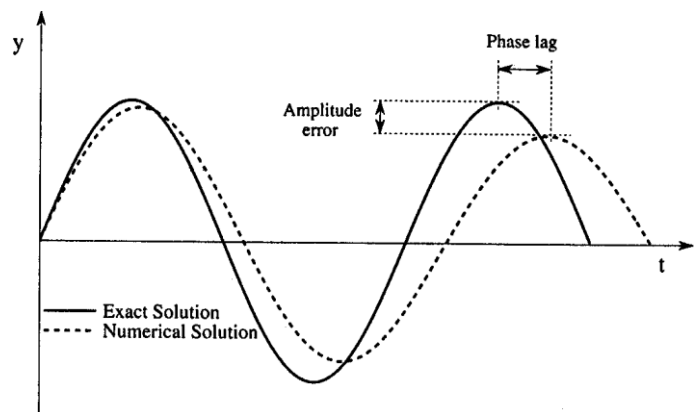
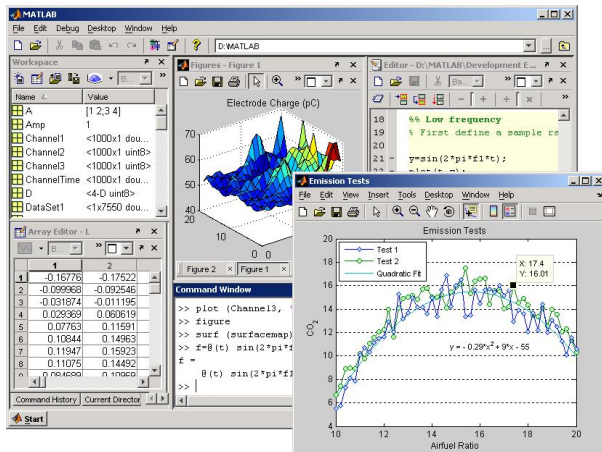
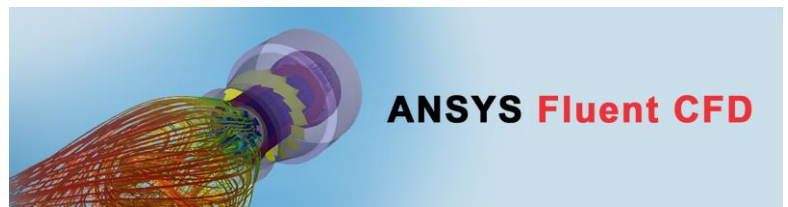




## UE 703 Méthodes Numériques pour la Mécanique

### POLYCOPE DE TP – M1 ENERGIE



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Lectures web page : <https://arche.univ-lorraine.fr/course/view.php?id=6745>



# Analyse Numérique - Master MEPP

Enoncés des TP - Introduction à MATLAB pour le calcul numérique

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## 0.1 Organisation des TP

Les TP d'analyse numérique déroulent en 8 séances de 3h de la manière suivante :

- 5 séances d'analyse numérique avec MATLAB ,
- 3 TP d'initiation à FLUENT, le logiciel commercial de simulation en mécanique des fluides et thermique.

Si vous n'avez aucune connaissance en MATLAB , il est recommandé de réaliser les TP d'introduction dont l'énoncé se trouve dans ce poly. Pour réaliser les TP d'analyse numérique, vous devez télécharger sur ARCHE (<http://arche.univ-lorraine.fr>) l'archive 'tpmatlab\_m1mepp.zip' du cours *UE104* et l'extraire sur votre PC. Cette archive contient la documentation et les programmes MATLAB à utiliser et compléter lors des séances.

## 0.2 Notation des travaux pratiques

Les TP seront notées d'une part par une de votre travail et assiduité durant les séances, et sur la base de compte-rendus individuels que vous devez rendre à la fin du semestre.

lieu à 2 notes : une pour l'EC '*Introduction à MATLAB* ', et une note de TP pour l'EC '*Analyse numérique*'.

- Les compte-rendu devront répondre aux questions de l'énoncé, et comporter 6 pages au maximum. Les réponses doivent être rédigées de façon précise et succincte (ne recopiez pas l'énoncé !), et seront illustrées par les graphiques que vous avez réalisés avec MATLAB . Dans certains cas, il vous sera demandé de rendre le listing de votre programme MATLAB .
- Les compte-rendu doivent être rédigés avec le traitement de texte WORD ou convertis en fichiers PDF. Ne rendez pas de fichiers MATLAB . N'oubliez pas de mettre votre nom, votre parcours licence et l'intitulé du TP sur les compte-rendu.

# Analyse Numérique - Master MEPP

## Introduction à MATLAB pour le calcul numérique

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## 1 Prise en main de MATLAB

### 1.1 Introduction

MATLAB (abréviation de *MATrix LABoratory*) est un système interactif, basé sur le calcul matriciel, qui permet de réaliser des calculs scientifiques sans nécessiter d'écrire de longs programmes informatiques. Le but de ces séances de TP est de vous introduire aux fonctionnalités de base de MATLAB dans l'optique de son utilisation en analyse numérique. Néanmoins, MATLAB vous sera utile non seulement en calcul scientifique, mais aussi pour l'exploitation de données expérimentales.

### 1.2 Premiers pas dans MATLAB

Pour réaliser ces TP, vous devez télécharger sur ARCHE (<http://arche.univ-lorraine.fr>) l'archive 'tpmatlab\_m1mepp.zip' du cours *UE 104* et l'extraire sur votre PC. Cette archive contient la documentation et les programmes MATLAB à utiliser et compléter lors des séances.

Sur votre PC, allez dans le répertoire de travail '*TP0 INITIATION MATLAB*' qui contient les programmes MATLAB (fichiers *\*.m*) pour les séances d'initiation. Le fichier '*Tp0\_matlab.m*' contient les commandes (ou *instructions*) MATLAB utiles au premier TP. Ouvrez ce fichier par double-clic, et l'interface graphique représentée sur la figure 1 apparaîtra à l'écran.

L'interface graphique de MATLAB se décompose en :

- **Command Window** : pour entrer les commandes MATLAB (après le prompt `>>`) et voir les résultats s'afficher.
- **MATLAB Editor** : pour éditer les fichiers *\*.m* contenant les commandes et fonctions MATLAB.
- **Current Folder** : affiche l'ensemble des fichiers *\*.m* contenus dans le répertoire de travail.
- **Workspace** : affiche les données, matrices, *etc* ... que vous créez durant la séance de travail.
- **Command History** : historique des commandes de la séance de travail.

### 1.3 Exécution d'un programme MATLAB

Il y a 3 manières équivalentes d'exécuter le programme '*Tp0\_matlab.m*' :

- Taper `>> Tp0_matlab` dans la **Command Window**,
- Cliquer sur le bouton  dans la barre du menu **EDITOR** de l'interface graphique,
- Placer la souris dans le **MATLAB Editor** du programme '*Tp0\_matlab.m*' et taper F5 au clavier.

Le résultat du programme s'affiche dans la **Command Window**. Il est à noter que le programme '*Tp0\_matlab.m*' utilise une fonction définie dans le fichier '*fonction\_ex0.m*' qui doit nécessairement se trouver dans le répertoire de travail, sinon MATLAB ne trouvera pas la fonction et affichera un message d'erreur.

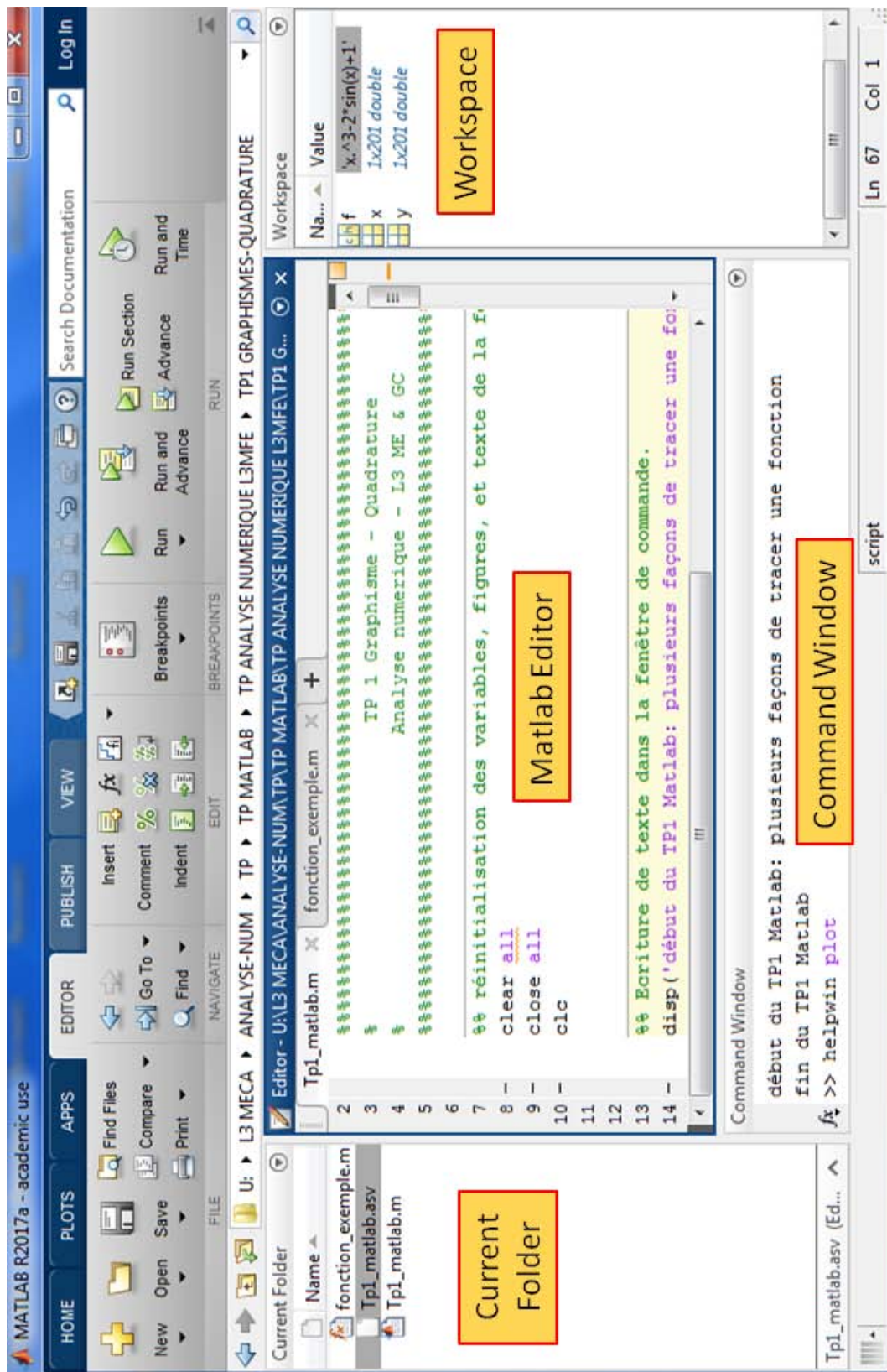


Figure 1: Interface graphique de MATLAB .

## 2 TP d'initiation à MATLAB

### 2.1 Introduction

Vous devez effectuer en 3 séances les exercices donnés ci-dessous. Pour cela, vous devez lire en premier le polycopié '*Introduction à MATLAB*' et exécuter tous les exemples qu'il contient. Il sera souvent nécessaire de rechercher des informations complémentaires au poly dans l'aide en ligne de MATLAB :

>>helpwin commande

### 2.2 Enoncés des exercices d'initiation

#### Exercice 1. (*Algèbre linéaire et calcul matriciel*)

On considère la matrice rectangulaire  $\mathcal{M}$  définie par :

$$\mathcal{M} = \begin{bmatrix} 4 & 8 & 12 \\ 3 & 8 & 13 \\ 2 & 9 & 18 \\ -1 & 0 & 1 \end{bmatrix}$$

- 1) Créez la matrice  $\mathcal{M}$  dans MATLAB . Donnez le nombre de lignes et de colonnes (commande `size`).
- 2) Extrayez de  $\mathcal{M}$  la matrice carrée :

$$\mathcal{A} = \begin{bmatrix} 4 & 8 & 12 \\ 3 & 8 & 13 \\ 2 & 9 & 18 \end{bmatrix}$$

- 3) On considère le système linéaire  $\mathcal{A}X = F$ , tel que le second membre  $F$  est défini par :

$$F = \begin{pmatrix} 4 \\ 5 \\ 11 \end{pmatrix}$$

Montrez que le système possède une solution unique en calculant son déterminant et résolvez le système linéaire.

- 4) Vérifiez que la solution donnée par MATLAB est exacte en calculant le résidu  $R = F - \mathcal{A}X$ .

#### Exercice 2. (*Graphisme sous MATLAB*)

Cet exercice a pour objet d'initier aux multiples possibilités graphiques de MATLAB . Utilisez l'aide en ligne (`helpwin plot`) pour le détail des options disponibles.

- 1) Sur une même fenêtre graphique (commande `subplot`) tracez les trois graphes représentant dans  $[-\pi, \pi] \times [-1, 1]$  la fonction  $x \mapsto \sin(x)$  et ses développements de Taylor autour de  $x = 0$ , i.e.  $x \mapsto x$  et  $x \mapsto x - x^3/3$ . Vous devrez respecter la mise en forme de la figure 2, c'est-à-dire que chaque graphe ait une couleur et un trait différent, et que les figures soient correctement légendées.
- 2) Utilisez la barre des tâches de la fenêtre graphique MATLAB pour exporter les graphes sous format *jpeg* ou *png*, ainsi que pour faire un zoom autour du point  $x = 0$ .

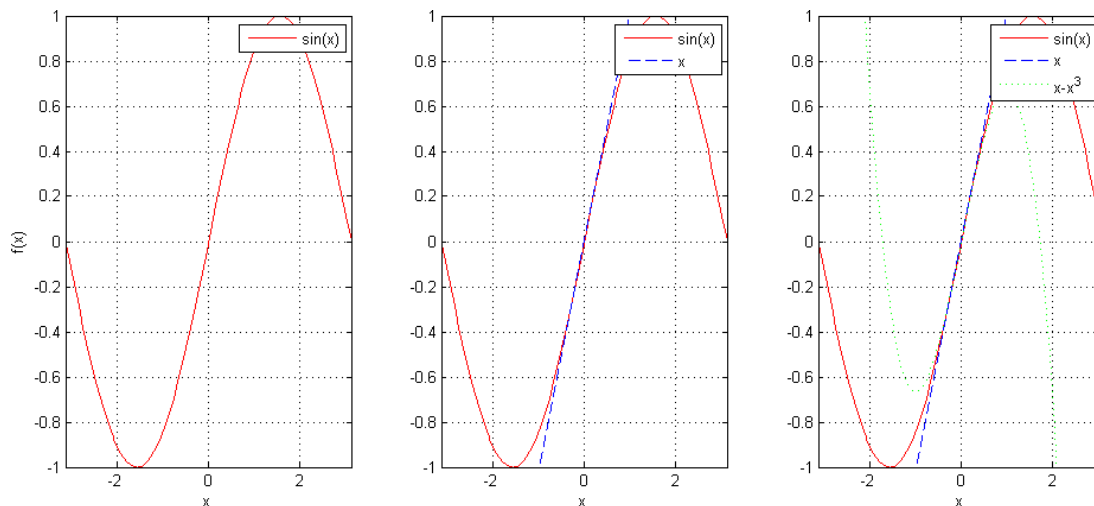


Figure 2: Graphisme obtenu sous Matlab

**Exercice 3. (Fonctions MATLAB et boucles de contrôle)**

- 1) A l'aide de la boucle `for`, Ecrivez la fonction MATLAB `F=fact(n)` qui calcule la factorielle de l'entier  $n$  (Faites un test conditionnel (boucle `if`) qui calcule  $n!$  seulement si  $n \geq 0$ ).  
*Application* : calculez la factorielle de -1, 0 et 3.

- 2) Ecrivez la fonction `R=residumat(A,X,F)` qui calcule le résidu du système linéaire  $AX = F$  de taille  $n \times m$ , soit :

$$\text{Pour } i = 1, \dots, n, \quad R_i = F_i - \sum_{j=1}^m A_{i,j} X_j$$

Testez-la sur l'exemple de l'exercice 1.

- 3) Ecrivez la fonction MATLAB `[r1,r2]=racines_poly(a,b,c)` qui calcule les racines *réelles* du polynôme du second degré  $ax^2 + bx + c$ , en faisant un test suivant la valeur du discriminant  $\Delta$ . Appliquez-le au calcul des racines de  $x^2 + x - 2$ .

**Exercice 4. (Intégration numérique par méthode du trapèze)**

- 1) Ecrivez la fonction MATLAB `trapeze(a,b,f,n)` qui calcule l'intégrale  $\int_a^b f(x)dx$  par la formule du trapèze composite avec  $n$  sous intervalles de taille uniforme<sup>1</sup>.
- 2) Validez la fonction que vous avez écrite en vérifiant qu'elle donne l'intégrale exacte d'une fonction linéaire (polynôme de degré 1).

<sup>1</sup>On rappelle que, d'après le cours, cette formule s'écrit :

$$I_h = \frac{h}{2} \sum_{k=0}^{n-1} [f(x_k) + f(x_{k+1})],$$

avec  $h = (b - a)/n$  et  $x_k = a + kh$ .

3) Vous allez maintenant utiliser votre fonction `trapeze(a,b,f,n)` pour calculer l'intégrale

$$I = \int_1^{\pi} \frac{\sin x}{2x^3} dx \simeq 0.198557 \dots \quad (1)$$

qui a été vue en TD d'analyse numérique. Il est possible de calculer la valeur exacte de cette intégrale à l'aide de la *toolbox* de calcul formel symbolic de MATLAB, en tapant ces lignes :

```
>> syms x
>> I=double(int(f_exo_trapeze(x),1,pi))
```

- a) Pour un nombre de sous-intervalles  $n$  suffisamment grand, calculez le facteur de réduction de la méthode du trapèze et vérifiez qu'on retrouve le résultat théorique.
- b) Pour  $n \in [50, 200]$ , représentez en échelle semi-logarithmique (commande `semilogy`) l'erreur numérique  $E_n = |I_n - I|$  de la méthode du trapèze et montrez qu'elle est au second-ordre en traçant sur le même graphe la fonction  $n \mapsto 1/n^2$ .



# MATLAB Quick Reference

## Operators

| Matrix Operations |                                                 | Array or Element by Element |                                                         |
|-------------------|-------------------------------------------------|-----------------------------|---------------------------------------------------------|
| +                 | Addition                                        |                             |                                                         |
| -                 | Subtraction                                     |                             |                                                         |
| *                 | Matrix Multiplication                           | .*                          | Element by Element Multiplication                       |
| /                 | Right Matrix Division $b/A = bA^{-1}$           | ./                          | Element by Element Right Division                       |
| \                 | Left Matrix Division $A \setminus b = A^{-1}b$  | .\                          | Element by Element Left Division $A \setminus B = B./A$ |
| ^                 | Raise Matrix to a power                         | .^                          | Raise each Element to a power                           |
| '                 | Transpose matrix (conjugate if complex)<br>$A'$ | .'                          | Transpose matrix (no complex conjugation) $A.'$         |

| Relational Operators |                       | Logical Operators |     |
|----------------------|-----------------------|-------------------|-----|
| <                    | Less Than             | &                 | And |
| <=                   | Less Than or Equal    |                   | Or  |
| >                    | Greater Than          | ~                 | NOT |
| >=                   | Greater Than or Equal |                   |     |
| ==                   | Equal                 |                   |     |
| ~=                   | Not Equal             |                   |     |

## Special Number Symbols

|     |                     |
|-----|---------------------|
| pi  | $\pi$               |
| inf | $\infty$            |
| NaN | Not a number ie 0/0 |

|   |             |
|---|-------------|
| i | $\sqrt{-1}$ |
| j | $\sqrt{-1}$ |

## Entering Matrices

|                                   |                                                                                                      |
|-----------------------------------|------------------------------------------------------------------------------------------------------|
| $x = [1\ 2\ 3; 4\ 5\ 6; 7\ 8\ 9]$ | Comma or space between elements.<br>Semicolon or return for new row.<br>Enclosed in square brackets. |
| $x = [1,2,3,4,5,6,7,8,9]$         |                                                                                                      |
| $a = [\exp(0)\ \sqrt{4}\ 1+2]$    | Each element can be an expression.                                                                   |
| $C = [A\ B]$                      | A and B are matrices with the same number of rows.                                                   |
| $C = [A;\ B]$                     | A and B are matrices with the same number of columns.                                                |

## Indexing

| Vector Indexing |                                                    |
|-----------------|----------------------------------------------------|
| variable(V)     | V is a vector of indexes                           |
| Examples        |                                                    |
| A(3)            | The third element in vector A.                     |
| A([3 8 11])     | The third, eighth and eleventh elements of A.      |
| A(3:11)         | The third to eleventh elements of A.               |
| A(11:end)       | All elements from the eleventh to the last element |

| Matrix Indexing |                                                      |
|-----------------|------------------------------------------------------|
| variable(R,C)   | R is a vector of rows and C a vector of Columns      |
| Examples        |                                                      |
| M(2,3)          | The element in the second row and third column of M. |
| M(2, [3 8 11])  | Second row, third, eighth and eleventh column.       |
| M(4,3:11)       | Fourth row, third to eleventh column.                |
| M(11:end,2)     | Eleventh to the last row, second column.             |
| M(5,:)          | Fifth row, all columns.                              |
| M(11:end,:)     | Eleventh to the last row, all columns.               |

## Output Display Format

|                 |                                        |                |                                         |
|-----------------|----------------------------------------|----------------|-----------------------------------------|
| format short    | 4 decimal places                       | format long    | 15 decimal places                       |
| format          | Same as above                          |                |                                         |
| format shortE   | Scientific notation, 4 decimal places  | format longE   | Scientific notation, 15 decimal places  |
| format shortEng | Engineering notation, 4 decimal places | format longEng | Engineering notation, 15 decimal places |

## Generating Vectors

|                                                                                              |                                                     |
|----------------------------------------------------------------------------------------------|-----------------------------------------------------|
| <b>&lt;start&gt; : &lt;end&gt;</b>                                                           | <b>x = 1 : 5</b> generates x = [ 1 2 3 4 5]         |
| <b>&lt;start&gt; : &lt;separation&gt; : &lt;end&gt;</b>                                      | <b>y = 0:5:20</b> generates y = [ 0 5 10 15 20]     |
| <b>linspace(start,end,n)</b> n = number of elements                                          | <b>linspace(0,10,5)</b> generates [ 0 2.5 5 7.5 10] |
| <b>logspace(d1,d2,n)</b> n elements logarithmically spaced between $10^{d1}$ and $10^{d2}$ . | <b>logspace(-1,2,4)</b> generates [0.1 1 10 100]    |

## Utility Matrices

|                     |                                           |
|---------------------|-------------------------------------------|
| <b>zeros(n)</b>     | n by n matrix where each element is zero. |
| <b>zeros(m,n)</b>   | m by n matrix where each element is zero. |
| <b>zeros(a,b,c)</b> | 3 dimensional array, a by b by c.         |
| <b>ones(m,n)</b>    | m by n matrix where each element is one.  |
| <b>rand(m,n)</b>    | m by n matrix of random numbers.          |
| <b>eye(n)</b>       | n by n identity matrix.                   |

## Variable Control

|                               |                                          |
|-------------------------------|------------------------------------------|
| <b>who</b>                    | List all variables in memory.            |
| <b>whos</b>                   | Same as above but with more information. |
| <b>clear</b>                  | Remove all variables from memory.        |
| <b>clear &lt;variable&gt;</b> | Remove specified variables from memory.  |

## File Control Commands

|                                |                                                 |
|--------------------------------|-------------------------------------------------|
| <b>dir</b>                     | List contents of current directory.             |
| <b>ls</b>                      | List contents of current directory.             |
| <b>what</b>                    | List the Matlab files in the current directory. |
| <b>cd &lt;directory&gt;</b>    | Change the current directory.                   |
| <b>type &lt;filename&gt;</b>   | Display the contents of a text or .m file.      |
| <b>delete &lt;filename&gt;</b> | Delete a file.                                  |
| <b>diary &lt;filename&gt;</b>  | Record all commands and results to a file.      |
| <b>diary off</b>               | Stop above.                                     |

## Help

|                              |                                         |
|------------------------------|-----------------------------------------|
| <b>help</b>                  | Display help topics.                    |
| <b>help &lt;function&gt;</b> | Help on a particular function.          |
| <b>lookfor &lt;word&gt;</b>  | Look for word in function descriptions. |
| <b>doc &lt;function&gt;</b>  | Full documentation on function.         |

## In Built Functions

Hit **fx** icon next to prompt for function browser. Only selected functions shown here.

|                 |                                    |                 |                               |
|-----------------|------------------------------------|-----------------|-------------------------------|
| <b>abs(x)</b>   | The absolute value. Modulus        | <b>round(x)</b> | Round to the nearest integer. |
| <b>sqrt(x)</b>  | The square root.                   | <b>ceil(x)</b>  | Round up.                     |
| <b>exp(x)</b>   | The exponential base e. $e^x$      | <b>floor(x)</b> | Round down.                   |
| <b>log(x)</b>   | The natural logarithm. $\log_e(x)$ | <b>fix(x)</b>   | Round towards zero.           |
| <b>log10(x)</b> | The log base 10. $\log_{10}(x)$    | <b>rem(x,b)</b> | Remainder of x divided by b.  |

## Trigonometry

|                                                                      |                                             |
|----------------------------------------------------------------------|---------------------------------------------|
| <b>sin(x)</b>                                                        | Sine. x in radians.                         |
| <b>sind(x)</b>                                                       | Sine. x in degrees.                         |
| <b>asin(x)</b>                                                       | The arcsine. The inverse of sin(x). Radians |
| <b>asind(x)</b>                                                      | The arcsine. Degrees                        |
| <b>sinh(x)</b>                                                       | Hyperbolic Sine.                            |
| <b>asinh(x)</b>                                                      | The inverse Hyperbolic Sine.                |
| The above variations are also available for the following functions. |                                             |
| <b>cos(x)</b>                                                        | Cosine                                      |
| <b>tan(x)</b>                                                        | Tangent                                     |
| <b>sec(x)</b>                                                        | Secant                                      |
| <b>cot(x)</b>                                                        | Cotangent                                   |
| <b>csc(x)</b>                                                        | Cosecant                                    |

