

PARALLEL IMPLEMENTATION OF BREADTH-FIRST SEARCH

Markus Püschel

Department of Computer Science
ETH Zürich
Zürich, Switzerland

ABSTRACT

Describe in concise words what you do, why you do it (not necessarily in this order), and the main result. The abstract has to be self-contained and readable for a person in the general area. You should write the abstract last.

1. INTRODUCTION

Motivation. Breadth-first search (BFS) is a graph traversing algorithm which is used in many applications. Some examples are finding connected components, determining a shortest-path for navigation systems or computing the maximum flow in a network. As real-world graphs can get very large, for example the US road network graph contains roughly 24 million vertices and 57 million edges, the need for fast implementations is high.

In this paper, we will discuss multiple approaches and compare their performance on three different platforms.

Related work. In this section we will give a brief overview of related work. In the first part we will discuss the work covering the top-down-bottom-up hybrid approach and in the second part an approach that focuses on avoiding atomic operations.

Beamer et al.[1] report a different algorithm to deal with the performance issues encountered when designing a BFS algorithm. The proposed hybrid algorithm combines the usual top down approach with a new bottom up part. In the bottom up part a level is processed by searching a parent for all unvisited vertices where a parent is only valid if it is a neighbour of the unvisited vertex. This is advantageous for small-world graphs because it saves accesses and data processing when a large fraction of the vertices are in the frontier. To get optimal results a hybrid algorithm is proposed where a heuristic switching criteria controls the use of top down or bottom up step depending on the size of the frontier and a predicted size of the next frontier. Yasui et al.[2] describe an implementation of such a hybrid algorithm for kronecker and R-MAT graphs as well as a detailed description of the heuristic switching parameters.

The author thanks Jelena Kovacevic. This paper is a modified version of the template she used in her class.

Berrendorf[3] describes a technique to avoid atomic operations in a generalized scenario. The scenario is given as an if-statement followed by some operations that change a state, where multiple threads might execute the predicate and execute the operations afterwards. The operations need to change the state to the same value if executed multiple times otherwise there exists a race condition, i.e. the change of the distance of a visited vertex to the value of the level or the addition of a vertex to the next frontier. The trade-off is that doing a BFS this way can result in additional work, since any unvisited vertex may get added multiple times.

2. BACKGROUND: BREADTH-FIRST SEARCH

In this section we will give a brief overview over the idea of breadth-first search and its sequential asymptotic runtime cost and a short discussion of different graphs and graph properties.

Breadth-first search.

Breadth-first search (BFS) is a graph traversal algorithm which starts at a source and either travels until it finds a specific vertex or until it has explored all connected vertices. In the first step all vertices adjacent to the source are explored and stored in some data structure (called frontier or next) as well as marked as visited. In the second step the newly visited vertices become the new sources (called neighbours or current) from which the search continues by repeating this step. Doing a traversal in this way assures that all the nodes at the same distance to the source are explored on the same level before any vertices with greater distance can be explored.

The desired output of a BFS can differ depending on where it is applied. For example with minimal modifications BFS can deliver a predecessor map, where every vertex points to only one parent, or a distance map, where the distance to the source for every vertex is saved. Since the predecessor map is not necessarily unique, we choose to return a distance map as our output to make verification of correctness simpler.

The sequential version of a BFS can be implemented using a single queue and has a theoretical asymptotic runtime

of $\mathcal{O}(|V| + |E|)$ where $|V|$ is the number of vertices and $|E|$ is the number of edges of the connected graph.

Graphs.

3. YOUR PROPOSED METHOD

We implemented many different algorithms and multiple variants for most of them. They return a distance map from one source vertex to all reachable vertices in the graph.

All our implementations are based on OpenMP for synchronization.

Topdown. A lot of our approaches are based on a simple topdown algorithm. The idea is to do a level-synchronous traversal of the graph by keeping record of the vertices in the current as well as those in the next level (“frontier” and “neighbour”) in two data structures of the same type and setting the frontier to the current neighbours after each level. This leads to an implicit barrier, as the work done by the different threads has to be synchronized. We experimented with different data structures and synchronization methods.

The naive version of the topdown algorithm has a global standard vector for the frontier to allow easy dynamic splitting between the threads in each level. This balances the load between threads, however it has a significant overhead because it relies on a critical section (`OMP critical`) for checking whether the vertices were already visited and inserting them into the neighbour data structure.

To improve the naive implementation, we used atomics, the built-in `__sync_val_compare_and_swap` (CAS) to atomically check whether a vertex was visited and set the correct distance. It also uses a local neighbourhood data structure (standard vector) to prevent needing a critical section. Only at the end of each level, a lock (`omp_lock_t`) is used to combine the local neighbourhoods to a global one, which can then be distributed between the threads for the next level. This is an idea adapted from Berrendorf [3].

An extended version of the algorithm before first uses a non-atomic check if a vertex visited before each CAS.

Also inspired by Berrendorf [3] is the idea to remove atomics altogether. This results in a race condition where additional work for the next level is created, as some vertices might be added to the neighbours multiple times. However, expensive atomics are omitted, which results in a trade-off between the additional time from the added work and the faster runtime by leaving out the CAS. This algorithm still uses a local neighbourhood and a lock when combining them.

