<table>
<thead>
<tr>
<th>Version</th>
<th>Description</th>
</tr>
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<tr>
<td>3.5</td>
<td>April 2001</td>
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<td></td>
<td>Original Printing.</td>
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Determining how much memory is available on your system

I/O bound

## Procedure 4: Determining if the code is I/O bound

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Preface

This publication documents techniques you can use to optimize Fortran 90, C++, or C code on Cray SV1 systems. This manual discusses using the Cray CF90 compiler, the Cray Standard C compiler, the Cray C++ compiler, and various performance tools to help you analyze and optimize your code.

This is a guide for programmers with working knowledge of the CF90, Cray Standard C, or Cray C++ compilers.

Related Publications

The following documents contain additional information that may be helpful:

• CF90 Commands and Directives Reference Manual
• Cray Standard C and Cray C++ Reference Manual
• Application Programmer’s Library Reference Manual
• Intrinsic Procedures Reference Manual
• Scientific Library Reference Manual

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Customers who subscribe to the CRInform program can order software release packages electronically by using the Order Cray Software option.

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Conventions

The following conventions are used throughout this document:

<table>
<thead>
<tr>
<th>Convention</th>
<th>Meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td>command</td>
<td>This fixed-space font denotes literal items (such as commands, files, routines, pathnames, signals, messages, programming language structures, and e-mail addresses) and items that appear on the screen.</td>
</tr>
<tr>
<td>variable</td>
<td>Italic typeface denotes variable entries and words or concepts being defined.</td>
</tr>
<tr>
<td>user input</td>
<td>This bold, fixed-space font denotes literal items that the user enters in interactive sessions. Output is shown in nonbold, fixed-space font.</td>
</tr>
<tr>
<td>[ ]</td>
<td>Brackets enclose optional portions of a command or directive line.</td>
</tr>
<tr>
<td>...</td>
<td>Ellipses indicate that a preceding element can be repeated.</td>
</tr>
</tbody>
</table>

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We value your comments and will respond to them promptly.
This document provides information about techniques you can use to optimize Fortran, C, or C++ code on Cray SV1 systems. This is a guide for programmers with working knowledge of either the UNICOS or UNIX operating system and experience in working with CF90, Cray C++, or Cray Standard C compilers.

This document makes the following assumptions about code to be optimized:

- The code has been debugged and is running on a Cray SV1 system.
- When running the code, you have used the best compiler options and data types for your program. For example, you have used aggressive optimization options (-O scalar3,vector3 for the Fortran f90 command or -hscalar3,vector3 for the C or C++ cc or CC command) and, if you do not need 96 bits of floating-point precision, you have used single precision (f90 -dp or the float or double type specifier instead of long double type specifier within C++ code).

**Note:** Unless otherwise noted, all discussions of C++ code optimizations also apply to the C language.

### 1.1 About this manual

The following information is provided in this manual:

- Section 1.2, page 2, gives an overview of the optimization process.
- Section 1.3, page 4, provides an overview of the Cray SV1 hardware.
- Chapter 2, page 11, describes how to evaluate code and determine where to focus optimization efforts.
- Chapter 3, page 25, describes multi-streaming for Cray SV1 systems.
- Chapter 4, page 39, describes vectorization.
- Chapter 5, page 47, describes how to optimize memory use.
- Chapter 6, page 67, describes how to optimize input and output.
1.2 Optimization overview

Optimization is the process of changing a program or the environment in which it runs to improve its performance. Performance gains generally fall into one of two categories of measured time:

- **User CPU time.** Time accumulated by a user process when it is attached to a CPU and executing. When running on a single CPU, CPU time is a fraction of elapsed time. When multitasked, CPU time is a multiple of elapsed time.

- **Elapsed (wall-clock) time.** The amount of time that passes between the start and termination of a user process. Elapsed time includes the following:
  - User CPU time
  - UNICOS system CPU time
  - I/O wait time
  - Sleep or idle time

The flowchart in Figure 1, page 3, shows a comprehensive view of the CPU optimization process described in this document. This is an iterative method that requires you to determine when to stop optimizing the code.

**Note:** During the optimization process you will need to execute your code repeatedly to assess performance and measure performance gain. To save execution time, we encourage you to work with smaller, sample data sets that exercise all of the code within your program.

Each area of optimization involves a three-step method, which includes the following general steps:

1. Identify a problem.
2. Evaluate the problem.
3. Apply a technique.

Check your answers after each code modification to avoid regression, then compare the new performance against the initial performance to decide where to proceed. If you want to continue optimizing for CPU performance, return to the initial analysis.

**Note:** If none of these techniques brings you closer to the desired computation rates for the CPU for your system, you may want to seek help from Cray. See your local account manager or service representative for more information.
Introduction

Sufficiently optimized?

Yes

No

- Identify where execution time is being spent
- Determine whether you can make the code run faster
- Apply optimization technique

- Identify files using large amounts of processing time
- Evaluate I/O alternatives
- Restructure I/O

Evaluate the code

- Identify large amounts of memory wait or system CPU time
- Evaluate dynamic memory alternatives
- Restructure code or insert loader directives

Check answers

- Sufficiently optimized?

Program is optimized for single-CPU performance. Can the program benefit from Autotasking?

Figure 1. Optimization overview
1.3 Hardware overview

The Cray SV1 computer is significantly different from previous Cray vector machines in that it provides a cache for the data resulting from scalar, vector, and instruction buffer memory references. Like its predecessors, the Cray SV1 system achieves high bandwidth to memory for both unit and non-unit stride memory references.

The Cray SV1 system is configured with between 4 and 32 CPUs. Each CPU has 2 add and 2 multiply functional units, allowing them to deliver 4 floating point results per CPU clock cycle. With the 300 MHz CPU clock, the peak floating point rate per CPU is 1.2 Gflops and 38.4 Gflops for the system.

The memory architecture is uniform access, shared central memory. Uniform memory access (UMA) means that the access time for any CPU memory reference to any location in memory is the same. Memory capacity for the system ranges from a minimum of 4 GBytes up to a maximum of 32 GBytes.

The Cray SV1 system has two module types: processor and memory. The system must be configured with eight memory modules and one to eight processor modules. Each processor module has four CPUs. A Cray J90 processor module can be upgraded with a Cray SV1 processor module and the Cray J90 system can be configured with both processor module types, Cray J90 or Cray SV1.

The following figure shows the block diagram for a single processor:
Control and/or data from other processors.
The second vector logical functional unit and the vector integer multiply functional unit share hardware with the floating-point multiply functional unit.
The vector pop/parity/leading zero functional unit shares hardware with the floating-point reciprocal approximation functional unit.
The bit matrix multiply unit shares hardware with the floating-point add functional unit.

Figure 2. Block Diagram
1.3.1 The processor

The Cray SV1 processor uses a custom CMOS chip. The processor is implemented using two chip types: cpu and cache.

The cpu chip contains the vector and scalar units. Scalar registers, scalar functional units, and the instruction buffers reside in the scalar unit, while the vector unit contains vector registers and the vector functional units. As in previous Cray vector systems, the processor contains eight vector (V) registers, eight scalar (S) registers backed by 64 T registers, and eight address (A) registers backed up by 64 B registers. A parallel job also has access to eight shared B and eight shared T registers, which are used for low overhead data passing and synchronization between processors.

A vector functional unit contains two pipes, each capable of producing a result every CPU clock cycle. This results in a peak rate for a functional unit of two results per clock cycle. The maximum vector length, or VL, is 64. The combine floating point functional units, add and multiply, deliver 4 results per CPU clock cycle. With the CPU clock rate of 300 MHz a peak, a floating point rate of 1.2 Gflops for the processor is achieved.

In addition to the add and multiply units, the other vector functional units are reciprocal, integer add, shift, pop/parity/leading zero, bit matrix multiply, and logical.

The vector units are capable of full chaining and tailgating. Chaining is reading from a V register that is still being written to, and tailgating is writing to a V register that is still being read from a prior vector instruction. Scalar floating point operations are executed using the vector functional units. This is different from the Cray J90 system, which has separate floating point functional units for scalar operations.

Two data paths or ports are provided to move data between CPU registers and memory via cache. In any given clock cycle, two memory requests can be active and consist of two dual-port reads or one dual-port read and one dual-port write. If there are no read requests, there can be only one write request active. The processor can access up to 32 Gbytes of memory, but an application is limited to a 16-Gbyte address space.

Instructions are decoded by the scalar unit; when vector instructions are encountered, they are dispatched to the vector unit, which maintains an independent instruction queue. Barring instruction dependency, the two units can execute independently.
There are 32 performance counters in four groups of eight each. Only one group can be active at a time, with software providing user access to the data. The groups are labeled 0 through 3. The collection order, based on how useful the performance information is to the user, is 0, 3, 2 and 1. For an example of using the hardware performance monitor, see Section 2.1.1, page 13.

Note: In addition to the CPU clock, there is a system clock that runs at the rate of 100 MHz. When using the CPU instruction to return the count of clock ticks, the tick count is generated by the system clock rate.

1.3.2 Cache

Cache lies between memory and a processor’s registers (see Figure 3, page 8). Its purpose is to speed up loads from memory.

Sacrificing the performance improvements to be gained from cache can eat into the performance gains you may be realizing with other optimizations. To know how to get the best out of both, you must first know something about how cache works.

Data read from memory is accessed through a high-speed, 32-Kword cache. Each of the processors involved in multi-streaming has its own cache.

Cache is a four-way set associative temporary storage area, meaning each line in cache is divided into four places, or sets. See the representation of cache in Figure 3, page 8.

Moving data from cache to a register is 2-to-3 times faster than moving data directly from memory to a register (ideally, 32 clock periods (CPs) compared to about 102 CPs). Having the data items you are going to use available in cache can represent a significant optimization in itself.

The following example shows the advantage of cache, how it helps move data quickly between memory and processor registers. The data being read in the example is from an array named A that is accessed as follows in a Fortran program:

```fortran
DO I = 1, N
  ... = A(I)
ENDDO
```

A single-streaming processor (SSP) in a multi-streaming program would be assigned its own part of the array. Each would use the following procedure to move data from memory to its registers:
Procedure 1: Moving data from memory

1. A register requests the value of $A(1)$ from cache.

2. Cache does not have $A(1)$. It requests $A(1)$ from memory.

3. Cache receives 8 64-bit words ($A(1)$ through $A(8)$) from memory and stores them in one of four sets.

4. The register receives $A(1)$ from cache. The state of the data at this point is as illustrated in the following figure:

![Figure 3. Using Cache](image_url)
5. When the register needs \( A(2) \) through \( A(8) \), it finds them in cache.

6. Meanwhile, cache is prefetching \( A(9) \) through \( A(16) \) from memory.

7. By the time the register needs \( A(9) \), it is again available in cache.

In this way, a constant flow of data is set up between memory and the processor registers. Whenever an array item is needed, it will be available in cache.

On non-cached Cray vector systems, performance on vector constructs generally increased as a predictable function of vector length. This is not always the case on the Cray SV1 system, since long vectors can lead to a reduction in data cache efficiency. In general, it is better to use blocked algorithms (similar to those commonly used on microprocessors), balancing the vector length against any potential data reuse that can be exploited via data cache.

### 1.3.3 I/O

Disk drives, interfaces to other networks, and other peripherals are connected to the Cray SV1 system using the high-speed GigaRing I/O system. The double-ring product is illustrated in the following illustration.

![GigaRing I/O](image)

Figure 4. GigaRing I/O

Each of the rings has a maximum transfer rate of 500 Mbytes/s, which provides an effective total bandwidth of 800 Mbytes/s. Since the two rings rotate in different directions, the shortest path to the target node is selected for each transfer.
The first step in optimizing a program is evaluating its overall performance. This allows you to decide where to focus optimization efforts.

When you compile and execute a program on a Cray SV1 system, it usually will be dominated by one of the following general activities:

- CPU computation (see Section 2.1, page 11).
- Memory management (see Section 2.2, page 19).
- Input/output (I/O) processing (see Section 2.3, page 23).

A program is considered to be memory bound, I/O bound, or CPU bound because its performance is limited by the dominant activity. Use the job accounting (ja) and hardware performance monitor (hpm) tools as directed in this chapter to identify where in your code to obtain the greatest potential performance gain.

Figure 5, page 12, summarizes the recommended initial analysis.

2.1 CPU-bound programs

A CPU-bound code spends most of its elapsed-time performing CPU calculations. The hpm(1) report will tell you how effectively the code is using vector registers, instruction buffers, and memory ports.

The following sections provide information and steps for determining CPU-related code bottlenecks. Section 2.1.1, page 13, describes the types of reports that you can generate with the hpm(1) command. Procedure 2, page 13, describes the use of the hpm(1) command to determine whether code is performing at peak levels on Cray SV1 systems.
Begin evaluating code using ja and hpm

Memory bound?

I/O bound?

CPU bound

Low computational intensity?

Low MFLOPS and MIPS?

High fetch rate?

Done with single-CPU optimization. Can you benefit from Autotasking?

Memory management

I/O optimization

CPU optimization

Use ja report to determine:
- Excessive unlocked wait time?
- Efficient I/O requests?
- Expected I/O transfer rate?

Figure 5. Evaluating code
2.1.1 Using the hpm Command to Issue Reports

The hpm(1) command allows you to access the hardware performance monitor (HPM) and obtain overall program timing information. The hpm tool can issue any one of the following four types of reports:

- Program summary (HPM counter group 0)
- Hold-issue conditions (HPM counter group 1)
- Memory use (HPM counter group 2)
- Vectorization (HPM counter group 3)

For complete information on the hpm utility, see the hpm(1) man page or the Guide to Parallel Vector Applications.

Procedure 2: Using the hpm command to determine CPU performance statistics

Use the following procedure to access the code’s CPU performance statistics and to determine whether the code is optimized for single-CPU speed:

1. Produce an hpm report by first compiling the program, then issuing the hpm command with the program name, as shown in the following example:

   ```
   f90 yourcode.f90 or CC yourcode.C
   hpm .a.out
   ```

   By default, this hpm command generates a program summary (group 0) report, the most commonly run report. Also, by default, the hpm command writes report-style output to the stderr file, allowing the program’s default output to appear on the screen.

   The following is a sample hpm group 0 report.

   ```
   Group 0: CPU seconds : 31.74961 CP executing : 952488354
   Million inst/sec (MIPS) : 42.72 Instructions : 1356230656
   Avg. clock periods/inst : 7.02 CP holding issue : 7743041828
   % CP holding issue : 81.29 Inst.buf. fetches : 7680895
   Inst.buffer fetches/sec : 0.24M F.P. adds : 2491024131
   Floating adds/sec : 78.46M F.P. multiplies : 2630592592
   Floating multiplies/sec : 82.85M F.P. reciprocals : 10201
   Cache hits/sec : 92.75M Cache hits : 2944803625
   Floating reciprocal/sec : 0.00M CPU mem. references : 4436364260
   ```
2. Determine whether the code is dominated by scalar or vector operations. To do this, use the values shown in the hpm report to perform the following steps:

   a. Find the millions of instructions per second (MIPS) for the code by looking at the Million inst/sec (MIPS) row of the hpm report.

   b. Find the floating point operations per second (FLOPS) for the code by looking at the Floating ops/CPU second row of the hpm report.

   c. Use the following table to determine whether the MIPS and FLOPS for the code are low or high, based on the hardware expectations.

      The code is not likely to achieve peak FLOPS or MIPS rates, but some codes are capable of performing at substantial fractions of peak speed.

