Breadth First Search Optimization

High Performance Programming with Multicore & GPUs

By Chen-Yu Chang

Graph Traversal

Definition

Graph traversal refers to the process of visiting each vertex in a graph. Such traversals are classified by the order in which the vertices are visited.

Procedure

Graph

- → Adjacency List
- → Adjacency Matrix
- → Undirected Graph

Find Connections

→ Match from parent nodes

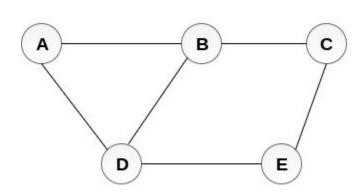
Check Visited

- → Record when the node is visited
- → Prevent from repeating

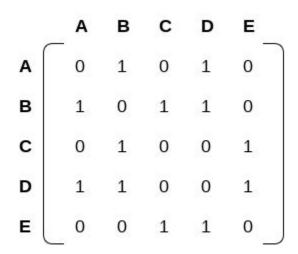
How Fast: O(|V|+|E|) (Worst Case)

where V (vertices), E (Edges)

Graph Traversal



Undirected Graph



Adjacency Matrix

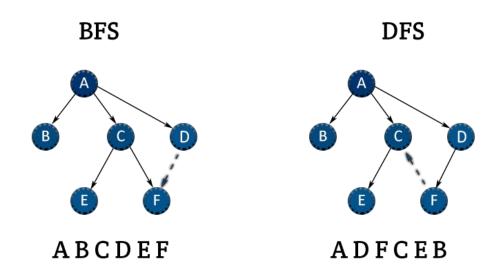
Breadth First Search

Definition

BFS is an algorithm for traversing or searching tree or graph data structures. It starts at the tree root, and explores all of the neighbor nodes at the present depth prior to moving on to the nodes at the next depth level.

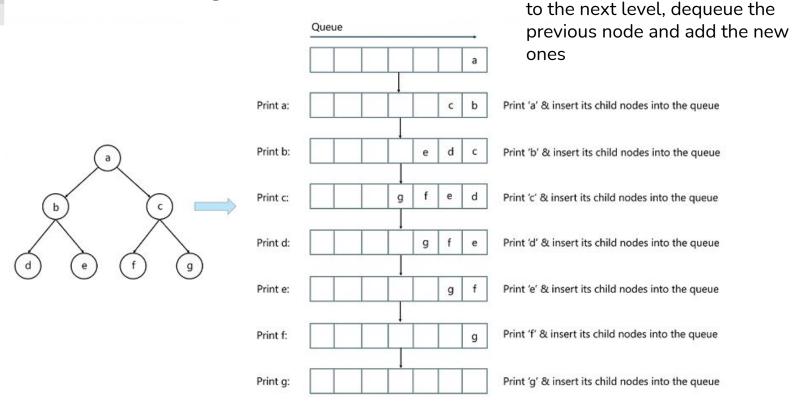
Compare to DFS

Unlike BFS, DFS explores the node branch as far as possible before being forced to backtrack and expand other nodes.



Serial Method

Method: Queue



Key: When the node is visited,

add to the queue. After going

Image Source: https://www.edureka.co/blog/breadth-first-search-algorithm/

Challenges

- ★ Need to set a high capacity for he queue if we operate BFS for huge graph
- ★ In huge graph, BFS will operate slowly since all the other levels need to wait for the previous levels
- \bigstar Memory access is another constraint : O(V²)
 - Including Parents, Visited, Queue, and Graph arrays

Hybrid Algorithm

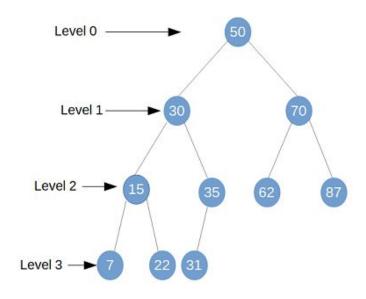
Definition

Set the targeted node and find the nodes that are adjacent (same distance to the root node). Group them based on breadth.

