

COMP/CS 605: Introduction to Parallel Computing

Topic : MPI: Trapezoid Function

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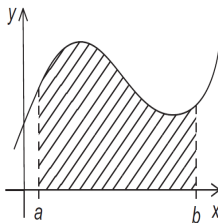
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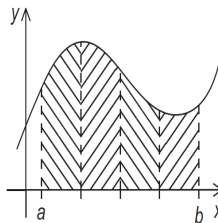
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The Trapezoid Rule for Numerical Integration

Solve the Integral: $\int_a^b F(x)dx$



(a)



(b)

Where $F(x)$ can be any function of x : $f(x^2)$, $f(x^3)$

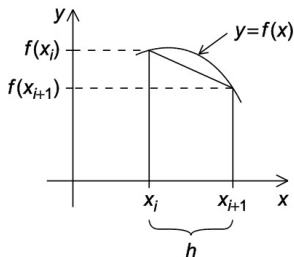
See Pacheco IPP (2011), Ch3.

Trapezoid Equations

Integral: $\int_a^b f(x) dx$

Area of 1 trapezoid: $= \frac{h}{2} [f(x_i) + f(x_{i+1})]$

Base: $h = \frac{b-a}{n}$



Endpoints: $x_0 = a, \quad x_1 = a + h, \quad x_2 = a + 2h, \dots, \quad x_{n-1} = a + (n-1)h, \quad x_n = b$

Sum of Areas: $Area = h \left[\frac{f(x_0)}{2} + f(x_1) + f(x_2) + \dots + f(x_{n-1}) + \frac{f(x_n)}{2} \right]$

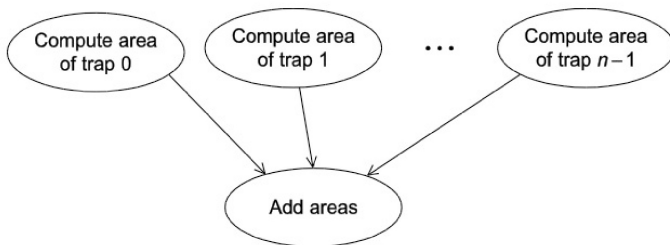
Pseudocode for a Serial Algorithm

```
/*  Input: a ,b, n  */  
h = (b-a)/n ;  
approx = (F(a) + F(b))/2.0  
for (i=0; i<= n-1; i++) {  
    x_i = a + i*H;  
    approx += f(x_i);  
}  
approx = h* approx
```

Paralellizing the Trapezoidal Rule

PCAM Approach

- Partition problem solution into tasks.
- Identify communication channels between tasks.
- Aggregate tasks into composite tasks.
- Map composite tasks to cores.



Two types of tasks:

Compute area of 1 trapezoid

Compute area sums

Parallel Pseudocode

```
Get a,b,n;
h = b*a/n ;
local_n = n/comm_sz;
local_a = a + my_rank*local_n*h;
local_b = local_a + local_n*h;
local_integral = Trap(local_a , local_b , local_n , h);
if (my_rank != 0)
    Send local integral to process 0;
else /* my_rank==0 */
    total_integral = local_integral;
    for (proc = 1; proc < comm_sz; proc++)
    {
        Receive local integral from proc;
        total integral += local integral;
    }
}

if (my_rank == 0)
    print result;
```


Datatype Example: Pacheco code: mpi-trap1.c

```
/* File:      mpi_trap4.c
 * Purpose:   Use MPI to implement a parallel version of the trapezoidal rule.
 *           This version uses collective communications and MPI derived datatypes
 *           to distribute the input data and compute the global sum.
 *
 * Input:     The endpoints of the interval of integration and the number of trapezoids
 * Output:    Estimate of the integral from a to b of f(x) using the trapezoidal rule and n trapezoids.
 *
 * Compile:   mpicc -g -Wall -o mpi_trap4 mpi_trap.c
 * Run:      mpiexec -n <number of processes> ./mpi_trap
 *
 * Algorithm:
 *   1. Each process calculates "its" interval of integration.
 *   2. Each process estimates the integral of f(x) over its interval using the trapezoidal rule.
 *   3a. Each process != 0 sends its integral to 0.
 *   3b. Process 0 sums the calculations received from the individual processes and prints the result.
 *
 * Note: f(x) is all hardwired.
 * IPP:  Section 3.2.2 (pp. 96 and ff.)
 */
#include <stdio.h>
#include <mpi.h>      /* MPI routines, definitions, etc. */

/* Calculate local integral */
double Trap(double left_endpt, double right_endpt, int trap_count, double base_len);

/* Function we're integrating */
double f(double x);
```