      Low FLOPS is any single-digit rate up to 20% of peak rates for that system. High FLOPS are generally anything above 66% of peak rates for that system.

      Low MIPS rates are at or near the bottom of the range for that system. High MIPS rates are generally anything near 50% of peak MIPS for that system.
Table 1. Single-CPU hardware expectations

<table>
<thead>
<tr>
<th></th>
<th>Cray SV1</th>
<th>Cray SV1e</th>
</tr>
</thead>
<tbody>
<tr>
<td>Peak FLOPS</td>
<td>1200M</td>
<td>1800M</td>
</tr>
<tr>
<td>MIPS range</td>
<td>30 to 300</td>
<td>30 to 450</td>
</tr>
</tbody>
</table>

d. Use the following table to determine whether the code is dominated by scalar or vector operations.

Table 2. Determining if the code is dominated by scalar or vector operations

<table>
<thead>
<tr>
<th>FLOPS</th>
<th>MIPS</th>
<th>Determination</th>
</tr>
</thead>
<tbody>
<tr>
<td>High or Medium</td>
<td>Low</td>
<td>Code is most likely dominated by vector operations. The code is performing well.</td>
</tr>
<tr>
<td>Medium</td>
<td>Medium</td>
<td>Code is most likely a mix of scalar and vector operations. CPU optimization might help improve its CPU performance.</td>
</tr>
<tr>
<td>Medium or Low</td>
<td>High</td>
<td>Code is most likely dominated by scalar operations (is not vector code). CPU optimization will help improve its CPU performance.</td>
</tr>
<tr>
<td>Low</td>
<td>Medium</td>
<td>Code is dominated by scalar operations. CPU optimization will help improve its CPU performance.</td>
</tr>
<tr>
<td>Low</td>
<td>Low</td>
<td>Code is performing poorly, whether it is scalar or vector. It has a CPU performance problem. CPU optimization will help improve its CPU performance.</td>
</tr>
</tbody>
</table>

3. Determine if the instruction buffer fetches per second (as shown by the Inst.buffer fetches/sec row of the hpm report) are close to or greater than 0.1 million per second. If yes, see the following note on analyzing the CPU-bound code. If no, go to the next step.
Note: Excessive instruction buffer fetches tend to slow down the code. If the code has a high rate (approaching 0.1 million per second), the code may have excessive jumping (as with go to or if-then-else constructs) or excessive calls to subprograms. CPU optimization probably will help the overall performance.

4. Determine the computational intensity ratio of the code. Use the values in the hpm report, as follows:

   a. Divide the number in the Floating ops/CPU second row by the number in the CPU mem. references/sec row. This is the computational intensity of the code.

   The computational intensity ratio is the ratio of the floating-point operation rate to the memory access rate. This ratio should reflect the floating-point operations in the code (for example, $a=b+c$ has 1 Floating ops/CPU second and 3 CPU mem. references/sec, or a computational intensity of 0.33). Any ratio less than 1/3 makes excessive use of the CPU memory ports while the remainder of the processor idles. This is usually caused by memory-to-memory traffic ($a=b$), highly scalar code, or a hidden performance problem.

   b. If the computational intensity ratio of the code is less than 1/3, the program is not making effective use of the CPU. If the computational intensity ratio of the code is 1/3 or more, go to the next step.

5. If you have already determined that the code is not memory or I/O bound, the code is probably optimized for single-CPU performance. However, if your program still does not achieve the performance expected from your system after you have exhausted all optimization techniques, Cray can offer assistance for further optimization on a fee basis. See your local account manager or service representative for more information.

6. If you have completed single-CPU optimization of your program and are satisfied with its single-CPU performance, you might want to run your program in parallel on multiple CPUs. The multi-streaming processor (MSP) feature is the most convenient method to achieve parallel execution. To determine whether your program can benefit from multi-streaming, go to Chapter 3, page 25.

2.1.2 Analyzing multi-streaming code

The following other tools are available to analyze multi-streaming code:

- The prof(1) and profview(1) commands (see Section 3.6.1, page 33).
• The MSP_STATS environment variable (see Section 3.6.2, page 36).
• The ps(1) command (see Section 2.1.2.1, page 17).
• The jstat(1) command (see Section 2.1.2.2, page 18).
• The cpu(8) command (see Section 2.1.2.3, page 18).
• The sysconf(1) (see Section 2.2.2.1, page 22).
• The ja(1) (see Section 2.2.2, page 19).

2.1.2.1 The ps(1) command

The –l option of the ps command reports the process connection state in the WCHAN field of its output.

If you are running a single-processor job, an SSP process connected to CPU 5 would have cpu-05 in the WCHAN field. When running on an MSP, ps displays both the MSP number and the number of each of the SSPs involved. The following convention applies:

msp.n.i

n: MSP number.

i: SSP number.

For example, ps –el | grep msptest, while multi-streaming the executable program msptest, produces output similar to the following:

```
F S UID PID PPID C PRI NI ADDR SZ WCHAN TTY TIME CMD
101 R 1330 4711 4689 0 996 24 250164 8504 msp0.2 p002 0:07 msptest
101 R 1330 4709 4689 0 996 24 250164 8504 msp0.0 p002 0:07 msptest
101 R 1330 4712 4689 0 996 24 250164 8504 msp0.1 p002 0:07 msptest
101 R 1330 4710 4689 0 995 24 250164 8504 msp0.3 p002 0:07 msptest
```

Output for a multitasking and multi-streaming program running on two MSPs is as follows:

```
F S UID PID PPID C PRI NI ADDR SZ WCHAN TTY TIME CMD
101 R 0 2246 2203 0 995 24 264064 16504 msp0.1 p003 0:17 msptest
101 R 0 2245 2203 0 996 24 264064 16504 msp0.0 p003 0:17 msptest
101 R 0 2247 2203 0 995 24 264064 16504 msp0.2 p003 0:17 msptest
101 R 0 2248 2203 0 995 24 264064 16504 msp0.3 p003 0:17 msptest
101 R 0 2249 2203 0 996 24 264064 16504 msp1.0 p003 0:17 msptest
101 R 0 2250 2203 0 996 24 264064 16504 msp1.1 p003 0:17 msptest
```
2.1.2.2 The jstat(1) command

The jstat command displays information pertaining to active jobs. Using the -j option, you can view the CPU connection state of all processes within a job. The state field of the jstat -j id output identifies the MSP and SSP on which each executing process within the job is running. For example:

<table>
<thead>
<tr>
<th>pid</th>
<th>status</th>
<th>state</th>
<th>utime</th>
<th>stime</th>
<th>size</th>
<th>addr</th>
<th>system call</th>
<th>command</th>
</tr>
</thead>
<tbody>
<tr>
<td>6647</td>
<td>running</td>
<td>msp0.1</td>
<td>117</td>
<td>0</td>
<td>16504</td>
<td>90116</td>
<td>nosys</td>
<td>msptest</td>
</tr>
<tr>
<td>6652</td>
<td>running</td>
<td>msp0.0</td>
<td>117</td>
<td>0</td>
<td>16504</td>
<td>90116</td>
<td>nosys</td>
<td>msptest</td>
</tr>
<tr>
<td>6651</td>
<td>running</td>
<td>msp0.2</td>
<td>117</td>
<td>0</td>
<td>16504</td>
<td>90116</td>
<td>nosys</td>
<td>msptest</td>
</tr>
<tr>
<td>6649</td>
<td>running</td>
<td>msp0.3</td>
<td>117</td>
<td>0</td>
<td>16504</td>
<td>90116</td>
<td>nosys</td>
<td>msptest</td>
</tr>
</tbody>
</table>

The order of the SSPs is random.

2.1.2.3 The cpu(8) command

The -i option of the /etc/cpu command displays information for each configured CPU. Among the items displayed is an indicator identifying participation in an MSP. The following example shows output for a 12 CPU machine with one MSP configured:

<table>
<thead>
<tr>
<th>cpu</th>
<th>type</th>
<th>state</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>1</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>2</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>3</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>4</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>5</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>6</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>7</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>8</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>9</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>10</td>
<td>svl</td>
<td>up</td>
</tr>
<tr>
<td>11</td>
<td>svl</td>
<td>up</td>
</tr>
</tbody>
</table>
2.2 Memory and cache

There are two ways of optimizing memory on the Cray SV1 system:

- By getting the most out of cache. Cache is a high-speed memory located between main memory and each MSP. The time it takes a register to access data from cache is two or three times faster than the time it takes to access data from memory. See Section 2.2.1, page 19.

- By minimizing the size of your program in memory. When the compiled program begins execution, it becomes a UNICOS user process in memory. The process resides in memory during execution and it interacts with the operating system. Its size might expand or decrease during execution. If the process spends excessive elapsed time managing memory, it is considered a memory-bound process. See Section 2.2.2, page 19.

2.2.1 Cache

(Need help here. How can a user measure things like cache hits versus cache misses?)

2.2.2 Code size

A general rule for the size of a code’s process in memory is “smaller is better.” The advantages of a smaller process are reductions in swap frequency, wait-time, and time-to-load. Although there are that provide exceptions to this rule, for now you need only to identify whether your process is memory bound.

The ja utility reports job or session-related accounting information provided by the job accounting daemon. It can provide you with a report that contains an overall view of memory usage, I/O, and system overhead. This report can help you determine whether your code is CPU bound, memory bound, or I/O bound.

For more information on the job accounting tool, see the ja(1) man page.

Procedure 3: Determining if the code is memory bound

To determine if the code is memory bound, perform the following steps:

1. To obtain a report that shows an overall view of memory usage, I/O, and system overhead, run the job accounting utility, ja -c lth (single-tasked), with the code, as shown in the following example. (To save an extra step, you can also run ja(1) concurrently with hpm(1), the results of which you will use later).

   f90 -o bigio bigio.f90
Compiles and loads the code

`ja`

Turns on job accounting

`.bigio`

Executes the code as a UNICOS user process

`ja -clth > ja.rpt`

Stores a 132-column report in `ja.rpt` file

The following shows a sample `ja` report.

```
cut -c1-72 ja.rpt
```

```
Job Accounting - Command Report
===============================
Command Started Elapsed User CPU Sys CPU I/O Wait I/O Wait
Name At Seconds Seconds Seconds Sec Lck Sec Unlock
ja 11:29:50 0.1906 0.0014 0.0137 0.18 0.0000
```

```
cut -c1-9,73-132 ja.rpt
```

```
Job Accounting - Command Report
Command CPU MEM I/O WMem Kwords Log I/O Memory Ex
Name Avg Mwds Avg Mwds Xferred Request HiWater St Ni Fl SBU'
ja 0.3556 0.4297 0.00 0 880 0 20 0
```

20
Note: To ensure a representative reading, the user process you are measuring with the ja utility should have an execution time that is greater than 1 second.

2. Find the Memory HiWater column of the ja report. Perform the following steps to determine if the memory high-water mark for the process is a significant fraction of available user memory.

Note: A significant fraction of available user memory on your system is a subjective measure based on the chances of the code finding room in memory as it competes with other users’ jobs. Whether the code finds room in memory depends on how many jobs are competing for memory, the job size, the priority of the process, memory latency, and other factors.

a. Multiply the number in the Memory HiWater column of the ja report by 512. This gives you the maximum size of your process as measured in Cray words.

b. Determine the amount of available user memory and the number of available processors on your system by using the sysconf(1) command (see Section 2.2.2.1, page 22). Use the USRMEM value in the SOFTWARE report of the sysconf command as the amount of available user memory for your system. Use the NCPU value in the HARDWARE report of the sysconf command as the number of available processors.

c. Although you might be able to execute a process as large as the size stated in USRMEM of the sysconf report, you are probably sharing your system with other users. Also, your system most likely has multiple CPUs. Therefore, use the following rules of thumb to determine whether the memory high-water mark for the process is a significant fraction of available user memory.

• If the system on which the code is running has 4 CPUs or less, divide the maximum size of your process by user memory (USRMEM on the sysconf output). If the result is equal to or greater than .333, your process will probably benefit from optimization of memory management. If the result is less than .333, go to the next step.

• If the system on which the code is running has more than 4 CPUs, divide the user memory (USRMEM on the sysconf output) by the number of processors (NCPU on the sysconf output). Compare this number to the maximum size of your process. If the maximum size of your process is bigger than this number, your process will probably benefit from optimization of memory management. If the
maximum size of your process is not bigger than this number, go to the next step.

3. If the number in the Sys CPU Seconds column in the ja report for the code is greater than 10% of the number in the User CPU Seconds column, you have memory-bound code. Inefficient memory management is one cause for excessive elapsed time.

If the number in the Sys CPU Seconds column is not greater than 10% of the number in the User CPU Seconds column, your code is not memory bound.

2.2.2.1 Determining how much memory is available on your system

To find out how much user memory is available on your Cray SV1 system, use the `sysconf`, `target`, or `sar -M` commands, or ask your system administrator.

The `sysconf(1)` command reports a number for user memory (USRMEM) in the SOFTWARE portion of its report, as shown in the following sample portion of a `sysconf` command output.

```
HARDWARE: SERIAL= SN9617  MTYPE= Cray SV1  MFSUBTYPE= SV1
        NCPU= 20  NSSP= 8  NMSP= 3  CPCYCLE= 10.0000 ns
        MEM= 4294964992  NBANKS= 1024  CHIPSZ= 67108864
        AVL= YES  BDM= YES  EMA= YES  HPM= YES  BMM= YES
        SSD= 0  SSDRINGS= 0  IOS= MODEL_F  RINGS= 4

SOFTWARE: RELEASE= 10.00  POSIX VERSION= 199009  SECURE SYS= ON
        SYSMEM= 60092416 WRDS  USRMEM= 4234872576 WRDS
        OS_HZ= 60  CLK_TCK= 10000000
        JOB_CONTROL= YES  SAVED_IDS= YES  SCTRACE= ON
        UID_MAX= 16777215  PID_MAX= 100000
        NDISK= 32  SDS= 0  NBUF= 8000
        POSIX_PRIV= ON  SECURE_MLSDIR= SECURE  SECURE_MAC= OFF
        PRIV_SU= ON  PRIV_TFM= OFF
```

The `target(1)` command reports the absolute memory size of your system in the memsize value as shown in the following sample portion of a `target` command output. Available user memory should be greater than 90% of this value.

```
Primary machine type is: CRAY SV1
banks        = 1024
```
numcpus = 20
ibufsize = 32
memsize = 4294967296
memspeed = 33
clocktim = 10000
numclstr = 21
bankbusy = 14

The sar(8) command with the -M option specified provides a dynamic report of available user memory.

For more information on the sysconf(1), target(1), and sar(8) commands, see the man pages.

2.3 I/O bound

If a program spends most of its elapsed time performing I/O, it is considered I/O bound. I/O optimization can offer a significant savings in elapsed time. If the design of a program requires large amounts of I/O, you should optimize for I/O performance. However, if the code runs 24 hours and performs only 1 hour of I/O, there are other areas of optimization that will probably have a greater impact on overall code performance.

Procedure 4: Determining if the code is I/O bound

To determine whether the code is I/O bound, perform the following steps:

1. Run the job accounting utility, ja -clth (single-tasked), with the code to get a report for an overall view of memory usage, I/O, and system overhead. You can use the same report you created in Procedure 3, page 19, without running the code again.

   Note: To ensure a representative reading, the user process you are measuring with the ja utility should have an execution time that is greater than 1 second.