Advantage

★ Instead of node by node, the whole algorithm categorize the nodes into groups

★ Potential improvement can be operating multiple nodes at the same time



Top-Down Method: O(E)

```
\label{eq:function top-down-step} \begin{split} & \textbf{for } v \in \text{frontier } \textbf{do} \\ & \textbf{for } n \in \text{neighbors}[v] \ \textbf{do} \\ & \textbf{if } parents[n] = -1 \ \textbf{then} \\ & parents[n] \leftarrow v \\ & next \leftarrow next \cup \{n\} \\ & \textbf{end if} \\ & \textbf{end for} \\ & \textbf{end for} \end{split}
```

- → Efficient on small frontier
- → Loop the current frontier and its neighbors
- → If not visited, add to the next frontier

Down-Top Method: O(NV+E)

```
function bottom-up-step(vertices, frontier, next, parents)

for v \in \text{vertices do}

if parents[v] = -1 then

for n \in \text{neighbors}[v] do

if n \in \text{frontier then}

parents[v] \leftarrow n

next \leftarrow next \cup \{v\}

break

end if

end for

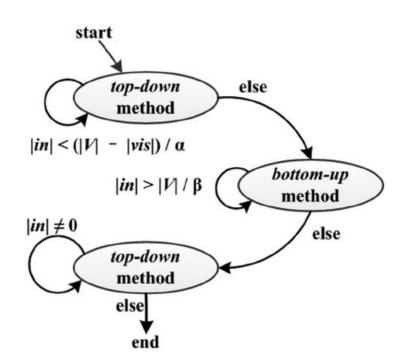
end if

end for
```

- → Efficient on large frontier
- → Find the nodes' parents that are not in the current frontier

Hybrid

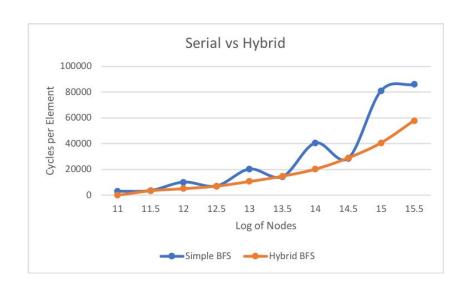
```
function hybrid-bfs(vertices, source)
   frontier \leftarrow {source}
   next \leftarrow \{\}
   parents \leftarrow [-1,-1,\cdots,-1]
   while frontier \neq \{\} do
       if next-direction() = top-down then
           top-down-step (vertices, frontier, next, parents)
       else
           bottom-up-step (vertices, frontier, next, parents)
        end if
       frontier \leftarrow next
       next \leftarrow \{\}
   end while
   return parents
end function
```



Data Comparison

By comparing the cycles per element, we gain the data from using the serial code & hybrid code.

→ Overall, the hybrid approach seems to decrease the iterations and shows a better control on frontiers



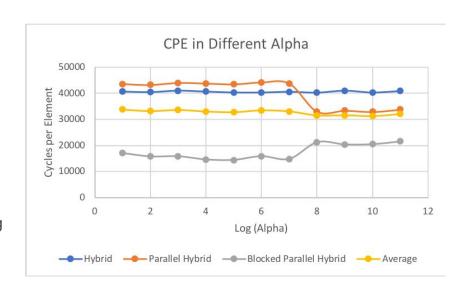
a (Alpha)

Determine when the frontier is too large

From the data:

- → When there are 32768 nodes, the optimized Alpha is at about 1024.
- → It also shows the logical result in the graph on the right.

Usage: Compensate for bottom-up finishing



β (Beta)

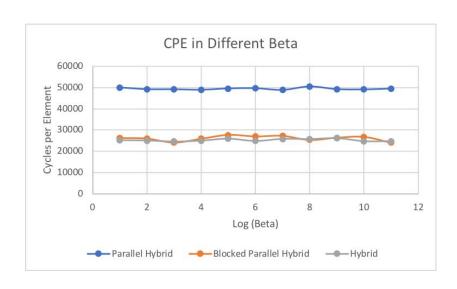
Less important comparing to alpha

Check whether the frontier has appropriate size

From the data:

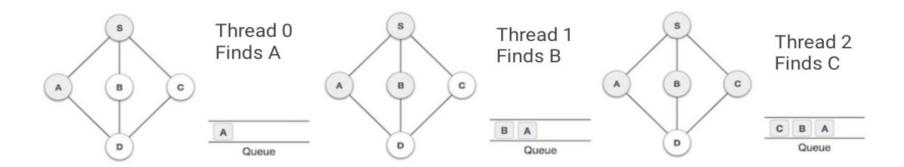
→ When there are 32768 nodes, the optimized Alpha is at about 16.

