

# Fortran Programming Guide

FORTRAN 77 5.0 — Fortran 90 2.0

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## **Preface**

This guide combines the essential information programmers need to develop efficient applications using the two Sun Fortran compilers, £77 (FORTRAN 77 version 5.0) and £90 (Fortran 90 version 2.0). It presents issues relating to input/output, program development, use and creation of software libraries, program analysis and debugging, numerical accuracy, porting, performance, optimization, parallelization, and the C/Fortran interface.

Discussion of the compiler command-line options and their use can be found in the companion book, the *Fortran User's Guide*.

**Note -** This guide covers the Sun FORTRAN 77 and Fortran 90 compilers. The text uses "f77/f90" and "Fortran" to indicate information that is common to *both* compilers.

# Who Should Use This Book

This guide is intended for scientists, engineers, and programmers who have a working knowledge of the Fortran language and wish to learn how to use the Sun Fortran compilers effectively. Familiarity with the Solaris operating environment or UNIX $^{\otimes}$  $^{\odot}$  in general is also assumed.

# How This Book Is Organized

This guide is organized into the following chapters:

- Chapter 1, "Introduction," briefly describes the features of the compilers.
- Chapter 2, "Fortran Input/Output," discusses how to use I/O efficiently.
- Chapter 3, "Program Development," demonstrates how program management tools like SCCS, make, and Teamware can be helpful.
- Chapter 4, "Libraries," explains use and creation of software libraries.
- Chapter 5, "Program Analysis and Debugging," describes use of dbx and other analysis tools.
- Chapter 6, "Floating Point Arithmetic," introduces important issues regarding numerical computation accuracy.
- Chapter 7, "Porting," considers porting programs to Sun compilers.
- Chapter 8, "Performance Profiling," describes techniques for performance measurement.
- Chapter 9, "Performance and Optimization," indicates ways to improve execution performance of Fortran programs.
- *Chapter 10, "Parallelization,"* explains the multiprocessing features of the compilers.
- Chapter 11, "C-Fortran Interface," describes how C and Fortran routines can call each other and pass data.

# Multiplatform Release

**Note -** The name of the latest Solaris operating environment release is Solaris 7 but some documentation and path or package path names may still use Solaris 2.7 or SunOS 5.7.

The Sun Fortran documentation covers the release of the Fortran compilers on a number of operating environments and hardware platforms:

FORTRAN 77 5.0 is released for:

- Solaris 2.5.1, 2.6, and Solaris 7 environments on:
  - lacksquare architectures based on the SPARC  $^{^{TM}}$  microprocessor
  - x86-based architectures, where x86 refers to the Intel $^{®TM}_{TM}$  implementation of one of the following: Intel 80386 , Intel 80486 , Pentium , or the equivalent

Fortran 90 2.0 is released for:

■ Solaris 2.5.1, 2.6, and Solaris 7 environments on SPARC processors only.

**Note -** The term "x86" refers to the Intel 8086 family of microprocessor chips, including the Pentium, Pentium Pro, and Pentium II processors and compatible microprocessor chips made by AMD and Cyrix. In this document, the term "x86" refers to the overall platform architecture. Features described in this book that are particular to a specific platform are differentiated by the terms "SPARC" and "x86" in the text.

## **Related Books**

The following books augment this manual and provide essential information:

- Fortran User's Guide—provides information on command line options and how to use the compilers.
- FORTRAN 77 Language Reference. Complete language reference.
- Fortran Library Reference-gives details on the language and routines.
- Sun Performance WorkShop Fortran Overview gives a high-level outline of the Fortran package suite.

## Other Programming Books

- *C User's Guide*—describes compiler options, pragmas, and more.
- Numerical Computation Guide—details floating-point computation and numerical accuracy issues.
- Sun WorkShop Performance Library Reference-discusses the library of subroutines and functions to perform useful operations in computational linear algebra and Fourier transforms.

### Other Sun WorkShop Books

- Sun WorkShop Quick Install-provides installation instructions.
- Sun WorkShop Installation Reference-provides supporting installation and licensing information.
- *Sun Visual WorkShop C++ Overview*-gives a high-level outline of the C++ package suite.
- Using Sun WorkShop—gives information on performing development operations through Sun WorkShop.

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- *Debugging a Program With dbx*—provides information on using dbx commands to debug a program.
- Analyzing Program Performance with Sun WorkShop—describes the profiling tools; LoopTool, LoopReport, LockLint utilities; and the Sampling Analyzer to enhance program performance.
- Sun WorkShop TeamWare User's Guide—describes how to use the Sun WorkShop TeamWare code management tools.

### Solaris Books

The following Solaris manuals and guides provide additional useful information:

- The Solaris Linker and Libraries Guide—gives information on linking and libraries.
- The Solaris *Programming Utilities Guide*—provides information for developers about the special built-in programming tools available in the SunOS system.

# **Ordering Sun Documents**

The SunDocs<sup>SM</sup> program provides more than 250 manuals from Sun Microsystems, Inc. If you live in the United States, Canada, Europe, or Japan, you can purchase documentation sets or individual manuals using this program.

For a list of documents and how to order them, see the catalog section of the  $SunExpress^{TM}$  Internet site at http://www.sun.com/sunexpress.

# **Accessing Sun Documents Online**

Sun WorkShop documentation is available online from several sources:

- The docs.sun.com Web site
- AnswerBook2<sup>TM</sup> collections
- HTML documents
- Online help and release notes

### Using the docs.sun.com Web site

The docs.sun.com Web site enables you to access Sun technical documentation online. You can browse the docs.sun.com archive or search for a specific book title or subject. The URL is http://docs.sun.com.

## Accessing AnswerBook2 Collections

The Sun WorkShop documentation is also available using AnswerBook2 software. To access the AnswerBook2 collections, your system administrator must have installed the AnswerBook2 documents during the installation process (if the documents are not installed, see your system administrator or Chapter 3 of Sun WorkShop Quick Install for installation instructions). For information about accessing AnswerBook2 documents, see Chapter 6 of Sun WorkShop Quick Install, Solaris installation documentation, or your system administrator.

**Note** - To access AnswerBook2 documents, Solaris 2.5.1 users must first download AnswerBook2 documentation server software from a Sun Web page. For more information, see Chapter 6 of *Sun WorkShop Quick Install*.

### **Accessing HTML Documents**

The following Sun Workshop documents are available online only in HTML format:

- Tools.h++ Class Library Reference
- Tools.h++ User's Guide
- Numerical Computation Guide
- Standard C++ Library User's Guide
- Standard C++ Class Library Reference
- Sun WorkShop Performance Library Reference Manual
- Sun WorkShop Visual User's Guide
- Sun WorkShop Memory Monitor User's Manual

To access these HTML documents:

### 1. Open the following file through your HTML browser:

install-directory/SUNWspro/DOC5.0/lib/locale/C/html/index.html

Replace *install-directory* with the name of the directory where your Sun WorkShop software is installed (the default is /opt).

The browser displays an index of the HTML documents for the Sun WorkShop products that are installed.

2. Open a document in the index by clicking the document's title.

# Accessing Sun WorkShop Online Help and Release Notes

This release of Sun WorkShop includes an online help system as well as online manuals. To find out more see:

- Online Help. A help system containing extensive task-oriented, context-sensitive help. To access the help, choose Help Help Contents. Help menus are available in all Sun WorkShop windows.
- Release Notes. The Release Notes contain general information about Sun WorkShop and specific information about software limitations and bugs. To access the Release Notes, choose Help Release Notes.
- You can view the latest release information regarding the Fortran compilers by invoking the f77 or f90 compiler with the -xhelp=readme flag.

# What Typographic Changes Mean

The following table describes the typographic changes used in this book.

TABLE P-1 Typographic Conventions

Typeface or Symbol	Meaning	Example
AaBbCc123	The names of commands, files, and directories; on-screen computer	Edit your .login file. Use ls -a to list all files.
	output	machine_name% You have mail.
AaBbCc123	What you type, contrasted with on-screen computer output	machine_name% <b>su</b>
	- •	Password:

TABLE P-1 Typographic Conventions (continued)

Typeface or Symbol	Meaning	Example
AaBbCc123	Command-line placeholder: replace with a real name or value	To delete a file, type rm filename.
AaBbCc123	Book titles, new words or terms, or words to be emphasized	Read Chapter 6 in <i>User's Guide</i> . These are called <i>class</i> options.  You <i>must</i> be root to do this.

# Shell Prompts in Command Examples

The following table shows the default system prompt and superuser prompt for the C shell, Bourne shell, and Korn shell.

TABLE P-2 Shell Prompts

Shell	Prompt
C shell prompt	machine_name%
C shell superuser prompt	machine_name#
Bourne shell and Korn shell prompt	\$
Bourne shell and Korn shell superuser prompt	#

# Other Conventions Used in This Book

The following conventions appear in the text of this book:

■ Code listings and examples appear in boxes:

WRITE( \*, \* ) "Hello world"

- The symbol ¤ stands for a blank space where a blank is significant: ¤¤36.001
- FORTRAN 77 examples appear in tab format, while Fortran 90 examples appear in free format. Examples common to both FORTRAN 77 and 90 use tab format except where indicated.
- Uppercase characters are generally used to show Fortran keywords and intrinsics (PRINT), and lowercase or mixed case is used for variables (TbarX).
- The Sun Fortran compilers are referred to by their command names, either £77 or £90. "£77/£90" indicates information that is common to both the FORTRAN 77 and Fortran 90 compilers.
- References to online man pages appear with the topic name and section number. For example, a reference to GETENV will appear as *getenv*(3F), implying that the man command to access this page would be: man -s 3F getenv
- System Administrators may install the Sun Fortran compilers and supporting material at: <install\_point>/SUNWspro/SC5.0/ where <install\_point> is usually /opt for a standard install. This is the location assumed in this book.

# Introduction

The Sun Fortran compilers, £77 and £90, described in this book (and the companion *Fortran User's Guide*) are available under the Solaris operating environment on the various hardware platforms that Solaris supports. The compilers themselves conform to published Fortran language standards, and provide many extended features, including multiprocessor parallelization, sophisticated optimized code compilation, and mixed C/Fortran language support.

## **Standards Conformance**

- £77 was designed to be compatible with the ANSI X3.9-1978 Fortran standard and the corresponding International Organization for Standardization (ISO) 1539-1980, as well as standards FIPS 69-1, BS 6832, and MIL-STD-1753.
- f90 was designed to be compatible with the ANSI X3.198-1992 standard, and ISO/IEC 1539:1991
- Floating-point arithmetic for both compilers is based on IEEE standard 754-1985, and international standard IEC 60559:1989.
- On SPARC processors, both compilers provide support for the optimization-exploiting features of SPARC V8, and SPARC V9, including the UltraSPARC implementation. These features are defined in the SPARC Architecture Manuals, Version 8 (ISBN 0-13-825001-4), and Version 9 (ISBN 0-13-099227-5), published by Prentice-Hall for SPARC International.
- In this document, "Standard" means conforming to the versions of the standards listed above. "Non-standard" or "Extension" refers to features that go beyond these versions of these standards.

The responsible standards bodies may revise these standards from time to time. The versions of the applicable standards to which these compilers conform may be revised or replaced, resulting in features in future releases of the Sun Fortran compilers that create incompatibilities with earlier releases.

# Features of the Fortran Compilers

Sun Fortran compilers provide the following features or extensions:

- £77: Global program checking across routines for consistency of arguments, commons, parameters, and the like
- *SPARC:* Multiprocessor support, including automatic and explicit loop parallelization tightly integrated with optimization

**Note -** Parallelization features of the Fortran compilers are available only with Sun Performance WorkShop .

- £77: Many VAX/VMS Fortran extensions, including:
  - NAMELIST
  - DO WHILE
  - Structures, records, unions, maps
  - Variable format expressions
  - Recursion
  - Pointers
  - Double-precision complex
  - *SPARC*: Quadruple-precision real
  - *SPARC*: Quadruple-precision complex
- Cray-style parallelization directives, including TASKCOMMON, with extensions on f90.
- Global, peephole, and potential parallelization optimizations produce high-performance applications. Benchmarks show that optimized applications can run significantly faster when compared to unoptimized code.
- Common calling conventions on Solaris systems permit routines written in C or C++ to be combined with Fortran programs.
- Support for 64-bit Solaris 7 environments on UltraSPARC processors.
- Fortran 95 features in £90 include the attributes PURE and ELEMENTAL, and enhanced forms of MAXVAL and MINVAL.
- Call-by-value, %VAL, implemented in both £77 and £90.

■ Interoperability between FORTRAN 77 and Fortran 90 programs and object binaries.

# Fortran Input/Output

This chapter discusses the input/output features provided by Sun Fortran compilers.

# Accessing Files From Within Fortran Programs

Data is transferred between the program and devices or files through a Fortran *logical unit*. Logical units are identified in an I/O statement by a logical unit number, a nonnegative integer from 0 to the maximum 4-byte integer value (2,147,483,647).

The character \* can appear as a logical unit identifier. The asterisk stands for standard input file when it appears in a READ statement; it stands for *standard output file* when it appears in a WRITE or PRINT statement.

A Fortran logical unit can be associated with a specific, named file through the OPEN statement. Also, certain "preconnected" units are automatically associated with specific files at the start of program execution.

### **Accessing Named Files**

The OPEN statement's FILE= specifier establishes the association of a logical unit to a named, physical file at runtime. This file can be pre-existing or created by the program. See the Sun *FORTRAN 77 Language Reference Manual* for a full discussion of the OPEN statement.

The FILE= specifier on an OPEN statement may specify a simple file name (FILE='myfile.out') or a file name preceded by an absolute or relative directory

path (FILE='../Amber/Qproj/myfile.out'). Also, the specifier may be a character constant, variable, or character expression.

Library routines can be used to bring command-line arguments and environment variables into the program as character variables for use as file names in OPEN statements. (See man page entries for getarg(3F) and getenv(3F) for details; these and other useful library routines are also described in the *Fortran Library Reference*).

The following example (GetFilNam.f) shows one way to construct an absolute path file name from a typed-in name. The program uses the library routines GETENV, LNBLNK, and GETCWD to return the value of the \$HOME environment variable, find the last non-blank in the string, and determine the current working directory:

```
CHARACTER F*128, FN*128, FULLNAME*128
      PRINT*, "ENTER FILE NAME:"
      READ *, F
      FN = FULLNAME( F )
      PRINT *, "PATH IS: ",FN
      CHARACTER*128 FUNCTION FULLNAME( NAME )
      CHARACTER NAME*(*), PREFIX*128
С
          This assumes C shell.
           Leave absolute path names unchanged.
С
            If name starts with "~/", replace tilde with home
С
            directory; otherwise prefix relative path name with
С
           path to current directory.
      IF ( NAME(1:1) .EQ. "/" ) THEN
           FULLNAME = NAME
      ELSE IF ( NAME(1:2) .EQ. "\sim/" ) THEN
           CALL GETENV( "HOME", PREFIX )
            FULLNAME = PREFIX(:LNBLNK(PREFIX)) //
&
                          NAME (2:LNBLNK(NAME))
      ELSE
           CALL GETCWD( PREFIX )
            FULLNAME = PREFIX(:LNBLNK(PREFIX)) //
&
                         "/" // NAME(:LNBLNK(NAME))
      ENDIF
      RETURN
      END
```

Compiling and running GetFilNam.f results in:

```
demo% pwd
/home/users/auser/subdir
demo% f77 -silent -o getfil GetFilNam.f
demo% getfil
anyfile
/home/users/auser/subdir/anyfile
demo%
```

### Opening Files Without a Name

The OPEN statement need not specify a name; the runtime system supplies a file name according to several conventions.

### Opened as Scratch

Specifying STATUS='SCRATCH' in the OPEN statement opens a file with a name of the form tmp. FAAAxnnnnn, where nnnnn is replaced by the current process ID, AAA is a string of three characters, and x is a letter; the AAA and x make the file name unique. This file is deleted upon termination of the program or execution of a CLOSE statement, unless (with f77) STATUS='KEEP' is specified in the CLOSE statement.

### Already Open

If the file has already been opened by the program, you can use a subsequent OPEN statement to change some of the file's characteristics; for example, BLANK and FORM. In this case, you would specify only the file's logical unit number and the parameters to change.

### **Preconnected Units**

Three unit numbers are automatically associated with specific standard I/O files at the start of program execution. These preconnected units are standard input, standard output, and standard error:

- Standard input is logical unit 5 (also Fortran 90 unit 100)
- Standard output is logical unit 6 (also Fortran 90 unit 101)
- Standard error is logical unit 0 (also Fortran 90 unit 102)

Typically, standard input receives input from the workstation keyboard; standard output and standard error display output on the workstation screen.

In all other cases where a logical unit number but no FILE= name is specified on an OPEN statement, a file is opened with a name of the form fort. n, where n is the logical unit number.

### Opening Files Without an OPEN Statement

Use of the OPEN statement is optional in those cases where default conventions can be assumed. If the first operation on a logical unit is an I/O statement other than OPEN or INQUIRE, the file fort. n is referenced, where n is the logical unit number (except for 0, 5, and 6, which have special meaning).

These files need not exist before program execution. If the first operation on the file is not an OPEN or INQUIRE statement, they are created.

Example: The WRITE in the following code creates the file fort.25 if it is the first input/output operation on that unit:

```
demo% cat TestUnit.f
    IU=25
    WRITE( IU, "(I4)" ) IU
    END
demo%
```

The preceding program opens the file fort. 25 and writes a single formatted record onto that file:

```
demo% f77 -silent -o testunit TestUnit.f
demo% testunit
demo% cat fort.25
   25
demo%
```

## Passing File Names to Programs

The file system does not have any automatic facility to associate a logical unit number in a Fortran program with a physical file.

However, there are several satisfactory ways to communicate file names to a Fortran program.

### Via Runtime Arguments and GETARG

The library routine *getarg*(3F) can be used to read the command-line arguments at runtime into a character variable. The argument is interpreted as a file name and used in the OPEN statement FILE= specifier:

### Via Environment Variables and GETENV

Similarly, the library routine *getenv*(3F) can be used to read the value of any environment variable at runtime into a character variable that in turn is interpreted as a file name:

```
demo% cat testenv.f
         CHARACTER outfile*40
C Get $OUTFILE as output file name for unit 51
         CALL getenv("OUTFILE", outfile)
         OPEN(51,FILE=outfile)
         WRITE(51,*) "Writing to file: ", outfile
demo% f77 -silent -o tstenv testenv.f
demo% setenv OUTFILE EnvFileName
demo% tstenv
demo% cat EnvFileName
Writing to file: EnvFileName
demo%
```

When using getarg or geteny, care should be taken regarding leading or trailing blanks. (FORTRAN 77 programs can use the library function LNBLNK; Fortran 90 programs can use the intrinsic function TRIM.) Additional flexibility to accept relative path names can be programmed along the lines of the FULLNAME function in the example at the beginning of this chapter.

### f77: Logical Unit Preattachment Using IOINIT

The library routine IOINIT can also be used with £77 to attach logical units to specific files at runtime. IOINIT looks in the environment for names of a user-specified form and then opens the corresponding logical unit for sequential formatted I/O. Names must be of the general form PREFIXnn, where the particular *PREFIX* is specified in the call to IOINIT, and *nn* is the logical unit to be opened. Unit numbers less than 10 must include the leading 0. See the Sun Fortran Library Reference, and the IOINIT(3F) man page. (The IOINIT facility is not implemented for f90.)

Example: Associate physical files test.inp and test.out in the current directory to logical units 1 and 2:

First, set the environment variables.

```
demo$ TST01=ini1.inp
demo$ TST02=ini1.out
demo$ export TST01 TST02
```

With ksh or sh:

#### With csh:

The program inil.f reads 1 and writes 2:

With environment variables and ioinit, inil.f reads inil.inp and writes to inil.out:

```
demo% cat inil.inp
  12 3.14159012 6
demo% f77 -silent -o tstinit inil.f
demo% tstinit
demo% cat inil.out
  12 3.14159 6
demo%
```

IOINIT is adequate for most programs as written. However, it is written in Fortran specifically to serve as an example for similar user-supplied routines. Retrieve a copy from the following file, a part of the FORTRAN 77 package installation:

/opt/SUNWspro/SC5.0/src/ioinit.f

### Command-Line I/O Redirection and Piping

Another way to associate a physical file with a program's logical unit number is by redirecting or piping the preconnected standard I/O files. Redirection or piping occurs on the runtime execution command.

In this way, a program that reads standard input (unit 5) and writes to standard output (unit 6) or standard error (unit 0) can, by redirection (using <, >, >>, >&, |, |&, |&, |&> 2>&1 on the command line), read or write to any other named file.

This is shown in the following table:

TABLE 2-1 csh/sh/ksh Redirection and Piping on the Command Line

Action	Using C Shell	Using Bourne or Korn Shell
Standard input —read from mydata	myprog < mydata	myprog < mydata
Standard output —write (overwrite) myoutput	myprog > myoutput	myprog > myoutput
Standard output — write/ append to myoutput	myprog >> myoutput	myprog >> myoutput
Redirect standard error to a file	myprog >& errorfile	myprog 2> errorfile
Pipe standard output to input of another program	myprog1   myprog2	myprog1   myprog2
Pipe standard error and output to another program	myprog1  & myprog2	myprog1 2>&1   myprog2

See the csh, ksh, and sh man pages for details on redirection and piping on the command line.

# f77: VAX / VMS Logical File Names

If you are porting from VMS FORTRAN to FORTRAN 77, the VMS-style logical file names in the INCLUDE statement are mapped to UNIX path names. The environment variable LOGICALNAMEMAPPING defines the mapping between the logical names and the UNIX path name. If the environment variable LOGICALNAMEMAPPING is set and the -vax, -xl or -xld compiler options are used, the compiler interprets VMS logical file names on the INCLUDE statement.

The compiler sets the environment variable to a string with the following syntax:

"lname1=path1; lname2=path2; "

Each *lname* is a logical name, and each *path* is the path name of a directory (without a trailing /). All blanks are ignored when parsing this string. Any trailing /list or /nolist is stripped from the file name in the INCLUDE statement. Logical names in a file name are delimited by the first colon in the VMS file name. The compiler converts file names of the form:

Iname1: file

to:

path1/file

Uppercase and lowercase are significant in logical names. If a logical name is encountered on the INCLUDE statement that was not specified by LOGICALNAMEMAPPING, the file name is used unchanged.

## Direct I/O

Direct or random I/O allows you to access a file directly by record number. Record numbers are assigned when a record is written. Unlike sequential I/O, direct I/O records can be read and written in any order. However, in a direct access file, all records must be the same fixed length. Direct access files are declared with the ACCESS='DIRECT' specifier on the OPEN statement for the file.

A logical record in a direct access file is a string of bytes of a length specified by the OPEN statement's RECL= specifier. READ and WRITE statements must not specify logical records larger than the defined record size. (Record sizes are specified in bytes.) Shorter records are allowed. Unformatted, direct writes leave the unfilled part of the record undefined. Formatted, direct writes cause the unfilled record to be padded with blanks.

Direct access READ and WRITE statements have an extra argument, REC=n, to specify the record number to be read or written.

Example: Direct access, unformatted:

This program opens a file for direct access, unformatted I/O, with a fixed record length of 200 bytes, then reads the thirteenth record into X and Y.

Example: Direct access, *formatted*:

```
OPEN( 2, FILE="inven.db", ACCESS="DIRECT", RECL=200,
        FORM="FORMATTED", ERR=90 )
READ( 2, FMT="(I10,F10.3)", REC=13, ERR=30 ) X, Y
```

This program opens a file for direct access, formatted I/O, with a fixed record length of 200 bytes. It then reads the thirteenth record and converts it with the format (I10,F10.3).

For formatted files, the size of the record written is determined by the FORMAT statement. In the preceding example, the FORMAT statement defines a record of 20 characters or bytes. More than one record can be written by a single formatted write if the amount of data on the list is larger than the record size specified in the FORMAT statement. In such a case, each subsequent record is given successive record numbers.

Example: Direct access, formatted, multiple record write:

```
OPEN( 21, ACCESS="DIRECT", RECL=200, FORM="FORMATTED")
WRITE(21, "(10F10.3)", REC=11) (X(J), J=1, 100)
```

The write to direct access unit 21 creates 10 records of 10 elements each (since the format specifies 10 elements per record) these records are numbered 11 through 20.

## **Internal Files**

An internal file is an object of type CHARACTER such as a variable, substring, array, element of an array, or field of a structured record. Internal file READ can be from a constant character string. I/O on internal files simulates formatted READ and WRITE statements by transferring and converting data from one character object to another data object. No file I/O is performed.

When using internal files:

- The name of the character object receiving the data appears in place of the unit number on a WRITE statement. On a READ statement, the name of the character object source appears in place of the unit number.
- A constant, variable, or substring object constitutes a single record in the file.
- With an array object, each array element corresponds to a record.
- f77: f77 extends direct I/O to internal files. (The ANSI standard includes only sequential formatted I/O on internal files.) This is similar to direct I/O on external

files, except that the number of records in the file cannot be changed. In this case, a record is a single element of an array of character strings.

■ Each sequential READ or WRITE statement starts at the beginning of an internal file.

Example: Sequential formatted read from an internal file (one record only):

```
demo% cat intern1.f
    CHARACTER X*80
    READ( *, "(A)" ) X
    READ( X, "(I3,I4)" ) N1, N2 ! This codeline reads the internal file X
    WRITE( *, * ) N1, N2
    END
demo% f77 -silent -o tstintern intern1.f
demo% tstintern
12 99
12 99
demo%
```

Example: Sequential formatted read from an internal file (three records):

Example: Direct access read from an internal file (one record) (f77 only):

```
demo% cat intern3.f
      CHARACTER LINE(4)*16
                             ! This is our "internal file"
                      12341234
      DATA LINE(1) / " 81 81 "
     DATA LINE(2) / " 82 82 " /
      DATA LINE(3) / " 83 83 " /
      DATA LINE(4) / " 84 84 " /
      READ ( LINE, FMT=20, REC=3 ) M, N
20
       FORMAT( 14, 14 )
      PRINT *, M, N
      END
demo% f77 -silent intern3.f
demo% a.out
  83 83
demo%
```

# Tape I/O (f77 Only)

Most typical Fortran I/O is done to disk files. However, by associating a logical unit number to a physically mounted tape drive via the OPEN statement, it is possible to do I/O directly to tape.

It could be more efficient to use the  $\mathtt{TOPEN}(\ )$  routines rather than Fortran I/O statements to do I/O on magnetic tape.

### Using TOPEN Routines

With the nonstandard tape I/O package (see *topen*(3F)) you can transfer blocks between the tape drive and buffers declared as Fortran character variables. You can then use internal I/O to fill and empty these buffers. This facility does not integrate with the rest of Fortran I/O and even has its own set of tape logical units. Refer to the man pages for complete information.

# Fortran Formatted I/O for Tape

The Fortran I/O statements provide facilities for transparent access to *formatted*, sequential files on magnetic tape. There is no limit on formatted record size, and records may span tape blocks.

## Fortran Unformatted I/O for Tape

Using the Fortran I/O statements to connect a magnetic tape for *unformatted* access is less satisfactory. The implementation of unformatted records implies that the size of a record (plus eight characters of overhead) cannot be bigger than the buffer size.

As long as this restriction is complied with, the I/O system does not write records that span physical tape blocks, writing short blocks when necessary. This representation of unformatted records is preserved (even though it is inappropriate for tapes) so that files can be freely copied between disk and tapes.

Since the block-spanning restriction does not apply to tape reads, files can be copied from tape to disk without any special considerations.

### Tape File Representation

A Fortran data file is represented on tape by a sequence of data records followed by an endfile record. The data is grouped into blocks, with maximum block size determined when the file is opened. The records are represented in the same way as records in disk files: formatted records are followed by newlines; unformatted records are preceded and followed by character counts. In general, there is no relation between Fortran records and tape blocks; that is, records can span blocks, which can contain parts of several records.

The only exception is that Fortran does not write an unformatted record that spans blocks; thus, the size of the largest unformatted record is eight characters less than the block size.

### The dd Conversion Utility

An end-of-file record in Fortran maps directly into a tape mark. In this respect, Fortran files are the same as tape system files. But since the representation of Fortran files on tape is the same as that used in the rest of UNIX, naive Fortran programs cannot read 80-column card images on tape. If you have an existing Fortran program and an existing data tape to read with it, translate the tape using the dd(1) utility, which adds newlines and strips trailing blanks.

Example: Convert a tape on mt0 and pipe that to the executable ftnprg:

demo% dd if=/dev/rmt0 ibs=20b cbs=80 conv=unblock | ftnprg

### The getc Library Routine

As an alternative to dd, you can call the *getc*(3F) library routine to read characters from the tape. You can then combine the characters into a character variable and use internal I/O to transfer formatted data. See also *TOPEN*(3F).

### **End-of-File**

The end-of-file condition is reached when an end-of-file record is encountered during execution of a READ statement. The standard states that the file is positioned after the end-of-file record. In real life, this means that the tape read head is poised at the beginning of the next file on the tape. Although it seems as if you could read the next file on the tape, this is not strictly true, and is not covered by the ANSI FORTRAN 77 Language Standard.