2. If the sum of the number in the I/O Wait Sec Lck column plus the number in the I/O Wait Sec Unlck column of the ja report is close to or greater than 50% of the number in the User CPU-Seconds column, the code is probably I/O bound.
Note: If a program has frequent or large formatted I/O requests, it is possible for it to be I/O bound even if the test in this step indicates that it is not. Formatted I/O (specified by using a Fortran `FORMAT` statement or a C++ `scanf` or `printf` command) consumes user CPU seconds to translate between ASCII and internal binary representation. This form of I/O time is not reflected in the two I/O Wait time columns of the ja report. A program that contains frequent or large formatted I/O requests might benefit from I/O optimization.
Multi-streaming is a feature that lets you schedule four dedicated Cray SV1 CPUs as one multi-streaming processor (MSP) for a Fortran, C, or C++ program. It automatically divides loop iterations among the four CPUs. You may get speedup factors of up to four on loops to which this technique can be applied.

Figure 6, page 25, illustrates the loop iterations that each CPU will operate on for the following loop. (An MSP is constructed of four CPUs, and an SSP is a single CPU.)

```fortran
DO I = 1, 20
    A(I) = A(I) * 3.14
ENDDO
```

<table>
<thead>
<tr>
<th>CPU0</th>
<th>I = 1, 5</th>
</tr>
</thead>
<tbody>
<tr>
<td>CPU1</td>
<td>I = 6, 10</td>
</tr>
<tr>
<td>CPU2</td>
<td>I = 11, 15</td>
</tr>
<tr>
<td>CPU3</td>
<td>I = 16, 20</td>
</tr>
</tbody>
</table>

Figure 6. Dividing Loop Iterations Among CPUs

Multi-streaming is an optional feature. To determine whether or not multi-streaming is enabled on your Cray SV1 system, enter the `sysconf(1)` command at your system prompt. The `HARDWARE` output field contains the `NMSP` field, which shows the number of multi-streaming processors configured. The relevant field is in boldface type in the following example of `sysconf` output. Having the `NMSP` field set to 2, as in this example, means two MSPs are configured for multi-streaming.

```
HARDWARE: SERIAL= SN3202   MTYPE= CRAY-SV1   MFSUBTYPE= SV14XX
NCPU= 32   NSSP= 24   NMSP= 2   CPCYCLE= 10.0000 ns
MEM= 1073739520   NBANKS= 1024   CHIPSZ= 16777216
```
Multi-streaming on a Cray SV1 system is similar to autotasking in distributing loop iterations across processors. However, multi-streaming causes gang scheduling of all requested processors, meaning they are attached to the program whether they are actually executing code or not. Autotasking schedules processors as they are needed. Processor utilization efficiency is directly proportional to the extent that you are able to multi-stream the program. For information on how to invoke and tune your program to enhance multi-streaming, see the following section.

### 3.1 Compiler options and directives

You can control multi-streaming within your program by means of command-line options and compiler directives.

#### 3.1.1 Fortran compiler options

To enable multi-streaming in a CF90 program, include either the `-O stream1`, `-O stream2`, or `-O stream3` option on the `f90(1)` command line. The `-O stream2` and `-O stream3` options are more aggressive than `-O stream1`. For more information on using these options with Fortran, see the CF90 Commands and Directives Reference Manual.

In the following example, the `f90(1)` command turns on streaming:

```bash
% setenv NCPUS 1
% f90 -O stream2 -O nothreshold streamer.f90
% ./a.out
```
The `setenv` command specifies one multi-streaming processor, meaning your program gets four SSPs.

**Note:** You may get slight numeric differences when you select the most aggressive streaming option (stream3). Check your program output to see if the answers generated are acceptable.

### 3.1.2 C and C++ compiler options

In C and C++, use the `-h stream1`, `-h stream2`, and `-h stream3` compiler options to enable multi-streaming. The `-h stream2` and `-h stream3` options are more aggressive than `-h stream1`. See the *Cray Standard C and Cray C++ Reference Manual* for more information on these options.

The following example specifies conservative streaming. By default, the program runs on one MSP. The `-h report=m` option generates multi-streaming optimization messages.

```sh
% setenv NCPUS 1
% cc -h stream1 -h report=m -h nothreshold streams.c
% ./a.out
```

### 3.1.3 Directives

The following directives allow you to enhance the performance of multi-streaming:

- **CONCURRENT**, see the following subsection.
- **PREFERSTREAM, STREAM, and NOSTREAM.** See Section 3.1.3.2, page 28.

#### 3.1.3.1 CONCURRENT directive

The `CONCURRENT` directive lets you communicate to the compiler that a loop is parallel when the compiler cannot make that determination on its own. Placing the following in front of a loop instructs the compiler to multi-stream that loop.

**Fortran:**

```fortran
!DIR$ CONCURRENT
```

**C and C++:**

```c
#pragma _CRI concurrent
```
This directive is especially important for C programs in which the presence of pointers can make parallelizing loops difficult for the compiler.

The ivdep directive serves this purpose for vectorization, but it is not a powerful enough assertion to allow the compiler to multi-stream a loop. Converting all ivdeps to concurrent in your program both enables multi-streaming improves the performance of vector code.

Applying the concurrent directive to all the loops that can be multitasked gives the compiler the best chance of generating optimal code. For information on the kinds of loops that can be multi-streamed, see Section 3.2, page 29.

### 3.1.3.2 PREFERSTREAM, STREAM, and NOSTREAM directives

You can designate or skip loops using the multi-streaming directives. The Fortran directives are as follows:

```fortran
!DIR$ preferstream
!DIR$ nostream
!DIR$ stream
```

The preferstream directive selects which loop, among two or more nested loops, to be multi-streamed. Insert the directive immediately in front of the loop to be multi-streamed. If the compiler has determined that it is safe to multi-stream the loop, it will do so.

The nostream and stream directives toggle multi-streaming off and on in Fortran. The nostream directive turns multi-streaming off in your program until a stream directive is encountered.

The following Fortran example turns streaming on for the inner loop:

```fortran
DO I=1,N1
  !DIR$ PREFERSTREAM
  DO J=1,N2
    A(I,J) = B(J,I)
  ENDDO
ENDDO
```

The C and C++ directives are as follows:

```c
#pragma _CRI nostream
#pragma _CRI preferstream
```
Unlike Fortran, there is no \#pragma_CRI stream directive. The nostream directive has different meanings in Fortran and C. The C and C++ version turns multi-streaming off only for the loop that immediately follows the directive.

In cases in which the compiler could multi-stream more than one loop in a loop nest, the preferstream directive instructs it to choose the one immediately following the directive.

### 3.2 Loops that are multi-streamed by the compiler

The compiler uses certain criteria to judge whether or not a loop can be automatically multi-streamed.

At the point the compiler evaluates a program, the code has already gone through an initial restructuring, as it would for vectorization or multitasking. Optimizations such as loop interchange (switching an inner loop with an outer loop) and loop splitting (changing a single loop into two or more loops) may have been done.

If a loop passes the following tests, it will be multi-streamed:

- Its iterations can be divided among different processors without delivering incorrect results. The loop can contain private arrays if their values are not used outside the loop.

- There are no function or subroutine calls within the loop. (The autotasking CNCALL directive is ignored.)

- The loop contains no bit matrix multiply (BMM) operations.

- A scatter operation is ordered as opposed to random. That is, there must be a constant stride through the array to be scattered; taking random elements from an array cannot be multi-streamed.

- The loop is not contained in a vector loop.

- It is not prefaced by the nostream directive.

- It has a trip count of at least two.

- The \-O threshold CF90 option or the \-h threshold C and C++ option is not in effect, and the resulting compile-time or run-time evaluation determines that there is enough work to make multi-streaming worthwhile. Because threshold is the default, you must specify \-O nothreshold or \-h nothreshold to turn it off.
When two loops that qualify to be multi-streamed are nested, the following criteria are used in the following order to choose which loop to multi-stream:

1. The outermost loop that is prefaced by the preferstream directive.

2. The loop that the compiler estimates will have the greatest amount of work after the loop is interchanged to its outermost valid position.

3. The outermost loop after initial restructuring.

Once a loop has been selected, it is moved to its outermost valid position before multi-streaming. If the loop is not prefaced by a preferstream directive and has a run-time trip count, two versions of the loop are generated, one multi-streamed and the other not. Which version is executed at run time depends on the trip count.

### 3.3 Bit Matrix Multiply (BMM) and multi-streaming

Your view of the Bit Matrix Multiply (BMM) programming model on the Cray SV1 system is identical to that used on all previous Cray platforms: a single BMM register, 64 elements deep and wide, and the intrinsic functions. The Fortran functions are m@ld, m@mx, m@ldmx, m@ul, and m@clr, and the C and C++ functions are _mclr_mld, _mmx, _mldmx, and _mul.

The compiler automatically utilizes the multiple BMM registers available on an MSP and divides the BMM work between them.

There are two types of BMM loops: those that directly contain BMM operations and those that indirectly contain BMM operations (that is, those that contain direct or indirect BMM loops). The following list uses these terms in describing how the compiler applies BMM operations to the SSPs of an MSP:

- Loops that directly contain an m@ld or _mld that is used outside the loop need to be executed redundantly across all SSPs. Loops that indirectly contain an m@ld or _mld can and should be multi-streamed as usual.

- Loops that directly contain m@ldmx or _mldmx can be multi-streamed, but partitions (except the last) must be multiples of 64. Loops that indirectly contain m@ldmx or _mldmx can be multi-streamed with no restrictions on partition size.

- Loops that contain m@mx or _mmx (directly or indirectly) can be multi-streamed with no restrictions on partition size.
Multi-streaming

• Loops that directly contain _mul or _mul cannot be multi-streamed. They can, however, be executed redundantly across SSPs when containing a feeding _mul or _mld, as in the following loop:

```c
CDIR$ shortloop
    do i=0, vl-1
        m@bmm(i) = m@ld( m@ul() )
    enddo
```

• Loops that contain _mx or _mmx with a reaching _mld or _mld (or a _mldmx or _mldmx) outside the loop and imported from a loop that is not redundantly multi-streamed, first broadcast the BMM value from SSP 0 to the others before multi-streaming. (All loops will be multi-streamed so that SSP 0 will execute the last iteration of the loop. Consequently, SSP 0’s BMM register will always contain the value that is indicated by the BMM programming model to be in the user’s BMM register at the end of a loop.)

• Loops that directly or indirectly contain _clr or _mclr can be multi-streamed. If an _clr or _mclr is not in a multi-streamed loop, it should be executed redundantly across all SSPs.

3.4 Tasking and multi-streaming

Multi-streaming and multitasking can coexist in the same program. See the following sections for more information:

• Multitasking on MSPs with multi-streaming (see the following section).
• Multitasking on MSPs without multi-streaming (see Section 3.4.2, page 32).

Multi-streaming offers the following performance advantages over multitasking:

• The startup time for multi-streaming is faster.
• Communication between multi-streaming processors is more efficient than between multitasking processors.

Still, some programs may run as fast or even faster when they use multitasking. Running your program both ways will tell you which method is better for you.

3.4.1 Multitasking on MSPs with multi-streaming

Multitasking and multi-streaming can be combined within an application to provide better performance than either model alone. Within a function, the
compiler will do either multi-streaming or multitasking, but not both. If the function contains multitasking directives, the compiler will do multitasking; otherwise, it will do multi-streaming.

Usually, the best approach to combining the models is to apply multitasking through user directives to parallel loops that contain subroutine calls, and apply multi-streaming within the called subroutines.

If you include both multitasking and multi-streaming options on the compiler command line (for instance, `-O task2, -O stream2` in Fortran and `-h task2, -h stream2` in C and C++), the value of the NCPUS environment variable (see Section 3.1, page 26) will refer to MSPs. If you set NCPUS to 2 before compiling your program, you will have two MSPs, or a total of eight SSPs to work with.

In the following example, the loop will be multitasked by the compiler, and the iterations of the loop will be split between the two MSPs requested by the programmer:

```
!CMIC$ DOALL PRIVATE (I), SHARED (A,B,C,N)
   DO I = 1, N
       CALL FOO(A,B,C,I)
   ENDDO
```

The best CPU utilization is achieved if the work within a call to FOO is effectively distributed through multi-streaming across the SSPs in the MSP.

### 3.4.2 Multitasking on MSPs without multi-streaming

If you enable multitasking on the compiler command line (for instance, `-O task3` or `-h task3`) and do not specify multi-streaming, the compiler will not multi-stream for the compilation. The value of the NCPUS environment variable (see Section 3.1, page 26) now refers to SSPs instead of MSPs.

You may see improved performance for a multitasking program on the Cray SV1 system without making changes to the code. A multitasking program will compete with multi-streaming programs for processors if you set the MP_DEDICATED environment variable and enter the `cpu(8)` command as in the following example. MP_DEDICATED relieves some of the overhead of task scheduling. Notice that you must set the NCPUS environment variable to four times the value of the `cpu -a` argument.

```
% setenv NCPUS 8
% setenv MP_DEDICATED 1
```
3.5 Vectorization and multi-streaming

Once multi-streaming divides the iterations of a loop among the SSPs working on your program, a number of single-processor optimizations created by the compiler continue to improve the performance of your program. The most important is vectorization.

Vectorization usually yields a greater speedup than any of the other single-processor optimizations. The more you can get out of vectorization, the closer you can come to realizing the full potential of your Cray SV1 system.

For more information on vectorization, see Chapter 4, page 39.

The major strategies you can follow in order to enhance vectorization are:

- Keep the stride through the array small; ideally, use a stride of 1.
- Use long vector lengths; the more iterations, the better.
- Avoid less efficient vector code, such as loops that contain variant IFs and reductions.

3.6 Analyzing the performance of a multi-streaming program

The following tools are available for analyzing a multi-streaming program:

- The prof(1) command (see Section 3.6.1, page 33).
- The MSP STATS environment variable (see Section 3.6.2, page 36).

3.6.1 The prof command

The prof(1) command returns timing information at the function and subroutine level for each processor involved in a program. To generate and view performance data for a Fortran program, use the prof library and command and the profview(1) visualization tool, as follows:

```
% f90 -l prof -O stream2 -V -v streamer.f
% setenv PROF_WPB 1
% ./a.out
% prof -x a.out >prof.report
% profview prof.report &
```
For a C or C++ program, use the following commands:

% cc -l prof -h stream2 -V myprog.c
% setenv PROF_WPB 1
% ./a.out
% prof -x a.out >prof.report
% profview prof.report &

The following is an example of a pie chart view generated by profview. This represents an aggregation of data collected from all four MSPs involved in executing the program. It shows you the percentage of time consumed by the longest-running routines:
When a program is multi-streamed but not multitasked, the amount of time spent in barriers will depend on the percentage of the program that can be multi-streamed. When the program is executing in multi-streaming mode but is not currently processing a loop, only one SSP is executing the program; the others are waiting at a barrier until the next loop. The above example shows that wait time in the _MsBarrier routine.