## Complex Numbers

|                |                             |                 |                          |
|----------------|-----------------------------|-----------------|--------------------------|
| <b>real(z)</b> | The real part of z.         | <b>imag(z)</b>  | The imaginary part of z. |
| <b>abs(z)</b>  | The modulus of z.           | <b>angle(z)</b> | The phase angle of z.    |
| <b>conj(z)</b> | The complex conjugate of z. |                 |                          |

## Matrix

|                     |                                      |                 |                                |
|---------------------|--------------------------------------|-----------------|--------------------------------|
| <b>det(A)</b>       | Determinant                          | <b>sqrtm(A)</b> | The matrix square root.        |
| <b>norm(A)</b>      | Norm                                 | <b>expm(A)</b>  | The matrix exponential base e. |
| <b>inv(A)</b>       | Inverse                              | <b>logm(A)</b>  | The matrix natural logarithm.  |
| <b>[v,d]=eig(A)</b> | d = Eigenvalues,<br>v = Eigenvectors |                 |                                |

## Statistics

|               |         |                  |         |               |                    |
|---------------|---------|------------------|---------|---------------|--------------------|
| <b>max(x)</b> | Maximum | <b>median(x)</b> | Median  | <b>var(x)</b> | Variance           |
| <b>min(x)</b> | Minimum | <b>mean(x)</b>   | Average | <b>std(x)</b> | Standard Deviation |

## Polynomials

|                                                                                       |                                                                  |
|---------------------------------------------------------------------------------------|------------------------------------------------------------------|
| $p = [1 \ 2 \ 3 \ 4 \ 5]$ ; can represent the polynomial $x^4 + 2x^3 + 3x^2 + 4x + 5$ |                                                                  |
| <b>y = polyval(p,x)</b>                                                               | Evaluate polynomial for each value in x.                         |
| <b>roots(p)</b>                                                                       | Roots of polynomial.                                             |
| <b>p = poly( &lt;roots&gt;)</b>                                                       | Polynomial with given roots.                                     |
| <b>p = polyfit(x,y,n)</b>                                                             | Best fit of x,y data points to $n^{\text{th}}$ order polynomial. |

## Saving, Exporting and Importing Data

|                                                      |                                                                        |
|------------------------------------------------------|------------------------------------------------------------------------|
| <b>save &lt;filename&gt;</b>                         | Save all variable to the file <i>filename.mat</i>                      |
| <b>load &lt;filename&gt;</b>                         | Load in variables from the file <i>filename.mat</i>                    |
| <b>save &lt;filename&gt; &lt;variable&gt;</b>        | Save only the variable <i>variable</i> to the file <i>filename.mat</i> |
| <b>save &lt;filename&gt; &lt;variable&gt; -ascii</b> | Save <i>variable</i> to the text file <i>filename</i>                  |
| <b>load &lt;filename&gt;.&lt;ext&gt;</b>             | Load from the text file, to a variable called <i>filename</i>          |

## Misc

|                                    |                                          |
|------------------------------------|------------------------------------------|
| <b>x = input( 'What is x ? ' )</b> | Ask the user for a number.               |
| <b>[x,y] = ginput(1)</b>           | Graph coordinates of a clicked on point. |
| <b>pause</b>                       | Wait for the user to press a key.        |
| <b>pause(5)</b>                    | Wait for 5 seconds                       |
| <b>! command</b>                   | Execute an operating system command      |
| <b>display(a)</b>                  | Suppose a = 5. This would print "a = 5"; |
| <b>disp(a)</b>                     | Same as above, but with out the "a = ".  |

## fprintf Formatting

|           |                              |               |                       |           |                   |
|-----------|------------------------------|---------------|-----------------------|-----------|-------------------|
| <b>%f</b> | Fixed point                  | <b>%5f</b>    | 5 characters wide     | <b>\n</b> | New line          |
| <b>%e</b> | Exponential Notation         | <b>%5.2f</b>  | 2 decimal places (dp) | <b>\t</b> | Horizontal tab    |
| <b>%g</b> | auto chooses %f or %e        | <b>%-5.2f</b> | Left justify          | <b>\\</b> | Back slash        |
| <b>%s</b> | String of Characters         | <b>%+5.2f</b> | Print sign (+ or -)   | <b>%%</b> | Percent character |
| <b>%d</b> | Integer (Whole numbers only) |               |                       |           |                   |

| Examples                              | What is printed     | Comment                                         |
|---------------------------------------|---------------------|-------------------------------------------------|
| <b>fprintf(' x = %f ',pi)</b>         | <b>x = 3.141593</b> | No new line after                               |
| <b>fprintf(' x = %f \n',pi)</b>       | <b>x = 3.141593</b> | Moves to a new line after                       |
| <b>fprintf(' L%6.2fR \n',pi)</b>      | <b>L 3.14R</b>      | Number is 6 chars wide with 2 dp                |
| <b>fprintf('%s L%-6.2fR','Hi',pi)</b> | <b>Hi L3.14 R</b>   | The string 'Hi', then left justify number       |
| <b>S = sprintf(' x = %f ',pi);</b>    | <b>NOTHING</b>      | <b>S</b> is a string containing ' x = 3.141593' |

## Graph Commands

|                                    |                                                        |
|------------------------------------|--------------------------------------------------------|
| <b>plot(y)</b>                     | Plot y against index number.                           |
| <b>plot(x,y)</b>                   | Plot y against x                                       |
| <b>plot(x1,y1,x2,y2)</b>           | Plot y1 against x1 and y2 against x2.                  |
| <b>plot(x,y,'r+')</b>              | Plot y against x using red plus signs.                 |
| <b>plot(x1,y1,'r+',x2,y2,'go')</b> | Red plus signs for x1 and y1, Green circles for x2 y2. |

| Symbol | Line Type or Mark | Symbol | Colour  |
|--------|-------------------|--------|---------|
| .      | Point             | r      | Red     |
| o      | Circle            | g      | Green   |
| x      | X mark            | b      | Blue    |
| +      | Plus sign         | y      | Yellow  |
| *      | Stars             | m      | Magenta |
| -      | Solid line        | c      | Cyan    |
| :      | Dotted line       | w      | White   |
| .-     | Dash dot line     | k      | Black   |
| --     | Dash Line         |        |         |

| Other Types of Plot   |                  |
|-----------------------|------------------|
| <b>fill(x,y,'r')</b>  | Red filled graph |
| <b>bar(x,y)</b>       | Bar graph        |
| <b>loglog(x,y)</b>    | x & y log scale  |
| <b>semilogx(x,y)</b>  | x log, y linear  |
| <b>semilogy(x,y)</b>  | x linear, y log  |
| <b>polar(theta,r)</b> | Polar plot       |
| <b>surf(x,y,z)</b>    | 3D surface       |
| <b>mesh(x,y,z)</b>    | 3D mesh          |
| <b>plot3(x,y,z)</b>   | 3D line plot     |

## Other Graphics Commands

|                                                 |                                                                 |                |                |
|-------------------------------------------------|-----------------------------------------------------------------|----------------|----------------|
| <b>title('Title')</b>                           | Graph title.                                                    |                |                |
| <b>xlabel('X axis')</b>                         | Label the x-axis.                                               |                |                |
| <b>ylabel('Y axis')</b> <b>zlabel('Z axis')</b> | Label y and z axis.                                             |                |                |
| <b>text(x,y,'My Text')</b>                      | Place text at coordinates x,y.                                  |                |                |
| <b>grid</b>                                     | Place a grid on the graph.                                      |                |                |
| <b>hold on</b>                                  | Add any new plot to the current graph.                          |                |                |
| <b>hold off</b>                                 | Replace current plot with any new plot.                         |                |                |
| <b>subplot(r,c,n)</b>                           | Split figure into r rows by c columns of subplots.              |                |                |
|                                                 | subplot(2,3,1)                                                  | subplot(2,3,2) | subplot(2,3,3) |
|                                                 | subplot(2,3,4)                                                  | subplot(2,3,5) | subplot(2,3,6) |
| <b>axis([minX maxX minY maxY])</b>              | Set the limits of the graph in X and Y. (Note 4 element vector) |                |                |
| <b>h = figure</b>                               | New graphics window.                                            |                |                |
| <b>figure(h)</b>                                | Change to plotting in figure h.                                 |                |                |
| <b>delete(h)</b>                                | Delete figure h.                                                |                |                |
| <b>clf</b>                                      | Clear current figure.                                           |                |                |
| <b>drawnow</b>                                  | Force the graph to update now.                                  |                |                |

# Programming

## Enumerated Loops (for)

The general form of a **for** loop is :-

```
for <variable> = <vector>
    <statement>
    <statement>
    etc
end
```

Examples

```
A = rand(1,5);
for a = A
    disp(a)
end
```

```
for k = [ 1 7 3 pi i]
    disp(k)
end
```

## Precondition Loops (while)

Below is the general form of a **while** loop.

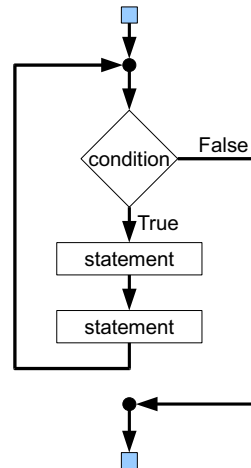
```
while(<condition>)
    <statement>
    <statement>
    etc
end
```

Example

```
A = 7;      %Find the square root of A.
x=1;       %First guess
err=1;     %Set error to get started

% Newton Raphson Iteration
while(err>0.0001)
    x = (x.^2+A)./(2*x);
    % calculate the error
    err = abs(A-x.^2);
end

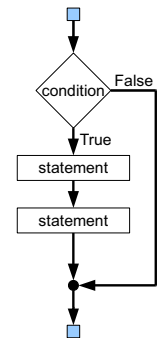
disp(x);
```



## Conditional Execution (if)

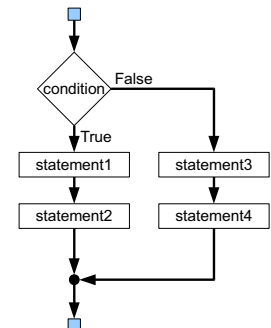
General form of a simple **if** statement.

```
if <condition>
    <statement>
    <statement>
    etc
end
```



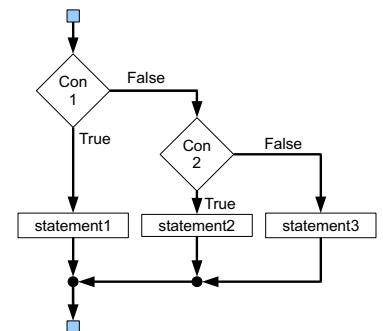
The form of **if** statement with **else**

```
if <condition>
    <statement1>
    <statement2>
    etc
else
    <statement3>
    <statement4>
    etc
end
```



You can also have as many **elseif** parts as you like.

```
if <condition1>
    <statement1>
    etc
elseif <condition2>
    <statement2>
    etc
else
    <statement3>
    etc
end
```



# Switch

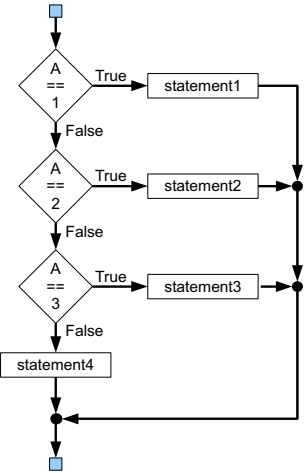
Execution depends on the value of a variable.

```
switch <variable>
case 1
  <statement1>
  etc
case 2
  <statement2>
  etc
case 3
  <statement3>
  etc
otherwise
  <statement4>
  etc
end
```

The case values can be any value that the switch variable can take. You can also put multiple values after the case.

```
case {1,2,3,4,5}
  <statement>
  etc
```

On the right A is the switch variable.



# Functions

| Function Definition                                                                                                                                                      | Using the Function                                                |
|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-------------------------------------------------------------------|
| <pre>function sum = myadd(a,b)  % The first block of comments % defines what is printed out when % you type    help myadd  % Comment not part of help sum = a + b;</pre> | <pre>s = myadd(&lt;expression&gt;,&lt;expression&gt;)</pre>       |
| <pre>function [sum,diff] = add_sub(a,b)  sum = a + b; diff = a - b;</pre>                                                                                                | <pre>[s,d] = add_sub(&lt;expression&gt;,&lt;expression&gt;)</pre> |

# Introduction à Matlab

## 1 Introduction

Matlab (*MATrix LABoratory*) est un logiciel pour le calcul scientifique, orienté vers le vecteurs et les listes de données. Matlab est un langage interprété, chaque ligne d'un programme Matlab est lue, interprétée et exécutée.

Pour entrer dans Matlab cliquez sur **Start/Programmes/Matlab** sous l'environnement Windows, ou tapez **matlab** au prompt de la shell sous Unix. Matlab se présente avec un environnement interactif et un prompt (généralement **>>**) dans lequel on peut introduire des *commandes*. Par exemple, pour sortir de Matlab tapez la commande:

```
>> quit
```

## 2 Aide en ligne

Matlab offre une aide en ligne très complète. Tapez:

```
>> help commande
```

pour une explication de l'utilisation de la commande Matlab **commande**

```
>> lookfor sujet
```

pour une liste des commandes Matlab qui concernent le sujet **sujet**

```
>> helpwin
```

pour ouvrir un guide des commandes Matlab sur différents sujets.

## 3 Déclaration des variables

Matlab permet de créer et d'initialiser des variables. La déclaration des variables en Matlab suit les règles suivantes:

- toutes les variables sont des *matrices*
- pas de déclaration de type

```
>> a=5 variable scalaire (1 × 1)
```

```
>> b=[4 6] vecteur ligne (1 × 2)
```

```
>> c=[-5; 2] vecteur colonne (2 × 1)
```

```
>> d=[2 3; -1 7] matrice carrée (2 × 2)
```

Dans ces exemples on a utilisé les opérateurs suivants:

- séparateur de ligne: point virgule ou return
- séparateur de colonne: virgule ou espace blanc

Il faut noter que Matlab montre les nombres avec quatre chiffres après la virgule, tandis que la représentation interne est faite avec 16 chiffres. Pour changer la façon de montrer les nombres en Matlab, on utilise la commande **format**. Par exemple, si avant de taper **>> pi** (en Matlab la variable **pi** est une approximation de  $\pi$ ), nous écrivons

```
>> format long nous obtiendrons 3.1416;
```

```
>> format short nous obtiendrons 3.14159265358979;
```

```
>> format long e nous obtiendrons 3.141592653589793e+00;
```

```
>> format short e nous obtiendrons 3.1416e+00.
```

## 4 Workspace

Matlab permet de connaître plusieurs informations sur les variables déclarées. Tapez:

```
>> who           pour afficher toutes les variables.

>> whos         pour afficher toutes les variables
                  avec indication sur leur taille.

>> size(a)       pour afficher les dimensions de la
                  matrice a.

>> clear var     pour effacer la variable var.

>> clear (ou >> clear all) pour effacer toutes les variables.
```

Pour ces opérations vous pouvez aussi utiliser l'interface graphique de Matlab.

## 5 Opérations fondamentales

Matlab peut effectuer plusieurs opérations entre matrices. Les opérations fondamentales peuvent être partagées en deux catégories.

### Opérations matricielles

Les opérations matricielles usuelles sont définies par  $+$ ,  $-$ ,  $*$ ,  $/$ ,  $^$

```
>> C = A + B      est la somme matricielle,  $C_{ij} = A_{ij} + B_{ij}$ 

>> C = A * B      est le produit matriciel,  $C_{ij} = \sum_k A_{ik} B_{kj}$ 

>> C = A / B      est la division matricielle,  $C = AB^{-1}$ 

>> C = A^3        est la troisième puissance matricielle (C = A*A*A)
```

Il faut remarquer que ces opérations sont bien définies seulement si les matrices ont des *dimensions cohérentes*. Pour l'addition  $A+B$ ,  $A$  et  $B$  doivent avoir la même taille; pour le produit  $A*B$ , le nombre des colonnes de  $A$  et le nombre des lignes de  $B$  doivent être égaux; les opérations  $A/B$  et  $B^3$  demandent une matrice  $B$  carrée. Par exemple:

```
>> A = [1 2 3; 4 5 6]; B = [7 8 9; 10 11 12]; C = [13 14; 15 16; 17 18];
>> A + B
ans =

     8     10     12
    14     16     18

>> A + C
```

```
??? Error using ==> +
Matrix dimensions must agree.
```

```
>> A * C
ans =

     94     100
    229     244
```

```
>> A * B
??? Error using ==> *
Inner matrix dimensions must agree.
```

Notez que Matlab renvoie un message d'erreur si le dimensions des matrices ne s'accordent pas avec l'opération commandée.

### Opérations élément par élément

Pour exécuter des opérations entre matrices élément par élément il faut faire précéder l'opérateur d'un point. Les opérateurs élément par élément sont donc  $.*$ ,  $./$ ,  $.^$

```
>> C = A .* B      est le produit élément par élément,  $C_{ij} = A_{ij} B_{ij}$ 

>> C = A ./ B      est la division élément par élément,  $C_{ij} = \frac{A_{ij}}{B_{ij}}$ 

>> C = A.^3        est la troisième puissance élément par élément,  $C_{ij} = A_{ij}^3$ 
```

## 6 Manipulation des matrices

Après avoir défini la matrice  $A$  (par exemple la matrice carrée  $A$  de taille  $n$ ), Matlab permet les opérations suivantes sur les *sous-matrices* de  $A$ .

### Extraction de sous-matrices

```
>> A(2,3)          extraction de l'élément  $A_{23}$ 

>> A(:,5)          extraction de la colonne  $[A_{13}; \dots; A_{n3}]$ 

>> A(1:4,3)        extraction de la sous-colonne  $[A_{13}; \dots; A_{43}]$ 

>> A(1,:)          extraction de la ligne  $[A_{1j}, \dots, A_{1n}]$ 

>> diag(A)         extraction de la diagonale  $[A_{11}; \dots; A_{nn}]$ 
```

### Construction de matrices particulières

Matlab permet, en plus, de définir des matrices particulières (une fois définis les entiers **n**, **m** et le vecteur **v**).

```
>> A = eye(n)           matrice identité  $n \times n$ 

>> A = diag(v)          matrice diagonale avec v comme diagonale

>> A = zeros(n,m)       matrice de zéros avec n lignes et m colonnes

>> A = ones(n,m)        matrice de un avec n lignes et m colonnes

>> A = rand(n,m)        matrice aléatoire avec n lignes et m colonnes

>> x = [but:pas:fin]     vecteur ligne de points équidistribués
                        avec pas pas entre but et fin
```

### Fonctions matricielles

Soit **A** une matrice, Matlab permet d'effectuer directement les opérations suivantes.

```
>> C = A'               transposée de A,  $C_{ij} = A_{ji}$ 

>> C = inv(A)           inverse de A (matrice carrée),  $C = A^{-1}$ 

>> d = det(A)           déterminant de A (matrice carrée)

>> r = rank(A)          rang de A

>> nrm = norm(A)        norme 2 de A

>> v = eig(A)           valeurs propres de A (matrice carrée)
```

Soit **A** une matrice carrée de taille **n** et soit **b** un vecteur de taille **n**, alors le vecteur **x**, solution du système linéaire **Ax = b** est défini par

```
>> x = A^-1 * b
```

Cette commande est théoriquement correcte, cependant si on est intéressé seulement à la solution **x** et pas à l'inverse **A^-1**, il est préférable d'utiliser un *backslash*

```
>> x = A \ b
```

Cette commande résout le système linéaire avec des algorithmes fiables et optimaux.

### 7 Graphisme 2D

Matlab offre plusieurs possibilités pour tracer un graphe en 2D. On va en présenter deux: la commande **plot** et la commande **fplot**.

Avant d'expliquer ces deux commandes en détail, on souligne qu'avec **plot** on doit toujours utiliser des vecteurs, alors qu'avec **fplot** non.

On considère maintenant la fonction  $f(x) = x^3 - 2 \sin x + 1$  et on voit comment on peut tracer son graphe dans l'intervalle  $[-1, 1]$ , en utilisant les commandes **plot** et **fplot**.

#### Commande plot

Pour tracer le graphe de  $f(x)$  il faut passer par les étapes suivantes :

- Définir la fonction  $f(x)$  :

```
>> f = 'x.^3 - 2*sin(x) + 1';
```

le mot **f** contient la définition de la fonction choisie (les guillemets font partie de la définition dans la syntaxe Matlab).

- Définir un *vecteur de points* dans l'intervalle donné:

```
>> x = [-1:0.1:+1];
```

on a défini un vecteur de 21 points equidistribués dans l'intervalle donné, avec un pas de 0.1.

- Évaluer la fonction **f** dans l'intervalle  $[-1, 1]$ , en utilisant la commande **eval**:

```
>> y = eval(f);
```

on note que dans la définition de la fonction  $f$  on a dû écrire  $.^$  (puissance élément par élément) et pas  $^$  (puissance matricielle), parce qu'il faut évaluer la fonction en chaque élément du *vecteur* **x**. C'est-à-dire:

$$y(i) = x(i)^3 - 2 \sin(x(i)) + 1, \quad \text{pour } i = 1, \dots, 21;$$

La variable **y** est donc un *vecteur* qui contient les valeurs de **f** pour chaque entrée du vecteur **x**.

- Tracer le graphe :

```
>> plot(x,y); grid;
```

cette commande ouvre une fenêtre avec le graphe. La commande **grid** dessine une grille de repère.

Le résultat de ces commandes est représenté dans la Figure 1.



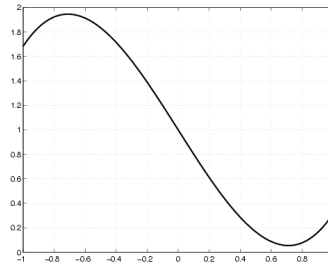


Figure 1: Graphe de la fonction  $f(x) = x^3 - 2\sin(x) + 1$

### Commande `fplot`

Pour tracer le graphe avec la commande `fplot` il faut passer par les étapes suivantes :

- Définir la fonction  $f(x)$  :

```
>> f = 'x^3 - 2*sin(x) + 1';
```

- Tracer le graphe :

```
>> fplot(f, [-1, 1]); grid;
```

La commande `fplot` demande la définition de la fonction `f` et de l'intervalle où il faut la tracer. La définition des vecteurs `x` et `y` n'est plus nécessaire et on note que dans la définition de `f` il suffit d'écrire `f='x^3 - ...'` au lieu de `f = x.^3 - ...'`, comme on avait fait pour utiliser `plot`. Le graphe obtenu est encore montré dans la Figure 1.

**Exemple** Tracer le graphe des fonctions  $f_1(x) = \sin(2x) - 1 + x$  et  $f_2(x) = x^3 - 2\sin(x)$  dans l'intervalle  $[-1, 1]$  sur la même fenêtre.

En utilisant la commande `fplot`, il faut procéder de la façon suivante:

- Définir les fonctions:

```
>> f1 = 'sin(2*x) - 1 + x';
>> f2 = 'x^3 - 2*sin(x)';
```

- Tracer le graphe:

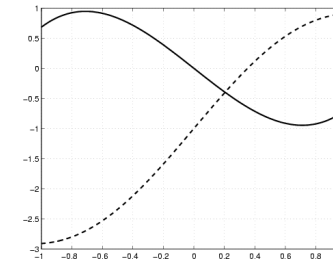


Figure 2: Résultat de l'exemple

```
>> fplot(f1, [-1, 1]); grid; hold on;
```

La commande `hold on` affiche tous les prochains graphes sur la fenêtre qu'on vient d'ouvrir. Donc le graphe:

```
>> fplot(f2, [-1, 1]);
```

est affiché sur la même fenêtre, comme on peut le voir dans la Figure 2.

La commande `hold off` arrête cet effet.

Pour plus de détails sur les fonctions `plot` et `fplot` (par exemple comment changer la couleur du graphe) tapez `help plot` ou `help fplot`.

## 8 Boucles de contrôle

Matlab offre, comme plusieurs autres langages, quelques boucles de contrôle. Ces commandes sont spécialement utiles pour écrire des programmes Matlab.

### Boucle `for`

Si l'on veut exécuter des instructions pour chaque valeur

$$i = m, m + 1, \dots, n$$

d'une certaine variable  $i$  entre deux bornes assignés  $m$  et  $n$ , on utilise `for`. Par exemple pour calculer le produit scalaire `ps` entre deux vecteurs `x` et `y` de taille `n` on utilise:

```
>> ps = 0;
>> for i = 1:n;
>>     ps = ps + x(i)*y(i);
>> end;
```

Cette boucle est équivalente au le produit matriciel entre le vecteur transposé de **x** et le vecteur **y**, en supposant que les facteurs soient deux vecteurs colonne:

```
>> ps = x'*y;
```

**Boucle while**

Si l'on veut exécuter plusieurs fois des instructions tandis qu'une expression logique est vraie, on utilise **while**. Par exemple, le même calcul qu'on a considéré avec la boucle **for** peut être exécuté avec

```
>> ps = 0;
>> i = 0;
>> while (i < n);
>>     i = i + 1;
>>     ps = ps + x(i)*y(i);
>> end;
```

**Instruction conditionnelle if**

Si l'on veut exécuter des instructions seulement si une expression logique est vraie, on utilise **if**. Par exemple, si l'on veut calculer la racine carrée d'une variable scalaire **r** seulement si **r** n'est pas négative:

```
>> if (r >= 0)
>>     racine = sqrt(r);
>> end;
```

On a plusieurs *opérateurs logiques* à disposition:

| Opérateur | Action logique |
|-----------|----------------|
| &         | and            |
|           | or             |
| ~         | not            |
| ==        | equal to       |

Taper par exemple **help &** pour une explication de la liste des opérateurs.

**9 Scripts et fonctions**

Matlab permet d'écrire des programmes. Les outils principaux sont les script files et les fonctions.

**Script files**

Un script file est une suite de commandes Matlab. Les noms des script files doivent avoir l'extension **“.m”**. Pour exécuter un script file tapez son nom, sans extension, dans le prompt Matlab.

**Fonctions**

Une fonction Matlab est une suite de commandes qui nécessite un input pour être exécutée et qui renvoie un output. La déclaration des fonctions en Matlab suit les règles suivantes.

