Basic Pollard Rho Algorithm Implementation On CUDA Device

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1 Introduction

The factorisation problem of a huge number resolved into massive parallel solutions. A Large number of algorithms are currently state-of-art and are continuously developed into better forms. This short article is focused on implementation of the Pollard-Rho algorithm on CUDA device. In the first part, there is a definition of algorithm. Next, the article focuses on options of parallelism on CUDA device. Results are measured in multiple instances and compared in graphs.

2 Definition of the algorithm

The ρ algorithm (named after the shape of curves symbolising two functions trying to reach themselves in the projective space) is based on finding the cycle. In t random numbers of x_1, x_2, \ldots, x_t in range [1, n] will contain repetition with probability of P > 0.5 if $t > 1.777n^{\frac{1}{2}}$

The ρ algorithm uses g(x) with modulo n as a generator for pseudo-random sequence. In this article function $x^2+k;\ k\in\mathbb{N}$. We are assuming that $n=p\cdot q$ that also mean $p<\sqrt{n}$ and $q<\sqrt{n}$. Algorithm actually generates sequence of $x_1=g(2),\ x_2=g(g(2)),\ \dots$ in two separated sequences running in same time. One sequence is generated as $x_1=g(x_0)\mod n$ and second as $x_1=g(g(x_0))\mod n$. Since we know that $p<\sqrt{n}$ the faster sequence is likely to cycle faster then the sequence generated just in in application of a g function. The repetition of the p will be detected with p which will divide p without any residue [2].

```
 \begin{array}{l} x \! = \! 2; \; y \! = \! 2; \; d \! = \! 1 \\ \text{while } d \; \text{is } 1 \colon \\ x \! = \! g(x) \\ y \! = \! g(g(y)) \\ d \! = \! \gcd(|x - y| \;, \; n) \\ \text{if } d \! = \! n \colon \\ \text{return } failure \\ \text{else} \colon \\ \text{return } d \end{array}
```

Figure 1: Sequential pseudo-code of algorithm

While the implementation in Python is basically just a copy of pseudo-code, in C there is plenty struggle with fix size integer size. Another problem will be with the transformation of libraries like GNU MP, that's why the basic focus on sequential solution was on basic mathematical operations. For the basic adding and subtracting there was no bigger effort than just creating basic implementation with carry bits. Much more time was spent on operations like division and modulo.

2.1 Division

The first idea on how to divide numbers was to use school *Long division* algorithm which proved to be a little bit expensive but was behaving in the same manners for every input numbers.

```
\begin{array}{l} Q \,=\, 0 \\ R \,=\, 0 \\ \text{for } i \,=\, n \,-\, 1 \,\ldots\, 0 \,\text{ do} \\ R \,=\, R \,<<\, 1 \\ R(0) \,=\, N(\,i\,) \\ \text{if } R \,>=\, D \,\text{ then} \\ R \,=\, R \,-\, D \\ Q(\,i\,) \,=\, 1 \end{array}
```

Figure 2: Long division [1]

Another much faster division was to use *Knuth's division* [3] which took a considerably shorter time to compute. Yet there was no luck in implementation. Current version which was considered is from 11vm [4]. But it failed with some badly shifted bits in residuum hence was not used.

2.2 GCD

Since the division wasn't fast enough, at least GCD was implemented with shifts and small bit checks. This may lead to more divergent behaviour in CUDA implementation.

```
inline void gcd(unsigned int * A, unsigned int * B){
    unsigned int t [SIZE];
    unsigned int shift;
    if(zeroNum(B)){
        return;
    if (zeroNum(A)) {
        copyNum(A, B);
        return;
    for (shift = 0; ((A[0] \mid B[0]) \& 1) == 0; ++ shift)
        shiftRightNum(A);
        shiftRightNum(B);
    while ((A[0] \& 1) == 0){
        shiftRightNum(A);
    do{
        while ((B[0] \& 1) = 0){
            shiftRightNum(B);
        if(bigger(A, B) == 1){
            copyNum(t, B);
            copyNum(B, A);
            copyNum(A, t);
        subNum(B, A);
    } while (! zeroNum(B));
    shiftLeftNumBy(A, shift);
}
```

Figure 3: GCD algorithm

2.3 Measurements

To make some basic idea on how fast the algorithm is, there was picked large number composed with big prime numbers $N = 0x0fd42d4eb2c4b7b1 = 1140586421661513649 = 1067982407 \cdot 1067982407$. Whole run of sequential implementation shows that there is only 54443 iterations to finish the task.

Figure 4: Command time output on sequential run

CPU which was used was AMD Phenom(tm) II X4 960T Processor.

