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Parallelizing the Symbolic Manipulation Program FORM *

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Abstract

After an introduction to the sequential version of FORM and the mechanisms behind it we report on the status of our ongoing project of its parallelization. An analysis of the parallel platforms used is given and the structure of a parallel prototype of FORM is explained.

1. The Sequential Version of FORM

FORM [1] is a program for symbolic manipulation of algebraic expressions specialized to handle very large expressions of millions of terms in an efficient and reliable way. It is used non-interactively by executing a program that contains several parts called modules. The execution of each module is again divided into three steps:

- **Compilation:** the input is translated into an internal representation.
- **Generating:** for each term of the input expressions the statements of the module are executed. This in general generates a lot of terms for each input term.
- **Sorting:** all the output terms that have been generated are sorted and equivalent terms are summed up.

FORM only allows local operations on single terms, like replacing parts of a term or multiplying something to it. Together with a sophisticated pattern matcher this seemingly strong limitation allows the formulation of general and efficient algorithms. The limitation to local operations makes it possible to handle expressions as "streams" of terms, that can be read in sequentially from a file and be worked on one at a time. The generation of terms is done in a way that the output terms drop out term by term also. These output terms are stored in two intermediate buffers and a temporary sortfile in a staged procedure: When the smaller buffer gets full, its content is sorted and copied to the larger buffer. Consequently the small buffer is free to be filled with the next patch of output terms. If the larger buffer gets full, its content is again sorted and copied to the temporary file, freeing the large buffer. At the end of the module all of the existing sorted patches residing in the two buffers and the sortfile have to be merged stage by stage to one single sorted output stream of terms which is written to file and used as input source of terms for the next module. For the sorting of the small buffer a slightly modified merge sort algorithm is used; merging sorted patches of terms in the other stages of sorting is done with a tree of losers [2]. This results in a tree-like structure for the generation of terms as well as for the sorting.

Of course the first step before parallelizing the program was to profile and optimize the sequential code. One of the main achievements was a speedup of about a factor 2 on the 64-bit architecture of the alpha processors by changing the internal used word-length from 16 (which results in 32-bit arithmetic operations) to 32 bits (64-bit arithmetics). Another speedup of about 1.5 was simply achieved by experimenting with the compiler optimizations. Profiling of FORM in typical applications shows that the time needed for compiling the program text into the internal representation is not dependent on the size of the problem and usually neglectible. Most of the time is spent in generating and sorting of terms.

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2. Evaluation of Different Parallel Platforms

One of the suppositions for the parallelizing of FORM has been not to limit the approach by using a too specialized hardware, but instead to use standard message passing libraries (MPI(CH) [3] and PVM [4]) for the parallelization that are available on a wide class of architectures and should be portable to new and more powerful systems. Usually both message passing libraries use specialized device drivers underneath to yield maximum performance. During the first part of our project the following combinations of hard– and software have been used:

- DEC alpha workstation cluster running DEC UNIX 4.0D,
  8 processors, 600MHz, 512MB RAM and 2×4GB disk each,
  interconnection: Fast (100 MBit) & Gigabit (1000 MBit) Ethernet,
  1.26 GBit Myrinet [5], 1.26 GBit ParaStation II [6],
  message passing: MPI(CH) and PVM (over IP and ParaStation II).

- IBM SP2 running AIX 4.2.1,
  168 (in total 256) processors, 120MHz, 512MB RAM,
  interconnection: special low latency switches,
  message passing: MPI.

- ALR Quad6 SMP machine running Solaris 2.6,
  4 processors, 200MHz PentiumPro, 512MB RAM and 2×4GB disks,
  interconnection: shared memory,
  message passing: MPICH (over shared memory).

In the case of the DEC alpha cluster different IP drivers and different implementations of device drivers for the messages libraries exist. For a thorough understanding of the parallel systems behaviour it is of course crucial to compare the performance of the messages passing libraries. The MPI(CH) libraries have been examined using the Pallas MPI benchmarks (PMB) [8]. A shell script is provided that can be used to measure the throughput and latency of the message passing operations under different circumstances without further interaction. We only present the results for the simplest possible single transfer benchmark, a ping-pong test: One process(or) sends a message of n bytes to another process(or), which immediately sends that message back. In contrast to more complicated benchmarks there is no concurrency with other messages passing activity during this test. In other words, the bandwidth and latency are measured under optimal conditions.

