

CSC4005 – Distributed and Parallel Computing

Prof. Yeh-Ching Chung

School of Data Science
Chinese University of Hong Kong, Shenzhen





Outline

- Introduction to Parallel Computers
- Message Passing Computing and Programming
- Multithreaded Programming
- CUDA Programming
- OpenMP Programming
- Embarrassingly Parallel Computations
- Partitioning and Divide-and-Conquer Strategies
- Pipelined Computations
- Synchronous Computations
- Load Balancing and Termination Detection
- Sorting Algorithms





Embarrassingly Parallel Computations (1)

A computation that can be divided into a number of completely independent parts, each of which can be executed by a separate processor.

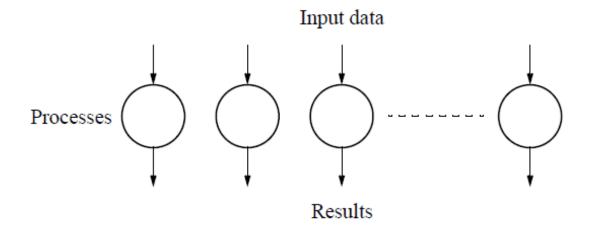


Figure 3.1 Disconnected computational graph (embarrassingly parallel problem).





Embarrassingly Parallel Computations (2)

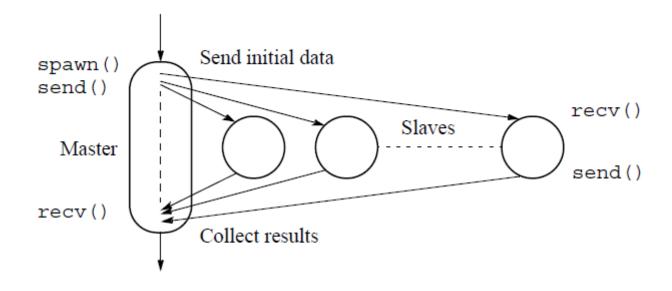


Figure 3.2 Practical embarrassingly parallel computational graph with dynamic process creation and the master-slave approach.



Embarrassingly Parallel Examples (1)

Low level image operations:

(a) Shifting

Object shifted by Δx in the x-dimension and Δy in the y-dimension:

$$x' = x + \Delta x$$

$$y' = y + \Delta y$$

where x and y are the original and x' and y' are the new coordinates.

(b) Scaling

Object scaled by a factor S_x in the x-direction and S_y in the y-direction:

$$\chi' = \chi S_{\gamma}$$

$$y' = yS_y$$

(c) Rotation

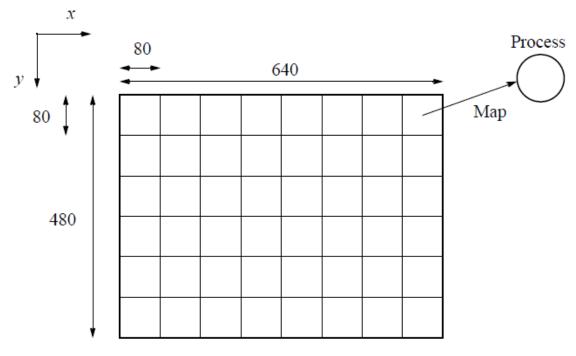
Object rotated through an angle θ about the origin of the coordinate system:

$$x' = x \cos \theta + y \sin \theta$$

$$y' = -x \sin\theta + y \cos\theta$$



Embarrassingly Parallel Examples (2)

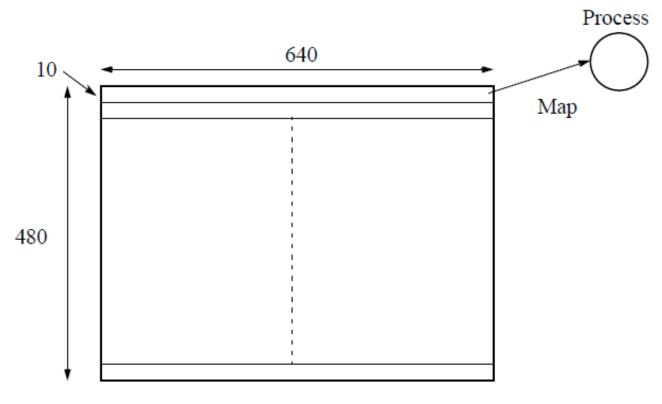


(a) Square region for each process

Figure 3.3 Partitioning into regions for individual processes.