- Une fonction est contenue dans un fichier **“.m”** avec le même nom que la fonction.
- Le fichier qui définit une fonction doit commencer par:  
`function [output arguments] = nom_fonction(input arguments)`  
Par exemple, l'en-tête d'une fonction pour la méthode de dichotomie pourrait être:

```
function [zero,res,niter]=bisection(fun,a,b,tol,nmax,varargin)
%BISECTION Find function zeros.
%   ZERO=BISECTION(FUN,A,B,TOL,NMAX) tries to find a zero ZERO
%   of the continuous function FUN in the interval [A,B]
%   using the bisection method.
.
.
.   % instructions
.
zero = ...; res = ...; niter = ...;
.
.   % instructions
.
.
return           % fin du corps de la fonction
```

- Les lignes de commentaires qui éventuellement suivent cet en-tête constituent le help de la fonction. En Matlab les lignes de commentaire sont introduites par **%**.
- Toutes les variables internes à la fonction sont locales.

Par exemple, supposons qu'on aie écrit sur le fichier **my\_function.m** le lignes suivantes:

```
function f = my_function(x);
f = x.^3 - 2*sin(x) + 1;
return;
```

Il est possible d'utiliser la fonction `my_function.m` de la même façon que les autres commandes Matlab. Donc les commandes

```
>> x = 0;
>> y = x.^3 - 2*sin(x) + 1
```

et

```
>> x = 0;
>> y = my_function(x)
```

sont équivalentes et renvoient le même résultat:

```
y =
    1
```

### Fonctions de plusieurs variables

Pour définir des fonctions dépendantes d'un ou plusieurs paramètres nous pouvons utiliser les fonctions *inline*. Pour définir une fonction *inline* `f` qui dépend de la variable `x` on écrit:

```
>> f = inline('log(x) - 1','x');
```

Si `f` dépend de de plusieurs variables, nous écrirons par exemple

```
>> f = inline('log(x) - p','x','p');
```

Pour évaluer une fonction `f` définie grâce à *inline*, il faut utiliser la commande `feval` en donnant les valeurs des paramètres comme montré dans l'exemple suivant:

```
>> f = inline('x.^3 + a.*x','x','a');
>> xval = 1; aval = 2;
>> fval = feval(f, xval, yval)
fval =
    3
```

## 10 Polynômes en Matlab

Considérons un polynôme de degré  $n$ :

$$f(x) = a_n x^n + a_{n-1} x^{n-1} + \dots + a_1 x + a_0.$$

Matlab représente les polynômes de degré  $n$  sous la forme d'un vecteur  $p = [a_n, a_{n-1}, \dots, a_0]$  de taille  $n+1$  qui contient les coefficients en ordre décroissant par rapport au degré associé. Par exemple, le vecteur associé au polynôme  $f(x) = 3x^3 - 4x^2 + x$  est

```
>> p = [3, -4, 1, 0];
```

| Mois | Heure de son réveil (moyenne) | Mois | Heure de son réveil (moyenne) |
|------|-------------------------------|------|-------------------------------|
| 1    | 8.0                           | 7    | 8.0                           |
| 2    | 7.5                           | 8    | 10.0                          |
| 3    | 7.5                           | 9    | 8.0                           |
| 4    | 7.5                           | 10   | 7.5                           |
| 5    | 7.6                           | 11   | 7.5                           |
| 6    | 7.8                           | 12   | 7.7                           |

Table 1: Heure de son réveil (de 0 à 24 avec décimes d'heure)

Pour évaluer ce polynôme on utilise la commande `polyval`:

```
>> x = [0:.1:1];
>> y = polyval(p, x);
```

Dans ce cas, les éléments du vecteur `y` sont les valeurs du polynôme calculées pour chaque élément du vecteur `x`.

Pour obtenir le vecteur associé au polynôme de degré  $n$  approximant un jeu de données on utilise `polyfit`. Si la taille des données est plus grande de  $n+1$  on a le polynôme approximant au sens des moindres carrés; si elle est égal à  $n+1$  on a le polynôme interpolant. Par exemple, si on veut calculer le polynôme de degré 8 qui approxime le données de la Table 10 (au sens des moindre carrées) et en tracer le graphe, il faut utiliser les commandes qui suivent;

```
x = [1 2 3 4 5 6 7 8 9 10 11 12];
y = [8 7.5 7.5 7.5 7.6 7.8 8 10 8 7.5 7.5 7.7];
p = polyfit(x,y,8); % Calculons le polynome interpolant (degre=8)
plot(x, y, 'or');
axis([1 12 6 12])
hold on;
f = inline('polyval(p,x)','x','p');
fplot(f, [1 12], [], [], [], p); % Trace le graphe de f(x,p)
                                % pour x entre 0 et 12, en donnant
                                % p comme parametre de f.
```

Il faut noter que `f` est déclarée comme fonction *inline* avec les arguments `x` et `p` (le vecteur associé au polynôme interpolant); la commande `fplot` (voir `help fplot`) accepte le paramètre `p` à donner à `f`.

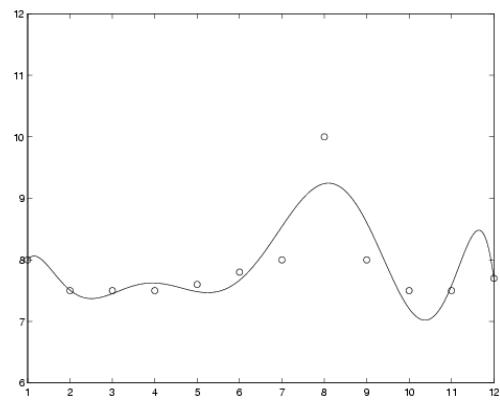


Figure 3: Données et polynôme d'interpolation

# MATLAB Primer

## Third Edition

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### ON THE THIRD EDITION

The Third Edition of the MATLAB Primer is based on version 4.0/4.1 of MATLAB. While this edition reflects an extensive general revision of the Second Edition, most significant is the new information to help one begin to use the major new features of version 4.0/4.1, the sparse matrix and enhanced graphics capabilities.

The plain TeX source and corresponding PostScript file of the latest printing of the MATLAB Primer are always available via anonymous ftp from:

Address: math.ufl.edu    Directory: pub/matlab    Files: primer.tex, primer.ps

You are advised to download anew each term the latest printing of the Primer since minor improvements and corrections may have been made in the interim. If ftp is unavailable to you, the Primer can be obtained via `listserv` by sending an email message to `listserv@math.ufl.edu` containing the single line `send matlab/primer.tex`.

Also available at this ftp site are both English (`primer35.tex`, `primer35.ps`) and Spanish (`primer35sp.tex`, `primer35sp.ps`) versions of the Second Edition of the Primer, which was based on version 3.5 of MATLAB. The Spanish translation is by Celestino Montes, University of Seville, Spain. A Spanish translation of the Third Edition is under development.

Users of the Primer usually appreciate the convenience and durability of a bound copy with a cover, copy center style.

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## INTRODUCTION

MATLAB is an interactive, matrix-based system for scientific and engineering numeric computation and visualization. You can solve complex numerical problems in a fraction of the time required with a programming language such as Fortran or C. The name MATLAB is derived from MATrix LABoratory.

The purpose of this Primer is to help you begin to use MATLAB. It is not intended to be a substitute for the User's Guide and Reference Guide for MATLAB. The Primer can best be used hands-on. You are encouraged to work at the computer as you read the Primer and freely experiment with examples. This Primer, along with the on-line help facility, usually suffice for students in a class requiring use of MATLAB.

You should liberally use the on-line help facility for more detailed information. When using MATLAB, the command `help functionname` will give information about a specific function. For example, the command `help eig` will give information about the eigenvalue function `eig`. By itself, the command `help` will display a list of topics for which on-line help is available; then `help topic` will list those specific functions under this topic for which help is available. The list of functions in the last section of this Primer also gives most of this information. You can preview some of the features of MATLAB by first entering the command `demo` and then selecting from the options offered.

The scope and power of MATLAB go far beyond these notes. Eventually you will want to consult the MATLAB User's Guide and Reference Guide. Copies of the complete documentation are often available for review at locations such as consulting desks, terminal rooms, computing labs, and the reserve desk of the library. Consult your instructor or your local computing center to learn where this documentation is located at your institution.

MATLAB is available for a number of environments: Sun/Apollo/VAXstation/HP workstations, VAX, MicroVAX, Gould, PC and AT compatibles, 80386 and 80486 computers, Apple Macintosh, and several parallel machines. There is a relatively inexpensive Student Edition available from Prentice Hall publishers. The information in these notes applies generally to all of these environments.

MATLAB is licensed by The MathWorks, Inc., 24 Prime Park Way, Natick, MA 01760, (508)653-1415, Fax: (508)653-2997, Email: [info@mathworks.com](mailto:info@mathworks.com).

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## 1. Accessing MATLAB.

On most systems, after logging in one can enter MATLAB with the system command `matlab` and exit MATLAB with the MATLAB command `quit` or `exit`. However, your local installation may permit MATLAB to be accessed from a menu or by clicking an icon.

On systems permitting multiple processes, such as a Unix system or MS Windows, you will find it convenient, for reasons discussed in section 14, to keep both MATLAB and your local editor active. If you are working on a platform which runs processes in multiple windows, you will want to keep MATLAB active in one window and your local editor active in another.

You should consult your instructor or your local computer center for details of the local installation.

## 2. Entering matrices.

MATLAB works with essentially only one kind of object—a rectangular numerical matrix with possibly complex entries; all variables represent matrices. In some situations, 1-by-1 matrices are interpreted as scalars and matrices with only one row or one column are interpreted as vectors.

Matrices can be introduced into MATLAB in several different ways:

- Entered by an explicit list of elements,
- Generated by built-in statements and functions,
- Created in a diskfile with your local editor,
- Loaded from external data files or applications (see the User's Guide).

For example, either of the statements

```
A = [1 2 3; 4 5 6; 7 8 9]
```

and

```
A = [  
1 2 3  
4 5 6  
7 8 9 ]
```

creates the obvious 3-by-3 matrix and assigns it to a variable *A*. Try it. The elements within a row of a matrix may be separated by commas as well as a blank. When listing a number in exponential form (e.g. 2.34e-9), blank spaces must be avoided.

MATLAB allows complex numbers in all its operations and functions. Two convenient ways to enter complex matrices are:

```
A = [1 2;3 4] + i*[5 6;7 8]  
A = [1+5i 2+6i;3+7i 4+8i]
```

When listing complex numbers (e.g. 2+6i) in a matrix, blank spaces must be avoided. Either *i* or *j* may be used as the imaginary unit. If, however, you use *i* and *j* as variables and overwrite their values, you may generate a new imaginary unit with, say, `ii = sqrt(-1)`.

Listing entries of a large matrix is best done in an ASCII file with your local editor, where errors can be easily corrected (see sections 12 and 14). The file should consist of a rectangular array of just the numeric matrix entries. If this file is named, say, `data.ext` (where `.ext` is any extension), the MATLAB command `load data.ext` will read this file to the variable `data` in your MATLAB workspace. This may also be done with a script file (see section 12).

The built-in functions `rand`, `magic`, and `hilb`, for example, provide an easy way to create matrices with which to experiment. The command `rand(n)` will create an  $n \times n$  matrix with randomly generated entries distributed uniformly between 0 and 1, while `rand(m,n)` will create an  $m \times n$  one. `magic(n)` will create an integral  $n \times n$  matrix which is a magic square (rows, columns, and diagonals have common sum); `hilb(n)` will create the  $n \times n$  Hilbert matrix, the king of ill-conditioned matrices ( $m$  and  $n$  denote, of course, positive integers). Matrices can also be generated with a for-loop (see section 6 below).

Individual matrix and vector entries can be referenced with indices inside parentheses in the usual manner. For example, `A(2,3)` denotes the entry in the second row, third column of matrix *A* and `x(3)` denotes the third coordinate of vector *x*. Try it. A matrix or a vector will only accept *positive* integers as indices.

## 3. Matrix operations, array operations.

The following matrix operations are available in MATLAB:

|   |                     |
|---|---------------------|
| + | addition            |
| - | subtraction         |
| * | multiplication      |
| ^ | power               |
| ' | conjugate transpose |
| \ | left division       |
| / | right division      |

These matrix operations apply, of course, to scalars (1-by-1 matrices) as well. If the sizes of the matrices are incompatible for the matrix operation, an error message will result, except in the case of scalar-matrix operations (for addition, subtraction, and division as well as for multiplication) in which case each entry of the matrix is operated on by the scalar.

The “matrix division” operations deserve special comment. If *A* is an invertible square matrix and *b* is a compatible column, resp. row, vector, then

$x = A \backslash b$  is the solution of  $A * x = b$  and, resp.,  
 $x = b / A$  is the solution of  $x * A = b$ .

In left division, if *A* is square, then it is factored using Gaussian elimination and these factors are used to solve  $A * x = b$ . If *A* is not square, it is factored using Householder orthogonalization with column pivoting and the factors are used to solve the under- or over- determined system in the least squares sense. Right division is defined in terms of left division by  $b / A = (A' \backslash b)'$ .

### Array operations.

The matrix operations of addition and subtraction already operate entry-wise but the other matrix operations given above do not—they are *matrix* operations. It is important to observe that these other operations, `*`, `^`, `\`, and `/`, can be made to operate entry-wise by preceding them by a period. For example, either `[1,2,3,4].*[1,2,3,4]` or `[1,2,3,4].^2` will yield `[1,4,9,16]`. Try it. This is particularly useful when using Matlab graphics.

### 4. Statements, expressions, and variables; saving a session.

MATLAB is an *expression* language; the expressions you type are interpreted and evaluated. MATLAB statements are usually of the form

*variable* = *expression*, or simply  
*expression*

Expressions are usually composed from operators, functions, and variable names. Evaluation of the expression produces a matrix, which is then displayed on the screen and assigned to the variable for future use. If the variable name and = sign are omitted, a variable **ans** (for answer) is automatically created to which the result is assigned.

A statement is normally terminated with the carriage return. However, a statement can be continued to the next line with three or more periods followed by a carriage return. On the other hand, several statements can be placed on a single line if separated by commas or semicolons.

If the last character of a statement is a semicolon, the printing is suppressed, but the assignment is carried out. This is essential in suppressing unwanted printing of intermediate results.

MATLAB is case-sensitive in the names of commands, functions, and variables. For example, **solveUT** is not the same as **solveut**.

The command **who** (or **whos**) will list the variables currently in the workspace. A variable can be cleared from the workspace with the command **clear** *variablename*. The command **clear** alone will clear all nonpermanent variables.

The permanent variable **eps** (epsilon) gives the machine unit roundoff—about  $10^{-16}$  on most machines. It is useful in specifying tolerances for convergence of iterative processes.

A runaway display or computation can be stopped on most machines without leaving MATLAB with CTRL-C (CTRL-BREAK on a PC).

### Saving a session.

When one logs out or exits MATLAB all variables are lost. However, invoking the command **save** before exiting causes all variables to be written to a non-human-readable diskfile named **matlab.mat**. When one later reenters MATLAB, the command **load** will restore the workspace to its former state.

### 5. Matrix building functions.

Convenient matrix building functions are

|                       |                                       |
|-----------------------|---------------------------------------|
| <code>eye</code>      | identity matrix                       |
| <code>zeros</code>    | matrix of zeros                       |
| <code>ones</code>     | matrix of ones                        |
| <code>diag</code>     | create or extract diagonals           |
| <code>triu</code>     | upper triangular part of a matrix     |
| <code>tril</code>     | lower triangular part of a matrix     |
| <code>rand</code>     | randomly generated matrix             |
| <code>hilb</code>     | Hilbert matrix                        |
| <code>magic</code>    | magic square                          |
| <code>toeplitz</code> | see <b>help</b> <code>toeplitz</code> |

For example, **zeros(m,n)** produces an *m*-by-*n* matrix of zeros and **zeros(n)** produces an *n*-by-*n* one. If *A* is a matrix, then **zeros(size(A))** produces a matrix of zeros having the same size as *A*.

If *x* is a vector, **diag(x)** is the diagonal matrix with *x* down the diagonal; if *A* is a square matrix, then **diag(A)** is a vector consisting of the diagonal of *A*. What is **diag(diag(A))**? Try it.

Matrices can be built from blocks. For example, if *A* is a 3-by-3 matrix, then

**B = [A, zeros(3,2); zeros(2,3), eye(2)]**

will build a certain 5-by-5 matrix. Try it.

### 6. For, while, if — and relations.

In their basic forms, these MATLAB flow control statements operate like those in most computer languages.

#### For.

For example, for a given *n*, the statement

**x = []; for i = 1:n, x=[x,i^2], end**

or

**x = [];**  
**for i = 1:n**  
    **x = [x,i^2]**  
**end**

will produce a certain *n*-vector and the statement

**x = []; for i = n:-1:1, x=[x,i^2], end**

will produce the same vector in reverse order. Try them. Note that a matrix may be empty (such as **x = []**).



The statements

```
for i = 1:m
    for j = 1:n
        H(i, j) = 1/(i+j-1);
    end
end
H
```

will produce and print to the screen the  $m$ -by- $n$  hilbert matrix. The semicolon on the inner statement is essential to suppress printing of unwanted intermediate results while the last `H` displays the final result.

The `for` statement permits *any* matrix to be used instead of `1:n`. The variable just consecutively assumes the value of each column of the matrix. For example,

```
s = 0;
for c = A
    s = s + sum(c);
end
```

computes the sum of all entries of the matrix  $A$  by adding its column sums (Of course, `sum(sum(A))` does it more efficiently; see section 8). In fact, since `1:n = [1,2,3,...,n]`, this column-by-column assignment is what occurs with “`if i = 1:n,...`” (see section 11).

#### While.

The general form of a `while` loop is

```
while relation
    statements
end
```

The statements will be repeatedly executed as long as the relation remains true. For example, for a given number  $a$ , the following will compute and display the smallest nonnegative integer  $n$  such that  $2^n \geq a$ :

```
n = 0;
while 2^n < a
    n = n + 1;
end
n
```

#### If.

The general form of a simple `if` statement is

```
if relation
    statements
end
```

The statements will be executed only if the relation is true. Multiple branching is also possible, as is illustrated by

```
if n < 0
    parity = 0;
```

```
elseif rem(n,2) == 0
    parity = 2;
else
    parity = 1;
end
```

In two-way branching the `elseif` portion would, of course, be omitted.

#### Relations.

The relational operators in MATLAB are

|    |                       |
|----|-----------------------|
| <  | less than             |
| >  | greater than          |
| <= | less than or equal    |
| >= | greater than or equal |
| == | equal                 |
| ~= | not equal.            |

Note that “`=`” is used in an assignment statement while “`==`” is used in a relation. Relations may be connected or quantified by the logical operators

|   |      |
|---|------|
| & | and  |
|   | or   |
| ~ | not. |

When applied to scalars, a relation is actually the scalar 1 or 0 depending on whether the relation is true or false. Try entering `3 < 5`, `3 > 5`, `3 == 5`, and `3 == 3`. When applied to matrices of the same size, a relation is a matrix of 0's and 1's giving the value of the relation between corresponding entries. Try `a = rand(5)`, `b = triu(a)`, `a == b`.

A relation between matrices is interpreted by `while` and `if` to be true if each entry of the relation matrix is nonzero. Hence, if you wish to execute *statement* when matrices  $A$  and  $B$  are equal you could type

```
if A == B
    statement
end
```

but if you wish to execute *statement* when  $A$  and  $B$  are not equal, you would type

```
if any(any(A ~= B))
    statement
end
```

or, more simply,

```
if A == B else
    statement
end
```

Note that the seemingly obvious

```
if A ~= B, statement, end
```

will not give what is intended since *statement* would execute only if *each* of the corresponding entries of *A* and *B* differ. The functions **any** and **all** can be creatively used to reduce matrix relations to vectors or scalars. Two **any**'s are required above since **any** is a vector operator (see section 8).

## 7. Scalar functions.

Certain MATLAB functions operate essentially on scalars, but operate element-wise when applied to a matrix. The most common such functions are

|     |      |                   |      |       |
|-----|------|-------------------|------|-------|
| sin | asin | exp               | abs  | round |
| cos | acos | log (natural log) | sqrt | floor |
| tan | atan | rem (remainder)   | sign | ceil  |

## 8. Vector functions.

Other MATLAB functions operate essentially on a vector (row or column), but act on an  $m$ -by- $n$  matrix ( $m \geq 2$ ) in a column-by-column fashion to produce a row vector containing the results of their application to each column. Row-by-row action can be obtained by using the transpose; for example, **mean(A')'**. A few of these functions are

|      |      |        |     |
|------|------|--------|-----|
| max  | sum  | median | any |
| min  | prod | mean   | all |
| sort |      | std    |     |

For example, the maximum entry in a matrix *A* is given by **max(max(A))** rather than **max(A)**. Try it.

## 9. Matrix functions.

Much of MATLAB's power comes from its matrix functions. The most useful ones are

|       |                                        |
|-------|----------------------------------------|
| eig   | eigenvalues and eigenvectors           |
| chol  | cholesky factorization                 |
| svd   | singular value decomposition           |
| inv   | inverse                                |
| lu    | LU factorization                       |
| qr    | QR factorization                       |
| hess  | hessenberg form                        |
| schur | schur decomposition                    |
| rref  | reduced row echelon form               |
| expm  | matrix exponential                     |
| sqrtm | matrix square root                     |
| poly  | characteristic polynomial              |
| det   | determinant                            |
| size  | size                                   |
| norm  | 1-norm, 2-norm, F-norm, $\infty$ -norm |
| cond  | condition number in the 2-norm         |
| rank  | rank                                   |

MATLAB functions may have single or multiple output arguments. For example,

```
y = eig(A), or simply eig(A)
```

produces a column vector containing the eigenvalues of *A* while

```
[U,D] = eig(A)
```

produces a matrix *U* whose columns are the eigenvectors of *A* and a diagonal matrix *D* with the eigenvalues of *A* on its diagonal. Try it.

## 10. Command line editing and recall.

The command line in MATLAB can be easily edited. The cursor can be positioned with the left/right arrows and the Backspace (or Delete) key used to delete the character to the left of the cursor. Other editing features are also available. On a PC try the Home, End, and Delete keys; on a Unix system or a PC the Emacs commands Ctl-a, Ctl-e, Ctl-d, and Ctl-k work; on other systems see **help cedit** or **type cedit**.

A convenient feature is use of the up/down arrows to scroll through the stack of previous commands. One can, therefore, recall a previous command line, edit it, and execute the revised command line. For small routines, this is much more convenient than using an M-file which requires moving between MATLAB and the editor (see sections 12 and 14). For example, **flopcounts** (see section 15) for computing the inverse of matrices of various sizes could be compared by repeatedly recalling, editing, and executing

```
a = rand(8); flops(0), inv(a); flops
```

If one wanted to compare plots of the functions  $y = \sin mx$  and  $y = \sin nx$  on the interval  $[0, 2\pi]$  for various  $m$  and  $n$ , one might do the same for the command line:

```
m=2; n=3; x=0:.01:2*pi; y=sin(m*x); z=cos(n*x); plot(x,y,x,z)
```

## 11. Submatrices and colon notation.

Vectors and submatrices are often used in MATLAB to achieve fairly complex data manipulation effects. "Colon notation" (which is used both to generate vectors and reference submatrices) and subscripting by integral vectors are keys to efficient manipulation of these objects. Creative use of these features to vectorize operations permits one to minimize the use of loops (which slows MATLAB) and to make code simple and readable. *Special effort should be made to become familiar with them.*

The expression **1:5** (met earlier in **for** statements) is actually the row vector **[1 2 3 4 5]**. The numbers need not be integers nor the increment one. For example,

```
0.2:0.2:1.2
```

gives **[0.2, 0.4, 0.6, 0.8, 1.0, 1.2]**, and

```
5:-1:1 gives [5 4 3 2 1].
```

The following statements will, for example, generate a table of sines. Try it.

```
x = [0.0:0.1:2.0]';
y = sin(x);
[x y]
```

Note that since `sin` operates entry-wise, it produces a vector  $y$  from the vector  $x$ .

The colon notation can be used to access submatrices of a matrix. For example,

`A(1:4,3)` is the column vector consisting of the first four entries of the third column of  $A$ .

A colon by itself denotes an entire row or column:

`A(:,3)` is the third column of  $A$ , and `A(1:4,:)` is the first four rows.

Arbitrary integral vectors can be used as subscripts:

`A(:, [2 4])` contains as columns, columns 2 and 4 of  $A$ .

Such subscripting can be used on both sides of an assignment statement:

`A(:, [2 4 5]) = B(:, 1:3)` replaces columns 2,4,5 of  $A$  with the first three columns of  $B$ . Note that the *entire* altered matrix  $A$  is printed and assigned. Try it.

Columns 2 and 4 of  $A$  can be multiplied on the right by the 2-by-2 matrix `[1 2; 3 4]`:

`A(:, [2,4]) = A(:, [2,4])*[1 2; 3 4]`

Once again, the entire altered matrix is printed and assigned.

If  $x$  is an  $n$ -vector, what is the effect of the statement `x = x(n:-1:1)`? Try it. Also try `y = flipr(x)` and `y = flipud(x')`.

To appreciate the usefulness of these features, compare these MATLAB statements with a Pascal, FORTRAN, or C routine to effect the same.

## 12. M-files.

MATLAB can execute a sequence of statements stored in diskfiles. Such files are called “M-files” because they must have the file type of “.m” as the last part of their filename. Much of your work with MATLAB will be in creating and refining M-files. M-files are usually created using your local editor.

There are two types of M-files: *script files* and *function files*.

### Script files.

A script file consists of a sequence of normal MATLAB statements. If the file has the filename, say, `rotate.m`, then the MATLAB command `rotate` will cause the statements in the file to be executed. Variables in a script file are global and will change the value of variables of the same name in the environment of the current MATLAB session.

Script files may be used to enter data into a large matrix; in such a file, entry errors can be easily corrected. If, for example, one enters in a diskfile `data.m`

```
A = [
1 2 3 4
5 6 7 8
];
```

then the MATLAB statement `data` will cause the assignment given in `data.m` to be carried out. However, it is usually easier to use the MATLAB function `load` (see section 2).

An M-file can reference other M-files, including referencing itself recursively.

### Function files.

Function files provide extensibility to MATLAB. You can create new functions specific to your problem which will then have the same status as other MATLAB functions. Variables in a function file are by default local. A variable can, however, be declared global (see `help global`).