The standard also says that a BACKSPACE or REWIND statement can be used to reposition the file. Consequently, after reaching end-of-file, you can backspace over

the end-of-file record and further manipulate the file-for example, writing more records at the end, rewinding the file, and rereading or rewriting it.

## Multifile Tapes

The name used to open the tape file determines certain characteristics of the connection, such as the recording density and whether the tape is automatically rewound when opened and closed.

To access a file on a tape with multiple files, first use the mt(1) utility to position the tape to the needed file. Then open the file as a no-rewind magnetic tape such as /dev/nrmt0. Referencing the tape with this name prevents it from being repositioned when it is closed. By reading the file until end-of-file and then reopening it, a program can access the next file on the tape. Any program subsequently referencing the same tape can access it where it was last left, preferably at the beginning of a file, or past the end-of-file record.

However, if your program terminates prematurely, it may leave the tape positioned anywhere. Use the SunOS mt(1) command to reposition the tape appropriately.

# Fortran 90 I/O Considerations

Fortran 90 2.0 and FORTRAN 77 5.0 are I/O compatible. Executables containing intermixed £77 and £90 compilations can do I/O to the same unit from both the £77 and £90 parts of the program.

However, Fortran 90 provides some additional features:

- A file opened with FORM='BINARY' enables nonstandard I/O of raw data without record marks. This has nearly the same effect as FORM='UNFORMATTED', except that no record lengths are embedded in the file. As a result, it is not possible to backspace a FORM='BINARY' file. A READ statement on such a 'BINARY' file reads as much data as needed to fill all the variables on the input list
- ADVANCE='NO' enables nonadvancing I/O, as in:

```
write(*,"(a)",ADVANCE="NO") "Enter size= "
    read(*,*) n
```

- NAMELIST input features:
  - f90 allows the group name to be preceded by \$ or & on input. The Fortran 90 standard accepts only & and this is what a NAMELIST write outputs.

- f90 accepts \$ as the symbol terminating an input group unless the last data item in the group is CHARACTER, in which case the \$ is treated as input data.
- f90 allows NAMELIST input to start in the first column of a record.
- ENCODE and DECODE are recognized and implemented by £90 just as they are by £77.

# **Program Development**

This chapter briefly introduces two powerful program development tools, make and SCCS, that can be used very successfully with Fortran programs.

A number of good, commercially published books on using make and SCCS are currently available, including *Managing Projects with make*, by Andrew Oram and Steve Talbott, and *Applying RCS and SCCS*, by Don Bolinger and Tan Bronson. Both are from O'Reilly & Associates.

# Facilitating Program Builds With the make Utility

The make utility applies intelligence to the task of program compilation and linking. Typically, a large application consists of a set of source files and INCLUDE files, requiring linking with a number of libraries. Modifying any one or more of the source files requires recompilation of that part of the program and relinking. You can automate this process by specifying the interdependencies between files that make up the application along with the commands needed to recompile and relink each piece. With these specifications in a file of directives, make ensures that only the files that need recompiling are recompiled and that relinking uses the options and libraries you need to build the executable. The following discussion provides a simple example of how to use make. For a summary, see <code>make(1)</code>.

#### The Makefile

A file called makefile tells make in a structured manner which source and object files depend on other files. It also defines the commands required to compile and link the files.

For example, suppose you have a program of four source files and the makefile:

```
demo% ls
makefile
commonblock
computepts.f
pattern.f
startupcore.f
demo%
```

Assume both pattern.f and computepts.f have an INCLUDE of commonblock, and you wish to compile each.f file and link the three relocatable files, along with a series of libraries, into a program called pattern.

The makefile looks like this:

The first line of makefile indicates that making pattern depends on pattern.o, computepts.o, and startupcore.o. The next line and its continuations give the command for making pattern from the relocatable.o files and libraries.

Each entry in makefile is a rule expressing a target object's dependencies and the commands needed to make that object. The structure of a rule is:

target: dependencies-listTAB build-commands

- **Dependencies**. Each entry starts with a line that names the target file, followed by all the files the target depends on.
- **Commands**. Each entry has one or more subsequent lines that specify the Bourne shell commands that will build the target file for this entry. Each of these command lines must be indented by a tab character.

#### make Command

The make command can be invoked with no arguments, simply:

```
demo% make
```

The make utility looks for a file named makefile or Makefile in the current directory and takes its instructions from that file.

The make utility:

- Reads makefile to determine all the target files it must process, the files they depend on, and the commands needed to build them.
- Finds the date and time each file was last changed.
- Rebuilds any target file that is older than any of the files it depends on, using the commands from makefile for that target.

#### **Macros**

The make utility's *macro* facility allows simple, parameterless string substitutions. For example, the list of relocatable files that make up the target program pattern can be expressed as a single macro string, making it easier to change.

A macro string definition has the form:

```
NAME = string
```

Use of a macro string is indicated by:

\$(NAME)

which is replaced by make with the actual value of the macro string.

This example adds a macro definition naming all the object files to the beginning of  ${\tt makefile}$ :

```
OBJ = pattern.o computepts.o startupcore.o
```

Now the macro can be used in both the list of dependencies as well as on the f77 link command for target pattern in makefile:

```
pattern: $(OBJ)
     f77 $(OBJ) --lcore77 --lcore --lsunwindow \
     --lpixrect --o pattern
```

For macro strings with single-letter names, the parentheses may be omitted.

## Overriding of Macro Values

The initial values of make macros can be overridden with command-line options to make. For example:

Now a simple make command without arguments uses the value of FFLAGS set above. However, this can be overridden from the command line:

```
demo% make "FFLAGS=-u -O"
```

Here, the definition of the FFLAGS macro on the make command line overrides the makefile initialization, and both the -O flag and the -u flag are passed to £77. Note that "FFLAGS=" can also be used on the command to reset the macro to a null string so that it has no effect.

#### Suffix Rules in make

To make writing a makefile easier, make will use its own default rules depending on the suffix of a target file. Recognizing the <code>.f</code> suffix, make uses the <code>f77</code> compiler, passing as arguments any flags specified by the <code>FFLAGS</code> macro, the <code>-c</code> flag, and the name of the source file to be compiled.

The example below demonstrates this rule twice:

make uses default rules to compile computepts.f and startupcore.f.

Similarly, suffix rules for .f90 files will also invoke the f90 compiler.

# Version Tracking and Control With SCCS

SCCS stands for Source Code Control System. SCCS provides a way to:

- Keep track of the evolution of a source file—its change history
- Prevent a source file from being simultaneously changed by other developers
- Keep track of the version number by providing version stamps

The basic three operations of SCCS are:

- Putting files under SCCS control
- Checking out a file for editing
- Checking in a file

This section shows you how to use SCCS to perform these tasks, using the previous program as an example. Only basic SCCS is described and only three SCCS commands are introduced: create, edit, and delget.

# Controlling Files With SCCS

Putting files under SCCS control involves:

- Making the SCCS directory
- Inserting SCCS ID keywords into the files (this is optional)
- Creating the SCCS files

#### Making the SCCS Directory

To begin, you must create the SCCS subdirectory in the directory in which your program is being developed. Use this command:

demo% mkdir SCCS

SCCS must be in uppercase.

## **Inserting SCCS ID Keywords**

Some developers put one or more SCCS ID keywords into each file, but that is optional. These keywords are later identified with a version number each time the files are checked in with an SCCS get or delget command. There are three likely places to put these strings:

- Comment lines
- Parameter statements
- Initialized data

The advantage of using keywords is that the version information appears in the source listing and compiled object program. If preceded by the string @(#), the keywords in the object file can be printed using the what command.

Included header files that contain only parameter and data definition statements do not generate any initialized data, so the keywords for those files usually are put in comments or in parameter statements. In some files, like ASCII data files or makefiles, the SCCS information will appear in comments.

SCCS keywords appear in the form %keyword% and are expanded into their values by the SCCS get command. The most commonly used keywords are:

 $\$  expands to the identifier string  $\$  (#) recognized by the what command. %M% expands to the name of the source file. %I% expands to the version number of this SCCS maintained file. %E% expands to the current date.

For example, you could identify the makefile with a make comment containing these keywords:

```
# %Z%%M% %I% %E%
```

The source files, startupcore.f, computepts.f, and pattern.f, can be identified by initialized data of the form:

```
CHARACTER*50 SCCSID
DATA SCCSID/"%Z%%M% %I% %E%\n"/
```

When this file is processed by SCCS, compiled, and the object file processed by the SCCS what command, the following is displayed:

```
demo% f77 -c pattern.f
...
demo% what pattern
pattern:
    pattern.f 1.2 96/06/10
```

You can also create a PARAMETER named  $\tt CTIME$  that is automatically updated whenever the file is accessed with  $\tt get.$ 

```
CHARACTER*(*) CTIME
PARAMETER ( CTIME="%E%")
```

INCLUDE files can be annotated with a Fortran comment containing the SCCS stamp:

C

Now you can put these files under control of SCCS with the SCCS create command:

```
demo% sccs create makefile commonblock startupcore.f \
   computepts.f pattern.f
demo%
```

# Checking Files Out and In

Once your source code is under SCCS control, you use SCCS for two main tasks: to *check out* a file so that you can edit it, and to *check in* a file you have finished editing.

Check out a file with the sccs edit command. For example:

```
demo% sccs edit computepts.f
```

SCCS then makes a writable copy of computepts.f in the current directory, and records your login name. Other users cannot check the file out while you have it checked out, but they can find out who has checked it out.

When you have completed your editing, check in the modified file with the sccs delget command. For example:

```
demo% sccs delget computepts.f
```

This command causes the SCCS system to:

- Make sure that you are the user who checked out the file by comparing login names
- Prompt for a comment from you on the changes
- Make a record of what was changed in this editing session
- Delete the writable copy of computepts.f from the current directory
- Replace it by a read-only copy with the SCCS keywords expanded

The sccs delget command is a composite of two simpler SCCS commands, delta and get. The delta command performs the first three tasks in the list above; the get command performs the last two tasks.

# Libraries

This chapter describes how to use and create libraries of subprograms. Both *static* and *dynamic* libraries are discussed.

# **Understanding Libraries**

A software *library* is usually a set of subprograms that have been previously compiled and organized into a single binary *library file*. Each member of the set is called a library *element* or *module*. The linker searches the library files, loading object modules referenced by the user program while building the executable binary program. See *ld*(1) and the Solaris *Linker and Libraries Guide* for details.

There are two basic kinds of software libraries:

- Static library. A library in which modules are bound into the executable file before
  execution. Static libraries are commonly named libname.a. The .a suffix refers to
  archive.
- **Dynamic library**. A library in which modules can be bound into the executable program at runtime. Dynamic libraries are commonly named libname.so. The .so suffix refers to *shared object*.

Typical system libraries that have both static (.a) and dynamic (.so) versions are:

- FORTRAN 77 libraries: libF77, libM77
- Fortran 90 libraries: libfsu, libfui, libfai, libfai2, libfsunai, libfprodai, libfminlai, libfmaxlai, libminvai, libmaxvai, libf77compat
- VMS Fortran libraries: libv77
- C libraries: libc

There are two advantages to the use of libraries:

- There is no need to have source code for the library routines that a program calls.
- Only the needed modules are loaded.

Library files provide an easy way for programs to share commonly used subroutines. You need only name the library when linking the program, and those library modules that resolve references in the program are linked and merged into the executable file.

# **Specifying Linker Debugging Options**

Summary information about library usage and loading can be obtained by passing additional options to the linker through the LD\_OPTIONS environment variable. The compiler calls the linker with these options (and others it requires) when generating object binary files.

Using the compiler to call the linker is always recommended over calling the linker directly because many compiler options require specific linker options or library references, and linking without these could produce unpredictable results.

```
demo% setenv LD_OPTIONS '-m -Dfiles'
demo% f77 -o myprog myprog.f
```

Example: Using LD\_OPTIONS to create a load map:

Some linker options do have compiler command-line equivalents that can appear directly on the £77 or £90 command. These include

-Bx, -dx, -G, -h name, -R path, and -z t ext. See the f77(1) and f90(1) man pages or the Fortran User's Guide for details.

More detailed examples and explanations of linker options and environment variables can be found in the Solaris *Linker and Libraries Guide*.

## Generating a Load Map

The linker -m option generates a load map that displays library linking information. The routines linked during the building of the executable binary program are listed together with the libraries that they come from.

Example: Using -m to generate a load map:

```
demo% setenv LD_OPTIONS '-m'
demo% f77 any.f
any.f:
    MAIN:
    LINK EDITOR MEMORY MAP
```

```
output
         input
                 virtual
section section address
                                size
                 100d4
.interp
                               11 (null)
          .interp 100d4
.hash
                100e8
                              2e8
          .hash
                  100e8
                               2e8 (null)
.dynsym
                 103d0
                               650
          .dynsym 103d0
                               650 (null)
.dynstr
                10a20
                               366
           .dynstr 10a20
                               366 (null)
                              1e70
.text
                10c90
                 10c90 00 /opt/SUNWspro/SC5.0/lib/crti.o
.text
                 10c90
                         f4 /opt/SUNWspro/SC5.0/lib/crt1.o
                 10d84 00 /opt/SUNWspro/SC5.0/lib/values-xi.o
.text
.text
                 10d88
                         d20 sparse.o
```

# **Listing Other Information**

Additional linker debugging features are available through the linker's -Dkeyword option. A complete list can be displayed using -Dhelp.

Example: List linker debugging aid options using the -Dhelp option:

```
demo%

1d -Dhelp

...

debug: args display input argument processing debug: bindings display symbol binding; debug: detail provide more information debug: entry display entrance criteria descriptors

...

demo%
```

For example, the -Dfiles linker option lists all the files and libraries referenced during the link process:

See the Linker and Libraries Guide for further information on these linker options.

# Consistent Compiling and Linking

Ensuring a consistent choice of compiling and linking options is critical whenever compilation and linking are done in separate steps. Compiling any part of a program with any of the following options requires linking with the same options:

```
-a, -autopar, -Bx, -fast, -G, -Lpath, -lname, -mt, -nolib, -norunpath, -p, -pg, -xlibmop
```

Example: Compiling sbr.f with -a and smain.f without it, then linking in separate steps (-a invokes toov old style profiling):

```
demo% f77 -c -a sbr.f
demo% f77 -c smain.f
demo% f77 -a sbr.o smain.o {link step; pass -a to the linker}
```

Also, a number of options require that *all* source files be compiled with that option. These include:

```
-autopar, -cg92, -dx, -dalign, -dbl, -explicitpar, -f, -misalign, -native, -parallel,
```

See the f77(1) and f90(1) man pages and the Fortran User's Guide for details on all compiler options.

# Setting Library Search Paths and Order

The linker searches for libraries at several locations and in a certain prescribed order. Some of these locations are standard paths, while others depend on the compiler options <code>-Rpath</code>, <code>-llibrary</code>, and <code>-Ldir</code> and the environment variable <code>LD\_LIBRARY\_PATH</code>.

## Search Order for Standard Library Paths

The standard library search paths used by the linker are determined by the installation path, and they differ for static and dynamic loading. <code><install-point></code> is the path to where the Fortran compilers have been installed. In a standard install of the software this is <code>/opt</code>.

#### Static Linking

While building the executable file, the static linker searches for any libraries in the following paths (among others), in the specified order:

<pre><install-point>/SUNWspro/lib</install-point></pre>	Sun shared libraries
/usr/ccs/lib/	Standard location for SVr4 software
/usr/lib	Standard location for UNIX software

These are the *default* paths used by the linker.

#### **Dynamic Linking**

The dynamic linker searches for *shared* libraries at runtime, in the specified order:

- Paths specified by user with -Rpath
- <install-point>/SUNWspro/lib/
- /usr/lib standard UNIX default

The search paths are built into the executable.

# LD\_LIBRARY\_PATH Environment Variable

Use the LD\_LIBRARY\_PATH environment variable to specify directory paths that the linker should search for libraries specified with the -library option.

Multiple directories can be specified, separated by a colon. Typically, the  ${\tt LD\_LIBRARY\_PATH}$  variable contains two lists of colon-separated directories separated by a semicolon:

dirlist1; dirlist2

The directories in dirlist1 are searched first, followed by any explicit -Ldir directories specified on the command line, followed by dirlist2 and the standard directories.

That is, if the compiler is called with any number of occurrences of -L, as in:

```
f77 ... -Lpath1 ... -Lpathn ...
```

then the search order is:

dirlist1 path1 ... pathn dirlist2 standard\_paths

When the LD\_LIBRARY\_PATH variable contains only one colon-separated list of directories, it is interpreted as *dirlist2*.

In the Solaris 7 operating environment, a similar environment variable,  $LD\_LIBRARY\_PATH\_64$  can be used to override  $LD\_LIBRARY\_PATH$  when searching for 64-bit dependencies. See the Solaris *Linker and Libraries Guide* and the Id(1) man page for details.

- On a 32-bit SPARC processor and linking with Solaris 7, LD\_LIBRARY\_PATH\_64 is ignored.
- If only LD LIBRARY PATH is defined, it is used for both 32-bit and 64-bit linking.
- If both LD\_LIBRARY\_PATH and LD\_LIBRARY\_PATH\_64 are defined, 32-bit linking will be done using LD\_LIBRARY\_PATH, and 64-bit linking with LD\_LIBRARY\_PATH\_64.

**Note -** Use of the LD\_LIBRARY\_PATH environment variable with production software is strongly discouraged. Although useful as a temporary mechanism for influencing the runtime linker's search path, *any* dynamic executable that can reference this environment variable will have its search paths altered. You might see unexpected results or a degradation in performance.

# Library Search Path and Order—Static Linking

Use the -llibrary compiler option to name additional libraries for the linker to search when resolving external references. For example, the option -lmylib adds the library libmylib.so or libmylib.a to the search list.

The linker looks in the standard directory paths to find the additional libmylib library. The -L option (and the LD\_LIBRARY\_PATH environment variable) creates a list of paths that tell the linker where to look for libraries outside the standard paths.

Were libmylib.a in directory /home/proj/libs, then the option -L/home/proj/libs would tell the linker where to look when building the executable:

demo% f77 -o pgram part1.o part2.o -L/home/proj/libs -lmylib

## Command-Line Order for -1 library Options

For any particular unresolved reference, libraries are searched only once, and only for symbols that are undefined at that point in the search. If you list more than one library on the command line, then the libraries are searched in the order in which they are found on the command line. Place -1 library options as follows:

- Place the -1 *library* option after any .f, .for, .F, .f90, or .o files.
- If you call functions in libx, and they reference functions in liby, then place -1x before -1y.

#### Command-Line Order for -Ldir Options

The -Ldir option adds the dir directory path to the library search list. The linker searches for libraries first in any directories specified by the -L options and then in the standard directories. This option is useful only if it is placed *preceding* the -1 library options to which it applies.

# Library Search Path and Order—Dynamic Linking

With dynamic libraries, changing the library search path and order of loading differs from the static case. Actual linking takes place at runtime rather than build time.

#### Specifying Dynamic Libraries at Build Time

When *building* the executable file, the linker records the paths to shared libraries in the executable itself. These search paths can be specified using the  $\neg \mathbb{R}$  path option. This is in contrast to the  $\neg \mathbb{L}$  dir option which indicates at buildtime where to find the library specified by a  $\neg \mathbb{L}$  library option, but does not record this path into the binary executable.

The directory paths that were built in when the executable was created can be viewed using the dump command.

Example: List the directory paths built into a.out:

```
demo% f77 program.f -R/home/proj/libs -L/home/proj/libs -lmylib
demo% dump -Lv a.out | grep RPATH
[5] RPATH /home/proj/libs:/opt/SUNWspro/lib
```

## Specifying Dynamic Libraries at Runtime

At *runtime*, the linker determines where to find the dynamic libraries that an executable needs from:

- The value of LD\_LIBRARY\_PATH at runtime
- The paths that had been specified by -R at the time the executable file was built

As noted earlier, use of  $\mbox{LD\_LIBRARY\_PATH}$  can have unexpected side-effects and is not recommended.

## Fixing Errors During Dynamic Linking

When the dynamic linker cannot locate a needed library, it issues this error message:

```
ld.so: proq: fatal: libmylib.so: can't open file:
```

The message indicates that the libraries are not where they are supposed to be. Perhaps you specified paths to shared libraries when the executable was built, but the libraries have subsequently been moved. For example, you might have built a .out with your own dynamic libraries in /my/libs/, and then later moved the libraries to another directory.

Use 1dd to determine where the executable expects to find the libraries:

```
demo% ldd a.out
    libsolib.so => /export/home/proj/libsolib.so
    libF77.so.4 => /opt/SUNWspro/lib/libF77.so.4
    libc.so.1 => /usr/lib/libc.so.1
    libdl.so.1 => /usr/lib/libdl.so.1
```

If possible, move or copy the libraries into the proper directory or make a soft link to the directory (using ln -s) in the directory that the linker is searching. Or, it could be that LD\_LIBRARY\_PATH is not set correctly. Check that LD\_LIBRARY\_PATH includes the path to the needed libraries at runtime.

# **Creating Static Libraries**

Static library files are built from precompiled object files (.0 files) using the ar(1) utility.

The linker extracts from the library any elements whose entry points are referenced within the program it is linking, such as a subprogram, entry name, or COMMON block initialized in a BLOCKDATA subprogram. These extracted elements (routines) are bound permanently into the a .out executable file generated by the linker.

## Tradeoffs for Static Libraries

There are three main issues to keep in mind regarding static, as compared to dynamic, libraries and linking:

Static libraries are more self-contained but less adaptable.

If you bind an a .out executable file *statically*, the library routines it needs become part of the executable binary. However, if it becomes necessary to update a static library routine bound into the a .out executable, the entire a .out file must be relinked and regenerated to take advantage of the updated library. With *dynamic* libraries, the library is not part of the a .out file and linking is done at runtime. To

take advantage of an updated dynamic library, all that is required is that the new library be installed on the system.

■ The "elements" in a static library are individual compilation units, .o files.

Since a single compilation unit (a source file) can contain more than one subprogram, these routines when compiled together become a single module in the static library. This means that *all* the routines in the compilation unit are loaded together into the a.out executable, even though only one of those subprograms was actually called. This situation can be improved by optimizing the way library routines are distributed into compilable source files. (Still, only those library modules actually referenced by the program are loaded into the executable.)

Order matters when linking static libraries.

The linker processes its input files in the order in which they appear on the command line—left to right. When the linker decides whether or not to load an element from a library, its decision is determined by the library elements that it has already processed. This order is not only dependent on the order of the elements as they appear in the library file but also on the order in which the libraries are specified on the compile command line.

Example: If the Fortran program is in two files, main.f and crunch.f, and only the latter accesses a library, it is an error to reference that library before crunch.f or crunch.o:

# Creation of a Simple Static Library

Suppose that you can distribute all the routines in a program over a group of source files and that these files are wholly contained in the subdirectory test\_lib/.

Suppose further that the files are organized in such a way that they each contain a single principal subprogram that would be called by the user program, along with any "helper" routines that the subprogram might call but that are called from no other routine in the library. Also, any helper routines called from more than one library routine are gathered together into a single source file. This gives a reasonably well-organized set of source and object files.

Assume that the name of each source file is taken from the name of the first routine in the file, which in most cases is one of the principal files in the library:

The lower-level "helper" routines are gathered together into the file etc.f. The other files can contain one or more subprograms.

First, compile each of the library source files, using the -c option, to generate the corresponding relocatable .o files:

```
demo% f77 -c *.f
delte.f:
    delte:
    q_fixx:
dropx.f:
    dropx:
etc.f:
    q_fill:
    q_step:
    q_node:
    q_warn:
 ...etc
demo% ls
total 42
 2 dropx.f
                4 etc.o
                             2 linkz.f 4 markx.o
                4 dropx.o 2 evalx.f 2 etc.f 4 evalx.o
                                           4 linkz.o
2 markx.f
 2 delte.f
                                                            2 point.f
 4 delte.o
                                                             4 point.o
demo%
```

Now, create the static library testlib.a using ar:

```
demo% ar cr testlib.a *.o
```

To use this library, either include the library file on the compilation command or use the -1 and -L compilation options. The example uses the .a file directly:

Notice that the main program calls only two of the routines in the library. You can verify that the uncalled routines in the library were not loaded into the executable file by looking for them in the list of names in the executable displayed by nm:

```
demo% nm trylib | grep FUNC | grep delte demo% nm trylib | grep FUNC | grep markx demo% ..etc
```

In the preceding example, grep finds entries in the list of names only for those library routines that were actually called.

Another way to reference the library is through the -1 library and -1 path options. Here, the library's name would have to be changed to conform to the 1 ibname. a convention:

```
demo% mv test_lib/testlib.a test_lib/libtestlib.a
demo% f77 -o trylib trylib.f -Ltest_lib -ltestlib
trylib.f:
    MAIN:
```

The -llibrary and -Lpath options are used with libraries installed in a commonly accessible directory on the system, like /usr/local/lib, so that other users can reference it. For example, if you left libtestlib.a in /usr/local/lib, other users could be informed to compile with the following command:

```
demo% f77 -o myprog myprog.f -L/usr/local/lib -ltestlib
```

#### Replacement in a Static Library

It is not necessary to recompile an entire library if only a few elements need recompiling. The -r option of ar permits replacement of individual elements in a static library.

Example: Recompile and replace a single routine in a static library:

```
demo% f77 --c point.f
demo% ar r testlib.a point.o
demo%
```

## Ordering Routines in a Static Library

To order the elements in a static library when it is being built by ar, use the commands *lorder*(1) and *tsort*(1):

```
demo% ar cr mylib.a 'lorder exg.o fofx.o diffz.o | tsort'
```

# **Creating Dynamic Libraries**

Dynamic library files are built by the linker 1d from precompiled object modules that can be bound into the executable file *after* execution begins.

Another feature of a dynamic library is that modules can be used by other executing programs in the system *without* duplicating modules in each program's memory. For this reason, a dynamic library is also a *shared* library.

A dynamic library offers the following features:

- The object modules are *not* bound into the executable file by the linker during the compile-link sequence; such binding is deferred until runtime.
- A shared library module is bound into system memory when the first running program references it. If any subsequent running program references it, that reference is mapped to this first copy.
- Maintaining programs is easier with dynamic libraries. Installing an updated dynamic library on a system immediately affects all the applications that use it without requiring relinking of the executable.

## Tradeoffs for Dynamic Libraries

Dynamic libraries introduce some additional tradeoff considerations:

■ Smaller a.out file

Deferring binding of the library routines until execution time means that the size of the executable file is less than the equivalent executable calling a static version of the library; the executable file does not contain the binaries for the library routines.

Possibly smaller process memory utilization

When several processes using the library are active simultaneously, only one copy of the memory resides in memory and is shared by all processes.

Possibly increased overhead

Additional processor time is needed to load and link-edit the library routines during runtime. Also, the library's position-independent coding might execute more slowly than the relocatable coding in a static library.

■ Possible overall system performance improvement

Reduced memory utilization due to library sharing should result in better overall system performance (reduced I/O access time from memory swapping).

Performance profiles among programs vary greatly from one to another. It is not always possible to determine or estimate in advance the performance improvement

(or degradation) between dynamic versus static libraries. However, if both forms of a needed library are available to you, it would be worthwhile to evaluate the performance of your program with each.

## Position-Independent Code and -pic

Position-independent code (PIC) can be bound to any address in a program without requiring relocation by the link editor. Such code is inherently sharable between simultaneous processes. Thus, if you are building a dynamic, shared library, you must compile the component routines to be position-independent (by using compiler options -pic or -PIC).

In position-independent code, each reference to a global item is compiled as a reference through a pointer into a global offset table. Each function call is compiled in a relative addressing mode through a procedure linkage table. The size of the global offset table is limited to 8 Kbytes on SPARC processors. The -PIC compiler option is similar to -pic, but -PIC allows the global offset table to span the range of 32-bit addresses.

Version 5.0 of £77 and version 2.0 of £90 introduce a more flexible compiler flag, -xcode=v, for specifying the code address space of a binary object. With this compiler flag, 32-, 44-, or 64-bit absolute addresses can be generated, as well as small and large model position-independent code. -xcode=pic13 is equivalent to -pic, and -xcode=pic32 is equivalent to -PIC. See the f77(1) and f90(1) man pages, or the Fortran User's Guide, for details.

# **Binding Options**

You can specify dynamic or static library binding when you compile. These options are actually linker options, but they are recognized by the compiler and passed on to the linker.

- -Bdynamic | -Bstatic
- -Bdynamic sets the preference for shared, dynamic binding whenever possible.
- -Bstatic restricts binding to static libraries only.

When both static and dynamic versions of a library are available, use this option to toggle between preferences on the command line:

```
f77 prog.f -Bdynamic -lwells -Bstatic -lsurface
```

-dy | -dn

Allows or disallows dynamic linking for the entire executable. (This option may appear on the command line only once.)

-dy allows dynamic, shared libraries to be linked. -dn does not allow linking of dynamic libraries.

#### Binding in 64-Bit Environments

Some static system libraries, such as libm.a and libc.a, are not available on 64-bit environments with Solaris 7. These are supplied as dynamic libraries only. Use of -dn in these environments will result in an error indicating that some static system libraries are missing. Also, ending the compiler command line with -Bstatic will have the same effect.

