### Modular Programming (PHY1610 Lecture 4)

Ramses van Zon

Winter 2022



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### Section 1

### Modularity



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# Why does modularity matter?

Modularity? Who cares?

- Scientific software can be large, complex and subtle.
- If each section uses the internal details of other sections, you must understand the entire code at once to understand what the code in a particular section is doing.

(This is why global variables are bad bad bad!)

• Interactions grow as (number of lines of code)<sup>2</sup>.



# Example: Monolythic code for hydrogen's ground state

```
#include <iostream>
#include <fstream>
const int n = 4913;
double m[n][n], a[n], b = 0.0;
void pw() {
    double q[n] = \{0\};
    for (int i = 0; i < n; i++)
        for (int j = 0; j < n; j++)
            q[i] += m[i][j]*a[j];
    for (int i = 0; i < n; i++)
        a[i] = a[i]:
double en() {
    double e = 0.0, z = 0.0;
    for (int i = 0; i < n; i++) {
        z += a[i]*a[i]:
        for (int j = 0; j < n; j++)
            e += a[i]*m[i][j]*a[j];
    ጉ
    return b + e/z;
```

```
int main() {
    for (int i = 0; i < n; i++) {</pre>
        a[i] = 1.0;
        for (int j = 0; j < n; j++) {</pre>
            m[i][j] = H(i,j,n);
    for (int i = 0; i < n; i++)
        if (m[i][i] > b)
            b = m[i][i]:
    for (int i = 0; i < n; i++)
        m[i][i] -= b;
    for (int p = 0; p < 20; p++)
        ;()wq
    std::cout<<"Ground state energy="<<en()<<"\n";</pre>
    std::ofstream f("data.txt"):
    for (int i = 0; i < n; i++)
        f << a[i] << std::endl;</pre>
    std::ofstream g("data.bin", std::ios::binary);
    g.write((char*)(a), sizeof(a));
    return 0;
```

# What is wrong with that code?

The hydrogen.cpp code uses functions. Is that not modular? A few bad things:

- Global variables that all of the code can modify.
- All code in one file.
- No comments.
- Not clear what part does what, or what part needs which variables.
- Cryptic variable and function names.
- Hard-coded filenames and parameters.
- Automatic arrays.

Who cares, you might say, as long as it runs? But:

- Code is not written for a computer but for humans.
- Code almost never a one-off.

# What to do: Use modularity

- You must enforce **boundaries** between sections of code so that you have self-contained modules of functionality.
- This is not just for your own sanity. There are added benefits:
- Each section can then be tested individually, which is significantly easier.
- Makes rebuilding software more efficient.
- Makes version control more powerful.
- Makes changing the code easier.



## But it's more work up-front

- Think about the **blocks of functionality** that you are going to need.
- How are the routines within these blocks going to be used?
- Think about what you might want to use these routines for; only then design the interface.
- The interfaces to your routines may change a bit in the early stages of your code development, but if it changes a lot you should stop and rethink things you're not using the functionality the way you expected to.
- More work up-front but results in higher productivity in the long run.

Developing good infrastructure is always time well spent.



The code writes out the array in binary and text formats. Let's start with putting those parts in functions.

```
//hydrogen.cpp
#include <string>
```

```
void writeBinary(const std::string& s, int n, const double x[]) {
   std::ofstream g(s, std::ios::binary);
   g.write((char*)(x), n*sizeof(x[0]));
   g.close();
void writeText(const std::string& s, int n, const double x[]) {
   std::ofstream f(s):
   for (int i=0: i<n: i++)
      f << a[i] << std::endl:</pre>
   f.close():
int main() {
   writeText("data.txt", n, a);
   writeBinarv("data.bin", n, a):
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```

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The code writes out the array in binary and text formats. Let's extract function declarations.

```
//hydrogen.cpp
#include <string>
```

```
void writeBinary(const std::string& s, int n, const double x[]) {
   std::ofstream g(s, std::ios::binary);
   g.write((char*)(x), n*sizeof(x[0]));
   g.close();
void writeText(const std::string& s, int n, const double x[]) {
   std::ofstream f(s):
   for (int i=0: i<n: i++)</pre>
      f << a[i] << std::endl:</pre>
   f.close():
int main() {
   writeText("data.txt", n, a);
   writeBinarv("data.bin", n, a):
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```

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The code writes out the array in binary and text formats. Let's extract declarations.

```
#include <string>
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void writeText(const std::string& s, int n, const double x[]);
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   g.write((char*)(x), n*sizeof(x[0]));
   g.close();
void writeText(const std::string& s, int n, const double x[]) {
   std::ofstream f(s):
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```
#include <string>
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   writeText("data.txt", n, a);
   writeBinarv("data.bin", n, a):
```

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The code writes out the array in binary and text formats. Function definitions can be moved.

