack: Base 🔮	⊞ 🗆 🖯 🗗 🗖
$1 - Sn_old = 0;$	
2 - Sn_new = 1;	
$3 - Sn_new = 0;$	
4 - N = 1e4;	
5 %	
6 - for n = 1:N	
$7 - Sn = Sn_old + 1/n^2;$	
$8 - dS = Sn - Sn_old;$	
9 - if abs(dS) < 1e-6	
10 - break 11 - end	
11 - end 12 %	
13 - Sn old = Sn;	
14 - end	
15 - disp(['S = S', num2str(n), ' = ', num2str(Sn)]);	
script Ln 10	Col 14 OVR

Figure 10.2: Example of Matlab code with "lint": statement $Sn_new = 1$; is superfluous and should have been removed.

in the command window.

Finally: the **Debug** menu in the Editor window should also be consulted. This menu contains the same operations as can also be carried out using the debugger icons.

The debugger is useful for understanding runtime errors, but even more so for finding out why your Matlab code does not return the answer that you expected. If you are serious about using Matlab as a tool, learning to use the debugger is mandatory.

10.2 De-linting the code

Developing complicated Matlab code in the Matlab editor will to some degree involve some trial and error: we try with some code, then change our mind and modify the code. In this development, it is very common that we forget to remove unnecessary code, or that we forget to define some variables. This problem is illustrated in the (not very realistic) example of fig. 10.2.

We can use the M-Lint tool to find such code: type Tools/Check Code with M-Lint, and the report in fig.10.3 is provided.

As the report in fig. 10.3 indicates, the statement on line 2 of the code in fig. 10.2 should be removed. (Clearly, Matlab can not guess that statement on line 3 is without interest.)

10.3 Structuring the file

From Matlab 7 of, it is possible to structure Matlab code in the editor by defining text cells. Essentially, a text cell is initiated by putting the symbol %% (double % character) in the first position of a line. The cell ends on the line preceding the next %% symbol. After the %% symbol, a space should be inserted, followed by a descriptive text. The Matlab editor marks a cell by framing it with a yellow box: when you put the cursor in a cell, the frame is shown. In order for this to work, the *Cell Mode* must have been *Enabled*, see the Cell menu of the Matlab editor. Figure 10.4 illustrates a simple script where cells have been defined.

A simple use of such cells is use it for browsing a long Matlab file: click on the Show cell titles icon, fig. 10.5

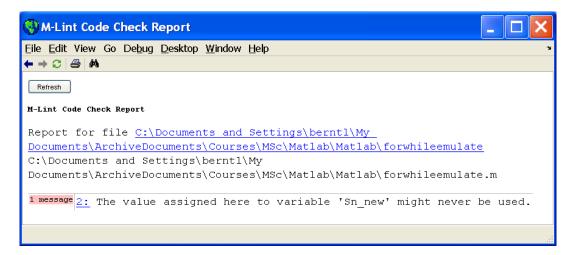


Figure 10.3: Report from checking the code of fig. 10.2 with M-Lint.

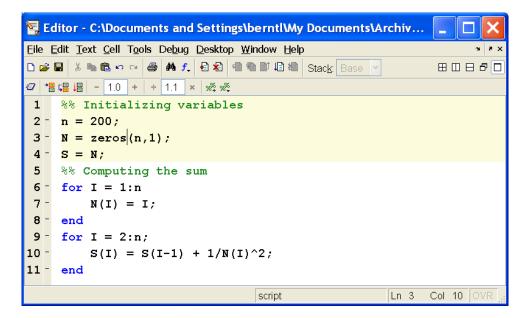


Figure 10.4: Example of text cells in the Matlab editor.

🔄 Editor -	C:\Docur	nents a	and Setti	ngs\berntl\My	Documents\Archiv	v 💶 🗖 🗙
<u>Eile E</u> dit <u>T</u> e	xt <u>C</u> ell T <u>o</u>	ols De <u>l</u>	oug <u>D</u> eskto	op <u>W</u> indow <u>H</u> elp)	X 5 K
ا & 🖬 😂 🗅	h 🗈 ហ 🖓	<i>a</i> M	🖵 🗗 🕈		Stac <u>k</u> : Base 🍸	
🖉 🕂 📮	- 1.0 +	÷ 1.1	1 × %**	% •		
			Show	cell titles		
Tools for u	using ce	lls	CHOW			

Figure 10.5: Tools for taking advantage of cells in Matlab files. The Show cell titles tool can be used for browsing through the cells: clicking this tool, gives a list of the cells.

and a list of the cell headlines (descriptive text following %% symbol) pops up. By clicking on the relevant cell title, the cursor is positioned in the cell head. This allows for easy navigation of large Matlab files.

However, cells can be used for much more. It is e.g. possible to insert markup commands in the Matlab file, e.g. making text boldface, etc., and also markup text using T_EX syntax so that it can be interpreted as mathematics. The Matlab file can then be published as an HTML document, a IAT_EX document, a Word document, etc., in such a way that the published document documents the code with headlines, section heads, with computed results, etc.

Data structures in Matlab

Historically, the matrix was the only data type in Matlab. Vectors and scalars are special cases of the general matrix. Since Version 5, some important new data structures have arrived. One is a the multi-dimensional array, which just extends the matrix to more than two dimensions. More important newcomers are the structure and the cell array.

11.1 Structures

A *structure* is a data structure that can hold diverse data types, not necessarily numbers, and with named "data containers" called *fields*, similar to a record with fields in a database. Example:

```
>> tank1.diameter = 0.5;
>> tank1.height = 0.4;
>> tank1.type = 'cylinder';
```

This is useful if we want to collect several items of information about some object. We can now tell e.g. a function to calculate the volume: all it needs to know about tank1 by using the *tank1 structure* as function argument:

```
function V = volume(a)
switch a.type
    case 'cylinder'
        V = pi*a.diameter^2*a.height/4;
    case 'box'
        V = a.length*a.width*a.heigth;
    case 'sphere'
        V = pi*a.diameter^3/6;
    otherwise
        error('unknown type')
end
```

Here we have used a new flow control structure, switch ... case ... otherwise ... end.

```
>> volume(tank1)
ans =
    0.025
>> tank2.type = 'sphere';
>> tank2.diameter = 2;
>> volume(tank2)
ans =
    4.1888
```

Structures may be used to limit the number of arguments needed by a function. This can be achieved by collecting all parameters in a parameter structure.

11.2 Cell arrays

A *cell array* is a data structure that can hold diverse data types, not necessarily numbers. The File Import Wizard returns text and data as cell arrays, as we saw in Part I of the course. We can collect items into a cell array by enclosing them in curly brackets:

```
>> A = rand(4);
>> C = {A sum(A) prod(prod(A))}
C =
[4x4 double] [1x4 double] [8.7912e-006]
```

We access the contents of the individual cells by curly brackets:

```
>> b = C{2}
b =
2.1251 1.8055 2.8766 2.8019
```

The Matlab Help contains much more information about structures and cell arrays.

Handle graphics

As you have seen, simple plotting in Matlab is quite easy. At the same time, Matlab graphics is very flexible, and you have great freedom to get the exact graphics that you need. Here, we will demonstrate how to modify some properties of graphs. You can find out more by consulting Matlab's help system.

We consider the plotting of two Taylor series for $\cos x$, and compare them to the exact function:

$$\cos x \approx \begin{cases} T_1 = 1 - \frac{1}{2}x^2 + \frac{1}{24}x^4 \\ T_2 = -\left(x - \frac{\pi}{2}\right) + \frac{1}{6}\left(x - \frac{\pi}{2}\right)^3 \end{cases}$$

To make sure that we can distinguish between the three lines, we use markers for T_1 and T_2 :

x = linspace(0,pi/2,25); f = cos(x); T1 =1 - x.²/2 + x.⁴/24; T2 = -(x-pi/2)+(x-pi/2).³/6; figure(1) h = plot(x,f,x,T1,'*',x,T2,'+');

The result is displayed in fig. 12.1.

Note the form of the plot command. The variable h now contains *handles* to each of the three curves. In fact, h is an *object* with certain *properties*, and we can get a list of all of the properties for the first curve by typing get(h(1)):

>> get(h(1))

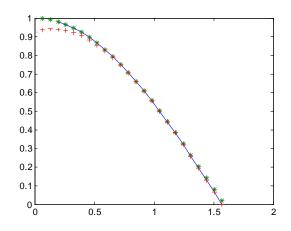


Figure 12.1: Comparison of $\cos x$, and two Taylor expansions of $\cos x$.

```
Color: [0 0 1]
         EraseMode: 'normal'
         LineStyle: '-'
         LineWidth: 2
            Marker: 'none'
        MarkerSize: 6
   MarkerEdgeColor: 'auto'
   MarkerFaceColor: 'none'
             XData: [1x25 double]
             YData: [1x25 double]
             ZData: [1x0 double]
      BeingDeleted: 'off'
     ButtonDownFcn: []
          Children: [Ox1 double]
          Clipping: 'on'
         CreateFcn: []
         DeleteFcn: []
        BusyAction: 'queue'
  HandleVisibility: 'on'
           HitTest: 'on'
     Interruptible: 'on'
          Selected: 'off'
SelectionHighlight: 'on'
               Tag: ''
              Type: 'line'
     UIContextMenu: []
          UserData: []
           Visible: 'on'
            Parent: 150.0048
       DisplayName: ''
         XDataMode: 'manual'
       XDataSource: ''
       YDataSource: ''
       ZDataSource: ''
```

We can now use the handles to change the properties of the curves. For instance, we may increase the line width of the first curve by;

>> set(h(1),'linewidth',2)

(note: Matlab is not case sensitive when it comes to property names).

The properties can also be modified interactively, as we have seen in Section 3.2.2 p. 44. Using handles enables us to specify the desired look from our scripts and functions. Once we get the script right, we get the right look every time. Furthermore, using handles makes it possible to automate the look in repetition loops.

As an example, we may even modify the data to be plotted in the graph. This enables the creation of *animations*. Example:

```
% animation example
x = linspace(0,1,100);
y = 0*sin(x*pi);
figure(2);
h=plot(x,y);
set(h,'erasemode','xor') % Redraw only curve, not full figure
axis([0 1 -1 1])
for i = 1:1000
```

```
pause(0.01); % Wait 0.01 seconds to slow things down
y = sin(i/100*pi)*sin(x*pi);
set(h,'ydata',y) % Enter new y-data
drawnow % Update graph
```

end

For more information on graphics and graphics properties, see the interactive help. (Matlab help or Full product help, from the help menu.)

Exercise 12.1 Make a plot of $\sin x$ and $\cos x$ on $[0, \pi]$.

1. Write a script that changes the line style of the first curve to '-.' and the second to ':', and the colors of both to black ('k'). You can get handles to all line objects of the current axes by the command:

h = findobj(gca,'Type','line')

2. Repeat the exercise, this time using the Figure Properties Editor. This is accessible from the Edit menu on the figure window.

This kind of modification may be useful if you want to copy your paper using black and white copiers. The Matlab Help System contains much more information on how to tailor your graphs to suit your needs.

Problems

In the following problems, use the debugger and Matlab's help facilities to sort out the troubles you are sure to encounter:

Problem 13.1 An integer k is called a proper divisor of n if 0 < k < n, and if the fraction n/k is an integer. Write a function **propdiv(n)** that finds all proper divisors of the positive integer n. A simple algorithm for this is to test every integer $k \le n/2$ to find those that give no remainder when n is divided by k. Use the standard Matlab function **rem** to do this.

Problem 13.2 Write a function perfect(n) that finds the first perfect number > n. A perfect number is the sum of all its proper divisors. The first three perfect numbers are 1, 6 (= 1 + 2 + 3), and 28 (= 1 + 2 + 4 + 7 + 14).

The following problems relate to graphics in Matlab.

Problem 13.3 Let t = linspace(0,2*pi). Let $x1 = \sin(t)$, $y1 = \cos(t)$, $y2 = \sin(2*t)$, $y3 = \cos(3*t)$, and $y4 = \sin(4*t)$. Create a figure containing a 2 by 2 subplot array. In subplot 1, plot y1 vs. x. In subplot 2, plot y2 vs. x, and so on. Use the interactive plotting tools to change line colors, add markers, and stretch or move the subplot axes. You may also change background colors and make other modifications as you like.

Problem 13.4 Use the generate m-file menu choice in the figure's file menu to save the figure set-ups as an m-file function, which you may call createfigure.m. Use createfigure to make a new figure identical to the first, only now the x-axis data for each figure is x2 = cos(t).

Problem 13.5 Modify createfigure to add xlabels and ylabels to each subplot. The ylabel for subplot 1 may be " $\cos(t)$ " and so on.

Problem 13.6 In this problem, do *not* use the interactive plotting tools. Make a new figure, this time showing y1 through y4 vs. x1 in the same axes. Let the plot command return handles to each graph, and use the handles to add markers and modify colors and line types of each graph.

To solve the next problem, it is necessary to study Appendix D.

Problem 13.7 Write the Matlab commands: x=0:0.25:2; y=0:0.5:5; [X,Y] = meshgrid(x,y). Examine X and Y to see the effect of the meshgrid function. Let $Z = X.^2 + sin(Y)$. Plot Z as a function of X and Y using the following commands: surf, mesh, contour, contourf. Write labels on the axes. Rotate and move the graphs using the various interactive tools, including those on the camera toolbar.

Part III

Becoming a Matlab Master

The once and future Matlab Master

14.1 Revisiting Part II

Numerical methods with Matlab

• Repetition of functions from Part I, and some new functions.

Writing your own functions

- Introduction through example: distance to horizon
 - Basic use of functions
 - Returning more than one result
- Workspaces and variables: local value of variables within functions
- Relational and logical expressions
- Program flow control
- Loops vs. matrix operations
- Errors and the debugger

Data structures in Matlab

- Structures
- Cell arrays

Handle graphics

- Figures and handles
- Modification of figures through handles and setting properties
- Animation

14.2 Overview of new material

- Functions and function handles
- Calling functions through their handles
- Anonymous functions
- Function functions

- Writing your own Newton solver
- Generalizing the newton solver
- ODE solvers
 - Writing your own Euler solver
 - Generalizing the Euler solver
 - Built-in Matlab solvers

Functions and function handles

15.1 Function handles

A *function handle* is a Matlab *value* which can be used for calling a function *indirectly*. Basically, the function handle is an address to the function.