We first illustrate with a simple example of a function file.

```
function a = randint(m,n)
%RANDINT Randomly generated integral matrix.
%   randint(m,n) returns an m-by-n such matrix with entries
%   between 0 and 9.
a = floor(10*rand(m,n));
```

A more general version of this function is the following:

```
function a = randint(m,n,a,b)
%RANDINT Randomly generated integral matrix.
%   randint(m,n) returns an m-by-n such matrix with entries
%   between 0 and 9.
%   rand(m,n,a,b) return entries between integers a and b.
if nargin < 3, a = 0; b = 9; end
a = floor((b-a+1)*rand(m,n)) + a;
```

This should be placed in a diskfile with filename `randint.m` (corresponding to the function name). The first line declares the function name, input arguments, and output arguments; without this line the file would be a script file. Then a MATLAB statement `z = randint(4,5)`, for example, will cause the numbers 4 and 5 to be passed to the variables  $m$  and  $n$  in the function file with the output result being passed out to the variable  $z$ . Since variables in a function file are local, their names are independent of those in the current MATLAB environment.

Note that use of `nargin` (“number of input arguments”) permits one to set a default value of an omitted input variable—such as  $a$  and  $b$  in the example.

A function may also have multiple output arguments. For example:

```
function [mean, stdev] = stat(x)
% STAT Mean and standard deviation
%   For a vector x, stat(x) returns the mean of x;
%   [mean, stdev] = stat(x) both the mean and standard deviation.
%   For a matrix x, stat(x) acts columnwise.
[m n] = size(x);
if m == 1
    m = n; % handle case of a row vector
end
mean = sum(x)/m;
stdev = sqrt(sum(x.^2)/m - mean.^2);
```

Once this is placed in a diskfile `stat.m`, a MATLAB command `[xm, xd] = stat(x)`, for example, will assign the mean and standard deviation of the entries in the vector  $x$  to

$xm$  and  $xd$ , respectively. Single assignments can also be made with a function having multiple output arguments. For example, `xm = stat(x)` (no brackets needed around  $xm$ ) will assign the mean of  $x$  to  $xm$ .

The `%` symbol indicates that the rest of the line is a comment; MATLAB will ignore the rest of the line. Moreover, the first few contiguous comment lines, which document the M-file, are available to the on-line help facility and will be displayed if, for example, `help stat` is entered. Such documentation should *always* be included in a function file.

This function illustrates some of the MATLAB features that can be used to produce efficient code. Note, for example, that `x.^2` is the matrix of squares of the entries of  $x$ , that `sum` is a vector function (section 8), that `sqrt` is a scalar function (section 7), and that the division in `sum(x)/m` is a matrix-scalar operation. Thus all operations are vectorized and loops avoided.

If you can't vectorize some computations, you can make your `for` loops go faster by preallocating any vectors or matrices in which output is stored. For example, by including the second statement below, which uses the function `zeros`, space for storing  $E$  in memory is preallocated. Without this MATLAB must resize  $E$  one column larger in each iteration, slowing execution.

```
M = magic(6);
E = zeros(6,50);
for j = 1:50
    E(:,j) = eig(M^i);
end
```

Some more advanced features are illustrated by the following function. As noted earlier, some of the input arguments of a function—such as `tol` in this example, may be made optional through use of `nargin` (“number of input arguments”). The variable `nargout` can be similarly used. Note that the fact that a relation is a number (1 when true; 0 when false) is used and that, when `while` or `if` evaluates a relation, “nonzero” means “true” and 0 means “false”. Finally, the MATLAB function `feval` permits one to have as an input variable a string naming another function. (Also see `eval`.)

```
function [b, steps] = bisect(fun, x, tol)
%BISECT Zero of a function of one variable via the bisection method.
%   bisect(fun,x) returns a zero of the function. fun is a string
%   containing the name of a real-valued MATLAB function of a
%   single real variable; ordinarily functions are defined in
%   M-files. x is a starting guess. The value returned is near
%   a point where fun changes sign. For example,
%   bisect('sin',3) is pi. Note the quotes around sin.
%
%   An optional third input argument sets a tolerance for the
%   relative accuracy of the result. The default is eps.
%   An optional second output argument gives a matrix containing a
%   trace of the steps; the rows are of form [c f(c)].
```

```
% Initialization
if nargin < 3, tol = eps; end
trace = (nargout == 2);
if x ~= 0, dx = x/20; else, dx = 1/20; end
a = x - dx; fa = feval(fun,a);
b = x + dx; fb = feval(fun,b);

% Find change of sign.
while (fa > 0) == (fb > 0)
    dx = 2.0*dx;
    a = x - dx; fa = feval(fun,a);
    if (fa > 0) ~= (fb > 0), break, end
    b = x + dx; fb = feval(fun,b);
end
if trace, steps = [a fa; b fb]; end

% Main loop
while abs(b - a) > 2.0*tol*max(abs(b),1.0)
    c = a + 0.5*(b - a); fc = feval(fun,c);
    if trace, steps = [steps; [c fc]]; end
    if (fb > 0) == (fc > 0)
        b = c; fb = fc;
    else
        a = c; fa = fc;
    end
end
end
```

Some of MATLAB's functions are built-in while others are distributed as M-files. The actual listing of any non-built-in M-file—MATLAB's or your own—can be viewed with the MATLAB command `type functionname`. Try entering `type eig`, `type vander`, and `type rank`.

### 13. Text strings, error messages, input.

Text strings are entered into MATLAB surrounded by single quotes. For example,

```
s = 'This is a test'
```

assigns the given text string to the variable `s`.

Text strings can be displayed with the function `disp`. For example:

```
disp('this message is hereby displayed')
```

Error messages are best displayed with the function `error`

```
error('Sorry, the matrix must be symmetric')
```

since when placed in an M-File, it aborts execution of the M-file.

In an M-file the user can be prompted to interactively enter input data with the function `input`. When, for example, the statement

```
iter = input('Enter the number of iterations: ')
```

is encountered, the prompt message is displayed and execution pauses while the user keys in the input data. Upon pressing the return key, the data is assigned to the variable `iter` and execution resumes.

#### 14. Managing M-files.

While using MATLAB one frequently wishes to create or edit an M-file with the local editor and then return to MATLAB. One wishes to keep MATLAB active while editing a file since otherwise all variables would be lost upon exiting.

This can be easily done using the `!`-feature. If, while in MATLAB, you precede it with an `!`, any system command—such as those for editing, printing, or copying a file—can be executed without exiting MATLAB. If, for example, the system command `ed` accesses your editor, the MATLAB command

```
>> !ed rotate.m
```

will let you edit the file named `rotate.m` using your local editor. Upon leaving the editor, you will be returned to MATLAB just where you left it.

However, as noted in section 1, on systems permitting multiple processes, such as one running Unix or MS Windows, it may be preferable to keep both MATLAB and your local editor active, keeping one process suspended while working in the other. If these processes can be run in multiple windows, you will want to keep MATLAB active in one window and your editor active in another.

You should consult your instructor or your local computing center for details of the local installation.

Many debugging tools are available. See `help dbtype` or the list of functions in the last section.

When in MATLAB, the command `pwd` will return the name of the present working directory and `cd` can be used to change the working directory. Either `dir` or `ls` will list the contents of the working directory while the command `what` lists only the M-files in the directory. The MATLAB commands `delete` and `type` can be used to delete a diskfile and print an M-file to the screen, respectively. While these commands may duplicate system commands, they avoid the use of an `!`. You may enjoy entering the command `why` a few times.

M-files must be in a directory accessible to MATLAB. M-files in the present working directory are always accessible. On most mainframe or workstation network installations, personal M-files which are stored in a subdirectory of one's home directory named `matlab` will be accessible to MATLAB from any directory in which one is working. The current list of directories in MATLAB's search path is obtained by the command `path`. This command can also be used to add or delete directories from the search path. See `help path`.

#### 15. Comparing efficiency of algorithms: `flops`, `tic` and `toc`.

Two measures of the efficiency of an algorithm are the number of floating point operations (`flops`) performed and the elapsed time.

The MATLAB function `flops` keeps a running total of the flops performed. The command `flops(0)` (not `flops = 0!`) will reset `flops` to 0. Hence, entering `flops(0)` immediately before executing an algorithm and `flops` immediately after gives the flop count for the algorithm. For example, the number of flops required to solve a given linear system via Gaussian elimination can be obtained with:

```
flops(0), x = A\b; flops
```

The elapsed time (in seconds) can be obtained with the stopwatch timers `tic` and `toc`; `tic` starts the timer and `toc` returns the elapsed time. Hence, the commands

```
tic, any statement, toc
```

will return the elapsed time for execution of the statement. The elapsed time for solving the linear system above can be obtained, for example, with:

```
tic, x = A\b; toc
```

You may wish to compare this time—and flop count—with that for solving the system using `x = inv(A)*b`. Try it.

It should be noted that, on timesharing machines, elapsed time may not be a reliable measure of the efficiency of an algorithm since the rate of execution depends on how busy the computer is at the time.

#### 16. Output format.

While all computations in MATLAB are performed in double precision, the format of the displayed output can be controlled by the following commands.

|                             |                                                 |
|-----------------------------|-------------------------------------------------|
| <code>format short</code>   | fixed point with 4 decimal places (the default) |
| <code>format long</code>    | fixed point with 14 decimal places              |
| <code>format short e</code> | scientific notation with 4 decimal places       |
| <code>format long e</code>  | scientific notation with 15 decimal places      |
| <code>format rat</code>     | approximation by ratio of small integers        |
| <code>format hex</code>     | hexadecimal format                              |
| <code>format bank</code>    | fixed dollars and cents                         |
| <code>format +</code>       | +, -, blank                                     |

Once invoked, the chosen format remains in effect until changed.

The command `format compact` will suppress most blank lines allowing more information to be placed on the screen or page. The command `format loose` returns to the non-compact format. These commands are independent of the other format commands.

## 17. Hardcopy.

Hardcopy is most easily obtained with the **diary** command. The command

```
diary filename
```

causes what appears subsequently on the screen (except graphics) to be written to the named diskfile (if the filename is omitted it will be written to a default file named **diary**) until one gives the command **diary off**; the command **diary on** will cause writing to the file to resume, etc. When finished, you can edit the file as desired and print it out on the local system. The **!**-feature (see section 14) will permit you to edit and print the file without leaving MATLAB.

## 18. Graphics.

MATLAB can produce planar plots of curves, 3-D plots of curves, 3-D mesh surface plots, and 3-D faceted surface plots. The primary commands for these facilities are **plot**, **plot3**, **mesh**, and **surf**, respectively. An introduction to each of these is given below.

To preview some of these capabilities, enter the command **demo** and select some of the graphics options.

### Planar plots.

The **plot** command creates linear x-y plots; if  $x$  and  $y$  are vectors of the same length, the command **plot(x,y)** opens a graphics window and draws an x-y plot of the elements of  $x$  versus the elements of  $y$ . You can, for example, draw the graph of the sine function over the interval -4 to 4 with the following commands:

```
x = -4:.01:4; y = sin(x); plot(x,y)
```

Try it. The vector  $x$  is a partition of the domain with meshsize 0.01 while  $y$  is a vector giving the values of sine at the nodes of this partition (recall that **sin** operates entrywise).

You will usually want to keep the current graphics window (“figure”) exposed—but moved to the side—and the command window active.

One can have several graphics figures, one of which will at any time be the designated “current” figure where graphs from subsequent plotting commands will be placed. If, for example, figure 1 is the current figure, then the command **figure(2)** (or simply **figure**) will open a second figure (if necessary) and make it the current figure. The command **figure(1)** will then expose figure 1 and make it again the current figure. The command **gcf** will return the number of the current figure.

As a second example, you can draw the graph of  $y = e^{-x^2}$  over the interval -1.5 to 1.5 as follows:

```
x = -1.5:.01:1.5; y = exp(-x.^2); plot(x,y)
```

Note that one must precede **^** by a period to ensure that it operates entrywise (see section 3).

MATLAB supplies a function **fplot** to easily and efficiently plot the graph of a function. For example, to plot the graph of the function above, one can first define the function in an M-file called, say, **expnormal.m** containing

```
function y = expnormal(x)
y = exp(-x.^2);
```

Then the command

```
fplot('expnormal', [-1.5,1.5])
```

will produce the graph. Try it.

Plots of parametrically defined curves can also be made. Try, for example,

```
t=0:.001:2*pi; x=cos(3*t); y=sin(2*t); plot(x,y)
```

The graphs can be given titles, axes labeled, and text placed within the graph with the following commands which take a string as an argument.

|               |                                         |
|---------------|-----------------------------------------|
| <b>title</b>  | graph title                             |
| <b>xlabel</b> | x-axis label                            |
| <b>ylabel</b> | y-axis label                            |
| <b>gtext</b>  | place text on the graph using the mouse |
| <b>text</b>   | position text at specified coordinates  |

For example, the command

```
title('Best Least Squares Fit')
```

gives a graph a title. The command **gtext('The Spot')** allows one to interactively place the designated text on the current graph by placing the mouse pointer at the desired position and clicking the mouse. To place text in a graph at designated coordinates, one would use the command **text** (see **help text**).

The command **grid** will place grid lines on the current graph.

By default, the axes are auto-scaled. This can be overridden by the command **axis**. Some features of **axis** are:

|                                    |                                            |
|------------------------------------|--------------------------------------------|
| <b>axis([xmin,xmax,ymin,ymax])</b> | set axis scaling to prescribed limits      |
| <b>axis(axis)</b>                  | freezes scaling for subsequent graphs      |
| <b>axis auto</b>                   | returns to auto-scaling                    |
| <b>v = axis</b>                    | returns vector $v$ showing current scaling |
| <b>axis square</b>                 | same scale on both axes                    |
| <b>axis equal</b>                  | same scale and tic marks on both axes      |
| <b>axis off</b>                    | turns off axis scaling and tic marks       |
| <b>axis on</b>                     | turns on axis scaling and tic marks        |

The **axis** command should be given *after* the **plot** command.

Two ways to make multiple plots on a single graph are illustrated by

```
x=0:.01:2*pi;y1=sin(x);y2=sin(2*x);y3=sin(4*x);plot(x,y1,x,y2,x,y3)
```

and by forming a matrix  $Y$  containing the functional values as columns

```
x=0:.01:2*pi; Y=[sin(x)', sin(2*x)', sin(4*x)']; plot(x,Y)
```

Another way is with **hold**. The command **hold on** freezes the current graphics screen so that subsequent plots are superimposed on it. The axes may, however, become rescaled. Entering **hold off** releases the “hold.”

One can override the default linetypes, pointtypes and colors. For example,

```
x=0:.01:2*pi; y1=sin(x); y2=sin(2*x); y3=sin(4*x);
plot(x,y1,'--',x,y2,':',x,y3,'+')
```

renders a dashed line and dotted line for the first two graphs while for the third the symbol **+** is placed at each node. The line- and mark-types are

Linetypes: solid (**-**), dashed (**--**), dotted (**:**), dashdot (**-.**)  
 Marktypes: point (**.**), plus (**+**), star (**\***), circle (**o**), x-mark (**x**)

Colors can be specified for the line- and mark-types.

Colors: yellow (**y**), magenta (**m**), cyan (**c**), red (**r**)  
 green (**g**), blue (**b**), white (**w**), black (**k**)

For example, `plot(x,y,'r--')` plots a red dashed line.

The command `subplot` can be used to partition the screen so that several small plots can be placed in one figure. See `help subplot`.

Other specialized 2-D plotting functions you may wish to explore via `help` are:

`polar`, `bar`, `hist`, `quiver`, `compass`, `feather`, `rose`, `stairs`, `fill`

### Graphics hardcopy

A hardcopy of the current graphics figure can be most easily obtained with the MATLAB command `print`. Entered by itself, it will send a high-resolution copy of the current graphics figure to the default printer.

The `printopt` M-file is used to specify the default setting used by the `print` command. If desired, one can change the defaults by editing this file (see `help printopt`).

The command `print filename` saves the current graphics figure to the designated filename in the default file format. If *filename* has no extension, then an appropriate extension such as `.ps`, `.eps`, or `.jet` is appended. If, for example, PostScript is the default file format, then

```
print lissajous
```

will create a PostScript file `lissajous.ps` of the current graphics figure which can subsequently be printed using the system `print` command. If *filename* already exists, it will be overwritten unless you use the `-append` option. The command

```
print -append lissajous
```

will append the (hopefully different) current graphics figure to the existing file `lissajous.ps`. In this way one can save several graphics figures in a single file.

The default settings can, of course, be overwritten. For example,

```
print -deps -f3 saddle
```

will save to an Encapsulated PostScript file `saddle.eps` the graphics figure 3 — even if it is not the current figure.

### 3-D line plots.

Completely analogous to `plot` in two dimensions, the command `plot3` produces curves in three dimensional space. If *x*, *y*, and *z* are three vectors of the same size, then the command `plot3(x,y,z)` will produce a perspective plot of the piecewise linear curve in 3-space passing through the points whose coordinates are the respective elements of *x*, *y*, and *z*. These vectors are usually defined parametrically. For example,

```
t=.01:.01:20*pi; x=cos(t); y=sin(t); z=t.^3; plot3(x,y,z)
```

will produce a helix which is compressed near the *x-y* plane (a “slinky”). Try it.

Just as for planar plots, a title and axis labels (including `zlabel`) can be added. The features of `axis` command described there also hold for 3-D plots; setting the axis scaling to prescribed limits will, of course, now require a 6-vector.

### 3-D mesh and surface plots.

Three dimensional wire mesh surface plots are drawn with the command `mesh`. The command `mesh(z)` creates a three-dimensional perspective plot of the elements of the matrix *z*. The mesh surface is defined by the *z*-coordinates of points above a rectangular grid in the *x-y* plane. Try `mesh(eye(10))`.

Similarly, three dimensional faceted surface plots are drawn with the command `surf`. Try `surf(eye(10))`.

To draw the graph of a function  $z = f(x, y)$  over a rectangle, one first defines vectors *xx* and *yy* which give partitions of the sides of the rectangle. With the function `meshgrid` one then creates a matrix *x*, each row of which equals *xx* and whose column length is the length of *yy*, and similarly a matrix *y*, each column of which equals *yy*, as follows:

```
[x,y] = meshgrid(xx,yy);
```

One then computes a matrix *z*, obtained by evaluating *f* entrywise over the matrices *x* and *y*, to which `mesh` or `surf` can be applied.

You can, for example, draw the graph of  $z = e^{-x^2-y^2}$  over the square  $[-2, 2] \times [-2, 2]$  as follows (try it):

```
xx = -2:.2:2;
yy = xx;
[x,y] = meshgrid(xx,yy);
z = exp(-x.^2 - y.^2);
mesh(z)
```

One could, of course, replace the first three lines of the preceding with

```
[x,y] = meshgrid(-2:.2:2, -2:.2:2);
```

Try this plot with `surf` instead of `mesh`.

As noted above, the features of the `axis` command described in the section on planar plots also hold for 3-D plots as do the commands for titles, axes labelling and the command `hold`.

The color shading of surfaces is set by the `shading` command. There are three settings for shading: `faceted` (default), `interpolated`, and `flat`. These are set by the commands

shading faceted, shading interp, or shading flat

Note that on surfaces produced by `surf`, the settings `interpolated` and `flat` remove the superimposed mesh lines. Experiment with various shadings on the surface produced above. The command `shading` (as well as `colormap` and `view` below) should be entered *after* the `surf` command.

The color profile of a surface is controlled by the `colormap` command. Available pre-defined colormaps include:

`hsv` (default), `hot`, `cool`, `jet`, `pink`, `copper`, `flag`, `gray`, `bone`

The command `colormap(cool)` will, for example, set a certain color profile for the current figure. Experiment with various colormaps on the surface produced above.

The command `view` can be used to specify in spherical or cartesian coordinates the viewpoint from which the 3-D object is to be viewed. See `help view`.

The MATLAB function `peaks` generates an interesting surface on which to experiment with shading, `colormap`, and `view`.

Plots of parametrically defined surfaces can also be made. The MATLAB functions `sphere` and `cylinder` will generate such plots of the named surfaces. (See `type sphere` and `type cylinder`.) The following is an example of a similar function which generates a plot of a torus.

```
function [x,y,z] = torus(r,n,a)
%TORUS Generate a torus
%   torus(r,n,a) generates a plot of a torus with central
%   radius a and lateral radius r.  n controls the number
%   of facets on the surface.  These input variables are optional
%   with defaults r = 0.5, n = 30, a = 1.
%
%   [x,y,z] = torus(r,n,a) generates three (n+1)-by-(n+1)
%   matrices so that surf(x,y,z) will produce the torus.
%
%   See also SPHERE, CYLINDER

if nargin < 3, a = 1; end
if nargin < 2, n = 30; end
if nargin < 1, r = 0.5; end
theta = pi*(0:2:2*n)/n;
phi = 2*pi*(0:2:n)/n;
xx = (a + r*cos(phi))*cos(theta);
yy = (a + r*cos(phi))*sin(theta);
zz = r*sin(phi)*ones(size(theta));
if nargin == 0
    surf(xx,yy,zz)
    ar = (a + r)/sqrt(2);
    axis([-ar,ar,-ar,ar,-ar,ar])
else
    %
```

```
x = xx; y = yy; z = zz;
end
```

Other 3-D plotting functions you may wish to explore via `help` are:

`meshz`, `surf`, `surfl`, `contour`, `pcolor`

## Handle Graphics.

Beyond those described above, MATLAB's graphics system provides low level functions which permit one to control virtually all aspects of the graphics environment to produce sophisticated plots. Enter the command `set(1)` and `gca,set(ans)` to see some of the properties of figure 1 which one can control. This system is called Handle Graphics, for which one is referred to the MATLAB User's Guide.

## 19. Sparse Matrix Computations.

In performing matrix computations, MATLAB normally assumes that a matrix is dense; that is, any entry in a matrix *may* be nonzero. If, however, a matrix contains sufficiently many zero entries, computation time could be reduced by avoiding arithmetic operations on zero entries and less memory could be required by storing only the nonzero entries of the matrix. This increase in efficiency in time and storage can make feasible the solution of significantly larger problems than would otherwise be possible. MATLAB provides the capability to take advantage of the sparsity of matrices.

Matlab has two storage modes, full and sparse, with full the default. The functions `full` and `sparse` convert between the two modes. For a matrix *A*, full or sparse, `nnz(A)` returns the number of nonzero elements in *A*.

A sparse matrix is stored as a linear array of its nonzero elements along with their row and column indices. If a full tridiagonal matrix *F* is created via, say,

```
F = floor(10*rand(6)); F = triu(tril(F,1),-1);
```

then the statement `S = sparse(F)` will convert *F* to sparse mode. Try it. Note that the output lists the nonzero entries in column major order along with their row and column indices. The statement `F = full(S)` restores *S* to full storage mode. One can check the storage mode of a matrix *A* with the command `issparse(A)`.

A sparse matrix is, of course, usually generated directly rather than by applying the function `sparse` to a full matrix. A sparse banded matrix can be easily created via the function `spdiags` by specifying diagonals. For example, a familiar sparse tridiagonal matrix is created by

```
m = 6; n = 6; e = ones(n,1); d = -2*e;
T = spdiags([e,d,e],[-1,0,1],m,n)
```

Try it. The integral vector `[-1,0,1]` specifies in which diagonals the columns of `[e,d,e]` should be placed (use `full(T)` to view). Experiment with other values of *m* and *n* and, say, `[-3,0,2]` instead of `[-1,0,1]`. See `help spdiags` for further features of `spdiags`.



The sparse analogs of `eye`, `zeros`, `ones`, and `randn` for full matrices are, respectively,

`speye`, `sparse`, `spones`, `sprandn`

The latter two take a matrix argument and replace only the nonzero entries with ones and normally distributed random numbers, respectively. `randn` also permits the sparsity structure to be randomized. The command `sparse(m,n)` creates a sparse zero matrix.

The versatile function `sparse` permits creation of a sparse matrix via listing its nonzero entries. Try, for example,

```
i = [1 2 3 4 4 4]; j = [1 2 3 1 2 3]; s = [5 6 7 8 9 10];
S = sparse(i,j,s,4,3), full(S)
```

In general, if the vector  $s$  lists the nonzero entries of  $S$  and the integral vectors  $i$  and  $j$  list their corresponding row and column indices, then

```
sparse(i,j,s,m,n)
```

will create the desired sparse  $m \times n$  matrix  $S$ . As another example try

```
n = 6; e = floor(10*rand(n-1,1)); E = sparse(2:n,1:n-1,e,n,n)
```

The arithmetic operations and most MATLAB functions can be applied independent of storage mode. The storage mode of the result? Operations on full matrices always give full results. Selected other results are ( $S=\text{sparse}$ ,  $F=\text{full}$ ):

Sparse:  $S+S$ ,  $S*S$ ,  $S.*S$ ,  $S.*F$ ,  $S^{\wedge}n$ ,  $S \backslash S$

Full:  $S+F$ ,  $S*F$ ,  $S \backslash F$ ,  $F \backslash S$

Sparse:  $\text{inv}(S)$ ,  $\text{chol}(S)$ ,  $\text{lu}(S)$ ,  $\text{diag}(S)$ ,  $\text{max}(S)$ ,  $\text{sum}(S)$

For sparse  $S$ ,  $\text{eig}(S)$  is full if  $S$  is symmetric but undefined if  $S$  is unsymmetric; `svd` requires a full argument. A matrix built from blocks, such as  $[A,B;C,D]$ , is sparse if any constituent block is sparse.

You may wish to compare, for the two storage modes, the efficiency of solving a tridiagonal system of equations for, say,  $n = 20, 50, 500, 1000$  by entering, recalling and editing the following two command lines:

```
n=20;e=ones(n,1);d=-2*e; T=spdiags([e,d,e],[-1,0,1],n,n); A=full(T);
b=ones(n,1);s=sparse(b);tic,T\s;sparsetime=toc, tic,A\b;fulltime=toc
```

## 20. Reference.

There are many MATLAB features which cannot be included in these introductory notes. Listed below are some of the MATLAB functions and operators available, grouped by subject area<sup>1</sup>. Use the on-line help facility or consult the Reference Guide for more detailed information on the functions.