Embarrassingly Parallel Examples (3)



(b) Row region for each process





Pseudocode to Perform Image Shift (1)

Master



Embarrassingly Parallel Examples (2)

Slave





Embarrassingly Parallel Examples (3)

Analysis Sequential

$$t_s = n^2 = \mathcal{O}(n^2)$$

Parallel

Communication

$$t_{\text{comm}} = t_{\text{startup}} + mt_{\text{data}}$$
$$t_{\text{comm}} = p(t_{\text{startup}} + 2t_{\text{data}}) + 4n^2(t_{\text{startup}} + t_{\text{data}}) = O(p + n^2)$$

Computation

$$t_{\text{comp}} = 2\left(\frac{n^2}{p}\right) = O(n^2/p)$$

Overall Execution Time

$$t_p = t_{\rm comp} + t_{\rm comm}$$

For constant p, this is $O(n^2)$. However, the constant hidden in the communication part far exceeds those constants in the computation in most practical situations.



Mandelbrot Set Computation (1)

Set of points in a complex plane that are quasi-stable (will increase and decrease, but not exceed some limit) when computed by iterating the function

$$z_{k+1} = z_k^2 + c$$

where z_{k+1} is the (k+1)th iteration of the complex number z = a + bi and c is a complex number giving the position of the point in the complex plane.

The initial value for z is zero.

Iterations continued until magnitude of z is greater than 2 or number of iterations reaches arbitrary limit. Magnitude of z is the length of the vector given by

$$z_{\text{length}} = \sqrt{a^2 + b^2}$$



Mandelbrot Set Computation (2)

Computing the complex function, $z_{k+1} = z_k^2 + c$, is simplified by recognizing that

$$z^2 = a^2 + 2abi + bi^2 = a^2 - b^2 + 2abi$$

or a real part that is $a^2 - b^2$ and an imaginary part that is 2ab.

The next iteration values can be produced by computing:

$$z_{\text{real}} = z_{\text{real}}^2 - z_{\text{imag}}^2 + c_{\text{real}}$$

$$z_{\rm imag} = 2z_{\rm real}z_{\rm imag} + c_{\rm imag}$$



Mandelbrot Set Computation (3)

Seq. Routine computing value of one pt, returning no of iterations

```
structure complex {
  float real;
  float imag;
};
int cal pixel(complex c)
int count, max;
complex z;
float temp, lengthsq;
\max = 256;
z.real = 0; z.imag = 0;
                                  /* number of iterations */
count = 0;
do {
  temp = z.real * z.real - z.imag * z.imag + c.real;
  z.imag = 2 * z.real * z.imag + c.imag;
  z.real = temp;
  lengthsq = z.real * z.real + z.imag * z.imag;
  count++;
} while ((lengthsq < 4.0) && (count < max));</pre>
return count;
```





Mandelbrot Set Computation (4)

Scaling Coordinate System

```
For computational efficiency, let

scale_real = (real_max - real_min)/disp_width;
scale_imag = (imag_max - imag_min)/disp_height;
```

Including scaling, the code could be of the form

```
for (x = 0; x < disp_width; x++) /* screen coordinates x and y */
  for (y = 0; y < disp_height; y++) {
    c.real = real_min + ((float) x * scale_real);
    c.imag = imag_min + ((float) y * scale_imag);
    color = cal_pixel(c);
    display(x, y, color);
}</pre>
```

where display() is a routine to display the pixel (x, y) at the computed color.





Mandelbrot Set Computation (5)

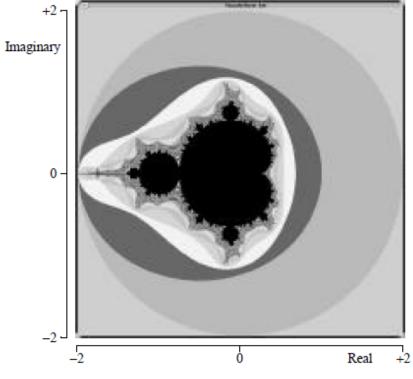


Figure 3.4 Mandelbrot set.