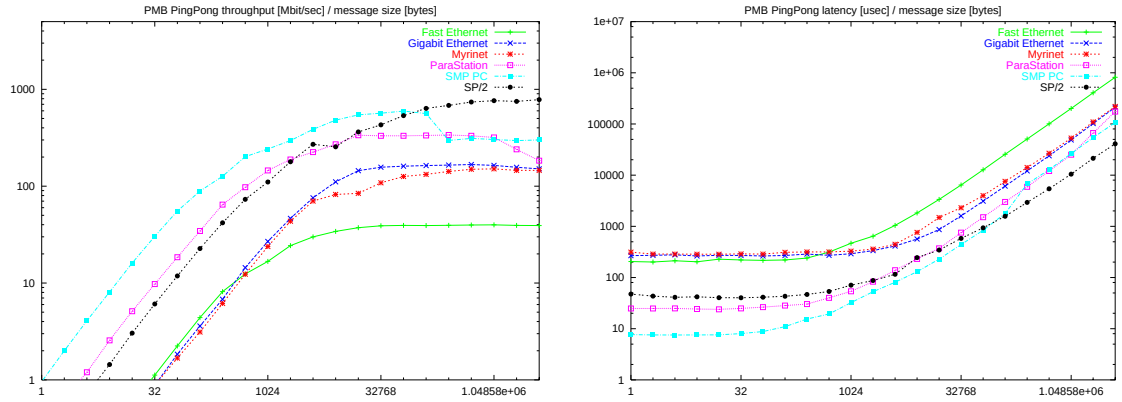

Figure 1: MPI(CH) bandwidth and latency results for the PingPong benchmark.

As can be seen from figure 1 (left) the bandwidth increases as the message size does. For large message sizes there is a saturation effect. The maximum throughput for the Fast Ethernet is about 40 MBit/s, for the Gigabit Ethernet and the Myrinet the maximum bandwidth is about 200 Mbit/s. The ParaStation II software provides a special device driver for MPI(CH) on the Myrinet hardware that does not use the IP protocol and adds much less protocol overhead. With this spezialized software a bandwidth of about 300 MBit/s is achieved. The IBM SP2 reaches a maximum MPI bandwidth of 800 MBit/s, which is even higher than the maximum MPI(CH) throughput on the SMP machine that uses shared memory segments for the data transfer. In figure 1 (right) the measured MPI(CH) latency for the ping-pong benchmark is shown. The latency ranges from a few microseconds for small message sizes up to a second for large message sizes. Using MPI(CH) over the IP protocol results in a minimum latency of more than 200 microseconds. Lower latencies can be achieved with special device drivers only. It turns out that the minimum latency on the ParaStation II (about 20 microseconds) is even smaller than on the IBM SP2. Of course the SMP machine has the smallest latency—less than ten microseconds. Note that for our particular interest the latency at larger message sizes is most important. In this region again the IBM SP2 provides optimal performance.

3. The Parallelization of FORM

Allowing local operations only makes FORM very well suited for a straightforward parallelization: distribute the input terms among the available processors, let each of them perform the local operations on its input terms and generate and sort the arising output terms. At the end of a module the sorted streams of terms from all processors have to be merged into one final output stream again. The compiling of the program-text to the internal representation was considered not to be worth parallelizing. This concept indicates to use a master-slave structure for the parallelization, where the master would store the expressions and distribute and recollect all the terms of each expression. For the implementation of this raw concept we used a four step strategy:

- one process(or) generates terms, a second process(or) sorts the output terms
- instead of only one process arbitrary many processes perform the sorting.
- the input terms are distributed and the term generation is also done in parallel
- final optimizations: avoid or handle worst cases, load leveling, fault tolerance

This approach has several advantages, the most important being that having working versions in every stage gives us a good idea of how good the parallelization is and the possibility of doing realistic tests even in a very early stage.