A function handle can be created as follows:

>> mysin = @sin mysin = @sin

Here, sin is a function name — the built-in Matlab function to compute the sine of the argument. By preceding the function name by the handle indicator **@**, **@sin** is the handle for the sine function, and points to the built-in sine function. Finally: with the statement mysin = **@sin**, we *declare* variable mysin to be the function handle **@sin**.

From Matlab 7 of, we can use the variable name mysin containing the function handle of the sin function, just as if it was the sin function:

```
>> x = linspace(0,2*pi);
>> plot(x, mysin(x))
```

In older versions of Matlab, other and more arcane techniques must be used.¹

The function handle is a mapping to the memory address of the function. We can treat function handles as any variable, e.g. copying the handle:

```
>> myothersin = mysin;
```

Next, we can use myothersin just as we used mysin, and

```
>> plot(x, myothersin(x));
```

will give the same result as plot(x, mysin(x)). Function handles can e.g. be stored in cell arrays:

```
>> mysin = @sin;
>> mycos = @cos;
>> mytrig = {mysin, mycos}
```

¹In older versions of Matlab, we would have to write the plot statement below as plot(x, feval(mysin, x)).

mytrig =

[sin] [cos]

Then we can use mytrig to plot the sine function:

```
>> x = linspace(0,2*pi);
>> plot(x, mytrig{1}(x))
```

A very important feature of the function handle, is that we can transfer the handle as an input argument to a function, and then use the function handle from within the new function — because of this, we talk about function handles as "function functions".

15.2 Function functions

We have already seen examples of many built-in Matlab functions. With some of the built-in Matlab functions, the user has to supply an additional function. As an example, Matlab has built-in functions for performing numerical integration, e.g. the quad function². But the user has to supply the function that he/she wants to integrate. In the simplest case, the arguments of quad are: (i) the function handle for the function we want to integrate, (ii) lower integration boundary, and (iii) upper integration boundary.

Suppose we want to find

$$\int_0^\pi \sin x dx$$

by numeric integration. This can be done as follows:

>> quad(@sin,0,pi)

ans =

2.0000

The point here is that Matlab function quad is a general function that can integrate any function numerically, and we, the user must supply the function handle for the function which we want to integrate — @sin in this case. Because we supply the function handle @sin, Matlab function quad can internally make calls to the specified function sin.

Similarly, we can compute

$$\int_0^{\frac{\pi}{4}} \cos x dx$$

as follows:

```
>> quad(@cos,0,pi/4)
```

ans =

0.7071

Some Matlab functions that require a user specified function are shown in Table 15.1.

We have previously seen how we can define our own functions in m-files, where the function name and the file name coincide. We can then transfer the reference to such functions using the function handle technique, e.g. to the built-in functions in Table 15.1. It is also possible to define functions in a simpler way, without writing specific m-files for them. Such functions are denoted *anonymous functions*.

²Numerical integration is often called *quadrature*.

Function	Comment
quad, quadl	Numerical integration, quadrature
fzero	Find a zero of a function of one variable
fminbnd	Find a minimum of a function of one variable
fminsearch	Find a minimum of a function of several variables
fplot	Plot a function
ode23, ode45	Solve an ODE system
ode15s	Solve an ODE or DAE system

Table 15.1: Some built-in Matlab functions that require a function reference (function handle) as input argument.

15.3 Anonymous functions

Anonymous functions give you a quick way to define functions without writing m-files. In mathematics and in computer algebra systems such as Maple, Mathematica, MuPAD, etc., we can define functions as follows:

$$f: x \to y$$

Here, the f before the colon : is the name of the function, x is the input argument, and y is the mapping of x. As an example, we can define a function named g as

$$q: x \to \sin^2 x.$$

Figure 15.1 illustrates how such a function definition can be carried out in a Computer Algebra System such as MuPAD.

In Matlab³, we can define such a function using the following syntax:

>> g = @(x) (sin(x)).^2;

Here, we notice that we have prepared g(x) for accepting array inputs x. Then we can compute g(x) for selected values of x as follows:

Similarly, we can plot g(x) as follows:

```
>> plot(x,g(x))
```

Let us also consider another example:

sinc: $x \to \frac{\sin x}{x}$.

Using Matlab, we find:

>> sinc = $O(x) \sin(x)./x;$

³Matlab 7 and later versions.

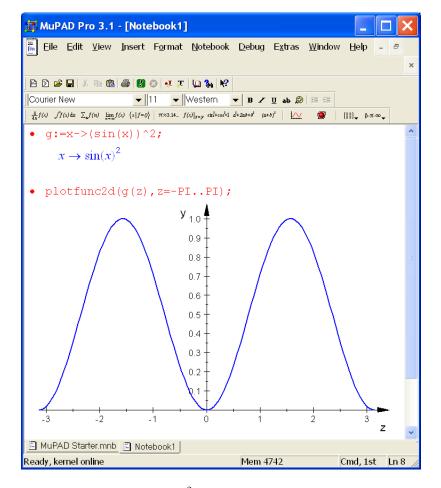


Figure 15.1: Definition of function $g: x \to \sin^2 x$ in Computer Algebra System (CAS) MuPAD. Note in particular that the mapping g is independent of what we choose to name the *argument*.

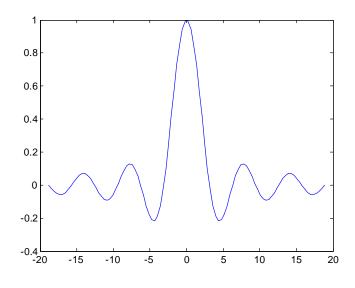


Figure 15.2: Plot of sinc $x = \frac{\sin x}{x}, x \in [-6\pi, 6\pi]$.

Here, we have prepared our sinc function to accept array input argument \mathbf{x} . We can now use the sinc function handle as follows:

>> x = linspace(-6*pi,6*pi);
>> plot(x, sinc(x))

to find the well known plot of the sinc x function, fig. 15.2.

We can also use function handles of anonymous functions as input arguments to the functions in Table 15.1. As an example, we could consider the problem of computing

$$\int_0^{6\pi} \operatorname{sinc} x dx.$$

In the simplest case, the arguments of quad are: (i) a function handle for the function we want to integrate, (ii) lower integration boundary, and (iii) upper integration boundary. We have previously done this for the sin x function, and in principle, we could carry out a similar integration with the anonymous function handle sinc in the interval $x \in [0, 6\pi]$:

```
>> quad(sinc, 0, 6*pi)
Warning: Divide by zero.
> In @(x) sin(x)./x
    In quad at 62
ans =
        1.5180
```

We see that Matlab has a problem since $\sin x/x$ is not handled very well at x = 0. This illustrates a problem with anonymous functions. The general syntax is O(args) expr, where args is a list of comma-separated input arguments, and expr is a one line Matlab expression which depends on the input arguments. In the case of the sinc x function, we would really like to have more than one Matlab expression in the definition, e.g. O(x) if x == 0, 1; else, sin(x)/x; end; this is, however, more than one expression, hence is not a valid expression for the anonymous function.

One more example of anonymous functions: we have in previous chapters searched for the root of $\sin x = 0.5$. We can find this root by combining the idea of anonymous functions with function **fzero** in Table 15.1, specifying 1 as the initial guess of the root x:

As indicated, anonymous functions can be functions of more than one argument:

```
>> sin2 = @(x,y) sin(x)*sin(y);
>> sin2(1,2)
ans =
        0.7651
```

15.4 Function functions: motivating example

15.4.1 Implementation of Newton method

We want solve the equations $\sin x = 0.5$, using Newton's metod. We use the following code and store it as a script in file newton0.m:

```
x0 = 1;
xi = x0;
fi = sin(xi)-0.5;
Ji = cos(xi);
while abs(fi) > 1e-10
    xip1 = xi-fi/Ji;
    disp(['xi = ', num2str(xi)]);
    xi = xip1;
    fi = sin(xi)-0.5;
    Ji = cos(xi);
end
disp(['sin(xi) - 0.5 = ', num2str(sin(xi)-0.5)]);
```

Here is a brief description of how the Newton method works. We want to solve the equations

 $f\left(x\right) = 0,$

where in the example case above,

$$f(x) = \sin x - 0.5.$$

Here are the steps in the Newton method:

Algorithm 1 Newton method:

- 1. Let x^* denote the correct solution such that $f(x^*) = 0$; x^* is currently unknown. We want to find an approximation x_{∞} of x^* through iteration.
- 2. First, we have to guess an initial value x_0 . The closer x_0 is to the correct solution x^* , the faster we will find the our approximate estimate x_∞ of x^* . Set counter i = 0.
- 3. Next, we find a first order Taylor series approximation of f(x) in the point x_i :

$$f(x) \approx f(x_i) + \left. \frac{df(x)}{dx} \right|_{x=x_i} (x-x_i)$$

15.4. FUNCTION FUNCTIONS: MOTIVATING EXAMPLE

4. We now let the next iterate x_{i+1} be defined by $f(x_{i+1}) \approx 0$:

$$0 = f(x_i) + \frac{df(x)}{dx} \Big|_{x=x_i} (x_{i+1} - x_i)$$

$$\downarrow$$

$$x_{i+1} = x_i - \frac{1}{\frac{df(x)}{dx}} f(x_i).$$

5. Hopefully, by iterating infinitely many times, $x_{\infty} \approx x^*$. Since we do not have infinitely much time, we need to terminate the approximation after a finite number of approximation. We thus need tests to see whether we are close to x_{∞} . One possible tests is:

$$|f(x_{i+1})| \le \varepsilon_f,$$

and if this test is satisfied, we set

$$x^* \approx x_{i+1},$$

and terminate the iteration. But other tests are possible, e.g.:

$$|f(x_{i+1}) - f(x_i)| \le \varepsilon_{\Delta f},$$

which will indicate that we do not improve our approach to the correct solution in the ordinate direction. When satisfied, we set

 $x^* \approx x_{i+1},$

and terminate the iteration. Or we could require that

$$|x_{i+1} - x_i| \le \varepsilon_{\Delta x},$$

which will indicate that we do not improve our approach to the correct solution in the abscissa direction. When satisfied, we set

 $x^* \approx x_{i+1},$

and terminate the iteration.

In reality, we may stop the iteration when one of the test above are satisfied, or we may require that each of them are satisfied. In practice, it may also be necessary to introduce a maximal number of iterations i_{max} , in case the Newton method doesn't work. Other refinements may also be necessary.

6. If the iteration is not terminated, we increase the iteration counter by 1: i = i + 1, and return to step 3.

Let us illustrate the method by plotting $f(x) = \sin x - 0.5$ together with the linear approximation

$$f(x) \approx f(1) + \left. \frac{df(x)}{dx} \right|_{x=1} (x-1) = \sin 1 - 0.5 + (\cos 1) (x-1),$$

see fig. 15.3.

Here is the result of running the Matlab code above:

>> newton0
xi = 1
xi = 0.368
xi = 0.51831
xi = 0.52359
sin(xi) - 0.5 = -1.596e-011

We see how the iterates approach the correct solution $x^* = \arcsin(0.5) = 0.523599$.

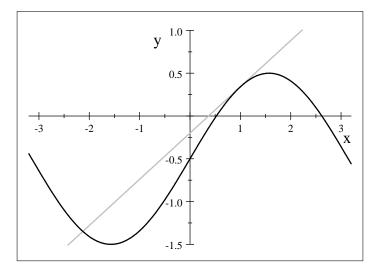


Figure 15.3: Plot of function $f(x) = \sin x - 0.5$ (black), and linear approximation at $x = 1 : f(x) \approx \sin 1 - 0.5 + (x - 1) \cos 1$ (gray). We want to find x : f(x) = 0.

15.4.2 Generalizing the Newton method

Natural we observe that the Newton method is not restricted to the problem $\sin x = 0.5$. Furthermore, we observe that the Newton method is the same for any problem f(x) = 0, and that the only thing that is special is the form of the function f(x) and the form of the Jacobian J(x) = df(x)/dx. It is thus natural to consider separating the Newton algorithm itself from the form of the functions f(x) and J(x).

We could consider the following strategy:

- 1. Define the function f(x) in a Matlab function myfunction, with input argument x, and output argument f(x).
- 2. Define the Jacobian J(x) = df(x)/dx in Matlab function myjacobian, with input argument x and output argument J(x).
- 3. Write a Newton solver function mynewtonsolver, with input argument x_0 , and output argument x^* (or an approximation of x^*).

Here is how our code works:

In order to get this result, we have written the function mynewtonsolver for a general Newton solver:

```
function xsolution = mynewtonsolver(x0)
%
xi = x0;
fi = myfunction(xi);
Ji = myjacobian(xi);
while abs(fi) > 1e-10
    xip1 = xi-fi/Ji;
    xi = xip1;
    fi = myfunction(xi);
    Ji = myjacobian(xi);
```

end xsolution = xi;

In addition, we have written the special function myfunction for defining $f(x) = \sin x - 0.5$:

```
function fx = myfunction(x)
%
fx = sin(x) - 0.5;
```

and the special function myjacobian for defining $J(x) = df(x)/dx = \cos x$.

```
function Jx = myjacobian(x)
%
Jx = cos(x);
```

15.4.3 Some final thoughts

We have now seen how we can separate a general algorithm (the Newton method) from a special case (finding the root of $f(x) = \sin x - 0.5$) by writing a general function for the Newton method, and then implement the specific functions f(x) and J(x) in separate Matlab functions.

Still, our implementation is *not satisfactory*:

- With the current implementation, we can define one and only one problem f(x) = 0 per directory on our computer! This is so since we are only allowed to have a single file with the name myfunction, and a single file named myjacobian.
- Clearly, this is too limiting. We need to be able to define many problems for each directory.
- The solution is to change the code of the Newton solver so that the user can define the file names where f(x) and J(x) are stored.

This modification of our Newton solver is the topic of the next section.

15.5 Function handles as function arguments

To make our Newton solver more general, i.e., such that the user can choose the name of the functions where f(x) and J(x) = df(x)/dx, we need to transfer information about the function names to the mynewtonsolver. The best way to transfer the names, is as input arguments to the Newton solver.

In Matlab, we do not transfer the names of the functions, but rather so-called *function handles*, which are pointers to the memory address of the functions. In order to specify that we mean the function *handles*, we precede the function names by the *function handle indicator* @. Here is a typical call structure to our new Newton solver, which we name mynewtsolver:

```
>> x0 = 1;
>> mynewtsolver(@mysinfunc, @mysinjacobian, x0)
```

Here, we notice:

- 1. The new function (which we have yet to code/write), is named mynewtsolver, and is quite similar to mynewtonsolver.
- 2. Compared to our previous solver mynewtonsolver, the new solver requires two additional input arguments: the function handles to the function where f(x) is defined (in this case function: mysinfunc) and where J(x) = df(x)/dx is defined (in this case function: mysinjacobian). We have to write both of these new functions.