Figure 7. Profview Pie Chart
3.6.2 The MSP_STATS environment variable

The MSP_STATS environment variable provides statistics on multi-streaming for Fortran, C, and C++ programs. To enable the generation of statistics, set MSP_STATS as follows:

```bash
% setenv MSP_STATS 1
```

Enter the following commands to run your Fortran program with aggressive multi-streaming, on one multi-streaming processor, and with MSP_STATS enabled:

```bash
% setenv MSP_STATS 1
% setenv NCPUS 1
% setenv NEW_ENTRY
% f90 -O stream2 -l libmsx.a myprog.f
% ./a.out
```

Running MSP_STATS for a C or C++ program involves the same set of commands:

```bash
% setenv MSP_STATS 1
% setenv NCPUS 1
% setenv NEW_ENTRY
% cc -h stream2 -l libmsx.a myprog.c
% ./a.out
```

MSP_STATS provides the following output:

<table>
<thead>
<tr>
<th>MSP</th>
<th>SSP</th>
<th>UserSecs</th>
<th>MsLibSecs</th>
<th>#MsEnts</th>
<th>#Parks</th>
<th>#Barrs</th>
<th>#CInvs</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>0</td>
<td>35.612</td>
<td>2.029</td>
<td>777737</td>
<td>783438</td>
<td>5401</td>
<td>788839</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>0.826</td>
<td>36.815</td>
<td>777737</td>
<td>783438</td>
<td>5401</td>
<td>788839</td>
</tr>
<tr>
<td>0</td>
<td>2</td>
<td>0.826</td>
<td>36.815</td>
<td>777737</td>
<td>783438</td>
<td>5401</td>
<td>788839</td>
</tr>
<tr>
<td>0</td>
<td>3</td>
<td>0.856</td>
<td>36.785</td>
<td>777737</td>
<td>783438</td>
<td>5401</td>
<td>788839</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>MSP</th>
<th>SSP</th>
<th>WaitIters</th>
<th>Mispredicts</th>
<th>OverFlows</th>
<th>Reconnect</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>0</td>
<td>3316440</td>
<td>0</td>
<td>1</td>
<td>24</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>205831159</td>
<td>10207</td>
<td>0</td>
<td>22</td>
</tr>
<tr>
<td>0</td>
<td>2</td>
<td>206147477</td>
<td>10207</td>
<td>0</td>
<td>23</td>
</tr>
<tr>
<td>0</td>
<td>3</td>
<td>205925417</td>
<td>10207</td>
<td>0</td>
<td>23</td>
</tr>
</tbody>
</table>
The column headers have the following meanings:

<table>
<thead>
<tr>
<th>MSP</th>
<th>Number of the multi-streaming processor. Since only one MSP is used, its number is 0.</th>
</tr>
</thead>
<tbody>
<tr>
<td>SSP</td>
<td>Number of each SSP. SSP 0 is the master processor.</td>
</tr>
<tr>
<td>UserSecs</td>
<td>Wall-clock time spent outside of multi-streamed code for each SSP. The master processor spent most of its time outside multi-streamed code.</td>
</tr>
<tr>
<td>MsLibSecs</td>
<td>Wall-clock time spent inside the multi-streaming library. When an SSP is inside the library, it is either waiting for a multi-streamed section of code to be entered or resumed (SSPs are parked while waiting at a barrier) or an SSP is triggering the beginning or resumption of a multi-streamed section of code.</td>
</tr>
<tr>
<td>#MsEnts</td>
<td>The number of times multi-streamed procedures were entered. Entire procedures are multi-streamed; within these procedures, SSPs 1-3 skip past code that must be executed by one SSP.</td>
</tr>
<tr>
<td>#Parks</td>
<td>The number of times the MSP was parked so that SSP 0 could execute a single-streamed piece of code. Parking and unparking is the way SSPs 1-3 skip around single-streamed code. The value is for the entire program.</td>
</tr>
<tr>
<td>#Barrs</td>
<td>The number of times the MSP encountered a streaming barrier in the program. Barriers are often needed for data synchronization or cache invalidation.</td>
</tr>
<tr>
<td>#CInvs</td>
<td>The number of cache invalidations. This must be less than or equal to the number of parks plus the number of barriers. Cache is invalidated for safety reasons.</td>
</tr>
<tr>
<td>WaitIters</td>
<td>SSPs execute <em>idle wait</em> loops while they are inside of the multi-streaming library and waiting for other SSPs to synchronize with them. <em>WaitIters</em> is the total number of iterations each SSP spent in these loops. The numbers should correlate roughly with <em>MsLibSecs</em>.</td>
</tr>
</tbody>
</table>
**Mispredicts**

The streaming library attempts to predict where SSPs 1-3 will be needed next after being parked, and it advances execution speculatively toward that point. The library safely prefetches data and instructions into the cache and may execute instructions leading up to the unpark location. **Mispredicts** is the number of times this prediction was incorrect. It should be much less than the number of parks for best performance.

**Overflows**

The number of times the streaming library detected a common stack overflow. This should be nonzero only for SSP 0.

**Reconnect**

The approximate number of times each SSP was reconnected by the operating system. The values for the SSPs making up an MSP should be small and roughly equal, indicating long connect times and gang scheduling.

The statistics are gathered whether or not you have set the **MSP_STATS** environment variable. There is little overhead associated with gathering the statistics.
Vectorized constructs perform up to a factor of 20 times faster compared with non-vector constructs. However, vectorization on the Cray SV1 system does not always perform in the same way it did on earlier Cray systems.

To make the best use of vectorization, it is first necessary to understand it. The following section describes how it provides you with performance gains. If you already understand vectorization, continue on with Section 4.2, page 41.

4.1 What is vectorization?

Vectorization is similar to an assembly line in manufacturing. To illustrate, Figure 8, compares two ways to carry out a five-step process for making chairs, in which each step takes one minute:

- In one process, one assembly must go through all five steps before the next assembly can begin the first step. One chair is produced every five minutes.

- With the use of an assembly line, each step in the sequence is performed as soon as the previous step is complete. In this way, the first chair is completed after five minutes, and one new chair is completed each minute after that.
Scalar processing: one at a time

Each incoming seat must wait for processing until previous assembly completes all five steps. One new chair is completed every five minutes.

Vector processing: assembly line

One chair is completed each minute.

Figure 8. Scalar versus vector, illustrated

Like manufacturing a chair, processing an array in a loop is a multistep process that can be performed one step at a time, or by a process resembling an assembly line.

To process an array in either vector or scalar mode, a loop must do the following:
• Load elements from memory to a register
• Process the elements, placing the results in another register
• Store the results back to memory

Conventional (scalar) code is a one-at-a-time process: each element must finish the final step of processing before the next element can begin the first step; that is, each loop iteration begins when the previous iteration ends. But with vector code, as in an assembly line, each element begins the first step when the previous element finishes the first step, rather than the final step.

4.2 Vectorization inhibitors

The Cray CF90 compiler will generate a loop mark listing showing which loops in a program vectorize or parallelize. This can be enabled with the \(-r m\) option on the \(f90\) command line. Loops with a very large number of lines (hundreds) may require the \(-O\) \(aggress\) optimization option on the compiler line; this allows for larger internal tables for compiler analysis.

Vectorization inhibitors within DO loops include:

• CALL statements
• I/O statements
• Backward branches
• Statement numbers with references from outside the loop
• References to character variables
• External functions that do not vectorize
• RETURN, STOP, or PAUSE statements
• Dependencies (see the following section)

Many of these can be addressed through slight modifications of the source code.

4.3 Dependencies

Dependencies are code constructs that prevent the compiler from vectorizing a loop. Dependencies can be real or potential. For example, the compiler will not cleanly vectorize the following loop.
Note: The compiler will mark this loop as vectorized, but the method used has extra overhead needed to check for repeated index values.

\[
\text{do } i = 1, n \\
\quad A(\text{index}(i)) = A(\text{index}(i)) + b(i) \\
\text{enddo}
\]

In this case, there is a potential dependency on the array \( A \). A dependency exists if the index array has repeated values within a vector length (64 elements). If, for example, the programmer knows that \( \text{index}(i) \) is monotonically increasing, or that \( \text{index}(i) \) is unique for each value of \( i \), then no dependency exists. You can assert that fact through a compiler directive:

\[
!\text{dir$ ivdep} \\
\text{do } i = 1, n \\
\quad A(\text{index}(i)) = A(\text{index}(i)) + b(i) \\
\text{enddo}
\]

Adding this directive to a safe loop can improve performance by up to a factor of 10.

Four indirect update cases in which an index list is used have also been identified. A set of routines has been developed by the Cray benchmarking group that will vectorize four flavors of indirect update loops by removing any vector dependency regarding the storing index. The routines will exceed the performance of Fortran 90 generated code, providing the storing index changes infrequently compared to the number of times the indirect update loops are executed. The four cases covered are as follows:

- The \texttt{ind_update.f} routine vectorizes the loop:
  \[
  \text{do } j=1,N \\
  \quad A( I(j) ) = A( I(j) ) + B(j) \\
  \text{enddo}
  \]

- The \texttt{ind_vpv.f} routine vectorizes the loop:
  \[
  \text{do } j=1,N \\
  \quad A( I(j) ) = A( I(j) ) + B(K(j)) \\
  \text{enddo}
  \]

- The \texttt{ind_vpvu.f} routine vectorizes the loop:
  \[
  \text{do } j=1,N \\
  \quad A( I(j) ) = A( I(j) ) + B(j) \times C(K(j)) \\
  \text{enddo}
  \]
- The `ind_vpvv.f` routine vectorizes the loop:

```fortran
    do j=1,N
        A(I(j)) = A(I(j)) + B(K(j))*C(L(j))
    enddo
```

The above routines are available from the Cray benchmarking group upon request.

In many cases, the dependency is real but can be programmed around with some loop restructuring. For example, in the following loop there is a vector dependency in both the inner and outer loops:

```fortran
    do i=1,n
        do j=1,m
            temp = .25*(x(i,j-1)+x(i-1,j)+x(i+1,j)+x(i,j+1)) - x(i,j)
            x(i,j) = x(i,j) + omega * temp
            if (abs(temp) .gt. err1) err1 = abs(temp)
        enddo
    enddo
```

In this case, `x(i,j)` is defined using its north, south, east, and west neighbors. In the current form of the loop, `x(i-1,j)` is needed to update `x(i,j)`, which creates a dependency in the `i` direction. Switching the loops gives us a similar problem in the `j` direction. This loop runs at about 20 Mflops on the Cray SV1.

The stencil and resulting dependency is shown in the following figure:

A solution to this problem is to change the vectors to run diagonally through the matrix `X`. This can be coded as follows:

```fortran
    do jd=2,n+m
        !dir$ ivdep
        do j=max(1,jd-m),min(n,jd-1)
            i = jd - j
            temp = .25*(x(i,j-1) + x(i-1,j) + x(i+1,j) + x(i,j+1)) - x(i,j)
            x(i,j) = x(i,j) + omega * temp
            if (abs(temp) .gt. err1) err1 = abs(temp)
        enddo
    enddo
```

[The figure that will go here is not yet complete.]
The leading dimension of the matrix should now be even because the stride is the leading dimension of array A-1. This loop now vectorizes and runs at over 260 Mflop/s on a 1000x1000 grid. Cache hit rates run at about 50% because two of the four stencil points lie in the recently updated vector and hence are in cache.

4.4 Examples

The following sections show vectorization examples and how you can optimize them.

4.4.1 Using odd–numbered strides

Odd numbered strides through a loop are optimal on Cray SV1 systems. The stride can be either positive or negative, so long as it is an odd number. In addition, all strides within a loop should match in order to minimize cache conflicts.

If you must use an even stride, avoid powers of two. If there is only a single factor of two, the access will only be slowed slightly (about 20%). If it contains a factor of four, the best you can get is half speed. Multiples of eight run at 1/4 speed.

4.4.2 A problem with unrolling

The following code fragment is from the X-Ray crystallography program SHELXL. As written here, it performs at 153 Mflops on the Cray SV1 computer:

```plaintext
REAL A(N), B(N), C(N), D(N), E(N)
M=(N/4)*4
DO 1 I=1,M,4
   A(I)=A(I)+C(I)*D(I)
   B(I)=B(I)+C(I)*E(I)
   A(I+1)=A(I+1)+C(I+1)*D(I+1)
   B(I+1)=B(I+1)+C(I+1)*E(I+1)
   A(I+2)=A(I+2)+C(I+2)*D(I+2)
   B(I+2)=B(I+2)+C(I+2)*E(I+2)
   A(I+3)=A(I+3)+C(I+3)*D(I+3)
   B(I+3)=B(I+3)+C(I+3)*E(I+3)
1  CONTINUE
IF (M+1.GT.N) GOTO 2
   A(M+1)=A(M+1)+C(M+1)*D(M+1)
```

S–2312–35
This is a simple loop that has been unrolled by 4. This was a common technique to speed up programs on non-vector systems, although today the unrolling task is better handled automatically by compilers. Unrolling the loop by 4 causes the following problems on the Cray SV1:

- The loop still vectorizes, but a stride of 4 has been introduced.
- The effective vector length has been reduced from \( N \) to \( N/4 \).

The code is simplified to get rid of these two problems by re-rolling the loop as follows:

```
DO 1 I=1,M
    A(I)=A(I)+C(I)*D(I)
    B(I)=B(I)+C(I)*E(I)
1    CONTINUE
```

The resulting code improves to 357 Mflops.

### 4.4.3 Loop ordering

When presented with a nest of loops, the compiler makes choices about which loop to vectorize and which loop to parallelize or multi-stream based on stride and vector length. When incomplete information is presented to the compiler, it can result in less than optimal performance. In addition, the compiler does not make loop order choices based on temporal locality considerations. Consider the following loop nest:

```plaintext
real a(69,92,115)
do k = 2,115
    do j = 2,92
        do i = 2,69
            c(i,j,k) = 2.5*(a(i,j,k)-a(i,j,k-1))
            * (b(i,j,k)-b(i,j,k-1))
        end do
    end do
end do
```

In this case, the compiler vectorized the \( k \) loop because it has the largest iteration count. Because array \( a \) has an even middle dimension and because \( k \) is the last index, this results in an even-strided vectorized loop. The code runs at 109 Mflops.

```fortran
25. V------< do k = 2,115
26. V 2----< do j = 2,92
27. V 2 3--< do i = 2,69
28. V 2 3 -> c(i,j,k) =
29. V 2 3 --> enddo
30. V 2----> enddo
31. V------> enddo
```

The \( j \) loop (second dimension) will give us the longest vector length that is not a multiple of two, so it makes sense to choose it for the innermost loop. Next, we see that a temporal reuse opportunity exists on the \( k \) index because both \( k \) and \( k-1 \) are referenced for arrays \( a \) and \( b \). For this reason, \( k \) should be the next loop in the nest. That leaves the \( i \) loop as the outermost. In addition, we can force the compiler to unroll the \( k \) loop into the \( j \) loop with an unroll directive that facilitates register reuse opportunities. Finally, the \texttt{prefer vector} directive tells the compiler to vectorize the \( j \) loop. The resulting code is as follows:

```fortran
real a(69,92,115)
do i = 2,69
!dir$ unroll(4)
do k = 2,115
!dir$ prefer vector
do j = 2,92
  c(i,j,k) = 2.5*(a(i,j,k)-a(i,j,k-1))
$ *(b(i,j,k)-b(i,j,k-1))
  enddo
enddo
enddo
```

The performance improves to 247 Mflops.