To link with static versions of specific libraries, use a command line that looks something like:

```
f77 -o prog prog.f -Bstatic -labc -lxyz -Bdynamic
```

Here the user's libabc.a and libxyz.a file are linked (rather than libabc.so or libxyz.so), and the final -Bdynamic insures that the remaining libraries, including system libraries, and dynamically linked.

In more complicated situations, it may be necessary to explicitly reference each system and user library on the link step with the appropriate <code>-Bstatic</code> or <code>-Bdynamic</code> as required. First use <code>LD\_OPTIONS</code> set to '<code>-Dfiles</code>' to obtain a listing of all the libraries needed. Then perform the link step with <code>-nolib</code> (to suppress automatic linking of system libraries) and explicit references to the libraries you need. For example:

```
f77 -xarch=v9 -o cdf -nolib cdf.o-Bstatic -lF77 -lM77 -lsunmath -Bdynamic -lm -lc
```

## Naming Conventions

To conform to the dynamic library naming conventions assumed by the link loader and the compilers, assign names to the dynamic libraries that you create with the prefix lib and the suffix .so. For example, libmyfavs.so could be referenced by the compiler option -lmyfavs.

The linker also accepts an optional version number suffix: for example, libmyfavs.so.1 for version *one* of the library.

The compiler's -h name option records name as the name of the dynamic library being built.

# A Simple Dynamic Library

Building a dynamic library requires a compilation of the source files with the -pic or -PIC option and linker options -G, -ztext, and -h name. These linker options are available through the compiler command line.

You can create a dynamic library with the same files used in the static library example.

Example: Compile with -pic and other linker options:

```
demo% f77 -o libtestlib.so.1 -G -pic -ztext -hlibtestlib.so.1 *.f
delte.f:
      delte:
     q_fixx:
dropx.f:
     dropx:
etc.f:
      q_fill:
     q_step:
     q_node:
      q_warn:
evalx.f:
      evalx:
linkz.f:
     linkz:
markx.f:
     markx:
point.f:
     point:
Linking:
```

- -G tells the linker to build a dynamic library.
- -ztext warns you if it finds anything other than position-independent code, such as relocatable text.

Example: Make an executable file a . out using the dynamic library:

```
demo% f77 -o trylib -R'pwd' trylib.f libtestlib.so.1
trylib.f:
MAIN main:
demo% file trylib
trylib:ELF 32--bit MSB executable SPARC Version 1, dynamically linked, not stripped
demo% ldd trylib
    libtestlib.so.1 => /export/home/U/Tests/libtestlib.so.1
    libF77.so.4 => /opt/SUNWspro/lib/libF77.so.4
    libc.so.1 => /usr/lib/libc.so.1
    libdl.so.1 => /usr/lib/libdl.so.1
```

Note that the example uses the  $-\mathbb{R}$  option to bind into the executable the path (the current directory) to the dynamic library.

The file command shows that the executable is dynamically linked.

The ldd command shows that the executable, trylib, uses some shared libraries, including libtestlib.so.1; libf77, libdl, and libc are included by default by f77.

# Libraries Provided with Sun Fortran Compilers

The table shows the libraries installed with the compilers.

TABLE 4-1 Major Libraries Provided With the Compilers

Library	Name	Options Needed
£77 functions, nonmath	libF77	None
£77 functions, nonmath, multithread safe	libF77_mt	-parallel
f77 math library	libM77	None
f90 support intrinsics	libfsu	None
f90 interface	libfui	None
f90 array intrinsics libraries	libf*ai	None
f90/f77 I/O compatibility library	libf77compat	-lf77compat
VMS library	libV77	-1 <b>v</b> 77
Library used with Pascal, Fortran, and C	libpfc	None
Library of Sun math functions	libsunmath	None
POSIX bindings	libFposix	-lFposix

TABLE 4-1 Major Libraries Provided With the Compilers (continued)

f90 POSIX interface libposix9 -lposix9

POSIX bindings for extra runtime checking libFposix\_c -lFposix\_c

See also the math\_libraries README file for more information.

## VMS Library

The  $\mbox{lib}\mbox{V77}$  library is the VMS library, which contains two special VMS routines, idate and time.

To use either of these routines, include the -1V77 option.

For idate and time, there is a conflict between the VMS version and the version that traditionally is available in UNIX environments. If you use the -1V77 option, you get the VMS compatible versions of the idate and time routines.

See the Fortran Library Reference Manual and the FORTRAN 77 Language Reference Manual for details on these routines.

# POSIX Library

There are two versions of POSIX bindings provided with FORTRAN 77:

- libFposix, which is just the bindings (-lFposix)
- libFposix\_c, which does some runtime checking to make sure you are passing correct handles (-lFposix\_c)

If you pass bad handles:

- libFposix\_c returns an error code (ENOHANDLE).
- libFposix core dumps with a segmentation fault.

Of course, the checking is time-consuming, and  $\mbox{libFposix\_c}$  is several times slower.

Both POSIX libraries come in static and dynamic forms.

The POSIX bindings provided are for IEEE Standard 1003.9-1992.

IEEE 1003.9 is a binding of 1003.1-1990 to FORTRAN (X3.8-1978).

For more information, see these POSIX.1 documents:

■ ISO/IEC 9945-1:1990

- IEEE Standard 1003.1-1990
- IEEE Order number SH13680
- IEEE CS Catalog number 1019

To find out precisely what POSIX is, you need both the 1003.9 and the POSIX.1 documents.

The POSIX library for f90 is libposix9.

# Shippable Libraries

If your executable uses a Sun dynamic library that is listed in the runtime.libraries README file, your license includes the right to redistribute the library to your customer.

This README file is located in the READMEs directory:

<install-point>/SUNWspro/READMEs/

Do not redistribute or otherwise disclose the header files, source code, object modules, or static libraries of object modules in any form.

Refer to your software license for more details.

# Program Analysis and Debugging

This chapter presents a number of Sun Fortran compiler features that facilitate program analysis and debugging.

# **Global Program Checking**

The f77 compiler's -Xlistx options provide a valuable way to analyze a source program for inconsistencies and possible runtime problems. The analysis performed by the compiler is *global*, across subprograms.

- -Xlistx reports errors in alignment, agreement in number and type for subprogram arguments, common block, parameter, and various other kinds of errors.
- -Xlistx also can be used to make detailed source code listings and cross-reference tables.

**Note -** The £90 compiler provides only a subset of the --Xlist options described here. A conventional cross-reference map is produced, but complete global program checking is not performed.

#### **GPC** Overview

Global program checking (GPC), invoked by the -Xlistx option, does the following:

■ Enforces type-checking rules of Fortran more stringently than usual, especially between separately compiled routines

- Enforces some portability restrictions needed to move programs between different machines or operating systems
- Detects legal constructions that nevertheless might be suboptimal or error-prone
- Reveals other potential bugs and obscurities

In particular, global checking reports problems such as:

- Interface problems
  - Conflicts in number and type of dummy and actual arguments
  - Wrong types of function values
  - Possible conflicts due to data type mismatches in common blocks between different subprograms
- Usage problems
  - Function used as a subroutine or subroutine used as a function
  - Declared but unused functions, subroutines, variables, and labels
  - Referenced but not declared functions, subroutines, variables, and labels
  - Usage of unset variables
  - Unreachable statements
  - Implicit type variables
  - Inconsistency of the named common block lengths, names, and layouts

# How to Invoke Global Program Checking

The -xlist option on the command line invokes the compiler's global program analyzer. There are a number of -xlistx suboptions, as described in the sections that follow.

Example: Compile three files for basic global program checking:

```
demo% f77 -Xlist any1.f any2.f any3.f
```

In the preceding example, the compiler:

Produces output listings in the file any1.1st

■ Compiles and links the program if there are no errors

#### Screen Output

Normally, output listings produced by -Xlistx are written to a file. To display directly to the screen, use -Xlisto to write the output file to /dev/tty.

Example: Display to terminal:

demo% f77 -Xlisto /dev/tty any1.f

#### **Default Output Features**

The -Xlist option provides a combination of features available for output. With no other -Xlist options, you get the following by default:

- The listing file name is taken from the first input source or object file that appears, with the extension replaced by .lst
- A line-numbered source listing
- Error messages (embedded in listing) for inconsistencies across routines
- Cross-reference table of the identifiers
- Pagination at 66 lines per page and 79 columns per line
- No call graph
- No expansion of include files

## File Types

The checking process recognizes all the files in the compiler command line that end in .f, .f90, .for, .F, or .o. The .o files supply the process with information regarding only global names, such as subroutine and function names.

## Analysis Files (.fln Files)

Programs compiled with --xlist options have their analysis data built into the binary files automatically. This enables global program checking over programs in libraries.

Alternatively, the compiler will save individual source file analysis results into files with a .fln suffix if the -Xlistfln dir option is also specified. dir indicates the directory to receive these files.

# Some Examples of --Xlist and Global Program Checking

Here is a listing of the Repeat.f source code used in the following examples:

```
demo% cat Repeat.f
       PROGRAM repeat
         pn1 = REAL( LOC ( rp1 ) )
        CALL subr1 ( pn1 )
        CALL nwfrk ( pn1 )
        PRINT *, pn1
       END ! PROGRAM repeat
       SUBROUTINE subr1 ( x )
         IF ( \times .GT. 1.0 ) THEN
         CALL subr1 ( x * 0.5 )
        END IF
       END
       SUBROUTINE nwfrk( ix )
         EXTERNAL fork
         INTEGER prnok, fork
        PRINT *, prnok ( ix ), fork ( )
       INTEGER FUNCTION prnok ( x )
        prnok = INT (x) + LOC(x)
       SUBROUTINE unreach_sub()
        CALL sleep(1)
       END
```

#### Example: Use -XlistE to show errors and warnings:

```
repeat/nwfrk//prnok in line #21 but set as real
                by repeat in line #2
subroutine subr1
                 CALL subr1 ( x * 0.5 )
**** WAR #348: recursive call for "subr1". See dynamic calls:
                "Repeat.f" line #3
subroutine nwfrk
                 PRINT *, prnok ( ix ), fork ( )
**** ERR #418: argument "ix" is integer*4, but dummy argument
                is real
                See: "Repeat.f" line #20
subroutine unreach_sub
   24 SUBROUTINE unreach_sub()
**** WAR #338: subroutine "unreach_sub" isn"t called from program
        Wed Feb 24 10:40:32 1999
Files:
Lines:
             2 (Sources: 1; libraries: 1)
             26 (Sources: 26; Library subprograms:2)
Routines:
            5 (MAIN: 1; Subroutines: 3; Functions: 1)
            5 (Errors: 3; Warnings: 2)
Messages:
demo%
```

# Compiling the same program with --Xlist also produces a cross-reference table on standard output:

```
CROSS REFERENCE TABLE
 Source file: Repeat.f
   Definition/Declaration
D
IJ
     Simple use
Μ
     Modified occurrence
     Actual argument
Subroutine/Function call
Α
C
I
     Initialization: DATA or extended declaration
     Occurrence in EQUIVALENCE
E
      Occurrence in NAMELIST
     PROGRAM FORM
Program
-----
                     D
repeat
          <repeat>
Functions and Subroutines
     int*4 <nwfrk> DC 15:D 16:D 17:C
fork
int intrinsic
           loc
     intrinsic
          2:C
           <repeat> C 4:C
nwfrk
```

		<nwfrk></nwfrk>	D	14:D	
prnok	int*4	<nwfrk> <prnok></prnok></nwfrk>	DC DM	16:D 20:D	17:C 21:M
real	intrin	sic <repeat></repeat>	С	2:C	
sleep		<unreach_sub></unreach_sub>		С	25:C
subr1		<repeat> <subr1></subr1></repeat>	C DC	3:C 8:D	10:C
unreach_	sub	<unreach_sub></unreach_sub>		D	24:D

#### Output from compiling f77 -Xlist Repeat.f (Continued)

ix	int*4	dummy <nwfrk></nwfrk>	DA	14:D	17:A		
pn1	real*4	<repeat></repeat>	UMA	2:M	3:A	4:A	5 <b>:</b> U
rp1	real*4	<repeat></repeat>	А	2:A			
x	real*4	dummy <subr1> <prnok></prnok></subr1>	DU DUA	8:D 20:D	9:U 21:A	10:U 21:U	

Date: Tue Feb 22 13:15:39 1995
Files: 2 (Sources: 1; libraries: 1)
Lines: 26 (Sources: 26:7:7) Lines: 26 (Sources: 26; Library subprograms:2)
Routines: 5 (MAIN: 1; Subroutines: 3; Functions: 1)
Messages: 5 (Errors: 3; Warnings: 2)

demo%

#### In the cross-reference table in the preceding example:

■ ix is a 4-byte integer:

Variables and Arrays

- Used as an argument in the routine nwfrk
- At line 14, used as a declaration of argument
- At line 17, used as an actual argument
- pn1 is a 4-byte real in the routine repeat:
  - At line 2, modified
  - At line 3, argument
  - At line 4, argument

- At line 5. used
- rp1 is a 4-byte real in the routine, repeat. At line 2, it is an argument.
- x is a 4-byte real in the routines subr1 and prnok:
  - In subr1, at line 8, defined; used at lines 9 and 10
  - In prnok, at line 20, defined; at line 21, used as an argument

# **Suboptions for Global Checking Across Routines**

The basic global cross-checking option is -Xlist with no suboption. It is a combination of suboptions, each of which could have been specified separately.

The following sections describe options for producing the listing, errors, and cross-reference table. Multiple suboptions may appear on the command line.

#### **Suboption Syntax**

Add suboptions according to the following rules:

- Append the suboption to -Xlist.
- Put no space between the -Xlist and the suboption.
- Use only one suboption per -Xlist.

#### --Xlist and its Suboptions

Combine suboptions according to the following rules:

- The most general option is -Xlist (listing, errors, cross-reference table).
- Specific features can be combined using -Xlistc, -XlistE, -XlistL, or -XlistX.
- Other suboptions specify further details.

Example: Each of these two command lines performs the same task:

```
demo% f77 -Xlistc -Xlist any.f

demo% f77 -Xlistc any.f
```

The following table shows the reports generated by these basic --Xlist suboptions alone:

TABLE 5-1 Xlist Suboptions

Generated Report	Option
Errors, listing, cross-reference	-Xlist
Errors only	-Xlist <b>E</b>
Errors and source listing only	-Xlist <b>L</b>
Errors and cross-reference table only	-Xlist <b>X</b>
Errors and call graph only	-Xlist <b>c</b>

The following table summarizes all  ${\tt -Xlist}$  suboptions.

TABLE 5-2 Summary of --Xlist Suboptions

Option	Action
-Xlist (no suboption)	Shows errors, listing, and cross-reference table
-Xlist <b>c</b>	Shows call graphs and errors (f77 only)
-Xlist <b>E</b>	Shows errors
- Xlist <b>err</b> [nnn]	Suppresses error <i>nnn</i> in the verification report
-Xlist <b>f</b>	Produces fast output (f77 only)
-Xlist <b>fln</b> dir	Puts the .fln files in dir (f77 only)
-Xlist <b>h</b>	Shows errors from cross-checking stop compilation (f77 only)
-Xlist <b>I</b>	Lists and cross-checks include files

TABLE 5-2 Summary of --Xlist Suboptions (continued)

-Xlist <b>L</b>	Shows the listing and errors
-Xlist <b>l</b> n	Sets page breaks
-Xlist <b>o</b> name	Renames the -Xlist output report file
-Xlist <b>s</b>	Suppresses unreferenced symbols from cross-reference (f77 only)
-Xlist <b>v</b> n	Sets checking "strictness" level (f77 only)
-Xlistw[nnn]	Sets the width of output lines
- Xlist <b>war</b> [ <i>nnn</i> ]	Suppresses warning nnn in the report
-Xlist <b>X</b>	Shows just the cross-reference table and errors

## -Xlist Suboption Reference

This section describes the --Xlist suboptions. As noted, some are only available with £77.

#### -f77: -Xlistc — Show call graphs and cross-routine errors

Used alone, --Xlistc does not show a listing or cross-reference. It produces the call graph in a tree form, using printable characters. If some subroutines are not called from MAIN, more than one graph is shown. Each BLOCKDATA is printed separately with no connection to MAIN.

The default is *not* to show the call graph.

#### --XlistE - Show cross-routine errors

Used alone, --XlistE shows only cross-routine errors and does not show a listing or a cross-reference.

#### - Xlisterr[nnn] - Suppress error nnn

Use --Xlisterr to suppress a numbered error message from the listing or cross-reference.

For example: -Xlisterr338 suppresses error message 338. If *nnn* is not specified, all error messages are suppressed. To suppress additional specific errors, use this option repeatedly.

#### f77: -Xlist-f - Produce faster output

Use ——Xlistf to produce source file listings and a cross-checking report and to verify sources, but without generating object files.

The default without this option is to generate object files.

#### f77: -Xlist-flndir - Put .fln files into dir directory

Use --Xlistfln to specify the directory to receive .fln source analysis files. The directory specified (dir) must already exist. The default is to include the source analysis information directly within the object .o files (and not generate .fln files).

#### f77: -Xlist-h - Halt on errors

With --Xlisth, compilation stops if errors are detected while cross-checking the program. In this case, the report is redirected to stdout instead of the \*.lst file.

#### -Xlist-I - List and cross-check include files

If -XlistI is the only suboption used, include files are shown or scanned along with the standard -Xlist output (line numbered listing, error messages, and a cross-reference table).

- **Listing**—If the listing is not suppressed, then the include files are listed in place. Files are listed as often as they are included. The files are:
  - Source files
  - #include files
  - INCLUDE files
- **Cross-Reference Table**—If the cross reference table is not suppressed, the following files are all scanned while the cross reference table is generated:
  - Source files
  - #include files

INCLUDE files

The default is not to show include files.

#### -Xlist-L - Show listing and cross routine errors

Use ——XlistL to produce only a listing and a list of cross routine errors. This suboption by itself does not show a cross reference table. The default is to show the listing and cross reference table.

#### -Xlist-ln - Set the page length for pagination to n lines

Use --Xlistl to set the page length to something other than the default page size. For example, -Xlistl45 sets the page length to 45 lines. The default is 66.

With n=0 (-Xlist10) this option shows listings and cross-references with no page breaks for easier on-screen viewing.

#### -Xlist-o name - Rename the -Xlist output report file

Use --Xlisto to rename the generated report output file. (A space between o and name is required.) With --Xlisto name, the output is to name.lst.

To display directly to the screen, use the command: -Xlisto /dev/tty

#### f77: -Xlist-s - Suppress unreferenced identifiers

Use --Xlists to suppress from the cross reference table any identifiers defined in the include files but not referenced in the source files.

This suboption has no effect if the suboption -XlistI is used.

The default is not to show the occurrences in #include or INCLUDE files.

### f77: -Xlist-vn - Set level of checking strictness

n is 1,2,3, or 4. The default is 2 (-Xlistv2):

■ -Xlistv1

Shows the cross-checked information of all names in summary form only, with no line numbers. This is the lowest level of checking strictness—syntax errors only.

■ -Xlistv2

Shows cross-checked information with summaries and line numbers. This is the default level of checking strictness and includes argument inconsistency errors and variable usage errors.

#### ■ -Xlistv3

Shows cross-checking with summaries, line numbers, and common block maps. This is a high level of checking strictness and includes errors caused by incorrect usage of data types in common blocks in different subprograms.

#### ■ -Xlistv4

Shows cross-checking with summaries, line numbers, common block maps, and equivalence block maps. This is the strictest level of checking with maximum error detection.

#### -Xlist-w[nnn] - Set width of output line to n columns

Use ——Xlistw to set the width of the output line. For example, —Xlistw132 sets the page width to 132 columns. The default is 79.

#### -Xlist-war[nnn] - Suppress warning nnn in the report

Use ——Xlistwar to suppress a specific warning message from the output reports. If *nnn* is not specified, then all warning messages are suppressed from printing. For example, —Xlistwar338 suppresses warning message number 338. To suppress more than one, but not all warnings, use this option repeatedly.

# -Xlist-X - Show cross-reference table and cross routine errors

--XlistX produces a cross reference table and cross routine error list but no source listing.

# Some Examples Using Suboptions

Example: Use -Xlistwarnnn to suppress two warnings from a preceding example:

```
by repeat in line #2
subroutine nwfrk
                   PRINT *, prnok ( ix ), fork ( )
    17
**** ERR #418: argument "ix" is integer*4, but dummy argument
                 is real
                 See: "Repeat.f" line #20
         Wed Feb 24 10:40:32 1999
Lines:
Files:
              2 (Sources: 1; libraries: 1)
              26 (Sources: 26; Library subprograms:2)
Routines: 5 (MAIN: 1; Subroutines: 3; Functions: 1)
Messages: 5 (Errors: 3; Warnings: 2)
demo%
Example: Explain a message and find a type mismatch in program ShoGetc.f:
demo% cat ShoGetc.f
      CHARACTER*1 c
      i = getc(c)
      END
                                      Compile program
demo% f77 -silent ShoGetc.f
demo% a.out
               Type "Z" on keyboard. This causes run-time message. Why?
                               Program waits for input...
Note: IEEE floating-point exception flags raised:
  Invalid Operation;
 See the Numerical Computation Guide, ieee_flags(3M)
demo% f77 -XlistE -silent ShoGetc.f Compile with Global Program Checking
                                      and view listing
demo% cat ShoGetc.lst
FILE "ShoGetc.f"
program MAIN
    2
                i = getc(c)
**** WAR #320: variable "i" set but never referenced
              i = getc(c)
**** ERR #412: function "getc" used as real but declared as
                 integer*4
        Here is the error - function must be declared INTEGER.
                i = getc(c)
**** WAR #320: variable "c" set but never referenced
demo% cat ShoGetc.f
                       Modify program to declare getc INTEGER and run again.
      CHARACTER*1 c
      INTEGER getc
      i = getc(c)
     END
```

Type "Z" on keyboard Now no error.

demo% f77 -silent ShoGetc.f

demo% a.out

demo%

# **Special Compiler Options**

Some compiler options are useful for debugging. They check subscripts, spot undeclared variables, show stages of the compile-link sequence, display versions of software, and so on.

The Solaris linker has additional debugging aids. See ld(1), or run the command 1d -Dhelp at a shell prompt to see the online documentation.

# Subscript Bounds (-C)

The -C option adds checks for out-of-bounds array subscripts.

If you compile with  $\neg C$ , the compiler adds checks at runtime for out-of-bounds references on each array subscript. This action helps catch some situations that cause segmentation faults.

Example: Index out of range:

```
demo% cat indrange.f
    REAL a(10,10)
    k = 11
    a(k,2) = 1.0
    END

demo% f77 --C --silent indrange.f
demo% a.out
Subscript out of range on file indrange.f, line 3, procedure MAIN.
Subscript number 1 has value 11 in array a.
Abort (core dumped)
demo%
```

### f77: Undeclared Variable Types (-u)

The -u option checks for any undeclared variables. (Not available with £90.)

The -u option causes all variables to be initially identified as undeclared, so that all variables that are not explicitly declared by type statements, or by an <code>IMPLICIT</code> statement, are flagged with an error. The -u flag is useful for discovering mistyped variables. If -u is set, all variables are treated as undeclared until explicitly declared. Use of an undeclared variable is accompanied by an error message.

### Version Checking (-V)

The -V option causes the name and version ID of each phase of the compiler to be displayed. This option can be useful in tracking the origin of ambiguous error messages and in reporting compiler failures, and to verify the level of installed compiler patches.

# Interactive Debugging With dbx and Sun WorkShop

The Sun WorkShop provides a tightly integrated development environment for building and browsing, as well as debugging applications written in Fortran, C, and

The Sun WorkShop debugging facility is a window-based interface to dbx, while dbx itself is an interactive, line-oriented, source-level symbolic debugger. Either can be used to determine where a program crashed, to view or trace the values of variables and expressions in a running code, and to set breakpoints.

Sun WorkShop adds a sophisticated graphical environment to the debugging process that is integrated with tools for editing, building, and source code version control. It includes a data visualization capability to display and explore large and complex datasets, simulate results, and interactively steer computations.

For details, see the Sun manuals Using Sun WorkShop and Debugging a Program With Sun WorkShop, and the dbx(1) man pages.

The dbx program provides event management, process control, and data inspection. You can watch what is happening during program execution, and perform the following tasks:

- Fix one routine, then continue executing without recompiling the others
- Set watchpoints to stop or trace if a specified item changes
- Collect data for performance tuning
- Graphically monitor variables, structures, and arrays
- Set breakpoints (set places to halt in the program) at lines or in functions
- Show values—once halted, show or modify variables, arrays, structures
- Step through a program, one source or assembly line at a time
- Trace program flow—show sequence of calls taken
- Invoke procedures in the program being debugged
- Step over or into function calls; step up and out of a function call

- Run, stop, and continue execution at the next line or at some other line
- Save and then replay all or part of a debugging run
- Examine the call stack, or move up and down the call stack
- Program scripts in the embedded Korn shell
- Follow programs as they fork(2) and exec(2)

To debug optimized programs, use the dbx fix command to recompile the routines you want to debug:

- 1. Compile the program with the appropriate -0n optimization level.
- 2. Start the execution under dbx.
- 3. Use fix -g any.f without optimization on the routine you want to debug.
- 4. Use continue with that routine compiled.

Some optimizations will be inhibited by the presence of --g on the compilation command. For example, --g suppresses the automatic inlining usually obtained with -04. --g cancels any parallelization option (--autopar, --explicitpar, --parallel), as well as --depend and --reduction. Debugging is facilitated by specifying --g without any optimization options. See the dbx documentation for details.

# f 77: Viewing Compiler Listing Diagnostics

Use the error utility program to view compiler diagnostics merged with the source code. error inserts compiler diagnostics above the relevant line in the source file. The diagnostics include the standard compiler error and warning messages, but *not* the -Xlist error and warning messages.

**Note -** The error utility rewrites your source files and does not work if the source files are read-only, or are in a read only directory.

error(1) is included as part of a "developer" installation of the Solaris operating environment; it can also be installed from the package, SUNWbtool.

Facilities also exist in the Sun WorkShop for viewing compiler diagnostics. Refer to *Using Sun WorkShop*.

# Floating-Point Arithmetic

This chapter considers floating-point arithmetic and suggests strategies for avoiding and detecting numerical computation errors.

For a detailed examination of floating-point computation on SPARC and x86 processors, see the Sun *Numerical Computation Guide*.

# Introduction

Sun's floating-point environment on SPARC and x86 implements the arithmetic model specified by the IEEE Standard 754 for Binary Floating Point Arithmetic. This environment enables you to develop robust, high-performance, portable numerical applications. It also provides tools to investigate any unusual behavior by a numerical program.

In numerical programs, there are many potential sources for computational error:

- The computational model could be wrong.
- The algorithm used could be numerically unstable.
- The data could be ill-conditioned.
- The hardware could be producing unexpected results.

Finding the source of the errors in a numerical computation that has gone wrong can be extremely difficult. The chance of coding errors can be reduced by using commercially available and tested library packages whenever possible. Choice of algorithms is another critical issue. Using the appropriate computer arithmetic is another.

This chapter makes no attempt to teach or explain numerical error analysis. The material presented here is intended to introduce the IEEE floating-point model as implemented by Sun's Fortran compilers.

# **IEEE Floating-Point Arithmetic**

IEEE arithmetic is a relatively new way of dealing with arithmetic operations that result in such problems as invalid, division by zero, overflow, underflow, or inexact. The differences are in rounding, handling numbers near zero, and handling numbers near the machine maximum.

The IEEE standard supports user handling of exceptions, rounding, and precision. Consequently, the standard supports interval arithmetic and diagnosis of anomalies. IEEE Standard 754 makes it possible to standardize elementary functions like *exp* and *cos*, to create high precision arithmetic, and to couple numerical and symbolic algebraic computation.

IEEE arithmetic offers users greater control over computation than does any other kind of floating-point arithmetic. The standard simplifies the task of writing numerically sophisticated, portable programs. Many questions about floating-point arithmetic concern elementary operations on numbers. For example:

- What is the result of an operation when the infinitely precise result is not representable in the computer hardware?
- Are elementary operations like multiplication and addition commutative?

Another class of questions concerns floating-point exceptions and exception handling. What happens if you:

- Multiply two very large numbers with the same sign?
- Divide nonzero by zero?
- Divide zero by zero?

In older arithmetic models, the first class of questions might not have the expected answers, while the exceptional cases in the second class might all have the same result: the program aborts on the spot or proceeds with garbage results.

The standard ensures that operations yield the mathematically expected results with the expected properties. It also ensures that exceptional cases yield specified results, unless the user specifically makes other choices.