There are many functions beyond these. There exist, in particular, several “toolboxes” of functions for specific areas<sup>2</sup>. Included among such are signal processing, control systems, robust-control, system identification, optimization, splines, chemometrics,  $\mu$ -analysis and synthesis, state-space identification, neural networks, image processing, symbolic math (Maple kernel), and statistics. These can be explored via the command `help`.

| MANAGING COMMANDS AND FUNCTIONS |                                                    |
|---------------------------------|----------------------------------------------------|
| <b>help</b>                     | help facility                                      |
| <b>what</b>                     | list M-files on disk                               |
| <b>type</b>                     | list named M-file                                  |
| <b>lookfor</b>                  | keyword search through the help entries            |
| <b>which</b>                    | locate functions and files                         |
| <b>demo</b>                     | run demonstrations                                 |
| <b>path</b>                     | control MATLAB's search path                       |
| <b>cedit</b>                    | set parameters for command line editing and recall |
| <b>version</b>                  | display MATLAB version you are running             |
| <b>whatsnew</b>                 | display toolbox README files                       |
| <b>info</b>                     | info about MATLAB and The MathWorks                |
| <b>why</b>                      | receive flippant answer                            |

| MANAGING VARIABLES AND THE WORKSPACE |                                           |
|--------------------------------------|-------------------------------------------|
| <b>who</b>                           | list current variables                    |
| <b>whos</b>                          | list current variables, long form         |
| <b>save</b>                          | save workspace variables to disk          |
| <b>load</b>                          | retrieve variables from disk              |
| <b>clear</b>                         | clear variables and functions from memory |
| <b>pack</b>                          | consolidate workspace memory              |
| <b>size</b>                          | size of matrix                            |
| <b>length</b>                        | length of vector                          |
| <b>disp</b>                          | display matrix or text                    |

<sup>1</sup> Source: MATLAB Reference Guide, version 4.1

<sup>2</sup> The toolboxes, which are optional, may not be installed on your system.

| WORKING WITH FILES AND THE OPERATING SYSTEM |                                                 |
|---------------------------------------------|-------------------------------------------------|
| <b>cd</b>                                   | change current working directory                |
| <b>pwd</b>                                  | show current working directory                  |
| <b>dir, ls</b>                              | directory listing                               |
| <b>delete</b>                               | delete file                                     |
| <b>getenv</b>                               | get environment variable                        |
| <b>!</b>                                    | execute operating system command                |
| <b>unix</b>                                 | execute operating system command; return result |
| <b>diary</b>                                | save text of MATLAB session                     |

| CONTROLLING THE COMMAND WINDOW |                                        |
|--------------------------------|----------------------------------------|
| <b>clc</b>                     | clear command window                   |
| <b>home</b>                    | send cursor home—to top of screen      |
| <b>format</b>                  | set output format                      |
| <b>echo</b>                    | echo commands inside script commands   |
| <b>more</b>                    | control paged output in command window |

| STARTING AND QUITTING FROM MATLAB |                                        |
|-----------------------------------|----------------------------------------|
| <b>quit</b>                       | terminate MATLAB                       |
| <b>startup</b>                    | M-file executed when MATLAB is started |
| <b>matlabrc</b>                   | master startup M-file                  |

| MATRIX OPERATORS                     | ARRAY OPERATORS          |
|--------------------------------------|--------------------------|
| <b>+</b> addition                    | <b>+</b> addition        |
| <b>—</b> subtraction                 | <b>—</b> subtraction     |
| <b>*</b> multiplication              | <b>.*</b> multiplication |
| <b>^</b> power                       | <b>.^</b> power          |
| <b>/</b> right division              | <b>./</b> right division |
| <b>\</b> left division               | <b>.\</b> left division  |
| <b>'</b> conjugate transpose         |                          |
| <b>.'</b> transpose                  |                          |
| <b>kron</b> Kronecker tensor product |                          |

| RELATIONAL AND LOGICAL OPERATORS |                       |              |              |
|----------------------------------|-----------------------|--------------|--------------|
| <b>&lt;</b>                      | less than             | <b>&amp;</b> | and          |
| <b>&lt;=</b>                     | less than or equal    | <b> </b>     | or           |
| <b>&gt;</b>                      | greater than          | <b>~</b>     | not          |
| <b>&gt;=</b>                     | greater than or equal | <b>xor</b>   | exclusive or |
| <b>==</b>                        | equal                 |              |              |
| <b>~=</b>                        | not equal             |              |              |

| SPECIAL CHARACTERS |                                                                               |
|--------------------|-------------------------------------------------------------------------------|
| <b>=</b>           | assignment statement                                                          |
| <b>[]</b>          | used to form vectors and matrices; enclose multiple function output variables |
| <b>( )</b>         | arithmetic expression precedence; enclose function input variables            |
| <b>.</b>           | decimal point                                                                 |
| <b>..</b>          | parent directory                                                              |
| <b>...</b>         | continue statement to next line                                               |
| <b>,</b>           | separate subscripts, function arguments, statements                           |
| <b>;</b>           | end rows, suppress printing                                                   |
| <b>%</b>           | comments                                                                      |
| <b>:</b>           | subscripting, vector generation                                               |
| <b>!</b>           | execute operating system command                                              |

| SPECIAL VARIABLES AND CONSTRAINTS |                                         |
|-----------------------------------|-----------------------------------------|
| <b>ans</b>                        | answer when expression not assigned     |
| <b>eps</b>                        | floating point precision                |
| <b>realmax</b>                    | largest floating point number           |
| <b>reallmin</b>                   | smallest positive floating point number |
| <b>pi</b>                         | $\pi$                                   |
| <b>i, j</b>                       | imaginary unit                          |
| <b>inf</b>                        | infinity                                |
| <b>NaN</b>                        | Not-a-Number                            |
| <b>flops</b>                      | floating point operation count          |
| <b>nargin</b>                     | number of function input arguments      |
| <b>nargout</b>                    | number of function output arguments     |
| <b>computer</b>                   | computer type                           |

| TIME AND DATE   |                           |
|-----------------|---------------------------|
| <b>date</b>     | current date              |
| <b>clock</b>    | wall clock                |
| <b>etime</b>    | elapsed time function     |
| <b>tic, toc</b> | stopwatch timer functions |
| <b>cputime</b>  | elapsed CPU time          |

| SPECIAL MATRICES |                                      |
|------------------|--------------------------------------|
| <b>zeros</b>     | matrix of zeros                      |
| <b>ones</b>      | matrix of ones                       |
| <b>eye</b>       | identity                             |
| <b>diag</b>      | diagonal                             |
| <b>toeplitz</b>  | Toeplitz                             |
| <b>magic</b>     | magic square                         |
| <b>compan</b>    | companion                            |
| <b>linspace</b>  | linearly spaced vectors              |
| <b>logspace</b>  | logarithmically spaced vectors       |
| <b>meshgrid</b>  | array for 3-D plots                  |
| <b>rand</b>      | uniformly distributed random numbers |
| <b>randn</b>     | normally distributed random numbers  |
| <b>hilb</b>      | Hilbert                              |
| <b>invhilb</b>   | inverse Hilbert (exact)              |
| <b>vander</b>    | Vandermonde                          |
| <b>pascal</b>    | Pascal                               |
| <b>hadamard</b>  | Hadamard                             |
| <b>hankel</b>    | Hankel                               |
| <b>rosser</b>    | symmetric eigenvalue test matrix     |
| <b>wilkinson</b> | Wilkinson's eigenvalue test matrix   |
| <b>gallery</b>   | two small test matrices              |

| MATRIX MANIPULATION |                                         |
|---------------------|-----------------------------------------|
| <b>diag</b>         | create or extract diagonals             |
| <b>rot90</b>        | rotate matrix 90 degrees                |
| <b>fliplr</b>       | flip matrix left-to-right               |
| <b>flipud</b>       | flip matrix up-to-down                  |
| <b>reshape</b>      | change size                             |
| <b>tril</b>         | lower triangular part                   |
| <b>triu</b>         | upper triangular part                   |
| <b>.'</b>           | transpose                               |
| <b>:</b>            | convert matrix to single column; $A(:)$ |

| LOGICAL FUNCTIONS |                                         |
|-------------------|-----------------------------------------|
| <b>exist</b>      | check if variables or functions exist   |
| <b>any</b>        | true if any element of vector is true   |
| <b>all</b>        | true if all elements of vector are true |
| <b>find</b>       | find indices of non-zero elements       |
| <b>isnan</b>      | true for NaNs                           |
| <b>isinf</b>      | true for infinite elements              |
| <b>finite</b>     | true for finite elements                |
| <b>isieee</b>     | true for IEEE floating point arithmetic |
| <b>isempty</b>    | true for empty matrix                   |
| <b>issparse</b>   | true for sparse matrix                  |
| <b>isstr</b>      | true for text string                    |
| <b>strcmp</b>     | compare string variables                |

| CONTROL FLOW  |                                                         |
|---------------|---------------------------------------------------------|
| <b>if</b>     | conditionally execute statements                        |
| <b>else</b>   | used with <b>if</b>                                     |
| <b>elseif</b> | used with <b>if</b>                                     |
| <b>end</b>    | terminate <b>if</b> , <b>for</b> , <b>while</b>         |
| <b>for</b>    | repeat statements for a specific number of times        |
| <b>while</b>  | repeat statments while condition is true                |
| <b>break</b>  | terminate execution of <b>for</b> or <b>while</b> loops |
| <b>return</b> | return to invoking function                             |
| <b>error</b>  | display message and abort function                      |

| PROGRAMMING     |                                             |
|-----------------|---------------------------------------------|
| <b>input</b>    | prompt for user input                       |
| <b>keyboard</b> | invoke keyboard as if it were a script file |
| <b>menu</b>     | generate menu of choices for user input     |
| <b>pause</b>    | wait for user response                      |
| <b>function</b> | define function                             |
| <b>eval</b>     | execute string with MATLAB expression       |
| <b>feval</b>    | evaluate function specified by string       |
| <b>global</b>   | define global variables                     |
| <b>nargchk</b>  | validate number of input arguments          |

| TEXT AND STRINGS |                                             |
|------------------|---------------------------------------------|
| <b>string</b>    | about character strings in MATLAB           |
| <b>abs</b>       | convert string to numeric values            |
| <b>blanks</b>    | a string of blanks                          |
| <b>eval</b>      | evaluate string with MATLAB expression      |
| <b>num2str</b>   | convert number to string                    |
| <b>int2str</b>   | convert integer to string                   |
| <b>str2num</b>   | convert string to number                    |
| <b>isstr</b>     | true for string variables                   |
| <b>strcmp</b>    | compare string variables                    |
| <b>upper</b>     | convert string to uppercase                 |
| <b>lower</b>     | convert string to lowercase                 |
| <b>hex2num</b>   | convert hex string to floating point number |
| <b>hex2dec</b>   | convert hex string to decimal integer       |
| <b>dec2hex</b>   | convert decimal integer to hex string       |

| DEBUGGING       |                                |
|-----------------|--------------------------------|
| <b>dbstop</b>   | set breakpoint                 |
| <b>dbclear</b>  | remove breakpoint              |
| <b>dbcont</b>   | remove execution               |
| <b>dbdown</b>   | change local workspace context |
| <b>dbstack</b>  | list who called whom           |
| <b>dbstatus</b> | list all breakpoints           |
| <b>dbstep</b>   | execute one or more lines      |
| <b>dbtype</b>   | list M-file with line numbers  |
| <b>dbup</b>     | change local workspace context |
| <b>dbdown</b>   | opposite of <b>dbup</b>        |
| <b>dbquit</b>   | quit debug mode                |

| SOUND PROCESSING FUNCTIONS |                             |
|----------------------------|-----------------------------|
| <b>saxis</b>               | sound axis scaling          |
| <b>sound</b>               | convert vector to sound     |
| <b>auread</b>              | Read Sun audio file         |
| <b>auwrite</b>             | Write Sun audio file        |
| <b>lin2mu</b>              | linear to mu-law conversion |
| <b>mu2lin</b>              | mu-law to linear conversion |

| ELEMENTARY MATH FUNCTIONS |                                     |
|---------------------------|-------------------------------------|
| <b>abs</b>                | absolute value or complex magnitude |
| <b>angle</b>              | phase angle                         |
| <b>sqrt</b>               | square root                         |
| <b>real</b>               | real part                           |
| <b>imag</b>               | imaginary part                      |
| <b>conj</b>               | complex conjugate                   |
| <b>gcd</b>                | greatest common divisor             |
| <b>lcm</b>                | least common multiple               |
| <b>round</b>              | round to nearest integer            |
| <b>fix</b>                | round toward zero                   |
| <b>floor</b>              | round toward $-\infty$              |
| <b>ceil</b>               | round toward $\infty$               |
| <b>sign</b>               | signum function                     |
| <b>rem</b>                | remainder                           |
| <b>exp</b>                | exponential base e                  |
| <b>log</b>                | natural logarithm                   |
| <b>log10</b>              | log base 10                         |

| TRIGONOMETRIC FUNCTIONS       |                                                                    |
|-------------------------------|--------------------------------------------------------------------|
| <b>sin, asin, sinh, asinh</b> | sine, arcsine, hyperbolic sine, hyperbolic arcsine                 |
| <b>cos, acos, cosh, acosh</b> | cosine, arccosine, hyperbolic cosine, hyperbolic arccosine         |
| <b>tan, atan, tanh, atanh</b> | tangent, arctangent, hyperbolic tangent, hyperbolic arctangent     |
| <b>cot, acot, coth, acoth</b> | cotangent, arccotangent, hyperbolic cotan., hyperbolic arccotan.   |
| <b>sec, asec, sech, asech</b> | secant, arcsecant, hyperbolic secant, hyperbolic arcsecant         |
| <b>csc, acsc, csch, acsch</b> | cosecant, arccosecant, hyperbolic cosecant, hyperbolic arccosecant |

| SPECIAL FUNCTIONS |                                |
|-------------------|--------------------------------|
| <b>bessel</b>     | bessel function                |
| <b>beta</b>       | beta function                  |
| <b>gamma</b>      | gamma function                 |
| <b>rat</b>        | rational approximation         |
| <b>rats</b>       | rational output                |
| <b>erf</b>        | error function                 |
| <b>erfinv</b>     | inverse error function         |
| <b>ellipke</b>    | complete elliptic integral     |
| <b>ellipj</b>     | Jacobian elliptic integral     |
| <b>expint</b>     | exponential integral           |
| <b>log2</b>       | dissect floating point numbers |
| <b>pow2</b>       | scale floating point numbers   |

| MATRIX DECOMPOSITIONS AND FACTORIZATIONS |                                                   |
|------------------------------------------|---------------------------------------------------|
| <b>inv</b>                               | inverse                                           |
| <b>lu</b>                                | factors from Gaussian elimination                 |
| <b>rref</b>                              | reduced row echelon form                          |
| <b>chol</b>                              | Cholesky factorization                            |
| <b>qr</b>                                | orthogonal-triangular decomposition               |
| <b>nls</b>                               | nonnegative least squares                         |
| <b>lsconv</b>                            | least squares in presence of known covariance     |
| <b>null</b>                              | null space                                        |
| <b>orth</b>                              | orthogonalization                                 |
| <b>eig</b>                               | eigenvalues and eigenvectors                      |
| <b>hess</b>                              | Hessenberg form                                   |
| <b>schur</b>                             | Schur decomposition                               |
| <b>cdf2rdf</b>                           | complex diagonal form to real block diagonal form |
| <b>rsf2csf</b>                           | real block diagonal form to complex diagonal form |
| <b>balance</b>                           | diagonal scaling for eigenvalue accuracy          |
| <b>qz</b>                                | generalized eigenvalues                           |
| <b>polyeig</b>                           | polynomial eigenvalue solver                      |
| <b>svd</b>                               | singular value decomposition                      |
| <b>pinv</b>                              | pseudoinverse                                     |

| MATRIX CONDITIONING |                                               |
|---------------------|-----------------------------------------------|
| <b>cond</b>         | condition number in 2-norm                    |
| <b>rcond</b>        | LINPACK reciprocal condition number estimator |
| <b>condest</b>      | Hager/Higham condition number estimator       |
| <b>norm</b>         | 1-norm, 2-norm, F-norm, OC-norm               |
| <b>normest</b>      | 2-norm estimator                              |
| <b>rank</b>         | rank                                          |

| ELEMENTARY MATRIX FUNCTIONS |                                                     |
|-----------------------------|-----------------------------------------------------|
| <b>expm</b>                 | matrix exponential                                  |
| <b>expm1</b>                | M-file implementation of <b>expm</b>                |
| <b>expm2</b>                | matrix exponential via Taylor series                |
| <b>expm3</b>                | matrix exponential via eigenvalues and eigenvectors |
| <b>logm</b>                 | matrix logarithm                                    |
| <b>sqrtn</b>                | matrix square root                                  |
| <b>fmm</b>                  | evaluate general matrix function                    |
| <b>poly</b>                 | characteristic polynomial                           |
| <b>det</b>                  | determinant                                         |
| <b>trace</b>                | trace                                               |

| POLYNOMIALS     |                                           |
|-----------------|-------------------------------------------|
| <b>poly</b>     | construct polynomial with specified roots |
| <b>roots</b>    | polynomial roots—companion matrix method  |
| <b>roots1</b>   | polynomial roots—Laguerre's method        |
| <b>polyval</b>  | evaluate polynomial                       |
| <b>polyvalm</b> | evaluate polynomial with matrix argument  |
| <b>conv</b>     | multiply polynomials                      |
| <b>deconv</b>   | divide polynomials                        |
| <b>residue</b>  | partial-fraction expansion (residues)     |
| <b>polyfit</b>  | fit polynomial to data                    |
| <b>polyder</b>  | differentiate polynomial                  |

| COLUMN-WISE DATA ANALYSIS |                                |
|---------------------------|--------------------------------|
| <b>max</b>                | largest component              |
| <b>min</b>                | smallest component             |
| <b>mean</b>               | average or mean value          |
| <b>median</b>             | median value                   |
| <b>std</b>                | standard deviation             |
| <b>sort</b>               | sort in ascending order        |
| <b>sum</b>                | sum of elements                |
| <b>prod</b>               | product of elements            |
| <b>cumsum</b>             | cumulative sum of elements     |
| <b>cumprod</b>            | cumulative product of elements |
| <b>hist</b>               | histogram                      |

| SIGNAL PROCESSING |                                                 |
|-------------------|-------------------------------------------------|
| <b>abs</b>        | complex magnitude                               |
| <b>angle</b>      | phase angle                                     |
| <b>conv</b>       | convolution and polynomial multiplication       |
| <b>deconv</b>     | deconvolution and polynomial division           |
| <b>corrcoef</b>   | correlation coefficients                        |
| <b>cov</b>        | covariance matrix                               |
| <b>filter</b>     | one-dimensional digital filter                  |
| <b>filter2</b>    | two-dimensional digital filter                  |
| <b>cplxpair</b>   | sort numbers into complex pairs                 |
| <b>unwrap</b>     | remove phase angle jumps across 360° boundaries |
| <b>nextpow2</b>   | next higher power of 2                          |
| <b>fft</b>        | radix-2 fast Fourier transform                  |
| <b>fft2</b>       | two-dimensional FFT                             |
| <b>ifft</b>       | inverse fast Fourier transform                  |
| <b>ifft2</b>      | inverse 2-D FFT                                 |
| <b>fftshift</b>   | zero-th lag to center of spectrum               |

| FINITE DIFFERENCES AND DATA INTERPOLATION |                                       |
|-------------------------------------------|---------------------------------------|
| <b>diff</b>                               | approximate derivatives               |
| <b>gradient</b>                           | approximate gradient                  |
| <b>del2</b>                               | five point discrete Laplacian         |
| <b>subspace</b>                           | angle between two subspaces           |
| <b>spline</b>                             | cubic spline interpolation            |
| <b>interp1</b>                            | 1-D data interpolation                |
| <b>interp2</b>                            | 2-D data interpolation                |
| <b>interpft</b>                           | 1-D data interpolation via FFT method |
| <b>griddata</b>                           | data gridding                         |

| NUMERICAL INTEGRATION |                                    |
|-----------------------|------------------------------------|
| <b>quad</b>           | adaptive 2-panel Simpson's Rule    |
| <b>quad8</b>          | adaptive 8-panel Newton-Cotes Rule |
| <b>trapz</b>          | trapezoidal method                 |

| DIFFERENTIAL EQUATION SOLUTION |                                           |
|--------------------------------|-------------------------------------------|
| <b>ode23</b>                   | 2nd/3rd order Runge-Kutta method          |
| <b>ode23p</b>                  | solve via <b>ode23</b> , displaying plot  |
| <b>ode45</b>                   | 4th/5th order Runge-Kutta-Fehlberg method |

| NONLINEAR EQUATIONS AND OPTIMIZATION |                                                                                                |
|--------------------------------------|------------------------------------------------------------------------------------------------|
| <b>fmin</b>                          | minimize function of one variable                                                              |
| <b>fmin</b> s                        | minimize function of several variables                                                         |
| <b>fsolve</b>                        | solution to a system of nonlinear equations<br>(find zeros of a function of several variables) |
| <b>fzero</b>                         | find zero of function of one variable                                                          |
| <b>fplot</b>                         | plot graph of a function                                                                       |

| TWO DIMENSIONAL GRAPHS |                          |
|------------------------|--------------------------|
| <b>plot</b>            | linear plot              |
| <b>loglog</b>          | log-log scale plot       |
| <b>semilogx</b>        | semilog scale plot       |
| <b>semilogy</b>        | semilog scale plot       |
| <b>fill</b>            | draw filled 2-D polygons |
| <b>polar</b>           | polar coordinate plot    |
| <b>bar</b>             | bar graph                |
| <b>stairs</b>          | stairstep plot           |
| <b>errorbar</b>        | error bar plot           |
| <b>hist</b>            | histogram plot           |
| <b>rose</b>            | angle histogram plot     |
| <b>compass</b>         | compass plot             |
| <b>feather</b>         | feather plot             |
| <b>fplot</b>           | plot function            |

| GRAPH ANNOTATION |                            |
|------------------|----------------------------|
| <b>title</b>     | graph title                |
| <b>xlabel</b>    | x-axis label               |
| <b>ylabel</b>    | y-axis label               |
| <b>zlabel</b>    | z-axis label for 3-D plots |
| <b>grid</b>      | grid lines                 |
| <b>text</b>      | text annotation            |
| <b>gtext</b>     | mouse placement of text    |
| <b>ginput</b>    | graphical input from mouse |

| FIGURE WINDOW/AXIS CREATION AND CONTROL |                                     |
|-----------------------------------------|-------------------------------------|
| <b>figure</b>                           | create figure (graph window)        |
| <b>gcf</b>                              | get handle to current figure        |
| <b>clf</b>                              | clear current figure                |
| <b>close</b>                            | close figure                        |
| <b>hold</b>                             | hold current graph                  |
| <b>ishold</b>                           | return hold status                  |
| <b>subplot</b>                          | create axes in tiled positions      |
| <b>axes</b>                             | create axes in arbitrary positions  |
| <b>gca</b>                              | get handle to to current axes       |
| <b>axis</b>                             | control axis scaling and appearance |
| <b>caxis</b>                            | control pseudocolor axis scaling    |

| GRAPH HARDCOPY AND STORAGE |                                   |
|----------------------------|-----------------------------------|
| <b>print</b>               | print graph or save graph to file |
| <b>printopt</b>            | configure local printer defaults  |
| <b>orient</b>              | set paper orientation             |

| THREE DIMENSIONAL GRAPHS |                                                    |
|--------------------------|----------------------------------------------------|
| <b>mesh</b>              | 3-D mesh surface                                   |
| <b>meshc</b>             | combination mesh/contour plot                      |
| <b>meshz</b>             | 3-D mesh with zero plane                           |
| <b>surf</b>              | 3-D shaded surface                                 |
| <b>surfc</b>             | combination surface/contour plot                   |
| <b>surf1</b>             | 3-D shaded surface with lighting                   |
| <b>plot3</b>             | plot lines and points in 3-D space                 |
| <b>fill3</b>             | draw filled 3-D polygons in 3-D space              |
| <b>contour</b>           | contour plot                                       |
| <b>contour3</b>          | 3-D contour plot                                   |
| <b>clabel</b>            | contour plot elevation labels                      |
| <b>contourc</b>          | contour plot computation (used by <b>contour</b> ) |
| <b>pcolor</b>            | pseudocolor (checkerboard) plot                    |
| <b>quiver</b>            | quiver plot                                        |
| <b>image</b>             | display image                                      |
| <b>waterfall</b>         | waterfall plot                                     |
| <b>slice</b>             | volumetric visualization plot                      |

| 3-D GRAPH APPEARANCE |                                   |
|----------------------|-----------------------------------|
| <b>view</b>          | 3-D graph viewpoint specification |
| <b>viewmtx</b>       | view transformation matrices      |
| <b>hidden</b>        | mesh hidden line removal mode     |
| <b>shading</b>       | color shading mode                |
| <b>axis</b>          | axis scaling and appearance       |
| <b>caxis</b>         | pseudocolor axis scaling          |
| <b>specular</b>      | specular reflectance              |
| <b>diffuse</b>       | diffuse reflectance               |
| <b>surfnorm</b>      | surface normals                   |
| <b>colormap</b>      | color lookup table (see below)    |
| <b>brighten</b>      | brighten or darken color map      |
| <b>spinmap</b>       | spin color map                    |
| <b>rgbplot</b>       | plot colormap                     |
| <b>hsv2rgb</b>       | hsv to rgb color map conversion   |
| <b>rgb2hsv</b>       | rgb to hsv color map conversion   |

| COLOR MAPS    |                                         |
|---------------|-----------------------------------------|
| <b>hsv</b>    | hue-saturation-value (default)          |
| <b>jet</b>    | variant of <b>hsv</b>                   |
| <b>gray</b>   | linear gray-scale                       |
| <b>hot</b>    | black-red-yellow-white                  |
| <b>cool</b>   | shades of cyan and magenta              |
| <b>bone</b>   | gray-scale with tinge of blue           |
| <b>copper</b> | linear copper tone                      |
| <b>pink</b>   | pastel shades of pink                   |
| <b>flag</b>   | alternating red, white, blue, and black |

| 3-D OBJECTS     |                       |
|-----------------|-----------------------|
| <b>sphere</b>   | generate sphere       |
| <b>cylinder</b> | generate cylinder     |
| <b>peaks</b>    | generate demo surface |

| MOVIES AND ANIMATION |                               |
|----------------------|-------------------------------|
| <b>moviein</b>       | initialize movie frame memory |
| <b>getframe</b>      | get movie frame               |
| <b>movie</b>         | play recorded movie frames    |

| HANDLE GRAPHICS OBJECTS |                               |
|-------------------------|-------------------------------|
| <b>figure</b>           | create figure window          |
| <b>axes</b>             | create axes                   |
| <b>line</b>             | create line                   |
| <b>text</b>             | create text                   |
| <b>patch</b>            | create patch                  |
| <b>surface</b>          | create surface                |
| <b>image</b>            | create image                  |
| <b>uicontrol</b>        | create user interface control |
| <b>uimenu</b>           | create user interface menu    |

| HANDLE GRAPHICS OPERATIONS |                               |
|----------------------------|-------------------------------|
| <b>set</b>                 | set object properties         |
| <b>get</b>                 | get object properties         |
| <b>reset</b>               | reset object properties       |
| <b>delete</b>              | delete object                 |
| <b>drawnow</b>             | flush pending graphics events |

| SPARSE MATRIX FUNCTIONS |                                                     |
|-------------------------|-----------------------------------------------------|
| <b>spdiags</b>          | sparse matrix formed from diagonals                 |
| <b>speye</b>            | sparse identity matrix                              |
| <b>sprandn</b>          | sparse random matrix                                |
| <b>spones</b>           | replace nonzero entries with ones                   |
| <b>sprandsym</b>        | sparse symmetric random matrix                      |
| <b>spfun</b>            | apply function to nonzero entries                   |
| <b>sparse</b>           | create sparse matrix; convert full matrix to sparse |
| <b>full</b>             | convert sparse matrix to full matrix                |
| <b>find</b>             | find indices of nonzero entries                     |
| <b>spconvert</b>        | convert from sparse matrix external format          |
| <b>issparse</b>         | true if matrix is sparse                            |
| <b>nnz</b>              | number of nonzero entries                           |
| <b>nonzeros</b>         | nonzero entries                                     |
| <b>nzmax</b>            | amount of storage allocated for nonzero entries     |
| <b>spalloc</b>          | allocate memory for nonzero entries                 |
| <b>spy</b>              | visualize sparsity structure                        |
| <b>gplot</b>            | plot graph, as in “graph theory”                    |
| <b>colmmd</b>           | column minimum degree                               |
| <b>colperm</b>          | order columns based on nonzero count                |
| <b>dmlperm</b>          | Dulmage-Mendelsohn decomposition                    |
| <b>randperm</b>         | random permutation vector                           |
| <b>symmmd</b>           | symmetric minimum degree                            |
| <b>symrcm</b>           | reverse Cuthill-McKee ordering                      |
| <b>condest</b>          | estimate 1-norm condition                           |
| <b>normest</b>          | estimate 2-norm                                     |
| <b>sprank</b>           | structural rank                                     |
| <b>spaaugment</b>       | form least squares augmented system                 |
| <b>spparms</b>          | set parameters for sparse matrix routines           |
| <b>symbfact</b>         | symbolic factorization analysis                     |
| <b>parsefun</b>         | sparse auxillary functions and parameters           |



## B Travaux pratiques d'analyse numérique avec MATLAB

### B.1 Introduction

Vous disposez de 5 séances pour réaliser les 4 travaux pratiques ci-dessous. Les TP 1 et 3 doivent s'effectuer en 2 séances environ, le TP 2 ne doit nécessiter qu'une séance.