3.1 The Two-Processor Version

This first step turns out to be useful, since it not only gives a deep insight into how changes to the source code of FORM have to be made without affecting the efficiency of the well optimized sequential code. The two-processor version also serves as a check of whether and how the concept can possibly lead to a decent speedup, because we basically add communication overhead (no speedup can be expected from seperating the generation and the sorting of terms only). It turns out that for parallelizing software on a cluster of very fast workstations the importance of avoiding communication overhead cannot be overstated. Moreover these experiments show that the communication has to be done in a buffered way, since sending single terms increases the runtimes of the two-processor code up to a factor 20. With the buffered version the increase in runtime was limited to about a factor of 1.5 of the sequential code with two workstation connected by a (slow) 10 Mbit/s Ethernet using the PVM and MPI(CH) libraries.

3.2 Parallel Sorting

The second step is to distribute the output terms among arbitrary many processors and do the sorting in parallel. Since this part of the sorting relies strongly on communication between the processors, it most probably sets the limits of parallel speedup. Therefore already in this stage it could be predicted quite reliably whether a speedup of the whole program could be achieved or not. A first try was to map the "tree of losers" used in FORM to merge sorted patches onto the processors. While it would distribute the workload in an optimal way this approach adds too much communication overhead. This is why in the end a much simpler communication structure was used, where all slaves send their sorted terms to the master process and this process uses a local "tree of losers" to merge the output streams of the slave-processes. Additional effort was made to overlap the work on the master process with the sorting done on the slaves, which caused a much deeper interference with the sequential code, but finally resulted in a very fast and stable implementation.

Of course the scaling of this approach with a large number of processors is a possible limitation, but tests on the 256 node IBM SP2 parallel computer showed that, at least for such a specialized network, the performance does not dramatically break down up to 32 processors. If problems should occur with more processors or a different architecture, an intermediate layer of "foremen" could be used to circumvent this possible bottleneck.

Since we are only testing the sorting, any problem that produces a sufficient number of intermediate terms can be used to test the parallel speedup. We wrote a very simple program, that expanded the expression $(a_1 + \cdots + a_n)^2$ and then replaced a_1 by $-(a_4 + \cdots + a_n)$ which results in a short, easy to check result and, by choosing different values of n, can be scaled in an easy way. The runtimes we could achieve with this version on the alpha cluster with different combinations of communication soft/hardware are shown in figure 2 for a medium ($n = 3000$, ~ 500 MB) and a large ($n = 5000$, ~ 1400 MB) problem (the sizes in MB are for the 64-bit architecture of the alphas, they are half as large for 32 bit systems (a more efficient compression to reduce disk access is in progress). They show that only for sufficiently large problems a satisfying speedup can be achieved without sophisticated network soft/hardware.

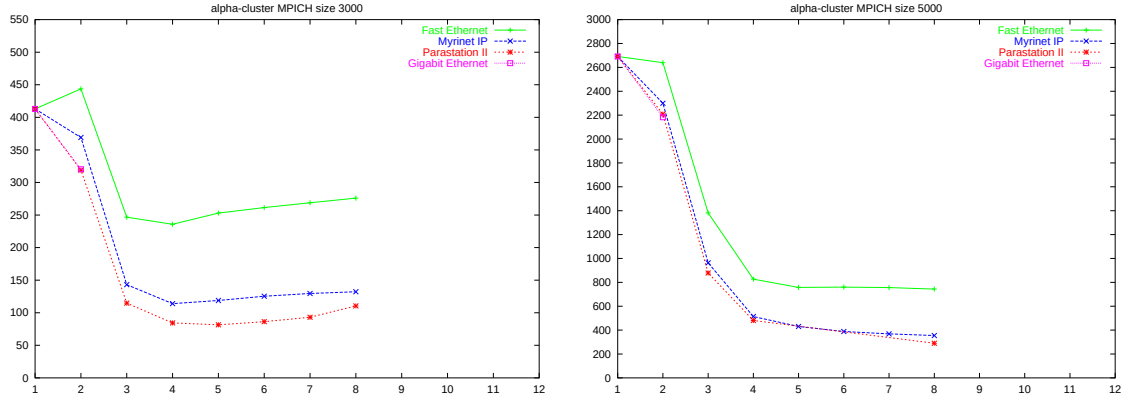

Figure 2: Runtimes [sec] for parallel sorting on the alpha cluster, shown is the runtime vs. number of processors for different communication soft/hardware.

The parallel sorting was implemented in a way that can take advantage of the non-blocking unbuffered MPI message passing functions by using a set of cyclic buffers which eliminates any unnecessary overhead and also minimizes the time consumed for synchronizing. Choosing the number of cyclic buffers to be one results in using the send/receive function in blocking mode. On a 4 processor SMP machine we could see a speedup of about 10% for some cases using this feature, while on a IBM SP2

there are no differences between the two versions (figure 3).