For example, the exceptional values +Inf, -Inf, and NaN are introduced intuitively:

Also, five types of floating-point exception are identified:

- **Invalid.** Operations with mathematically invalid operands—for example, 0.0/0.0, sqrt(-1.0), and log(-37.8)
- **Division by zero.** Divisor is zero and dividend is a finite nonzero number—for example, 9.9/0.0
- **Overflow.** Operation produces a result that exceeds the range of the exponent—for example, MAXDOUBLE+0.00000000001e308
- Underflow. Operation produces a result that is too small to be represented as a normal number—for example, MINDOUBLE \* MINDOUBLE
- **Inexact.** Operation produces a result that cannot be represented with infinite precision—for example, 2.0 / 3.0, log(1.1) and 0.1 in input

The implementation of the IEEE standard is described in the Sun *Numerical Computation Guide.* 

### -ftrap=mode Compiler Options

The <code>-ftrap=mode</code> option enables trapping for floating-point exceptions. If no signal handler has been established by an <code>ieee\_handler()</code> call, the exception terminates the program with a memory dump core file. See <code>Fortran User's Guide</code> for details on this compiler option. For example, to enable trapping for overflow, division by zero, and invalid operations, compile with <code>-ftrap=common</code>.

**Note -** You must compile the application's main program with -ftrap= for trapping to be enabled.

# Floating-Point Exceptions and Fortran

Programs compiled by £77 automatically display a list of accrued floating-point exceptions on program termination. In general, a message results if any one of the invalid, division-by-zero, or overflow exceptions have occurred. Inexact exceptions do not generate messages because they occur so frequently in real programs.

£90 programs do not automatically report on exceptions at program termination. An explicit call to *ieee\_retrospective*(3M) is required.

You can turn off any or all of these messages with ieee\_flags() by clearing exception status flags. Do this at the end of your program.

# **Handling Exceptions**

Exception handling according to the IEEE standard is the default on SPARC and x86 processors. However, there is a difference between detecting a floating-point exception and generating a signal for a floating-point exception (SIGFPE).

Following the IEEE standard, two things happen when an untrapped exception occurs during a floating-point operation:

- The system returns a default result. For example, on 0/0 (*invalid*), the system returns NaN as the result.
- A flag is set to indicate that an exception is raised. For example, 0/0 (*invalid*), the system sets the "invalid operation" flag.

# Trapping a Floating-Point Exception

£77 and £90 differ significantly in the way they handle floating-point exceptions.

With £77, the default on SPARC and x86 systems is *not* to automatically generate a signal to interrupt the running program for a floating-point exception. The assumptions are that signals could degrade performance and that most exceptions are not significant as long as expected values are returned.

The default with £90 is to automatically trap on division by zero, overflow, and invalid operation.

The f77 and f90 command-line option -ftrap can be used to change the default. In terms of -ftrap, the default for f77 is -ftrap=%none. The default for f90 is -ftrap=common.

To enable exception trapping, compile the main program with one of the -ftrap options—for example: -ftrap=common.

### SPARC: Nonstandard Arithmetic

One aspect of standard IEEE arithmetic, called *gradual underflow*, can be manually disabled. When disabled, the program is considered to be running with nonstandard arithmetic.

The IEEE standard for arithmetic specifies a way of handling underflowed results gradually by dynamically adjusting the radix point of the significand. In IEEE floating-point format, the radix point occurs before the significand, and there is an implicit leading bit of 1. Gradual underflow allows the implicit leading bit to be cleared to 0 and shifts the radix point into the significant when the result of a floating-point computation would otherwise underflow. With a SPARC processor this result is not accomplished in hardware but in software. If your program generates many underflows (perhaps a sign of a problem with your algorithm) and you run on a SPARC processor, you may experience a performance loss.

Gradual underflow can be disabled either by compiling with the -fns option or by calling the library routine nonstandard arithmetic() from within the program to turn it off. Call standard\_arithmetic() to turn gradual underflow back on.

**Note -** To be effective, the application's main program must be compiled with -fns. See the Fortran User's Guide.

For legacy applications, take note that:

- The standard\_arithmetic() subroutine replaces an earlier routine named gradual\_underflow().
- The nonstandard\_arithmetic() subroutine replaces an earlier routine named abrupt\_underflow().

Note - The -fns option and the nonstandard\_arithmetic() library routine are effective only on some SPARC systems. On x86 platforms, gradual underflow is performed by the hardware.

# **IEEE Routines**

The following interfaces help people use IEEE arithmetic. These are mostly in the math library libsunmath and in several .h files.

- ieee\_flags(3m)—Controls rounding direction and rounding precision; query exception status; clear exception status
- ieee\_handler(3m)—Establishes an exception handler routine
- ieee\_functions(3m)—Lists name and purpose of each IEEE function
- ieee\_values(3m)—Lists functions that return special values
- Other libm functions described in this section:
  - ieee\_retrospective
  - nonstandard\_arithmetic
  - standard arithmetic

The SPARC processors conform to the IEEE standard in a combination of hardware and software support for different aspects, x86 processors conform to the IEEE standard entirely through hardware support.

The newest SPARC processors contain floating-point units with integer multiply and divide instructions and hardware square root.

Best performance is obtained when the compiled code properly matches the runtime floating-point hardware. The compiler's -xtarget= option permits specification of

the runtime hardware. For example, -xtarget=ultra would inform the compiler to generate object code that will perform best on an UltraSPARC processor.

On SPARC platforms: The utility fpversion displays which floating-point hardware is installed and indicates the appropriate -xtarget value to specify. This utility runs on all Sun SPARC architectures. See *fpversion*(1), the Sun *Fortran User's Guide* (regarding -xtarget) and the *Numerical Computation Guide* for details.

# Flags and ieee\_flags()

The ieee\_flags function is used to query and clear exception status flags. It is part of the libsunmath library shipped with Sun compilers and performs the following tasks:

- Controls rounding direction and rounding precision
- Checks the status of the exception flags
- Clears exception status flags

The general form of a call to ieee\_flags is:

```
flags = ieee_flags( action, mode, in, out )
```

Each of the four arguments is a string. The input is *action, mode*, and *in*. The output is *out* and *flags*. ieee\_flags is an integer-valued function. Useful information is returned in *flags* as a set of 1-bit flags. Refer to the man page for *ieee\_flags*(3m) for complete details.

Possible parameter values are shown in the following table:

 TABLE 6-1
 ieee\_flags(action, mode, in, out)Argument Values

action	mode	in, out
get	direction	nearest
set	precision	tozero
clear	exception	negative
clearall		positive
		extended
		double
		single
		inexact
		division
		underflow
		overflow
		invalid
		all
		common

The precision *mode* is available only on x86 platforms.

Note that these are literal character strings, and the output parameter out must be at least CHARACTER\*9. The meanings of the possible values for in and out depend on the action and mode they are used with. These are summarized in the following table:

 TABLE 6-2 ieee\_flags Argument Meanings

Value of in and out	Refers to
nearest, tozero, negative, positive	Rounding direction
extended, double, single	Rounding precision

TABLE 6-2 ieee\_flags Argument Meanings (continued)

Value of in and out	Refers to
<pre>inexact, division, underflow, overflow, invalid</pre>	Exceptions
all	All five exceptions
common	Common exceptions: invalid, division, overflow

For example, to determine what is the highest priority exception that has a flag raised, pass the input argument *in* as the null string:

```
CHARACTER *9, out
ieeer = ieee_flags( "get", "exception", "", out )
PRINT *, out, " flag raised"
```

Also, to determine if the overflow exception flag is raised, set the input argument in to overflow. On return, if out equals overflow, then the overflow exception flag is raised; otherwise it is not raised.

```
ieeer = ieee_flags( "get", "exception", "overflow", out )
IF ( out.eq. "overflow") PRINT *,"overflow flag raised"
```

#### Example: Clear the invalid exception:

```
ieeer = ieee_flags( "clear", "exception", "invalid", out )
```

#### Example: Clear all exceptions:

```
ieeer = ieee_flags( "clear", "exception", "all", out )
```

#### Example: Set rounding direction to zero:

```
ieeer = ieee_flags( "set", "direction", "tozero", out )
```

#### Example: Set rounding precision to double:

```
ieeer = ieee_flags( "set", "precision", "double", out )
```

#### Turning Off All Warning Messages With ieee\_flags

Calling ieee\_flags with an *action* of clear, as shown in the following example, resets any uncleared exceptions. Put this call before the program exits to suppress system warning messages about floating-point exceptions at program termination.

```
Example: Clear all accrued exceptions with ieee_flags():
    i = ieee_flags("clear", "exception", "all", out)
```

### Detecting an Exception With ieee\_flags

The following example demonstrates how to determine which floating-point exceptions have been raised by earlier computations. Bit masks defined in the system include file f77\_floatingpoint.h are applied to the value returned by ieee\_flags.

In this example, <code>DetExcFlg.F</code>, the include file is introduced using the <code>#include</code> preprocessor directive, which requires us to name the source file with a <code>.F</code> suffix. Underflow is caused by dividing the smallest double-precision number by 2.

Example: Detect an exception using ieee\_flags and decode it:

```
#include "f77_floatingpoint.h"
        CHARACTER*16 out
        DOUBLE PRECISION d_max_subnormal, x
        INTEGER div, flgs, inv, inx, over, under

    Cause underflow

      x = d_{max\_subnormal()} / 2.0
      flgs=ieee_flags("get","exception","",out) ! Which are raised?
      inx = and(rshift(flgs, fp_inexact) , 1) ! Decode
      div = and(rshift(flgs, fp_division) , 1) ! the value
                                                   ! returned
      under = and(rshift(flgs, fp_underflow), 1)
                                                    ! by
      over = and(rshift(flgs, fp_overflow) , 1)
       inv = and(rshift(flgs, fp_invalid) , 1)
                                                       ! ieee_flags
      PRINT *, "Highest priority exception is: ", out
      PRINT *, " invalid divide overflo underflo inexact"
      PRINT "(5i8)", inv, div, over, under, inx
       PRINT *, "(1 = exception is raised; 0 = it is not)"
       i = ieee_flags("clear", "exception", "all", out) ! Clear all
      END
```

Example: Compile and run the preceding example (DetExcFlg.F):

```
demo% f77 -silent DetExcFlg.F
demo% a.out
```

### **IEEE Extreme Value Functions**

The compilers provide a set of functions that can be called to return a special IEEE extreme value. These values, such as *infinity* or *minimum normal*, can be used directly in an application program.

Example: A convergence test based on the smallest number supported by the hardware would look like:

```
IF ( delta .LE. r_min_normal() ) RETURN
```

The values available are listed in the following table:

TABLE 6-3 Functions Returning IEEE Values

IEEE Value	Double Precision	Single Precision
infinity	d_infinity()	r_infinity()
quiet NaN	d_quiet_nan()	r_quiet_nan()
signaling NaN	d_signaling_nan()	r_signaling_nan()
min normal	d_min_normal()	r_min_normal()
min subnormal	d_min_subnormal()	r_min_subnormal()
max subnormal	d_max_subnormal()	r_max_subnormal()
max normal	d_max_normal()	r_max_normal()

The two NaN values (quiet and signaling) are *unordered* and should not be used in comparisons such as IF(X.ne.r\_quiet\_nan())THEN... To determine whether some value is a NaN, use the function ir\_isnan(r) or id\_isnan(d).

The Fortran names for these functions are listed in these man pages:

- libm\_double(3f)
- libm\_single(3f)
- ieee\_functions(3m)

#### Also see:

- ieee\_values(3m)
- The f77\_floatingpoint.h header file

# Exception Handlers and ieee\_handler()

Typical concerns about IEEE exceptions are:

- What happens when an exception occurs?
- How do I use ieee\_handler() to establish a user function as an exception handler?
- How do I write a function that can be used as an exception handler?
- How do I locate the exception—where did it occur?

Exception trapping to a user routine begins with the system generating a signal on a floating-point exception. The standard UNIX name for *signal: floating-point exception* is SIGFPE. The default situation on SPARC and x86 platforms is *not* to generate a SIGFPE when an exception occurs. For the system to generate a SIGFPE, exception trapping must first be enabled, usually by a call to ieee\_handler().

### **Establishing an Exception Handler Function**

To establish a function as an exception handler, pass the name of the function to ieee\_handler(), together with the name of the exception to watch for and the action to take. Once you establish a handler, a SIGFPE signal is generated whenever the particular floating-point exception occurs, and the specified function is called.

The form for invoking ieee\_handler() is shown in the following table:

TABLE 6-4 Arguments for ieee\_handler(action, exception, handler)

Argument	Туре	Possible Values
action	character	get, set, or clear
exception	character	invalid, division, overflow, underflow, or inexact
handler	Function name	The name of the user handler function or SIGFPE_DEFAULT, SIGFPE_IGNORE, or SIGFPE_ABORT
Return value	integer	0 =OK

The routine that calls ieee\_handler() should also declare:

#include 'f77\_floatingpoint.h'

For £90 programs, declare:

#include 'f90/floatingpoint.h'

The special arguments SIGFPE\_DEFAULT, SIGFPE\_IGNORE, and SIGFPE\_ABORT are defined in f77\_floatingpoint.h and can be used to change the behavior of the program for a specific exception:

SIGFPE_DEFAULT or SIGFPE_IGNORE	No action taken when the specified exception occurs.
SIGFPE_ABORT	Program aborts, possibly with dump file, on exception.

### Writing User Exception Handler Functions

The actions your exception handler takes are up to you. However, the routine must be an integer function with three arguments specified as shown:

handler\_name( sig, sip, uap)

- *handler\_name* is the name of the integer function.
- sig is an integer.
- sip is a record that has the structure siginfo.
- uap is not used.

#### Example: An exception handler function:

```
INTEGER FUNCTION hand( sig, sip, uap )
INTEGER sig, location
STRUCTURE /fault/
    INTEGER address
    INTEGER trapno
END STRUCTURE
STRUCTURE /siginfo/
    INTEGER si_signo
    INTEGER si_code
     INTEGER si_errno
    RECORD /fault/ fault
END STRUCTURE
RECORD /siginfo/ sip
location = sip.fault.address
... actions you take ...
```

This £77 example would have to be modified to run on SPARC V9 architectures (-xarch=v9 or v9a) by replacing all INTEGER declarations within each STRUCTURE with INTEGER\*8.

If the handler routine enabled by ieee\_handler() is in Fortran as shown in the example, the routine should not make any reference to its first argument (sig). This first argument is passed by value to the routine and can only be referenced as loc(sig). The value is the signal number.

#### Detecting an Exception by Handler

The following examples show how to create handler routines to detect floating-point exceptions.

Example: Detect exception and abort:

```
demo% cat DetExcHan.f
     EXTERNAL myhandler
      REAL r / 14.2 /, s / 0.0 /
      i = ieee_handler ("set", "division", myhandler )
      t = r/s
      END
      INTEGER FUNCTION myhandler(sig,code,context)
      INTEGER sig, code, context(5)
      CALL abort()
      END
demo% f77 -silent DetExcHan.f
demo% a.out
abort: called
Abort (core dumped)
demo%
```

SIGFPE is generated whenever that floating-point exception occurs. When the SIGFPE is detected, control passes to the myhandler function, which immediately aborts. Compile with -g and use dbx to find the location of the exception.

#### Locating an Exception by Handler

Example: Locate an exception (print address) and abort:

```
demo% cat LocExcHan.F
#include "f77_floatingpoint.h"
     EXTERNAL Exhandler
     INTEGER Exhandler, i, ieee_handler
     REAL r / 14.2 /, s / 0.0 /, t
C Detect division by zero
     i = ieee_handler( "set", "division", Exhandler )
     t = r/s
     END
     INTEGER FUNCTION Exhandler( sig, sip, uap)
      INTEGER sig
      STRUCTURE /fault/
                 INTEGER address
      END STRUCTURE
      STRUCTURE /siginfo/
                  INTEGER si_signo
                 INTEGER si code
                 INTEGER si_errno
                 RECORD /fault/ fault
     END STRUCTURE
     RECORD /siginfo/ sip
     WRITE (*,10) sip.si_signo, sip.si_code, sip.fault.address
         FORMAT("Signal ",i4," code ",i4," at hex address ", Z8 )
10
     CALL abort()
     END
demo% f77 -silent -g LocExcHan.F
demo% a.out
Signal 8 code
                3 at hex address 11230
abort: called
Abort (core dumped)
demo%
```

In SPARC V9 environments, replace the INTEGER declarations within each STRUCTURE with INTEGER\*8, and the i4 formats with i8.

In most cases, knowing the actual *address* of the exception is of little use, except with dbx:

```
demo% dbx a.out
(dbx) stopi at 0x11230 Set breakpoint at address
(2) stopi at &MAIN+0x68
(dbx) run Running: a.out
(process id 18803)
stopped in MAIN at 0x11230
MAIN+0x68: fdivs %f3, %f2, %f2
(dbx) where Shows the line number of the exception
=>[1] MAIN(), line 7 in "LocExcHan.F"
```

Of course, there are easier ways to determine the source line that caused the error. However, this example does serve to show the basics of exception handling.

#### **Disabling All Signal Handlers**

With £77, some system signal handlers for trapping interrupts, bus errors, segmentation violations, or illegal instructions are automatically enabled by default.

Although generally you would not want to turn off this default behavior, you can do so by compiling a C program that sets the global C variable f77\_no\_handlers to 1 and linking into your executable program:

```
demo% cat NoHandlers.c
    int f77_no_handlers=1;
demo% cc -c NoHandlers.c
demo% f77 NoHandlers.o MyProgram.f
```

Otherwise, by default, f77\_no\_handlers is 0. The setting takes effect just before execution is transferred to the user program.

This variable is in the global name space of the program; do not use f77 no handlers as the name of a variable anywhere else in the program.

With £90, no signal handlers are on by default.

### **Retrospective Summary**

The ieee\_retrospective function queries the floating-point status registers to find out which exceptions have accrued and a message is printed to standard error to inform you which exceptions were raised but not cleared. This function is automatically called by Fortran 77 programs at normal program termination (CALL EXIT). The message typically looks like this; the format may vary with each compiler release:

```
Note: IEEE floating-point exception flags raised:
    Division by Zero;
IEEE floating-point exception traps enabled:
    inexact; underflow; overflow; invalid operation;
See the Numerical Computation Guide, ieee_flags(3M),
    ieee_handler(3M)
```

Fortran 90 programs do not call ieee\_retrospective automatically. A Fortran 90 program would need to call ieee\_retrospective explicitly (and link with -lf77compat).

# **Debugging IEEE Exceptions**

In most cases, the only indication that any floating-point exceptions (such as overflow, underflow, or invalid operation) have occurred is the retrospective summary message at program termination. Locating *where* the exception occurred requires exception trapping be enabled. This can be done by either compiling with the <code>-ftrap=common</code> option or by establishing an exception handler routine with <code>ieee\_handler()</code>. With exception trapping enabled, run the program from <code>dbx</code> or the Sun WorkShop, using the <code>dbx catch FPE</code> command to see where the error occurs.

The advantage of recompiling with -ftrap=common is that the source code need not be modified to trap the exceptions. However, by calling  $ieee\_handler()$  you can be more selective as to which exceptions to look at.

Example: Recompiling with -ftrap=common and using dbx:

```
demo% f77 -g -ftrap=common -silent myprogram.f
demo% dbx a.out
Reading symbolic information for a.out
Reading symbolic information for rtld /usr/lib/ld.so.1
Reading symbolic information for libF77.so.3
Reading symbolic information for libc.so.1
Reading symbolic information for libdl.so.1
(dbx) catch FPE
(dbx) run
Running: a.out
(process id 19739)
signal FPE (floating point divide by zero) in MAIN at line 212 in file "myprogram.f"
  212
                   Z = X/Y
(dbx) print Y
y = 0.0
(dbx)
```

If you find that the program terminates with overflow and other exceptions, you can locate the first overflow specifically by calling <code>ieee\_handler()</code> to trap just

overflows. This requires modifying the source code of at least the main program, as shown in the following example.

Example: Locate an overflow when other exceptions occur:

```
demo% cat myprog.F
#include ``f77_floatingpoint.h''
       program myprogram
      ier = ieee_handler('set','overflow',SIGFPE_ABORT)
demo% f77 -g -silent myprog.F
demo% dbx a.out
Reading symbolic information for a.out
Reading symbolic information for rtld /usr/lib/ld.so.1
Reading symbolic information for libF77.so.3
Reading symbolic information for libc.so.1
Reading symbolic information for libdl.so.1
(dbx) catch FPE
(dbx) run
Running: a.out
(process id 19793)
signal FPE (floating point overflow) in MAIN at line 55 in file "myprog.F"
                    w = rmax * 200.
                                                         ! Cause of the overflow
                                             ! Continue execution to completion
(dbx) cont
 Note: IEEE floating-point exception flags raised:
   Inexact; Division by Zero; Underflow; ! There were other exceptions
 IEEE floating-point exception traps enabled:
   overflow;
 See the Numerical Computation Guide...
execution completed, exit code is 0
(dbx)
```

To be selective, the example introduces the <code>#include</code>, which required renaming the source file with a <code>.F</code> suffix and calling <code>ieee\_handler()</code>. You could go further and create your own handler function to be invoked on the overflow exception to do some application-specific analysis and print intermediary or debug results before aborting.

# **Further Numerical Adventures**

This section addresses some real world problems that involve arithmetic operations that may unwittingly generate invalid, division by zero, overflow, underflow, or inexact exceptions.

For instance, prior to the IEEE standard, if you multiplied two very small numbers on a computer, you could get zero. Most mainframes and minicomputers behaved that way. With IEEE arithmetic, *gradual underflow* expands the dynamic range of computations.

For example, consider a machine with 1.0E-38 as the machine's *epsilon*, the smallest representable value on the machine. Multiply two small numbers:

```
a = 1.0E-30

b = 1.0E-15

x = a * b
```

In older arithmetic, you would get 0.0, but with IEEE arithmetic and the same word length, you get 1.40130E-45. Underflow tells you that you have an answer smaller than the machine naturally represents. This result is accomplished by "stealing" some bits from the mantissa and shifting them over to the exponent. The result, a *denormalized number*, is less precise in some sense, but more precise in another. The deep implications are beyond this discussion. If you are interested, consult *Computer*, January 1980, Volume 13, Number 1, particularly J. Coonen's article, "Underflow and the Denormalized Numbers."

Most scientific programs have sections of code that are sensitive to roundoff, often in an equation solution or matrix factorization. Without gradual underflow, programmers are left to implement their own methods of detecting the approach of an inaccuracy threshold. Otherwise they must abandon the quest for a robust, stable implementation of their algorithm.

For more details on these topics, see the Sun Numerical Computation Guide.

# **Avoiding Simple Underflow**

Some applications actually do a lot of computation very near zero. This is common in algorithms computing residuals or differential corrections. For maximum numerically safe performance, perform the key computations in extended precision arithmetic. If the application is a single-precision application, you can perform key computations in double precision.

Example: A simple dot product computation in single precision:

If a(i) and b(i) are very small, many underflows occur. By forcing the computation to double precision, you compute the dot product with greater accuracy and do not suffer underflows:

On SPARC platforms: You can force a SPARC processor to behave like an older system with respect to underflow (Store Zero) by adding a call to the library routine nonstandard\_arithmetic() or by compiling the application's main program with the -fns option.

### Continuing With the Wrong Answer

You might wonder why you would continue a computation if the answer is clearly wrong. IEEE arithmetic allows you to make distinctions about what kind of wrong answers can be ignored, such as NaN or Inf. Then decisions can be made based on such distinctions.

For an example, consider a circuit simulation. The only variable of interest (for the sake of argument) from a particular 50-line computation is the voltage. Further, assume that the only values that are possible are +5v, 0, -5v.

It is possible to carefully arrange each part of the calculation to coerce each sub-result to the correct range:

```
if computed value is greater than 4.0, return 5.0 if computed value is between -4.0 and +4.0, return 0 if computed value is less than -4.0, return -5.0
```

Furthermore, since Inf is not an allowed value, you need special logic to ensure that big numbers are not multiplied.

IEEE arithmetic allows the logic to be much simpler. The computation can be written in the obvious fashion, and only the final result need be coerced to the correct value—since Inf can occur and can be easily tested.

Furthermore, the special case of 0/0 can be detected and dealt with as you wish. The result is easier to read and faster in executing, since you don't do unneeded comparisons.

### SPARC: Excessive Underflow

If two very small numbers are multiplied, the result underflows.

If you know in advance that the operands in a multiplication (or subtraction) may be small and underflow is likely, run the calculation in double precision and convert the result to single precision later.

For example, a dot product loop like this:

```
real sum, a(maxn), b(maxn)
...
```

```
do i =1, n
    sum = sum + a(i)*b(i)
enddo
```

where the a(\*) and b(\*) are known to have small elements, should be run in double precision to preserve numeric accuracy:

```
real a(maxn), b(maxn)
double sum
...
do i =1, n
    sum = sum + a(i)*dble(b(i))
enddo
```

Doing so may also improve performance due to the software resolution of excessive underflows caused by the original loop. However, there is no hard and fast rule here; experiment with your intensely computational code to determine the most profitable solutions.

# **Porting**

This chapter discusses the porting of programs from other dialects of Fortran to Sun compilers. VAX VMS Fortran programs compile almost exactly as is with Sun £77; this is discussed further in the chapter on VMS extensions in the *FORTRAN 77 Language Reference Manual*.

**Note -** Porting issues bear mostly upon FORTRAN 77 programs. The Sun Fortran 90 compiler, £90, incorporates few nonstandard extensions, and these are described in the *Fortran User's Guide*.

# Time and Date Functions

Library functions that return the time of day or elapsed CPU time vary from system to system.

The following time functions are not supported directly in the Sun Fortran libraries, but you can write subroutines to duplicate their functions:

- Time-of-day in 10h format
- Date in A10 format
- Milliseconds of job CPU time
- Julian date in ASCII

The time functions supported in the Sun Fortran library are listed in the following table:

TABLE 7–1 Sun Fortran Time Functions

Name	Function	Man Page
time	Returns the number of seconds elapsed since January, 1, 1970	time(3F)
date	Returns date as a character string	date(3F)
fdate	Returns the current time and date as a character string	fdate(3F)
idate	Returns the current month, day, and year in an integer array	idate(3F)
itime	Returns the current hour, minute, and second in an integer array	itime(3F)
ctime	Converts the time returned by the time function to a character string	ctime(3F)
ltime	Converts the time returned by the ${\tt time}$ function to the local time	ltime(3F)
gmtime	Converts the time returned by the ${\tt time}$ function to Greenwich time	gmtime(3F)
etime	Single processor: Returns elapsed user and system time for program execution Multiple processors: Returns the wall clock time	etime(3F)
dtime	Returns the elapsed user and system time since last call to $\ensuremath{\operatorname{dtime}}$	dtime(3F)
date_and_time	Returns date and time in character and numeric form	date_and_time(3F)

For details, see Fortran Library Reference Manual or the individual man pages for these functions. The routines listed in the following table provide compatibility with VMS Fortran system routines idate and time. To use these routines, you must include the -1V77 option on the f77 command line, in which case you also get these VMS versions instead of the standard f77 versions.

TABLE 7-2 Summary: Nonstandard VMS Fortran System Routines

Name	Definition	Calling Sequence	Argument Type
idate	Date as day, month, year	call idate( d, m, y )	integer
time	Current time as hhmmss	call time( t )	character*8

**Note** - The *date*(3F) routine and the VMS version of *idate*(3F) cannot be Year 2000 safe because they return 2-digit values for the year. Programs that compute time duration by subtracting dates returned by these routines will compute erroneous results after December 31, 1999. The Fortran 90 routine *date\_and\_time*(3F) is available for both FORTRAN 77 and Fortran 90 programs, and should be used instead. See the *Fortran Library Reference Manual* for details.

The error condition subroutine errsns is *not* provided, because it is totally specific to the VMS operating system.

Here is a simple example of the use of these time functions (TestTim.f):

```
subroutine startclock
      common / myclock / mytime
      integer mytime, time
      mytime = time()
      return
      end
      function wallclock
      integer wallclock
      common / myclock / mytime
      integer mytime, time, newtime
      newtime = time()
      wallclock = newtime -- mytime
      mytime = newtime
     return
      end
      integer wallclock, elapsed
      character*24 greeting
     real dtime, timediff, timearray(2)
      print a heading
С
      call fdate( greeting )
                   Hello, Time Now Is: ", greeting
     print*, "
                  "See how long "sleep 4" takes, in seconds"
      print*,
      call startclock
      call system( "sleep 4" )
      elapsed = wallclock()
     print*, "Elapsed time for sleep 4 was: ", elapsed," seconds"
      now test the cpu time for some trivial computing
      timediff = dtime( timearray )
      q = 0.01
      do 30 i = 1, 1000
```

Running this program produces the following results:

# **Formats**

Some £77 and £90 format edit descriptors can behave differently on other systems. Here are some format specifiers that £77 treats differently than some other implementations:

- A Alphanumeric conversion. Used with character type data elements. In FORTRAN 77, this specifier worked with any variable type. £77 supports the older usage, up to four characters to a word.
- \$ Suppresses newline character output.
- R Sets an arbitrary radix for the I formats that follow in the descriptor.
- SU Selects unsigned output for following I formats. For example, you can convert output to either hexadecimal or octal with the following formats, instead of using the Z or O edit descriptors:

```
10 FORMAT( SU, 16R, I4 )
20 FORMAT( SU, 8R, I4 )
```

# Carriage-Control

Fortran carriage-control grew out of the capabilities of the equipment used when Fortran was originally developed. For similar historical reasons, an operating system derived from the UNIX operating system, does not have Fortran carriage control, but you can simulate it in two ways.