- 3. In addition, we need to specify an initial guess x0, i.e. x_0 .
- 4. Notice that the order of the input arguments is arbitrary we can choose the order as we want. However, when we have chosen the function order, we must stick to it!

Below is an example of how we can run the new code:

```
>> x0 = 1;
>> mynewtsolver(@mysinfunc, @mysinjacobian, x0)
ans =
        0.5236
```

Here is what the new function mynewtsolver looks like:

```
function xsolution = mynewtsolver(myfunc, myjac, x0)
%
% Prepared for Matlab 7
%
xi = x0;
fi = myfunc(xi);
Ji = myjac(xi);
while abs(fi) > 1e-10\quad \quad % Here, we could introduce more stop criteria
    xip1 = xi-fi/Ji;
    xi = xip1;
    fi = myfunc(xi);
    Ji = myjac(xi);
end
xsolution = xi;
```

The functions mysinfunc and mysinjacobian are defined just as before (but with other names):

```
function fx = mysinfunc(x)
%
fx = sin(x) - 0.5;
```

and:

```
function Jx = mysinjacobian(x)
%
Jx = cos(x);
```

What have we gained from this change? The important thing is that we can now use the Newton solver with new function names *without* changing the code within the Newton solver. This is very important: never write functions that the user is supposed to modify!

Suppose we want to use the code to solver another function, e.g.

$$f(x) = \cos x - 0.5$$

This function is displayed in fig. 15.4. The Jacobian of this function is

$$J\left(x\right) = -\sin x,$$

and let us use $x_0 = 0.5$ as an initial guess. We have to write two new functions, where we define f(x) and J(x). Function mycosfunc is:

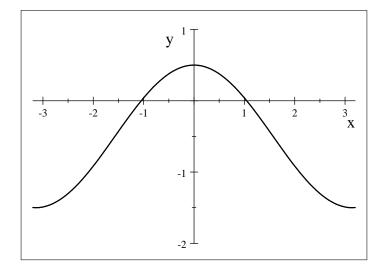


Figure 15.4: Function $f(x) = \cos x - 0.5$.

```
function fx = mycosfunc(x)
%
fx = cos(x) - 0.5;
```

and function mycosjacobian is:

```
function Jx = mycosjacobian(x)
%
Jx = -sin(x);
```

We solve the equation $f(x) = \cos x - 0.5 = 0$ as follows:

This idea of using function handles generalizes to any application that we may want to write. Here, it could be mentioned that there exist code that can automatically find the Jacobian of a vector function; such methods are denoted *automatic differentiation*.

15.6 Exercises

Exercise 15.1 In the previous sections, we have solved $\sin x = 0.5$ ($f(x) = \sin x - 0.5 = 0$) with initial guess $x_0 = 0.5$.

- Use the functions that we have developed with initial values $x_0 = 1.3$, $x_0 = 1.4$, $x_0 = 1.5$. Comment on the results.
- Explain the result with the various initial guesses.

Exercise 15.2 Redo the case $f(x) = \sin x - 0.5 = 0$, with $x_0 = \pi/2$ and with $x_0 = 0.999 \cdot \pi/2$.

• Explain the results.

- Why may it be necessary to put restrictions on the maximal number of iterations, i_{max} ?
- Modify the Newton solver mynewtsolver such that there is a limit to the number of iterations, and let your code write a message to the user if this maximum number is reached.

Exercise 15.3 Use the developed code to find roots for the functions:

• $f(x) = \sqrt{x} - 1$,

•
$$f(x) = x^5 - 1.5$$
,

• $f(x) = \tan x - 0.5$.

Exercise 15.4 For some functions, it may be a hassle to find an analytic expression for the Jacobian J(x) = df(x)/dx. A possible alternative is to develop a function for computing the Jacobian numerically.

• Write a function for computing a numerical approximation for J(x) based on the central difference

$$J(x) \approx \frac{f(x + \Delta x) - f(x - \Delta x)}{\Delta x}$$

(or other approximations of the derivative). The code should automatically choose a relevant value for Δx , e.g. as $\Delta x = x/100$, $\Delta x = 10^{-5}$, etc.

(**Hint**: Make sure that you avoid the possible problem that x may have the value x = 0, e.g. by setting $\Delta x = \max\left(\frac{x}{100}, 10^{-5}\right)$.)

• Modify your Newton solver to automatically detect whether the user has specified a function for the Jacobian. If not, call the numerical approximation.

(Hint: Use the function nargin to find the number of arguments used by the user.)

Exercise 15.5 Write a function grad(f,x) that finds the numerical gradient of a function of several variables. Assume that the function has form f(x) where x is a vector of several variables.

(**Hint**: To find the *i*-th component of the gradient, you must evaluate f at $x + \Delta x$ etc. where element i of Δx differs from zero, while all other elements of Δx are zero. Test grad on the function $\sin(x_1) + \cos(x_2)$.)

Exercise 15.6 Write a Newton solver that can solve the equation f(x) = 0 for the general case, when both x and f(x) are vectors. It can be assumed that the number of unknowns x equals the number of equations f(x) = 0.

(Hint: You may assume that function f returns both the vector of function values, and the Jacobian, i.e. has the Matlab form [f,J] = f(x). Remember that one Newton step in Matlab has the scalar case form [f,J] = f(x); x = x - f/J; In the multivariable case, we need to change the scalar case line x = x - f/J; $(x := x - \frac{f}{J})$ to the multivariable version $x = x - J \setminus f$; $(x := x - J^{-1}f)$. Iterate until the norm of x-J/f is less than 10^{-5} . Make sure that the iteration will stop eventually, even if the iteration does not converge.

Exercise 15.7 Test your Newton solver in the previous exercise, on the function

$$f_1(x) = x_1^2 + x_2^2 - 4$$

$$f_2(x) = \exp(x_1) + x_2 - 1.$$

Exercise 15.8 Write a function jacobian(f,x) that finds the numerical Jacobian matrix of a vector function f as a function of a vector x of variables. Thus: $x \in \mathbb{C}^n$ and $f(x) \in \mathbb{C}^m$.

Exercise 15.9 Modify the Newton solver so that if the user has not specified the Jacobian of the system, then the numerical Jacobian is computed using Jacobian.

Exercise 15.10 Write a simple function myint for finding the integral $I = \int_{a}^{b} f(x) dx$, where the integral is approximated by

$$I \approx \sum_{i=0}^{N-1} f\left(a+i \cdot \frac{b-a}{N}\right) \cdot \Delta x.$$

Test function myint with functions $f(x) = x^2$ and $f(x) = \sin x$, and compare the result with what you get using the built-in Matlab function quad.

Chapter 16

Simulation of dynamic systems

16.1 Solution of ODEs using Euler integration

We consider the damped mass-spring system in fig. 16.1. A model for the system in fig. 16.1 is:

$$\frac{dx}{dt} = v \tag{16.1}$$

$$\frac{dv}{dt} = -\frac{k}{m}(x - x_0) - \frac{\mu}{m}v - g + \frac{F}{m}.$$
(16.2)

The parameters, initial conditions, and input signals in Table 16.1 are to be used for this system.

Notice that we have not stated any reference position for x! It is natural to assume that the unloaded spring hangs downwards. Since Table 16.1 specifies the unloaded spring position to be $x_0 = 1 \text{ m}$, this must mean that the spring hangs from a position > 1 m (the x-axis points upwards!). Furthermore, since it is specified that $x|_{t=0} = 1.5 \text{ m}$, this means that the spring at t_0 is compressed relative to its unloaded length.

In addition to the relationships above, we can express the kinetic energy K, the potential energy P, and the total energy E as follows:

$$K = \frac{1}{2}mv^{2}$$

$$P = mgx + \frac{k}{2}(x - x_{0})^{2}$$

$$E = K + P.$$

$$\frac{dy}{dt} \approx \frac{y(t + h) - y(t)}{h}.$$

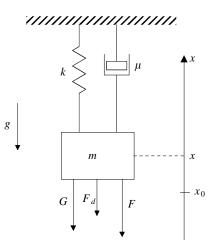


Figure 16.1: Sketch of damped mass-spring system.

Numerical values				
g = 9.81 m/s ²		gravitational acceleration		
k = 2	J/m^2	spring constant		
$\mu = 0.5$	N/(ms)	damping constant		
$x_0 = 1$	m	unloaded spring position		
m = 5	kg	mass		
$x _{t=0} = 1.5$	m	initial value for x		
$v _{t=0} = 0$	m/s	initial value for v		
F = 0	Ν	external force		

Table 16.1: Numerical values for damped mass-spring system.

We will now see how we can use the Explicit Euler Method to solve the ODE for x and v. The Explicit Euler Method works as follows:

Algorithm 2 In the Explicit Euler Method, we consider equations of form

$$\frac{dy}{dt} = f\left(y, u, t\right)$$

- 1. Assume that $y(t_0)$ is known, together with $u(t_0), u(t_0+h), u(t_0+2h), \ldots$
- 2. Use the approximation

$$\frac{dy}{dt} \approx \frac{y\left(t_i + h\right) - y\left(t_i\right)}{h},$$

and rewrite the ODE to the following approximation:

$$y(t_i + h) = y(t_i) + h \cdot f(y(t_i), u(t_i), t_i).$$

3. With $y(t_0)$ given, together with t_0 and $u(t_0)$, compute $y(t_0 + h)$. Then iterate to compute $y(t_i)$.

The following script file eulersim.m will solve these equations, and compute x(t), v(t), as well as K(t), P(t), and E(t).

- 1. Run the file by typing >>eulersim.
- 2. Plot the result using commands >> plot(T, Y(:, 1)), >> plot(T, Y(:, 2)) and >> plot(T, E).
- 3. Explain the figures.

Here is an example of what the file eulersim.m can look like:

```
% Script for simulating a damped mass-spring system
% using the Explicit Euler Method
% Defining system parameters
data;
% Specifying simulation time Tfin, step length h,
% and computing the necessary number of steps N
Tfin = 150;
h = 0.1;
N = ceil(Tfin/h); % Rounds the necessary steps up to the nearest integer
% Allocating space for the time vector T, as well as the state matrix Y
% and the energy function E
T = zeros(N,1);
Y = zeros(N,2);
E = zeros(N,2);
```

```
% Initial values for the states
Y(1,:) = [1.5, 0];
% Running the simulation
for I = 1: (N-1),
   x = Y(I,1);
   v = Y(I,2);
   fx = v;
   fv = -my*v/m - k*(x-x0)/m - g + F/m;
   Y(I+1,1) = x + h*fx;
   Y(I+1,2) = v + h*fv;
   T(I+1) = T(I) + h;
end
%Post processing to find the time response of energy E
K = 0.5*m*Y(:,2).*Y(:,2);
P = m*g*Y(:,1) + 0.5*k*(Y(:,1)-x0).^2;
E = K + P;
```

Note the command data. This command runs a script file (data.m) which holds the numerical values for parameters, etc.:

g = 9.81;	\qquad \qquad % m/s^2, gravitational acceleration
k = 2;	% J/m ² , spring constant
my = 0.5;	\qquad \qquad % N/m/s, damping constant
x0 = 1;	% m, unloaded spring position
m = 5;	% kg, masse hanging in the spring
F = 0;	% N, external force on mass

Also, notice the following post processing step at the end of script file eulersim.m:

Y(:,2).*Y(:,2) (Y(:,1)-x0).^2

Here we have used array processing to compute the energy.

16.2 Generalizing our Euler solver

The script eulersim.m is unfortunate in the sense that we have mixed a general algorithm (the Euler method) with a specific problem (the damped spring-mass system). This unfortunate formulation is similar to a previous case, where we mixed a general method (the Newton method) with a specific problem (solving $\sin x = 0.5$).

Just as in the case of developing a general Newton solver, we want to develop a general Euler solver where we separate the general method from the specific problem. Here is what we need to do:

- Separate the Euler solver from the specific model (the ODE), by specifying the ODE in separate functions.
- Allow for the user to specify the function name where the ODE is defined, with a free choice of the name of this function.

The following listing of file odexeu.m gives a relatively general ODE solver¹ based on the Explicit Euler Method:

¹ODE = ordinary differential equation. In fact, the ODE solvers require that the system is in state space form, i.e. ODEs of form dy/dt = f(t, y) where y may be a vector, $y \in \mathbb{R}^n$.

```
function [T,Y] = odexeu(filehandle,Tspan,y0,h);
%
% Function call: [T,Y] = odexeu(filename,Tspan,y0,h)
%
% The function utilizes the Explicit Euler Method
% to solve the ordinary differential equation (ODE) dy/dt = f(t,y)
% where y may be a vector, and f(.,.) denotes the
% vector field of the ODE
%
% Interpretation of variables:
\% filehandle: the function handle of the file where the vector field is defined
\% Tspan: a vector [t0,tf] specifying the initial and final time of the simulation
% y0: the vector of initial values, y(t0)
% h: the step length of the simulation
% T: the resulting vector of time instances where y is computed
\% Y: a matrix of the time response of the states, one response fore each column of Y
%
% As an example, the time response of state 1 can be plotted using
% command: >>plot(T,Y(:,1))
% Prepared for Matlab 7
% Computing of the number of time steps in the Euler method:
t0 = Tspan(1);
tf = Tspan(2);
N = ceil(tf/h); \quad \guad \ \rightarrow the iteration number upwards to nearest integer
% Allocates space to store the results
n = length(y0);
T = zeros(N, 1);
Y = zeros(N,n);
% Inserts initial values in T and Y
T(1) = t0;
Y(1,:) = y0';
% Computes y(t) in a for loop:
for I=1:(N-1);
    y = Y(I,:)';
    t = T(I);
    fy = filehandle(t,y);
    Y(I+1,:) = (y + h*fy)';
    T(I+1) = T(I) + h;
end
```

The following script file (e.g. named dampms.m — damped mass spring) simulates the damped massspring system.

