Final Project BDDs Re-ordering

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https://github.com/cheng-hsiang-chiu/ECE6740-CAD



Problem

• Minimize the node sizes of the BDDs by re-ordering the variable in the BDDs.

Steps

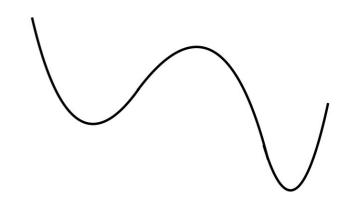
Read in an BLIF file

Convert BLIF format into BDD

 Pick the re-ordering variable that results in a minimum number of nodes in the BDD by simulated annealing algorithm.

Algorithm - Simulated Annealing (SA)

```
while(TEMPERATURE < FROZEN TEMPERATURE) {
iterate(MAX ITERATIONS PER TEMPERATURE) {
    generate_prop_ordering()
     delta_cost = prop_node_size - curr_node_size
     if(delta cost < 0) {
         curr_ordering = prop_ordering
         if(prop_node_size < best_node_size) {</pre>
             best_ordering = prop_ordering
     else {
         if(exp(-delta_cost/TEMPERATURE) > rand()) {
             curr_ordering = prop_ordering
 TEMPERATURE *= DEGRADE
```



Generate proposed ordering

Randomly pick a strategy

- Strategy 1: shuffle variables in a random range.

Strategy 2: swap two random variables.

- Strategy 3: reverse the current ordering.

Strategy 4: swap first half and second half.

Experimental Environments

- Ubuntu 20.04
- Intel(R) Xeon(R) Gold 6138 CPU @ 2.00GHz
- L1d cache: 1.3 MiB
- L1i cache: 1.3 MiB
- L2 cache: 40 MiB
- L3 cache: 55 MiB
- CPU(s): 80
- Memory: 256GiB
- C/C++17

Experimental Results - 1

• TEMPERATURE = 1000, MAX_ITERATIONS_PER_TEMPERATURE = 1000, FROZEN_TEMPERATURE = 0.1, DEGRADE = 0.9

	Time (secs)	My_ node_ size	Memor y
test_L14	18	5	42 MB
testXOR	18	4	42 MB
adder8	29	82	44 MB
mult6	111	1116	43 MB
c499	152122	45922	2.1 GB
c880	54187	51863	1.2 GB

	Default_ node_ size	ABC_ node_ size
test_L14	6	6
testXOR	4	4
adder8	1267	104
mult6	1158	444
c499	45922	800
c880	346660	654

Experimental Results - 2

• TEMPERATURE = 1000, MAX_ITERATIONS_PER_TEMPERATURE = 100, FROZEN_TEMPERATURE = 0.1, DEGRADE = 0.9

	Time (secs)	My_ node_ size
test_L14	2	5
testXOR	2	4
adder8	3	95
mult6	11	1116
c499	14077	45922
c880	5309	85242

	Default_ node_ size	ABC_ node_ size
test_L14	6	6
testXOR	4	4
adder8	1267	104
mult6	1158	444
c499	45922	800
c880	346660	654

Experimental Results - 3

• TEMPERATURE = 1000, MAX_ITERATIONS_PER_TEMPERATURE = 10, FROZEN_TEMPERATURE = 0.1, DEGRADE = 0.9

	Time (secs)	My_ node_ size
test_L14	0.2	5
testXOR	0.2	4
adder8	0.3	174
mult6	2.1	1158
c499	1533	45922
c880	531	92524

	Default_ node_ size	ABC_ node_ size
test_L14	6	6
testXOR	4	4
adder8	1267	104
mult6	1158	444
c499	45922	800
c880	346660	654

Experimental Results - Performance vs Execution time

• TEMPERATURE = 1000, MAX_ITERATIONS_PER_TEMPERATURE = ?, FROZEN_TEMPERATURE = 0.1, DEGRADE = 0.9

	1000	100	10
test_L14	5 / 81 sec	5 / 2sec	5 / 0.2 sec
testXOR	4 / 77 sec	4 / 2 sec	4 / 0.2 sec
adder8	82 / 117 sec	95 / 3 sec	174 / 0.3 sec
mult6	1116 / 111 sec	1116 / 11 sec	1158 / 2.1 sec
c499	45922 / 152122 sec	45922 / 14077 sec	45922 / 1533 sec
c880	51863 / 54187 sec	85242 / 5309 sec	92524 / 531 sec

Experimental Results - Performance Improvement

	Default	ABC	Mine
test_L14	0%	0 %	16.67%
testXOR	0%	0%	0%
adder8	0%	91.79%	93.53%
mult6	0%	61.66%	3.76%
c499	0%	98.26%	0 %
c880	0%	99.81%	84.37%

Conclusions

 SA parameters, such as TEMPERATURE, DEGRADE, and FROZEN_TEMPERATURE, could be learned by more test cases.

More iterations may not give more satisfying results.

 To examine more solution space and save more execution time, the work could be implemented in a parallel manner with OpenMP, CUDA, OpenCL, etc.

Contribution

• Contribute lines 56 – 470 in ./src/main.c

 Some lines in functions initialize() and calculate_node_size() are copied from original github project.

Thank You