- Use the asa filter to transform Fortran carriage-control conventions into the UNIX carriage-control format (see the *asa* (1) man page) before printing files with the lpr command.
- f77: For simple jobs, use OPEN(N, FORM='PRINT') to enable single or double spacing, formfeed, and stripping off of column one. It is legal to reopen unit 6 to change the form parameter to PRINT. For example:

```
OPEN( 6, FORM="PRINT")
```

You can use lp(1) to print a file that is opened in this manner.

# Working With Files

Early Fortran systems did not use named files, but did provide a command line mechanism to equate actual file names with internal unit numbers. This facility can be emulated in a number of ways, including standard UNIX redirection.

Example: Redirecting stdin to redir.data (using csh(1)):

# **Porting From Scientific Mainframes**

If the application code was originally developed for 64-bit (or 60-bit) mainframes such as CRAY or CDC, you might want to compile these codes with the

```
f77 -xtypemap option: "-xtypemap=real:64,double:128,integer:64"
```

to preserve the expected precision of the original. This option automatically promotes all default REAL variables to REAL\*8, default DOUBLE to REAL\*16, and COMPLEX to COMPLEX\*16. Only undeclared variables or variables declared as simply REAL or DOUBLE or DOUBLE PRECISION or COMPLEX are promoted; variables declared explicitly (for example, REAL\*4) are not promoted.

On x86, or for better performance on SPARC, use: "-xtypemap=real:64,double:64,integer:64"

which does not promote default DOUBLE PRECISION.

The -xtypemap option, is preferred over -dbl and -r8 and -i2. See the *Fortran User's Guide* and the f77(1) man pages for details.

To further recreate the original mainframe environment, it is probably preferable to stop on overflows, division by zero, and invalid operations. Compile the main program with <code>-ftrap=common</code> to ensure this.

# **Data Representation**

The FORTRAN 77 Language Reference Manual, Fortran User's Guide, and the Sun Numerical Computation Guide discuss in detail the hardware representation of data objects in Fortran. Differences between data representations across systems and hardware platforms usually generate the most significant portability problems.

The following issues should be noted:

- Sun adheres to the IEEE Standard 754 for floating-point arithmetic. Therefore, the first four bytes in a REAL\*8 are not the same as in a REAL\*4.
- The default sizes for reals, integers, and logicals are described in the FORTRAN 77 standard, except when these default sizes are changed by the -xtypemap= option (or by -i2, -dbl, or -r8).
- Character variables can be freely mixed and equivalenced to variables of other types, but be careful of potential alignment problems.
- £77 IEEE floating-point arithmetic does raise exceptions on overflow or divide by zero but does not signal SIGFPE or trap by default. It does deliver IEEE indeterminate forms in cases where exceptions would otherwise be signaled. This is explained in the *Floating Point Arithmetic* chapter of this Guide.
- The extreme finite, normalized values can be determined. See *libm\_single(3F)* and *libm\_double(3F)*. The indeterminate forms can be written and read, using formatted and list-directed I/O statements.

# Hollerith Data

Many "dusty-deck" Fortran applications store Hollerith ASCII data into numerical data objects. With the 1977 Fortran standard (and Fortran 90), the CHARACTER data type was provided for this purpose and its use is recommended. You can still initialize variables with the older Fortran Hollerith (nH) feature, but this is not standard practice. The following table indicates the maximum number of characters that will fit into certain data types. (In this table, boldfaced data types indicate default types subject to promotion by the f77 command-line flags <code>-dbl</code>, <code>-r8</code>, or <code>-xtypemap=</code>).

TABLE 7-3 f 77: Maximum Characters in Data Types

	Maximum Number of Standard ASCII Characters				
Data Type	No -i2, -i4, -r8, -dbl	-i2	-i4	-r8	-dbl
BYTE	1	1	1	1	1
COMPLEX	8	8	8	16	16
COMPLEX*16	16	16	16	16	16
COMPLEX*32	32	32	32	32	32
DOUBLE COMPLEX	16	16	16	32	32
DOUBLE PRECISION	8	8	8	16	16
INTEGER	4	2	4	4	8
INTEGER*2	2	2	2	2	2
INTEGER*4	4	4	4	4	4
INTEGER*8	8	8	8	8	8
LOGICAL	4	2	4	4	8
LOGICAL*1	1	1	1	1	1
LOGICAL*2	2	2	2	2	2

TABLE 7-3 f 77: Maximum Characters in Data Types (continued)

	Maximum Number of Standard ASCII Characters				
Data Type	No -i2, -i4, -r8, -dbl	-i2	-i4	-r8	-dbl
LOGICAL*4	4	4	4	4	4
LOGICAL*8	8	8	8	8	8
REAL	4	4	4	8	8
REAL*4	4	4	4	4	4
REAL*8	8	8	8	8	8
REAL*16	16	16	16	16	16

When storing standard ASCII characters with normal Fortran:

- With -r8, unspecified size INTEGER and LOGICAL do not hold double.
- With -dbl, unspecified size INTEGER and LOGICAL *do* hold double.

The storage is allocated with both options, but it is unavailable in normal Fortran with -r8.

Options -i2, -r8, and -dbl are now considered obsolete; use -xtypemap instead.

Example: Initialize variables with Hollerith:

```
demo% cat FourA8.f
     double complex x(2)
     data x /16Habcdefghijklmnop, 16Hqrstuvwxyz012345/
     write( 6, "(4A8, "!")" ) x
     end

demo% f77 -silent -o FourA8 FourA8.f
demo% FourA8
abcdefghijklmnopqrstuvwxyz012345!
demo%
```

If you pass Hollerith constants as arguments, or if you use them in expressions or comparisons, they are interpreted as character-type expressions.

If needed, you can initialize a data item of a compatible type with a Hollerith and then pass it to other routines.

#### Example:

```
program respond
        integer yes, no
        integer ask
        data yes, no / 3hyes, 2hno /
        if ( ask() .eq. yes ) then
            print *, "You may proceed!"
            print *, "Request Rejected!"
        end
        integer function ask()
        double precision solaris, response
        integer yes, no
        data yes, no / 3hyes, 2hno / data solaris/ 7hSOLARIS/
10
        format( "What system? ", $ )
20
        format( a8 )
        write( 6, 10 )
        read (5, 20) response
        ask = no
        if ( response .eq. solaris ) ask = yes
```

# **Nonstandard Coding Practices**

As a general rule, porting an application program from one system and compiler to another can be made easier by eliminating any nonstandard coding. Optimizations or work-arounds that were successful on one system might only obscure and confuse compilers on other systems. In particular, optimized hand-tuning for one particular architecture can cause degradations in performance elsewhere. This is discussed later in the chapters on performance and tuning. However, the following issues are worth considering with regards to porting in general.

### Uninitialized Variables

Some systems automatically initialize local and COMMON variables to zero or some "not-a-number" (NaN) value. However, there is no standard practice, and programs should not make assumptions regarding the initial value of any variable. To assure maximum portability, a program should initialize all variables.

### **Aliasing Across Calls**

Aliasing occurs when the same storage address is referenced by more than one name. This happens when actual arguments to a subprogram overlap between themselves or between COMMON variables within the subprogram. For example, arguments X and Z refer to the same storage locations, as do B and H:

```
COMMON /INS/B(100)
REAL S(100), T(100)
...
CALL SUB(S,T,S,B,100)
...
SUBROUTINE SUB(X,Y,Z,H,N)
REAL X(N),Y(N),Z(N),H(N)
COMMON /INS/B(100)
```

Avoid aliasing in this manner in all portable code. The results on some systems and with higher optimization levels could be unpredictable.

# **Obscure Optimizations**

Legacy codes may contain source-code restructurings of ordinary computational DO loops intended to cause older vectorizing compilers to generate optimal code for a particular architecture. In most cases, these restructurings are no longer needed and may degrade the portability of a program. Two common restructurings are strip-mining and loop unrolling.

#### Strip-Mining

Fixed-length vector registers on some architectures led programmers to manually "strip-mine" the array computations in a loop into segments:

```
REAL TX(0:63)
...

DO IOUTER = 1,NX,64

DO IINNER = 0,63

TX(IINNER) = AX(IOUTER+IINNER) * BX(IOUTER+IINNER)/2.

QX(IOUTER+IINNER) = TX(IINNER)**2

END DO

END DO
```

Strip-mining is no longer appropriate with modern compilers; the loop can be written much less obscurely as:

```
DO IX = 1,N

TX = AX(I)*BX(I)/2.

QX(I) = TX**2
```

#### Loop Unrolling

Unrolling loops by hand was a typical source-code optimization technique before compilers were available that could perform this restructuring automatically. A loop written as:

```
K = 1, N-5, 6

J = 1, N
DO
  DO
     DO I = 1,N
        A(I,J) = A(I,J) + B(I,K) * C(K,J)
                         + B(I,K+1) * C(K+1,J)
                         + B(I,K+2) * C(K+2,J)
                         + B(I,K+3) * C(K+3,J)
                         + B(I,K+4) * C(K+4,J)
                         + B(I,K+5) * C(K+5,J)
     END DO
  END DO
END DO
        KK = K, N
DO
       J = 1, N
  DO
     DO I =1,N
        A(I,J) = A(I,J) + B(I,KK) * C(KK,J)
     END DO
   END DO
END DO
```

should be rewritten the way it was originally intended:

# **Troubleshooting**

Here are a few suggestions for what to try when programs ported to Sun Fortran do not run as expected.

# Results Are Close, but Not Close Enough

Try the following:

■ Pay attention to the size and the engineering units. Numbers very close to zero can appear to be different, but the difference is not significant, especially if this number is the difference between two large numbers, such as the distance across the continent in feet, as calculated on two different computers. For example, 1.9999999e-30 is very near -9.9992112e-33, even though they differ in sign.

VAX math is not as good as IEEE math, and even different IEEE processors may differ. This is especially true if the mathematics involves many trigonometric functions. These functions are much more complicated than one might think, and the standard defines only the basic arithmetic functions. There can be subtle differences, even between IEEE machines. Review the *Floating-Point Arithmetic* chapter in this Guide.

- Try running with a call nonstandard\_arithmetic(). Doing so can also improve performance considerably, and make your Sun workstation behave more like a VAX system. If you have access to a VAX or some other system, run it there also. It is quite common for many numerical applications to produce slightly different results on each floating-point implementation.
- Check for NaN, +Inf, and other signs of probable errors. See the *Floating-Point Arithmetic* chapter in this Guide, or the man page *ieee\_handler*(3m) for instructions on how to trap the various exceptions. On most machines, these exceptions simply abort the run.
- Two numbers can differ by  $6 \times 10^{29}$  and still have the same floating-point form. Here is an example of different numbers, with the same representation:

```
real*4 x,y
    x=9999990e+29
    y=9999996e+29
    write (*,10), x, x

0 format("99,999,990 x 10^29 = ", e14.8, " = ", z8)
    write(*,20) y, y

format("99,999,996 x 10^29 = ", e14.8, " = ", z8)
    end
```

#### The output is:

```
99,999,990 x 10^29 = 0.99999993E+37 = 7cf0bdc1
99,999,996 x 10^29 = 0.99999993E+37 = 7cf0bdc1
```

In this example, the difference is  $6 \times 10^{29}$ . The reason for this indistinguishable, wide gap is that in IEEE single-precision arithmetic, you are guaranteed only six decimal digits for any one decimal-to-binary conversion. You may be able to convert seven or eight digits correctly, but it depends on the number.

# **Program Fails Without Warning**

If the program fails without warning and runs different lengths of time between failures, then:

- Compile with minimal optimization (-O1). If the program then works, compile only selective routines with higher optimization levels.
- Understand that optimizers must make assumptions about the program. Nonstandard coding or constructs can cause problems. Almost no optimizer handles all programs at all levels of optimization.

# **Performance Profiling**

This chapter describes how to measure and display program performance. Knowing where a program is spending most of its compute cycles and how efficiently it uses system resources is a prerequisite for performance tuning.

# The time Command

The simplest way to gather basic data about program performance and resource utilization is to use the time (1) command or, in csh, the set time command.

Running the program with the time command prints a line of timing information on program termination.

```
demo% time myprog
    The Answer is: 543.01
6.5u 17.1s 1:16 31% 11+21k 354+210io 135pf+0w
demo%
```

#### The interpretation is:

user system wallclock resources memory I/O paging

- *user* 6.5 seconds in user code, approximately
- *system* 17.1 seconds in system code for this task, approximately
- *wallclock* 1 minute 16 seconds to complete
- resources 31% of system resources dedicated to this program
- memory 11 kilobytes of shared program memory, 21 kilobytes of private data memory

- *I/O* 354 reads, 210 writes
- paging 135 page faults, 0 swapouts

### Multiprocessor Interpretation of time Output

Timing results are interpreted in a different way when the program is run in parallel in a multiprocessor environment. Since /bin/time accumulates the user time on different threads, only wall clock time is used.

Since the user time displayed includes the time spent on all the processors, it can be quite large and is not a good measure of performance. A better measure is the real time, which is the wall clock time. This also means that to get an accurate timing of a parallelized program you must run it on a quiet system dedicated to just your program.

# The gprof Profiling Command

The <code>gprof(1)</code> command provides a detailed postmortem analysis of program timing at the subprogram level, including how many times a subprogram was called, who called it, whom it called, and how much time was spent in the routine and by the routines it called.

To enable gprof profiling, compile and link the program with the -pg option:

```
demo% f77 -o Myprog -fast -pg Myprog.f ...
demo% Myprog
demo% gprof Myprog
```

The program must complete normally for <code>gprof</code> to obtain meaningful timing information.

At program termination, the file <code>gmon.out</code> is automatically written in the working directory. This file contains the profiling data that will be interpreted by <code>gprof</code>.

Invoking gprof produces a report on standard output. An example is shown on the next pages. Not only the routines in your program are listed but also the library procedures and the routines they call.

The report is mostly two profiles of how the total time is distributed across the program procedures: the call graph and the flat profile. They are preceded by an explanation of the column labels, followed by an index. (The <code>gprof</code> -b option eliminates the explanatory text; see the <code>gprof(1)</code> man page for other options that can be used to limit the amount of output generated.)

In the graph profile, each procedure (subprogram, procedure) is presented in a call-tree representation. The line representing a procedure in its call-tree is called the *function line*, and is identified by an index number in the leftmost column, within square brackets; the lines above it are the *parent lines*; the lines below it, the *descendant* lines.

```
parent line caller 1
parent line caller 2
....

[n] time function line function name
descendant line called 1
descendant line called 2
....
```

The call graph profile is followed by a flat profile that provides a routine-by-routine overview. An (edited) example of gprof output follows.

**Note -** User-defined functions appear with their Fortran names followed by an underscore. Library routines appear with leading underscores.

#### The call graph profile:

granularity: each sample hit covers 2 byte(s) for 0.08% of 12.78 seconds

index	%time	self des	cendents		parents name index children
[3]	99.1	0.92 0.62 0.11 0.02 0.00	12.66 10.99 0.00 0.00 0.00 0.00	2000/2001 1000/1000 1000/1000 1/1 2/2	
[4]	93.2	0.92 1.11 1.29	10.99 4.52 2.91	1000 3000/3000 3000/6000	MAIN_ [3] diffr_ [4] deriv_ [7] cheb1_ [5] dissip_ [8]
[5]	65.7	1.29 2.58 5.81	2.91 5.81 0.00	3000/6000 6000	fftb_[6]

		0.00	0.00	128/192	sin [279]
[6]	45.5	5.81 5.81 0.00 0.00	0.00 0.00 0.00 0.00	6000/6000 6000 64/321 64/192	cheb1_ [5] fftb_ [6] cos [21] sin [279]

#### The flat profile overview:

granularity: each sample hit covers 2 byte(s) for 0.08% of 12.84 seconds

%	cumulative	self		self	total		
time	seconds	seconds	calls	ms/call	ms/call	name	
45.2	5.81	5.81	6000	0.97	0.97	fftb_	[6]
20.1	8.39	2.5	6000	0.43	1.40	cheb1_	[5]
9.1	9.56	1.17	3000	0.39	0.39	dissip_	[8]
8.6	10.67	1.11	3000	0.37	1.88	deriv_	[7]
7.1	11.58	0.92	1000	0.92	11.91	diffr_	[4]
4.8	12.20	0.62	2001	0.31	0.31	code_	[9]
2.5	12.53	0.33	69000	0.00	0.00	exp	[10]
0.9	12.64	0.11	1000	0.11	0.11	shock_	[11]

#### **■** Function Line.

The function line [5] in the preceding example reveals that:

- cheb1 was called 6000 times— 3000 from deriv, 3000 from diffr.
- 2.58 seconds were spent in cheb1 itself.
- 5.81 seconds were spent in routines called by cheb1.
- 65.7% of the execution time of the program was within cheb1.

#### **■** Parent Lines.

The parent lines above [5] indicate that cheb1 was called from two routines, deriv and diffr. The timings on these lines show how much time was spent in cheb1 when it was called from each of these routines.

#### Descendant Lines.

The lines below the function line indicate the routines called from cheb1, fftb, sin, and cos. The library sine function is called indirectly.

#### **■** Flat Profile.

Function names appear on the right. The profile is sorted by percentage of total execution time.

#### **Overhead Considerations**

Profiling (compiling with the -pg option) may greatly increase the running time of a program. This is due to the extra overhead required to clock program performance and subprogram calls. Profiling tools like gprof attempt to subtract an approximate overhead factor when computing relative runtime percentages. All other timings shown may not be accurate due to UNIX and hardware timekeeping inaccuracies.

Programs with short execution times are the most difficult to profile because the overhead may be a significant fraction of the total execution time. The best practice is to choose input data for the profiling run that will result in a realistic test of the program's performance. If this is not possible, consider enclosing the main computational part of the program within a loop that effectively runs the program N times. Estimate actual performance by dividing the profile results by N.

The Fortran library includes two routines that return the total time used by the calling process. See the man pages for *dtime*(3F) and *etime*(3F).

# The toov Profiling Command

The tcov(1) command, when used with programs compiled with the -a, -xa, or -xprofile=tcov options, produces a statement-by-statement profile of the source code showing which statements executed and how often. It also gives a summary of information about the basic block structure of the program.

There are two implementations of tcov coverage analysis. The original tcov is invoked by the -a or -xa compiler options. Enhanced statement level coverage is invoked by the -xprofile=tcov compiler option and the -x tcov option. In either case, the output is a copy of the source files annotated with statement execution counts in the margin. Although these two versions of tcov are essentially the same as far as the Fortran user is concerned (most of the enhancements apply to C++ programs), there will be some performance improvement with the newer style.

# "Old Style" toov Coverage Analysis

Compile the program with the -a (or -xa) option. This produces the file \$TCOVDIR/file.d for each source .f file in the compilation. (If environment variable \$TCOVDIR is not set at compile time, the .d files are stored in the current directory.)

Run the program (execution must complete normally). This produces updated information in the .d files. To view the coverage analysis merged with the individual source files, run toov on the source files. The annotated source files are named TCOVDIR/file.

The output produced by tcov shows the number of times each statement was actually executed. Statements that were not executed are marked with ####-> to the left of the statement.

Here is a simple example:

```
demo% f77 -a -o onetwo -silent one.f two.f
demo% onetwo
       ... output from program
demo% tcov one.f two.f
demo% cat one.tcov two.tcov
                      program one
                      do i=1,10
     10 ->
                            call two(i)
                      end do
     1 ->
                      end
             Top 10 Blocks
                       Count
             Line
               3
                            10
                2
                             1
                5
                              1
         3
                 Basic blocks in this file
         3
                 Basic blocks executed
                 Percent of the file executed
      100.00
                  Total basic block executions
          12
         4.00
                   Average executions per basic block
                      subroutine two(i)
     10 ->
                      print*, "two called", i
                      return
             Top 10 Blocks
             Line
                     Count
         1
                 Basic blocks in this file
         1
               Basic blocks executed
                Percent of the file executed
      100.00
          10
                   Total basic block executions
        10.00
                 Average executions per basic block
demo%
```

# "New Style" Enhanced toov Analysis

To use new style tcov, compile with -xprofile=tcov. When the program is run, coverage data is stored in *program*.profile/tcovd, where *program* is the name of the executable file. (If the executable were a.out, a.out.profile/tcovd would be created.)

Run tcov -x *dirname source\_files* to create the coverage analysis merged with each source file. The report is written to *file*.tcov in the current directory.

#### Running a simple example:

Environment variables \$SUN\_PROFDATA and \$SUN\_PROFDATA\_DIR can be used to specify where the intermediary data collection files are kept. These are the \*.d and toovd files created by old and new style toov, respectively.

Each subsequent run accumulates more coverage data into the tcovd file. Data for each object file is zeroed out the first time the program is executed after the corresponding source file has been recompiled. Data for the entire program is zeroed out by removing the tcovd file.

These environment variables can be used to separate the collected data from different runs. With these variables set, the running program writes execution data to the files in \$SUN\_PROFDATA\_DIR/\$SUN\_PROFDATA/.

Similarly, the directory that tcov reads is specified by tcov -x \$SUN\_PROFDATA. If \$SUN\_PROFDATA\_DIR is set, tcov will prepend it, looking for files in \$SUN\_PROFDATA\_DIR/\$SUN\_PROFDATA/, and not in the working directory.

For the details, see the *tcov*(1) man page.

# f77 I/O Profiling

You can obtain a report about how much data was transferred by your program. For each Fortran unit, the report shows the file name, the number of I/O statements, the number of bytes, and some statistics on these items.

To obtain an I/O profiling report, insert calls to the library routines start\_iostats and end\_iostats around the parts of the program you wish to measure. (A call to end\_iostats is required if the program terminates with an END or STOP statement rather than a CALL EXIT.)

Note - The I/O statements profiled are: READ, WRITE, PRINT, OPEN, CLOSE, INQUIRE, BACKSPACE, ENDFILE, and REWIND. The runtime system opens stdin, stdout, and stderr before the first executable statement of your program, so you must explicitly reopen these units after the call to start\_iostats.

Example: Profile stdin, stdout, and stderr:

```
EXTERNAL start iostats
CALL start_iostats
OPEN(5)
OPEN(6)
OPEN(0)
```

If you want to measure only part of the program, call end\_iostats to stop the process. A call to end\_iostats may also be required if your program terminates with an END or STOP statement rather than CALL EXIT.

The program must be compiled with the -pg option. When the program terminates, the I/O profile report is produced on the file name.io\_stats, where name is the name of the executable file.

#### Here is an example:

```
demo% f77 -o myprog -pg -silent myprog.f
demo% myprog
 ... output from program
demo% cat myprog.io_stats
            INPUT REPORT
1. unit 2. file name
                        input data
                                           4. map
                      total avg std dev
                   cnt
                                            (cnt)
          stderr 0 0 0 0
    Ω
                                            No
                       0
8
                                  0
                    0
                 2
1
0 0
0 0
8 48
4 48
9 48
                   2
                                4
    5
           stdin
                                         0
                                             No
                                8
                                         0
    6
           stdout
                                  0
                                         0
                                0
                                         Ω
   19
           fort.19
                                 6 4.276
                                 12
                                         0
   20
           fort.20
                                      4.276
                                 12
                   8
                          48
   21
           fort.21
                                      4.276
                                 6
                                             No
                          48
                                 12
                                         0
                          48
                                 6
   22
           fort.22
                                      4.276
                                             No
                          48
                                 12
                                         0
             OUTPUT REPORT
1. unit
              5. output data
                              6. blk size 7. fmt 8. direct
            total avg std dev
       cnt.
                                               (rec len)
            ______
              40 10 0 -1 Yes seq
```

	1	40	40	0			
5	0	0	0	0	-1	Yes	seq
	0	0	0	0			
6	26	248	9.538	1.63	-1	Yes	seq
	6	248	41.33	3.266			
19	8	48	6	4.276	500548	Yes	seq
	4	48	12	0			
20	8	48	6	4.276	503116	No	seq
	4	48	12	0			
21	8	48	6	4.276	503116	Yes	dir
	4	48	12	0			( 12)
22	8	48	6	4.276	503116	No	dir
	4	48	12	0			( 12)

Each pair of lines in the report displays information about an I/O unit. One section shows input operations and another shows output. The first line of a pair displays statistics on the number of data elements transferred before the unit was closed. The second row of statistics is based on the number of I/O statements processed.

In the example, there were 6 calls to write a total of 26 data elements to standard output. A total of 248 bytes was transferred. The display also shows the average and standard deviation in bytes transferred per I/O statement (9.538 and 1.63, respectively), and the average and standard deviation per I/O statement call (42.33 and 3.266, respectively).

The input report also contains a column to indicate whether a unit was memory mapped or not. If mapped, the number of mmap() calls is recorded in parentheses in the second row of the pair.

The output report indicates block sizes, formatting, and access type. A file opened for direct access shows its defined record length in parentheses in the second row of the pair.

**Note -** Compiling with environment variable LD\_LIBRARY\_PATH set might disable I/O profiling, which relies on its profiling I/O library being in a standard location.

# Performance and Optimization

This chapter considers some optimization techniques that may improve the performance of numerically intense Fortran programs. Proper use of algorithms, compiler options, library routines, and coding practices can bring significant performance gains. This discussion does not discuss cache, I/O, or system environment tuning. Parallelization issues are treated in the next chapter.

Some of the issues considered here are:

- Compiler options that may improve performance
- Compiling with feedback from runtime performance profiles
- Use of optimized library routines for common procedures
- Coding strategies to improve performance of key loops

The subject of optimization and performance tuning is much too complex to be treated exhaustively here. However, this discussion should provide the reader with a useful introduction to these issues. A list of books that cover the subject much more deeply appears at the end of the chapter.

Optimization and performance tuning is an art that depends heavily on being able to determine *what* to optimize or tune.

# **Choice of Compiler Options**

Choice of the proper compiler options is the first step in improving performance. Sun compilers offer a wide range of options that affect the object code. In the default case, where no options are explicitly stated on the compile command line, most options are off. To improve performance, these options must be explicitly selected.

Performance options are normally off by default because most optimizations force the compiler to make assumptions about a user's source code. Programs that conform to standard coding practices and do not introduce hidden side effects should optimize correctly. However, programs that take liberties with standard practices might run afoul of some of the compiler's assumptions. The resulting code might run faster, but the computational results might not be correct.

Recommended practice is to first compile with all options off, verify that the computational results are correct and accurate, and use these initial results and performance profile as a baseline. Then, proceed in steps—recompiling with additional options and comparing execution results and performance against the baseline. If numerical results change, the program might have questionable code, which needs careful analysis to locate and reprogram.

If performance does not improve significantly, or degrades, as a result of adding optimization options, the coding might not provide the compiler with opportunities for further performance improvements. The next step would then be to analyze and restructure the program at the source code level to achieve better performance.

# Performance Option Reference

The compiler options listed in the following table provide the user with a repertoire of strategies to improve the performance of a program over default compilation. Only some of the compilers' more potent performance options appear in the table. A more complete list can be found in the *Fortran User's Guide*.

TABLE 9-1 Some Effective Performance Options

Action	Option
Uses various optimization options together	-fast
Sets compiler optimization level to $n$	-on (-o = -o3)
Specifies target hardware	-xtarget= <i>sys</i>
Optimizes using performance profile data (with -05)	- xprofile=use
Unrolls loops by <i>n</i>	-unroll=n
Permits simplifications and optimization of floating-point	-fsimple=1 2
Performs dependency analysis to optimize loops	-depend

Some of these options increase compilation time because they invoke a deeper analysis of the program. Some options work best when routines are collected into files along with the routines that call them (rather than splitting each routine into its own file); this allows the analysis to be global.

#### -fast

This single option selects a number of performance options that, working together, produce object code optimized for execution speed without an excessive increase in compilation time.

The options selected by -fast are subject to change from one release to another, and not all are available on each platform:

- -native generates code optimized for the host architecture.
- -04 sets optimization level.
- -libmil inlines calls to some simple library functions.
- -fsimple=1 simplifies floating-point code (SPARC only).
- -dalign uses faster, double word loads and stores (SPARC only).
- -xlibmopt use optimized libm math library (SPARC only).
- -fns -ftrap=%none turns off all trapping.
- -depend analyze loops for data dependencies (SPARC only).
- -nofstore disables forcing precision on expressions (x86 only).

-fast provides a quick way to engage much of the optimizing power of the compilers. Each of the composite options may be specified individually, and each may have side effects to be aware of (discussed in the Fortran User's Guide). Following -fast with additional options adds further optimizations. For example:

sets the optimization to level 5 instead of 4.

**Note -** -fast includes -dalign and -native. These options may have unexpected side effects for some programs.

#### -0*n*

No compiler optimizations are performed by the compilers unless a -O option is specified explicitly (or implicitly with macro options like -fast). In nearly all cases, specifying an optimization level for compilation improves program execution performance. On the other hand, higher levels of optimization increase compilation time and may significantly increase code size.

For most cases, level -03 is a good balance between performance gain, code size, and compilation time. Level -04 adds automatic inlining of calls to routines contained in the same source file as the caller routine, among other things. Level -05 adds more aggressive optimization techniques that would not be applied at lower levels. In general, levels above -03 should be specified only to those routines that make up the most compute-intensive parts of the program and thereby have a high certainty of improving performance. (There is no problem linking together parts of a program compiled with different optimization levels.)