% Script for simulating damped mass-spring system

```
% Necessary data
Tfin = 150;
h = 0.1;
y0 = [1.5;0];
% Simulating system
[T,Y] = odexeu(@dampedspring,[0,Tfin],y0,h);
% Computes the energy
data;
K = 0.5*m*Y(:,2).*Y(:,2);
```

P = m*g*Y(:,1) + 0.5*k*(Y(:,1)-x0).^2; E = K + P;

The data for the system is given in the file data.m, see p. 143. Also, we define the "vector field" for the differential equation in the function dampedspring.m:

```
function fy = dampedspring(t,y)
%
% Defining data
data;
% Defining vector field
x = y(1);
v = y(2);
fx = v;
fv = - my*v/m - k*(x-x0)/m - g + F/m;
% Setting up vector field
fy = [fx;fv];
```

In order to run the program, issue the Matlab command >>dampms. The result is plotted using the command >>plot(T,E), etc.

Exercise 16.1 With h = 0.1 s, simulate the system in the interval $t \in [0, 150]$ s (as indicated in the script file dampms.m. Plot the time response of x (Y(:,1)) and v (Y(:,2)) as functions of time (T).

Exercise 16.2 Plot the time response of the total energy E (E) as a function of the time. From the relationship $dE/dt = -\mu v^2 < 0$ (assuming F = 0), it follows that E must decrease monotonously with the time t. Why is this theoretical fact not supported by the plot of E(t)?

Exercise 16.3 Experiment with various values on the step length h, and see if this changes the time response of E. What is the largest step length you can use, and still get a relatively good result from the simulations?

(**Hint**: Even if E(t) oscillates, we will get a relatively good picture of E by averaging the energy over a few oscillation periods. For what value of h will the energy E(t) "take off"?)

Exercise 16.4 Experiment with other values on the parameters (g, k, μ, x_0, m) and the initial values $(x|_{t=0} = 1.5, v|_{t=0} = 0)$. Will the changes in the parameter values or the initial values affect the requirement on step length h such that the simulation gives a good approximation of E(t)?

Exercise 16.5 Study the undamped case $\mu = 0$ when F = 0. In this case, the energy E should be constant according to theory, dE/dt = 0. Simulate the system with $\mu = 0$ and step length h = 0.1 (it suffices to simulate the system for 50 s, i.e. set Tfin=50). What happens with E? Is the theoretically correct result that dE/dt = 0, satisfied?

Exercise 16.6 In the previous exercise, we simulated the undamped case ($\mu = 0$), and in this case, E should theoretically be constant. With h = 0.1, it turns out that the the energy increases, which implies that the approximation we introduced is imperfect. One possible way to improve the simulation is to use a better approximation for dy/dt. Alternatively, we can attempt to rectify the error in the approximation by introducing an artificial damping, i.e. increasing μ until the energy on the average is constant.

What value μ must be used such that the energy E averaged over some oscillation periods becomes constant?

ODE solver	Comment
ode45	Fourth order (explicit) Runge-Kutta method.
ode23	Second order (explicit) Runge-Kutta method.
ode113	Variable order Adams-Bashforth-Moulton method.
ode15s	Variable-order multistep solver (1st - 5th order).
ode23s	Second order modified Rosenbrock single step method.
ode23t	Second order trapezoidal method.
ode23tb	Second order implicit Runge-Kutta method.

Table 16.2: Table Caption

16.3 ODE solvers in Matlab

Matlab comes with a number of ODE solvers as standard. All of them are named odeXXX, where the addition XXX specifies what type of method is used in the particular solver. This should be compared to our solver odexeu from the previous section, where XXX is equal to xeu denoting eXplicit EUler method. A list of available solvers is shown in Table 16.2.

As a general comment, the methods are more accurate the higher their order. However, higher order methods require more computing time to find the solution. If a solution reaches a constraint (e.g. if $y \leq y_{\text{max}}$, such as when a vessel becomes full), then the accuracy is lost when $y = y_{\text{max}}$.

Most methods are single step methods, meaning that for an ODE of type

$$\frac{dy}{dt}=f\left(t,y\right) ,$$

 y_i is computed from the knowledge of y_{i-1} . Two methods in Table 16.2 are multistep solvers (ode113 and ode15s), meaning that y_i is computed from the knowledge of y_{i-1} , y_{i-2} , etc. Multistep solvers are usually more efficient than single-step solvers, but this is not true if the solver needs to be reset often — as an example, at start-up, only $y(t_0)$ is known, and multistep solvers need to use a single-step method in the beginning.

Finally, the first three solvers in Table 16.2 are non-stiff solvers, while the four following methods are stiff solvers. As a general rule, stiff problems have modes with very different time constants. For such problems, non-stiff solvers are very slow at finding the solution, while stiff solvers may find the solution rapidly. As an example, for some stiff problems, solver ode15s may be several hundred times faster than ode45.

We have seen that we call our odexeu solver as follows:

[T,Y] = odexeu(@dampedspring,[0,Tfin],y0,h);

Here, the final input argument is the time increment **h**. In fact, it is very important to find a good choice of the time increment **h**. With too large values of **h**, the approximation may become very poor, even unstable. For too small values of **h**, the solution becomes inefficient: it takes too much time to find the solution.

All of the Matlab solvers in Table 16.2 have algorithms for automatically computing the time increment **h**. As an example: for ode45, it is postulated that the error in the simulation is equal to the difference between the fourth order solution and the fifth order solution, and then the time increment **h** is adjusted to achieve sufficient accuracy in the solution.

Consequently, it is not necessary to inform Matlab about the time increment — in fact, Matlab varies **h** from iteration number to iteration number. *Apart* from this modification, the input arguments to the ODE solvers in Table 16.2 are *identical* to the arguments for **odexeu**. If we want to use solver **ode45** to solve the damped mass-spring system, we need to modify the script file **dampms** into the following:

% Script for simulating damped mass-spring system

```
% Necessary data
Tfin = 150;
h = 0.1;
```

Parameter/variable	Symbol	Nominal value
Volume of reactor	V	1001
Volumetric feed flow	q	$100 \mathrm{l} \mathrm{min}^{-1}$
Feed concentration of A	c_{Af}	$1 \mathrm{mol}\mathrm{l}^{-1}$
Feed temperature	T_f	$350\mathrm{K}$
Volumetric coolant flow	q_c	$100 \mathrm{l} \mathrm{min}^{-1}$
Coolant inlet temperature	T_{cf}	$350\mathrm{K}$
Densities	$ ho, ho_c$	$1000 \mathrm{g}\mathrm{l}^{-1}$
Specific heat capacities	c_v, c_{vc}	$1 \mathrm{cal} \mathrm{g}^{-1} \mathrm{K}^{-1}$
Preexponential factor	k_0	$7.2 \times 10^{10} \mathrm{min}^{-1}$
Exponential factor	E/R	$9.98 \times 10^3 \mathrm{K}$
Reaction order	m	1
Reaction enthalpy	$-\Delta \tilde{H}_r$	$2.0 \times 10^5 \mathrm{cal mol}^{-1}$
Overall heat transfer factor	hA	$7 \times 10^5 \mathrm{cal min^{-1} K^{-1}}$
Discretization time	Δt	$0.01 \min$
Initial value, c_A	$c_A \left(t = 0 \right)$	$8.235 \times 10^{-2} \mathrm{mol} \mathrm{l}^{-1}$
Initial value, T	T(t=0)	441.81 K
Nominal value, q_c	$q_{c}\left(t\right)$	$100 \mathrm{l} \mathrm{min}^{-1}$

Table 16.3: Parameters and nominal operating conditions for CSTR.

```
y0 = [1.5;0];
% Simulating system
tic;
%[T,Y] = odexeu(@dampedspring,[0,Tfin],y0,h);
[T,Y] = ode45(@dampedspring,[0,Tfin],y0);
%[T,Y] = ode15s(@dampedspring,[0,Tfin],y0);
toc
% Computes the energy
data;
K = 0.5*m*Y(:,2).*Y(:,2);
P = m*g*Y(:,1) + 0.5*k*(Y(:,1)-x0).^2;
E = K + P;
```

```
Here, we have also inserted the commands tic/toc to measure the time it takes to solve the ODE using the various methods. For this simple case, we will find that the Euler method by far is the most efficient method. However, the other methods are more accurate, and with the built-in Matlab solvers, we do not have to find/choose the step-length h.
```

Finally, we should mention that we can set some parameters for the ODE solvers via the function odeset.

16.4 Exercises

Exercise 16.7 Consider the following reactor model:

$$\frac{dc_A}{dt} = \frac{q}{V} (c_{Af} - c_A) - k_0 \exp\left(\frac{-E}{RT}\right) c_A^m$$
$$\frac{dT}{dt} = \frac{q}{V} (T_f - T) + \frac{\left(-\Delta \tilde{H}_r\right)}{\rho \hat{c}_{v,S}} \frac{k_0}{m} \exp\left(\frac{-E}{RT}\right) c_A^m + \frac{Q}{\rho V \hat{c}_{v,S}},$$

where supplied heat Q is expressed as:

$$Q = \rho_c q_c \hat{c}_{vc} \left[1 - \exp\left(-\frac{hA}{\rho_c q_c \hat{c}_{vc}}\right) \right] \left(T_{cf} - T\right).$$

Numerical values for parameters and nominal operating conditions are given in Table 16.3.

- Implement the reactor model and use the explicit Euler solver odexeu developed in this chapter.
- How much time does it take to solve the reactor model when you simulate 8 min of operation.
- Use the built-in Matlab solvers ode45 and ode15s, and compare the execution time for simulating 8 min of operation with each other, and with odexeu.
- Change the simulation such that after 1 min of simulated time, the coolant flow:
 - Increases by 10%.
 - Decreases by 10%.

What do you observe?

Part IV Closing

Chapter 17

Conclusions

These lectures notes give a relatively simple introduction to Matlab, and how Matlab can be used to do engineering computations. The notes have been updated to cover Matlab 7.0. Matlab is constantly being improved: the user interface changes, new data structures are introduced, and so on. This means that it is difficult to keep lecture notes for Matlab up to date. Fortunately, there is also a large number of books on Matlab: both introductory books, specialized books (e.g. how to use Matlab do develop numerical algorithms, how to design graphical user interfaces (GUI) for your Matlab code), etc. As with computer books at large, some of the Matlab books are good, and others are less useful.

The first part of the lecture notes covers Arrays: what they are, how to name them, and how to create and manipulate them. Next, basic plotting is treated. A chapter on simple data analysis is provided, and finally the automation of tasks is introduced, with basic string handling, the basic repetition loop (the **for** loop), as well as the concept of Matlab scripts and the Matlab editor.

In the second part, more techniques for program flow control is introduced, and then the important concept of functions is introduced. Followed by this is a discussion of more advanced use of the Matlab editor such as the debugger. Finally some more advanced data structures are introduced (structures, cell arrays), and some basic information about objects are given (handle graphics).

In the third part, functions and function handles are introduced, their usefulness are illustrated through examples (a Newton solver), and finally, there is an overview of how ODE solvers work, as an introduction to the use of built-in Matlab ODE solvers.

In appendices, more details are given. In the first appendix, there is a discussion of vectors and matrices; this should be contrasted with the concept of arrays. In the second appendix, there is a technical discussion of the intricacies of indexing arrays. The third appendix gives an example of how hexadecimal numbers relate to decimal numbers; hexadecimal numbers are the native numbers of computers. Appendix four discusses three dimensional plotting, as well as the housekeeping of plots. In appendix five, a more advanced discussion of string manipulation is given. In appendix six, some numerical methods for vectors and matrices are discussed; Matlab is particular useful for handling linear algebra problems. Appendix seven discusses how anonymous functions can be used to handle extra parameters in function calls. Finally, appendix eight discusses how differential algebraic equations (DAEs) can be solved in Matlab.

To really learn to use Matlab, it is necessary to use it, explore the help browser, but also to study Matlab in more detail. Books such as Hanselman & Littlefield (2005) and Higham & Higham (2000) will provide a good start for becoming a Matlab master.

Part V Appendices

Appendix A

Arrays vs. vectors and matrices*

A.1 Vectors

Definition A.1 A vector is an element of a vector space.

Definition A.2 A vector space is a set \mathcal{V} together with an addition on \mathcal{V} : u + v, $u, v \in \mathcal{V}$, and a scalar multiplication on \mathcal{V} : av, $a \in \mathbb{F}$ and $v \in \mathcal{V}$, such that the following properties hold:

- 1. Commutativity: u + v = v + u for all $u, v \in \mathcal{V}$.
- 2. Associativity: (u + v) + w = u + (v + w) and (ab) v = a (bv) for all $u, v, w \in \mathcal{V}$ and all $a, b \in \mathbb{F}$.
- 3. Additive identity: there exists an element $0 \in \mathcal{V}$ such that v + 0 = v for all $v \in \mathcal{V}$.
- 4. Additive inverse: for every $v \in V$, there exists $w \in V$ such that v + w = 0. Normally, w is denoted -v.
- 5. Multiplicative identity: there exists an element $1 \in \mathbb{F}$ such that 1v = v for all $v \in \mathcal{V}$.
- 6. Distributive properties: a(u+v) = au + av and (a+b)u = au + bu for all $a, b \in \mathbb{F}$ and all $u, v \in \mathcal{V}$.

What does this mean? The vector space is a set, and the operation of addition between vectors, and multiplication of vectors with a scalar must be defined. The operation of addition and multiplication must satisfy the requirements above.

Example A.1 Let us consider the set of *n*-tuples of real numbers, $\mathcal{V} = \mathbb{R}^n$, i.e. the ordered list (v_1, \ldots, v_n) where each $v_i \in \mathbb{R}$. In addition, let us define addition of the elements of the set, v + u, as

$$v+u=(v_1+u_1,\ldots,v_n+u_n)$$

Thus, the operation of addition is well defined. Furthermore, let us define $\mathbb{F} = \mathbb{I} = \{1, 2, 3, ...\}$, i.e. the positive integers, and let us define scalar multiplication as

$$av = (a \cdot v_1, \ldots, a \cdot v_n).$$

Thus, the operation of scalar multiplication is well defined. It is easy to see that Properties 1, 2, 6 are fulfilled. Furthermore, let us define $0 \triangleq (0, \dots, 0)$. Here it is vital to realize that the vector 0 (which is a list of *n* scalar zeros) is different from the scalar 0. Obviously, $0 \in \mathbb{R}^n$, so Property 3 is fulfilled. Furthermore, obviously Property 4 is fulfilled. Finally, scalar $1 \in \mathbb{I} = \mathbb{F}$, and thus Property 5 is fulfilled.