Instead of using a standard vector for the frontier and the neighbour, one of our implementation relies on a bool array for all data structures (visited vertices, frontier, neighbour). This makes it possible to remove all critical sections, as the insertion of an element into a vector was what made them necessary in the first place. Similarly to the algorithm be-

fore, there might be additional work due to avoiding atomics. The downside of this implementation is that in each level, you have to loop through all the vertices, not just the current frontier, to be able to explore from there. Depending on the structure of the graph, this can be much more work.

4. EXPERIMENTAL RESULTS

Here you evaluate your work using experiments. You start again with a very short summary of the section. The typical structure follows.

Experimental setup. We run experiments on three different platforms.

The first platform is the EULER cluster which is operated by the HPC Group of ETH. We had access to one node with a 12-core Intel Xeon E5-2697v2 processors (2.7 GHz nominal, 3.0-3.5 GHz peak). It supports hyper-threading, so we ran our algorithms with up to 24 threads. On this platform, we used the gcc compiler with the `-O2` flag.

The next platform we ran our algorithms on is the Xeon Phi provided through the class.

Lastly, we also used an AMD FX-8350 (4GHz x8, 8Gb, W2k3).

Results. Next divide the experiments into classes, one paragraph for each. In each class of experiments you typically pursue one questions that then is answered by a suitable plot or plots. For example, first you may want to investigate the performance behavior with changing input size, then how your code compares to external benchmarks.

For some tips on benchmarking including how to create a decent viewgraph see pages 22–27 in

Comments:

- Create very readable, attractive plots (do 1 column, not 2 column plots for this report) with readable font size. However, the font size should also not be too large; typically it is smaller than the text font size. An example is in Fig. 1 (of course you can have a different style).
- Every plot answers a question. You state this question and extract the answer from the plot in its discussion.
- Every plot should be referenced and discussed.

5. CONCLUSIONS

Here you need to summarize what you did and why this is important. *Do not take the abstract* and put it in the past tense. Remember, now the reader has (hopefully) read the report, so it is a very different situation from the abstract. Try to highlight important results and say the things you really want to get across such as high-level statements (e.g., we believe that is the right approach to Even though

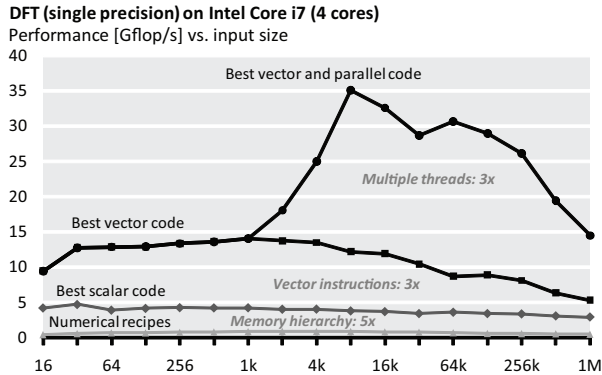


Fig. 1. Performance of four single precision implementations of the discrete Fourier transform. The operations count is roughly the same. The labels in this plot are maybe a little bit too small.

we only considered x, the technique should be applicable) You can also formulate next steps if you want. Be brief. After the conclusions there are only the references.

6. FURTHER COMMENTS

Here we provide some further tips.

Further general guidelines.

- For short papers, to save space, I use paragraph titles instead of subsections, as shown in the introduction.
- It is generally a good idea to break sections into such smaller units for readability and since it helps you to (visually) structure the story.
- The above section titles should be adapted to more precisely reflect what you do.
- Each section should be started with a very short summary of what the reader can expect in this section. Nothing more awkward as when the story starts and one does not know what the direction is or the goal.
- Make sure you define every acronym you use, no matter how convinced you are the reader knows it.
- Always spell-check before you submit (to us in this case).
- Be picky. When writing a paper you should always strive for very high quality. Many people may read it and the quality makes a big difference. In this class, the quality is part of the grade.
- Books helping you to write better:

- Conversion to pdf (latex users only):

```
dvips -o conference.ps -t letter -Ppdf -G0 conference.dvi
and then
ps2pdf conference.ps
```

Graphics. For plots that are not images *never* generate the bitmap formats jpeg, gif, bmp, tif. Use eps, which means encapsulate postscript. It is scalable since it is a vector graphic description of your graph. E.g., from Matlab, you can export to eps.

The format pdf is also fine for plots (you need pdflatex then), but only if the plot was never before in the format jpeg, gif, bmp, tif.

7. REFERENCES

- [1] Scott Beamer, Krste Asanovi, and David A Patterson, "Searching for a parent instead of fighting over children: A fast breadth-first search implementation for graph500," 2011.
- [2] Y. Yasui, K. Fujisawa, and K. Goto, "Numa-optimized parallel breadth-first search on multicore single-node system," in *Big Data, 2013 IEEE International Conference on*, Oct 2013, pp. 394–402.
- [3] R. Berrendorf, "A technique to avoid atomic operations on large shared memory parallel systems," *International Journal on Advances in Software*, vol. 7, no. 1 & 2, pp. 197–210, 2014.