A good general rule of thumb is to vectorize on the longest odd dimension, and then look for temporal reuse opportunities in the next level loop.
This chapter is organized as follows:
• Overview of memory (see the following section)
• Optimizing cache use (see Section 5.2, page 51)
• Managing memory (see Section 5.3, page 55)

5.1 Overview of Memory

This section is organized as follows:
• Central memory (see the following section)
• Cache (see Section 5.1.2, page 49)

5.1.1 Central Memory

The tables in this section give latencies and bandwidths for the Cray SV1 system. The first lists typical processor latencies. The time is given in 300 Mhz clock periods:

<table>
<thead>
<tr>
<th>Operation</th>
<th>Time</th>
</tr>
</thead>
<tbody>
<tr>
<td>V register to cache</td>
<td>25</td>
</tr>
<tr>
<td>S register to cache</td>
<td>22</td>
</tr>
<tr>
<td>V register to memory</td>
<td>109</td>
</tr>
<tr>
<td>Floating point add</td>
<td>8</td>
</tr>
<tr>
<td>Floating point multiply</td>
<td>9</td>
</tr>
<tr>
<td>Floating point reciprocal</td>
<td>16</td>
</tr>
<tr>
<td>Jump</td>
<td>6</td>
</tr>
</tbody>
</table>

The next table is a summary of the configuration characteristics and peak rates for memory module types:
Memory module type  | mem128 | mem512 |
---|---|---|
Number of CPUs  | 32 | 32 |
Number of memory sections  | 8 | 8 |
Number of memory subsections  | 8 | 8 |
Number of pseudo bank pairs  | 512 | 512 |
Pseudo bank pair busy clocks  | 8 | 6 |
Memory size in Gbytes  | 8 | 32 |
CPU to cache rate in Gbytes per second  | 9.6 | 9.6 |
CPU to memory rate in Gbytes per second  | 3.2 | 3.2 |
Memory rate in Gbytes per second  | 51.2 | 68.3 |

To measure the sustainable rate between the CPU and memory, the STREAM benchmark was used. Because of its memory reference patterns, the performance of the benchmark is not affected by cache. In the following table, results from the benchmark are presented with varying processor counts and varying processor assignments to modules. The results are in Gbytes per second.

<table>
<thead>
<tr>
<th>CPUs</th>
<th>Modules</th>
<th>mem512</th>
<th>mem128</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>1</td>
<td>2.52</td>
<td>2.51</td>
</tr>
<tr>
<td>2</td>
<td>1</td>
<td>4.71</td>
<td>4.68</td>
</tr>
<tr>
<td>2</td>
<td>2</td>
<td>5.00</td>
<td>4.97</td>
</tr>
<tr>
<td>4</td>
<td>1</td>
<td>5.08</td>
<td>5.07</td>
</tr>
<tr>
<td>4</td>
<td>2</td>
<td>9.25</td>
<td>9.10</td>
</tr>
<tr>
<td>4</td>
<td>4</td>
<td>9.89</td>
<td>9.76</td>
</tr>
<tr>
<td>8</td>
<td>2</td>
<td>9.95</td>
<td>9.74</td>
</tr>
<tr>
<td>8</td>
<td>4</td>
<td>16.76</td>
<td>15.00</td>
</tr>
<tr>
<td>8</td>
<td>8</td>
<td>18.89</td>
<td>18.26</td>
</tr>
<tr>
<td>12</td>
<td>4</td>
<td>14.22</td>
<td>13.51</td>
</tr>
<tr>
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<td>21.72</td>
<td>19.58</td>
</tr>
<tr>
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<td>8</td>
<td>22.93</td>
<td>20.62</td>
</tr>
<tr>
<td>16</td>
<td>4</td>
<td>17.69</td>
<td>16.30</td>
</tr>
</tbody>
</table>
### 5.1.2 Cache

The Cray SV1 cache size is 256 KBytes. It is 4-way set associative, write allocate, and write through; the least-recently-used data element is replaced. The data resulting from vector, scalar, and instruction buffer memory references is cached. The cache line size is 8 bytes, or 1 word long.

Cache is located between the CPU and the interface to main memory. The interface between cache and the CPU consists of four 64-bit read data paths and two 64-bit write data paths. The same number and types of paths exist between the cache and the interface to memory. For an example of how cache works on a read, see Section 1.3.2, page 7.

Cray SV1 cache differs from those on mainstream microprocessors in several important ways.

<table>
<thead>
<tr>
<th>CPUs</th>
<th>Modules</th>
<th>mem512</th>
<th>mem128</th>
</tr>
</thead>
<tbody>
<tr>
<td>16</td>
<td>8</td>
<td>25.04</td>
<td>21.66</td>
</tr>
<tr>
<td>20</td>
<td>8</td>
<td>24.85</td>
<td>21.65</td>
</tr>
<tr>
<td>24</td>
<td>8</td>
<td>24.98</td>
<td>21.84</td>
</tr>
<tr>
<td>28</td>
<td>8</td>
<td>25.37</td>
<td>21.83</td>
</tr>
<tr>
<td>32</td>
<td>8</td>
<td>25.44</td>
<td>21.93</td>
</tr>
</tbody>
</table>

**cache line width**

The cache-line width for vector references is a single word (8 bytes). Unlike most microprocessors, contiguous memory references are not required in order to achieve full cache (memory) bandwidth.

**bandwidth**

Data cache has very high memory bandwidth (up to 4 words per clock period).

**size**

The Cray SV1 cache is 256 Kbytes, which is small by modern standards. You may need smaller blocking factors than on many microprocessor-based systems.

**write-through**

This means that any memory stores go all the way to memory. Because system memory bandwidth is often a limiting resource, it pays to minimize unnecessary data stores.
coherence  Due to the need for binary compatibility, Cray SV1 cache does not maintain coherency with the other processors. This means that if another processor performs a store to memory, the new data value is not in cache. When the system synchronizes multiple processors, cache is invalidated (essentially, cleared). The implication is that better parallel performance is achieved with larger granularity.

Cache on the Cray SV1 system is 4-way set associative, which means the cache is organized into sets. Each set is composed of 4 separate ways or cache lines. (For an illustration, see Figure 3, page 8.)

A memory address is mapped in a manner that results in the data represented by that address being placed in one particular cache set. That is, in an $N$-way set associative cache, an expression such as the following indicates to which set in cache the address maps:

$$\text{modulo } (\text{memory_address}, \frac{\text{cache_size}}{N})$$

Since memory is much larger than cache, each set has many memory addresses that map into it. The four ways allow up to four memory addresses to map into the same cache set. If it becomes necessary to map a new address into a fully allocated cache set, the way for the least recently used address will be used and its data will be overwritten.

**Note:** Direct-mapped cache and fully associative cache can be viewed as special end cases of the general $N$-way set-associative cache. For direct-mapped cache, $N$ equals 1, and for fully-associative cache, $N$ equals the number of cache lines.

The write allocate attribute of cache requires that the address and data for a CPU write request to memory be mapped and placed into cache, if the request generates a cache miss. Because the Cray SV1 system has a write-through cache, any value that is updated in cache will always be written through to memory at the same time. This is in contrast to a write back cache, in which a value is updated in memory when there is no longer room for it in cache (that is, the value will be written back during a read to another value, which makes reads more expensive).

Write through can mean that more writes to memory are made than are strictly necessary when data is updated several times while remaining cache-resident. The cache contains buffers capable of holding 384 outstanding references. This
is sufficient for the cache to handle 6 vector references of stride 1 (384 references = 6 vectors * 64 references per vector).

The 1-word cache line size is advantageous for non-unit vector strides, since it does not cause the overhead of unnecessary data traffic when referencing memory using larger cache line sizes.

The disadvantage is that a single scalar reference will not bring in surrounding data, thus potentially inhibiting a scalar code from taking advantage of spatial locality. For this reason, scalar references have a prefetch feature, whereby a scalar reference causes 8 words to be brought into cache. These 8 words are determined by addresses that match the reference, except for the lower 3 bits. The effect for scalar loads is similar to having a cache line size of 8 words.

The Cray SV1 system relies on software that maintains cache coherency between processes that share memory. Cache is invalidated (cleared) as part of the test-and-set instruction. The test-and-set instruction has been used for processor synchronization in previous generation Cray vector systems. Adding the cache invalidation feature to this instruction allows old Cray binaries to run in parallel on the Cray SV1 system with the data cache enabled.

Data and instructions are cached by default. To turn caching off, use a command such as the following:

```bash
% /etc/cpu -m ecfoff a.out
```

For more information, see the `cpu(8)` man page.

## 5.2 Optimizing Cache Use

The following cache optimization techniques are described in this section:

- Minimizing stores (see Section 5.2.1, page 51)
- Porting issues (see Section 5.2.2, page 54)

### 5.2.1 Minimizing stores

The Cray SV1 cache policy is write-allocate and write-through. This means that any store will consume memory bandwidth and cache space. In many cases, stores can be reduced through unrolling and a technique called outer loop vectorization.
The following matrix-vector multiply kernel illustrates the techniques. A matrix vector multiply of size N has \(2N^2\) floating-point operations and \(N^2\) data. From an algorithm perspective, a minimum of one memory operation will be required for every two floating-point operations, giving a maximum computational intensity of two (two FLOPS per memory operation).

To inhibit full compiler optimization, the loop is compiled as follows:

```fortran
% f90 -O nopattern,nointerchange -rm mxv.f

```subroutine mxv(a,lda,n,b,x)
real a(lda,n), b(lda), x(lda)
1----< do j = 1, n
1 Vr--< do i = 1, n
1 Vr x(i) = x(i) + a(i,j) *b(j)
1 Vr--> enddo
1----> enddo
return
end
```

From the listing file, the inner loop is vectorized (V) and unrolled (r). The unrolling here is on a vector chime basis (64 elements), not on an iteration-by-iteration basis. For example, if the compiler unrolls a vector loop by two, it means two vector chimes (or 128 elements) are processed before loop iterations are incremented.

HPM shows that this loop runs at 250 MFlops and is requesting operands at the rate of 374 Mwords/s (three words for every two flops, or a computational intensity of 2/3). Of this, 250 Mwords/s is satisfied by main memory and the remaining 124 Mwords/s is satisfied from data cache.

Since \(x(i)\) is updated for each pass of \(j\), we can unroll the \(j\) loop into the inner loop and reduce the number of times \(x(i)\) is updated. For example, if we unroll by four times, we should reduce loads and stores to \(x\) by a factor of four:

```fortran
```

subroutine mxvl(a,lda,n,b,x)
real a(lda,n), b(lda), x(lda)
1----< do j = 1, n, 4
1 V--< do i = 1, n
1 V x(i) = x(i) + a(i,j) *b(j)
1 V 1 + a(i,j+1)*b(j)
1 V 1 + a(i,j+2)*b(j)
1 V 1 + a(i,j+3)*b(j)
1 V--> enddo
1----> enddo

```
Performance has improved from 250 Mflops to 301 Mflops. In addition, the overall bandwidth consumed has dropped from 374 Mwords/s to 226 Mwords/s (38 Mwords/s from cache and 188 Mwords/s from main memory).

Another alternative is to switch the loop ordering and go to a dot-product formulation. This eliminates vector store traffic in the inner loop:

```fortran
subroutine mxv(a,lda,n,b,x)
  real a(lda,n), b(lda), x(lda)
  do i = 1, n
    do j = 1, n
      x(i) = x(i) + a(i,j) * b(j)
    enddo
  enddo
  return
end
```

This formulation runs at 298 Mflops, consuming 306 Mwords/s of total bandwidth (153 Mwords/s from cache and 153 Mwords/s from memory). In this case, loads to \( b(j) \) are cached and loads to \( a(i,j) \) are not. We have reached our algorithmic ideal ratio of two flops for every memory operation. Dot products, however, are not ideal on vector systems due to the final vector reduction operation, in which a vector register of operands is collapsed down to a single scalar value.

The ideal algorithm would hold 64 elements of \( x \) in a vector register until all updates are complete. The 64 elements of the completed vector \( x \) would be written to memory. This technique is called outer-loop vectorization. It can be achieved by writing the loop in a dot-product formulation and then inserting a `cdir$ prefer vector` directive before the outer loop:

```fortran
subroutine mxv(a,lda,n,b,x)
  real a(lda,n), b(lda), x(lda)
  cdir$ prefer vector
  do i = 1, n
    do j = 1, n
      x(i) = x(i) + a(i,j) * b(j)
    enddo
  enddo
  return
end
```
This loop now runs at 395 Mflops, consuming only 201 Mwords/s of total bandwidth (199 Mwords/s from memory and 2 Mwords/s from cache). Reuse has moved from cache into a vector register. In many cases the compiler will choose this formulation automatically.

The Cray SV1 memory is capable of delivering operands at the rate of 2.5 GBytes/s (312 Mwords/s). Since two flops are computed for every word of memory bandwidth, this algorithm has a peak potential rate of 624 Mflops (312 Mwords/s * 2) on the Cray SV1. When coded in assembly language, vector loads to a can be carefully scheduled to achieve maximum bandwidth:

```fortran
subroutine mxv(a,lda,n,b,x)
  real a(lda,n), b(lda), x(lda)
  CALL SGEMV ('n', N, N, 1., A(1,1), LDA, B(1), 1, 1., X(1), 1)
  return
end
```

The scientific libraries code runs this problem at 600 Mflops, consuming 305 Mwords/s of memory bandwidth (3 Mwords/s from cache and 302 Mwords/s from memory).

### 5.2.2 Porting Issues

The Cray SV1 system libraries for parallel processing have all been modified to invoke the test-and-set instruction when cache must be invalidated. Therefore, you should be able to port codes written for Autotasking, OpenMP, and MPI without modification.

SHMEM codes should also port without modification, with one exception. Autotasking and OpenMP use a test-and-set is issued at the beginning and end of parallel regions, at parallel loop iterations, and at critical regions (locks and guards).

For MPI programs, there is no shared data from the user’s view, and the libraries take care of managing consistency within themselves. Although it is better to maximize the granularity of parallel tasks and minimize synchronization on any architecture, there is additional incentive for the Cray SV1 system because it is best to avoid cache invalidations.

The Cray SV1 SHMEM library routines `shmem_barrier(3)`, `shmem_wait(3)`, and `shmem_udcflush(3)` perform cache invalidations. In order to avoid race
conditions, a `shmem_barrier` or `shmem_wait` is usually issued before remotely updated data is used.

That will take care of cache coherency considerations at the same time. Codes that were originally written for the Cray T3D computer may need to be modified if they make use of the `shmem_set_cache_inv` and `shmem_clear_cache_inv` routines. These routines invalidated cache on the Cray T3D system and are ignored on the Cray T3E computer, but they are not supported on the Cray SV1 system; you will get an unsatisfied external message when you try to load.

These codes will need to be reworked, probably by replacing each `shmem_set` or `clear_cache_inv` call with a `shmem_barrier` or `shmem_udcflush`.

### 5.3 Managing Memory

The Cray SV1 system contains fixed-size, real memory; it has no virtual memory capabilities. Large memory codes must fit within available user memory. User processes that dynamically expand and shrink during execution are managing memory, and they must compete with other processes for real memory space. Processes that spend excessive time managing memory are said to be memory bound.

Inefficient memory management within a process can increase its elapsed time, and this can affect other user processes in the system. If the process is large compared to the amount of available user memory, small performance problems can become worse. When a process attempts to dynamically change its size, it becomes a candidate for a swap out of memory onto an external device (such as a disk) before the operating system can find enough contiguous memory to swap it back in. Also, the code might make system memory calls without your knowledge. These situations can greatly increase elapsed time.

Optimizing memory-bound code reduces elapsed time and system CPU time, and it may have a side effect of reducing user CPU time.

**Note:** Most of the memory management techniques described in this chapter reduce elapsed time at the cost of increased memory usage.
Use the `procstat` and `procview` utilities to identify whether code spends large amounts of system CPU time processing memory requests.

- **Yes**
  - Evaluate alternatives
  - Apply memory technique
  - Check your answers
  - Satisfied?
  - **Yes**
    - Return to initial analysis phase
  - **No**
    - Satisfied?
  - **No**
  - **Yes**
- **No**
  - Code has large amount of memory wait time?
  - **Yes**
    - Apply memory technique
    - Check your answers
    - Satisfied?
    - **Yes**
      - Return to initial analysis phase
    - **No**
      - Satisfied?
      - **Yes**
      - Return to initial analysis phase
      - **No**

Continued from Figure 2

Figure 9.