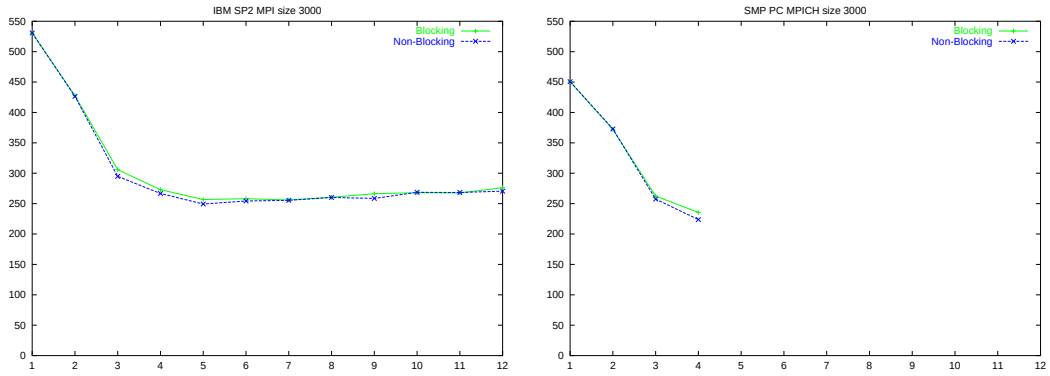

Figure 3: Runtimes [sec] for parallel sorting on two other architectures: IBM SP2 (left) and Intel SMP (right) shown are the version with blocking and nonblocking communication.

It is also quite interesting to see that obviously the master process can generate just enough terms to keep about 4 processes busy, almost independently of the problem size, the difference is that the cost of additional overhead for more than 4 processes is larger for the small problems, which of course is no surprise. It was know from the profiling of the sequential code that a large part of the runtime is spent in the generation routines, so the speedup we could see for very large problems is about the optimum we could expect to reach at this intermediate stage.

3.3 Parallel Generating

The last step towards a working prototype was the distribution of input terms among all processes before generation of the output terms. The slave-processes of course also need all the information necessary for the generation of the terms, which is at the moment realized by having them all read the program text file and compile their own internal representation and broadcasting only the rest of necessary information from the master process.

It was one of the main goals to get this version to run realistic problems as soon as possible, so the often used FORM-package MINCER [9], which can calculate certain types of Feynman diagrams up to three loops was used to serve as a testbed. First the necessary subset of FORM commands used in this package was made available and some easy standard integrals served as a test of the correctness of the parallel computations. After these preliminary tests the computation of some diagrams of an ongoing project, the calculation of the 3-loop triple-gluon vertexfunction, are now serving as "real world" tests.

Just as with the sorting the speedup is strongly dependent on the number of terms sent within one message. In the current implementation this number can be changed at the start of the program and results in very different speedups, as can be seen in figure 4. Of course choosing a too coarse grained distribution will result in the danger of running into worst cases, where all the work sits in only one of the input patches, and only one processor is busy. On the other hand, the fine grain distribution causes more overhead. The best setting turned out to be not only dependent on the underlying soft/hardware, but also strongly dependent on the problem that is run. The distribution is organized such that the master sends a patch of terms to each slave process at the beginning of the module and then waits for the slaves to ask for new terms whenever they are finished with their last patch of input terms. This actually turns the concept in that of a client-server situation which will also be useful to make the slaves receive any kind of global information when the need arises. It also produces a decent load levelling among the slaves, which can be controlled by the size of the

input-patches and could even be adjusted during runtime for further improvement.

It must be understood that for this specific problem there were now over hundred modules executed, where most of them were only seeing 1 to 10 terms, which makes them "worst cases" for the parallelization and there is a lot of room for further optimizations of this parallel program. Also the results are received from the FORM-code written for the sequential version of FORM without any modification or optimization in these algebraic packages, which corresponds to a perfect code reuse.