Usage: Compensate for up-bottom finishing



Parallelization

Serial Method with Parallelization: Threads



We see an adjacent node from S, which is A, B, C. For example, we find A and enqueue it.

Then, we enqueue and mark visited B that is adjacent from S.

Finally, the next adjacent node C is visited and enqueue.

Serial Method with Parallelization: Threads

The head of the queue is shared to all threads

The threads will split into several neighbors to visit

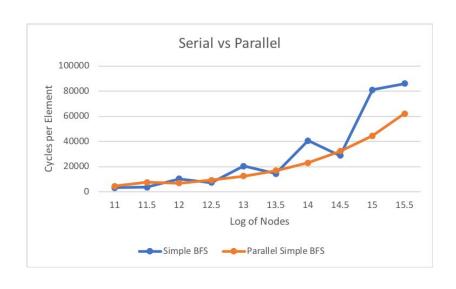
- The threads will wait for the next update of the queue; otherwise, it stays in rest
- Queue is locked

Data Comparison

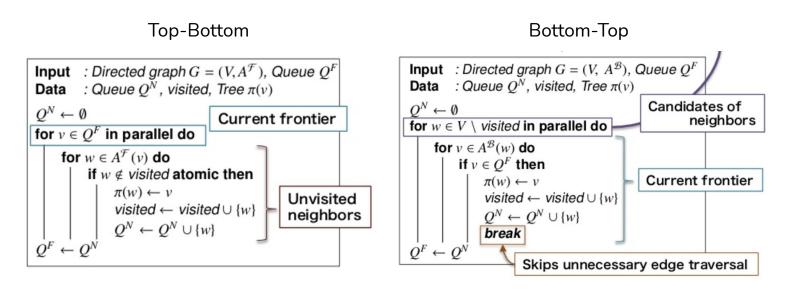
Overall, the performance of parallelization improves. However, the serial implementation is quite unstable, which sometimes performs better than the parallel implementation.

Reason:

It sometimes spends more time waiting.



Hybrid Algorithm with Parallelization



Make good use of barriers: each level of nodes

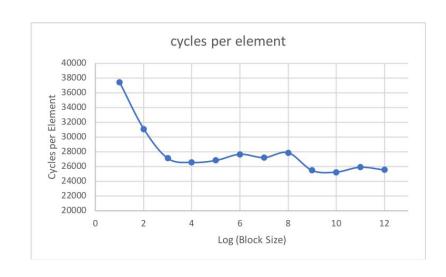
Blocked Parallel Hybrid

→ Temporal Locality

From the data:

→ We found the optimized blocked size is 1024 for 32768 nodes.

Overall, as blocked size increase, the number of iterations decreases.



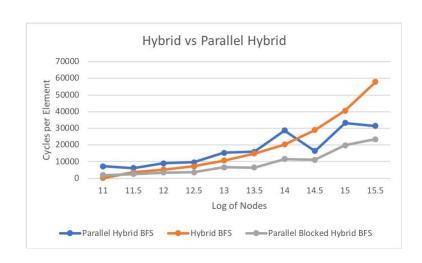
Data Comparison

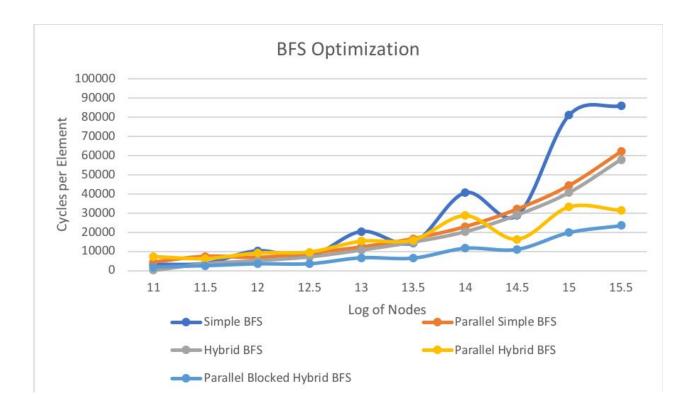
By comparing serial hybrid, parallel hybrid, and parallel blocked hybrid:

When there are large number of nodes:

- Multi Threads are efficient solution
- Parallel steps also helps

Overall, parallel blocked hybrid works the best.





