Dawn of GPU Era-Potentials of Chaos Theory

Saurabh Maniktala, Anisha Goel, A. B. Patki, and R. C. Meharde

Abstract—In the present era Chaos theory has tremendous potential in Computer Science Domain. The true potential of Chaos theory can be realized with the assistance of high performance computing aids such as GPU that have become available in present times. The main purpose is to develop a high performance experimental laboratory in academic institutions, for analysis of chaotic systems. In this paper we present the parallel implementation of One Dimensional Logistic Map on NVIDIA GeForce 9400 GT Graphical Processing Unit (GPU) which is a CUDA enabled GPU as a first step towards this direction. We report that the GPU version of Logistic Map executes 495.08 times faster than the CPU only version.

Index Terms-Chaos theory, CUDA, GPU, logistic map.

I. INTRODUCTION

The field of chaos has fascinated researchers from all over the world since its inception in the 1900s. Traditionally, in any English dictionary the word chaos is defined as complete disorder or anarchy. On the contrary, from a research perspective Chaos has been explored as a discipline which enables us to study the apparently random behavior of nonlinear dynamic systems. By the word apparently we stress the fact that datasets appearing random or chaotic can be generated from a deterministic mathematical equation. The key features of a chaotic system include:

- bounded
- nonlinearity
- mixing
- sensitive dependence on initial conditions (SDIC)

The milestones achieved in this field include electronic circuit by Leon O. Chua, the study of atmospheric convection by Edward N. Lorenz and chemical kinetics by Otto Rössler to name a few. Chaos theory finds applications in various scientific disciplines including computer science, population dynamics, aviation, politics, philosophy, robotics, finance, biology, and psychology.

In the present scenario Chaos theory is playing an even more prominent role for developing computer science based applications. Some of the works in this domain include preemptive queuing models with delay [1] and encryption techniques using Baptista Methodology [2].

In this paper, we have taken the first step towards setting up of an experimental laboratory in academic institutions,

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including AICTE approved engineering colleges, for analysis of chaotic systems. The systems can be divided into two broad categories of continuous and discrete time. We have presented a parallel implementation of one dimensional logistic map of discrete time system on the Graphical Processing Unit (GPU). The GPU was chosen for the implementation as it has an enormous parallel computational capability that makes analysis faster and efficient. We have achieved a speedup of 495.08 times over the sequential implementation—which we have also presented in this paper for comparison purposes.

We have carried out the experiment using the CUDA enabled NVIDIA GeForce 9400 GT and CUDA Toolkit version 3.0. Other details are mentioned in Table I.

Section II presents a brief introduction to one dimensional logistic map with its sequential C based implementation, section III elaborates the concepts of GPU and CUDA, section IV presents the pseudo-parallel implementation of one dimensional logistic map, section V includes the performance measurement of the parallel implementation and section VI presents conclusion and future work.

II. ONE DIMENSIONAL LOGISTIC MAP AND ITS SEQUENTIAL IMPLEMENTATION

The logistic equation as applied to Chaos models the growth curve of some natural phenomena. As an example, the curve could depict the variation in size of a population of a particular area over time. The size of the population is measured in discrete steps by using (1).

$$x_{t+1} = k x_t (1 - x_t) \tag{1}$$

TABLE I: DETAILS OF PLATFORM USED TO PERFORM EXPERIMENT

PLATFORM DETAILS		
CPU	Intel Pentium Dual Core	
OS	Windows XP	
GPU	NVIDIA GeForce 9400 GT	
Compute Capability	1.1	
Graphics Bus	PCI-Express	
CUDA Driver and Runtime	3.0	
Version		
Clock rate of GPU	1.4GHz	
Global Memory	1 GB	

The variable x_i denotes the current size of population, x_{i+1} represents future size of population and k is a constant known as the control parameter. The equation (1) is of a feedback nature as the size of population at some time interval is used to generate the size of population at the next time interval. The control parameter scales the size of population. In order to limit the size within realistic bounds, a factor of $(1-x_i)$ is multiplied. Since there is only one variable, (1) is more aptly

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referred to as one dimensional logistic equation. Equation (1), for simplicity, is normalized such that the size of population can only range from 0 to 1.