#### PRAGMA OPT=n

Use the C\$ PRAGMA SUN OPT=n directive to set different optimization levels for individual routines in a source file. This directive will override the -On flag on the compiler command line, but must be used with the -xmaxopt=n flag to set a maximum optimization level. See the f77(1) and f90(1) man pages for details.

#### Optimization With Runtime Profile Feedback

The compiler applies its optimization strategies at level O3 and above much more efficiently if combined with -xprofile=use. With this option, the optimizer is directed by a runtime execution profile produced by the program (compiled with -xprofile=collect) with typical input data. The feedback profile indicates to the compiler where optimization will have the greatest effect. This may be particularly important with -O5. Here's a typical example of profile collection with higher optimization levels:

```
demo% f77 -o prg -fast -xprofile=collect prg.f ...
demo% prg
demo% f77 -o prgx -fast -O5 -xprofile=use:prg.profile prg.f ...
demo% prgx
```

The first compilation in the example generates an executable that produces statement coverage statistics when run. The second compilation uses this performance data to guide the optimization of the program.

(See the Fortran User's Guide for details on -xprofile options.)

```
-dalign
```

With -dalign the compiler is able to generate double-word load/store instructions whenever possible. Programs that do much data motion may benefit significantly

when compiled with this option. (It is one of the options selected by <code>-fast.</code>) The double-word instructions are almost twice as fast as the equivalent single word operations.

However, users should be aware that using <code>-dalign</code> (and therefore <code>-fast</code>) may cause problems with some programs that have been coded expecting a specific alignment of data in COMMON blocks. With <code>-dalign</code>, the compiler may add padding to ensure that all double (and quad) precision data (either REAL or COMPLEX) are aligned on double-word boundaries, with the result that:

- COMMON blocks might be larger than expected due to added padding.
- All program units sharing COMMON must be compiled with -dalign if any one of them is compiled with -dalign.

For example, a program that writes data by aliasing an entire COMMON block of mixed data types as a single array might not work properly with <code>-dalign</code> because the block will be larger (due to padding of double and quad precision variables) than the program expects.

#### SPARC: -depend

Adding -depend to optimization levels -03 and higher (on the SPARC platform) extends the compiler's ability to optimize DO loops and loop nests. With this option, the optimizer analyzes inter-iteration loop dependencies to determine whether or not certain transformations of the loop structure can be performed. Only loops without dependencies can be restructured. However, the added analysis might increase compilation time.

#### -fsimple=2

Unless directed to, the compiler does not attempt to simplify floating-point computations (the default is <code>-fsimple=0</code>). With the <code>-fast</code> option, <code>-fsimple=1</code> is used and some conservative assumptions are made. Adding <code>-fsimple=2</code> enables the optimizer to make further simplifications with the understanding that this might cause some programs to produce slightly different results due to rounding effects. If <code>-fsimple</code> level 1 or 2 is used, all program units should be similarly compiled to ensure consistent numerical accuracy.

#### -unroll=n

Unrolling short loops with long iteration counts can be profitable for some routines. However, unrolling can also increase program size and might even degrade performance of other loops. With n=1, the default, no loops are unrolled automatically by the optimizer. With n greater than 1, the optimizer attempts to unroll loops up to a depth of n.

The compiler's code generator makes its decision to unroll loops depending on a number of factors. The compiler might decline to unroll a loop even though this option is specified with n>1.

If a DO loop with a variable loop limit can be unrolled, both an unrolled version and the original loop are compiled. A runtime test on iteration count determines whether or not executing the unrolled loop is inappropriate. Loop unrolling, especially with simple one or two statement loops, increases the amount of computation done per iteration and provides the optimizer with better opportunities to schedule registers and simplify operations. The tradeoff between number of iterations, loop complexity, and choice of unrolling depth is not easy to determine, and some experimentation might be needed.

The example that follows shows how a simple loop might be unrolled to a depth of four with -unroll=4 (the source code is not changed with this option):

```
Original Loop:
    DO I=1,20000
        X(I) = X(I) + Y(I)*A(I)
    END DO

Unrolled by 4 compiles as if it were written:
    DO I=1, 19997,4
        TEMP1 = X(I) + Y(I)*A(I)
        TEMP2 = X(I+1) + Y(I+1)*A(I+1)
        TEMP3 = X(I+2) + Y(I+2)*A(I+2)
        X(I+3) = X(I+3) + Y(I+3)*A(I+3)
        X(I) = TEMP1
        X(I+1) = TEMP2
        X(I+2) = TEMP3
    END DO
```

This example shows a simple loop with a fixed loop count. The restructuring is more complex with variable loop counts.

### -xtarget=*platform*

The performance of some programs might improve if the compiler has an accurate description of the target computer hardware. When program performance is critical, the proper specification of the target hardware could be very important. This is especially true when running on the newer SPARC processors. However, for most programs and older SPARC processors, the performance gain could be negligible and a generic specification might be sufficient.

The Fortran User's Guide lists all the system names recognized by -xtarget=. For any given system name (for example, ss1000, for SPARCserver 1000), -xtarget expands into a specific combination of -xarch, -xcache, and -xchip that properly matches that system. The optimizer uses these specifications to determine strategies to follow and instructions to generate.

The special setting -xtarget=native enables the optimizer to compile code targeted at the host system (the system doing the compilation). This is obviously useful when compilation and execution are done on the same system. When the execution system is not known, it is desirable to compile for a generic architecture. Therefore, -xtarqet=qeneric is the default, even though it might produce suboptimal performance.

# Other Performance Strategies

Assuming that you have experimented with using a variety of optimization options, compiling your program and measuring actual runtime performance, the next step might be to look closely at the Fortran source program to see what further tuning can be tried.

Focusing on just those parts of the program that use most of the compute time, you might consider the following strategies:

- Replace handwritten procedures with calls to equivalent optimized libraries.
- Remove I/O, calls, and unnecessary conditional operations from key loops.
- Eliminate aliasing that might inhibit optimization.
- Rationalize tangled, spaghetti-like code to use block IF.

These are some of the good programming practices that tend to lead to better performance. It is possible to go further, hand-tuning the source code for a specific hardware configuration. However, these attempts might only further obscure the code and make it even more difficult for the compiler's optimizer to achieve significant performance improvements. Excessive hand-tuning of the source code can hide the original intent of the procedure and could have a significantly detrimental effect on performance for different architectures.

# **Using Optimized Libraries**

In most situations, optimized commercial or shareware libraries perform standard computational procedures far more efficiently than you could by coding them by hand.

For example, the Sun Performance Library <sup>™</sup> is a suite of highly optimized mathematical subroutines based on the standard LAPACK, BLAS, FFTPACK, VFFTPACK, and LINPACK libraries. Performance improvement using these routines can be significant when compared with hand coding.

### **Eliminating Performance Inhibitors**

Use the profiling techniques described in Chapter 8 to identify the key computational parts of the program. Then, carefully analyze the loop or loop nest to eliminate coding that might either inhibit the optimizer from generating optimal code or otherwise degrade performance. Many of the nonstandard coding practices that make portability difficult might also inhibit optimization by the compiler.

Reprogramming techniques that improve performance are dealt with in more detail in some of the reference books listed at the end of the chapter. Three major approaches are worth mentioning here:

#### Removing I/O From Key Loops

I/O within a loop or loop nest enclosing the significant computational work of a program will seriously degrade performance. The amount of CPU time spent in the I/O library might be a major portion of the time spent in the loop. (I/O also causes process interrupts, thereby degrading program throughput.) By moving I/O out of the computation loop wherever possible, the number of calls to the I/O library can be greatly reduced.

### **Eliminating Subprogram Calls**

Subroutines called deep within a loop nest could be called thousands of times. Even if the time spent in each routine per call is small, the total effect might be substantial. Also, subprogram calls inhibit optimization of the loop that contains them because the compiler cannot make assumptions about the state of registers over the call.

Automatic inlining of subprogram calls (using -inline=x,y,..z, or -04) is one way to let the compiler replace the actual call with the subprogram itself (*pulling* the subprogram into the loop). The subprogram source code for the routines that are to be inlined must be found in the same file as the calling routine.

There are other ways to eliminate subprogram calls:

- Use statement functions. If the external function being called is a simple math function, it might be possible to rewrite the function as a statement function or set of statement functions. Statement functions are compiled in-line and can be optimized.
- Push the loop into the subprogram. That is, rewrite the subprogram so that it can be called fewer times (outside the loop) and operate on a vector or array of values per call.

### Rationalizing Tangled Code

Complicated conditional operations within a computationally intensive loop can dramatically inhibit the compiler's attempt at optimization. In general, a good rule to follow is to eliminate all arithmetic and logical IF's, replacing them with block IF's:

```
Original Code:
    IF(A(I)-DELTA) 10,10,11
10 XA(I) = XB(I)*B(I,I)
    XY(I) = XA(I) - A(I)
    GOTO 13
11 \quad XA(I) = Z(I)
    XY(I) = Z(I)
    IF(QZDATA.LT.0.) GOTO 12
    ICNT = ICNT + 1
    ROX(ICNT) = XA(I)-DELTA/2.
12 SUM = SUM + X(I)
13 SUM = SUM + XA(I)
Untangled Code:
    IF(A(I).LE.DELTA) THEN
      XA(I) = XB(I)*B(I,I)
      XY(I) = XA(I) - A(I)
    ELSE
      XA(I) = Z(I)
      XY(I) = Z(I)
      IF(QZDATA.GE.O.) THEN
        ICNT = ICNT + 1
        ROX(ICNT) = XA(I)-DELTA/2.
      ENDIF
      SUM = SUM + X(I)
    ENDIF
    SUM = SUM + XA(I)
```

Using block IF not only improves the opportunities for the compiler to generate optimal code, it also improves readability and assures portability.

# **Further Reading**

The following reference books provide more details:

- Numerical Computation Guide, Sun Microsystems, Inc.
- Analyzing Program Performance with Sun WorkShop, Sun Microsystems, Inc.
- FORTRAN Optimization, by Michael Metcalf, Academic Press 1985
- High Performance Computing, by Kevin Dowd, O'Reilly & Associates, 1993

# **SPARC:** Parallelization

This chapter presents an overview of multiprocessor parallelization and describes the capabilities of Sun's Fortran compilers. Implementation differences between £77 and £90 are noted.

**Note -** Parallelization features are available only on SPARC platforms with Solaris 2.5.1, 2.6, and Solaris 7 operating environments, and require a Sun WorkShop license.

# **Essential Concepts**

Parallelizing (or *multithreading*) an application recasts the compiled program to run on a multiprocessor system. Parallelization enables single tasks, such as a DO loop, to run over multiple processors with a potentially significant execution speedup.

Before an application program can be run efficiently on a multiprocessor system like the Ultra 60, Enterprise 450, or Ultra HPC 1000, it needs to be multithreaded. That is, tasks that can be performed in parallel need to be identified and reprogrammed to distribute their computations.

Multithreading an application can be done manually by making appropriate calls to the libthread primitives. However, a significant amount of analysis and reprogramming might be required. (See the Solaris *Multithreaded Programming Guide* for more information.)

Sun compilers can automatically generate multithreaded object code to run on multiprocessor systems. The Fortran compilers focus on DO loops as the primary language element supporting parallelism. Parallelization distributes the computational work of a loop over several processors without requiring modifications to the Fortran source program.

The choice of which loops to parallelize and how to distribute them can be left entirely up to the compiler (-autopar), determined explicitly by the programmer with source code directives (-explicitpar), or done in combination (-parallel).

**Note** - Programs that do their own (explicit) thread management should *not* be compiled with any of the compiler's parallelization options. Explicit multithreading (calls to libthread primitives) cannot be combined with routines compiled with these parallelization options.

Not all loops in a program can be profitably parallelized. Loops containing only a small amount of computational work (compared to the overhead spent starting and synchronizing parallel tasks) may actually run more slowly when parallelized. Also, some loops cannot be safely parallelized at all; they would compute different results when run in parallel due to dependencies between statements or iterations.

Only explicit Fortran 90 DO loops are candidates for parallelization with £90.

Sun compilers can detect loops that might be safely and profitably parallelized automatically. However, in most cases, the analysis is necessarily conservative, due to the concern for possible hidden side effects. (A display of which loops were and were not parallelized can be produced by the <code>-loopinfo</code> option.) By inserting source code directives before loops, you can explicitly influence the analysis, controlling how a specific loop is (or is not) to be parallelized. However, it then becomes your responsibility to ensure that such explicit parallelization of a loop does not lead to incorrect results.

# Speedups—What to Expect

If you parallelize a program so that it runs over four processors, can you expect it to take (roughly) one fourth the time that it did with a single processor (a fourfold *speedup*)?

Probably not. It can be shown (by Amdahl's law) that the overall speedup of a program is strictly limited by the fraction of the execution time spent in code running in parallel. This is true *no matter how many processors are applied*. In fact, if c is the percentage of the execution time run in parallel, the theoretical speedup limit is 100/(100-c); therefore, if only 60% of a program runs in parallel, the *maximum* increase in speed is 2.5, independent of the number of processors. And with just four processors, the theoretical speedup for this program (assuming maximum efficiency) would be just 1.8 and not 4. With overhead, the actual speedup would be less.

As with any optimization, choice of loops is critical. Parallelizing loops that participate only minimally in the total program execution time has only minimal effect. To be effective, the loops that consume the *major* part of the runtime *must* be parallelized. The first step, therefore, is to determine which loops are significant and to start from there.

Problem size also plays an important role in determining the fraction of the program running in parallel and consequently the speedup. Increasing the problem size increases the amount of work done in loops. A triply nested loop could see a cubic increase in work. If the outer loop in the nest is parallelized, a small increase in problem size could contribute to a significant performance improvement (compared to the unparallelized performance).

# Steps to Parallelizing a Program

Here is a very general outline of the steps needed to parallelize an application:

- 1. **Optimize**. Use the appropriate set of compiler options to get the best serial performance on a single processor.
- 2. **Profile**. Using typical test data, determine the performance profile of the program. Identify the most significant loops.
- 3. **Benchmark**. Determine that the serial test results are accurate. Use these results and the performance profile as the benchmark.
- 4. **Parallelize**. Use a combination of options and directives to compile and build a parallelized executable.
- 5. **Verify**. Run the parallelized program on a single processor and check results to find instabilities and programming errors that might have crept in.
- 6. Test. Make various runs on several processors to check results.
- 7. **Benchmark**. Make performance measurements with various numbers of processors on a dedicated system. Measure performance changes with changes in problem size (scalability).
- 8. **Repeat** steps 4 to 7. Make improvements to parallelization scheme based on performance.

# **Data Dependency Issues**

Not all loops are parallelizable. Running a loop in parallel over a number of processors may result in iterations executing out of order. Or the multiple processors executing the loop in parallel may interfere with each other. These situations arise whenever there are data dependencies in the loop.

#### Recurrence

Variables that are set in one iteration of a loop and used in a subsequent iteration introduce cross-iteration dependencies, or *recurrences*. Recurrence in a loop requires that the iterations to be executed in the proper order. For example:

```
DO I=2,N

A(I) = A(I-1)*B(I)+C(I)

END DO
```

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requires the value computed for A(I) in the previous iteration to be used (as A(I-1)) in the current iteration. To produce results running each iteration in parallel that are the same as with single processor, iteration I must complete before iteration I+1 can execute.

#### Reduction

Reduction operations reduce the elements of an array into a single value. For example, summing the elements of an array into a single variable involves updating that variable in each iteration:

```
DO K = 1,N
SUM = SUM + A(I)*B(I)
END DO
```

If each processor running this loop in parallel takes some subset of the iterations, the processors will interfere with each other, overwriting the value in SUM. For this to work, each processor must execute the summation one at a time, although the order is not significant.

Certain common reduction operations are recognized and handled as special cases by the compiler.

#### **Indirect Addressing**

Loop dependencies can result from stores into arrays that are indexed in the loop by subscripts whose values are not known. For example, indirect addressing could be order dependent if there are repeated values in the index array:

```
DO L = 1,NW
A(ID(L)) = A(L) + B(L)
END DO
```

In the preceding, repeated values in ID cause elements in A to be overwritten. In the serial case, the last store is the final value. In the parallel case, the order is not determined. The values of A(L) that are used, old or updated, are order dependent.

### **Data Dependent Loops**

You might be able to rewrite a loop to eliminate data dependencies, making it parallelizable. However, extensive restructuring could be needed.

Some general rules are:

- A loop is data *independent* only if all iterations write to distinct memory locations.
- Iterations may read from the same locations as long as no one iteration writes to them

These are general conditions for parallelization. The compilers' automatic parallelization analysis considers additional criteria when deciding whether to parallelize a loop. However, you can use directives to explicitly force loops to be parallelized, even loops that contain inhibitors and produce incorrect results.

# Parallel Options and Directives Summary

The following table shows the  $\tt f77~5.0$  and  $\tt f90~2.0$  compilation options related to parallelization.

TABLE 10-1 Parallelization Options

Option	Flag
Automatic (only)	-autopar
Automatic and Reduction	-autopar -reduction
Explicit (only)	-explicitpar
Automatic and Explicit	-parallel
Automatic and Reduction and Explicit	-parallel -reduction
Show which loops are parallelized	-loopinfo
Show warnings with explicit	-vpara
Allocate local variables on stack	-stackvar
Use Sun-style MP directives	-mp=sun
Use Cray-style MP directives	-mp=cray

#### Notes on these options:

- -reduction requires -autopar.
- -autopar includes -depend and loop structure optimization.
- -parallel is equivalent to -autopar -explicitpar.
- -noautopar, -noexplicitpar, -noreduction are the negations.

- Parallelization options can be in any order, but they must be all lowercase.
- Reduction operations are not analyzed for explicitly parallelized loops.
- Use of any of the parallelization options requires a WorkShop license.

The following table shows the f77/f90 and f90 parallel directives.

TABLE 10-2 Parallel Directives

Parallel Directive	Purpose
C\$PAR TASKCOMMON	Declares a common block private
C\$PAR DOALL optional qualifiers	Parallelizes next loop, if possible
C\$PAR DOSERIAL	Inhibits parallelization of next loop
C\$PAR DOSERIAL*	Inhibits parallelization of loop nest

### **Number of Processors**

The PARALLEL environment variable controls the maximum number of processors available to the program. The following example shows how to set it:

In this example, setting PARALLEL to four enables the execution of a program using at most four threads. If the target machine has four processors available, the threads will map to independent processors. If there are fewer than four processors available, some threads could run on the same processor as others, possibly degrading performance.

The SunOS command psrinfo(1M) displays a list of the processors available on a system:

```
demo% psrinfo
0    on-line    since 03/18/96 15:51:03
1    on-line    since 03/18/96 15:51:03
2    on-line    since 03/18/96 15:51:03
3    on-line    since 03/18/96 15:51:03
```

### Stacks, Stack Sizes, and Parallelization

The executing program maintains a main memory stack for the parent program and distinct stacks for each thread. Stacks are temporary memory address spaces used to hold arguments and AUTOMATIC variables over subprogram invocations.

The default size of the main stack is about 8 megabytes. The Fortran compilers normally allocate local variables and arrays as STATIC (not on the stack). However, the <code>-stackvar</code> option forces allocation of *all* local variables and arrays on the stack (as if they were AUTOMATIC variables). Use of <code>-stackvar</code> is recommended with parallelization because it improves the optimizer's ability to parallelize CALLs in loops. <code>-stackvar</code> is required with explicitly parallelized loops containing subprogram calls. (See the discussion of <code>-stackvar</code> in the Fortran User's Guide.)

The limit command displays the current main stack size as well as setting it:

```
C shell example
demo% limit
cputime
             unlimited
filesize
              unlimited
datasize
             2097148 kbytes
stacksize
               8192 kbytes
                                       <- current main stack size
coredumpsize
                  0 kbvtes
descriptors
                 64
memorysize
                unlimited
demo% limit stacksize 65536
                                  <- set main stack to 64Mb
demo% limit stacksize
stacksize
               65536 kbytes
demo$ >ulimit -a
                         Korn Shell example
time(seconds)
                    unlimited
file(blocks)
                    unlimited
data(kbytes)
                    2097148
stack(kbytes)
                    8192
coredump(blocks)
nofiles(descriptors) 64
vmemory(kbytes)
                    unlimited
demo$ ulimit -s 65536
demo$ ulimit -s
65536
```

Each thread of a multithreaded program has its own *thread* stack. This stack mimics the main program stack but is unique to the thread. The thread's PRIVATE arrays and variables (local to the thread) are allocated on the thread stack. The default size is 256 kilobytes. The size is set with the STACKSIZE environment variable:

```
demo% setenv STACKSIZE 8192 <- Set thread stack size to 8 Mb C shell demo% STACKSIZE=8192 Bourne/Korn Shell demo% export STACKSIZE
```

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Setting the thread stack size to a value larger than the default may be necessary for most parallelized Fortran codes. However, it may not be possible to know just how large to set it, except by trial and error, especially if private/local arrays are involved. If the stack size is too small for a thread to run, the program will abort with a segmentation fault.

# **Automatic Parallelization**

With the <code>-autopar</code> and <code>-parallel</code> options, the compilers automatically find DO loops that can be parallelized effectively. These loops are then transformed to distribute their iterations evenly over the available processors. The compiler generates the thread calls needed to make this happen.

# **Loop Parallelization**

The compiler's dependency analysis transforms a DO loop into a parallelizable task. The compiler may restructure the loop to split out unparallelizable sections that will run serially. It then distributes the work evenly over the available processors. Each processor executes a different chunk of iterations.

For example, with four CPUs and a parallelized loop with 1000 iterations:

Processor 1 executing iterations	1	through	250
Processor 2 executing iterations	251	through	500
Processor 3 executing iterations	501	through	750
Processor 4 executing iterations	751	through	1000

Only loops that do not depend on the order in which the computations are performed can be successfully parallelized. The compiler's dependency analysis rejects loops with inherent data dependencies. If it cannot fully determine the data flow in a loop, the compiler acts conservatively and does not parallelize. Also, it may choose not to parallelize a loop if it determines the performance gain does not justify the overhead.

Note that the compiler always chooses to parallelize loops using a *chunk* distribution—simply dividing the work in the loop into equal blocks of iterations. Other distribution schemes may be specified using explicit parallelization directives described later in this chapter.

### Arrays, Scalars, and Pure Scalars

A few definitions, from the point of view of automatic parallelization, are needed:

- An array is a variable that is declared with at least one dimension.
- A scalar is a variable that is not an array.
- A *pure scalar* is a scalar variable that is not aliased—not referenced in an EQUIVALENCE or POINTER statement.

#### Example: Array/scalar:

```
dimension a(10)
real m(100,10), s, u, x, z
equivalence ( u, z )
pointer ( px, x )
s = 0.0
...
```

Both  $\mathfrak m$  and a are array variables; s is pure scalar. The variables s, s, s, and s are scalar variables, but not *pure* scalars.

### Automatic Parallelization Criteria

DO loops that have no cross-iteration data dependencies are automatically parallelized by -autopar or -parallel. The general criteria for automatic parallelization are:

- DO loops are parallelized, but not DO WHILE or Fortran 90 array operations.
- The values of *array* variables for each iteration of the loop must not depend on the values of *array* variables for any other iteration of the loop.
- Calculations within the loop must not *conditionally* change any pure scalar variable that is referenced after the loop terminates.
- Calculations within the loop must not change a scalar variable across iterations.
   This is called a loop-carried dependency.

### £77: Apparent Dependencies

The £77 compiler may automatically eliminate a reference that appears to create a dependency transforming the compiled code. One of the many such transformations makes use of private versions of some of the arrays. Typically, the compiler does this if it can determine that such arrays are used in the original loops only as temporary storage.

Example: Using -autopar, with dependencies eliminated by private arrays:

```
parameter (n=1000)
real a(n), b(n), c(n,n)
```

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```
<--Parallelized
```

```
do i = 1, 1000
  do k = 1, n
    a(k) = b(k) + 2.0
  end do
  do j = 1, n
    c(i,j) = a(j) + 2.3
  end do
end do
end
```

In the preceding example, the outer loop is parallelized and run on independent processors. Although the inner loop references to array a(\*) appear to result in a data dependency, the compiler generates temporary private copies of the array to make the outer loop iterations independent.

#### Inhibitors to Automatic Parallelization

Under automatic parallelization, the compilers do not parallelize a loop if:

- The DO loop is nested inside another DO loop that is parallelized.
- Flow control allows jumping out of the DO loop.
- A user-level subprogram is invoked inside the loop.
- An I/O statement is in the loop.
- Calculations within the loop change an aliased scalar variable.

#### **Nested Loops**

On multiprocessor systems, it is most effective to parallelize the outermost loop in a loop nest, rather than the innermost. Because parallel processing typically involves relatively large loop overhead, parallelizing the outermost loop minimizes the overhead and maximizes the work done for each processor. Under automatic parallelization, the compilers start their loop analysis from the outermost loop in a nest and work inward until a parallelizable loop is found. Once a loop within the nest is parallelized, loops contained within the parallel loop are passed over.

# Automatic Parallelization With Reduction Operations

A computation that transforms an array into a scalar is called a *reduction operation*. Typical reduction operations are the sum or product of the elements of a vector. Reduction operations violate the criterion that calculations within a loop not change a scalar variable in a cumulative way across iterations.

Example: Reduction summation of the elements of a vector:

```
s = 0.0

do i = 1, 1000

s = s + v(i)

end do

t(k) = s
```

However, for some operations, if the reduction is the only factor that prevents parallelization, it is still possible to parallelize the loop. Common reduction operations occur so frequently that the compilers are capable of recognizing and parallelizing them as special cases.

Recognition of reduction operations is not included in the automatic parallelization analysis unless the -reduction compiler option is specified along with -autopar or -parallel.

If a parallelizable loop contains one of the reduction operations listed in Table 10–3, the compiler will parallelize it if -reduction is specified.

#### **Recognized Reduction Operations**

The following table lists the reduction operations that are recognized by £77 and £90.

 TABLE 10-3
 Recognized Reduction Operations

Mathematical Operations	Fortran Statement Templates
Sum of the elements	s = s + v(i)
Product of the elements	s = s * v(i)
Dot product of two vectors	s = s + v(i) * u(i)
Minimum of the elements	s = amin(s, v(i))
Maximum of the elements	s = amax(s, v(i))
OR of the elements	do i = 1, n
	$b = b \cdot or \cdot v(i)$
	end do

 TABLE 10–3
 Recognized Reduction Operations (continued)

<b>Mathematical Operations</b>	Fortran Statement Templates
AND of nonpositive elements	b = .true.
	do i = 1, n
	if $(v(i)$ .le. 0) b=b .and. $v(i)$
	end do
Count nonzero elements	k = 0
	do i = 1, n
	if ( $v(i)$ .ne. 0 ) $k = k + 1$
	end do

All forms of the MIN and MAX function are recognized.

#### **Numerical Accuracy and Reduction Operations**

Floating-point sum or product reduction operations may be inaccurate due to the following conditions:

- The order in which the calculations were performed in parallel was not the same as when performed serially on a single processor.
- The order of calculation affected the sum or product of floating-point numbers. Hardware floating-point addition and multiplication are not associative. Roundoff, overflow, or underflow errors may result depending on how the operands associate. For example, (X\*Y)\*Z and X\*(Y\*Z) may not have the same numerical significance.

In some situations, the error may not be acceptable.

Example: Overflow and underflow, with and without reduction:

```
demo% cat t3.f
    real A(10002), result, MAXFLOAT
    MAXFLOAT = r_max_normal()
    do 10 i = 1 , 10000, 2
    A(i) = MAXFLOAT
    A(i+1) = -MAXFLOAT
10    continue

A(5001)=-MAXFLOAT
    A(5002)=MAXFLOAT
```

```
do 20 i = 1 ,10002
                                  !Add up the array
        RESULT = RESULT + A(i)
20
        continue
      write(6,*) RESULT
      end
demo% setenv PARALLEL 2
                                    {Number of processors is 2}
demo% f77 -silent -autopar t3.f
demo% a.out
                                    {Without reduction, 0. is correct}
   0.
demo% f77 -silent -autopar -reduction t3.f
demo% a.out
                                    {With reduction, Inf. is not correct}
  Inf
demo%
```

Example: Roundoff, get the sum of 100,000 random numbers between -1 and +1:

```
demo% cat t4.f
    parameter ( n = 100000 )
    double precision d_lcrans, lb / -1.0 /, s, ub / +1.0 /, v(n)
    s = d_lcrans ( v, n, lb, ub ) ! Get n random nos. between -1 and +1
    s = 0.0
    do i = 1, n
        s = s + v(i)
    end do
    write(*, '(" s = ", e21.15)') s
    end
demo% f77 -autopar -reduction t4.f
```

Results vary with the number of processors. The following table shows the sum of 100,000 random numbers between –1 and +1.

<b>Number of Processors</b>	Output
1	s = 0.568582080884714E+02
2	s = 0.568582080884722E+02
3	s = 0.568582080884721E+02
4	s = 0.568582080884724E + 02

In this situation, roundoff error on the order of  $10^{-14}$  is acceptable for data that is random to begin with. For more information, see the Sun *Numerical Computation Guide*.

## **Explicit Parallelization**

This section describes the source code directives recognized by £77 5.0 and £90 2.0 to explicitly indicate which loops to parallelize and what strategy to use.

Explicit parallelization of a program requires prior analysis and deep understanding of the application code as well as the concepts of shared-memory parallelization.

DO loops are marked for parallelization by directives placed immediately before them. The compiler options <code>-parallel</code> and <code>-explicitpar</code> must be used for DO loops to be recognized and parallel code generated. Take care when choosing which loops to mark for parallelization. The compiler generates threaded, parallel code for all loops marked with DOALL directives, even if there are data dependencies that will cause the loop to compute incorrect results when run in parallel.

If you do your own multithreaded coding using the libthread primitives, do *not* use any of the compilers' parallelization options—the compilers cannot parallelize code that has already been parallelized with user calls to the threads library.

## Parallelizable Loops

A loop is appropriate for explicit parallelization if:

- It is a DO loop, but not a DO WHILE or Fortran 90 array syntax.
- The values of array variables for each iteration of the loop do not depend on the values of array variables for any other iteration of the loop.
- If the loop changes a scalar, that scalar is not referenced after the loop terminates. Such scalar variables are not guaranteed to have a defined value after the loop terminates, since the compiler does not automatically ensure a proper storeback for them.
- For each iteration, any subprogram that is invoked inside the loop does not reference or change values of array variables for any other iteration.
- The DO loop index must be an integer.

## Scoping Rules: Private and Shared

A *private* variable or array is private to a *single iteration* of a loop. The value assigned to a private variable or array in one iteration is not propagated to any other iteration of the loop.

A *shared* variable or array is shared with all other iterations. The value assigned to a shared variable or array in an iteration is seen by other iterations of the loop.

If an explicitly parallelized loop contains shared references, then you must ensure that sharing does not cause correctness problems. The compiler does no synchronization on updates or accesses to shared variables.

If you specify a variable as private in one loop, and its only initialization is within some other loop, the value of that variable may be left undefined in the loop.

#### Default Scoping Rules for Sun-Style Directives

For Sun-style (C\$PAR) explicit directives, the compiler uses default rules to determine whether a scalar or array is shared or private. You can override the default rules to specify the attributes of scalars or arrays referenced inside a loop. (With Cray-style !MIC\$ directives, all variables that appear in the loop must be explicitly declared either shared or private on the DOALL directive.)

The compiler applies these default rules:

- All scalars are treated as *private*. A processor local copy of the scalar is made in each processor, and that local copy is used within that process.
- All array references are treated as *shared* references. Any write of an array element by one processor is visible to all processors. No synchronization is performed on accesses to shared variables.

If inter-iteration dependencies exist in a loop, then the execution may result in erroneous results. You must ensure that these cases do not arise. The compiler may sometimes be able to detect such a situation at compile time and issue a warning, but it does not disable parallelization of such loops.

Example: Potential problem through equivalence:

```
equivalence (a(1),y)
C$PAR DOALL
   do i = 1,n
        y = i
        a(i) = y
   end do
```

In the preceding example, since the scalar variable y has been equivalenced to a(1), it is no longer a private variable, even though the compiler treats it as such by the default scoping rule. Thus, the presence of the DOALL directive might lead to erroneous results when the parallelized  $\dot{\text{l}}$  loop is executed.

You can fix the example by using C\$PAR DOALL PRIVATE(y).

## Sun-Style Parallelization Directives

Parallelization directives are comment lines that tell the compiler to parallelize (or not to parallelize) the DO loop that follows the directive. Directives are also called *pragmas*.

A parallelization directive consists of one or more directive lines.

Sun-style directives are recognized by £77 and £90 by default (or with the -mp=sun option). A Sun-style directive line is defined as follows:

```
C$PAR Directive [ Qualifiers ] <- Initial directive line
C$PAR& [More_Qualifiers] <- Optional continuation lines
```

- The letters of a directive line are case-insensitive.
- The first five characters are C\$PAR, \*\$PAR, or !\$PAR.
- An *initial* directive line has a blank in column 6.
- A *continuation* directive line has a nonblank in column 6.
- Directives are listed in columns 7 and beyond.
- Qualifiers, if any, follow directives—on the same line or continuation lines.
- Multiple qualifiers on one line are separated by commas.
- Spaces before, after, or within a directive or qualifier are ignored.
- Columns beyond 72 are ignored unless the -e option is specified.

The parallel directives and their actions are as follows:

Directive	Action
TASKCOMMON	Declares COMMON block private
DOALL	Parallelizes the next loop
DOSERIAL	Does not parallelize the next loop
DOSERIAL*	Does not parallelize the next nest of loops

#### Examples of £77 parallel directives:

C\$PAR TASKCOMMON ALPHA
COMMON /ALPHA/BZ,BY(100)

C\$PAR DOALL

C\$PAR DOSERIAL

No qualifiers

```
C$PAR DOALL SHARED(I,K,X,V), PRIVATE(A)

This one-line directive is equivalent to the three-line directive that follows. 
C$PAR DOALL 
C$PAR& SHARED(I,K,X,V) 
C$PAR& PRIVATE(A)
```

#### TASKCOMMON Directive

The TASKCOMMON directive declares variables in a global COMMON block as *private*. Every variable declared in a task common block becomes a private variable. Only named COMMON blocks can be declared TASK COMMON.

The syntax of the directive is:

```
C$PAR TASKCOMMON common_block_name
```

The directive must appear immediately after the defining COMMON declaration.

This directive is effective only when compiled with <code>-explicitpar</code> or <code>-parallel</code>. Otherwise, the directive is ignored and the block is treated as a regular common block.

Variables declared in task common blocks are treated as private variables in all the DOALL loops they appear in explicitly, and in the routines called from a loop where the specified common block is in its scope.

It is an error to declare a common block as task common in some but not *all* compilation units where the block is defined. A check at runtime for task common consistency can be enabled by compiling the program with the -xcommonchk=yes flag. (Enable the runtime check only during program development, as it can degrade performance.)

#### DOALL Directive

The compilers will parallelize the DO loop following a DOALL directive (if compiled with the -parallel or -explicit par options).

**Note -** Analysis and transformation of reduction operations within loops is not done if they are explicitly parallelized.

Example: Explicit parallelization of a loop:

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```
do k = 1, m
x(k) = x(k) * z(k,k)
end do
...
demo% f77 -explicitpar t4.f
```

#### Subprogram Call in a Loop

A subprogram call in a loop (or in any subprograms called from within the called routine) may introduce data dependencies that could go unnoticed without a deep analysis of the data and control flow through the chain of calls. While it is best to parallelize outermost loops that do a significant amount of the work, these tend to be the very loops that involve subprogram calls.

Because such an interprocedural analysis is difficult and could greatly increase compilation time, automatic parallelization modes do not attempt it. With explicit parallelization, the compiler generates parallelized code for a loop marked with a DOALL directive that contains calls to subprograms. It is still the programmer's responsibility toeinsure that no data dependencies exist within the loop and all that the loop encloses, including called subprograms.

Multiple invocations of a routine from different processors can cause problems resulting from references to local static variables that interfere with each other. Making all the local variables in a routine *automatic* rather than *static* prevents this. Each invocation of a subprogram then has its own unique store of local variables maintained on the stack, and no two invocations will interfere with each other.

Local subprogram variables can be made automatic variables that reside on the stack either by listing them on an AUTOMATIC statement or by compiling the subprogram with the <code>-stackvar</code> option. However, local variables initialized in DATA statements must be rewritten to be initialized in actual assignments.

**Note -** Allocating local variables to the stack can cause stack overflow. See "Stacks, Stack Sizes, and Parallelization" on page 10-7 about increasing the size of the stack.

Data dependencies can still be introduced through the data passed down the call tree as arguments or through COMMON blocks. This data flow should be analyzed carefully before parallelizing a loop with subprogram calls.

### **DOALL Qualifiers**

All qualifiers on the DOALL directive are optional. The following table summarizes them:

TABLE 10-4 DOALL Qualifiers

Qualifier	Assertion	Syntax
PRIVATE	Do not share variables <i>u1</i> , between iterations	DOALL PRIVATE(u1,u2,)
SHARED	Share variables $v1$ , $v2$ , between iterations	DOALL SHARED( $v1, v2,$ )
MAXCPUS	Use no more than <i>n</i> CPUs	DOALL MAXCPUS(n)
READONLY	The listed variables are <i>not</i> modified in the DOALL loop	DOALL READONLY( $v1, v2,$ )
SAVELAST	Save the last $DO$ iteration values of all <i>private</i> variables	DOALL SAVELAST
STOREBACK	Save the last DO iteration values of variables $v1$ ,	DOALL STOREBACK(v1,v2,)
REDUCTION	Treat the variables $v1$ , $v2$ , as reduction variables.	DOALL REDUCTION( $v1, v2$ ,)
SCHEDTYPE	Set the scheduling type to <i>t</i> .	DOALL SCHEDTYPE( $t$ )

#### PRIVATE (varlist)

The PRIVATE (varlist) qualifier specifies that all scalars and arrays in the list varlist are private for the DOALL loop. Both arrays and scalars can be specified as private. In the case of an array, each thread of the DOALL loop gets a copy of the entire array. All other scalars and arrays referenced in the DOALL loop, but not contained in the private list, conform to their appropriate default scoping rules.

Example: Specify a private array:

```
C$PAR DOALL PRIVATE(a)
    do i = 1, n
        a(1) = b(i)
    do j = 2, n
        a(j) = a(j-1) + b(j) * c(j)
    end do
    x(i) = f(a)
end do
```

In the preceding example, the array a is specified as private to the i loop.

#### SHARED (varlist)

The SHARED(varlist) qualifier specifies that all scalars and arrays in the list varlist are shared for the DOALL loop. Both arrays and scalars can be specified as shared. Shared scalars and arrays are common to all the iterations of a DOALL loop. All other scalars and arrays referenced in the DOALL loop, but not contained in the shared list, conform to their appropriate default scoping rules.

Example: Specify a shared variable:

```
equivalence (a(1),y)
C$PAR DOALL SHARED(y)
do i = 1,n
     a(i) = y
end do
```

In the preceding example, the variable y has been specified as a variable whose value should be shared among the iterations of the i loop.

#### READONLY (varlist)

The READONLY (varlist) qualifier specifies that all scalars and arrays in the list varlist are read-only for the DOALL loop. Read-only scalars and arrays are a special class of shared scalars and arrays that are not modified in any iteration of the DOALL loop. Specifying scalars and arrays as READONLY indicates to the compiler that it does not need to use a separate copy of that variable or array for each thread of the DOALL loop.

Example: Specify a read-only variable:

```
x = 3
C$PAR DOALL SHARED(x), READONLY(x)
do i = 1, n
b(i) = x + 1
end do
```

In the preceding example, x is a shared variable, but the compiler can rely on the fact that it will not change over each iteration of the i loop because of its READONLY specification.

#### STOREBACK (varlist)

A STOREBACK variable or array is one whose value is computed in a DOALL loop. The computed value can be used after the termination of the loop. In other words, the last loop iteration values of storeback scalars and arrays may be visible outside of the DOALL loop.

Example: Specify the loop index variable as storeback:

```
C$PAR DOALL PRIVATE(x), STOREBACK(x,i)
    do i = 1, n
        x = ...
    end do
    ... = i
    ... = x
```

In the preceding example, both the variables x and i are STOREBACK variables, even though both variables are private to the i loop.

There are some potential problems for STOREBACK, however.

The STOREBACK operation occurs at the last iteration of the explicitly parallelized loop, even if this is the same iteration that last updates the value of the STOREBACK variable or array.

Example: STOREBACK variable potentially different from the serial version:

```
C$PAR DOALL PRIVATE(x), STOREBACK(x)
  do i = 1, n
      if (...) then
      x = ...
  end if
  end do
  print *,x
```

In the preceding example, the value of the STOREBACK variable x that is printed out might not be the same as that printed out by a serial version of the i loop. In the explicitly parallelized case, the processor that processes the last iteration of the i loop (when i=n) and performs the STOREBACK operation for x, might not be the same processor that currently contains the last updated value of x. The compiler issues a warning message about these potential problems.

In an explicitly parallelized loop, arrays are not treated *by default* as STOREBACK, so include them in the list *varlist* if such a storeback operation is desired—for example, if the arrays have been declared as private.

#### SAVELAST

The SAVELAST qualifier specifies that all private scalars and arrays are STOREBACK for the DOALL loop. A STOREBACK variable or array is one whose value is computed in a DOALL loop; this computed value can be used after the termination of the loop. In other words, the last loop iteration values of STOREBACK scalars and arrays may be visible outside of the DOALL loop.

Example: Specify SAVELAST:

In the preceding example, variables x, y, and i are STOREBACK variables.

#### REDUCTION (varlist)

The REDUCTION(varlist) qualifier specifies that all variables in the list varlist are reduction variables for the DOALL loop. A reduction variable (or array) is one whose partial values can be individually computed on various processors, and whose final value can be computed from all its partial values.

The presence of a list of reduction variables can aid the compiler in identifying if a DOALL loop is a reduction loop and in generating parallel reduction code for it.

Example: Specify a reduction variable:

```
C$PAR DOALL REDUCTION(x)
    do i = 1, n
        x = x + a(i)
    end do
```

In the preceding example, the variable x is a (sum) reduction variable; the i loop is a (sum) reduction loop.

#### SCHEDTYPE (t)

The SCHEDTYPE(t) qualifier specifies that the specific scheduling type t be used to schedule the DOALL loop.

### TABLE 10-5 DOALL SCHEDTYPE Qualifiers

Scheduling Type	Action
STATIC	Use <i>static scheduling</i> for this DO loop.
	Distribute all iterations uniformly to all available processors.
	Example: With 1000 iterations and 4 CPUs each CPU gets a single iteration in turn until all the iterations have been distributed.
SELF[(chunksize)	]Use self-scheduling for this DO loop.
	Distribute chunksize iterations to each available processor:
	• Repeat with the remaining iterations until all the iterations have been processed. • If <i>chunksize</i> is not provided, £77 selects a value.
	Example: With 1000 iterations and $\it chunksize$ of 4, distribute 4 iterations to each CPU.
FACTORING[(	Use factoring scheduling for this DO loop.
<i>m</i> )]	With $n$ iterations initially and $k$ CPUs, distribute $n/(2k)$ iterations uniformly to each processor until all iterations have been processed.
	$\bullet$ At least $m$ iterations must be assigned to each processor.
	• There can be one final smaller residual chunk.
	• If $m$ is not provided, £77 selects a value.
	Example: With 1000 iterations and $FACTORING(4)$ , and 4 CPUs, distribute 125 iterations to each CPU, then 62 iterations, then 31 iterations, and so on.
GSS[( m )]	Use <i>guided self-scheduling</i> for this DO loop.
	With $n$ iterations initially, and $k$ CPUs, then:
	• Assign $n/k$ iterations to the first processor.
	$\bullet$ Assign the remaining iterations divided by $k$ to the second processor, and so on until all iterations have been processed.
	Note:
	• At least <i>m</i> iterations must be assigned to each CPU.
	• There can be one final smaller residual chunk.
	• If $m$ is not provided, £77 selects a value.
	Example: With 1000 iterations and GSS(10), and 4 CPUs, distribute 250 iterations to the first CPU, then 187 to the second CPU, then 140 to the third CPU, and so on.

### **Multiple Qualifiers**

Qualifiers can appear multiple times with cumulative effect. In the case of conflicting qualifiers, the compiler issues a warning message, and the qualifier appearing last prevails.

Example: A three-line Sun-style directive:

```
C$PAR DOALL MAXCPUS(4), READONLY(S), PRIVATE(A,B,X), MAXCPUS(2) C$PAR DOALL SHARED(B,X,Y), PRIVATE(Y,Z) C$PAR DOALL READONLY(T)
```

Example: A one-line equivalent of the preceding three lines (note duplicate MAXCPUS and conflicting SHARED/PRIVATE):

```
C$PAR DOALL MAXCPUS(2), PRIVATE(A,Y,Z), SHARED(B,X), READONLY(S,T)
```

#### DOSERIAL Directive

The DOSERIAL directive disables parallelization of the specified loop. This directive applies to the one loop immediately following it (if you compile it with -explicitpar or -parallel).

Example: Exclude one loop from parallelization:

```
do i = 1, n
C$PAR DOSERIAL
    do j = 1, n
        do k = 1, n
        ...
    end do
    end do
    end do
end do
```

In the preceding example, the j loop is not parallelized, but the i or k loop can be.

#### DOSERIAL\* Directive

The DOSERIAL\* directive disables parallelization the specified nest of loops. This directive applies to the whole nest of loops immediately following it (if you compile with -explicitpar or -parallel).

Example: Exclude a whole nest of loops from parallelization:

```
do i = 1, n
C$PAR DOSERIAL*
do j = 1, n
```

```
\begin{array}{ccccc} \text{do } k = 1, \text{ n} \\ & \dots \\ & \text{end do} \\ & \text{end do} \\ & \text{end do} \end{array}
```

In the preceding loops, the j and k loops are not parallelized; the i loop could be.

#### Interaction Between DOSERIAL\* and DOALL

If both DOSERIAL and DOALL are specified, the last one prevails.

Example: Specifying both DOSERIAL and DOALL:

```
C$PAR DOSERIAL*
do i = 1, 1000
C$PAR DOALL
do j = 1, 1000
...
end do
end do
```

In the preceding example, the i loop is not parallelized, but the j loop is.

Also, the scope of the  ${\tt DOSERIAL*}$  directive does not extend beyond the textual loop nest immediately following it. The directive is limited to the same function or subroutine that it is in.

Example: DOSERIAL\* does not extend to a loop of a called subroutine:

```
program caller
  common /block/ a(10,10)
C$PAR DOSERIAL*
  do i = 1, 10
     call callee(i)
  end do
  end

subroutine callee(k)
  common /block/a(10,10)
  do j = 1, 10
     a(j,k) = j + k
  end do
  return
  end
```

In the preceding example, DOSERIAL\* applies only to the i loop and not to the j loop, regardless of whether the call to the subroutine callee is inlined.

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### Inhibitors to Explicit Parallelization

In general, the compiler parallelizes a loop if you explicitly direct it to. There are exceptions—some loops the compiler just cannot parallelize.

The following are the primary detectable inhibitors that might prevent explicitly parallelizing a DO loop:

- The DO loop is nested inside another DO loop that is parallelized.
  - This exception holds for indirect nesting, too. If you explicitly parallelize a loop that includes a call to a subroutine, then even if you parallelize loops in that subroutine, those loops are not run in parallel at runtime.
- A flow control statement allows jumping out of the DO loop.
- The index variable of the loop is subject to side effects, such as being equivalenced.

If you compile with -vpara, you may get a warning message if f77/f90 detects a problem with explicitly parallelizing a loop. f77/f90 may still parallelize the loop. The following list of typical parallelization problems shows those that are ignored by the compiler

TABLE 10-6 Explicit Parallelization Problems

Problem	Parallelized	Message
Loop is nested inside another loop that is parallelized.	No	No
Loop is in a subroutine, and a call to the subroutine is in a parallelized loop.	No	No
Jumping out of loop is allowed by a flow control statement.	No	Yes
Index variable of loop is subject to side effects.	Yes	No
Some variable in the loop keeps a loop-carried dependency.	Yes	Yes
I/O statement in the loop—usually unwise, because the order of the output is not predictable.	Yes	No

and those that generate messages with -vpara.

Example: Nested loops:

```
C$PAR DOALL
do 900 i = 1, 1000 ! Parallelized (outer loop)
```

```
do 200 j = 1, 1000 ! Not parallelized, no warning
...
200 continue
900 continue
...
demo% f77 -explicitpar -vpara t6.f
```

Example: A parallelized loop in a subroutine:

```
C$PAR DOALL

do 100 i = 1, 200

...

C$PAR DOALL

call calc (a, x)

do 1 m = 1, 1000

...

100 continue

1 continue

return

demo% f77 -explicitpar -vpara t.f

At runtime, the loop could run in parallel.

At runtime, both loops do not run in parallel.
```

In the preceding example, the loop within the subroutine is not parallelized because the subroutine itself is run in parallel.

Example: Jumping out of a loop:

Example: An index variable subject to side effects:

```
equivalence ( a(1), y ) !

" Source of possible side effects
...

C$PAR DOALL
do i = 1, 2000 !

" Parallelized: no warning, but not safe
y = i
a(i) = y
end do
...

demo% f77 -explicitpar -vpara t11.f
```

Example: A variable in a loop has a loop-carried dependency:

#### I/O With Explicit Parallelization

You can do I/O in a loop that executes in parallel, provided that:

- It does not matter that the output from different threads is interleaved (program output is nondeterministic.)
- You can ensure the safety of executing the loop in parallel.

Example: I/O statement in loop

```
C$PAR DOALL
    do i = 1, 10    ! Parallelized with no warning (not advisable)
        k = i
        call show ( k )
    end do
    end
    subroutine show( j )
    write(6,1) j
1    format("Line number ", i3, ".")
    end
demo% f77 -silent -explicitpar -vpara t13.f
demo% setenv PARALLEL 2
demo% a.out
(The output displays the numbers 1 through 10, but in a different order each time.)
```

Example: Recursive I/O:

```
do i = 1, 10
<-- Parallelized with no warning —unsafe
    k = i
    print *, list(k) <-- list is a function that does I/O
    end do
    end
    function list(j)
    write(6,"('Line number ', i3, '.')") j
    list = j
    end
demo% f77 -silent -mt t14.f
demo% setenv PARALLEL 2
demo% a.out</pre>
```

In the preceding example, the program may deadlock in libF77\_mt and hang. Press Control-C to regain keyboard control.

There are situations where the programmer might not be aware that I/O could take place within a parallelized loop. Consider a user-supplied exception handler that prints output when it catches an arithmetic exception (like divide by zero). If a parallelized loop provokes an exception, the implicit I/O from the handler may cause I/O deadlocks and a system hang.

#### In general:

- The library libF77\_mt is MT safe, but mostly not MT hot.
- You cannot do recursive (nested) I/O if you compile with -mt.

As an informal definition, an interface is MT safe if:

- It can be simultaneously invoked by more than one thread of control.
- The caller is not required to do any explicit synchronization before calling the function.
- The interface is free of data races.

A *data race* occurs when the content of memory is being updated by more than one thread, and that bit of memory is not protected by a lock. The value of that bit of memory is nondeterministic—the two threads *race* to see who gets to update the thread (but in this case, the one who gets there last, wins).

An interface is colloquially called *MT hot* if the implementation has been tuned for performance advantage, using the techniques of multithreading. For some formal definitions of multithreading technology, read the Solaris *Multithreaded Programming Guide*.

## **Cray-Style Parallelization Directives**

Parallel directives have two forms: Sun style and Cray style. The £77 and £90 default is Sun style (-mp=sun). To use Cray-style directives, you must compile with -mp=cray.

Mixing program units compiled with both Sun and Cray directives can produce different results.

A major difference between Sun and Cray directives is that Cray style *requires explicit* scoping of every scalar and array in the loop as either SHARED or PRIVATE.

The following table shows Cray-style directive syntax.

```
!MIC$ DOALL
!MIC$& SHARED( v1, v2, )
!MIC$& PRIVATE( u1, u2, )
...optional qualifiers
```

#### Cray Directive Syntax

A parallel directive consists of one or more *directive lines*. A directive line is defined as follows:

- The directive line is case insensitive.
- The first five characters are CMIC\$, \*MIC\$, or !MIC\$.
- An *initial* directive line has a blank in column 6.
- A *continuation* directive line has a nonblank in column 6.
- Directives are listed in columns 7 and beyond.
- Qualifiers, if any, follow directives—on the same line or continuation lines.
- Multiple qualifiers on a line are separated by commas.
- All variables and arrays are in qualifiers SHARED or PRIVATE.
- Spaces before, after, or within a directive or qualifier are ignored.
- Columns beyond 72 are ignored.

With f90 -free free-format, leading blanks can appear before !MIC\$.

## Qualifiers (Cray Style)

For Cray-style directives, the PRIVATE qualifier is required. Each variable within the DO loop must be qualified as private or shared, and the DO loop index must always be private. The following table summarizes available Cray-style qualifiers.

 TABLE 10-7
 DOALL Qualifiers (Cray Style)

Qualifier	Assertion
SHARED( <i>v1</i> , <i>v2</i> ,)	Share the variables $v1$ , $v2$ , between parallel processes. That is, they are accessible to all the tasks.
PRIVATE( <i>x1</i> , <i>x2</i> ,)	Do not share the variables $x1$ , $x2$ , between parallel processes. That is, each task has its own private copy of these variables.
SAVELAST	Save the values of $private$ variables from the last DO iteration.
MAXCPUS( n )	Use no more than $n$ CPUs.

For Cray-style directives, the <code>DOALL</code> directive allows a single scheduling qualifier, for example, <code>!MIC\$& CHUNKSIZE(100)</code>. Table 10–8 shows the Cray-style DOALL directive

TABLE 10-8 DOALL Cray Scheduling

Qualifier	Assertion	
GUIDED	Distribute the iterations by use of guided self-scheduling.	
	This distribution minimizes synchronization overhead, with acceptable dynamic load balancing.	
SINGLE	Distribute <i>one</i> iteration to each available processor.	
CHUNKSIZE( n	Distribute $n$ iterations to each available processor.	
)	n may be an expression. For best performance, $n$ must be an integer constant. Example: With 100 iterations and <code>CHUNKSIZE(4)</code> , distribute 4 iterations to each CPU.	
NUMCHUNKS( m	If there are $n$ iterations, distribute $n/m$ iterations to each available processor. There can be one smaller residual chunk.	
	m is an expression. For best performance, $m$ must be an integer constant. Example: With 100 iterations and <code>NUMCHUNKS(4)</code> , distribute 25 iterations to each CPU.	

#### scheduling qualifiers:

The  ${\tt f77}$  default scheduling type is the Sun-style  ${\tt STATIC}.$  The  ${\tt f90}$  default is GUIDED.

#### Inhibitors to £90 Explicit Parallelization

With the explicit parallelization situations listed in "Inhibitors to Explicit Parallelization" on page 10-26, the additional parallelization inhibitors for £90 include:

- The DO increment parameter, if specified, is a variable.
- There is an I/O statement in the loop.
- Parallelized loops in subprograms called from parallelized loops are, in fact, not run in parallel.

## **Debugging Parallelized Programs**

Compiling with the -g option cancels any of the parallelization options -autopar, -explicitpar, and -parallel, as well as -reduction and -depend. Some alternative ways to debug parallelized code are suggested in the following section.

## Debugging Without dbx

Debugging parallelized programs requires some cleverness. The following schemes suggest ways to approach the problem:

■ Turn off parallelization.

You can do one of the following:

- Turn off the parallelization options—Verify that the program works correctly by compiling with -03 or -04, but without any parallelization.
- Set the CPUs to one—run the program with the environment variable PARALLEL=1.

If the problem disappears, then you know it was due to parallelization.

Check also for out of bounds array references by compiling with -C.

Problems using -autopar may indicate that the compiler is parallelizing something it should not.

■ Turn off -reduction.

If you are using the -reduction option, summation reduction may be occurring and yielding slightly different answers. Try running without this option.

Reduce the number of compile options.

Compile with just -parallel -03 and check the results.

■ Use fsplit or f90split.

If you have a lot of subroutines in your program, use *fsplit*(1) to break them into separate files. (Use *f90split*(1) on Fortran 90 source codes.) Then compile some files with and without -parallel, and use f77 or f90 to link the .o files. You must specify -parallel on this link step as well. (See *Fortran User's Guide* section on consistent compiling and linking.)

Execute the binary and verify results.

Repeat this process until the problem is narrowed down to one subroutine.

You can proceed using a dummy subroutine or explicit parallelization to track down the loop that causes the problem.

■ Use -loopinfo.

Check which loops are being parallelized and which loops are not.

■ Use a dummy subroutine.

Create a dummy subroutine or function that does nothing. Put calls to this subroutine in a few of the loops that are being parallelized. Recompile and execute. Use <code>-loopinfo</code> to see which loops are being parallelized.

Continue this process until you start getting the correct results.

Then remove the calls from the other loops, compile, and execute to verify that you are getting the correct results.

■ Use explicit parallelization.

Add the C\$PAR DOALL directive to a couple of the loops that are being parallelized. Compile with -explicitpar, then execute and verify the results. Use -loopinfo to see which loops are being parallelized. This method permits the addition of I/O statements to the parallelized loop.

Repeat this process until you find the loop that causes the wrong results.

**Note -** If you need -explicitpar only (without -autopar), do *not* compile with -explicitpar and -depend. This method is the same as compiling with -parallel, which, of course, includes -autopar.

■ Run loops *backward* serially.

Replace DO I=1, N with DO I=N, 1, -1. Different results point to data dependencies.

Avoid using the loop index. It is safer to do so in the loop body, especially if the index is used as an argument in a call.

```
Replace:
    DO I=1,N
    ...
    CALL SNUBBER(I)
```

```
ENDDO

With:

DO I1=1,N
I=I1
...
CALL SNUBBER(I)
...
ENDDO
```

## Using dbx

To use dbx on a parallel loop, temporarily rewrite the program as follows:

- Isolate the body of the loop in a file and subroutine of its own.
- In the original routine, replace loop body with a call to the new subroutine.
- Compile the new subroutine with -g and no parallelization options.
- Compile the changed original routine with parallelization and no -g.