3 CUDA Solution

3.1 First solution with explicit barrier

First idea which was found at https://github.com/dghost/factor-cuda. Implementation prepared at host memory X, Y, C for the algorithm which was, later on, ran with multiple kernels and explicitly synchronised after each loop. For the parameters X, Y, C were used random numbers which could let to same sequences.

From this, there was an easy step to create deterministic X, Y, C just by giving them numbers from 1 till the number of running threads. This actually led to creating a kernel which prepared data directly at the global memory of device without the need of host CPU. Algorithm work as follow:

```
create R variable which should, later on, contain the result
copy input N to the device memory space
call kernel which creates input variables X, Y, C
   for each thread in each block
while R is zero:
   call kernel which is one iteration of Pollard Rho cycle
   copy R from device to host memory space
```

Figure 5: First version of Pollard-Rho on CUDA

The whole approach is based not on dividing the cycle into same size small problems but rather trying different starting positions and hoping some of them reach the result much faster. That would mean the algorithm could be in the worst case as same fast as the implementation in the single thread.

3.2 Measurements

For measurement in this case was chose small number:

```
N[0] = 0xb2c4b7b1;
N[1] = 0x0fd42d4e;
```

The number is really small and at the most complex operation like a power of two, it would take about 5 times 32 bits. That would also raise a question of bank conflicts. Yet with using a huge number of blocks and threads, this wouldn't need to be a problem. Another change was to force using L1 cache. Since number size is not huge, caching could provide huge speed up.

```
cudaFuncSetCacheConfig(prepareDataKernel, cudaFuncCachePreferL1);
cudaFuncSetCacheConfig(pollardKernel, cudaFuncCachePreferL1);
```

Figure 6: Changing caching policy for both kernels

With that, the average time of execution was about 11505.4 [ms]. Graph of all runs does not show any significant speed up in a particular number of blocks and number of threads. Only large slow down at blocks size of 10 with 10 threads which is obvious.

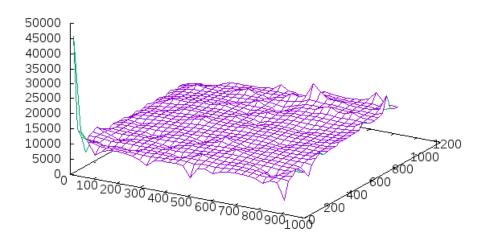


Figure 7: Graph of execution time

Some small speedups were noticed in cases where particular input X, Y, C were given. In these cases, the algorithm could end much faster because of simple luck of choosing good input. From opposite point of view, some cases were so unlucky with choosing initial values that some small slow downs were noticed as well.

3.2.1 Profiler output

Profiling was done on NVidia GeForce GTX 780 Ti at gpu-02.star.fit.cvut.cz.

Another problem would maybe occur in the divergence of threads in warp. The algorithm works with non-native mathematical operations which are used for big numbers. With big probability, the divergences occur at modulo operation. Because of large usage of while in GCD the operation could also cause a lot of divergences as well. For this reasons, profiler testing is provided showing about 50% of the runtime is given to waiting.

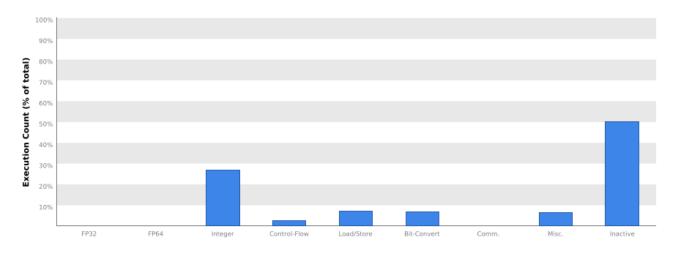


Figure 8: Inactivity of threads in warp

There are further indications of efficiency. Profiler provided the count of instructions which were used:

Cuda Fuctions :

pollardKernel(unsigned int const *, unsigned int*, unsigned int*)

Maximum instruction execution count in assembly: 12478876
Average instruction execution count in assembly: 716726
Instructions executed for the kernel: 1811166793
Thread instructions executed for the kernel: 31660047474
Non-predicated thread instructions executed for the kernel: 29193971935
Warp non-predicated execution efficiency of the kernel: 50,4%
Warp execution efficiency of the kernel: 54,6%

Figure 9: Count of instructions used by kernels

Some other techniques were used. For example $_$ restrict $_$ was given to data pointers and const was added to N which is untouched in the whole algorithm. Yet the biggest problem of diverging threads in warp wasn't fully solved.

For further detail see pollardKernel_v1.pdf.