*Optimizing*
5.4 Understanding Memory Management

Two basic types of dynamic memory management are available on Cray SV1 systems:

- Dynamic memory managed by the system heap routines
- Expandable dynamic common blocks

You cannot use both methods within the same code.

5.4.1 Dynamic Heap

When code is compiled and loaded, it is translated into Cray machine instructions (object code), logically linked together with library routines, and packaged in an executable file, named `a.out` by default. When you execute `a.out`, it is swapped into memory and becomes a user process. Most UNIX systems have at least the following three distinct areas defined in memory for each user process:

- User area
- Stack area
- Heap

The user area is where the object code (text area) and external and static variables (data area) reside. The stack area is commonly used to temporarily store context information for each routine that calls another subprogram. The heap is a dynamic area used for all other memory needs (except Fortran COMMON): dynamic variables, I/O buffers, flexible file input/output (FFIO) user cache, and so on.

The heap and the stack areas are allowed to grow dynamically, but the user area remains fixed. On a virtual memory system, both the heap and the stack area of a single process can grow independently with a virtual hole between them.

The UNICOS operating system implements dynamic allocation of heap and stack space differently from virtual systems. A UNICOS user process has a dynamically expandable heap, with stack space wholly contained within the heap. The stack space is managed without benefit of direct hardware support, and both heap and stack space must appear to grow and shrink independently.

Initial memory allocation for the heap and the stack space is established at the load step by the `ld(1)` utility (called by the `f90(1)`, `CC(1)`, and `cc(1)` commands).
The heap is dynamic and can be increased or decreased from an executing code by using calls to the library. Routines request space from the heap directly through calls to the UNICOS system.

Dynamic memory management is inherently expensive to a user process because it requires service from the operating system through system calls. An expansion of the heap might require the process to be relocated in memory. If there is no remaining space large enough in memory, the UNICOS operating system will swap the requesting process to a secondary device (such as a disk) until enough memory becomes available. This adds elapsed time and system CPU time.

Stack space must be allocated specifically in finite segments, or stack frames, within the heap. A stack frame is used as a place holder for procedure calls to save context, local variables, local arrays, and so on. Each procedure call pushes (allocates) a new frame on the stack, and pops (deallocates) that frame upon exit. The stack is a last-in, first-out model (LIFO) allocated inside the heap in finite-sized segments. Therefore, the stack frames compete with the rest of the dynamic memory requirements for space.

The following conditions can cause excessive system time associated with managing memory for your user process:

- Releasing heap blocks in a different order from that of their allocations can cause memory fragmentation for the process.
- Upon entry to a function, routine, or procedure in a user process, it might not be possible to allocate a stack frame within the boundaries of the current stack segment. This is a stack overflow condition (transparent to programmers) that causes attempts to allocate an additional stack segment within the heap.
- If a new stack segment cannot fit contiguously within the heap, total stack space becomes fragmented. This costs extra time for subroutine calls until that stack segment is no longer needed and returns to the heap (a stack underflow condition).
- stack thrashing is a rare condition in which a frequently called function allocates (stack overflow) and deallocates (stack underflow) both a stack segment and a heap block for every invocation.
- Fortran automatic arrays and C variable length arrays are allocated with compiler-generated heap requests upon entry to a routine and deallocated upon exit from the routine. These arrays cause hidden memory management from system calls.
• Multitasked processes share a heap between multiple slave processes, each using its own stack space in the heap.

5.4.2 Dynamic Common Blocks

The UNICOS operating system provides a second method to manage memory for Fortran 90 codes, the dynamic common block. To use this method, you must specify only one dynamic common block (which might be blank common) for the loader to place at the high end of the process memory space. For an example of an £90 command line that establishes a dynamic common block, see Example 3 in Section 5.7.1, page 62.

This technique requires your heap to be a fixed size. Heap expansion is not allowed because the dynamic common area is stored directly after the heap. Therefore, the initial size of the heap must be large enough to handle all requests for heap space. Generally, an initial heap size between 5,000 and 10,000 words is adequate.

Fortran codes that use this method typically overindex an array within a dynamic common block, but the programmer must be careful to avoid an operand range error. You can expand and contract the dynamic common block by using the SBREAK library routine. SBREAK expands the field length of the user process to provide more memory, and it also releases memory when it receives a negative argument.

All subroutines within the same code have access to its dynamic common block at any time during program execution. Its contents cannot be initialized at load or compile time.

5.5 Identifying Large Amounts of Memory Wait Time or System CPU Time

The procstat(1) utility can provide accurate memory information about your code. It gathers process-level memory statistics, such as elapsed time, number of calls to memory processor, number of memory declines, and total time to complete memory requests.

To create a report, execute the procstat utility with the name of the program to be analyzed listed as an argument.

For complete information on procstat, see the procstat(1) man page or Guide to Parallel Vector Applications.
Procedure 5: Creating a report

Use the following procedure to view a report of complete I/O information for your code:

1. Compile and run the code.
2. Run the `procstat` utility on the code.

The following examples show how to perform this procedure:

CF90 example:

```bash
f90 prog.f90
procstat -r -m a.out
```

C++ example:

```bash
cc prog.c
procstat -r -m a.out
```

5.6 Evaluating Dynamic Memory Alternatives and Applying a Technique

If your code is memory bound, it will probably exhibit one of the symptoms listed in the following sections. Each section lists a symptom followed by a recommended technique to reduce the elapsed time caused by memory requests. Evaluate the behavior of your code to see if it matches one or more of these symptoms, and select one of the corresponding techniques. After applying any optimization technique, check your answers and examine the new elapsed time for the code.

5.6.1 Large Number of System Calls

Check the number in the Number of Calls to Memory Processor row in the Procstat Process Report that you created. Does the code have a large number of system calls? If possible, reduce the number of system calls for memory within the source code.

5.6.2 Memory Expanded or Contracted in Small Increments

Check the number in the Number of Calls to Memory Processor row in the Procstat Process Report. Does the code still make many system requests to expand or contract memory? The requests may be too small.
If so, apply one of the following techniques:

- Reduce the number of memory requests and increase the size of the requests within the code.
- Ensure that the first system call requests sufficient memory for prolonged usage, and minimize system calls to shrink the size of the process.
- Initialize a larger heap, as described in Section 5.7, page 62.

### 5.6.3 Other Reasons for Excessive Memory Activity

If there are still a large number of calls to the memory processor in the Procstat Process report, it is caused by any of the following situations:

- Alternate requests and releases of memory from the heap. Apply one of the following techniques:
  - Within your source code, attempt to reuse existing heap space instead of releasing it. Use the Fortran 90 `ALLOCATE` and `DEALLOCATE` keywords or the C++ `new` and `delete` operators or the `malloc` or `free` library functions. These do not require the compiler to generate a system call.
  - In Fortran code, avoid using an `SBREAK` statement with a negative argument. In C++ code, avoid calling the `sbreak` function with a negative argument.
  - Initialize a larger heap as described in Section 5.7, page 62.
- Frequent stack overflows and underflows (or stack thrashing). You can check this condition by using the Fortran `STOP` statement or the C++ `stkstat(3)` system call to produce a report of stack overflows. To avoid stack overflows, use the `SEGLDR` directive to increase the value in the `Initial stack size` row of Figure 10, page 66, to the maximum stack size as displayed by the `STOP` statement output. To create the Load Map Program Statistics report, see Section 5.7.2, page 63.

The following `STOP` statement output shows that the program experiences a large number of stack overflows:

```f90
where.f90
./a.out < INPUT_FILE

STOP executed at line 261 in Fortran routine ‘CLACIER’
CP: 34.435s, Wallclock; 198.094s, 2.2% of 8-CPU Machine
HWM mem: 236775, HWM stack: 10048, Stack overflows: 750000
```
By indicating a stack size that matches the stack high water mark (shown by the HWM stack value), stack overflows are now lower, and CPU time improves from 34.4 seconds to 13.5 seconds:

```
f90 -Wl"-S10048" where.f90
./a.out < INPUT_FILE
```

STOP executed at line 261 in Fortran routine 'CLACIER'
CP: 13.543s, Wallclock: 58.227s, 2.9% of 8-CPU Machine
HWM mem: 209338, HWM stack: 10048, Stack overflows: 0

5.6.4 Temporary Memory Expansion of Significant Duration

When the process expands temporarily during execution, it runs the risk of being swapped to disk, costing excessive elapsed time. Use a SEGLDR directive (see Section 5.7, page 62) to initialize the process at its largest size to avoid a swap.

5.6.5 Heap Blocks Released in Different Order than Allocated

Examine the source code. Does the process release heap blocks in a different order than they were allocated? This can cause memory fragmentation for the process. Reorder the code to allocate and deallocate heap space in the opposite order (last allocated should be first deallocated).

5.7 Memory Initialization

When forced to obtain more memory within the code, it is more efficient to use a few large requests than to use many small requests. One way to do this is to directly modify source code to issue fewer, larger requests. Another way is to initialize the heap with a SEGLDR directive. The most effective way to minimize the elapsed time due to memory management is to start with enough memory in the first place.

5.7.1 Loader Directives

At the load step, SEGLDR, which is usually called by both the f90(1) and CC(1) commands, establishes initial memory allocation for both the heap and the stack space. You can specify any of these parameters with a command-line option for segldr, CC, or f90, as in the following examples.
Example 1:
The following SEGLDR command specifies an initial heap size of 150,000 words and a heap increment of 75,000 words for the object file, `file.o`:

```
segldr -H150000+75000 file.o
```

Example 2:
The following Cray C++ compiler command specifies an initial stack size of 100,000 words and stack increment of 50,000 words for the source file, `file.C`:

```
CC -dSTACK=100000+50000 file.C
```

Example 3:
The following CF90 compiler command names a dynamic common block (DYNAM), specifies an initial heap size of 10,000 words, and establishes a zero heap increment size for the source file, `code.f90`. The heap increment is set to zero to force a fixed heap size required by using a dynamic common block:

```
f90 -Wl"-H10000+0;DYNAMIC=DYNAM" code.f90
```

For more information on SEGLDR see the Segment Loader (SEGLDR) and ld Reference Manual.

### 5.7.2 Optimal Heap Size

An optimal heap size for your process is a size that is large enough to prevent system calls for memory, but not so large that it contains unused memory.

#### Procedure 6: Determining optimal heap size

To determine an optimal initial heap size for your SEGLDR directive, perform the following steps:

1. Generate a ja report or retrieve the ja report you used in your initial analysis in Procedure 3, page 19.
2. Create a Load Map Program Statistics report for the code by using SEGLDR directives.

Example:
The sample load map in Figure 10, page 66, was generated by using the following f90 command:

```
f90 -Wl"-H10000+9000 -S6000+5000 -Mmap.fil,stat" code.f90
```

This command line specifies the following:

- Load map statistics in a file called `map.fil`
- Initial heap size of 10 Kwords
- Heap increment of 9 Kwords
- Initial stack size of 6 Kwords
- Stack increment size of 5 Kwords

3. Look at the number in the Memory HiWater column of the ja report you created for the code (Procedure 3, page 19). This number is reported in blocks, and a block is equal to 512 (decimal) Cray memory words. Multiply this figure by 512 to determine the maximum process memory size for the code in Cray words.

Example:

Assume the Memory HiWater figure from the ja report and the load map in Figure 10, page 66, are from the same code, `bigio`. This user process grew to 8,128 blocks. Perform the following arithmetic to obtain decimal words for Memory HiWater: 8,128 x 512 = 4,161,536 words.

4. Look at the number listed in the Base address of managed memory/stack row of the Load Map Program Statistics report (Figure 10, page 66). This number is an octal representation of the user area size measured in words. Convert this number to decimal to determine the user area size for the code in decimal format.

Example:

The Base address of managed memory/stack figure is 376,372 octal. This is equivalent to 130,298 decimal.

5. Subtract the user area size (total from the preceding step) from the maximum process memory size (total from Procedure 6, step 3, page 64). The result should be a good estimate for the minimum heap size required to avoid system calls for more memory.
Example:

4,161,536 - 130,298 = 4,034,238

This represents the largest heap size for the process. Note that this might change for a different dataset, and it will expand with larger library I/O buffer and user cache sizes.

6. Use the result from the preceding step in a SEGLDR directive to set the initial heap size.

Example:

To avoid system requests for memory, reload the object file with the following segldr command:

```
segldr -H4035000+10000 bigio.o
```
Program Statistics

Non-segmented object module written to: a.out

Allocation order: text, data, bss

Program origin: 0 octal 0 decimal

Program length: 422012 octal 140298 decimal

Start entry point is '$START' at address 10a

Transfer is to entry point 'POINTERVECTOR' at address 324d

Managed memory statistics

Initial stack size: 13560 octal 6000 decimal
Stack increment size: 11610 octal 5000 decimal
Initial managed memory size: 23420 octal 10000 decimal
Managed memory increment size: 21450 octal 9000 decimal
Available managed memory: 706 octal 454 decimal
Base address of managed memory/stack: 376372 octal
Base address of pad area: 376132 octal

Default libraries included:
/lib/libc.a
/lib/libf.a
/lib/libfi.a
/lib/libm.a
/lib/libp.a
/lib/libsci.a
/lib/libu.a

Figure 10. Load map statistics
This chapter describes the following techniques for optimizing I/O-bound code:

- Optimizing formatted I/O (Section 6.1, page 67)
- Optimizing large, sequential, unformatted I/O requests (Section 6.2, page 71)
- Optimizing small, sequential, unformatted I/O requests (Section 6.3, page 74)
- Optimizing for direct access I/O (Section 6.4, page 75)
- Optimizing asynchronous I/O (Section 6.5, page 77)
- Using an optimal storage device (Section 6.6, page 79)
- Minimizing system calls (Section 6.7, page 82)

For more information on I/O, see the Application Programmer’s I/O Guide.

6.1 Optimizing Formatted I/O

Formatted I/O is the slowest I/O and is useful only when the files must be viewed by people or transferred to systems other than Cray systems. However, if you are transferring the data to a system other than a Cray system, you can easily send the unformatted (binary) version instead of the formatted ASCII version by using the Cray foreign file conversion facility provided by the Flexible File I/O (FFIO) library. Use the techniques described in the following sections to optimize code that contains formatted I/O.

6.1.1 Changing to Unformatted I/O

If possible, change formatted I/O to unformatted I/O by using one of the following methods:

- In Fortran code, remove references to the FORMAT statement label and modify the Fortran OPEN statement to include FORM=’UNFORMATTED’.

  Example:

  ```fortran
  OPEN (10, FORM=’UNFORMATTED’)
  ```

- For C++ codes, cout << and cin >> are formatted read and write member functions of the iostream class. Also, the scanf(3) and printf(3)
function calls (including `fscanf`, `sscanf`, `fprintf`, and `sprintf`) require formatting to human-readable ASCII. Convert these functions to call unformatted I/O functions such as `fread` and `fwrite` instead. You can access FFIO by using the `ffread(3)` and `fwrite(3)` functions in your code in conjunction with the `assign(1)` command.

- To access the I/O layers provided by the FFIO libraries, use the `-F` command-line option with the `assign` command. This will provide access to the automated foreign file conversion.

### 6.1.2 Reducing the Amount of Formatted I/O

If you cannot change formatted to unformatted I/O, reduce the quantity of formatted I/O. Show only small samples of the data by using the following techniques:

- Change the code to show final results instead of many intermediate results.
- Change the code to show a checksum instead of the data itself.
- If the program sends data to another computer system (or printer), revise the program so that only the final version of the data is formatted.
- If you need to view the data, consider shipping it unformatted to a graphics postprocessor.