The one dimensional logistic map is obtained from the one dimensional logistic equation by plotting the values of x_{t+1} as y coordinates against values of x_t as x coordinates on a graph. Since the map is using two dimensions, namely x and y, to plot the points representing the same variable, the size of population, it is also called pseudo two dimensional.

The initial size of the population is denoted by x_0 . As chaotic systems are sensitive to initial conditions x_0 plays an important role to enter into Chaos. Moreover, the control parameter *k* has to be tuned to the appropriate value for the onset of Chaos. The value of *k* governs the steepness and the height of the curve which is a parabola. As the values generated from the equation are normalized, the allowable range of k is (0,4).

The system shows varied behavior for different ranges of *k* as:

- For k<1, the system eventually converges to a zero size population, so in this case the attractor is zero.
- For 1<k<3, the attractor increases from point zero to 0.667. For 0<k<3 the population value sooner or later saturates irrespective of initial value.
- For $3 \le k < 3.57$, period doubling starts. At k=3.4, the attractor splits into two point attractor. For greater values of k in this range the attractor keeps on doubling to produce 4, 8, 16... point attractors. In this range the final values keep on oscillating.
- For $3.57 \le k < 4$, the system enters in chaotic region. In this case the attractor can be erratic with infinite number of points or stable.

We now present the sequential C based implementation of logistic map. We have developed a C function named *logisitcmap()* having the arguments x_0 as the initial value of population, *k* as the control parameter, *y* as the output array containing population sizes for consecutive iterations and *N* as the number of iterations. The function returns the output array *y*. Below we present the source code of *logisticmap()*.

```
float* logisticmap(float x0, float k, float
y[], int N)
{
    float temp=x0;
    for(int i=0;i<N;i++)
    {
       temp = k * temp * (1-temp);
       y[i] = temp;
    }
    return(y);}</pre>
```

Table II shows the output array for N=10 iterations, k=3.7and two initial values of x_0 as 0.01 and 0.011. The value of k=3.7 drives the system into chaos and apparently random successive output values are generated. Notice that the output values arising out of the two initial values of x_0 differing by decimal in the third place diverge as the index of iteration increases. This behavior of chaotic systems as referred to as Sensitive Dependence on Initial Conditions.

TABLE II: OUTPUT VALUES FOR TWO INITIAL VALUES IN CHAOS STATE

Index	<i>x</i> ₀ =0.01	<i>x</i> ₀ =0.011
1	0.036630	0.040252
2	0.130567	0.142939
3	0.420020	0.453276
4	0.901332	0.916923
5	0.329051	0.281850
6	0.816873	0.748919
7	0.553488	0.695746
8	0.914415	0.783228
9	0.289564	0.628192
10	0.761152	0.864197

III. GPU AND CUDA

The Graphical Processing Unit (GPU) was traditionally designed to offload intensive graphical computations, including scaling, rotating, and rendering, from the Central Processing Unit (CPU). It essentially functions as a coprocessor to the CPU which directs work to it. Hence, the GPU is also known as the Device whereas the CPU is known as the Host.

The traditional GPU architecture consisted of pipelined stages of combination of programmable vertex shader and pixel fragment processors which are used for generating graphics on the screen. Subsequently, the vertex shader and pixel fragment processors were combined into a unified GPU architecture for efficient working.

The advances in the technology of GPU encouraged researchers to step into the direction of General Purpose GPU (GPGPU) programming. The GPGPU concept focuses on using the computational capability of GPU for non-graphical applications. The current GPUs have developed into a programmable parallel processor offering high performance parallel computing capabilities far exceeding that of current multi-core CPUs. The design goal of GPU is to optimize execution of massive number of threads in parallel and hence a larger chip area is dedicated for floating point operations. However, as GPU works as a coprocessor for host CPU, it will not outperform on other jobs that CPU is designed.