Trapezoid Rule for Numerical Integration

First MPI Parallel Trapezoid Rule

Datatype Example: Pacheco code: mpi-trap1.c

```
int main(void) {
    int my_rank, comm_sz, n = 1024, local_n;
    double a = 0.0, b = 3.0, h, local_a, local_b;
    double local_int, total_int;
    int source;

    /* Let the system do what it needs to start up MPI */
    MPI_Init(NULL, NULL);

    /* Get my process rank */
    MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);

    /* Find out how many processes are being used */
    MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);

    h = (b-a)/n;          /* h is the same for all processes */
    local_n = n/comm_sz;  /* So is the number of trapezoids */

    /* Length of each process' interval of integration = local_n*h.  So my interval starts at: */
    local_a = a + my_rank*local_n*h;
    local_b = local_a + local_n*h;
    local_int = Trap(local_a, local_b, local_n, h);

    /* Add up the integrals calculated by each process */
    if (my_rank != 0) {
        MPI_Send(&local_int, 1, MPI_DOUBLE, 0, 0, MPI_COMM_WORLD);
    } else {
        total_int = local_int;
        for (source = 1; source < comm_sz; source++) {
            MPI_Recv(&local_int, 1, MPI_DOUBLE, source, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
            total_int += local_int;
        }
    }

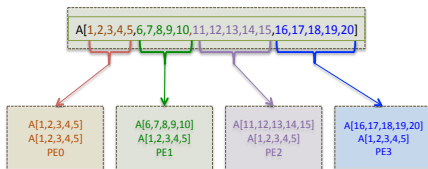
    /* Print the result */
    if (my_rank == 0) {
        printf("With n = %d trapezoids, our estimate\n", n);
        printf("of the integral from %f to %f = %.15e\n", a, b, total_int);
    }

    MPI_Finalize();          /* Shut down MPI */
}
```

Datatype Example: Pacheco code: mpi-trap1.c

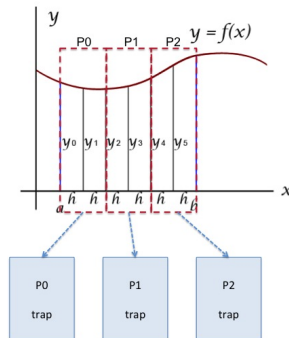
```
/*-----  
 * Function:    Trap  
 * Purpose:    Serial function for estimating a definite integral  
 *             using the trapezoidal rule  
 * Input args:  left_endpt  
 *             right_endpt  
 *             trap_count  
 *             base_len  
 * Return val:  Trapezoidal rule estimate of integral from  
 *             left_endpt to right_endpt using trap_count  
 *             trapezoids  
 */  
double Trap(  
    double left_endpt /* in */,  
    double right_endpt /* in */,  
    int trap_count /* in */,  
    double base_len /* in */) {  
    double estimate, x;  
    int i;  
  
    estimate = (f(left_endpt) + f(right_endpt))/2.0;  
    for (i = 1; i <= trap_count-1; i++) {  
        x = left_endpt + i*base_len;  
        estimate += f(x);  
    }  
    estimate = estimate*base_len;  
  
    return estimate;  
} /* Trap */  
  
/*-----  
 * Function:    f  
 * Purpose:    Compute value of function to be integrated  
 * Input args:  x  
 */  
double f(double x) {  
    return x*x;  
} /* f */
```

Data Distribution for Trap Function



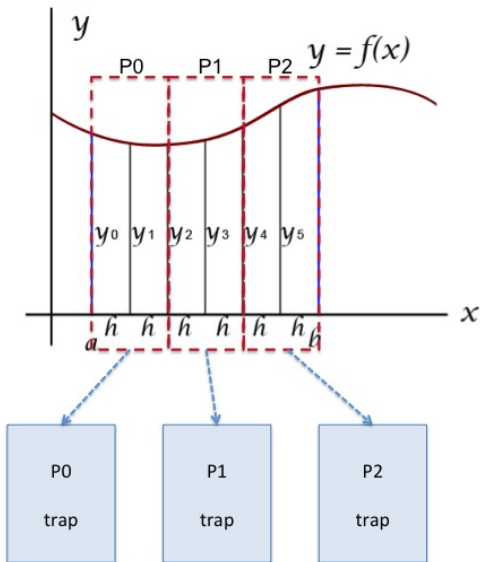
1D Vector Data Distribution

- Parallel *trap* distributes $[x_i]$ data points to different processors.
- Each processor runs the same code but solves for different $[x_i]$ data
- This is a *1-D* data decomposition
- This is a *1-D* (virtual) processor arrangement