Example: Manually transform a loop to allow using dbx in parallel:

```
Original code:
demo% cat loop.f
C$PAR DOALL
     DO i = 1,10
           WRITE(0,*) "Iteration ", i
      END DO
      END
Split into two parts: caller loop and loop body as a subroutine
demo% cat loop1.f
C$PAR DOALL
     DO i = 1,10
           k = i
            CALL loop_body ( k )
      END DO
      END
demo% cat loop2.f
      SUBROUTINE loop_body ( k )
      WRITE(0,*) "Iteration ", k
     RETURN
Compile caller loop with parallelization but no debugging
demo% f77 -O3 -c -explicitpar loop1.f
Compile the subprogram with debugging but not parallelized
demo% f77 -c -g loop2.f
Link together both parts into a.out
demo% f77 loop1.o loop2.o -explicitpar
Run a.out under dbx and put breakpoint into loop body subroutine
                          Various
demo% dbx a.out
dbx messages not shown
(dbx) stop in loop_body
(2) stop in loop_body
(dbx) run
Running: a.out
```

## **C-Fortran Interface**

This chapter treats issues regarding Fortran and C interoperability.

The discussion is inherently limited to the specifics of the Sun FORTRAN 77, Fortran 90, and C compilers.

**Note -** Material common to both Sun FORTRAN 77 and Fortran 90 is presented in examples that use FORTRAN 77.

## **Compatibility Issues**

Most C-Fortran interfaces must agree in all of these aspects:

- Function/subroutine: definition and call
- Data types: compatibility of types
- Arguments: passing by reference or value
- Arguments: order
- Procedure name: uppercase and lowercase and trailing underscore (\_)
- Libraries: telling the linker to use Fortran libraries

Some C-Fortran interfaces must also agree on:

- Arrays: indexing and order
- File descriptors and stdio
- File permissions

#### **Function or Subroutine?**

The word *function* has different meanings in C and Fortran. Depending on the situation, the choice is important:

- In C, all subprograms are functions; however, some may return a null (void) value.
- In Fortran, a function passes a return value, but a subroutine does not.

When a Fortran routine calls a C function:

- If the called C function returns a value, call it from Fortran as a function.
- If the called C function does not return a value, call it as a subroutine.

When a C function calls a Fortran subprogram:

- If the called Fortran subprogram is a *function*, call it from C as a function that returns a compatible data type.
- If the called Fortran subprogram is a *subroutine*, call it from C as a function that returns a value of int (compatible to Fortran INTEGER\*4) or void. A value is returned if the Fortran subroutine uses alternate returns, in which case it is the value of the expression on the RETURN statement. If no expression appears on the RETURN statement, and alternate returns are declared on SUBROUTINE statement, a zero is returned.

## Data Type Compatibility

The tables below summarize the data sizes and default alignments for FORTRAN 77 and Fortran 90 data types. In both tables, note the following:

- C data types int, long int, and long are equivalent (4 bytes). In a 64-bit environment and compiling with -xarch=v9 or v9a, long and pointers are 8 bytes. This is referred to as "LP64".
- REAL\*16 and COMPLEX\*32, (REAL(KIND=16) and COMPLEX(KIND=16)), are available only on SPARC platforms. In a 64-bit environment and compiling with -xarch=v9 or v9a, alignment is on 16-byte boundaries.
- Alignments marked 4/8 for SPARC indicate that alignment is 8-bytes by default, but on 4-byte boundaries in COMMON blocks. The maximum alignment in COMMON is 4-bytes.
- The elements and fields of arrays and structures must be compatible.
- You cannot pass arrays, character strings, or structures by value.
- You can pass arguments by value from £77 to C, but not from C to £77, since %VAL() is not allowed in a Fortran dummy argument list.

## FORTRAN 77 and C Data Types

Table 11–1 shows the sizes and allowable alignments for FORTRAN 77 data types. It assumes no compilation options affecting alignment or promoting default data sizes are applied. (See also the *FORTRAN 77 Language Reference Manual*).

TABLE 11-1 Data Sizes and Alignments—(in Bytes) Pass by Reference (£77 and cc)

FORTRAN 77 Data Type	C Data Type	Size	Default Alignment SPARC x86	
BYTE X	char x	1	1	1
CHARACTER X	unsigned char x	1	1	1
CHARACTER*n X	unsigned char x[n]	n	1	1
COMPLEX X	struct {float r,i;} x;	8	4	4
COMPLEX*8 X	struct {float r,i;} x;	8	4	4
DOUBLE	struct {double dr,di;}x;	16 16	4/8 4/8	4
COMPLEX X	struct {double dr,di;}x;	32	4/8/16	_
COMPLEX*16 X COMPLEX*32 X	<pre>struct {long double dr,di;} x;</pre>			
DOUBLE	double x	8	4/8	4
PRECISION X	float x	4	4	4
REAL X	float x	4 8	4 4/8	4
REAL*4 X	double x	16	4/8/16	_
REAL*8 X	long double x			
REAL*16 X				_

 TABLE 11–1
 Data Sizes and Alignments—(in Bytes) Pass by Reference (£77 and CC) (continued)

FORTRAN 77 Data Type	C Data Type	Size	Default Alignment SPARC x86	
INTEGER X	int x	4	4	4
INTEGER*2 X	short x	2	2	2
INIEGER"Z X	SHOLL X	4	4	4
INTEGER*4 X	int x	8	4	4
INTEGER*8 X	long long int x			
LOGICAL X	int x	4	4	4
LOGICAL*1 X	char x	1	1	1
LOGICAL I X	Char X	2	2	2
LOGICAL*2 X	short x	4	4	4
LOGICAL*4 X	int x	8	4	4
LOGICAL*8 X	long long int x			

## SPARC: Fortran 90 and C Data Types

The following table similarly compares the Fortran 90 data types with C.

 TABLE 11-2
 Data Sizes and Alignment—(in Bytes) Pass by Reference (£90 and cc)

Fortran 90 Data Type	C Data Type	<b>Sili</b> gnment
CHARACTER x	unsigned char x ;	1
CHARACTER (LEN=n) x	unsigned char $x[n]$ ;	h
COMPLEX x	struct {float r,i;} x;	8

 
 TABLE 11-2
 Data Sizes and Alignment—(in Bytes) Pass by Reference (f90 and cc)
 (continued)

Fortran 90 Data Type	C Data Type	<b>Ali</b> ignment
COMPLEX (KIND=4) x	struct {float r,i;} x;	8
COMPLEX (KIND=8) x	struct {double dr,di;} x;	<b>4</b> 68 <b>3</b> 28/16
COMPLEX (KIND=16) x	<pre>struct {long double, dr,di;} x;</pre>	
DOUBLE COMPLEX x	struct {double dr, di;} x;	468
DOUBLE PRECISION x	double x ;	8
REAL x	float x ;	4
REAL (KIND=4) x	float x ;	4
REAL (KIND=8) x	double x ;	<b>8</b> /8
REAL (KIND=16) x	long double x ;	468/16
INTEGER x	int x ;	4
INTEGER (KIND=1) x	signed char x ;	4
INTEGER (KIND=2) x	short x ;	<b>2</b> 4
INTEGER (KIND=4) x	int x ;	8
INTEGER (KIND=8) x	long long int x;	
LOGICAL x	int x ;	4
LOGICAL (KIND=1) x	signed char x ;	4
LOGICAL (KIND=2) x	short x ;	2
LOGICAL (KIND=4) x	int x ;	4 8
LOGICAL (KIND=8) x	long long int x;	

## Case Sensitivity

C and Fortran take opposite perspectives on case sensitivity:

- C is case sensitive—case matters.
- Fortran ignores case.

The £77 and £90 default is to ignore case by converting subprogram names to lowercase. It converts all uppercase letters to lowercase letters, except within character-string constants.

There are two usual solutions to the uppercase/lowercase problem:

- In the C subprogram, make the name of the C function all lowercase.
- Compile the £77 program with the -U option, which tells £77 to preserve existing uppercase/lowercase distinctions on function/subprogram names.

Use one of these two solutions, but not both.

Most examples in this chapter use all lowercase letters for the name in the C function, and do *not* use the £77 -U compiler option. (£90 does not have an equivalent option.)

#### **Underscores** in Routine Names

The Fortran compiler normally appends an underscore (\_) to the names of subprograms appearing both at entry point definition and in calls. This convention differs from C procedures or external variables with the same user-assigned name. If the name has exactly 32 characters, the underscore is not appended. All Fortran library procedure names have double leading underscores to reduce clashes with user-assigned subroutine names.

There are three usual solutions to the underscore problem:

- In the C function, change the name of the function by appending an underscore to that name.
- Use the £77 C() pragma to tell the FORTRAN 77 compiler to omit those trailing underscores.
- Use the f77 -ext\_names option to make external names without underscores.

Use only one of these solutions.

The examples in this chapter could use the FORTRAN 77 C() compiler pragma to avoid underscores. The C() pragma directive takes the names of external functions as arguments. It specifies that these functions are written in the C language, so the Fortran compiler does not append an underscore as it ordinarily does with external names. The C() directive for a particular function must appear before the first reference to that function. It must also appear in each subprogram that contains such a reference. The conventional usage is:

```
EXTERNAL ABC, XYZ !$PRAGMA C( ABC, XYZ )
```

If you use this pragma, the C function does not need an underscore appended to the function name.

Fortran 90 does not have equivalent methods for avoiding underscores. Trailing underscores are required in the names of C routines called from Fortran 90 routines.

## Argument-Passing by Reference or Value

In general, Fortran routines pass arguments by reference. In a call, if you enclose an argument with the £77 and £90 nonstandard function VAL(), the calling routine passes it by value.

In general, C passes arguments by value. If you precede an argument by the ampersand operator (&), C passes the argument by reference using a pointer. C always passes arrays and character strings by reference.

## **Argument Order**

Except for arguments that are character strings, Fortran and C pass arguments in the same order. However, for every argument of character type, the Fortran routine passes an additional argument giving the length of the string. These are long int quantities in C, passed by value.

The order of arguments is:

- Address for each argument (datum or function)
- A long int for each character argument (the whole list of string lengths comes after the whole list of other arguments)

#### Example:

```
This Fortran code fragment:
Is equivalent to this in C:
    CHARACTER*7 S
    INTEGER B(3)
    ...
    CALL SAM( S, B(2) )
    char s[7];
    long b[3];
    ...
    sam_( s, &b[1], 7L );
```

## Array Indexing and Order

Array indexing and order differ between Fortran and C.

#### **Array Indexing**

C arrays always start at zero, but by default Fortran arrays start at 1. There are two usual ways of approaching indexing.

- You can use the Fortran default, as in the preceding example. Then the Fortran element B(2) is equivalent to the C element b[1].
- You can specify that the Fortran array B starts at B(0) as follows:

```
INTEGER B(0:2)
```

This way, the Fortran element B(1) is equivalent to the C element b[1].

#### **Array Order**

Fortran arrays are stored in column-major order: A(3,2)

```
A(1,1) A(2,1) A(3,1) A(1,2) A(2,2) A(3,2) A(1,3) A(2,3) A(3,3)
```

C arrays in row-major order: A[3][2]

```
A[0][0] A[0][1] A[0][2] A[1][0] A[1][1] A[1][2] A[2][0] A[2][1] A[2][2]
```

For one-dimensional arrays, this is no problem. For two-dimensional and higher arrays, be aware of how subscripts appear and are used in all references and declarations—some adjustments might be necessary.

For example, it may be confusing to do part of a matrix manipulation in C and the rest in Fortran. It might be preferable to pass an *entire* array to a routine in the other language and perform *all* the matrix manipulation in that routine to avoid doing part in C and part in Fortran.

## File Descriptors and stdio

Fortran I/O channels are in terms of unit numbers. The I/O system does not deal with unit numbers but with *file descriptors*. The Fortran runtime system translates from one to the other, so most Fortran programs do not have to recognize file descriptors.

Many C programs use a set of subroutines, called  $standard\ I/O$  (or stdio). Many functions of Fortran I/O use standard I/O, which in turn uses operating system I/O calls. Some of the characteristics of these I/O systems are listed in in the following table.

TABLE 11-3 Comparing I/O Between Fortran and C

	Fortran Units	Standard I/O File Pointers	File Descriptors
Files Open	Opened for reading and writing	Opened for reading; or Opened for writing; or Opened for both; or Opened for appending; See <i>OPEN</i> (3S)	Opened for reading; or Opened for writing; or Opened for both
Attributes	Formatted or unformatted	Always unformatted, but can be read or written with format-interpreting routines	Always unformatted
Access	Direct or sequential	Direct access if the physical file representation is direct access, but can always be read sequentially	Direct access if the physical file representation is direct access, but can always be read sequentially
Structure	Record	Byte stream	Byte stream
Form	Arbitrary nonnegative integers from 0-2147483647	Pointers to structures in the user's address space	Integers from 0-1023

### File Permissions

C programs typically open input files for reading and output files for writing or for reading and writing. A f77 program can OPEN a file READONLY or with READWRITE='READ' or 'WRITE' or 'READWRITE'. f90 supports the READWRITE specifier, but not READONLY.

Fortran tries to open a file with the maximum permissions possible, first for both reading and writing, then for each separately.

This event occurs transparently and is of concern only if you try to perform a READ, WRITE, or ENDFILE operation but you do not have permission. Magnetic tape

operations are an exception to this general freedom, since you can have write permissions on a file, but not have a write ring on the tape.

# Libraries and Linking With the £77 or £90 Command

To link the proper Fortran and C libraries, use the f77 or f90 command to invoke the linker.

Example 1: Use f77 to link:

```
demo% cc -c RetCmplxmain.c
demo% f77 RetCmplx.f RetCmplxmain.o
" This command line does the linking.
demo% a.out
4.0 4.5
8.0 9.0
demo%
```

## Passing Data Arguments by Reference

The standard method for passing data between Fortran routines and C procedures is by reference. To a C procedure, a Fortran subroutine or function call looks like a procedure call with all arguments represented by pointers. The only peculiarity is the way Fortran handles character strings and functions as arguments and as the return value from a CHARACTER\*n function.

## Simple Data Types

For simple data types (not COMPLEX or CHARACTER strings), define or pass each associated argument in the C routine as a pointer:.

TABLE 11-4 Passing Simple Data Types

Fortran calls C	C calls Fortran
integer i	int i=100;
real r	float r;
external CSim	extern void fsim_(int *i, float *r);
i = 100	fsim_(&i, &r);
<pre>call CSim(i,r)</pre>	•••
	subroutine FSim(i,r)
<pre>void csim_(int *i, float *r)</pre>	integer i
{	real r
*r = *i;	r = i
}	return
1	end

#### **COMPLEX Data**

Pass a Fortran COMPLEX data item as a pointer to a C struct of two float or two double data types:

TABLE 11–5 Passing COMPLEX Data Types

Fortran calls C	C calls Fortran
complex w	struct cpx {float r, i;};
double complex z	struct cpx d1;
external CCmplx	struct cpx *w = &d1
call CCmplx(w,z)	struct dpx {double r, i;};
	struct dpx d2;
	struct dpx *z = &d2
<pre>struct cpx {float r, i;};</pre>	fcmplx_( w, z );
<pre>struct dpx {double r,i;};</pre>	
void ccmplx_(	
struct cpx *w,	subroutine FCmplx( w, z )
struct dpx *z)	complex w
{	double complex z
w -> r = 32.;	w = (32., .007)
w -> i = .007;	z = (66.67, 94.1)
z -> r = 66.67;	return
z -> i = 94.1;	end
}	

In 64-bit environments and compiling with  $\mbox{-xarch=v9},$  COMPLEX values are returned in registers.

## **Character Strings**

Passing strings between C and Fortran routines is not recommended because there is no standard interface. However, note the following:

■ All C strings are passed by reference.

■ Fortran calls pass an additional argument for every argument with character type in the argument list. The extra argument gives the length of the string and is equivalent to a C long int passed by value. (This is implementation dependent.) The extra string-length arguments appear after the explicit arguments in the call.

A Fortran call with a character string argument is shown in the next example with its C equivalent:

TABLE 11-6 Passing a CHARACTER string

Fortran call:	C equivalent:
CHARACTER*7 S	char s[7];
INTEGER B(3)	long b[3];
CALL CSTRNG( S, B(2) )	cstrng_( s, &b[1], 7L);

If the length of the string is not needed in the called routine, the extra arguments may be ignored. However, note that Fortran does not automatically terminate strings with the explicit null character that C expects. This must be added by the calling program.

#### **One-Dimensional Arrays**

Array subscripts in C start with 0.

TABLE 11-7 Passing a One-Dimensional Array

Fortran calls C	C calls Fortran
integer i, Sum	extern void vecref_
integer a(9)	( int[], int * );
external FixVec	
•••	int i, sum;
call FixVec ( a, Sum )	int v[9] =
	<pre>vecref_( v, ∑ );</pre>
<pre>void fixvec_ (</pre>	
int $v[9]$ , int *sum )	subroutine VecRef( v, total)
{	integer i, total, v(9)
int i;	total = 0
*sum = 0;	do i = 1,9
for ( i = 0; i <= 8; i++ )	total = total + v(i)
*sum = *sum + v[i];	end do
}	

# **Two-Dimensional Arrays**

Rows and columns between C and Fortran are switched.

TABLE 11-8 Passing a Two-Dimensional Array

Fortran calls C	C calls Fortran
REAL Q(10,20)	extern void
	<pre>qref_( int[][10], int *);</pre>
Q(3,5) = 1.0	
CALL FIXQ(Q)	int m[20][10] =;
	int sum;
<pre>void fixq_( float a[20][10] )</pre>	<pre>qref_( m, ∑ );</pre>
{	
a[5][3] = a[5][3] + 1.;	SUBROUTINE QREF(A,TOTAL)
	INTEGER A(10,20), TOTAL
}	DO I = 1,10
	DO J = 1,20
	TOTAL = TOTAL + A(I,J)
	END DO
	END DO

#### **Structures**

C and FORTRAN 77 structures and Fortran 90 derived types can be passed to each other's routines as long as the corresponding elements are compatible.

TABLE 11-9 Passing FORTRAN 77 STRUCTURE Records

Fortran calls C	C calls Fortran
STRUCTURE /POINT/	struct point {
REAL X, Y, Z	float x,y,z;
END STRUCTURE	};
RECORD / POINT / BASE	<pre>void fflip_ ( struct point *) ;</pre>
EXTERNAL FLIP	
	struct point d;
CALL FLIP( BASE )	struct point *ptx = &d
	<pre>fflip_ (ptx);</pre>
struct point {	
<pre>float x,y,z;</pre>	
};	SUBROUTINE FFLIP(P)
<pre>void flip_( struct point *v; )</pre>	STRUCTURE /POINT/
{	REAL X,Y,Z
float t;	END STRUCTURE
t = v -> x;	RECORD /POINT/ P
v -> x = v -> y;	REAL T
v -> y = t;	T = P.X
$v \rightarrow z = -2.*(v \rightarrow z);$	P.X = P.Y
}	P.Y = T
	P.Z = -2.*P.Z

TABLE 11–10 Passing Fortran 90 Derived Types

Fortran 90 calls C	C calls Fortran 90
TYPE point	struct point {
SEQUENCE	float x,y,z;
REAL :: x, y, z	};
END TYPE point	extern void fflip_ (
TYPE (point) base	struct point *);
EXTERNAL flip	
	struct point d;
CALL flip( base)	struct point *ptx = &d
	fflip_(ptx);
struct point {	
<pre>float x,y,z;</pre>	
};	SUBROUTINE FFLIP( P )
<pre>void flip_( struct point *v; )</pre>	TYPE POINT
{	REAL :: X, Y, Z
float t;	END TYPE POINT
t = v -> x;	TYPE (POINT) P
v -> x = v -> y;	REAL :: T
v -> y = t;	T = P%X
$v \rightarrow z = -2.*(v \rightarrow z);$	P%X = P%Y
}	P%Y = T
	P%Z = -2.*P%Z

#### **Pointers**

A FORTRAN 77 pointer can be passed to a C routine as a pointer to a pointer because the Fortran routine passes arguments by reference.

TABLE 11-11 Passing a FORTRAN 77 POINTER

Fortran calls C	C calls Fortran
REAL X	extern void fpass_;
POINTER (P2X, X)	
EXTERNAL PASS	float *x;
P2X = MALLOC(4)	float **p2x;
X = 0.	fpass_(p2x);
CALL PASS(X)	
	SUBROUTINE FPASS (P2X)
<pre>void pass_(x)</pre>	REAL X
int **x;	POINTER (P2X, X)
{	X = 0.
**x = 100.1;	
}	

C pointers are compatible with Fortran 90 (release 2.0) *scalar* pointers, but not *array* pointers.

# Passing Data Arguments by Value

Call by value is available only for simple data with FORTRAN 77, and only by Fortran routines calling C routines. There is no way for a C routine to call a Fortran routine and pass arguments by value. It is not possible to pass arrays, character strings, or structures by value. These are best passed by reference.

Use the nonstandard Fortran function %VAL(arg) as an argument in the call.

In the following example, the Fortran routine passes x by value and y by reference. The C routine incremented both x and y, but only y is changed.

#### Fortran 77 calls C

```
REAL x, y
x = 1.
y = 0.
PRINT *, x,y
CALL value( %VAL(x), y)
PRINT *, x,y
END
void value_( float x, float *y)
printf("%f, %f\n",x,*y);
x = x + 1.;
y = y + 1.;
printf("%f, %f\n",x,*y);
Compiling and running produces output:
1.00000 0. x and y from Fortran
1.000000, 0.000000 x and y from C
2.000000, 1.000000 new x and y from C
1.00000 1.00000 new x and y from Fortran
```

## Functions That Return a Value

A Fortran function that returns a value of type BYTE (*FORTRAN 77 only*), INTEGER, REAL, LOGICAL, DOUBLE PRECISION, or REAL\*16 (*SPARC only*) is equivalent to a C function that returns a compatible type (see Table 11–1 and Table 11–2). There are two extra arguments for the return values of character functions, and one extra argument for the return values of complex functions.

#### Returning a Simple Data Type

The following example returns a REAL or float value. BYTE, INTEGER, LOGICAL, DOUBLE PRECISION, and REAL\*16 are treated in a similar way:

TABLE 11-13 Functions Returning a REAL or float Value

Fortran calls C	C calls Fortran
real ADD1, R, S	float r, s;
external ADD1	extern float fadd1_();
R = 8.0	r = 8.0;
S = ADD1(R)	s = fadd1_( &r );
•••	
float add1_( pf )	real function fadd1 (p)
float *pf;	real p
{	fadd1 = p + 1.0
float f ;	return
f = *pf;	end
f++;	
return ( f );	
}	

## **Returning COMPLEX Data**

A Fortran function returning COMPLEX or DOUBLE COMPLEX is equivalent to a C function with an additional first argument that points to the return value in memory. The general pattern for the Fortran function and its corresponding C function is:

Fortran function	C function
COMPLEX FUNCTION CF(a1, a2,, an)	cf_ () return, a1, a2,, an
	struct { float r, i; } *;return

TABLE 11–14 Function Returning COMPLEX Data

Fortran calls C	C calls Fortran
COMPLEX U, V, RETCPX	struct complex { float r, i; };
EXTERNAL RETCPX	struct complex c1, c2;
U = ( 7.0, -8.0)	struct complex *u=&c1, *v=&c2
V = RETCPX(U)	extern retfpx_();
	u -> r = 7.0;
	u -> i = -8.0;
<pre>struct complex { float r, i; };</pre>	retfpx_( v, u );
<pre>void retcpx_( temp, w )</pre>	
struct complex *temp, *w;	
{	COMPLEX FUNCTION RETFPX(Z)
temp->r = $w->r + 1.0$ ;	COMPLEX Z
temp->i = w->i + 1.0;	RETFPX = Z + (1.0, 1.0)
return;	RETURN
}	END

In 64-bit environments and compiling with -xarch=v9, COMPLEX values are returned in floating-point registers: COMPLEX and DOUBLE COMPLEX in %f0 and %f1, and COMPLEX\*32 in %f0, %f1, %f2, and %f3.

#### Returning a CHARACTER String

Passing strings between C and Fortran routines is not encouraged. However, a Fortran character-string-valued function is equivalent to a C function with two additional first arguments—data address and string length. The general pattern for the Fortran function and its corresponding C function is:

Fortran function	C function
CHARACTER*n FUNCTION C(a1,, an)	void c_ (result, length, a1,, an)
	char result[];
	long length;

Here is an example:

TABLE 11-15 A Function Returning a CHARACTER String

Fortran calls C	C calls Fortran
CHARACTER STRING*16, CSTR*9	void fstr_( char *, int,
STRING = ''	char *, int *, int );
STRING = '123' // CSTR('*',9)	char sbf[9] = "123456789";
	char *p2rslt = sbf;
	<pre>int rslt_len = sizeof(sbf);</pre>
<pre>void cstr_( char *p2rslt,</pre>	char ch = '*';
int rslt_len,	int n = 4;
char *p2arg,	<pre>int ch_len = sizeof(ch);</pre>
int *p2n,	/* make n copies of ch in sbf
<pre>int arg_len )</pre>	*/
{ /* return n copies of arg */	fstr_( p2rslt, rslt_len,
int count, i;	&ch, &n, ch_len);
char *cp;	
count = *p2n;	
cp = p2rslt;	FUNCTION FSTR( C, N)
for (i=0; i <count; i++)="" td="" {<=""><td>CHARACTER FSTR*(*), C</td></count;>	CHARACTER FSTR*(*), C
*cp++ = *p2arg ;	FSTR = ''
}	DO I = 1,N
}	FSTR(I:I) = C
	END DO
	FSTR(N+1:N+1) = CHAR(0)
	END

In this example, the C function and calling C routine must accommodate two initial extra arguments (a pointer to the result string and the length of the string) and one

additional argument at the end of the list (length of character argument). Note that in the Fortran routine called from C, it is necessary to explicitly add a final null character.

# Labeled COMMON

Fortran labeled COMMON can be emulated in C by using a global struct.

TABLE 11-16 Emulating Labeled COMMON

Fortran COMMON Definition	C "COMMON" Definition
COMMON /BLOCK/ ALPHA,NUM	extern struct block {
	float alpha;
	int num;
	};
	extern struct block block_;
	main ()
	{
	blockalpha = 32.;
	blocknum += 1;
	}

Note that the external name established by the C routine must end in an underscore to link with the block created by the Fortran program. Note also that the C directive #pragma pack may be needed to get the same padding as with Fortran. Both £77 and £90 align data in COMMON blocks to at most 4-byte boundaries.

## Sharing I/O Between Fortran and C

Mixing Fortran I/O with C I/O (issuing I/O calls from both C and Fortran routines) is not recommended. It is better to do *all* Fortran I/O or *all* C I/O, not both.

The Fortran I/O library is implemented largely on top of the C standard I/O library. Every open unit in a Fortran program has an associated standard I/O file structure. For the stdin, stdout, and stderr streams, the file structure need not be explicitly referenced, so it is possible to share them.

If a Fortran main program calls C to do I/O, the Fortran I/O library must be initialized at program startup to connect units 0, 5, and 6 to stderr, stdin, and stdout, respectively. The C function must take the Fortran I/O environment into consideration to perform I/O on open file descriptors.

However, if a C main program calls a Fortran subprogram to do I/O, the automatic initialization of the Fortran I/O library to connect units 0, 5, and 6 to stderr, stdin, and stdout is lacking. This connection is normally made by a Fortran main program. If a Fortran function attempts to reference the stderr stream (unit 0) without the normal Fortran main program I/O initialization, output will be written to fort.0 instead of to the stderr stream.

The C main program can initialize Fortran I/O and establish the preconnection of units 0, 5, and 6 by calling the  $f_{init}()$  FORTRAN 77 library routine at the start of the program and, optionally,  $f_{exit}()$  at termination.

Remember: even though the main program is in C, you should link with £77.

#### Alternate Returns

Fortran's alternate returns mechanism is obsolescent and should not be used if portability is an issue. There is no equivalent in C to alternate returns, so the only concern would be for a C routine calling a Fortran routine with alternate returns.

The Sun Fortran implementation returns the int value of the expression on the RETURN statement. This is implementation dependent and its use should be avoided.

TABLE 11-17 Alternate Returns

C calls Fortran	Running the Example
<pre>int altret_ ( int * );</pre>	demo% cc -c tst.c
main ()	demo% f77 -o alt alt.f tst.o
{	alt.f:
int k, m;	altret:
k =0;	demo% <b>alt</b>
<pre>m = altret_( &amp;k );</pre>	1 2
<pre>printf( "%d %d\n", k, m); }</pre>	The C routine receives the return value 2 from the Fortran routine because it executed the
SUBROUTINE ALTRET( I, *, *)	RETURN 2 statement.
INTEGER I	
I = I + 1	
IF(I .EQ. 0) RETURN 1	
IF(I .GT. 0) RETURN 2	
RETURN	
END	

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