The greatest challenge that was faced was to use these numerous cores in an optimized fashion. In order to overcome such difficulties, NVIDIA Corporation developed a software environment named CUDA (Compute Unified Device Architecture) that extends the C language syntax [3]. [4].CUDA enables programmers to develop programs that can scale across the cores of a GPU while using the familiar syntax of C language. CUDA package includes a Runtime API and a Driver API. The CUDA runtime API extends C language syntax with function commands dedicated for GPU and is referred to as CUDA C. The Driver API, on the other hand, is a low level procedural API but is totally different from the graphical APIs such as Open GL and DirectX. The OpenGL and DirectX APIs are traditionally used for GPU programming for graphical applications. But the situation gets complicated when a programmer wants to deploy GPU to develop a parallel program for non graphical applications and had to put unnecessary effort of studying the graphic libraries is required. Thus CUDA C fulfills the programmer's needs. We explain the concepts of GPU in the context of CUDA programming.

Fig. 1. shows the architecture of an NVIDIA GeForce 8

series GPU. The basic unit shown is a Texture/Processor (TPC) which Cluster incorporates а Streaming Multiprocessor Controller (SMC), a texture memory unit, and 2 Streaming Multiprocessor (SM) blocks [5]. The Multiprocessor handles continuous flow of data streams and is thus named as a Streaming Multiprocessor. The SMC is responsible for work distribution amongst the two SMs. Each SM is responsible for executing hardware threads that perform computations. The texture memory unit inside the TPC is generally used for storing pixel values. Each SM houses 8 Streaming Processors (SP), 2 Special Function Units (SFUs), a Cache Unit and a 16 KB shared memory unit. The actual floating point computations take place in the SP which is essentially a single core handling a single hardware thread. A hardware thread is the smallest unit of computation. The SP utilizes cache memory for faster lookup of data, while the inter-SP communication in a SM takes place via the shared memory. The GPU actually consists of an array of TPCs. In addition to the above, the GPU also includes global memory unit and constant cache unit on the same chip. The number of TPCs and the size of global and constant cache memory depend on the GPU series.

The concurrent execution of the threads on the SPs of a SM forms the basis of parallel computation. Since a fixed number of threads can run concurrently on a SM, groups of 32 threads are made from the bank of threads and are scheduled to run in an ordered manner. From a hardware view point, these groups of 32 threads are known as Warps. At a particular instance of time only one warp can be active.

From the CUDA programming view point, the program is broken into sequential and parallel modules. For parallel modules, threads are organized into two level of hierarchies namely blocks and grids [6], [7]. The first hierarchy, block is a software programming concept wherein threads are organized in a 3 dimensional layout (as x, y, and zdimensional representation).



Fig. 1. Basic GPU architecture

The second hierarchy grid is also a software concept wherein the first hierarchy blocks are further organized in a two dimensional layout. The threads in a block execute cooperatively in parallel on the SM, whereas the blocks of given grid execute independent of each other across the GPU. On the contrary, the grids execute sequentially on a GPU. The independent execution of blocks forms the basis of coarse-grained parallelism and cooperative parallel execution of threads in block implements fine-grained parallelism. Each grid is formed while executing a kernel call. The CUDA kernel is the program module to be executed on the GPU and is typically invoked from host CPU. Multiple copies of the same program module are created and are assigned to each thread at the time of kernel call. We have to specify the dimensions of the grid and the block in the execution configuration at the time of the kernel call. The programmer is free to experiment with the dimension sizes of blocks and grids. The total number of threads in a grid and block is calculated by using (2).

$$T = grid_dimension \ x \ block_dimension$$
 (2)

where T is the total number of threads, grid_dimension denotes the total number of blocks in a grid and block_dimension denotes the total number of threads in a block. For example if grid is to have 6 blocks (grid_dimension=(3,2) organized as 3 blocks in x direction and 2 blocks in y direction) and further each block has to have 48 threads (block_dimension=(6,4,2) organized as 6 threads in x direction, 4 in y direction and 2 in z direction) totaling 48*6= 288 threads in all, then the kernel call will be given as:

dim3 grid_dimension (3,2) dim3 block_dimension(6,4,2) Another alternative allotment for same problem could be:



TABLE III: OUTPUT VALUES FOR TWO INITIAL VALUES IN CHAOS STATE IN PARALLEL IMPLEMENTATION

Index	<i>x</i> ₀ =0.01	<i>x</i> ₀ =0.011
1	0.036630	0.040252
2	0.130567	0.142939
3	0.420020	0.453276
4	0.901332	0.916923
5	0.329051	0.281850
6	0.816873	0.748919
7	0.553488	0.695746
8	0.914414	0.783228
9	0.289565	0.628192
10	0.761153	0.864197

TABLE IV: GPU AND CPU EXECUTION TIME		
No. of iterations	CPU(MSEC)	GPU(MSEC)
128	50.000	0.115
256	60.000	0.126
384	70.000	0.148
512	90.000	0.168
640	110.000	0.199

The performance issues resulting out of the above two alternative representations for 288 threads are beyond the scope of this paper. Fig. 2. illustrates the first alternative showing the arrangement of threads in blocks and blocks in grid. It also depicts the memory visible to each level of hierarchy. The local memory is only visible by its corresponding thread. Also, the threads in a block can communicate using shared memory. Lastly all the blocks in the grid can see the global memory.

IV. PARALLEL IMPLEMENTATION OF LOGISTIC MAP

In this section we present the parallel GPU based implementation of the one dimensional logistic map.

We have developed a kernel logisticmap() having the following arguments x_0 as the initial value of population, k as the control parameter, and y as the output array containing population sizes for consecutive iterations. Below we present the code snippet of logisticmap():

```
void logisticmap(float *y,
  global
\overline{float} \times 0, float k
  {
      shared
                 float temp;
    temp = x0;
    int i;
       int idx = blockIdx.x*blockDim.x +
       threadIdx.x;
    for(i=0;i<=idx;i++)</pre>
    {
      temp = k*temp*(1-temp);
    }
    y[idx] = temp;
  }
```

The execution configuration of the kernel call for a total of N iterations is given below:

dim3 block dimension(2);



Fig. 3. GPU versus CPU Execution Time

Table III shows the output array generated by the kernel *logisticmap()* for N=10, k=3.7 and initial values of x_0 as 0.01 and 0.011 for comparison with Table II. There is a difference in the sixth decimal place in last the 3 values for $x_0=0.01$ from the corresponding values in the sequential implementation due to round off scheme employed by the Device instead of truncation employed by Host. This difference can be overcome using a Device of compute capability 2.0.

V. PERFORMANCE MEASUREMENT

Now we present the comparison between the execution times of *logisticmap()* on GPU and CPU on the basis of the number of iterations performed. The execution times on GPU (parallel implementation) and CPU (sequential implementation) have been measured in milliseconds with precision of three decimal places and are enlisted in Table IV.

Fig. 3. depicts a graph generated using Table IV. The graph has been interpolated for the ease of understanding. As we observe from the figure the GPU execution time remains nearly constant over the entire range of iterations performed. On the contrary, the CPU execution time increases proportionally with the number of iterations performed. The most remarkable result that we have achieved is that on an average GPU computed the result 495.08 times faster than the CPU.

VI. CONCLUSION AND FUTURE WORK

The objective of this paper is to present a high performance parallel GPU based implementation of One Dimensional Logistic Map using the NVIDIA CUDA environment first step towards setting up of an experimental laboratory in academic institutions, including AICTE approved engineering colleges, for analysis of chaotic systems by developing one of the most fundamental tools for study. We have reported an enormous speedup of 495.08 times than the CPU only version. Towards this direction we have already explored potentials of Chaotic cryptography (Baptista's Methodology). Authors feel that this will be useful in offering an academic course with laboratory hands-on in high performance applications of Chaos.

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