Trapezoid Rule for Numerical Integration

First MPI Parallel Trapezoid Rule



Job Example: Serial Hello World

```
#include <stdio.h>
#include <string.h> /* For strlen */
#include <mpi.h> /* For MPI functions, etc */

const int MAX_STRING = 100;

int main(void) {
    char    greeting[MAX_STRING]; /* String storing message */
    int     comm_sz;              /* Number of processes */
    int     my_rank;              /* My process rank */

    /* Start up MPI */
    MPI_Init(NULL, NULL);

    /* Get the number of processes */
    MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);

    /* Get my rank among all the processes */
    MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);

    /* Print my message */
    printf("Greetings from process %d of %d!\n", my_rank, comm_sz);

    /* Shut down MPI */
    MPI_Finalize();

    return 0;
} /* main */
```

Job Example: Serial Hello World

```
[tuckoo] mthomas% mpirun -np 8 ./simp_mpi_hi
Greetings from process 2 of 8!
Greetings from process 0 of 8!
Greetings from process 1 of 8!
Greetings from process 3 of 8!
Greetings from process 6 of 8!
Greetings from process 5 of 8!
Greetings from process 7 of 8!
Greetings from process 4 of 8!
```

Output is random, or non-deterministic

To build more flexible code, you need to read in variable data. One approach is to use the STDIN.

- Most MPI implementations only allow process 0 in MPI_COMM_WORLD access to stdin.
- Process 0 must read the data (scanf) and send to the other processes.

Datatype Example: Pacheco code: mpi-trap2.c

```
/* File:      mpi_trap4.c
* Purpose:   Use MPI to implement a parallel version of the trapezoidal rule.
*           This version uses collective communications and MPI derived datatypes
*           to distribute the input data and compute the global sum.
*
* Input:     The endpoints of the interval of integration and the number of trapezoids
* Output:    Estimate of the integral from a to b of f(x) using the trapezoidal rule and n trapezoids.
*
* Compile:   mpicc -g -Wall -o mpi_trap4 mpi_trap.c
* Run:      mpiexec -n <number of processes> ./mpi_trap
*
* Algorithm:
*   1. Each process calculates "its" interval of integration.
*   2. Each process estimates the integral of f(x) over its interval using the trapezoidal rule.
*   3a. Each process != 0 sends its integral to 0.
*   3b. Process 0 sums the calculations received from the individual processes and prints the result.
*
* Note: f(x) is all hardwired.
* IPP:  Section 3.2.2 (pp. 96 and ff.)
*/
#include <stdio.h>
#include <mpi.h>      /* MPI routines, definitions, etc. */

/* Get the input values */
void Get_input(int my_rank, int comm_sz, double* a_p, double* b_p, int* n_p);

/* Calculate local integral */
double Trap(double left_endpt, double right_endpt, int trap_count, double base_len);

/* Function we're integrating */
double f(double x);
```

Datatype Example: Pacheco code: mpi-trap2.c

```
int main(void) {
    int my_rank, comm_sz, n, local_n, source;
    double a, b, h, local_a, local_b, local_int, total_int;

    /* Let the system do what it needs to start up MPI */
    MPI_Init(NULL, NULL);
    MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);      /* Get my process rank */
    MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);      /* Find out how many processes are being used */

    Get_input(my_rank, comm_sz, &a, &b, &n);

    h = (b-a)/n;          /* h is the same for all processes */
    local_n = n/comm_sz; /* So is the number of trapezoids */

    /* Length of each process' interval of integration = local_n*h. So my interval starts at: */
    local_a = a + my_rank*local_n*h;
    local_b = local_a + local_n*h;
    local_int = Trap(local_a, local_b, local_n, h);

    /* Add up the integrals calculated by each process */
    if (my_rank != 0) {
        printf("PE[%d] SEND: For %d trapezoids, local estimate=%f\n", my_rank, local_n, local_int);
        MPI_Send(&local_int, 1, MPI_DOUBLE, 0, 0, MPI_COMM_WORLD);
    }
    else {
        total_int = local_int;
        for (source = 1; source < comm_sz; source++) {
            MPI_Recv(&local_int, 1, MPI_DOUBLE, source, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
            printf("PE[%d] RECV from PE[%d]: local estimate=%f\n", my_rank, source, local_int);
            total_int += local_int;
        }
    }

    /* Print the result */
    if (my_rank == 0)
        printf("With n = %d trapezoids, our estimate of the integral from %f to %f = %.15e\n", n, a, b, total_int);

    MPI_Finalize();      /* Shut down MPI */
    return 0;
} /* main */
```