### 6.1.3 Increasing Formatted I/O Efficiency for Fortran Programs

Use the methods in the following sections to increase formatted I/O efficiency for Fortran programs.

#### 6.1.3.1 Minimizing the Number of Data Items in the I/O List

With the CF90 compiler you can increase formatted I/O efficiency by minimizing the number of data items in the I/O list. Consider the following example:

```fortran
DIMENSION X(20), Y(10), Z(5,30)
WRITE (6,101) M, (X(I), I=1,20), Z(M,J)
```

With vectorization turned off, this `WRITE` statement represents 22 data items. In this case, the `WRITE` operation would require 22 calls to the library routines that drive the `WRITE` statement. When vectorization is turned on, the compiler
treats each innermost implicit DO loop as a single data item, so that the preceding WRITE statement requires only 3 calls.

If you rewrite the statement as follows, the parameter list always represents 3 calls, even if all optimization is turned off:

\[
\text{WRITE (6,101) M, X, Z(M,J)}
\]

6.1.3.2 Using a Single READ, WRITE, or PRINT Statement

To increase formatted I/O efficiency for Fortran programs, read or write as much data as possible with a single READ, WRITE, or PRINT statement. Consider the following example:

\[
\begin{align*}
\text{DO } J &= 1, M \\
\text{DO } I &= 1, N \\
\quad \text{WRITE (42, 100) X(I,J)} \\
100 & \quad \text{FORMAT (E25.15)} \\
\end{align*}
\]

It is more efficient to write the entire array with a single WRITE statement, as follows:

\[
\text{WRITE (42, 100) ((X(I,J),I=1,N),J=1,M)}
\]

The following statement is even more efficient:

\[
\begin{align*}
\text{WRITE (42, 100) X} \\
100 & \quad \text{FORMAT (E25.15)}
\end{align*}
\]

Each of these three code fragments produce exactly the same output; however, the latter two examples are about twice as fast as the first. Also, the latter two examples are equivalent only if the implied DO loops write out the entire array, in order, and without omitting any items. You can use the format to control how much data is written per record.

6.1.3.3 Using Longer Records

To increase formatted I/O efficiency for Fortran programs, use longer records if possible. Because a certain amount of processing work is necessary to read or write each record, it is better to write fewer long records, rather than more short records. Consider the following example:
WRITE (42, 100) X
100 FORMAT (E25.15)

If you change it as follows, the resulting file will have 80% fewer records and, more importantly, the program will execute faster:

WRITE (42, 101) X
101 FORMAT (5E25.15)

Be careful to ensure that the resulting file does not contain records that are too long for the intended application. For example, certain text editors and utilities cannot process lines that are longer than a predetermined limit. Generally, lines that are not longer than 128 characters are safe to use in most applications.

6.1.3.4 Using Repeated Edit Descriptors

To increase formatted I/O efficiency for Fortran programs, use repeated edit descriptors whenever possible. For integers that fit in 4 digits (that is, less than 10000 and greater than —1000), avoid the following format:

200 FORMAT (16(X,I4))

Instead, use a format of the following form:

201 FORMAT (16I5)

6.1.3.5 Using Data Edit Descriptors that are the Same Width as the Character Data

To increase formatted I/O efficiency for Fortran programs, when reading and writing character data, use data edit descriptors that are the same width as the character data. For CHARACTER*n variables, the optimal data edit descriptor is A (or An). For Hollerith data in integer variables, the optimal data edit descriptor is A8 (or R8).

6.1.4 Increasing Formatted I/O Efficiency for C++ Programs

Calling a function increases overhead. To decrease overhead and increase formatted I/O efficiency for C++ programs, combine multiple calls to I/O functions into fewer calls. Consider the following example:

```c++
for (i=0; i<N; ) {
    fprintf(o1,"%d ",a[i];
    ++i;
    if (i%5 == 0) fprintf(o1,"\n");
    }
If you change it as follows, the resulting code will make 80% fewer calls to fprintf and the program will execute faster:

```c
for (i=0; i<N; i+=5) {
    fprintf(o2, "%d %d %d %d %d\n"
            a[i], a[i+1], a[i+2], a[i+3], a[i+4]);
}
```

### 6.1.5 Increasing Library Buffer Sizes for Formatted I/O Requests

For sequential-access formatted I/O files, the buffer size should be set equal to the length of a record or a multiple of that number. Generally, larger is better when buffering sequential access files.

To specify the library buffer size for Fortran, use the `assign` command with the following options:

```
assign -b size
```

For C++, use the `setvbuf` library function.

### 6.2 Optimizing Large, Sequential, Unformatted I/O Requests

*Sequential access* indicates that data items in a file have an implicit order. Unless the code issues positioning requests such as `fseek(3)` or `rewind(3)`, the system always accesses the next record automatically. If the code is issuing sequential, unformatted I/O requests larger than 1 Mword, use the techniques described in the following sections to optimize its I/O.

#### 6.2.1 Changing I/O File Format to Unbuffered and Unblocked

The default I/O file format for sequential unformatted Fortran I/O is COS blocked (`assign -s cos f: filename`), which means that the I/O request uses the library buffer and bypasses cache. Although the COS blocked file format helps fulfill the Fortran standard for sequential, unformatted I/O by marking (or blocking) record positions within a file, it is not the fastest I/O available for large, sequential transfers. COS blocked I/O requires user CPU time to create and insert (or interpret and remove) the control words.

If the code is issuing sequential, unformatted I/O requests larger than 1 Mword (8 Mbyte), change the I/O file format to unbuffered and unblocked by using the
-s u option, or by specifying the FFIO system, or syscall layer, as shown in
the following assign(1) command examples:

```
assign -s u f:filename
assign -F system f:filename
assign -F syscall f:filename
```

C++ codes can access the FFIO libraries by using the `ffread(3C)` and
`fwrite(3C)` I/O function calls in conjunction with the assign command.

Using the unbuffered, unblocked I/O file format requires you to construct
well-formed I/O requests in the code. These are simply I/O requests that begin
and end on disk sector boundaries, usually 512 words (4096 bytes) or a multiple
of 512 words. This unit of measurement is also known as a block or click. See
your system administrator to determine the sector size of the disks you are
using.

### 6.2.2 Converting to Asynchronous I/O

Converting to asynchronous I/O is a way to continue I/O activity in parallel
with the code’s CPU computation. If there are operations in the code that can
be executed while the code is waiting for I/O to complete, convert the code to
asynchronous I/O. For example, if the code contains any of the following
sequences, converting to asynchronous I/O might reduce elapsed time:

- Repetitive patterns of input, computation on that data, output, then input
  again
- I/O that appears in a loop

Most prominent sequential, unformatted I/O requests that consume a majority
of the code’s elapsed time will benefit from code conversion to asynchronous
I/O. You can convert to asynchronous I/O by using the assign(1) command
or by modifying your source code.

#### 6.2.2.1 Using the assign Command to Convert Code to Asynchronous I/O

The easiest way to convert code to asynchronous I/O is by using an FFIO layer,
either cachea or bufa, with the assign(1) command, as follows:

```
assign -F cachea:bs:nbufs f:filename
assign -F bufa:bs:nbufs f:filename
```
The `bs` argument specifies the size in 512-word blocks of each cache page or buffer. The `nbufs` argument specifies the number of cache pages (or buffers) to use. You can tune these arguments to better suit the I/O activities of the code.

If the code requires the use of COS blocked format, you can establish a specialized FFIO layer to provide asynchronous access by using the following `assign` command:

```
assign -F cos.async f:filename
```

### 6.2.2.2 Modifying Source Code to Convert Code to Asynchronous I/O

You can modify the source code to take better advantage of the asynchronous FFIO layer by breaking up a large I/O request into smaller iterative requests. Within the iterations, perform the necessary computation on that data. This technique is called *double-buffering*.

With double-buffering, two sets of data (buffers) are active at any given moment for each stream of input or output data. One buffer is active in CPU work, while the other is active in I/O (reading or writing). In a typical double buffer scheme, the I/O and CPU work sets are staggered, as in the following algorithm:

1. The first set of input data is read.
2. The second set of input data is read while the CPU works on the first set of input data.
3. The third set of input data is read while the CPU works on the second set of input data and the first set of data is output.
4. This sequence continues until all data is read. As the last data set is read, the next-to-last CPU work is in progress, and the third-from-last data set is output.
5. The CPU works on the last data set and the next-to-last data set is output.
6. The final data set is output.

**Example 1: C++ example of converting to asynchronous I/O**

In the following C++ example, the first input is the `reada(2)` system call in front of the `for` loop. Inside the loop, the first `recall(2)` system call synchronizes the previous `reada`, and the second `recall` synchronizes the previous `writea`. Notice that the `recall` system call is needed for synchronization. Generally, a second asynchronous system call to the same file descriptor will
not block execution for the first. This is called *queuing asynchronous I/O*, because each asynchronous request enters an I/O queue without blocking execution.

```c
#include <sys/types.h>
#include <sys/iosw.h>
#include <fcntl.h>
#define N 1001472
#define M 10
float a[N][2], b[N][2];
struct iosw sw[2], *prdsw[]={&sw[0]}, *pwrsw[]={&sw[1]};
int rfd, wfd, i, ird=0, iwk;
void work(float a[],float b[]);

main () {
    rfd = open("infile", O_RAW | O_RDONLY);
    wfd = open("outfile", O_RAW | O_CREAT | O_TRUNC | O_WRONLY, 0644);
    reada(rfd,(char *) &a[0][ird], N*sizeof(float), *prdsw, 0);
    for (i=0; i<M; i++) {
        iwk=ird; ird=(ird+1)%2;
        recall (rfd, 1, prdsw);
        if (i != M-1)
            reada (rfd, (char *) &a[0][ird], N*sizeof(float), *prdsw, 0);
        work(&a[0][iwk], &b[0][iwk]);
        if (i != 0) recall (wfd, 1, pwrsw);
        writea (wfd,(char *) &b[0][iwk], N*sizeof(float), *pwrsw, 0);
    }
}
```

### 6.2.3 Using Effective Library Buffer Sizes for Large, Sequential, Unformatted I/O

For large, sequential, unformatted I/O requests, enlarge the program’s library buffer to at least the size of its largest record, if possible. To specify the library buffer size for Fortran, use the `assign(1)` command with the following options:

```fortran
assign -b size f:filename
```

For C++, use the `setvbuf(3)` library function.

### 6.3 Optimizing Small, Sequential, Unformatted I/O Requests

If the code is issuing sequential, unformatted I/O requests that are 1 Mword or smaller, use the techniques described in the following sections to optimize I/O.
6.3.1 Using Effective Library Buffer Sizes for Small, Sequential, Unformatted I/O Requests

For small, sequential, unformatted I/O requests, use an effective library buffer size by ensuring that the library buffer is at least the size of the largest I/O request or a multiple of that size. For Fortran, use the `assign -b sz` command to specify the library buffer size. For C++, use the `setvbuf` library function.

6.3.2 Increasing I/O Request Size and Issuing Fewer Requests

To optimize small, sequential, unformatted I/O requests, increase the size of the I/O requests and issue fewer requests. This helps to reduce the overhead of both system and user CPU time and also may allow you to use the optimization techniques that apply to large I/O requests (see Section 6.2, page 71). You can use the following techniques to increase the size of the I/O requests:

- Read or write larger array sections instead of one element at a time.
- Combine read requests or write requests into a single read request or single write request.
- Extract I/O from inner loops.

6.3.3 Using the Memory-resident (MR) FFIO Layer

For small, sequential, unformatted I/O requests, if the file called by the code is heavily reused, the memory-resident (MR) layer in FFIO can improve performance over disk I/O by allowing the first portion of the file to reside in memory. For information on the MR layer, see Section 6.6.1, page 79.

6.4 Optimizing Techniques for Direct Access I/O

Direct access indicates that a program can access records or data at any point in the file. This also can be called nonsequential or random access I/O.

6.4.1 Fortran 90 Direct Access I/O

The Fortran 90 standard provides two types of access: sequential and direct. Sequential access restricts the program to reading from or writing to the I/O unit with records of any length in sequential order. Direct access divides the file associated with the I/O unit into fixed-length records, and allows the program to read or write records randomly. You can achieve Fortran direct access by opening a file with the `ACCESS=DIRECT` keyword on the `OPEN` statement and
specifying the fixed record size with the `RECL` keyword. All references to that file must specify the record number, `REC`, on subsequent `READ` and `WRITE` statements.

**Example 2: CF90 direct access example**

```fortran
OPEN (22, ACCESS='DIRECT', RECL=8000)
READ (22, REC=10) (DATA(I), I=1,1000)
WRITE (22, REC=2) (OUTNUM(J), J=1,150)
```

### 6.4.2 C++ Direct Access I/O

C++ programs do not use the I/O functions that transfer data to accomplish random access. C++ programs use the `fseek(3)` function or the `lseek(2)` system call to set the position in the file of the next input or output operation. The position is set in bytes, beginning at zero. Thus, C++ programmers are completely responsible for record keeping and indexing.

**Example 3: C++ direct access example**

```c++
stream = fopen ("file","r+");

bytes_per_word = 8;
nwords = 1000;

lrec = bytes_per_word*nwords;
fseek (stream,9*lrec,SEEK_SET);
fread (&data,bytes_per_word,nwords,stream);
fseek (stream,lrec,SEEK_SET);
fwrite (&outnum,bytes_per_word,150,stream);
fseek (stream, lrec - bytes_per_word*150,SEEK_CUR);
fread (&data,bytes_per_word,nwords,stream);
```

### 6.4.3 Optimizing Techniques for Direct Access Code

If the program is reading or writing files in direct access (as opposed to sequential access) you may be able to improve performance by using the following techniques:

- Ensure that the files are in binary file format and that they bypass the system cache by using the `assign -s bin` command.
- Ensure that the code is not using formatted or COS blocked file formats.
• Set the library buffer size as close to the length of a record (request) as possible without going under the length. This minimizes unnecessary data transfers, which detract from the performance of random I/O.

• For small, random I/O requests, use a smaller library buffer than the default. Limit its size to the record length of the code. This might improve performance by avoiding excessive unused data movement when filling the unused portion of the buffer.

• For large I/O requests, the library buffer size should be set equal to the length of the fixed-size record (request). To specify the library buffer size for Fortran, use the `assign -b sz` command. For C++, use the `setvbuf(3C)` library function.

• If the code makes repeated references to the same place in the data file, a memory-resident (MR) buffer might help if it can include the most frequently used area of data. For information on the MR layer, see Section 6.6.1, page 79.

• If the code uses word-addressable data, you can transfer the data faster with a binary file format (using the `assign -s bin` command), which also bypasses the system cache and forces use of the `GETWA` and `PUTWA` I/O routines without changing the source code. The `GETWA` and `PUTWA` I/O routines are among the fastest types of random-access I/O on Cray SV1 systems, but they place the burden of record keeping and indexing on you.

• Rearrange the data file so that the code can process it sequentially. Sequential I/O is usually faster than direct-access I/O. You might be able to use separate files to accomplish the same effect.

6.5 Optimizing Asynchronous I/O Requests

In most code, synchronous I/O is used more often than asynchronous I/O (also known as raw I/O). **Synchronous I/O** indicates that control is returned to the calling program after all requested data is transferred. The I/O transfer runs serially with respect to the CPU work.

**Asynchronous I/O** indicates that control is returned to the calling program after the I/O process has started, but before the I/O is completed. The I/O transfer runs in parallel with respect to the CPU work. The user program continues executing at the same time the I/O operation is executing.
If the code is using asynchronous I/O, use the techniques described in the following sections. Some of these methods increase CPU overhead but decrease total elapsed time if there is significant work to do during the I/O transfer.

### 6.5.1 Using Unblocked File Format for Asynchronous I/O Requests

To optimize asynchronous I/O requests, use unblocked file format if the code does not need to backspace, position the file pointer, read partial records, and so on. You can improve asynchronous I/O performance moderately by eliminating the overhead associated with record marking or blocking. This can be done in several ways, depending on the type of I/O and certain other characteristics.

For example, the following `assign` statements specify the unblocked file structure:

```plaintext
assign -s unblocked f:filename
assign -s u f:filename
assign -s bin f:filename
```

### 6.5.2 Avoiding Cache

For asynchronous I/O, avoid using cache by using the `assign -s u` command. This allows the data to transfer directly between the user process and the actual device without a stopover (with synchronization) in cache.

### 6.5.3 Using Effective Library Buffer Sizes for Asynchronous I/O Requests

If the program is using the default I/O file format for sequential unformatted Fortran I/O, which is COS blocked (with the `assign -F cos` command), to optimize asynchronous I/O requests, ensure that the largest record size is less than or equal to half the library buffer size. COS blocked I/O file format indicates that the I/O request uses the library buffer and bypasses cache.

Setting the library buffer size to an even number greater than 63 blocks causes COS blocked files to perform double-buffered asynchronous I/O by dividing the library buffer in half. When the library buffer size is an even number of disk sectors, each half of the buffer is well-formed. Thus, I/O requests for either half-buffer do not need to be rerouted through cache.

You can change the buffer size by using the `assign(1)` command to specify a special FFIO layer, as follows:

```plaintext
assign -F cos.async:size f:filename
```
6.5.4 Balancing Workload

Device I/O speeds are typically slower than CPU computation speeds by several orders of magnitude. If the code does not perform sufficient computation between I/O requests, it will spend most of its time waiting for I/O and lose the benefit of using asynchronous I/O. Try to balance both the I/O activity and the computation involving its data by moving as much of the CPU work as possible into the code that lies between asynchronous I/O requests.

6.5.5 Minimizing Required Synchronization

During asynchronous I/O processing, code reaches a synchronization point at which it has to wait for I/O completion before continuing. With an imbalance between CPU and I/O activity, this causes extended I/O wait time and an idle CPU. If this happens frequently, attempt to restructure the code to reduce required synchronization points.

6.5.6 Tune FFIO User Cache

If you are using asynchronous I/O through the cachea, bufa, or cos.async FFIO layers, you can adjust their sizes by using the assign(1) command. For complete information on controlling buffers and cache pages, see the Application Programmer's I/O Guide.

6.6 Using an Optimal Storage Device

If you have some flexibility with the storage devices your code uses, ensure that it uses the fastest devices available for the appropriate situations. The following sections describe storage devices and the situations in which they are best used.

6.6.1 Memory-Resident Files

Use memory-resident files for small requests, heavily reused files, or for large files in which most of the I/O activity occurs at the beginning of the file. The assign(l) command provides an option to declare certain files to be memory resident. This option causes these files to reside within the field length of the user’s process; its use can result in very fast access times.

To be most effective, this option should be used only with files that fit within the process’s field length limit. A program with a fixed-length heap and memory-resident files might deplete memory during execution. Sufficient space
for memory-resident files might exist, but might not exist for other run-time library allocations.

The memory-resident (MR) layer lets you declare that a file should reside in memory. The MR layer tries to allocate a buffer large enough to hold the entire file.

To use the MR layer for Fortran code I/O, use the following assign command and then rerun the program. To use the MR layer for C++ code I/O, use the FFIO system I/O calls in the code first (ffread, fwrite, and so on), and then use the assign command and rerun the program.

assign -F mr:size f:filename

The -F option invokes FFIO. The mr specification selects the memory-resident layer of FFIO. The size specification is the maximum size of the buffer in 512-word blocks. The filename specification is the name of the file.

6.6.2 Memory-Resident Predefined File Systems

Large memory systems might have predefined file systems resident in memory. Memory-resident file systems provide memory-to-memory speed, which is the fastest I/O available. Your system administrator can tell you which file systems are mounted in memory, and you might have access to create data files in those directories.

6.6.3 Disk Striping

If your file system is composed of partitions on more than one disk, using the disks at the same time can result in performance improvements. This technique is called disk striping. Disk striping can be accomplished through either hardware or software.

For example, if the file system spans three disks — partitions 0, 1, and 2 — it might be possible to increase performance by spreading the file over all three disks equally. Although 300 sequential writes might be required, assign only 100 to each disk and the disks can write simultaneously.

For hardware striping, your system administrator configures the disk and you can request that a file be opened on the striped disk during the OPEN system call.

Use software striping only for very large records, because all of the disk heads must do seeks on every transfer. Software striping can be useful for production
jobs that monopolize the system resources, but it can impede total system throughput if used on a heavily loaded multiuser system.

To specify software striping, consult with your system administrator to identify disk partitions for your file system. You can also use the df(1) command with the -p option, but always consult your system administrator before attempting this technique. You can achieve software striping by placing the desired file system partitions in the assign(1) command, as shown in the following two examples:

assign -p 0-2 -n 300:48 -b 144 f:filename
assign -p 0:1:2 -n 300:48 -F cos:144 f:filename

Factors such as channel capacity might limit the benefit of striping. Disk space on each partition should be contiguous and preallocated for maximum benefit.

6.6.4 Disk Arrays

Using disk arrays (for example, DA-60, DA-301, and so on) can be faster than single disk drives such as DD-60, DD-42, DD-301, and so on.

6.6.5 Disks

If possible, use disks only for files that are accessed one or two times or for saved files that are read at a later time. Try to use memory for most other activity.

Large disk files on a multiuser system can easily become fragmented across different disk cylinders and cause increases in I/O wait time from the device. Fragmentation causes a higher number of requests for disk space allocation and more physical seek requests. You can examine disk file fragmentation with the /etc/fck command. For more information, see the fck(1) man page.

Avoid disk file fragmentation by preallocating disk space for the data files that will be stored there. In Fortran 90, you can use the assign -n sz command. In C++, use the ialloc(2) system call after the file is opened. If it is necessary to allocate space prior to C++ program execution, use the setf(1) command. For more information, see the setf(1) man page.

The assign(1) and setf(1) commands will, by default, obtain contiguous allocation if it is available. If they do not succeed, you can force contiguous allocation by using the -c flag on either command, or by using the IA_CONT flag on the ialloc(2) system call.
6.6.6 Tapes

Consider tape to be a long-term storage device. Tape is both cost-effective and disaster-resistant. Before selecting tape, consider that it has slower access speed and that there is contention for the drive and delays for mounting. However, tape is appropriate for long-term archive storage of very large data files.

6.7 Minimizing System Calls

With few exceptions, system calls are required for all physical I/O requests and data movement to or from the library buffer. The following options minimize system calls:

- Ensure that I/O requests are as large as possible. For example, write whole arrays rather than one row at a time. Group multiple arrays into one write statement.

- Use larger buffers (or use cache) to capture many I/O requests in the user process space before the I/O library transfers the data out.

- Use scratch files for intermediate data that you no longer need after the code completes execution. This can eliminate unnecessary data movement and might avoid the device entirely.

  Scratch files are temporary and are deleted when they are closed. To create a Fortran scratch file, open a file with STATUS='SCRATCH' and use STATUS='DELETE'.

- Use the MR layer when appropriate (see Section 6.6.1, page 79).
alias
The alias shell command lets you define a synonym (a convenient name) for a command or command string. It lets you define a more mnemonic name for an existing command or a shorthand for a longer command string.

Autotasking
A trademarked process of Cray that automatically divides a program into individual tasks and organizes them to make the most efficient use of the computer hardware.

cache
A kind of temporary memory that speeds up load and store operations between memory and the registers of a processor.

chaining
A process of linking instructions together to save register storage time. Each instruction passes its results to the next linked instruction so that several operations may be done in approximately the same amount of time as one operation.

This function allows a vector register that is being used as a result register in one instruction also to be used as an operand register in a following instruction. By chaining vector instructions, overall speed is greatly increased; as soon as the first vector instruction has completed the function on element 0, that result is available to the second vector instruction as an operand. The first vector instruction does not have to complete processing on all vector elements before the second vector instruction can start processing.

chime
A sequence of vector operations that can be chained into a single pipeline. The limitation on such a sequence is that the same vector functional unit cannot be used twice in the same chain. Therefore, a loop that contains two vector adds, for example, contains at least two chimes because there is only one vector add functional unit.
**data dependence**
Occurs when the data that results from one segment of code depends on the
data that results from previous segments of code.

**disk striping**
Multiplexing or interleaving a disk file across two or more disk drives to
enhance I/O performance. The performance gain is function of the number of
drives and channels used.

**double-buffering**
Modifying source code with a technique by which you break up a large I/O
request into smaller iterative requests. Within the iterations, you perform the
necessary computations on that data.

**gather/scatter**
Either collecting data from multiple processors to a single processor (gather) or
dispersing data from a single processor to multiple processors (scatter). The
Fortran compiler multi-streams an ordered scatter (a scatter that involves
constant, as opposed to random, strides through the array being scattered).

**heap**
A section of memory within the user job area that provides a capability for
dynamic allocation. See the heap memory management routines in the library
documentation.

**induction variable**
A variable in a loop that controls the execution of that loop (for example, in
Fortran, \( i \) is the induction variable in \( \text{DO } I = 1, 100; \) \( \text{in C++}, i \text{ is the}
induction variable in \( \text{for } (i=0; i<100; i++) \)).

**inlining**
The process of replacing a user subroutine or function call with the subroutine
or function itself. This saves subprogram call overhead and may allow better
optimization of the inlined code. If all calls within a loop are inlined, the loop
becomes a candidate for vectorization, tasking, and multi-streaming.
load balancing
A process that ensures that each processor involved in a program performs roughly equal work.

loop counter
An integer variable that is incremented or decremented by an integer constant expression on each pass through the loop.

loop fusing
A code optimization technique by which two independent loops with the same iteration count are combined into one vector loop.

loop pushing
A code optimization technique by which a loop containing a subprogram call is moved into the called subprogram.

loop splitting
A code optimization technique by which a loop that contains both vectorizable work and scalar work is split into two loops: one that vectorizes, and one that does not.

loop unrolling
A code optimization technique in which the statements within a loop are replicated while reducing the trips through the loop.

loop unwinding
A code optimization technique in which a loop is unrolled so effectively that it is no longer a loop.

memory management
Management that allows memory to be allocated dynamically to programs while they are executing.

MSP
A multi-streaming processor capable of performing multi-streaming operations. It consists of four SSPs (single-streaming processors).
**multitasking**
An optimization method that incorporates multiple interconnected CPUs; these CPUs each run a part of a program simultaneously (in parallel) and share resources such as memory, storage devices, and printers. This term is used interchangeably with *parallel processing*.

**multi-streaming**
An optimization technique that automatically schedules one or more Cray SV1 multi-streaming processors (MSPs) for a program and divides the work for each MSP among four single-streaming processors (SSPS).

**parallel processing**
Processing in which multiple processors work on a single application simultaneously.

**parallel region**
An area within a program which multiple processors can execute in parallel. Its opposite is a *serial region*.

**pipelining**
A method of executing a sequence of instructions in a single processor so that subsequent instructions in the sequence can begin execution before previous instructions complete execution. This assembly-line approach to processing instructions is also called instruction pipelining or hardware pipelining.

**raw I/O**
A method of performing input/output in which the programmer must handle all of the I/O control. This is basically unformatted I/O. The opposite of raw I/O is *cooked I/O*.

**redundant code**
Code contained within a parallel region that is executed by all associated tasks, using the same data and generating the same results.

**scalar**
(1) In Fortran 90, a single object of any intrinsic or derived type. A structure is scalar even if it has a component that is an array. The rank of a scalar is 0. (2)
In C and C++, integral, floating, and pointer types are collectively called scalar types. (3) A nonvectorized, single numerical value that represents one aspect of a physical quantity and may be represented on a scale as a point. This term often refers to a floating-point or integer computation that is not vectorized; more generally, it also refers to logical and conditional (jump) computation.

**scalar processing**
A sequential operation in which one instruction produces one result; it starts an instruction, handles one operand or operand pair, and produces one result. Scalar processing complements vector processing by providing solutions to problems not readily adaptable to vector techniques.

**scalar temporary**
A simple variable defined and later referenced during each pass through a loop; it is not referenced outside the loop. The compiler can either replace a scalar temporary with a temporary vector, or it can eliminate it.

**stack**
(1) A data structure that provides a dynamic, sequential data list that can be accessed from either end; a last-in, first-out (push down, pop up) stack is accessed from just one end. (2) A dynamic area of memory used to hold information temporarily; a push/pop method of adding and retrieving information is used. (3) A portion of computer memory and/or registers used to hold information temporarily. The stack consists of stack frames that hold return locations for called routines, routine arguments, local variables, and saved registers.

**stack thrashing**
Frequent stack expansion (overflow) and contraction (underflow).

**stride**
An interval that is the same for all consecutive elements of a vector. On the Cray SV1 system, vectorization requires a constant, odd-numbered stride to perform at peak efficiency. An array is processed with a stride of 1, such as $A(1), A(2), A(3), ...$ is efficient. Avoid powers of 2, such as $B(2), B(4), B(6), ...$. A stride that is not constant is illustrated by a sequence such as $A(1), A(2), A(3), A(5), A(8), A(13)$. 
**tailgating**
Writing to a V register that is still being read from a prior vector instruction.

**user area**
A location at which the object code (text area) and external and static variables (data area) reside in memory.

**user CPU time**
Time accumulated by a user process when the process is attached to a CPU and executing. It is a fraction of wall-clock time, which measure the time from when you begin the execution of a program until execution completes. Wall-clock time includes the time a program is waiting for a CPU.

**user process**
An executable file becomes a user process after it is compiled and loaded. Your code is a UNICOS user process when it is executing.

**vector**
A computer vector is an array of numbers on which instructions operate; this can be an array or any subset of an array (such as a row, column, or diagonal). When arithmetic, logical, or memory operations are applied to vectors, it is referred to as vector processing.

**vector array reference**
An array element reference whose subscript expression is not a loop invariant. It is an array that is processed in vector registers.

**vector length**
The number of elements in a vector.

**vector processing**
A technique whereby iterative operations are performed on sets of ordered data. It provides results at rates exceeding the result rates of conventional scalar processing.
**vector register**
A vector (V) register is used for vector operations; successive elements from a V register enter a functional unit in successive clock periods.

**vectorizable expression**
An arithmetic or logical expression that consists of a combination of loop invariants, loop counters, vector array references, scalar temporaries, or a function with a vector version that has a vectorizable expression as an argument. This includes most Fortran intrinsic functions.

**vectorizable loop**
A loop that contains only vectorizable expressions (that is, expressions for which the compiler can produce vector code).

**vectorization**
Uses one instruction for the simultaneous performance of iterative operations on elements in sets of ordered data. It provides results at rates greatly exceeding those for conventional scalar processing, which works on only one element at a time.

**vectorized loop**
A source code loop that is processed with hardware vector registers.

**well-formed I/O requests**
I/O requests that begin and end on disk sector boundaries, usually 512 words (4096 bytes) or a multiple thereof.
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