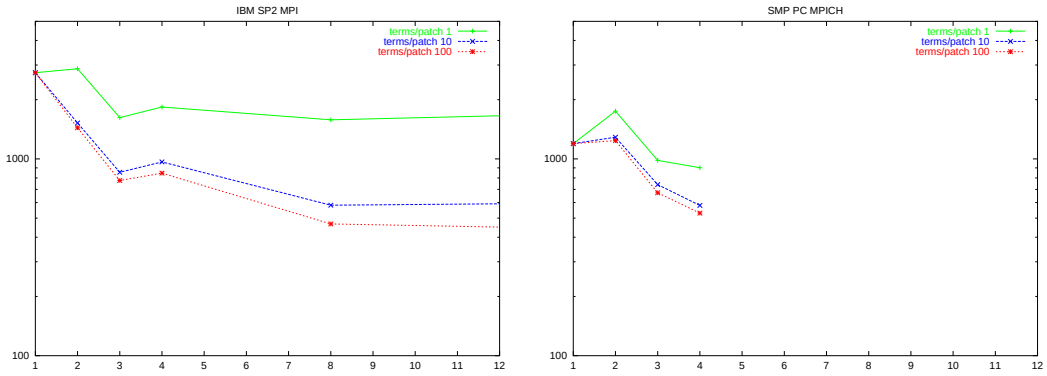

Figure 4: Runtimes [sec] vs. number of processors for a "simple" 3-loop Feynman-diagram on a IBM SP2 (left) and a 4 processor SMP machine (right) using its proprietary MPI-library. Shown are runtimes for different granularities of the input-term distribution.

As was expected, on the slow nodes of the IBM SP2 with fast connections a speedup is quite easy to obtain (see figure 4, left). Also it can be seen that for the fully parallelized program there is indeed a speedup up to 12 processors on the SP2. Another result is that the number of input terms that are distributed at once has a large influence on the speedup that can be achieved. Figure 4 (right) shows the runtimes that could be achieved on a 4 processor SMP machine. Also on that architecture the number of input terms sent at a time must be at least about 10 to see a speedup.

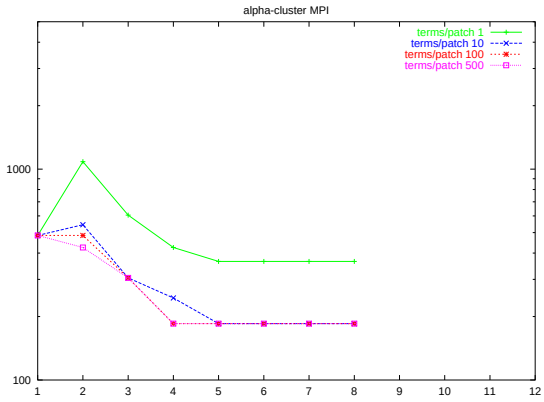

Figure 5: Runtimes [sec] vs. number of processors for a "simple" 3-loop Feynman-diagram on the DEC alpha cluster using MPI(CH) over the ParaStation II hard/software. Shown are runtimes for different granularities of the input-term distribution.

As can be seen in figure 5 this holds also for the DEC alpha cluster using MPI(CH) over the ParaStation II hard/software. Obviously a speedup is much harder to obtain in this case. Still the parallel

version on the IBM SP2 and the 4-processor SMP machine does hardly reach the runtimes measured with the sequential version on the DEC alpha cluster. From our experience with the parallel sorting and preliminary tests we expect a much larger speedup for larger problems, which are of course most interesting and investigated at the moment.

4. Conclusion & Outlook

The ongoing work is mainly to find and fix bugs to make the parallel version to run larger problems just as extremely reliable as the sequential one, which has been tested and improved for over ten years. As the runtimes show, there is also some work to be done to further optimize the parallel version, especially to insure that modules with only few terms will not cause a slowdown by unnecessary communication overhead. The next step will be the implementation of the full FORM version 2.3 standard, so that all existing software for that version can be run in parallel (which is basically all software that exists for FORM). Since the parallel program is actually based on version 3.0 which is in preparation by the author J. Vermaseren and offers some new and powerful features, we also will investigate if and how these new features can be implemented in the parallel version. Another field of current and planned activities is porting the program to other architectures. We are especially interested in taking better advantage of the possibilities of SMP machines. The main goal of all these efforts is to get from the current stage of a working prototype to an easy to use, powerful and reliable program that is not an end in itself, but a useful tool in real life applications on a wide variety of (parallel) architectures in the same manner as the current sequential version of FORM.

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