Example A.2 Consider the set of all *n*-th order polynomials $p^n = p_0 + p_1 x + \cdots + p_n x^n$, where normally $p_i \in \mathbb{R}$ or $p_i \in \mathbb{C}$. We define the addition of such polynomials as $p^n + q^n = (p_0 + q_0) + (p_1 + q_1)x + \cdots + (p_n + q_n)x^n$, thus addition is well defined. Furthermore, we define a set \mathbb{F} , e.g. $\mathbb{F} = \mathbb{I}$, and scalar multiplication as $ap^n = (ap_0) + (ap_1)x + \cdots + (ap_n)x^n$, hence scalar multiplication is well defined. It is straightforward to show that Properties 1–6 are fulfilled, hence the polynomials constitute a vector space, and e.g. $p^5 = 1 - x^3 + 2x^5$ is a vector.

The most commonly used vector space, is the space where $\mathcal{V} = \mathbb{R}^n$ (or $\mathcal{V} = \mathbb{C}^n$), and where $\mathbb{F} = \mathbb{R}$ (or $\mathbb{F} = \mathbb{C}$) and where addition and scalar multiplication is defined as in Example A.1. In fact, Matlab is tailor-made for the vector space $\mathcal{V} = \mathbb{C}^n$, $\mathbb{F} = \mathbb{C}$. Addition of vectors is performed with the operator +, and multiplication by scalar by the operator *, and vectors are represented by either row arrays or column arrays:

```
>> vec1 = [1,2,3]
vec1 =
     1
            2
                   З
>> vec2 = [4,5,6]
vec2 =
     4
            5
                   6
>> scalar = 3
scalar =
     З
>> vec1 + vec2
ans =
            7
     5
                   9
>> scalar*vec1
ans =
     3
            6
                   9
```

It should be noted that Matlab is strict in the sense that vector addition only works between two row arrays, or between two column arrays.

Note also that the operator + plays several rôles in Matlab: the plus operator can be used to add scalars. In addition, the plus operator can be used to add a vector and a scalar: in that case, the scalar is interpreted as follows: vector + scalar = vector + scalar*ones(size(scalar)), i.e., the scalar is first expanded into a vector, and then the vector addition is used.

In conclusion: vectors are more than simply arrays. But the most common vectors can be represented as row or column arrays with the addition of well defined vector addition and scalar multiplication. In fact, it can be shown that every vector space of dimension n in some sense is identical to the vector space given by $\mathcal{V} = \mathbb{C}^n$ and $\mathbb{F} = \mathbb{C}$ with addition and multiplication defined as above.

This means that Matlab is well equipped to compute with vectors. The relevant functions for performing such operations, belong to a course in linear algebra.

A.2 Matrices

Often, the word *matrix* is used as a synonym to *array*. However, since there has been developed an algebra of matrices, and furthermore, since matrices can be considered collections of vectors or as operators on vector spaces, we will use matrix to denote something more than arrays — in these notes, and array is simply and organization/storage of data as detailed in Section 2.2.

As an example, we often define matrix addition and the multiplication of matrices with scalar — much in the same way as we do with vector spaces. Also, several definitions of matrix multiplication exist:

Cayley product: Named after Arthur Cayley, who often is considered the founder of matrix algebra, the Cayley (or matrix) product of matrices A and B, $C = A \cdot B$, is defined as:

$$C_{i,j} = \sum_{k} A_{i,k} \cdot B_{k,j}.$$

It follows that the number of columns of A must equal the number of rows of B.

Schur product: The Schur- (or Direct- or Array-) product of matrices A and B, $C = A \times B$, is defined as

$$C_{i,j} = A_{i,j} \cdot B_{i,j}$$

It follows that A and B must have the same number of rows and columns.

Kronecker product: We have already met the Kronecker product of matrices A and B (see p. 162), $C = A \otimes B$, which is defined as

$$C = \begin{pmatrix} A_{1,1}B & A_{1,2}B & \cdots & A_{1,n}B \\ A_{2,1}B & A_{2,2}B & & A_{2,n}B \\ \vdots & & \ddots & \vdots \\ A_{m,1}B & A_{m,2}B & \cdots & A_{m,n}B \end{pmatrix}$$

It follows that there are no restrictions on the sizes of A and B.

Another use/interpretation of matrices is that each column may represent a column vector, or that each row may represent a row vector. With these interpretations, matrices are related to vector spaces.

A third use of matrices is as a mapping/an operator on vectors, typically written as Av where A is the matrix and v the vector. Usually, the standard (Cayley) matrix product is assumed in the notation Av, and in that case, it is necessary that v is considered a column vector/a column matrix.

In Matlab, the operator + can be used to add matrices in a natural way, and furthermore, the operator * works on arrays according to the Cayley matrix product:

```
>> A = rand(2,3);
>> B = rand(3,2);
>> C = ones(2,3);
>> A+C
ans =
    1.9501
              1.6068
                         1.8913
    1.2311
              1.4860
                         1.7621
>> A*B
ans =
    1.1771
              1.5018
    0.7405
               1.0054
```

Matlab also has a large number of built-in functions for treating matrices as collections of vector spaces, e.g. rank, null, orth, etc., or as operators, e.g. det, eig, etc. In addition, there are a large number of built-in functions for dealing with linear equations of form

$$Ax = b.$$

This means that Matlab is well equipped to compute with matrices. The relevant functions for performing such operations, belong to a course in linear algebra.

Appendix B

Manipulation of Arrays and Array Functions*

B.1 Indexing arrays

Suppose array A is given. We often wish to pick out element (i, j) from array A. Let us denote element (i, j) of array A by $A_{i,j}$, which in Matlab is given as A(i,j).

For array A, it can also be useful to be able to pick a subarray which has indices over the rectangle given by row i_1 to row i_2 , and column j_1 to j_2 . This submatrix can be denoted as $A_{i_1:i_2,j_1:j_2}$. In Matlab, this subarray is specified as $A(i1:i_2, j_1:j_2)$:

```
>> rand('state',0)
>> A = rand(3,4)
A =
      0.95013
                    0.48598
                                  0.45647
                                                 0.4447
      0.23114
                     0.8913
                                 0.018504
                                                0.61543
      0.60684
                     0.7621
                                  0.82141
                                                0.79194
>> A(1,3)
ans =
      0.45647
>> A(2:3,2:3)
ans =
       0.8913
                   0.018504
       0.7621
                    0.82141
```

Let us do some careful thinking here! We remember that the notation $i_1 : i_2$ in fact constructs a row array consisting of the following elements: $(i_1, i_1 + 1, \ldots, i_2)$, and similarly for $j_1 : j_2$. Let us therefore try to define the row arrays v_1 and v_2 of integers, and consider the notation A_{v_1,v_2} :

```
>> v1 = [1,3]; v2 = [4,2,2];
>> A(v1,v2)
ans =
```

0.4447 0.4860 0.4860 0.7919 0.7621 0.7621

With v_1 being a row array of length m, and v_2 being a row array of length n, we see that Matlab in fact works as follows:

$$A_{v_1,v_2} = \begin{pmatrix} A_{v_1(1),v_2(1)} & A_{v_1(1),v_2(2)} & \cdots & A_{v_1(1),v_2(n)} \\ A_{v_1(2),v_2(1)} & A_{v_2(2),v_2(2)} & & A_{v_2(2),v_2(n)} \\ \vdots & & \ddots & \\ A_{v_1(m),v_2(1)} & \cdots & & A_{v_1(m),v_2(n)} \end{pmatrix}.$$

What then if we only use one argument, e.g. a row array?

>> A(v1)

ans =

0.95013 0.60684

Now it is slightly more difficult to figure out what is going on. But we see that A assumes the shape of v_1 , and that

 $A(v_1) = (A(v_1(1)) \ A(v_1(2)) \ \cdots \ A(v_1(m)))),$

where one-dimensional addressing has been used. Similarly, if we use one argument where the argument is a two-dimensional array:

>> B = [1,3,2; 4,2, 2]

B = 1 3 2 4 2 2 >> A(B) ans =

0.95013	0.60684	0.23114
0.48598	0.23114	0.23114

Quite often, we are interested in picking out all rows in a column, or every column in a row. We can specify every row with the symbol :, e.g. $A_{:,v_2}$ where row array v_2 suggests which columns should be included, and every column as $A_{v_1,:}$ where row array v_1 indicates which rows of A should be included.

The smallest index in Matlab is always 1. The largest index in rows and columns can be specified with the symbol end in Matlab, e.g.: $A_{1:end,v_2} \triangleq A_{:,v_2}$. This notation is particularly useful if we want to partition an array into two (or more) parts, e.g. with arrays $A_{1:i,:}$ and $A_{i+1:end,:}$:

>> A(:,1:1)

ans =

0.9501 0.2311 0.6068

>> A(:,1+1:end)

ans =

0.4860	0.4565	0.4447
0.8913	0.0185	0.6154
0.7621	0.8214	0.7919

B.2 Building superarrays

B.2.1 Toeplitz and Hankel arrays

The structure of Toeplitz and Hankel arrays are best illustrated by examples. Let's first take a look at Toeplitz arrays:

```
>> c = [1,2,3]; r = [4,5,6,7];
>> toeplitz(c,r)
Warning: First element of input column does not match first element of input row.
         Column wins diagonal conflict.
(Type "warning off MATLAB:toeplitz:DiagonalConflict" to suppress this warning.)
> In C:\MATLAB6p5\toolbox\matlab\elmat\toeplitz.m at line 18
ans =
                        7
     1
           5
                 6
     2
                 5
                        6
           1
     3
           2
                        5
                 1
>> toeplitz(c)
ans =
     1
           2
                 3
     2
           1
                 2
     3
           2
                 1
>> toeplitz(r)
ans =
     4
           5
                 6
                        7
     5
           4
                 5
                        6
     6
           5
                 4
                        5
     7
           6
                 5
                        4
```

We see that Toeplitz arrays have constant elements along the *diagonals*. Notice also that if there is a conflict between the first element of c and the first element of r (these have to be equal!), the first element of c "wins".

Next, we look at Hankel arrays:

```
>> hankel(c,r)
Warning: Last element of input column does not match first element of input row.
        Column wins anti-diagonal conflict.
(Type "warning off MATLAB:hankel:AntiDiagonalConflict" to suppress this warning.)
> In C:\MATLAB6p5\toolbox\matlab\elmat\hankel.m at line 27
```

ans =				
1 2 3	2 3 5	3 5 6		
>> hankel(c	:)			
ans =				
1 2 3	2 3 0	3 0 0		
>> hankel(r)				
ans =				
4 5 6 7	5 6 7 0	6 7 0 0		

5 6 7

7 0 0

We see that Hankel arrays have constant elements along the *anti-diagonals*. Notice that if there is a conflict between the last element of c and the first element of r (these have to be equal!), then the last element of c "wins".

B.2.2 Kronecker product

The Kronecker product of two arrays (or matrices) A and B is denoted $A \otimes B$. If A is an array of size $m \times n$ and B is an array of size $k \times \ell$, then the product is of size $m \cdot k \times n \cdot \ell$, and the result generated as follows:

$$C = A \otimes B = \begin{pmatrix} A_{1,1}B & A_{1,2}B & \cdots & A_{1,n}B \\ A_{2,1}B & \ddots & & \vdots \\ \vdots & & \ddots & \vdots \\ A_{m,1}B & \cdots & \cdots & A_{m,n}B \end{pmatrix}.$$

Here, $A_{i,j}B$ means the value of element $A_{i,j}$ multiplied into every element of array B.

A common problem is to replicate array B in a certain pattern, e.g. we want to insert array B instead of a unit scalar in array A:

$$A = \left(\begin{array}{rrr} 1 & 0 & 1 \\ 1 & 1 & 0 \end{array}\right).$$

By using the Kronecker product which is achieved by kron(A,B), we can produce the concatenated superarray

$$C = A \otimes B = \left(\begin{array}{cc} B & 0 & B \\ B & B & 0 \end{array}\right).$$

The Matlab syntax for this operation is kron(A,B):

```
>> A = [1,0,1; 1,1,0]
```

A =

1 0 1 1 1 0>> B = rand(3,2)

B =					
0.4565 0.0185 0.8214	0.4447 0.6154 0.7919				
>> kron(A,B)					
ans =					
0.4565	0.4447	0	0	0.4565	0.4447
0.0185	0.6154	0	0	0.0185	0.6154
0.8214	0.7919	0	0	0.8214	0.7919
0.4565	0.4447	0.4565	0.4447	0	0
0.0185	0.6154	0.0185	0.6154	0	0
0.8214	0.7919	0.8214	0.7919	0	0

Obviously, we can use the Kronecker product to achieve the same effect as the function repmat(A,m,n) as follows: kron(ones(m,n), A). However, the repmat function is more efficient than using the kron function.

B.2.3 Block diagonal arrays

Function **blkdiag** is used for block diagonal concatenation of arrays:

```
>> A = rand(2,2);
>> B = randn(2,3);
>> C = blkdiag(A,B)
C =
    0.9218
               0.1763
                               0
                                          0
                                                     0
    0.7382
               0.4057
                               0
                                          0
                                                     0
          0
                    0
                         -0.4326
                                     0.1253
                                               -1.1465
         0
                         -1.6656
                    0
                                     0.2877
                                                1.1909
>> D = blkdiag(A,B,-1)
D =
    0.9218
               0.1763
                               0
                                          0
                                                     0
                                                                0
    0.7382
               0.4057
                                          0
                                                                0
                               0
                                                     0
          0
                    0
                         -0.4326
                                     0.1253
                                               -1.1465
          0
                    0
                         -1.6656
                                                1.1909
                                     0.2877
          0
                    0
                               0
                                          0
                                                     0
                                                          -1.0000
```

Note that it is not possible to specify a diagonal off the main diagonal in the case of block diagonal arrays.

0

0

Appendix C

Hexadecimal numbers*

Each byte (1 byte = 8 bits) is composed of 2 hexadecimal numbers, where the first hexadecimal number represents the first four bits, and the second hexadecimal number represents the four last bits of the byte. The 16 hexadecimal digits are $\{0, 1, 2, \ldots, 9, a, b, c, d, e, f\}$. The first half of the first byte of 1/3 is given by hexadecimal number 3_{16} , which in the decimal number system is 3_{10} — the subscript indicates the base of the number, i.e. 16 (hexadecimal) or 10 (decimal). The binary equivalent of this number is $0011_2 = (0 \cdot 2^3 + 0 \cdot 2^2 + 1 \cdot 2^1 + 1 \cdot 2^0)_{10} = 3_{10}$. The next half of the first byte is $f_{16} = 15_{10} = 1111_2$. In total, the first byte of the representation of 1/3 is thus 00111111. Similarly, the first half of the second byte is given by $d_{16} = 13_{10} = 1101_2$. Likewise, $5_{16} = 5_{10} = 0101_2$. Thus, the second byte is 11010101. The remaining 6 bytes are simply 01010101. We can thus write the binary representation of 1/3 as 00111111_11010101_201010101_3 \cdots 010101010_1_01010101_8 — this time, the subscript indicates the byte number. Let us number the bits from 1 to 64 as $b_1b_2\cdots b_{64}$, such that in this number, $b_1 = 0$, $b_2 = 0$, $b_3 = 1$, $b_4 = 1$, etc.