In summary, the efficiency: (Under Large Number of Nodes)

Parallel Blocked Hybrid > Parallel Hybrid > Hybrid > Simple Parallel > Simple

Dijkstra's Algorithm (GPU)

Definition

Dijkstra's algorithm found the shortest path between two given nodes, but a more common variant fixes a single node as the "source" node and finds shortest paths from the source to all other nodes in the graph.



	BFS	Dijkstra
Main Concept	Visit nodes level by level based on the closest to the source	In each step, visit the node with the lowest cost
Optimality	Gives an optimal solution for unweighted graphs or weighted ones with equal weights	Gives an optimal solution for both weighted and unweighted graphs
Queue Type	Simple queue	Priority queue
Time Complexity	O(V + E)	O(V + E(logV))

- Breadth-first search is just Dijkstra's algorithm with all edge weights equal to 1.
- Kernel call to initialize all node weights to infinity except for the source node. We mark the source node as settled after this point.
- Kernel Call to choose a node which is relaxed and settled and to relax the outgoing edges of each settled node.
- Kernel call to mark all the settled nodes.



- Kernel Call to initialize the values of the node weights to infinity except for the Source Node which is set to 0
- We mark the source Node to be settled after that and all other nodes unsettled
- Kernel Call to choose the node with minimum node weight whose edges are to be relaxed
- Kernel call to relax all the edges of a node

Algorithm 1 GPU implementation of Dijkstra's algorithm. CUDA kernels are delimited by <<< ... >>> .

```
 \begin{array}{lll} \text{1: } <<<&\text{initialize}>>> (U,F,\delta); & \textit{//Initialization} \\ \text{2: } & \textbf{while } (\Delta \neq \infty) & \textbf{do} \\ \text{3: } & <<<&\text{relax}>>> (U,F,\delta); & \textit{//Edge relaxation} \\ \text{4: } & \Delta =<<<&\text{minimum}>>> (U,\delta); & \textit{//Settlement step\_1} \\ \text{5: } & <<&\text{update}>>> (U,F,\delta,\Delta); & \textit{//Settlement step\_2} \\ \text{6: } & \textbf{end while} \\ \end{array}
```

Algorithm 2 Pseudo-code of a CUDA thread in relax kernel.

```
1: tid = thread.Id;

2: if (F[tid] == TRUE) then

3: for all j successor of tid do

4: if (U[j] == TRUE) then

5: BEGIN ATOMIC REGION

6: \delta[j] = \min\{\delta[j], \delta[tid] + w(tid, j)\};

7: END ATOMIC REGION

8: end if

9: end for

10: end if
```

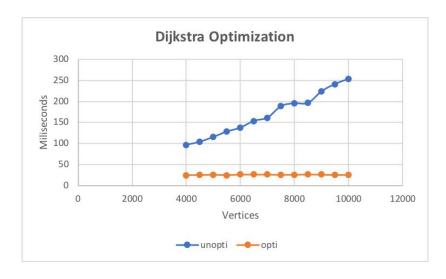
Algorithm 3 Pseudo-code of a CUDA thread in update kernel.

```
    tid = thread.Id;
    F[tid]= FALSE;
    if (U[tid]==TRUE and δ[tid] <= Δ) then</li>
    U[tid]= FALSE;
    F[tid]= TRUE;
    end if
```

Data Comparison

Under Edge per Nodes = 4000 & Weight of each Edge = 5:

After Optimization, the number of vertices does not seem to have influence on the runtime.



Implementation on GPU

Serial vs Parallel (GPU)

With 1024 threads, 10000000 vertices, 50000000 edges

Sequential Execution Time: 3.81 sec

Kernel Execution Time: 0.02636 sec

Speed up by: 144.554 times

Thank you