Parallelization of Mandelbrot Computation (1)

Static Task Assignment

for (i = 0, row = 0; i < 48; i++, row = row + 10)/* for each process*/

Master

```
/* send row no.*/
    send(&row, P;);
for (i = 0; i < (480 * 640); i++) {/* from processes, any order */}
    recv(&c, &color, P<sub>NV</sub>); /* receive coordinates/colors */
                             /* display pixel on screen */
    display(c, color);
}
Slave (process i)
recv(&row, Pmaster);
                             /* receive row no. */
for (x = 0; x < disp width; x++)/* screen coordinates x and y */
    for (y = row; y < (row + 10); y++) {
     c.real = min real + ((float) x * scale real);
      c.imag = min imag + ((float) y * scale imag);
     color = cal pixel(c);
      send(&c, &color, Pmaster);/* send coords, color to master */
```





Parallelization of Mandelbrot Computation (2)

Dynamic Task Assignment Work Pool/Processor Farms

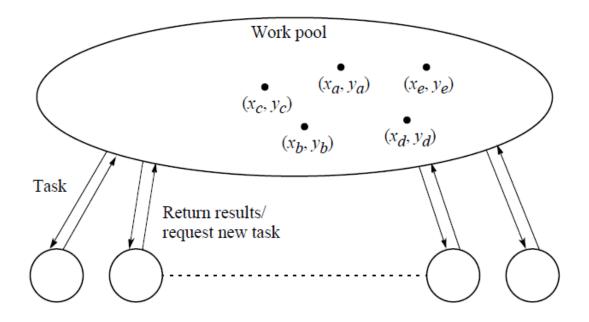


Figure 3.5 Work pool approach.



Parallelization of Mandelbrot Computation (3)

Coding for Work Pool Approach

```
Master
                                   /* counter for termination*/
count = 0;
                                  /* row being sent */
row = 0;
for (k = 0; k < procno; k++) { /* assuming procno<disp height */
    send(&row, Pk, data_tag); /* send initial row to process */
    count++;
                                  /* count rows sent */
                                  /* next row */
    row++;
do {
    recv (&slave, &r, color, P<sub>NNV</sub>, result tag);
    count--;
                                 /* reduce count as rows received */
    if (row < disp height) {</pre>
      send (&row, P<sub>slave</sub>, data_tag);
                                            /* send next row */
                                             /* next row */
      row++;
      count++;
    } else
      send (&row, P<sub>slave</sub>, terminator_tag); /* terminate */
    rows recv++;
```

/* display row */

display (r, color);

} while (count > 0);





Parallelization of Mandelbrot Computation (4)

Slave

```
recv(y, Pmaster, ANYTAG, source_tag);/* receive 1st row to compute */
while (source_tag == data_tag) {
    c.imag = imag_min + ((float) y * scale_imag);
    for (x = 0; x < disp_width; x++) {/* compute row colors */
        c.real = real_min + ((float) x * scale_real);
        color[x] = cal_pixel(c);
    }
    send(&i, &y, color, Pmaster, result_tag);/* row colors to master */
    recv(y, Pmaster, source_tag); /* receive next row */
};</pre>
```





Parallelization of Mandelbrot Computation (5)

Rows outstanding in slaves (count)

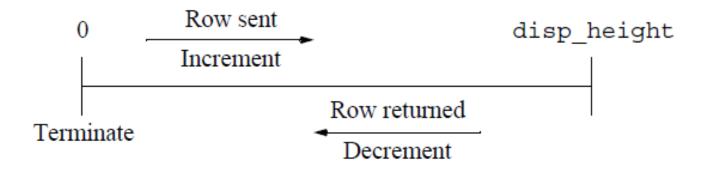


Figure 3.6 Counter termination.



Parallelization of Mandelbrot Computation (6)

Analysis

Sequential

 $t_{S} \leq \max \times n = O(n)$

Parallel program

Phase 1: Communication - Row number is sent to each slave

$$t_{\text{comm1}} = s(t_{\text{startup}} + t_{\text{data}})$$

Phase 2: Computation - Slaves perform their Mandelbrot computation in parallel

$$t_{\text{comp}} \le \frac{\max \times n}{s}$$

Phase 3: Communication - Results passed back to master using individual sends

$$t_{\text{comm2}} = \frac{n}{s} (t_{\text{startup}} + t_{\text{data}})$$

Overall

$$t_p \le \frac{\max \times n}{s} + \left(\frac{n}{s} + s\right)(t_{\text{startup}} + t_{\text{data}})$$



Monte Carlo Methods (1)

Basis of Monte Carlo methods is the use of random selections in calculations.