According to the IEEE standard, the representation of floating point numbers is composed of a mantissa part m, and an exponent part e, and the number is given as $m \times 2^e$, see (Gentle 1998). Typically, b_1 signifies the sign of the mantissa (0 means positive number, 1 means negative number), b_2-b_{12} (11 bits) gives the exponent (\tilde{e}), while $b_{13}-b_{64}$ gives the mantissa (52 bits). The mantissa is scaled such that the first bit in the mantissa is always 1. Since it is always 1, it is not included, and we have to add a bit with value 1 at the beginning of the mantissa. Thus, b_{13} is really the second bit in the mantissa. Finally, there is no sign bit in the exponent \tilde{e} . Instead, a fixed number \hat{e} is subtracted from the exponent to get $e = \tilde{e} - \hat{e}$ before we evaluate 2^e .

Let us consider the representation of 1/3. $b_1 = 0$, which means a positive sign. Next, the mantissa is given as

$$m = 1b_{13}b_{14}\cdots b_{64} = 10101_201010101_3\cdots 01010101_701010101_8$$

= $1\cdot 2^{-1} + b_{13}\cdot 2^{-2} + b_{14}\cdot 2^{-3} + \dots + b_{63}\cdot 2^{-52} + b_{64}\cdot 2^{-53}$
= $1\cdot 2^{-1} + 0\cdot 2^{-2} + 1\cdot 2^{-3} + 0\cdot 2^{-4} + 1\cdot 2^{-5}$
+ $\sum_{i=3}^{8} \left(0\cdot 2^{-6-(i-3)\cdot 8} + 1\cdot 2^{-7-(i-3)\cdot 8} + 0\cdot 2^{-8-(i-3)\cdot 8} + 1\cdot 2^{-9-(i-3)\cdot 8} + 0\cdot 2^{-10-(i-3)\cdot 8} + 1\cdot 2^{-11-(i-3)\cdot 8} + 0\cdot 2^{-12-(i-3)\cdot 8} + 1\cdot 2^{-13-(i-3)\cdot 8} \right)$

 $= 0.666\,666\,666\,666\,666\,629\,66.$

The exponent is given as

$$\tilde{e} = b_2 b_3 \cdots b_{12} = 0111111_1 1101$$

= $b_{12} \cdot 2^0 + b_{11} \cdot 2^1 + b_{10} \cdot 2^2 + \cdots + b_3 \cdot 2^9 + b_2 \cdot 2^{10}$
= $1 \cdot 2^0 + 0 \cdot 2^1 + 1 \cdot 2^2 + \cdots + 1 \cdot 2^9 + 0 \cdot 2^{10}$
= $1 \cdot 2^0 + 0 \cdot 2^1 + \sum_{i=2}^9 1 \cdot 2^i + 0 \cdot 2^{10}$
= $1021.$

To handle the sign of the mantissa, approximately half of the maximal exponent number is subtracted from the exponent before evaluating 2^e .: the maximal number is $\sum_{i=1}^{11} 2^{i-1} = 2047$, and half of this

is 2047/2 = 1023.5. In practice, the number $\hat{e} = 1022$ is subtracted. Thus, the real exponent is e = 1021 - 1022 = -1. The number represented by the 8 bytes is thus:

 $0.666\,666\,666\,666\,666\,629\,66\times 2^{-1} = 0.333\,333\,333\,333\,333\,333\,314\,83.$

We see that at the 16 first digits are correct.

Appendix D

Three Dimensional Plots and Plot Housekeeping*

D.1 Three dimensional plots

Let us consider the function

$$f(x,y) = \sin x \cos y \exp\left(-\frac{1}{10}(x^2 + y^2)\right),$$

which is displayed in fig. D.1. How can we plot this function using Matlab? We wish to plot the function in the region $(x, y) \in [-2\pi, 2\pi] \times [-2\pi, 2\pi]$. First we need to define the variation in x and y:

>> x = linspace(-2*pi,2*pi);
>> y = linspace(-2*pi,2*pi);

Remember that these variables (x, y) are row vectors.

The simplest way to compute the functional value of z = f(x, y) is by using array operations. We need to compute z for all possible combinations of x and y, i.e. for $[-2\pi, 2\pi] \times [-2\pi, 2\pi]$. We do this by generating arrays X and Y as follows:

>> [X,Y] = meshgrid(x,y);

Both X and Y are arrays of size $n_x \times n_y$, where $n_x = \text{length}(x)$ and $n_y = \text{length}(y)$. Array X is composed of n_y columns with copies of row array x, while array Y is composed of n_x rows with copies of row array y. The usefulness of this construction, is that if we take an arbitrary element X(i,j) and Y(i,j), then these elements are the ones we need for computing Z(i,j):

>> Z = sin(X).*cos(Y).*exp(-(X.^2 + Y.^2)/10);

Now, we have computed z for a selected number of values of x and y, and the result is found in array Z. There are a number of functions for plotting z as a function of x and y, depending on what we want to display. The most basic function is the **mesh** function:

>> mesh(X,Y,Z)

The result of this command is displayed in fig. D.2. Note that the default length of arrays generated using the **linspace** function, is 100. Thus, the graph in fig. D.2 is generated from $100 \cdot 100 = 10^4$ data points, and the resulting picture file behind fig. D.2 is relatively large. Figure D.3 illustrates how we can interactively rotate the graph, in order to look for a more interesting viewpoint.

Sometimes, we want to see contour lines for the 3D plot: these are like iso-curves in a map:

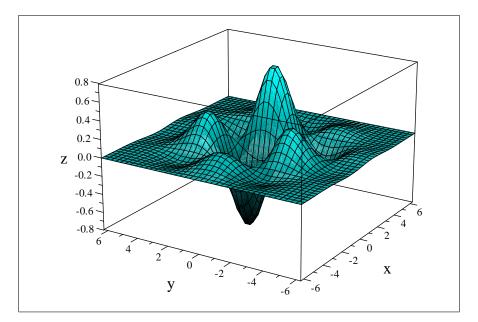


Figure D.1: Graph of function $\sin x \cos y \exp\left(-\frac{1}{10}(x^2+y^2)\right)$ generated using a CAS system (Computer Algebra System, here: MuPAD from within the word processor Scientific WorkPlace).

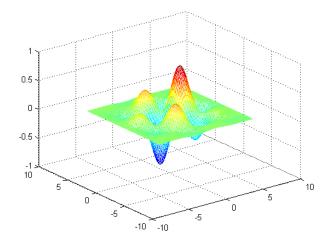


Figure D.2: Graph of function $f(x, y) = \sin x \cos x \exp\left(-\frac{1}{10}\left(x^2 + y^2\right)\right)$ generated in Matlab.

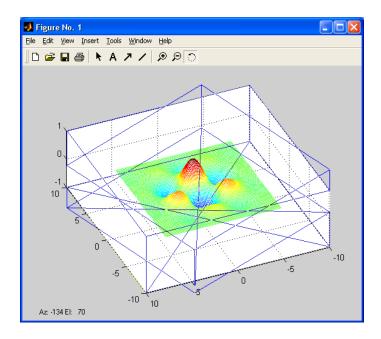


Figure D.3: In order to rotate 3D plots: click the rotation icon on the toolbar, click-and-hold in the plot, and "rotate" using the mouse.

>> meshc(X,Y,Z)

Note the small, but important difference between functions mesh and meshc. The resulting plot is shown in fig. D.4.

If we only want to see the contour lines, this is achieved using function contour:

>> contour(X,Y,Z,20)

draws contour lines at 20 levels, see fig. D.5.

Matlab function **surf** works similarly to function **mesh**.

Exercise D.1 Check out the help files for the following Matlab 3D functions: plot3, contour, contourf, contour3, mesh, meshc, meshz, surf, surfc, waterfall, bar3, bar3h, pie, fill3, comet3, scatter3, stem3.

It should be noted, that in many cases, it is advantageous to plot 3D results in two dimensions, e.g.:

>> plot(x,Z)

The result of this command is shown in fig. D.6, and should be compared to fig. D.2. Perhaps the current example is not the best example, but it is often difficult to understand 3D plots.

D.2 Housekeeping

Sometimes, it is necessary to clear the content of a figure. With the figure window selected, issue the command clf: the figure window is still open, but it is greyed out. In order to close a chosen figure window, issue command close.

It is possible to open more than one figure window at the same time. This is done by issuing the **figure** command. When this command is issued, the figure window is opened, and a so-called

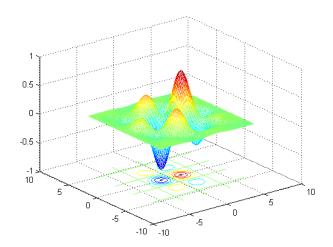


Figure D.4: Graph of function $f(x, y) = \sin x \cos x \exp\left(-\frac{1}{10}(x^2 + y^2)\right)$ generated in Matlab, with added contour lines.

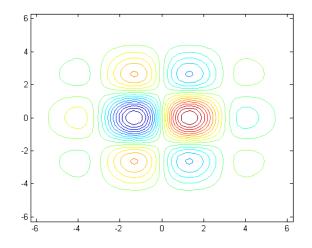


Figure D.5: Contour lines of function $f(x, y) = \sin x \cos x \exp\left(-\frac{1}{10}\left(x^2 + y^2\right)\right)$.

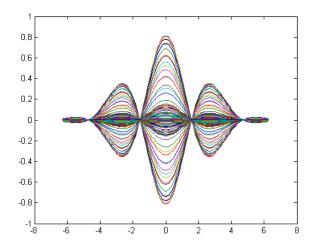


Figure D.6: Graph of function $f(x, y) = \sin x \cos x \exp\left(-\frac{1}{10}(x^2 + y^2)\right)$, mapped on the x-z plane.

handle for the figure is returned. This can later on be used to address the figure using the command figure(<handleno.>), which will make the figure window with the relevant handle the selected window. The handle for the figure can also be used to close a figure (close(<handleno.>) — clf can only clear the selected/current window. The following Matlab session illustrates the use of these commands:

>> f1 = figure; >> f2 = figure; >> f3 = figure; >> figure(f1); >> plot(x,y); >> figure(f3); >> plot(x,exp(-x)) >> figure(f2); >> plot(x,cos(x)); >> figure(f3); >> clf; >> close(f3) >> close(f2)

It is also possible to enforce a certain number on a figure window by using the command figure(N) where N is an integer no. This number can then also be used in place of the handle no.

It is useful to be able to save figures from the command line. This can be done using function saveas with the following arguments: saveas(H,'FILENAME','FORMAT'). Here, H is the handle of the figure, FILENAME is the file name with or without extension, and FORMAT is the file format. If the chosen format is fig, then it is also possible to open the figure for more manipulation in Matlab. For figures that are more or less finished, possible formats are bmp, jpg, etc.

The syntax for opening a figure, is **open** <**filename**>. The following example illustrates how the **saveas** function and the **open** commands may be used:

```
>> saveas(f1,'figure1','fig')
>> close(f1)
>> open figure1.fig
>> saveas(f1,'figure1','bmp')
>> saveas(f1,'figure1','jpg')
>> dir fig*
```

figure1.bmp figure1.fig figure1.jpg

Appendix E

Automation and String Operations*

E.1 Motivating example: generating sequence of plots

Based on the data in fig. 4.8 p. 68, suppose we want to produce a number of plots of the experimental data, and how well a polynomial model of order i fits the data when using the representation T(p), e.g. as in fig. 4.11 p. 71. Let us suppose that we also want to save the figures to files which are named according to its content. The following Matlab session would do the job:

```
Import Wizard created variables in the current workspace.
>> T = data(3:end,1);
>> p = data(3:end,2);
>> pp = linspace(min(p),max(p));
>> fig1 = figure;
>> \% ======= Ready to generate fig. for first order fit =======
>> plot(p,T,'kx')
>> xlabel('p [atm]')
>> ylabel('T [C^{\circ}]')
>> title('Temperature vs. pressure for saturated steam')
>> Tpar = polyfit(p,T,1);
>> hold on;
>> plot(pp,polyval(Tpar,pp),'k-')
>> saveas(fig1,'satvappfit1')
>> clf
>> % ====== Ready to generate fig. for second order fit =======
>> plot(p,T,'kx')
>> xlabel('p [atm]')
>> ylabel('T [C^{\circ}]')
>> title('Temperature vs. pressure for saturated steam')
>> Tpar = polyfit(p,T,2);
>> hold on;
>> plot(pp,polyval(Tpar,pp),'k-')
>> saveas(fig1,'satvappfit2')
>> dir satvappfit*
satvappfit1.fig satvappfit2.fig
>> % ====== Ready to generate fig. for third order fit =======
>> % ... etc. ...
```

Clearly, it is a major job to generate figures for every model order up to, say, 60: we have 61 data points, which means that a polynomial model of order 60 implies interpolation. There must be a simpler way, i.e. it must be possible to automate the generation of these plots.

In order to understand how to automate the generation of these plots, we need to discuss strings and string operations, as well as automatic repetition of operations.

	Table E.1: Some basic functions for operating on strings.	
Function	Description	
s = 'abc'	Creates a string \mathbf{s} containing the character sequence \mathbf{abc} .	
length(s)	The number of characters in string \mathbf{s} .	
blanks(n)	Creates a string of n blank elements (spaces).	
s(i)	If s is a string and i is a positive integer, $s(i)$ is the string ' $s(i)$ '.	
char(i)	If i is a an integer representing the ASCII value of a character, char(i) is	
	the resulting character string.	
[s1,s2]	Concatenates strings s1 and s2.	
<pre>strcat(s1,s1)</pre>	Concatenates strings s1 and s2, but neglects blank strings.	
deblank(s)	Strips trailing blanks from the end of strings.	
class(obj)	Responds with the type of the obj.	
<pre>isa(obj,'char')</pre>	If obj is a string of characters, the result is boolean TRUE; otherwise	
	FALSE.	
ischar(obj)	If obj is a string of characters, the result is boolean TRUE; otherwise	
	FALSE.	
<pre>isletter(s)</pre>	Creates a row array of length length(s), with value 1 for corresponding	
	letters in \mathbf{s} , and 0 otherwise.	
isspace(s)	Creates a row array of length length(s), with value 1 for corresponding	
	spaces in \mathbf{s} , and 0 otherwise.	
isequal(s1,s2)	If strings \$1 and \$2 are equivalent, the result is boolean TRUE; otherwise	
	FALSE.	
str2num(s)	If s is a string of characters representing a floating point number,	
	str2num(s) converts the string to the number.	
num2str(x)	If x is a floating point number, num2str(x) is the character string for the	
	number.	