### B.2 Enoncés des TP

#### B.2.1 Travaux pratiques 1: Approximation différences-finies des équations elliptiques et conditions aux Limites

Ce TP concerne la discrétisation différences-finies de problèmes elliptiques 1D, de type convection-diffusion stationnaire ou Helmholtz déjà abordés dans les exercices A.8 et A.9. Le répertoire de travail MATLAB du TP est '*TP MATLAB CONV DIFF STATIONNAIRE*'.

#### B.2.2 Equation de convection-diffusion stationnaire

Cette partie concerne le problème aux limites (A.1) de l'exercice A.8.

Le programme `Tp_matlabconv_diff_stationnaire.m` résoud cette EDP pour  $a = 1$  et  $\nu = 0.02$  (variables MATLAB `adv` et `nu` respectivement) par méthode de différences finies centrées (fonction `schema_centre_convdiff_stat.m`). Ensuite, il trace sur le même graphe la solution numérique (vecteur `unum`) et la solution exacte (définie dans la fonction `uex_convdiff`) aux points du maillage.

**Question 1.** *Exécutez le programme pour  $n = 11$  points et calculez le nombre de Péclet de maille  $Pe_h$ . Que remarquez-vous ? Calculez le nombre minimal de points pour que les problèmes numériques disparaissent. Exportez les graphes nécessaires pour le compte-rendu final.*

On veut maintenant vérifier numériquement que la méthode de différences finies centrées est au second-ordre, en calculant pour plusieurs valeurs du pas d'espace  $h$  la norme  $l^2$  de l'erreur :

$$\|u_{\text{ex}} - u_h\|_{l^2} \quad (\text{B.1})$$

où  $\|\cdot\|_{l^2}$  représente l'erreur quadratique moyenne aux points intérieurs du maillage, définie par

$$\|\phi\|_{l^2}^2 = \frac{1}{N} \sum_{i=1}^N \phi(x_i)^2. \quad (\text{B.2})$$

**Question 2.** *Tracez sur un graphe en coordonnées logarithmiques (commande `loglog`) l'erreur  $l^2$  de la méthode de différences-finies centrées obtenue pour des valeurs croissantes<sup>1</sup> de  $n$ , et montrez qu'elle est au second-ordre en traçant sur le même graphe la fonction  $n \mapsto 1/n^2$ .*

**Question 3.** *Programmez la méthode de différences finies upwind. Par expérimentation numérique, vérifiez que la solution numérique ne présente pas d'oscillations parasites quelle que soit la valeur de  $Pe_h$ .*

---

<sup>1</sup> Conseil : prenez des valeurs de  $n$  supérieures à 50.

### B.2.3 Equation de Helmholtz et conditions aux limites de Neumann

Le but de cette partie est d'établir numériquement l'ordre de précision de schémas aux différences finies centrés pour le problème de Helmholtz avec les conditions aux limites de type Neumann-Dirichlet :

$$\lambda u - \frac{d^2 u}{dx^2} = f(x), \quad \lambda \geq 0, \quad (B.3a)$$

$$u'(-1) = \sqrt{\lambda} \tanh^{-1}(\sqrt{\lambda}), \quad u(1) = 1, \quad (B.3b)$$

dans  $\Omega = ]-1, 1[$ , en utilisant différentes manières d'imposer une condition aux limites de Neumann, telle que cela a été vu dans l'exercice A.9.

#### Solution exacte du problème de Helmholtz

Pour  $f(x) = \lambda$ , la solution exacte de (B.3) s'écrit

$$u_\lambda(x) = 1 - \frac{\sinh \sqrt{\lambda} (1-x)}{\sinh 2\sqrt{\lambda}}. \quad (B.4)$$

On peut observer que  $u_\lambda(x) \sim 1$  lorsque  $\lambda \gg 1$  sauf dans un voisinage du bord  $x = -1$ , de taille  $\mathcal{O}(1/\sqrt{\lambda})$ , où la solution exhibe une couche limite.

#### Implémentation d'une condition de Neumann

On cherche à évaluer les estimations d'erreur (B.1) pour 3 types d'implémentations d'une C.L. de Neumann :

- I) Méthode de l'exemple 7.3 du cours,
- II) Méthode de l'exemple 7.4,
- III) Méthode vue en TD.

Qui va l'emporter ?

**Question 4.** Placez-vous dans le cas  $\lambda = 100$ . Programmez la méthode de différences-finies centrées avec les trois variantes de C.L. (faites une fonction MATLAB pour chaque méthode).

**Question 5.** Comparez les estimations d'erreur obtenue pour chaque méthode. Sont-elles consistantes avec la théorie ? Distinguez la meilleure de ces trois méthodes.

**Question 6.** En utilisant les graphes de la question précédente, estimez le nombre de nœuds nécessaires pour que le schéma du premier ordre fournisse une précision (en norme  $l^2$ ) comparable à celle donnée par la méthode III avec  $n = 100$  points.

## B.3 Travaux pratiques 2 : Schémas Numériques pour les Equations Hyperboliques

### B.3.1 Introduction

On s'intéresse à la résolution numérique de l'équation de transport :

$$\frac{\partial u}{\partial t} + a \frac{\partial u}{\partial x} = 0 \quad (B.1a)$$

dans le domaine  $\Omega = ]a_x, b_x[ \times ]0, t_f[$ , avec la condition initiale

$$u(x, t = 0) = u_0(x), \quad (\text{B.1b})$$

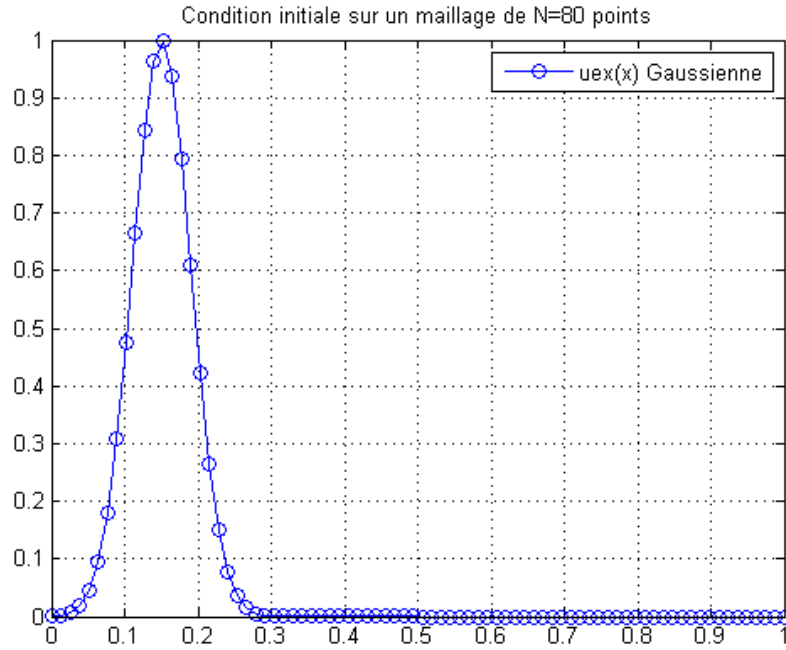
et la condition d'*inflow*

$$u(a_x, t) = g_-(t). \quad (\text{B.1c})$$

Le répertoire de travail MATLAB du TP est '*TP MATLAB EQT HYPERB*'.

### B.3.2 Condition initiale régulière

Dans cette première partie, on s'intéresse aux solutions de (B.1) qui restent continues en tout instant.



**Fig. B.1.** Condition initiale (B.2) avec  $\sigma = 0.04$  et  $x_0 = 0.15$ .

### Familiarisation avec le problème

On choisit le domaine  $\Omega = ]0, 1[$ , et le coefficient d'advection de l'E.D.P. est  $a = 1$  (variable "adv" dans le code MATLAB). La condition initiale est définie comme

$$u_0(x) = \exp\left(\frac{1}{2} \left[\frac{x - x_0}{\sigma}\right]^2\right), \quad (\text{B.2})$$

qui correspond à une gaussienne d'écart type  $\sigma$ , centrée en  $x_0 \in \Omega$  (cf. figure B.1). Plus la valeur de  $\sigma$  est grande, plus le problème est raide. On rappelle que la solution du problème (B.1) correspond à une translation uniforme de la condition initiale

*Remarque B.1.* Le problème (B.1) étant instationnaire, il faut fixer la borne supérieure d'intégration en temps  $T$ . Dans le code MATLAB, on définira

$$t_f \approx \frac{b_x - a_x}{a},$$

de façon que l'information toute entière soit sortie du domaine  $\Omega$ .

### L'approximation numérique

On va comparer qualitativement divers schémas pour la résolution de (B.1). Les schémas considérés sont les suivants :

- Schéma upwind (*cf.* cours),
- $\theta$ -schéma (*cf.* TD), pour  $\theta \in [0, 1]$  :

$$\frac{u_i^{n+1} - u_i^n}{\Delta t} + a [\theta \delta_x^0 u_i^{n+1} + (1 - \theta) \delta_x^0 u_i^n] = 0, \quad (\text{B.3})$$

- Schéma de Lax-Wendroff :

$$\frac{u_i^{n+1} - u_i^n}{\Delta t} + a \delta_x^0 u_i^n - a^2 \frac{\Delta t}{2} \delta_{xx}^0 u_i^n = 0. \quad (\text{B.4})$$

Le schéma upwind est déjà programmé (fonction `unew=schema_upwind(c,uold,ug)`). Ces schémas sont exprimés en fonction du nombre de courant  $C$  uniquement (variable “`cfl`” dans le code MATLAB). On rappelle la définition de  $C$  :

$$C = \frac{a\Delta t}{\Delta x}. \quad (\text{B.5})$$

*Remarque B.2.* Le problème (B.1) est bien posé avec l'unique condition aux limites amont (B.1c). Il est à noter que les schémas centrés, en particulier le  $\theta$ -schéma (B.3) et le schéma de Lax-Wendroff (B.4) nécessitent d'imposer une condition de sortie numérique en  $x = b_x$ , qui est donc artificielle.

### La comparaison des schémas

Le but de cette étude est d'observer par la pratique les notions qui ont été vues “théoriquement” en cours et en TD, *c.à.d* :

- Stabilité,
- Dissipation + dispersion numérique,
- Oscillations parasites.

Pour chacun des schémas, il s'agira d'évaluer qualitativement les propriétés énoncées ci-dessus et d'en effectuer une synthèse. A la différence du *TP* précédent, nous ne ferons pas de raffinement de maillage. On conseille ici de conserver  $N = 80$  points. Il est conseillé de prendre des notes manuscrites lors de l'évaluation de chacun des schémas.

**Question 1.** Le programme `MATLABTp_matlab_eqt_hyperboliques.m` utilise le schéma upwind pour résoudre l'équation de transport pour la condition initiale (B.2) avec  $\sigma = 0.04$  et  $x_0 = 0.15$ . Utilisez ce schéma avec  $C = 1.2$ . Que se passe-t-il ? Comment remédier au problème ? Observez et évaluez les phénomènes de dissipation et dispersion numérique.

*Remarque B.3.* Pour des valeurs de  $\Delta x$  et  $a$  données, il est équivalent de fixer la valeur de  $\Delta t$  ou celle de  $C$ , d'après la définition (B.5).

Désormais, la comparaison des différents schémas sera effectuée pour une même valeur du CFL.

**Question 2.** *Programmez le  $\theta$ -schéma de la manière qui a été définie en TD.*

**Question 3.** *Utilisez le  $\theta$ -schéma avec  $\theta = 0$ . Donnez le nom de ce schéma ? Quels sont les résultats obtenus ? Pouvez-vous le prévoir ?*

**Question 4.**  *$\theta$ -schéma avec  $\theta = 1/2$ . Le reconnaissez-vous ? Quel est son ordre de précision ? Utilisez le avec  $C = 0.8, 1.2, 2$ , etc ... Quelles sont ses propriétés de stabilité ? Pouvez-vous les prévoir ? Observez et évaluez les phénomènes de dissipation et dispersion numériques.*

**Question 5.**  *$\theta$ -schéma avec  $\theta = 1$ . Même questions que précédemment.*

**Question 6.** *Programmez le schéma de Lax-Wendroff. Rappelez l'ordre de précision de la méthode ? Retrouve-t'on les propriétés de stabilité théoriques ? Quelles sont les principales caractéristiques de la solution numérique par rapport à celle donnée par le schéma upwind ?*

**Question 7.** *Synthèse des résultats : parmi les schémas que vous venez d'utiliser, énoncez selon-vous les trois meilleurs et justifiez brièvement pourquoi vous rejetez les autres. Pour les trois meilleurs schémas, rappelez leurs propriétés et avantages et énoncez un classement subjectif selon ce que vous avez observé.*

### B.3.3 Condition initiale discontinue

Les problèmes de mécanique des fluides compressibles, modélisés par des équations hyperboliques, admettent des chocs, *i.e.* des discontinuités dans la solution. Comme prototype (linéaire), nous allons considérer le problème (B.1) avec la condition initiale de Riemann :

$$u_0(x) = \begin{cases} 1 & \text{si } x < x_0, \\ 0 & \text{si } x > x_0, \end{cases} \quad (\text{B.6})$$

qui va se propager à la vitesse  $a > 0$ .

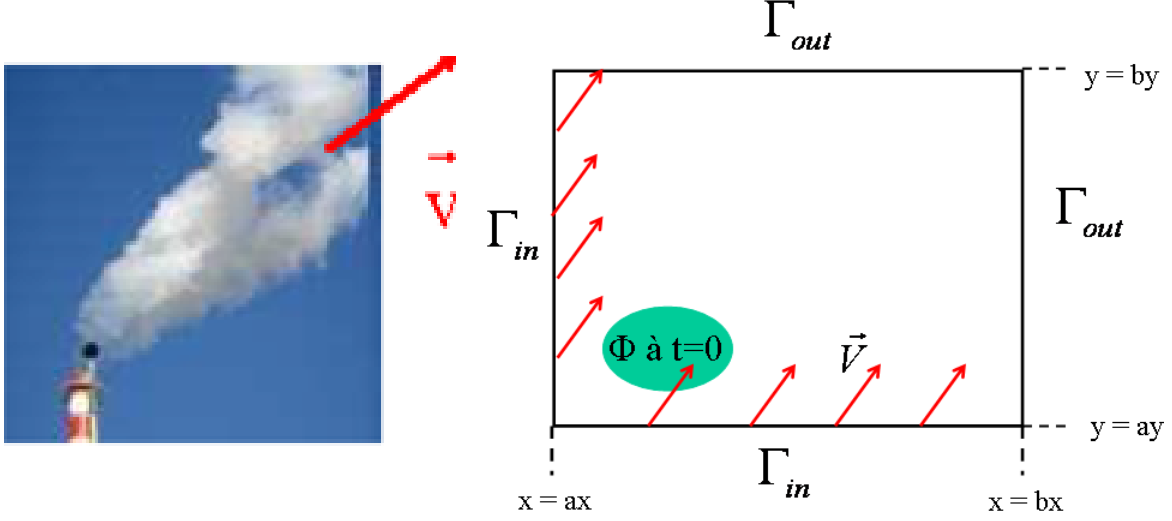
Il est, bien sûr, impossible d'obtenir une grande précision numérique au travers d'un choc. Il est néanmoins important de développer des méthodes numériques pour la mécanique des fluides telles que :

- une grande précision est obtenue loin du choc,
- il y a peu d'oscillations parasites près des discontinuités,
- La solution numérique conserve un profil de choc nettement défini, et non atténué par une trop forte viscosité artificielle.

**Question 8.** *Comparez les résultats des meilleurs schémas de la section B.3.2 lorsqu'ils sont appliqués à ce problème. Parmi les propriétés énoncées ci-dessus, lesquelles sont observées ? Conclusion + classement.*

## B.4 Travaux pratiques 3 : Modélisation du transport d'un polluant par la méthode de volumes finis

Ce TP concerne la modélisation d'un polluant dispersé par un fluide tel celui représenté sur la figure B.2 (gauche). Le transport de ce polluant est modélisé par une équation de convection-diffusion 2D qui sera discrétisée par la méthode des volumes finis. Le répertoire de travail MATLAB du TP est 'TP MATLAB 2D CONV DIFF'.



**Fig. B.2.** A gauche : dispersion de fumée dans l'atmosphère. A droite : Domaine de calcul considéré.

L'objectif du TP est de modéliser la dispersion de polluant (de fraction massique  $\phi(x, y, t)$ ) à la surface d'un domaine fluide représenté par le domaine de calcul  $\Omega = ]a_x, b_x[ \times ]a_y, b_y[$ . L'équation d'équilibre pour  $\phi$  traduit que la variation temporelle de la quantité de polluant  $\int_{\Omega} \rho \phi \, dV$  dans un volume élémentaire  $\Omega$  (de volume  $V$ , de surface  $\Gamma$ , et de normale extérieure  $\mathbf{n}$ ) est égale à un bilan de flux  $\int_{\Gamma} \mathbf{F} \cdot \mathbf{n} \, d\Gamma$  sur les faces  $\Gamma$  du volume élémentaire. Ce flux est modélisé par la somme d'un flux de diffusion :

$$\mathbf{F}_d = \lambda \nabla \phi, \quad \text{où } \lambda > 0 \text{ est le coefficient de diffusion,} \quad (\text{B.1})$$

et d'un flux de convection :

$$\mathbf{F}_c = -\rho \mathbf{v} \phi, \quad \text{tel que } \mathbf{v} \text{ est la vitesse du fluide.} \quad (\text{B.2})$$

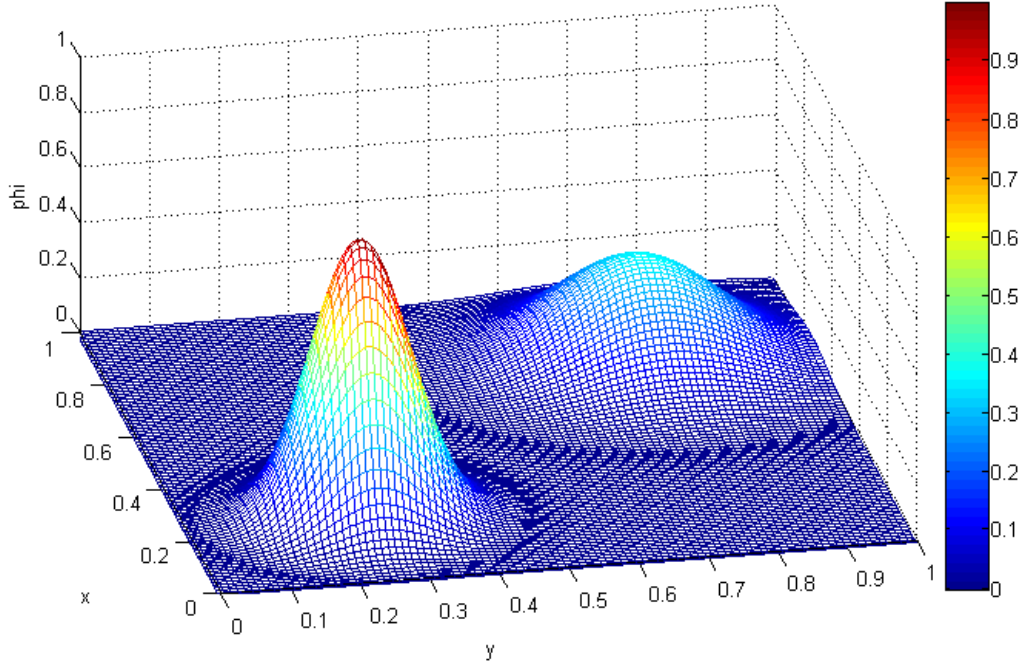
On suppose que la vitesse du fluide est telle que représentée sur la figure B.2, et on distinguera les frontières d'*inflow*  $\Gamma_{in}$  (telles que  $\mathbf{v} \cdot \mathbf{n} < 0$ ) où sont imposées des conditions de Dirichlet homogènes ( $\phi|_{\Gamma_{in}} = 0$ ), et les frontières d'*outflow*  $\Gamma_{out}$  (telles que  $\mathbf{v} \cdot \mathbf{n} > 0$ ) où le fluide transporte le polluant à l'extérieur du domaine de calcul, et où sont imposées des conditions de Neumann homogènes ( $\frac{\partial \phi}{\partial \mathbf{n}}|_{\Gamma_{out}} = 0$ ).

A l'instant initial ( $t = 0$ ), le polluant est placé à la position  $(x_0, y_0)$  du domaine de calcul et il est modélisé par :

$$\phi_0(x, y) = e^{-\left(\frac{x-x_0}{\sigma_0}\right)^2} e^{-\left(\frac{y-y_0}{\sigma_0}\right)^2} \quad (\text{B.3})$$

qui correspond à la une gaussienne d'amplitude 1 et de rayon  $\sigma_0$  représentée sur la figure B.3.

Le domaine de calcul  $\Omega = ]a_x, b_x[ \times ]a_y, b_y[$  est représenté par un maillage uniforme centré en cellules ( $\Omega_h = \{\Omega_{i,j}, i = 2, \dots, N_x, j = 2, \dots, N_y\}$ ), de centre  $\mathbf{x}_c(i, j) = (x_c(i), y_c(j))$ , de volume  $V_{i,j} = \Delta x \Delta y$ . Pour imposer les conditions aux limites, des cellules fantômes sont créées sur les parois West ( $\{\Omega_{1,j}, j = 2, \dots, N_y\}$ ), East ( $\{\Omega_{N_x+1,j}, j = 2, \dots, N_y\}$ ), South ( $\{\Omega_{i,1}, i = 2, \dots, N_x\}$ ) et North ( $\{\Omega_{i,N_y+1}, i = 2, \dots, N_x\}$ ).



**Fig. B.3.** Solution initiale et finale du transport d'une gaussienne placée en  $(x_0, y_0) = (1, 1)$  et de rayon  $\sigma_0 = 0.1$ .

**Question 1.** Ecrivez la forme conservative de l'équation d'équilibre de  $\phi$  dans une cellule  $\Omega_{i,j}$  du maillage, en notant  $\eta = \lambda/\rho$  le paramètre de diffusion.

**Question 2.** Faites un schéma du maillage complet en positionnant les variables discrètes  $\{\phi_{i,j}\}$ . Donnez la valeur que prend  $\phi_{i,j}$  dans les cellules fantômes selon le type de conditions aux limites.

**Question 3.** Effectuez une semi-discrétisation de l'équation de convection-diffusion pour  $\phi$  par méthode de volumes-finis centrés en espace. Achetez-la en utilisant le schéma d'Euler progressif (EP1). Ecrivez le schéma complet sous sa forme itérative.

Le programme MATLAB 'Tp\_matlabconv\_diff\_2D.m' crée le maillage  $\Omega_h$ , discrétise la condition initiale (B.3) (écrite dans le programme 'phi\_init.m') aux noeuds  $\mathbf{x}_c(i, j)$  du maillage, et la représente sur un graphe (fonction mesh(X,Y,Z)). Vous allez le compléter en écrivant la boucle en temps du schéma EP1 centré et réaliser une simulation complète de transport du polluant.

**Question 4.** Ecrivez la fonction MATLAB

```
phinew=schema_EP1_centre(advx,advy,eta,dt,dx,dy,ncvx,ncvy,phiold)
```

qui correspond à une itération en temps du schéma d'Euler progressif, tel que :

- les réels advx et advy correspondent aux composantes cartésiennes de la vitesse d'advection  $\mathbf{v}$ , supposée constante.
- Le tableau phiold de taille  $(ncvx + 1) \times (ncvy + 1)$  doit contenir la solution au temps  $t_n$  avec conditions aux limites dans les cellules fantômes.
- Le tableau phinew de taille  $(ncvx + 1) \times (ncvy + 1)$  donnera la solution au temps  $t_{n+1} = (n+1)\Delta t$ .

