(d) [15 points] If we modify the vector processor to *support chaining*, how many cycles would be required to execute the same program in part (c)? Explain.

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9 GPUs and SIMD [45 points]

We define the SIMD utilization of a program that runs on a GPU as the fraction of SIMD lanes that are kept busy with active threads during the run of the program. As we saw in lecture and practice exercises, the SIMD utilization of a program is computed across the complete run of the program.

The following code segment is run on a GPU. Each thread executes a single iteration of the shown loop. Assume that the data values of the arrays A and B are already in vector registers, so there are no loads and stores in this program. (Hint: Notice that there are 3 instructions in each iteration.) A warp in the GPU consists of 32 threads, and there are 32 SIMD lanes in the GPU.

Please answer the following six questions.

(a) [2 points] How many warps does it take to execute this program?

```
33 warps.  
Explanation:  
The number of warps is calculated as:  
\#Warps = \lceil \frac{\#Total\_threads}{\#Warp\_size} \rceil,  
where  
\#Total\_threads = 1025 = 2^{10} + 1 \text{ (i.e., one thread per loop iteration)},  
and  
\#Warp\_size = 32 = 2^5 \text{ (given)}.  
Thus, the number of warps needed to run this program is:  
\#Warps = \lceil \frac{2^{10}+1}{2^5} \rceil = 2^5 + 1 = 33.
```

(b) [10 points] What is the *maximum* possible SIMD utilization of this program? (Hint: The warp scheduler does *not* issue instructions when *no* threads are active).

 $\frac{1025}{1056}$

Explanation:

Even though all active threads in a warp follow the same execution path, the last warp will only have one active thread.

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