

Trapezoid Rule Integration using MPI

MPI is implemented to integrate the function outlined by the question using a cluster. The subintervals of the integration are equal to the number of processors one designates for the computation. When the trapezoid program is used to integrate the function over 10 processors it computes the following results:

```
-bash-4.2$ mpirun -np 10 Trapezoid_Rule_MPI
```

```
Area found in processor 7 = 77.2812
```

```
Area found in processor 3 = 43.2913
```

```
Area found in processor 9 = 31.2763
```

```
Area found in processor 2 = 26.0438
```

```
Area found in processor 5 = 73.2862
```

```
Area found in processor 1 = 11.2963
```

```
Area found in processor 8 = 62.0288
```

```
Area found in processor 6 = 80.0338
```

```
Area found in processor 4 = 60.0387
```

```
Area found in processor 0 = 2.04875
```

```
Trapezoidal Rule Total Area = 466.625
```

The function is split into segments and then sent numerically to each processor. The cluster does not necessarily finish the computations in numerical order.