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```
#include <string>
void writeBinary(const std::string& s, int n, const double x[]);
void writeText(const std::string& s, int n, const double x[]);
void writeBinary(const std::string& s, int n, const double x[]) {
void writeText(const std::string& s, int n, const double x[]) {
int main() {
   writeText("data.txt", n, a);
   writeBinarv("data.bin", n, a):
```

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The code writes out the array in binary and text formats. Function definitions can be moved.

```
#include <string>
void writeBinary(const std::string& s, int n, const double x[]);
void writeText(const std::string& s, int n, const double x[]);
int main() {
   writeText("data.txt", n, a);
  writeBinary("data.bin", n, a);
void writeBinarv(const std::string& s. int n. const double x[]) {
    // bunch of commands
}
void writeText(const std::string& s, int n, const double x[]) {
   // bunch of commands
```

The code writes out the array in binary and text formats. We're ready to make a module now!

```
#include <string>
void writeBinary(const std::string& s, int n, const double x[]);
void writeText(const std::string& s, int n, const double x[]);
int main() {
   writeText("data.txt", n, a);
  writeBinary("data.bin", n, a);
void writeBinarv(const std::string& s. int n. const double x[]) {
    // bunch of commands
}
void writeText(const std::string& s, int n, const double x[]) {
   // bunch of commands
```

## Creating the module

To create our own module, put the declarations for the functions in their own 'header' file.

//outputarray.h
#include <string>
void writeBinary(const std::string& s, int n, const double x[]);
void writeText(const std::string& s, int n, const double x[]);

The source code with the definitions of the functions shoule be put into its own separate file.

```
//outputarray.cpp
#include "outputarray.h"
#include <fstream>
void writeBinary(const std::string& s, int n, const double x[]) {
    // bunch of commands
}
void writeText(const std::string& s, int n, const double x[]) {
    // bunch of commands
}
```



# Using the module

The original code that uses these would look like:

```
//hydrogen.cpp
#include "outputarray.h"
//...
int main() {
    //...
    writeText("data.txt", data, n);
    writeBinary("data.bin", data, n);
    //...
}
```



# **Compiling + Linking = Building**

So how to compile this code?

- Before the full program can be compiled, all the **source** files (hydrogen.cpp, outputarray.cpp) must be **compiled**.
- outputarray.cpp doesn't contain a main function, so it can't be an executable (no program to run). Instead outputarray.cpp is compiled into a ".o" file, an **object file**, using the "-c" flag
- It is customary and advisable to compile all the code pieces into object files.
- After all the object files are generated, they are linked together to create the working executable.

\$ g++ -std=c++17 outputarray.cpp -c -o outputarray.o // compile \$ g++ -std=c++17 hydrogen.cpp -c -o hydrogen.o // compile \$ g++ -std=c++17 outputarray.o hydrogen.o -o hydrogen // link

If you leave out one of the needed .o files you will get a fatal linking error: "undefined reference to ...."



### Interface v. Implementation

By creating a header file, we separated the interface from the implementation.

- The implementation the actual code for writeBinary and writeText goes in the .cpp (or .cc or .cxx) or 'source' file. This is compiled on its own, separately from any program that uses its functions.
- The interface what the calling code needs to know goes in the .h or 'header' files. This is also called the API (Application Programming Interface).

//outputarray.h
#include <string>
void writeBinary(const std::string& s, int n, const double x[]);
void writeText(const std::string& s, int n, const double x[]);

This distinction is crucial for writing modular code.



### Interface v. implementation

So, to review:

• When hydrogen.cpp is being compiled, the header file outputarray.h is included to tell the compiler that there exists out there somewhere functions of the form

void writeBinary(const std::string& s, int n, const double x[]); void writeText(const std::string& s, int n, const double x[]);

- This allows the compiler to check the number and type of arguments and the return type for those functions (the interface).
- The compiler does not need to know the details of the implementation, since it's not compiling the implementation (the source code of the routine).
- The programmer of hydrogen.cpp also does not need to know the implementation, and is free to assume that writeBinary and writeText have been programmed correctly.



# Guards against multiple inclusion

Protect your header files!

- Header files can include other header files.
- It can be hard to figure out which header files are already included in the program.
- Including a header file twice will lead to doubly-defined entities, which results in a compiler error.
- The solution is to add a 'preprocessor guard' to every header file:

//outputarray.h
#ifndef OUTPUTARRAY\_H
#define OUTPUTARRAY\_H
#include <string>
void writeBinary(const std::string& s, int n, const double x[]);
void writeText(const std::string& s, int n, const double x[]);
#endif

We'll expect to see these in your homework.



## About the preprocessor

What do you mean by "preprocessor"?

- Before the compiler actually compiles the code, a "preprocessor" is run on the code.
- For our purposes, the preprocessor is essentially just a text-substitution tool.
- Every line that starts with "#" is interpreted by the preprocessor.
- The most common directives a beginner encounters are #include, #ifndef, #define, and #endif.

#### **Future Feature**

The new C++20 standard defines a new way to build modules than to rely on the preprocessor.

Implementations by compilers are not fully there yet, or are experimental. They also still vary a lot from compiler to compiler, to the extent that it changes how you build software.

So for now, let's stick with the #include technique.



## What goes into the interface (i.e. the header file)?

So what should one expect in a header file?

- At the very least, the **function declarations**.
- There may also be **constants** that the calling function and the routine need to agree on (error codes, for example) or **definitions of data structures**, classes, etc.
- **Comments**, which give a description of the module and its functions.

Further guidelines:

- There should really only be one header file per module. In theory there can be multiple source files.
- Not necessarily every function declaration is in the header file, just the public ones. Routines internal to the module are not in the public header file.



## What goes into the implementation (source file)?

What should one expect in a source file?

- Everything which is defined in the .h file which requires code that is not in the .h file. Particularly, **function definitions**.
- Internal routines which are used by the routines declared in the .h file.
- To ensure consistency, include the corresponding .h file at the top of the file.
- Everything that needs to be compiled and linked to code that uses the .h file.



## Modularization allows faster compilation

Consider a build tree with several source files, e.g.



 Each source file compilation into an .o file can be done simultaneously.

#### Parallel processing of code!

If only one .cc file has changed, only that file needs to be recompiled.

#### Fast rebuilds

But:

- Now you need to keep track of what depends upon what.
- You need to retype in the entire compilation command every time you need to recompile.
- It's easy to forget all of your compiler flags from one day to the next, as well as the location of external libraries.

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### Section 2

Make



### Make

- make is a build program that is used to build programs from multiple .cpp, .h, .o, and other files.
- It is actually a very general framework that is used to compile code, of any type.
- make takes a Makefile' as its input, which specifies what to do, and how.
- The Makefile contains variables, rules and dependencies.
- The Makefile specifies executables, compiler flags, library locations, ...
- Build programs are a crucial component of professional software development.

 $https://www.gnu.org/software/make/manual/html\_node/index.html$ 



### **Basic usage**

• Make is invoked with a list of target file names to build as *command-line arguments*:

```
$ make [TARGET ...]
```

- Without arguments, make builds the first target that appears in its makefile, which is traditionally a symbolic phony target named all.
- Make decides whether a target needs to be regenerated by comparing file modification times.
- This solves the problem of avoiding the building of files which are already up to date, as long as the timestamps are consistent and correct.



### **Makefiles**

- Make searches the current directory for the makefile to use, (e.g. makefile, Makefile, GNUmakefile and then runs the specified (or default) target(s) from (only) that file.
- It is possible to specify a different *makefile* by using the -f flag, e.g. sh \$ make -f myMakefile [TARGET ...]
- The makefile is a plain-text file, with a particular structure
- It may include rules and use commands from the shell.



### Rules

- A makefile consists of rules.
- Each rule begins with a textual dependency line which defines a target followed by a colon (:) and optionally an enumeration of components (files or other targets) on which the target depends.
- The dependency line is arranged so that the target (left of the colon) depends on "components" or "prerequisites" (to its right)
- Each command-line must start with a TAB character to be recognized as a command.

Unfortunately, as you can't easily see in your editor whether you have a TAB character or a set of spaces. If you have spaces instead of a TAB, make will print the unhelpful error: Makefile:3: \*\*\* missing separator. Stop.