La simulation que vous effectuerez prendra les paramètres suivants pour le domaine spatial :

$$a_x = 0, \quad b_x = 1, \quad a_y = 0, \quad b_y = 1.$$

La condition initiale prend les paramètres suivants :

$$\sigma = 0.1, \quad (x_0, y_0) = (1/4, 1/4),$$

et les paramètres de l'équation de convection diffusion sont :

$$\eta = 0.01, \quad \mathbf{v} = (1, 1).$$

**Question 5.** *Cette question va permettre de fixer les paramètres de l'intégration en temps.*

- Choisissez un intervalle de temps  $[0, t_f]$  telle que le polluant soit sorti du domaine à  $t = t_f$ .
- Etablissez les conditions de stabilité du schéma EP1 centré (cf. cours) pour les paramètres de la simulation.

**Question 6.** *Programmez dans 'Tp\_matlabconv\_diff\_2D.m' la boucle d'intégration en temps du schéma EP1 centré (conseil : inspirez vous de la boucle `while ... end` du TP2). Vous serez amené à faire une animation de la solution numérique (commandes `getframe` et `movie` du TP2). Des indications sur l'écriture de la boucle en temps sont données dans le programme MATLAB.*

**Question 7.** *Effectuez des simulations pour une valeur de  $\Delta t$  supérieure et inférieure à la condition de stabilité. Qu'observez-vous ? Dans le compte-rendu, montrez le graphe de la solution à des temps bien choisis, et rendez le listing de vos programmes MATLAB.*



## First steps with ANSYS FLUENT: The axisymmetric expansion flow

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### 1 Problem statement

The purpose of this lab exercise<sup>1</sup> is to learn the basics of CFD analysis with the commercial software ANSYS FLUENT<sup>2</sup> by computing the axisymmetric expansion flow presented in the lectures (Fig. 1).

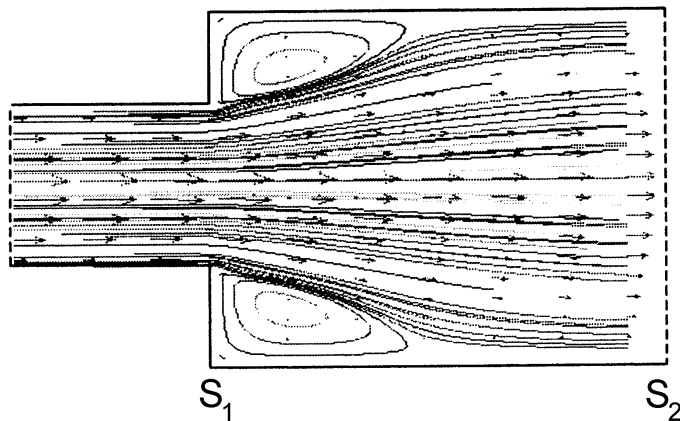


Figure 1: Velocity vectors at the vicinity of the sudden expansion (section  $S_1$  at  $x_1 = 1.0 m$ ). In the upstream pipe the flow is fully developed, then the singularity causes abrupt flow separation, which causes recirculation zones. In the downstream pipe, the flow reattaches at Section  $S_2$  and reaches fully developed regime.

This analysis will be performed with the ANSYS WorkBench platform, which integrates the following softwares:

- Geometry creation (DesignModeler )
- Mesh generation (Ansys Meshing )
- Flow setup and numerical solution (Fluent )

<sup>1</sup>The pdf version of this document and all necessary documents for completing the exercise are available online at the course webpage: *UE 203 - Computational Fluid Dynamics M1 MEPP* at the ARCHE website (<http://arche.univ-lorraine.fr/enrol/index.php?id=9396>) .

<sup>2</sup>This lab exercise has been realized with Version 15.0 of ANSYS WorkBench .

### Work to be performed

This document will introduce you to the basic functionality of the various WorkBench softwares and will guide you through the CFD analysis of the expansion flow.

Read carefully the document and execute with the WorkBench the tasks presented in the document. You should answer to the various questions of the document in a Word report. You will send me your report at the end of the CFD lectures as part of your evaluation.

### 2 The WorkBench platform

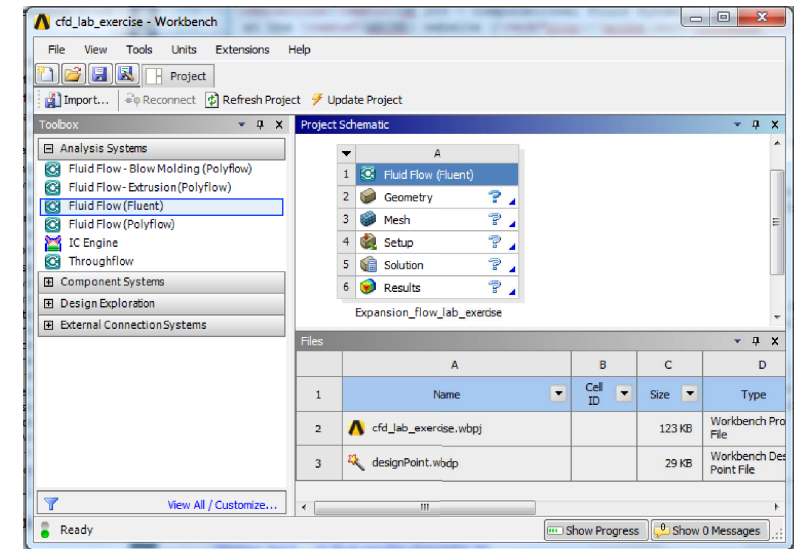


Figure 2: Graphical user interface (GUI) of ANSYS WorkBench .

Launch the ANSYS WorkBench software from your PC and create a **Fluid Flow (Fluent)** project. The **Fluent** workflow shown in Fig. 2 shows the consecutive steps of the CFD analysis. WorkBench will automatically manage the various data files generated by your **Fluent** project.

- Save now the WorkBench project (\*.wbproj file) on your disk with a meaningful and correct names for your file and working directory (no special characters, foreign accents, avoid spacings), or the software may malfunction and your data be lost!
- The language preferences of the various WorkBench softwares can be modified with the option **Regional and Language Options** from the menu bar: **Tools** → **Options...** You need to restart WorkBench to make the change effective.

### 3 Computational domain creation with DesignModeler

According to the symmetries of the flow domain of Fig. 1, we assume that the flow is asymmetric so that it is only required to create the computational domain shown in Fig. 3. These assumptions allow us to do 2D computations instead of 3D, that will save computational time and memory space.

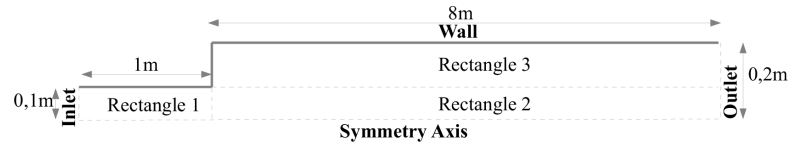


Figure 3: Computational domain and boundary conditions for the cylindrical expansion flow. The 2D computational domain is composed of 3 rectangular surfaces.



For axisymmetric flow computations with ANSYS Fluent, it is important to:

- Create the computational domain in the upper half XY plane (such as  $y > 0$ ) of DesignModeler.
- Ensure that the symmetry axis of the domain is the horizontal axis ( $y = 0$ ).

#### 3.1 Overview of DesignModeler

The graphics interface (GUI) of DesignModeler is shown in Fig. 4. The most relevant components are:

- **Graphics Window:** sketching area where the various geometrical components of the computational domain are viewed and manipulated with the mouse.
- **Menu Bar:** where you can read/write the geometry files, create new sketches, etc...
- **Toolbars:** The most important tools are:
  - **View toolbars** : icons for moving and zooming the geometry with the mouse in the graphics window.
  - **Selection toolbars** : Icons for mouse selection of the various elements of the geometry (vertex, edge, face and volume) in the graphics window.
  - **New sketch** : A “sketch” is a 2D drawing of the contours of the computational domain.
- **Mode Tabs:** there are 2 different modes:
  - **Sketching mode:** where you draw the contours of the geometry with the help of the various sketching tools shown in Fig. 5.

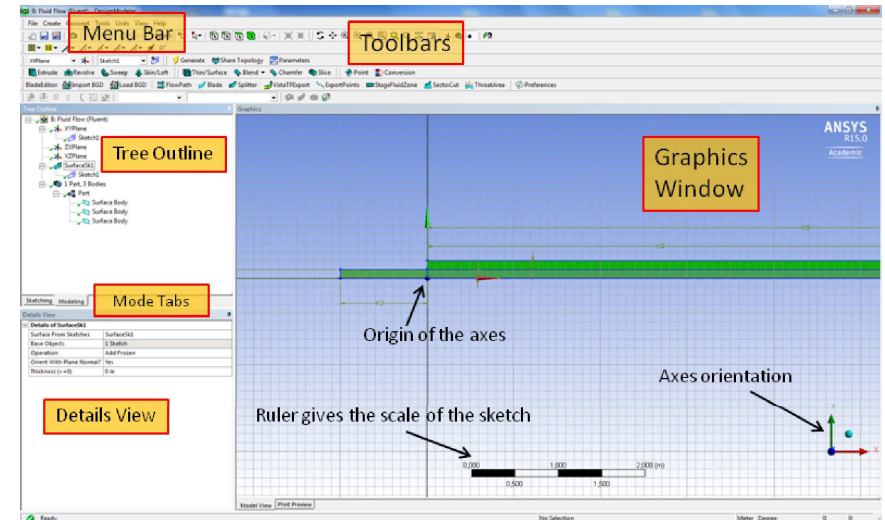


Figure 4: GUI of DesignModeler. Note that 2D geometries should be created in the XY plane. Click on the axes to change orientation of the graphics window.

- **Modeling mode:** that shows the Tree outline, where the basic operations for creating the geometry are shown sequentially: the default planes (where the sketches are stored), the geometry operations on the sketches, the list of created parts.
- **Details View:** where details of the selected geometry (dimensions, constraints, options, etc...) are displayed and can be modified. Note that grey colored details cannot be modified.

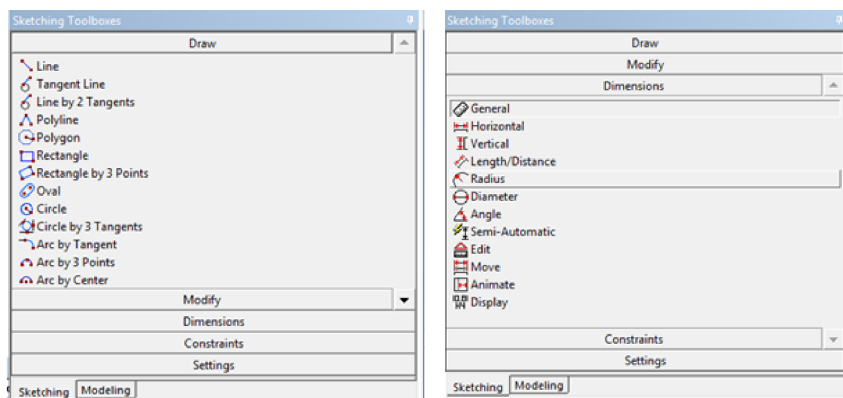


It is important that you memorize the structure of this GUI, since you will use it extensively to do the lab exercises. Note that the GUI of Ansys Meshing is structured similarly.

#### 3.2 Generation of the computational domain

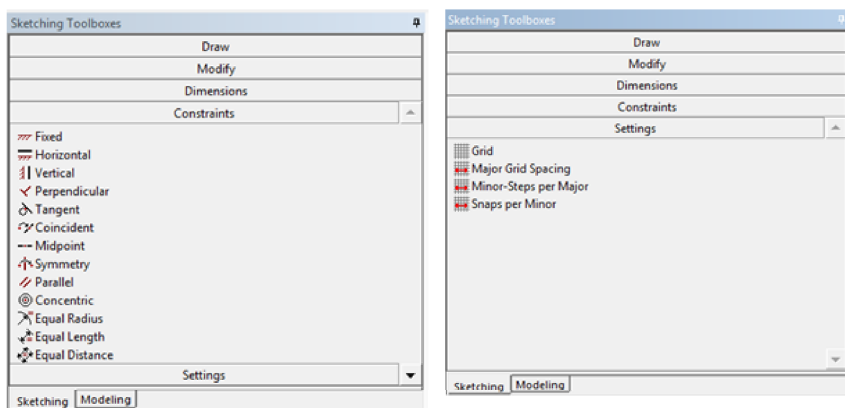
DesignModeler creates the computational domain with the following consecutive steps:

- Step 1) **Sketching mode:** draw the contours of the domain in one (or several) 2D sketches. It is easier to split the computational domain in several elementary entities (like the 3 rectangles of Fig. 3) and draw them separately.
- Step 2) **Modeling mode:** create “solid” bodies (in 2D, it is a surface body) from these contours.
- Step 3) **Modeling mode:** combine these bodies into a unique computational domain (which is called a part in DesignModeler).



(a)

(b)



(c)

(d)

Figure 5: Toolbox available for sketching. (a) **Draw**: tools for drawing elementary 2D entities (lines, rectangles, circles, ...), (b) **Dimensions**: tools for setting the dimensions of the entities (by mouse selection in the graphic windows), (c) **Constraints**: tools for imposing constraints on each entities to assemble the final geometry (alignment with the axes, *etc.*...), (d) **Settings**: tools for displaying grid lines in the graphic window to make sketching easier.

Apply now this strategy to create the domain of Fig. 4. In one sketch, you will draw separately the 3 rectangles shown in this figure, then assemble them as a unique **part** with 3 distinct **surface bodies**.

Note that the Undo/Redo tools are only available in the sketching mode. In the **Modeling Mode**, you can either **Suppress** a body<sup>3</sup> or **Delete** it permanently from the tree outline.

| Details View              |            |
|---------------------------|------------|
| Details of SurfaceSk1     |            |
| Surface From Sketches     | SurfaceSk1 |
| Base Objects              | 1 Sketch   |
| Operation                 | Add Frozen |
| Orient With Plane Normal? | Yes        |
| Thickness (>=0)           | 0 m        |

Figure 6: **Base Objects**: the 3 selected rectangles. **Add Frozen** (in French: **Ajouter Corps Bloqué**): it means that the three surfaces remain independent in the computational domain. You will be able to mesh each surface independantly in **Ansys Meshing**. Note that grey colored details cannot be modified.

- In the **Tree Outline** of the modeling mode, click on **XYPlane** and create a new sketch . Go to the **Settings** toolbox (Fig. 5 (d)) to show the grid and apply adequate parameters for viewing the geometry: **Major Grid spacing**: 1 m, **Minor-steps per Major**: 10.
- In the **sketching mode**, create **Rectangle 1** with the **Draw** toolbox of Fig. 3 and set its dimensions<sup>4</sup>. Repeat for the other 2 rectangles. Align them as shown in Fig. 4 by using the **Constraints** toolbox<sup>5</sup>.
- In the **Tree Outline** of the modeling mode, select the sketch and click on **Concept** → **Surfaces From Sketches** from the menu bar. A surface sketch **SurfaceSk1** is created in the tree outline, whose details are shown on Fig. 6. To finalize the operation click on **Generate**.
- You should have created 3 **parts** made from the 3 **surface bodies** you have drawn (look in the **Tree Outline**). Now, to assemble the 3 bodies in the same **part**, select them all and right-click on **Form New Part**. As a result, you have created the computational domain as 1 part made of 3 surfaces bodies, as in the tree outline of Fig. 4.
- The last step is to define the computational domain as “fluid”. To do this, set **Fluid/solid** to be **Fluid** in the **Details View** of the part.

If an element is highlighted with a yellow bolt in **WorkBench**, it means that the software waits for the user to update/generate this element.

<sup>3</sup>It will be ignored when the geometry is generated.

<sup>4</sup>Define them with **Dimensions** toolbox and set their values in the **Detail View**.

<sup>5</sup>Use the **Coincident** constraint to stitch the edges of each rectangle to the X or Y axes.

**Question 1.** Once the computational domain has been created, display it in your Word report and exit DesignModeler .

When you exit the software, the computational domain is automatically saved on disk in the DesignModeler database format (\*.agdb), and ready to be used by the next software in the WorkBench workflow. You can view all files generated during the lab exercise (and their location on your hard-disk) by toggling **View** → **Files** in the WorkBench menu bar.

## 4 Mesh generation with Ansys Meshing

Open now Ansys Meshing from the WorkBench . The GUI, shown in Fig. 7, is similar to DesignModeler with the same mouse functionality. In this first lab exercise, we will only consider the generation of the most simple case of quadrilateral mesh, *i.e.* the Cartesian mesh (see the CFD lectures). Ansys Meshing is able to generate a mesh automatically, with limited user's input. For example, after clicking on the Generate Mesh tool, you should have to create the uniform mesh shown in Fig. 7. The only user input needed is the cells' edge size (Max Face Size in the Details View of the mesh).

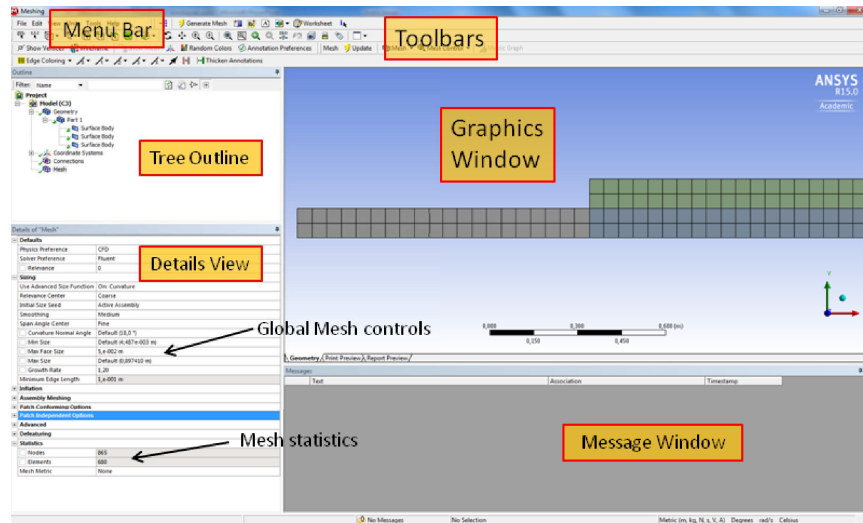


Figure 7: GUI of Ansys Meshing . The quad mesh shown in the graphics window has been generated with Max Face Size=0.05m and all other parameters set to default values.

**Question 2.** Try generate meshes with various cell sizes. Display the resulting meshes in your Word report with the corresponding number of cells (Elements) and vertices (Nodes) given in the Statistics detail view of the mesh.

### 4.1 Generation of a graded Cartesian mesh

The flaws of the “automatic” mesh of Fig. 7 is that the user does not control the mesh resolution in the vicinity of strong flow gradients, which are located in the boundary layers along solid walls and in the recirculation zone (see Fig. 1). Ansys Meshing allows various strategies for refining locally the mesh in selected areas of the computational domain. In this lab exercise, we will use the “bottom up” strategy to create a better quality mesh: first introduce a 1D meshing of the edges of the domain, then use a quad method to mesh the computational domain (which is called a Face, since it is 2D). The resulting mesh, shown in Fig. 8, is generated by the following steps:

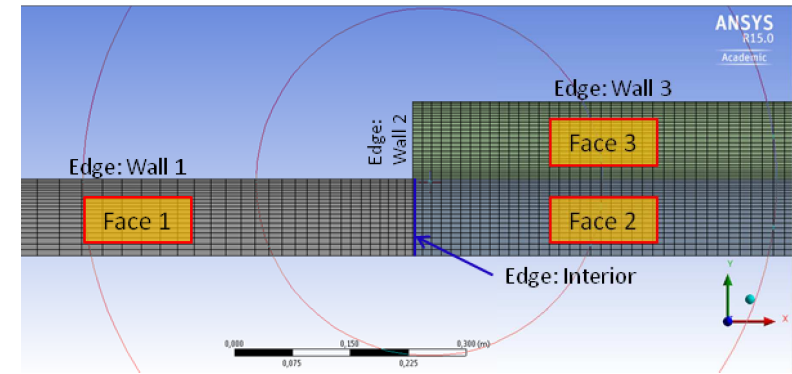


Figure 8: Zoom of the mesh near the singular expansion. Note that the mesh is refined near the walls and the salient corner. This mesh has 14300 cells.

- 1) Edge meshing (Tool for edge selection): In the Graphics Window, select one of the edges of Face 1 and insert an edge sizing method<sup>6</sup>. The Edge Sizing method appears in the Tree Outline of the mesh. Choose Type: Number of Divisions and enter the details given in Table 1. Repeat for the 3 other edges of Face 1.

| Edge           | Interior  | Wall 1    | Wall 2    | Wall 3    |
|----------------|-----------|-----------|-----------|-----------|
| # of divisions | 20        | 40        | 25        | 300       |
| Bias type      | - - - - - | - - - - - | - - - - - | - - - - - |
| Bias factor    | 3         | 5         | 2,5       | 4         |

Table 1: Sizing details for the edges shown in Fig. 8. Number of Divisions corresponds to the number of edge vertices  $N$ . Bias Type enables to create a graded mesh with geometric refinement of the form  $\Delta x_{i+1} = R\Delta x_i$ , where Bias Growth Rate  $R$  is the ratio between adjacent cells. Examples of graded meshes are given in Fig. 9. Bias Option: the option Smooth Transition needs the value of  $R$ , otherwise Bias Factor corresponds to  $\Delta x_{\max}/\Delta x_{\min} = R^{1/(N-1)}$ .

<sup>6</sup>Right-click on the moose: Insert → Sizing.



Figure 9: Examples of graded mesh with geometric refinement. At left: the mesh is refined on the west boundary only. At right: refinement at both west and east boundaries.

- 2) Face meshing (Tool  for face selection:) Complete the meshing of Face 1 by inserting a Mapped Face Meshing with Method: **Quadrilaterals**. Generate the mesh of Face 1 by right-clicking on **Generate Mesh on Selected Bodies**.
- 3) Preview the mesh by clicking on **Mesh** in the **Tree Outline**. If the mesh density is not similar to Fig. 8 or the grid lines are not parallel to the axes, you certainly made a mistake in some edge sizing details. In this case, delete the mesh<sup>7</sup>, review your edge sizings and generate again the mesh of Face 1.
- 4) Repeat these steps for Face 2 and 3.

## 4.2 Definition of the boundary conditions

Now that the meshing of the computational domain is completed, the last step before starting the CFD analysis with the **Fluent** solver is to label the boundary conditions as in Fig. 3. To perform this, click on **Model** in the **Tree Outline**, then select a boundary edge in the **Graphics Window** and insert a **Named Selection**<sup>8</sup>. In the **Tree Outline**, rename it as shown in Fig. 3. The label of the boundary edges will be recognized by **Fluent** and the corresponding boundary condition will be applied.

-  For an easy reviewing of the boundary conditions in **Fluent** by the user (think that an actual 3D mesh may have hundreds of boundary edges), it is worthwhile to group all edges with the same boundary conditions (*e.g.* the 3 edge walls) in a single named selection.

**Question 3.** Once the mesh is created and the boundary conditions labeled, display the mesh in your Word report, give the mesh statistics and exit **Ansys Meshing**.

## 5 CFD analysis with Fluent

Open **Fluent** from **WorkBench**. The **Fluent** launcher panel will open. Select **Dimension**: 2D, **Options**: **Double Precision** and leave the default settings for the other options. Note that:

- **Double precision**: if you choose this option, floating point operations will be made on 8 bits, which yields numerical results with a 16-digit precision. Using the default **Single Precision** (4 bits) will reduce the memory usage and yield faster computations, but will give less accurate results (8-digit accuracy) and the convergence of the iterative solution may be impacted. It is thus recommended to use the **Double Precision** solver.

<sup>7</sup>Right-click on **Mesh** in the **Tree Outline** and select **Clear Generated Data**. This operation only deletes the mesh, not the meshing methods you have defined.

<sup>8</sup>Right-click on the moose: **Insert** → **Named Selection**.

- **Serial/Parallel**: if your PC has a single processor, choose **Serial** (the default). If your PC has several processors (or cores), **Parallel** computations amounts to partitioning the mesh such that the computation on each mesh part is performed by a different processor. The interest of parallel computations is that the computational time of a serial computation is (ideally) divided by the number of processors used.

## 5.1 Overview of Fluent

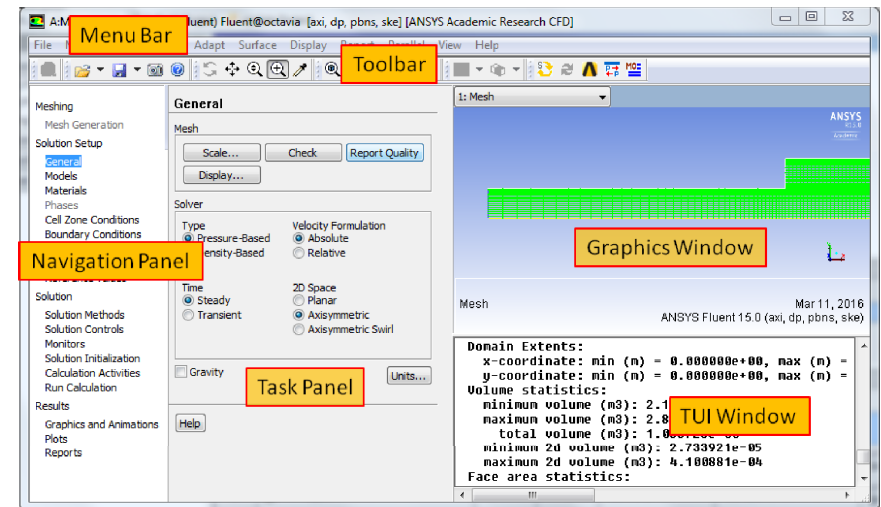


Figure 10: GUI of Fluent.

The graphics interface (GUI) of **Fluent** is shown in Fig. 10. The most relevant components are:

- **Toolbars**: As in **DesignModeler** and **Ansys Meshing**.
- **Graphics Window**: Graphical results are displayed here. The default mouse button functionality is<sup>9</sup>:
  - Left button translates.
  - Middle button zooms.
  - Right button selects/probes.
- **Navigation Panel**: The main stages of the CFD analysis are displayed here as items, organized as a tree from top to bottom:

<sup>9</sup>Note that they are different from the other softwares in **WorkBench**. They can be changed from **Display** → **Mouse buttons...** in the menu bar.