MPI_Send: Performs a blocking send

```
int MPI_Send(  
    const void *sendbuf,  
    int count,  
    MPI_Datatype datatype,  
    int dest,  
    int tag,  
    MPI_Comm comm )
```

Input Parameters:

- sendbuf: address of send buffer (choice)
- count: number of entries in buffer (integer)
- datatype: data type of buffer (handle)
- dest: rank of destination (integer)
- tag: message tag (integer)

MPI_Recv: Blocking receive for a message

```
int MPI_Recv(  
    const void *recvbuf,  
    int count,  
    MPI_Datatype datatype,  
    int source,  
    int tag,  
    MPI_Comm comm,  
    MPI_Status *status )
```

Input Parameters:

- count: number of entries in buffer (integer)
- datatype: data type of buffer (handle)
- dest: rank of destination (integer)
- tag: message tag (integer)

Output Parameters:

- recvbuf: address of receive buffer (choice)
- status: status object (Status)

Datatype Example: Pacheco code: mpi-trap2.c

```
/*-----  
* Function:      Get_input  
* Purpose:      Get the user input:  the left and right endpoints  
*              and the number of trapezoids  
* Input args:   my_rank:  process rank in MPI_COMM_WORLD  
*              comm_sz:  number of processes in MPI_COMM_WORLD  
* Output args:  a_p:  pointer to left endpoint  
*              b_p:  pointer to right endpoint  
*              n_p:  pointer to number of trapezoids  
*/  
void Get_input(int my_rank, int comm_sz, double* a_p, double* b_p, int* n_p)  
{  
    int dest;  
  
    if (my_rank == 0) {  
        printf("Enter a, b, and n\n");  
        scanf("%lf %lf %d", a_p, b_p, n_p);  
        for (dest = 1; dest < comm_sz; dest++) {  
            MPI_Send(a_p, 1, MPI_DOUBLE, dest, 0, MPI_COMM_WORLD);  
            MPI_Send(b_p, 1, MPI_DOUBLE, dest, 0, MPI_COMM_WORLD);  
            MPI_Send(n_p, 1, MPI_INT, dest, 0, MPI_COMM_WORLD);  
        }  
    } else { /* my_rank != 0 */  
        MPI_Recv(a_p, 1, MPI_DOUBLE, 0, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);  
        MPI_Recv(b_p, 1, MPI_DOUBLE, 0, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);  
        MPI_Recv(n_p, 1, MPI_INT, 0, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);  
    }  
} /* Get_input */
```

What happens if you are using the batch queuing system?
There is no way to capture *stdin*, so you have to read in data as
command line args or from a file.

Datatype Example: Pacheco code: mpi-trap2.c

```
main(int argc, char *argv[]) {
    long cmdvar;
    char cptr[100];
    int  comm_sz;           /* Number of processes */
    int  my_rank;          /* My process rank */

    /* Start up MPI */
    MPI_Init(NULL, NULL);

    /* Get the number of processes */
    MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);

    /* Get my rank among all the processes */
    MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);

    /* get a number from the command line */
    if (argc != 2) {
        printf("usage: hello-arg <integer number>\n");
        return 0;
    }
    cmdvar = strtol(argv[1], NULL, 10);

    /* Print my message */
    printf("Greetings from process %d of %d, cmd arg val=%ld\n", my_rank, comm_sz, cmdvar);

    /* Shut down MPI */
    MPI_Finalize();

    return 0;
} /* main */

"hello-arg.c" 41L, 980C written
```

Datatype Example: Pacheco code: mpi-trap2.c

```
[mthomas@tuckoo:~/pardev/hello] mpirun -np 5 ./hello-arg 837298
```

```
Greetings from process 4 of 5, cmd arg val=837298
```

```
Greetings from process 0 of 5, cmd arg val=837298
```

```
Greetings from process 1 of 5, cmd arg val=837298
```

```
Greetings from process 2 of 5, cmd arg val=837298
```

```
Greetings from process 3 of 5, cmd arg val=837298
```

```
[mthomas@tuckoo:~/pardev/hello] mpirun -np 8 ./hello-arg 10000000000000000000
```

```
Greetings from process 6 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 7 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 0 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 3 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 4 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 5 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 1 of 8, cmd arg val=9223372036854775807
```

```
Greetings from process 2 of 8, cmd arg val=9223372036854775807
```