```
TARGET: dependencies...
    [command 1]}
    :
    [command n]
TARGET1 [TARGET2 ...]: dep1 dep2 ...
    [command 1]
    :
    [command n]
```



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### **Rules - commands**

- Each command is executed by a separate shell or command-line interpreter instance.
- $\bullet\,$  The backslash  $\setminus\,$  can be used to have commands executed by the same shell, it represents line-continuation
- Commands can be separated by ;
- Comments are included using #
- A rule may have no command lines defined.

The dependency line can consist solely of components that refer to targets.

This means either there is nothing to do, or there is a predefined rule.



## Macros & Variables

- Macros are usually referred to as variables when they hold simple string definitions, like CXX = g++.
- Macros in makefiles may be overridden by the command-line arguments passed to the Make utility (e.g. "make CXX=icpc").
- Macros allow users to specify the programs invoked and other custom behavior during the build process.

For example, the macro CXX is used in makefiles to refer to the location of the C++ compiler  $% \left( {{{\rm{C}}_{{\rm{C}}}}_{{\rm{C}}}} \right)$ 

 To use variables, you need to use a dollar sign (\$) followed by the name of the variable in parenthesis or curly braces.

```
MACRO = definition
```

```
PACKAGE = package
VERSION = `date +"\%Y.\%m\%d"`
ARCHIVE = $(PACKAGE)-$(VERSION)
```

#### dist:

```
# Notice that only now macros are
# expanded for shell to interpret:
# tar -cf ../package-`date +"\%Y\%m\%d"`.tar
tar -cf ../$(ARCHIVE).tar .
```

Note: Environment variables are also available as macros.



# **Compilation and Linking**

# Example Makefile for the `hydrogen` program (after modularization) CXX=g++ CXXFLAGS=-std=c++17 -02 all: hydrogen

```
hydrogen: hydrogen.o outputarray.o initmatrix.o eigenvalue.o
        $(CXX) -o hydrogen hydrogen.o outputarray.o initmatrix.o eigenvalue.o
hydrogen.o: hydrogen.c outputarray.h initmatrix.h eigenvalue.h
    $(CXX) -c $(CXXFLAGS) -o hydrogen.o hydrogen.c
outputarray.o: outputarray.cpp outputarray.h
    $(CXX) -c $(CXXFLAGS) -o outputarray.o outputarray.c
initmatrix.o: initmatrix.cpp initmatix.h
    $(CXX) -c $(CXXFLAGS) -o initmatrix.o initmatrix.c
eigenvalue.o: eigenvalue.cpp eigenvalue.h
    $(CXX) -c $(CXXFLAGS) -o eigenvalue.o eigenvalue.c
clean:
   $(RM) eigenvalue.o initmatrix.o outputarray.o hydrogen.o
.PHONY: all clean
```

# **Compilation and Linking**

What happens when you type make?

- make will only recompile those dependencies that have source files that are newer then the library, thus only the code you are working on is modified.
- If a target is not a file, you should declare it 'PHONY'. Otherwise, should a file by that name exist, make thinks it's done already.
- It's good practice to put a clean rule in your Makefile that allows the whole compilation to restart.
- Several rules could be processed at the same time; you can tell make to try and use multiple processes when the dependencies allow it, but specifying a -j option, e.g.

\$ make -j 4



# **Special Variables**

- \$@: the target filename
- \$\*: the target filename without the file extension
- \$<: the first prerequisite filename
- \$^: the filenames of all the prerequisites, separated by spaces, discard duplicates.
- \$+: similar to \$^, but includes duplicates
- \$?: the names of all prerequisites that are newer than the target, separated by spaces



# **Compilation and Linking**

# Example Makefile for the `hydrogen` program (after modularization)
CXX=g++
CXXFLAGS=-std=c++17 -02
all: hydrogen

```
hydrogen: hydrogen.o outputarray.o initmatrix.o eigenvalue.o
    $(CXX) -o $@ $^
hydrogen.o: hydrogen.c outputarray.h initmatrix.h eigenvalue.h
    (CXX) - c (CXXFLAGS) - o 
outputarray.o: outputarray.cpp outputarray.h
    $(CXX) -c $(CXXFLAGS) -o $@ $<
initmatrix.o: initmatrix.cpp initmatix.h
    $(CXX) -c $(CXXFLAGS) -o $@ $<
eigenvalue.o: eigenvalue.cpp eigenvalue.h
    (CXX) - c (CXXFLAGS) - o 
clean:
    $(RM) eigenvalue.o initmatrix.o outputarrav.o hvdrogen.o
```

.PHONY: all clean

# Assignment 1

You will receive an existing, monolythic code that simulates a one-dimensional wave equation, and your task will be to convert it into a modular form, and supply it with a proper Makefile to compile the code.

Please see the course site for details later this Thursday.

Note that at the office hours on Friday (January 21, 2022), we will briefly go over this code as well.

The assignment will be due January 26th, 2022 at midnight. Upload all your code, the output, and a brief report (text format) on what you did.

