

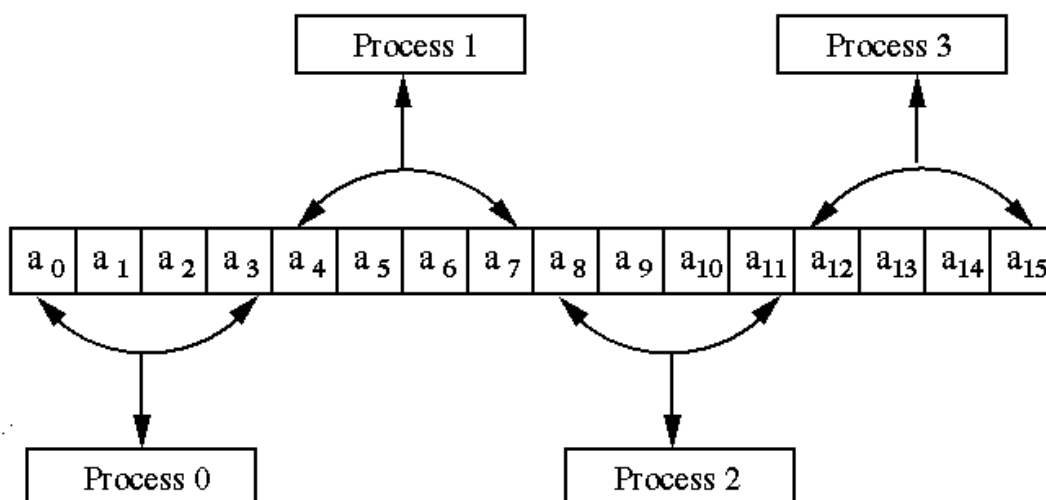
## VECTOR ADDITION USING MULTIPLE GPUS

### **Objective of the Lab:**

- To write a MPI - CUDA program, to compute the vector-vector addition on Multi-Core processor system with multi-GPU. Assume that each process computes the partial Vector Vector Addition using CUDA kernel.

### **Description:**

- The partitioning is called block-striped if each process is assigned contiguous elements. The process P0 gets the first  $n/p$  elements, P1 gets the next  $n/p$  elements and so on.
- The distribution of 16 elements of vector A on 4 processes is shown in the following below.



- Initially process with rank 0 distributes the input vectors using *MPI\_Scatter* on  $p$  processes. Each process will call a CUDA kernel which performs local

addition of the vectors and stores the partial addition. Now the process with rank 0 performs global reduction using *MPI\_Reduce* to get the final addition product of two vectors.

## **Implementation of Vector Vector Addition :**

### **Step 1 :**

Four vectors are required for computation. Two arrays for vector A and Vector B and the other two arrays for temporary storage data for two vector at the each node.

### **Step2 :**

Root process initializes the two vectors i.e Vector A and Vector B. The two vectors are constructed by assigning to each element one more than its index value.

### **Step 3 :**

Vector Size is broadcasted to all processes from the root process.

### **Step 4 :**

Memory is assigned for MyVectorA and MyVectorB on all nodes.

### **Step 5 :**

Process with rank 0 distributes the input vectors using *MPI\_Scatter* on to p processes.

### **Step 6 :**

Similar arrays are allocated on the device. The values of the arrays in the host machine are copied on to the arrays allocated on the device.

### **Step 7:**

Each node computes the partial addition value by calling VectorVectorAddition Cuda kernel.

### **Step 8 :**

Process with rank 0 performs global reduction using MPI\_Reduce to get the final addition of two vectors.

**Step 9 :**Process with rank 0 prints the addition value.

### **CUDA API used :**

To Allocate memory on device-GPU :

***cudaMalloc(void\*\* array, int size)***

To Free memory allocated on device-GPU:

***cudaFree(void\* array)***

To transfer from host-CPU to device-GPU:

***cudaMemcpy((void\*)device\_array, (void\*)host\_array, size, cudaMemcpyHostToDevice)***

To transfer from device-GPU to host-GPU:

***cudaMemcpy((void\*)host\_array, (void\*)device\_array, size, cudaMemcpyDeviceToHost)***

### **INPUT:**

The input to the problem is given as arguments in the command line. It should be given in the following format ; Suppose that the size of the vector is n and the number of nodes is m, then the program must be run as,

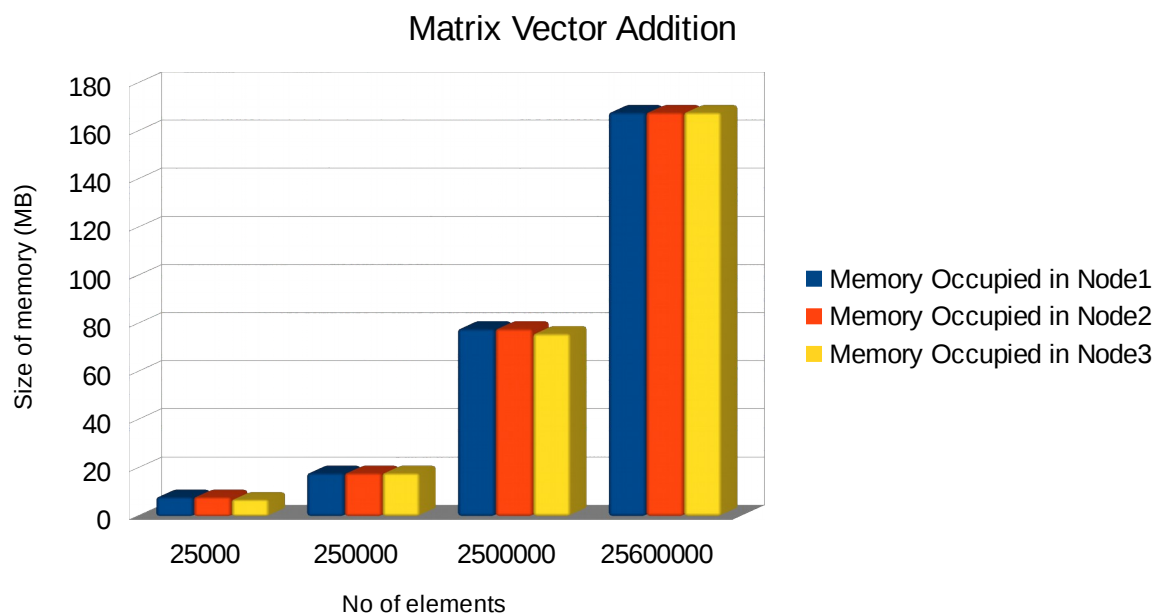
**mpirun -n m ./program\_name n**

Process 0 generates the two vectors i.e Vector A and Vector B

### **OUTPUT:**

Process 0 prints the final addition value of two vectors. The correctness of the output can be verified using the below formula The sum of the squares of the first n numbers is  $(n(n+1)(2n+1))/6$ .

Input Size	Memory Occupied ( Node 1 )	Memory Occupied ( Node 2 )	Memory Occupied ( Node 3 )
25000	8	8	8
250000	20	20	18
2500000	78	78	76
25600000	168	168	168



**POINT TO NOTE:**

Each node got an equal no of the elements and hence does an equal number of computation.