Table E 1. Sc mo basic functions for operating on strip

E.2 Strings and string operations

It is useful to have some knowledge of strings in Matlab, since these are important for some types of task automation. A string is a sequence of characters. When defining the string, the sequence of characters must be surrounded by an apostrophe ':

<pre>>> clear >> mystr1 = >> mystr2 = >> whos</pre>			
Name	Size	Bytes	Class
mystr1 mystr2	1x1 1x3		char array char array

Grand total is 4 elements using 8 bytes

We see that each character in a string takes up 2 bytes of memory.

Some basic functions for operating on strings are shown in Table E.1. Let us see how these functions work with a simple string:

```
>> mystring = 'John Donne';
>> length(mystring)
ans =
    10
>> find(mystring == ' ')
ans =
     5
```

```
>> mystr1 = mystring(1:5-1)
mystr1 =
John
>> mystr2 = mystring(5+1:end)
mystr2 =
Donne
>> [mystr1, ' ', mystr2]
ans =
John Donne
>> strcat(mystr1, ' ', mystr2)
ans =
JohnDonne
>> [mystr2, ', ', mystr1]
ans =
Donne, John
>> mystring(1)
ans =
J
>> mystring(2)
ans =
0
>> mystring(1:4)
ans =
John
>> mystring(end:-1:1)
ans =
ennoD nhoJ
```

In particular, note how we can change the sequence of characters in the string by using the command mystring(end:-1:1).

Let us also see how strings and numbers can be mixed and matched:

```
>> mynum = pi
mynum =
       3.1416
>> mynumstr = num2str(mynum)
mynumstr =
3.1416
>> class(mynum)
ans =
double
>> class(mynumstr)
ans =
char
>> isa(mynumstr,'char')
ans =
     1
>> isa(mynum,'char')
ans =
     0
>> ischar(mynumstr)
ans =
     1
>> isequal(mynum,mynumstr)
ans =
     0
>> isequal(mynumstr,'3.1416')
```

Tał	ble E.2: More advanced functions for operating on strings.
Function	Description
lower(s)	Converts any upper case character in string \mathbf{s} to lower case.
upper(s)	Converts any lower case character in string \mathbf{s} to upper case.
<pre>strjust(s,type)</pre>	Returns a justified string of string s according to type = 'right', 'left', or 'center'.
<pre>strrep(s,s1,s2)</pre>	Replaces any occurrence of string $s1$ in string s , with $s2$.
findstr(s1,s2)	Returns the starting indices of any occurrences of the shorter of the two strings s1, s2 in the longer.
<pre>strcmp(s1,s2)</pre>	Returns 1 if strings s1 and s2 are the same, and 0 otherwise.
<pre>strcmpi(s1,s2)</pre>	Returns 1 if strings s1 and s2 are the same except for case, and 0 otherwise.
<pre>strfind(s,s1)</pre>	Returns the starting indices of any occurrence of string s1 in string s .
<pre>strncmp(s1,s2,n)</pre>	Returns 1 if the first n characters of the strings s1 and s2 are the same, and 0 otherwise.
<pre>strncmpi(s1,s2,n)</pre>	Returns 1 if the first n characters of the strings s1 and s2 are the same
	except for case and 0 otherwise.
eval(s)	Execute string \mathbf{s} with MATLAB expression.

Table E.2: More advanced functions for operating on strings.

```
ans =
    1
>> isequal(mynum,pi)
ans =
    1
>> myprod = [mynumstr, '*2']
myprod =
3.1416*2
>> str2num(myprod)
ans =
    6.2832
```

Some more advanced functions for operating on strings are shown i Table E.2.

It is also possible to operate with *regular expressions* such as 'fig*', see >>help regexp. Some of the functions also accept other arguments — see the Matlab help system for more details.

E.3 Example: automatic generation of figures

Let us see how we can automate the task discussed in Section E.1. We assume that we manually import the data and define variables T and p, and we want to store the figures in files satvappfigI — saturated vapor pressure figure where I takes on values from 1 to 60. The script file displayed in fig. E.1 does the job of generating 60 figures.

The following Matlab command runs the script:

```
Import Wizard created variables in the current workspace.
>> tic; RepetitionDemoModelFit; toc
```

Note that the Matlab command tic resets an internal clock, while command toc tells the elapsed time (in seconds) since the clock was reset.

The response to this command is:

In C:\MyProjects\Documents\FAG\Matlab\Matlab\RepetitionDemoModelFit.m at line 11

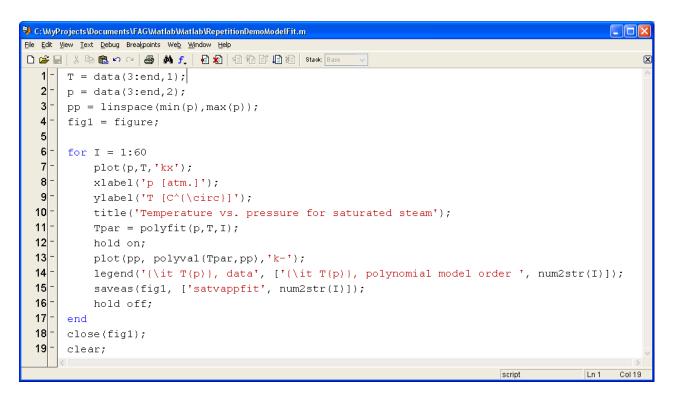


Figure E.1: Script file RepetitionDemoModelFit.m.

```
%
%... this warning is repeated a number of times for large values of I
%
elapsed_time =
4.0750
```

During the execution of the script file, the following files have been created:

```
>> dir satvappfit*
```

```
satvappfit1.fig
                 satvappfit24.fig satvappfit39.fig satvappfit53.fig
satvappfit10.fig satvappfit25.fig satvappfit4.fig
                                                     satvappfit54.fig
satvappfit11.fig satvappfit26.fig satvappfit40.fig satvappfit55.fig
satvappfit12.fig satvappfit27.fig satvappfit41.fig satvappfit56.fig
satvappfit13.fig satvappfit28.fig satvappfit42.fig satvappfit57.fig
satvappfit14.fig satvappfit29.fig satvappfit43.fig satvappfit58.fig
satvappfit15.fig satvappfit3.fig satvappfit44.fig satvappfit59.fig
satvappfit16.fig satvappfit30.fig satvappfit45.fig satvappfit6.fig
satvappfit17.fig satvappfit31.fig satvappfit46.fig satvappfit60.fig
{\tt satvappfit18.fig} \quad {\tt satvappfit32.fig} \quad {\tt satvappfit47.fig} \quad {\tt satvappfit7.fig}
satvappfit19.fig satvappfit33.fig satvappfit48.fig satvappfit8.fig
satvappfit2.fig satvappfit34.fig satvappfit49.fig satvappfit9.fig
satvappfit20.fig satvappfit35.fig satvappfit5.fig
satvappfit21.fig satvappfit36.fig satvappfit50.fig
satvappfit22.fig satvappfit37.fig satvappfit51.fig
satvappfit23.fig satvappfit38.fig satvappfit52.fig
```

We can check the fit of e.g. a 15-th order polynomial model by opening the figure:

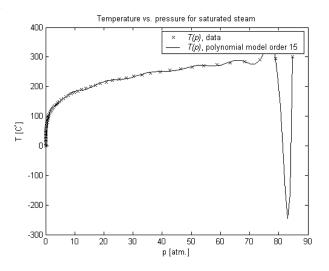


Figure E.2: The model fit of 15-th order polynomial model. Notice that although the model goes through most experimental data points, the model is very poor for interpolation at "high" pressures.

Function	Table E.3: Some functions for number conversion. Description
	Converts a decimal integer to a binary string. Leading zeros are dropped.
	Converts hexadecimal string \mathbf{s} (16 characters) to decimal integer.
nex2num(s)	Converts IEEE hexadecimal string to double precision number, where \mathbf{s} is
	a 16 character string.

>> open satvappfit15.fig

The result is shown in fig. E.2. Notice that although the model goes through most experimental data points, the model is very poor for interpolation at "high" pressures. For higher order models, the interpolation properties of the polynomial model is even poorer.

E.4 Example: hex 2 floating point

Let us consider the algorithm for converting hexadecimal numbers to floating point numbers, as indicated in the footnote p. 21. In doing so, it is convenient to introduce a handful of new Matlab functions, see Table E.3.

See also functions dec2base, base2dec, dec2hex, and bin2dec using the Matlab help system.

The following script does the job:

```
\% Assume that a hexadecimal number hexnum is given as a \% string of 16 hexadecimal numbers
```

% First we find the binary number binnum that represents hexnum:

```
iter = length(hexnum);
binnum = '';
for I = 1:iter,
   % We find the 4 bit equivalent of each hex number, with leading zeros:
   bytedummy = '0000';
   bytenum = dec2bin(hex2dec(hexnum(I)));
   blen = length(bytenum);
   bytedummy(5-blen:end) = bytenum;
```

```
bytenum = bytedummy;
    % We add 4 bit strings to previous string:
    binnum = [binnum, bytenum];
end
% Next we pick out the mantissa sign, the mantissa, and the exponent:
fsign = binnum(1);
fmantissa = binnum(13:64);
fexp1 = binnum(2:12);
% We then add the assumed binary digit "1" in front of the mantissa:
fmantissa = ['1', fmantissa];
\% Next, we convert the mantissa to decimal numbers:
mlen = length(fmantissa);
fman = 0;
for I = 1:mlen
    fman = fman + str2num(fmantissa(I))*2^(-I);
end
% We also need to include the sign of the mantissa:
fman = fman*sign(0.5 - str2num(fsign));
% Next, we compute the value of the temporary exponent:
elen = length(fexp1);
fexp1 = fexp1(elen:-1:1); % Reversing the bit order to simplify evaluation
fex1 = 0;
for I=1:elen.
    fex1 = fex1 + str2num(fexp1(I))*2^{(I-1)};
end
% Next, we subtract the offset of the exponent:
fexp = fex1 - 1022;
% Finally, we compute the floating point number:
fnum = fman*2^fexp
```

In this script, we assume that we have available the hexadecimal representation of the floating point number as 16 hexadecimal digits, where each digit is either of $\{0, 1, 2, \ldots, 9, a, b, c, d, f\}$. This hexadecimal representation must be stored in the variable hexnum.

Let us see how the script (named: hex2num_script) works:

We see that we found the same answer using hex2num_script, as we started with. Let us try another hex string where we do not know the answer:

```
>> hexnum = '701322c182aa1135'
hexnum =
701322c182aa1135
>> hex2num_script
fnum =
7.4272e+231
```

We can check the result by using the built-in Matlab function hex2num:

```
>> hex2num(hexnum)
ans =
7.4272e+231
```

The result appears to be identical to what was found using the script hex2num_script.

Appendix F

Numerical methods with Matlab*

F.1 Numerical functions

The functions in Table F.1 take arrays as arguments.

Table F.1: Some numerical functions in Matlab.		
Matlab function	Description	
diff	difference operator	
gradient	gradient using quadratic interpolating polynomials	
trapz	numeric integral of vector, using the trapeze method	
cumtrapz	As trapz, but returning vector of cumulative values	
polyfit	least squares fit of polynomial to data	
polyval	evaluate polynomial	
interp1	interpolation, various methods, 1D	
interp2	interpolation, various methods, 2D	
roots	roots of polynomial	

We met most of these functions in Part I, and will now apply them to some more examples. We shall learn something about the inner workings of several of these functions later in the course *Numerical Mathematics with Matlab*. For now, we treat them as "black boxes" and just assume that they do their job properly.

Exercise F.1 Generate the following vectors:

x = (0:0.2:10)';

- y = cos(x);
- Use diff to generate a vector dy containing approximate values of the derivative of $\cos x$.
- Plot dy together with $-\sin x$ in the same plot.

NB! Note that diff(y) has one element less than y. What should we use as abscissa axis variable (x-axis variable) when plotting dy?

Exercise F.2 Repeat the previous exercise using the **gradient** function to find approximate values of the derivative.

Exercise F.3 Repeat Exercise F.1, but generate approximate values of the integral instead of the derivative. Plot these together with $\sin x$.

Exercise F.4 Use interp1 and x and y from Exercise F.1 to find the value of sin(p/2).

Operation	Legal/Illegal	Interpretation
s1+s2, s1-s2	L	Addition/subtraction of scalars
s1*s2, s1/s2	\mathbf{L}	Standard scalar product/division
s+M, s-M	Ι	Interpreted as s*ones(size(M))+M , etc.
s*M	\mathbf{L}	Standard multiplication by scalar
s/M	Ι	
M/s	Ι	Interpretation: M./(s*ones(size(M)))
M1+M2, M1-M2	\mathbf{L}	M1 and M2 must have identical size
M1*M2	\mathbf{L}	No. of columns of $M1$ must equal no. rows of $M2$
M1/M2	Ι	_
Mîn	L	Legal if M is quare matrix (including scalar)

Table F.2: Basic binary operations among scalars s and matrices M.

F.2 Arrays vs. Vectors and Matrices

In Part I, we considered mainly arrays, which we "defined" as tables where the content of each element is of the same type. We also considered unary functions on array elements, and basic binary operations between arrays. In the case of binary operations, we required the two arrays, e.g. A, B, to be of the same size, e.g. $A \in \mathbb{C}^{m \times n}$ and $B \in \mathbb{C}^{m \times n}$. Furthermore, we defined addition, subtraction, multiplication, and division of arrays to be element-by-element operations. In Matlab, these are performed using the notation:

- array addition: A+B
- array subtraction: A-B
- array multiplication: A.*B
- array division: A./B

In Section A of Part I, we considered vector spaces and matrices. We saw that a wide-spread vectorspace is \mathbb{C}^n , which conveniently can be represented either by a row array $(\mathbb{C}^{1\times n})$ or by a column array $(\mathbb{C}^{n\times 1})$; for simplicity, we will denote these as a row vector and a column vector, respectively. For vectors, we need to define vector addition and subtraction — these fit well with the concepts of array addition and subtraction (+, -). Furthermore, we need the operation of scalar multiplication. Since a scalar and a vector, considered as arrays, have different sizes, we can not use array multiplication as the vector multiplication operation. Instead, Matlab supports the binary operation * among a scalar and a vector. Next: additive and multiplicative identities fits well with the array representation of Matlab. Finally, association, distribution, and commutativity is built into the way Matlab interprets parentheses.