3.3 Second solution with independent runners

Another idea was to omit using global memory at all. So the threads would run algorithm from different input values X, Y, C and after some time each one of them would check if the there is some result. If some thread found solution whole algorithm stop.

```
__global__ void pollardKernel(const unsigned int * N, unsigned int * result){
    unsigned int threadID = blockIdx.x * blockDim.x + threadIdx.x;
    unsigned int X[SIZE];
    unsigned int Y[SIZE];
    unsigned int C[SIZE];
    unsigned int G[SIZE];
    unsigned int N_tmp[SIZE];
    unsigned int abs_mxy[SIZE];
    cuda setZero(X);
    X[0] = 0x07;
    cuda setZero(C);
    C[0] = threadID + 1;
    cuda setZero(G);
    G[0] = 0x01;
    \operatorname{cuda}_{\operatorname{copyNum}}(Y, X);
    cuda fxfun(N, Y, C);
    unsigned int check = 0;
    while (cuda_isOne(G)){
         cuda_fxfun(N, X, C);
cuda_fxfun(N, Y, C);
cuda_fxfun(N, Y, C);
         if(cuda bigger(X, Y) == 1){
             cuda\_copyNum(abs\_mxy\,,\ X\,)\,;
             cuda subNum(abs mxy, Y);
         }else{
             cuda_copyNum(abs_mxy, Y);
             cuda subNum(abs mxy, X);
         cuda copyNum(G, abs mxy);
         cuda copyNum(N tmp, N);
         cuda gcd(G, N tmp);
         check ++;
         if ((check % (100) == 0) && !cuda zeroNum(result)){
             return:
    }
    cuda copyNum(result, G);
}
```

Figure 10: Kernel from second version

As you can see, after every 100 iteration thread check if some other thread hasn't solved the problem and stop.

3.4 Measurements

For measurement in this case was chossen small number:

```
N[0] = 0xb2c4b7b1;

N[1] = 0x0fd42d4e;
```

Again, the policy of preferring L1 was used. There was much less memory access since there is none global memory used. The only thing regarding bank conflict could be in accessing result which is actually provided to other threads as broadcast.

The average time of execution was 488627 [ms] which is much higher than in the last case.

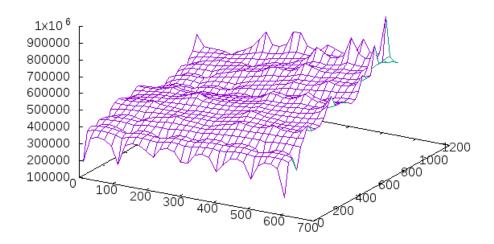


Figure 11: Graph of execution time

From this graph, we can observe that execution time is much higher.

3.4.1 Profiler output

Profiling was done on NVidia GeForce GTX 780 Ti at gpu-02.star.fit.cvut.cz.

In this case, the profiler didn't provide that much date with an explanation that it couldn't obtain them. Even after changing initial input value of N the profiler wasn't able to generate some result. But at least the ratio of which instructions were used was given. That actually showed how much diverged the algorithm was during the run.

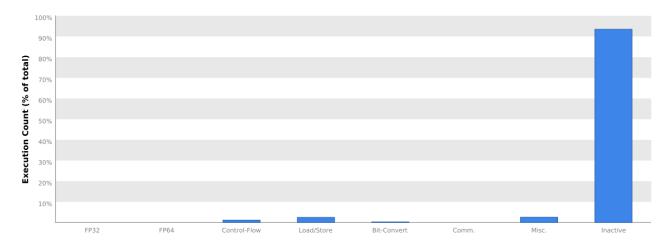


Figure 12: Inactivity of threads in warp

As the chart shows, divergence in this solution is huge, making it unusable for any usable results. A much worst case could appear with using bigger numbers.

4 Conclusion

The whole CUDA solution is based on pure luck finding some input values which actually converge faster to find some number dividing input N. To make the algorithm more deterministic, there is no random function used to fill input values. If random would be used, some threads could actually by some small probability count the same sequence. But on the other hand, some threads could converge much faster. CUDA contain library which can provide random numbers $curate{curate}$ which have big throughput, therefore there is no need to generate random numbers on host CPU.

In this solutions first CUDA implementation won. To compare for the same input N the speed up was 5.52 times. But! We have to notice that we picked block size 810 with a thread count 110. This solution was found with measuring going from 1 to 1024 for both block and thread count. Therefore, for new generic input number N there would be most strategic to get this numbers as big as local SM can serve.

Another much-needed improvement would be to work with some kind of GMP library which would contain a much better implementation of basic mathematical operations. Since this article is not about creating the fastest way to compute these operations, there is much to be done to improve them. With that, there should be some kind of relation with CUDA devices. If the number is big enough to rotate in memory banks, conflicts can occur making the whole implementation much slower. Same thing applies to the divergence of threads which was so much obvious in this implementation. If the mathematical operations would just work without any ifs and whiles, there should be much-needed speed up.

References

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