- **Solution Setup:** define the CFD problem, such as the fluid models (laminar, turbulent, heat transfer, multiphase, ...), type of materials (water, gas, solid, ...), boundary conditions.
- **Solution:** define the numerical inputs (such as discretization schemes, solution algorithms, initial condition, ...) and solve the problem.
- **Results:** once the solution is calculated, analyze (postprocess) the results.
- **Task Panel:** Selecting an item in the **Navigation Panel** opens the corresponding dialog box in this panel.
- **Menu Bar:** Additional items can be accessed from the menu bar.
- **TUI<sup>10</sup> Window:** display text results of the computation.<sup>11</sup>

## 5.2 Input/output files

There are 2 types of data files generated by **Fluent**.

- **Case File (\*.cas)** : stores the mesh and the setup parameters.
- **Data File (\*.dat)** : stores the last computed solution.

You can manage these files from the menu bar **[File]** → **Import** or **Export**. You can also import a mesh generated from a third-party meshing software. Since you are using **WorkBench**, the mesh generated previously with **Ansys Meshing** is imported at the start of **Fluent**, while **Case** and **Data** files are automatically managed when you start or exit **Fluent**.

## 5.3 Solution setup

All the **Fluent** tasks for set-up of the CFD problem are done in the **Solution Setup** panel. **Fluent** has predefined default settings for the various stages of the set-up, but it is important to set them correctly for the problem under consideration. Thus, it is important to have knowledge of the physics involved in your CFD problem for selecting the relevant flow simplifications, models and equations to be solved.

**Question 4.** We consider the flow of water ( $\rho = 998.2 \text{ kg/m}^3$ ,  $\mu = 0.001003 \text{ kg/ms}$ ) with inlet velocity  $u_{\text{inlet}} = 1.5 \times 10^{-3} \text{ m/s}$ . Compute the Reynolds number in the upstream pipe (based on its diameter). Infer the flow regime and the relevant flow simplifications.

### Check the mesh and select the CFD solver: **General** task page

It is important to check the suitability of the imported mesh for **Fluent** computations. Do it by clicking on **(Check)** on the Mesh subpanel. It is also possible to rescale the mesh.

<sup>10</sup>TUI: Text User Interface.

<sup>11</sup>For advanced user: all the tasks defined from graphic panels have a corresponding TUI command, in order to be able to run **Fluent** in batch mode for automating the solver set up/solution/postprocess stages with scripts and journal files.

As noted in the lectures, there is no universal numerical method that works for all flow regimes: incompressible flow are computed with the **Pressure-based Solver**, while the **Density-Based Solver** is used for compressible flows. The other options are related to the problem simplification mentioned in Section 1: **Time: Steady**, **2D Space: Axisymmetric**.

### Set up the flow physics: **Models** task page

The **Pressure-based** solver is configured with the relevant flow physics in the **Models** task page: turbulence models, flow with heat transfer (**Energy**, **Radiation**), combustion (**Species**), multiple phases (**Multiphase**), .... It is of the utmost importance to set up the form of the viscous tensor in the Navier-Stokes equations (**Viscous** item). Click on the **(Edit...)** button to review the various **Viscous** options:

- **Inviscid** : viscosity is assumed to be zero, which corresponds to the Euler equations, mainly used for external compressible flows.
- **Laminar** : corresponds to the classical Navier-Stokes equations with the Cauchy stress tensor. The form of the viscosity (constant, non-Newtonian or temperature dependent) is defined later with the material properties of the fluid.
- Other options are used for turbulent computations, and correspond to choosing the modified viscous tensor resulting from RANS<sup>12</sup> modeling of the Navier-Stokes equations. We can mention the well-know **k-epsilon** model which is widely used in CFD: using this model will add 2 transport equations (for the turbulent kinetic energy  $k$  and dissipation  $\epsilon$ ) to the NS equations. The parameters for these transport equations can be set from these this panel.

By default, **Fluent** uses the laminar model with all other models (**Multiphase**, **Energy**, *etc.*...) disabled.

### Set up the material properties: **Materials** task page

Materials refer to the various fluids or solids involved in your CFD problem. The materials available for your simulation are listed on the **Materials** task page. By default, **Fluent** selects the **Fluid** material as **air** with constant properties (density and viscosity). Click on **(Create/Edit...)** to add **water-liquid** to the list of available materials (use the **Fluent Database**, where a wide range of materials are predefined). You can edit the material properties if needed, for example if you wish to use a non-Newtonian fluid with shear-thinning viscosity.

### Set up the cell zone conditions

As stated in the lectures, a zone corresponds to a portion of the mesh defined as **fluid**, **solid** or **porous**, where distinct materials can be defined<sup>13</sup>. These zones have been defined previously in **DesignModeler** : in this first lab exercise, there is only one fluid zone filled with water. Click on

<sup>12</sup>RANS : *Reynolds Averaged Navier-Stokes Equations*, obtained through statistical modeling of turbulent flows.

<sup>13</sup>As an example, the computational domain for a conjugate heat transfer problem would be divided into fluid and solid zones.

**Edit...** in the **Cell Zone Conditions** task page to set this zone as **water-liquid** (by default **Fluent** fills all fluid zone with air).

*For advanced users:* In the **Fluent** pressure-based solver, the computed pressure (the **Gauge Pressure**  $p_{\text{gauge}}$ ) is defined up to a constant level (the **Operating Pressure**  $p_{\text{operating}}$ ), such that:

$$p(\mathbf{x}) = p_{\text{operating}} + p_{\text{gauge}}(\mathbf{x}). \quad (1)$$

The operating pressure can be set at a reference location in the domain by clicking on **Operating Conditions...**. By default, **Fluent** sets  $p_{\text{operating}}$  as the ambient pressure at the origin of the computational domain, and it is usually not necessary to change it. In addition, if it is required by your CFD problem, you can add gravity to your model with this item.

Note that the value of the operating pressure has no influence on the computed results since only pressure differences are relevant in incompressible flows.

### Set up the boundary conditions (BCs)

This task page defines the BCs settings at the boundary edges (in 2D) of the computational domain. The **Zone** panel displays the name of the boundary edges as you previously labelled them in **Ansys Meshing**<sup>14</sup>. You can visualize these edges by clicking on **Display Mesh...**. **Fluent** usually recognizes these labels and assigns them with a corresponding **BC Type**. Make sure the inlet boundary is defined as **velocity-inlet**, the symmetry axis as **axis**, and the outlet as **pressure-outlet** (see Fig 1). Some BCs needs additional settings which are defined by clicking on **Edit...**. For this lab exercise, you only have to set the inlet velocity given by Question 1.

### 5.4 Set up the numerical parameters of the computation

Now that the physical model, solver and boundary conditions have been set, the next step is to define the numerical parameters (algorithm for pressure-velocity coupling, spatial discretization, convergence criteria) of the pressure-based solver: this is performed with the **Solution** panel.

By default, **Fluent** chooses “robusts” numerical parameters, which yield converged solutions for a wide range of flows. Henceforth, the default discretization methods are usually low-order and dissipative (for stabilizing the solution) and should be modified to get accurate results.

### Set up the discretization methods: **Solution methods** task page

In this task page (shown in Fig 11) you first set up the iterative algorithm for solving the pressure-velocity coupling (SIMPLE<sup>15</sup> algorithm and its variants, see lecture notes). Use **SIMPLE** for steady problems such as this lab exercise (**PISO** is recommended for transient problems). Define next the various finite-volume (FV) discretization schemes for the NS equations<sup>16</sup> (**Gradient**: cell-face interpolation of velocity gradient used in the discretization of the viscous term and high-order discretization of the convection term, **Momentum**: discretization of convection term, **Pressure**: interpolation of cell-face pressure). The various pros and cons of each scheme has been discussed in the lecture notes. As a reminder:

<sup>14</sup>If some of the boundary edges are missing, it may be caused by a bug in the **WorkBench**. Ask advice to your instructor.

<sup>15</sup>Acronym for *Semi-Implicit Method for Pressure-Linked Equations*.

<sup>16</sup>When applicable, discretization schemes of the additional transport equations need to be defined too. Examples will be given for the turbulent computation of Section 6.

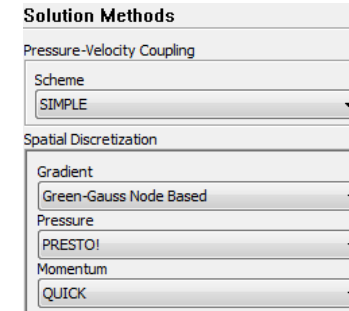


Figure 11: Recommended discretization methods for the expansion flow.

- Use **Green-Gauss Node-Based** interpolation for meshes with only tri/tet cells.
- Use **PRESTO!** for pressure interpolation or the solution may have convergence problems.
- For convection, use **First Order Upwind** only at the initial stages of the computation or if the solution exhibits convergence problems. Otherwise, second-order schemes such as **QUICK** (more accurate and less diffusive) are preferred.

### Set up the relaxation parameters of the iterative algorithm: **Solution Controls** task page

*Background on iterative algorithms for solving a system of algebraic equations:* for a (non-) linear system of the form:

$$\mathcal{A}\Phi = RHS, \quad (2)$$

where  $\Phi$  is the vector containing velocity or pressure mesh values, the matrix  $\mathcal{A}$  and vector  $RHS$  represent the FV discretization of the NS equation. Starting from the initial condition  $\Phi^0$  given by the user, an iterative algorithm seeks the solution as the series  $\{\Phi^k, k \geq 1\}$  such that:

$$\Phi^k = \Phi^{k-1} + \alpha \delta\Phi, \quad (3)$$

where  $k$  is the iteration number and  $\alpha \in [0, 1]$  is the **Under-Relaxation Factor**. The value of  $\alpha$ , chosen by the user, has impact on the rate of convergence of the algorithm. The closer is  $\alpha$  to 1 the faster is the convergence, but non-convergence (or solution blow-up) may occur because the lag  $\alpha \delta\Phi$  is too large. In the latter case, the user should decrease (*i.e.*, under-relax) the value of  $\alpha$  to get convergence, even though it may slow down the algorithm.

The under-relaxation factors are set-up in the **Solution Controls** task page for each flow equation. The default values given by **Fluent** are *robusts*, *i.e.* they give a solution for a wide range of problems. There is no need to modify them unless your computation undergoes convergence problems.



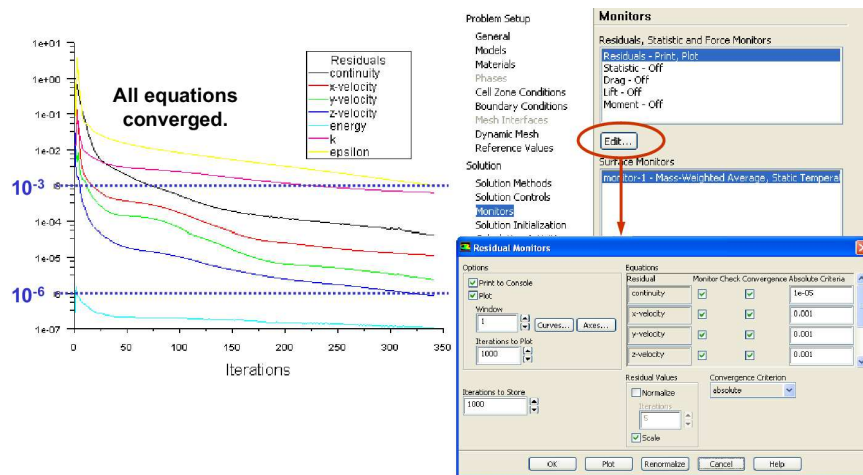


Figure 12: Example of residual monitoring plot for a turbulent flow computation.

## Monitoring the convergence of the iterative algorithm: Solution controls task page

Fluent considers that the iterative algorithm has converged when the residual  $R^k$  of Eq. 2, defined by  $R^k = RHS - A\Phi^k$ , falls below the convergence criterion:

$$\|R^k\| < \epsilon, \quad (4)$$

where the value of the tolerance threshold  $\epsilon$  is given by the user.

The set up of the convergence monitors for each scalar equation of the NS system is performed in the **Monitors** task page. By default, **Fluent** defines the convergence criteria to be  $\epsilon = 10^{-3}$  for all flow equations<sup>17</sup>. This value is usually not low enough to give accurate results. To change this criteria, click on **Edit...** in the **Residuals** panel (see Fig. 12). During the computation, the evolution of the residuals will be shown in the graphics window as in Fig. 12 if you choose the **Plot** option.

## 5.5 Perform the computation

Computations are performed in the **Solution** panel. You need first to define the initial condition  $\Phi^0$  of the iterative algorithm<sup>18</sup>. This is performed in the **Solution Initialization** task page. **Fluent** proposes several **Initialization** Methods for defining suitable initial conditions, notably from the values defined at the boundary conditions. For this lab exercise, you will start the algorithm from rest ( $\mathbf{v} = 0, p = 0$ ). To do this, choose **Standard Initialization** in the task page, **Compute From: [all-zones]**, and press **Initialize**.

To start the iterative algorithm, go to the **Run Calculation** task panel, enter **Number of Iterations: 10** and press **Calculate**. As a result, **Fluent** will write (in the TUI) and plot (in the graphics window) the residuals at each iterations<sup>19</sup>. Since after only 10 iterations the convergence criteria (4) is not met yet, set a larger number of iterations (say 1000) and click on **Calculate** to continue the iterative algorithm: **Fluent** will automatically stops the algorithm when convergence is reached.

Note that **Fluent** stores in memory the solution at the last iteration only. If you have pressed **Calculate** again, you will continue the algorithm from the latest saved iteration, i.e. the tenth. If you want to restart the algorithm from rest, you should do the **Solution Initialization** step first before clicking on **Calculate**.

You are now able to perform by yourself a full computation of the flow:

**Question 5.** Perform a computation of the expansion flow with the convergence threshold  $\epsilon = 10^{-5}$  for all flow equations. Display the residual monitoring plot in your Word report<sup>a</sup>. Save the converged solution to your PC disk (see Section 5.2 for details): give the name water\_inlet=0,0015 to the \*.cas and \*.dat files.

<sup>a</sup>For exporting a Fluent plot, right-click at the top of the graphics window and select **Copy to Clipboard**.

For advanced users: to perform periodically a back-up copy of the solution during the iterative algorithm, go to the **Calculation Activities** panel and enter the frequency back-up in the **Autosave Every (Iterations)** group box.

## 5.6 Postprocessing: Results panel

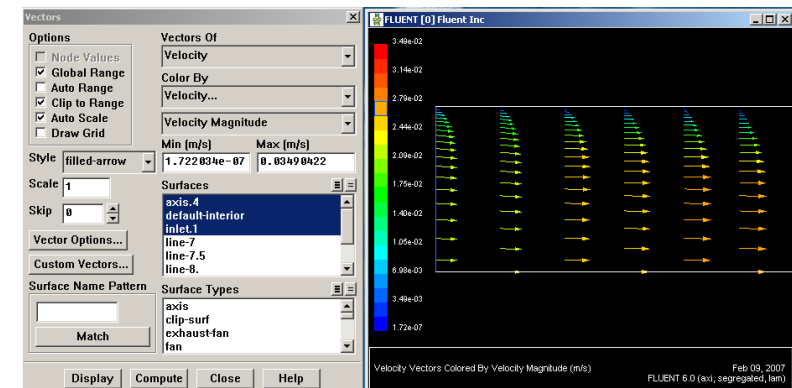


Figure 13: At left: Vectors dialog box. At right: 2D visualization of the velocity vectors upstream of the contraction.

<sup>19</sup>The frequency of the residual plots and writes can be changed with the **Reporting Interval** option.



Once the converged solution is obtained, **Fluent** proposes numerous postprocessing tools to analyze the solution in order to extract relevant flow information. This is performed under the **Results** panel. There are 3 main types of analysis:

- 2D (or 3D) solution visualization (**Graphics and Animation** task panel): for plotting the velocity vectors, contour levels of a flow variable (velocity, pressure) inside the computational domain.
- 1D visualization (**Plots** task panel): for profile plotting of the solution on selected sections (or lines) of the computational domain.
- Numerical reporting of mean flow quantities (**Reports** task panel): such as mass flow through selected sections, forces at boundaries, ...

### Velocity vector visualization: Graphics and Animation task panel

Choose **Vectors** in the **Graphics** group box and click on **Set Up...**. The **Vectors** dialog box of Fig. 13 (left) opens. From the **Surfaces** drop-down list choose **interior** to display the velocity vectors in the whole computational domain<sup>20</sup>, and press **Display**. You can use the mouse to zoom in/out in the graphics window.

**Question 6.** Zoom in the computational domain to observe the recirculation zone of the flow (see Fig. 1). Use the **Scale** options of the **Vectors** dialog box if the vectors are too small. Display this plot in your Word report.

### Isocontour visualization: Graphics and Animation task panel

Choose **Contours** in the **Graphics** group box and click on **Set Up...** to open the **Contours** dialog box. Select a flow variable in the **Contours** of drop-down list and press **Display**.

Most flow variables are predefined in **Fluent** with intuitive names. For example, for asymmetric flow:

$$u = \text{Axial Velocity}, \quad v = \text{Radial Velocity}, \quad ||v|| = \text{Velocity Magnitude}.$$

For pressure, note that **Static Pressure** and **Absolute Pressure** correspond respectively to  $p_{\text{gauge}}$  and  $p$  in Eq. (1), while the **Total Pressure** is :

$$p_{\text{total}} = p_{\text{gauge}} + \frac{1}{2}\rho ||v||^2. \quad (5)$$

By default, **Fluent** plots 20 iso-values between the min. and max. values of the visualized variable. In the **Contours** dialog box, you can set the number of levels with the **Levels** option, and modify the min. and max. levels by deselecting the **Auto Range** option and choosing manually the **Min** and **Max** values.

**Question 7.** Display the isobaric lines of the flow near the contraction. Export the graphics to your Word report. At what distance downstream of the singularity we get the isobaric lines of the Hagen-Poiseuille flow in a straight pipe ?

<sup>20</sup>Alternatively, you can visualize the flow on selected surfaces and boundaries from the **Surfaces** drop-down list.

### Velocity profiles on sections of the computational domain

You will learn now how to plot velocity vectors on selected surfaces (lines in 2D) of the computational domain, for example on the  $x = 7$  m section<sup>21</sup> near the outlet boundary, in order to verify that the flow is full developed there. To do this, select **Surface** → **Line/Rake...** from the menu bar to create the line  $x_0 = 7.0$ ,  $y_0 = 0$ ,  $x_1 = 7.0$ ,  $y_1 = 0.2$ . Give the line a meaningful label, such as **line-7.0** in **New Surface Name**. The surface **line-7.0** will now be available in the **Surfaces** drop-down list of any **Graphics and Visualization** dialog box.

**Question 8.** Plot the velocity profiles on Sections  $x = 6.0, 6.5, 7., 7.5$  and **pressure\_outlet.\*** (Give them a meaningful name!). Check that the Poiseuille velocity profile is obtained on these sections. Export the graphics to your Word report.

### Plot 1D profiles

The plotting of 1D profiles of predefined flow variables (axial or radial velocity, static or total pressure, ...) is performed on the **Plots** task panel. Choose **XY Plot** in the **Plots** group box and click on **Set Up...** to open the **Solution XY Plot** dialog box. For plotting along the radial direction  $y$ , choose the direction vector  $X = 0$ ,  $Y = 1$ ,  $Z = 0$  in the **Plot Direction** group box.

**Question 9.** Plot the velocity and (static) pressure along the sections of Question 8. Export the graphics to your Word report. Do these profiles correspond to a fully developed Poiseuille flow ?

### Define custom field functions

It is possible to define new flow variables (named “custom field functions”) from **Fluent** predefined variables, by using the “pocket calculator” provided by **Define** → **Custom Field Functions...** from the menu bar. This new variable can be visualized as previously.

**Question 10.** Define the new function **tau\_xy** that calculates the shear stress in cylindrical coordinates :

$$\tau_{xy} = \mu \left( \frac{\partial v}{\partial x} + \frac{\partial u}{\partial y} \right).$$

Plot the shear stress profile on the vertical sections of Question 8 and along the upper wall. Export the graphics to your Word report.

### Calculate mean values

For example, you can calculate the mean flow velocity in Section  $x = 8$  m of the pipe. In the **Reports** task panel, open the **Surface Integrals** dialog box and choose **Area Weighted Average** in the **Report Type** drop-down list.

**Question 11.** Verify that the mean flow velocity is constant on the sections of Question 8. Report these values in your Word report.

<sup>21</sup>Remember that the origin  $x = 0$  correspond to the section of the abrupt expansion.

## 5.7 Application : calculation of standard pressure drop

In a duct of diameter  $D$ , the pressure (or head) loss due to viscous friction between sections  $x_1$  and  $x_2$  reads :

$$H_1 - H_2 = \lambda \frac{\bar{V}^2}{2g} \frac{x_2 - x_1}{D}, \quad (6)$$

where  $\lambda$  is the (dimensionless) friction factor,  $\bar{V}$  is the mean flow velocity,  $g$  is the gravity,  $H_i = \bar{p}_{\text{tot}}(x_i)/\rho g$  is the mean head at Section  $x_i$ . This formula is valid when the streamlines are parallel to each others, *i.e.* when the flow is fully developed in the pipe.

**Question 12.** According to the previous questions, the flow is fully developed between the zones  $6.5 \lesssim x \leq 8$  of the pipe. Use **Fluent** to calculate the total pressure in various sections of the fully developed region, and give the corresponding value of  $\lambda$ . Compare your results with the empiric relations obtained in a straight pipe:

- Poiseuille formula for laminar flows,  $Re_D \lesssim 2000$  :

$$\lambda = \frac{64}{Re_D},$$

- Blasius formula for turbulent flows,  $4 \times 10^3 \lesssim Re_D \lesssim 10^5$  :

$$\lambda = 0.3164 Re_D^{-1/4}.$$

## 6 Numerical computations in the turbulent regime

### 6.1 Performing the computation

We consider now flow of water with inlet velocity  $u_{\text{inlet}} = 1.5 \times 10^{-1}$  m/s.

**Question 13.** Calculate the Reynolds number. Deduce the flow regime and the flow equations to solve.

If a turbulence model is needed, use **k-epsilon Standard**. In addition to the averaged Navier-Stokes equations, transport equations for  $k$  et  $\epsilon$  are now solved: you needed to define their solution settings, in particular:

- For inlet BCs, choose **Intensity and Hydraulic Diameter** in the drop-down list Turbulence Specification Method, and set a turbulent intensity of 5% and an hydraulic diameter of 0.2 m.
- Use 2nd-order discretization for the convective terms of these equations.
- Define the tolerance monitoring parameters for these equations.

**Question 14.** Starting from rest, compute the solution with residual threshold  $\epsilon = \times 10^{-6}$  for all equations. Export the residual history to your Word report.



When performing a high-Reynolds number with impulsive start from the rest, the iterative solution may undergo convergence problems: the residuals do not decrease monotonously, stall or increase until solution blow-up. If this happen, you need to modify the default solver settings for obtaining convergence. The “optimal” parameters for convergence are usually problems dependent, and fine-tuning of these parameters requires expertise and trial-and-errors. General guidelines for convergence troubleshooting are given in the lecture notes. Here is a quick summary:

- Initialize the solution with a previously calculated solution at lower Reynolds number, then gradually increases the inlet velocity.
- Use first-order methods in the initial stages of the iterative algorithm, then use more accurate methods.
- Decreases the values of the under-relaxation factors of the equation(s) that cause(s) difficulties (usually, it is the continuity equation).
- Improve the mesh quality and use finer cells in zones of high flow gradients.

However, for this lab exercise you shouldn't experience convergence problems: After an initial stage where the residuals seem to be stalled, the solution should converge after a few thousand iterations. Note however that this behaviour may depends on the discretization parameters you chose or the version of **Fluent** you are using.

### 6.2 Analyze the results

#### General flow features

**Question 15.** Verify that the flow is fully developed at the outlet. Display in the 2D domain the velocity profiles near inlet (several sections in the upstream pipe) and outlet (sections of Question 8). Compare with the laminar velocity profiles obtained previously. Report these results in your Word report.

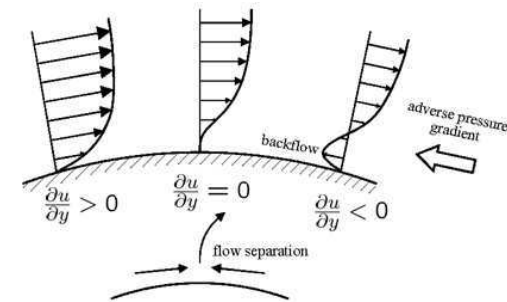


Figure 14: Velocity profiles at a wall boundary layer near a separation point.

### Calculation of singular head losses

At the vicinity of a geometrical singularity such as the sudden expansion at  $x_1 = 0 \text{ m}$  (cf. Fig. 1), the flow is no more unidirectional and the head loss equation (6) is not valid anymore. Experiments show that singular head losses are proportional to the mean kinetic energy:

$$\Delta H = \frac{\xi}{2g} \bar{V}^2,$$

where  $\xi$  is a dimensionless coefficient that depends on the type of singular geometry. For singular expansion flows, the value of the head-loss coefficient  $\xi$  can be estimated analytically at high flow rates with the following assumptions:

- Losses due to wall friction are negligible,
- The streamlines are parallel through Section  $S_1$  ( $x_1 = 1.0 \text{ m}$ ) and Section  $S_2$  (cf. Fig. 1), where the flow returns to fully developed regime. This implies that pressure distribution is uniform at these sections.

With these hypotheses, conservation of mass and momentum between  $S_1$  and  $S_2$  yields:

$$H_1 - H_2 = \frac{\xi}{2g} \bar{V}^2(x_1), \quad \text{with} \quad \xi = \left(1 - \frac{S_1}{S_2}\right)^2. \quad (7)$$

This formula can be found in any fluid dynamics textbook.

**Question 16.** *From the head losses given by your numerical solution, calculate the head-loss coefficient  $\xi$  and compare with the analytical value. The difficulty is to estimate the abscissa  $x_2$  where the streamlines do reattach. To estimate  $x_2$ , you may use the property that the shear stress is zero at a separation point (where flow separates or reattaches, see Fig. 14): the value of  $x_2$  is then obtain by plotting the computed shear stress along the upper wall (cf. Question 10).*