Matrices can also be represented by arrays (e.g. $\mathbb{C}^{m \times n}$), where matrix addition and subtraction are identical to array addition and multiplication. It is also convenient to expand the meaning of the operator for scalar-vector multiplication (*), to also denote the Cayley product of two matrices. In other words, the (Cayley) matrix product A*B is defined if the number of columns of matrix A is the same as the number of rows of matrix B, in other words: if $A \in \mathbb{C}^{m \times k}$ and $B \in \mathbb{C}^{k \times n}$, then the product AB is well defined, and $AB \in \mathbb{C}^{m \times n}$.

There is no division operation between vectors or matrices, thus symbol / does not really have a meaning when vectors and matrices are involved.

Finally, it should be noticed that for binary operations between two *scalars*, addition, subtraction, multiplication, and division is represented by the symbols +, -, *, and /, respectively. Furthermore, Matlab has a special interpretation for some operations that are theoretically illegal. The operations are summarized in Table F.2.

Note that although matrix addition commutes (i.e. A + B = B + A), in general, matrix multiplication does not commute (i.e. $A \cdot B \neq B \cdot A$).

Also note that matrix transpose M^T is achieved as follows: M.'. If we want to simultaneously find the transpose and take the complex conjugate of the matrix elements (often written as M^H or M^*), this is achieved as follows: M'.

Matlab function	Description
rank	Matrix rank, i.e. number of linearly independent rows/columns
det	Determinant, valid for square matrices.
cond	Condition number, measure of how close matrix is to singularity
inv	Matrix inverse, valid for square, non-singular matrices
	Solve linear equations using Gauss elimination
norm	Matrix/vector norm, extension of Pythagorean length of vectors
eig	Eigenvalues and eigenvectors

Table F.3: Basic Matrix operations

F.3 Matrix operations

F.3.1 Basic operations

Some basic matrix functions are given in Table F.3.

Exercise F.5 For the set of equations:

formulate the equations as a matrix equation of form Ax = b.

Exercise F.6 For matrix A of Exercise F.5:

- Find the rank of A.
- Find the determinant of A.
- Find the condition number of A.
- If possible, find the inverse of A.

Exercise F.7 For vector b and matrix A in Exercise F.5:

- Find the norm of vector *b*.
- Find the norm of matrix A.
- Experiment with other norms for b and A than the default norm (Hint: use the help system), and compare the results.

Exercise F.8 If possible:

- Find the solution x of the equation Ax = b in Exercise F.5 using the backslash operator /.
- Verify that you get the same result using the matrix inverse.

F.3.2 Matrix factorizations/decompositions¹

Matlab allows for the basic matrix factorizations/decompositions in Table F.4.

¹These methods are used extensively in linear algebra.

Factorization of A	Matlab command	Comment
$A = R^T R$	R=chol(A)	Cholesky factorization: A is positive definite, R is
PA = LU	[L,U,P] = lu(A)	upper triangular LU factorization: A is square, L is lower triangu-
		lar, U is upper triangular, P is permutation matrix $(P^T P = I)$
AE = QR	[Q,R,E] = qr(A)	QR decomposition: E is permutation matrix $(E^T E = I), Q$ is unitary matrix $(Q^H Q = I), R$ is
		upper triangular
$A = USV^H$	[U,S,V] = svd(A)	Singular value decomposition: U and V are uni- targe matrices $(U^{H}U) = U^{H}V = U$ and C is
		tary matrices $(U^H U = I, V^H V = I)$, and S is diagonal

Table F.4: Basic matrix factorizations in Matlab.

F.3.3 Vector space commands²

With $A \in \mathbb{C}^{m \times n}$ and $x \in \mathbb{C}^n$, Ax can be considered as the column vector $Ax = \sum_{i=1}^n x_i a_i$, where $a_i \in \mathbb{C}^m$ is column *i* of matrix A. The subspace consisting of all possible vectors Ax (i.e. by varying x) is known as the *column space* of matrix A, $\mathcal{R}(A)$. This subspace has the same number of so-called basis vectors, as the rank of A, rank A.

• The command orth(A) responds with a matrix with columns which constitute a complete set of basis vectors for $\mathcal{R}(A)$.

The particular matrix equation Ax = 0 may have many solutions. The set of possible solutions forms a subspace of \mathbb{C}^n which is known as the *nullspace* of A, denoted $\mathcal{N}(A)$. The nullspace can be characterized by a complete set of basis vectors. With $A \in \mathbb{C}^{m \times n}$, it can be shown that the number of basis vectors in $\mathcal{N}(A)$ is $n - \operatorname{rank} A$.

- The command null(A) responds with a matrix with columns which constitute a complete set of basis vectors for $\mathcal{N}(A)$.
- The command null(A,'r') has the same response as null(A), but with operation null(A,'r'), the resulting basis vectors are rational numbers.

The concepts column space $\mathcal{R}(A)$ and nullspace $\mathcal{N}(A)$ are used in order to assess the solvability of equations Ax = b. In particular:

- 1. A solution exists for equation Ax = b if and only if $b \in \mathcal{R}(A)$. We have that $b \in \mathcal{R}(A) \Leftrightarrow \operatorname{rank}[A, b] = \operatorname{rank} A$.
- 2. If a solution exists for Ax = b, then the solution is unique if and only if $\mathcal{N}(A) = \{0\}$ (i.e. the nullspace has zero basis vectors).

²These methods are used extensively in advanced linear algebra.

Appendix G

Extra parameters and anonymous functions*

Before going into more details about how to use function functions, let us consider how anonymous functions can simplify the *interface* between our own functions and built-in Matlab functions.

Many of the built-in Matlab functions, e.g. in Table 15.1, accept extra input parameters. As an example, the quad function accepts the more general form:

quad(fh1, a, b, tol, trace, P1, P2, ...)

Here, the first input argument is the function handle to the function we want to integrate (fh1), the second input argument is the lower integration limit (a), the third input argument is the upper integration limit (b), the fourth input argument is a user specified integration tolerance (tol), the fifth input argument specifies whether the user wants to see the intermediate results during the iteration for a solution (trace), while the remaining input arguments are extra parameters to the function referenced by the function handle (fh1). If we do not want to use parameters tol and trance, we can replace each of these by []:

```
quad(fh1, a, b, [], [], P1, P2, ...)
```

We can alternatively use extra parameters with **quad** by introducing an anonymous function as an interface:

fh2 = @(x) fh1(x, P1, P2, ...);
quad(fh2, a, b)

Using the anonymous function as an interface function, the notation and use of function functions with extra arguments becomes simpler.

Let us illustrate this idea with an example. We shoot up a projectile with an initial velocity v_0 at an angle φ with the horizon. The velocities in the horizontal an vertical directions are

$$v_x = v_0 \sin \varphi$$
$$v_y = v_0 \cos \varphi - gt$$

The speed is thus

$$v = \sqrt{v_x^2 + v_y^2} = \sqrt{v_0^2 + g^2 t^2 - 2v_0 g t \cos \varphi}$$

If the ground is in the horizon plane, the time before the projectile hits the ground is the solution for t > 0 of $\int_0^{t_{\text{max}}} v_y(t) dt = 0$ or $v_0 t_{\text{max}} \cos \varphi - \frac{1}{2}gt_{\text{max}}^2 = 0$:

$$t_{\max} = \frac{2v_0}{g}\sin\varphi$$

We now want to find the arc length of the trajectory, defined as

$$L = \int_0^{t_{\max}} v dt$$

If we were only interested in the result for one set of specified values for v_0 , φ , and g, then we could hard code those values in a function and integrate it using quad. But it is more flexible if we write a function with those parameters as input variables:

```
function v = speed(t, v0 , phi, g)
% Speed of projectile shot up with initial velocity v0 at angle
% phi with the horizon. g: gravitational acceleration
vx = v0*cos(phi);
vy = v0*sin(phi)-g*t;
v = sqrt(vx^2+vy.^2);
```

There are two ways of integrating this function using quad. Without using an anonymous function as interface, we have:

```
>> v0 = 10;g = 10;phi = pi/4;
>> tmax = 2*v0*sin(phi)/g;
>> L = quad(@speed, 0, tmax, [], [], v0, phi, g)
L =
11.4779
```

This works fine, but the user has to enter values (or empty arrays as placeholders) to ensure that the parameter arguments come at the right place in the function call. Another, perhaps more serious, problem is that the method for passing extra parameters when writing your own function functions is not so straightforward. In fact, most users find it a constant source of frustration.

In Matlab 7, we can alternatively use the anonymous function facility to define a new interface function:

```
>> f = @(t) speed(t, v0, phi, g)
f =
  @(t) speed(t,v0,phi,g)
```

Next, we can integrate **f** to find the arc length:

```
>> L = quad(f,0,tmax)
L =
11.4779
```

We can also define the function inside the call to quad. In this example, we change the angle a bit, just to demonstrate that the parameters do indeed influence the result:

```
>> phi = pi/3;
>> L = quad(@(t)speed(t,v0,phi,g),0,tmax)
L =
9.1949
>> L = quad(f,0,tmax)
L =
11.4779
```

Note that **f** is unaffected by the value change of the parameter.

Appendix H

DAE solvers in Matlab*

Two of the Matlab ODE solvers can also solve DAEs: ode15s and ode23t. Regarding the damped mass-spring system, we could alternatively have posed this as a DAE system:

$$\frac{dx}{dt} = v \tag{H.1}$$

$$\frac{dE}{dt} = Fv - \mu v^2 \tag{H.2}$$

$$P = mgx + \frac{k}{2} (x - x_0)^2$$
(H.3)

$$E = K + P \tag{H.4}$$

$$K = \frac{1}{2}mv^2. \tag{H.5}$$

However, this models is not entirely equivalent to the model in eqs. 16.1 – 16.2: with the given initial values, $(x, v)_{t=0} = (1.5, 0)$, we see that dx/dt = 0 and dE/dt = 0, hence the model in eqs. H.1 – H.5 will predict that the system stays at rest!

Instead, we choose to consider the following DAE:

$$\frac{dx}{dt} = v$$

$$\frac{dv}{dt} = -\frac{k}{m}(x - x_0) - \frac{\mu}{m}v - g + \frac{F}{m}$$

$$0 = P - mgx - \frac{k}{2}(x - x_0)^2$$

$$0 = K - \frac{1}{2}mv^2$$

$$0 = E - K - P.$$

This model obviously is equivalent: we simply use the same model for x and v, and add three algebraic equations in order to compute P, K and E. These equations constitute a DAE: a differential algebraic equation with 2 ordinary differential equation, and 3 additional algebraic equations. The standard form for DAEs in Matlab is

$$M\frac{dy}{dt} = f\left(t, y\right).$$

Here, we define

$$y = (x, v, P, K, E)^T.$$

In standard form, we have:

Notice that we have to specify F as a function of t or y in order to have the system in standard form. Notice also that we have a choice as to which of the algebraic equations belongs to dv/dt, dK/dt, and dP/dt. We can make that choice more or less as we like.

Since Matlab will attempt to compute y, notice that this formulation allows for automatic computation of all quantities of interest: x, E, v, K, P, and that no post processing is necessary.

The main problem with solving DAEs is that we need consistent conditions for y(0) and dy(0)/dt. Here, it is simplest to specify x(0) and v(0), and to compute E(0), K(0), and P(0) from the algebraic equations. Furthermore, we need a strategy for finding dy(0)/dt. We need to require that

$$M\frac{dy\left(0\right)}{dt} = f\left(0, y\left(0\right)\right)$$

which is a simple, linear algebra equation of type Ax = b, where A = M, the unknown is x = dy(0)/dt, and b = f(0, y(0)). We can solve this equation using the pseudo inverse of M, which in Matlab is computed through the use of function **pinv**.

Here is how we call DAE solvers for the damped mass-spring system (script file dampmsdae.m):

% Script for simulating damped mass-spring system

```
% Necessary data
Tfin = 150;
data;
xinit = 1.5;
v0 = 0;
PO = m*g*xinit + k*(xinit-x0)^2/2;
KO = m * v 0^2/2;
EO = KO + PO;
% Setting up initial values:
y0 = [xinit, v0, P0, K0, E0]';
% Setting up the M matrix in M*dy/dt = f(t,y):
M = zeros(length(y0));
M(1,1) = 1;
M(2,2) = 1;
% Computing consistent value for dy(t0)/dt:
MdyOdt = dmsdae(0,y0);
dyOdt = pinv(M)*MdyOdt;
% Specifying options for the ode solver:
opt = odeset('Mass',M,'MassSingular','yes', 'MstateDependence', 'none',...
    'InitialSlope', dyOdt, 'AbsTol', 1e-6, 'RelTol', 1e-3);
\% Simulating system - notice that we include the options opt
tic;
[T,Y] = ode15s(@dmsdae, [0,Tfin], y0, opt);
toc
```

Next, we need to define function dmsdae:

```
function fy = dmsdae(t,y)
%
% Defining data
data;
% Naming variables
x = y(1);
```

```
v = y(2);
P = y(3);
K = y(4);
E = y(5);
% Defining elements of the vector field
fx = v;
fv = -k*(x-x0)/m - mu*v/m - g + F/m;
fP = P - m*g*x - k*(x-x0)^2/2;
fK = K - m*v^2/2;
fE = E - K - P;
% Setting up the vector field
fy = [fx, fv, fP, fK, fE]';
```

The DAE is solved by running the script file dampmsdae. Solving the DAE (variables: x, v, P, K, E) takes approximately 3 times as long as solving the simple ODE (variables: x, v), but then we find more variables.

To sum up:

- Describing models as DAEs reduces¹ the necessary amount of formula manipulation.
- We can reduce/eliminate the necessary work to find auxiliary variables², if we describe models as DAEs.

¹Note: in this case, we didn't really save much manipulation, but normally we can save work by formulating models as DAEs. ²In this case, we eliminated the necessary post processing to find P, K, and E.

Part VI References

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