Example - To calculate π

A circle is formed within a square. Circle has unit radius so that square has sides 2×2 .

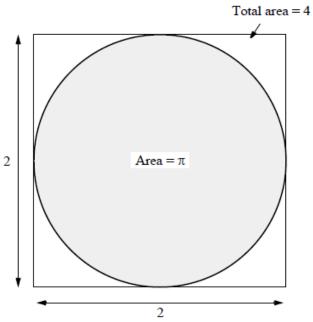


Figure 3.7 Computing π by a Monte Carlo method.



Monte Carlo Methods (2)

The ratio of the area of the circle to the square is given by

$$\frac{\text{Area of circle}}{\text{Area of square}} = \frac{\pi(1)^2}{2 \times 2} = \frac{\pi}{4}$$

Points within the square are chosen randomly and a score is kept of how many points happen to lie within the circle.

The fraction of points within the circle will be $\pi/4$, given a sufficient number of randomly selected samples.



Monte Carlo Methods (3)

Computing an Integral

One quadrant of the construction in Figure 3.7 can be described by the integral

$$\int_0^1 \sqrt{1 - x^2} \, dx \, = \, \frac{\pi}{4}$$

A random pair of numbers, (x_r, y_r) would be generated, each between 0 and 1, and then counted as in circle if $y_r \le \sqrt{1 - x_r^2}$; that is, $y_r^2 + x_r^2 \le 1$.

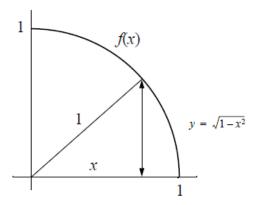


Figure 3.8 Function being integrated in computing π by a Monte Carlo method.



Monte Carlo Methods (4)

Alternative (better) Method

Use the random values of x to compute f(x) and sum the values of f(x):

Area =
$$\int_{x_1}^{x_2} f(x) dx = \lim_{N \to \infty} \frac{1}{N} \sum_{i=1}^{N} f(x_i)(x_2 - x_1)$$

where x_r are randomly generated values of x between x_1 and x_2 .



Monte Carlo Methods (5)

Example

Computing the integral

$$I = \int_{x_1}^{x_2} (x^2 - 3x) \, dx$$

Sequential Code

The routine randv(x1, x2) returns a pseudorandom number between x1 and x2.



Monte Carlo Methods (6)

Parallel Implementation

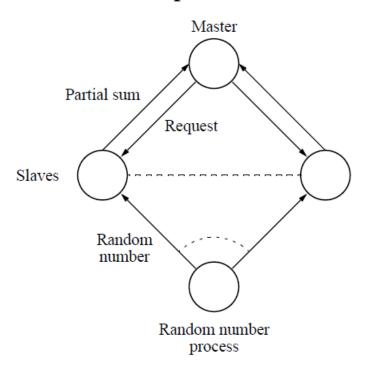


Figure 3.9 Parallel Monte Carlo integration.





Monte Carlo Methods (7)

Pseudocode

Master



Monte Carlo Methods (8)

Slave

```
sum = 0;
send(P<sub>master</sub>, req_tag);
recv(xr, &n, P<sub>master</sub>, source_tag);
while (source_tag == compute_tag) {
    for (i = 0; i < n; i++)
        sum = sum + xr[i] * xr[i] - 3 * xr[i];
    send(P<sub>master</sub>, req_tag);
    recv(xr, &n, P<sub>master</sub>, source_tag);
};
reduce_add(&sum, P<sub>qroup</sub>);
```



Random Number Generation

The most popular way of creating a pueudorandom number sequence:

$$x_1, x_2, x_3, \ldots, x_{i-1}, x_i, x_{i+1}, \ldots, x_{n-1}, x_n,$$

is by evaluating x_{i+1} from a carefully chosen function of x_i , often of the form

$$x_{i+1} = (ax_i + c) \mod m$$

where a, c, and m are constants chosen to create a sequence that has similar properties to truly random sequences.

- A good generator is with a = 16807, $m = 2^{31} 1$, and c = 0.
 - This generator creates a repeating sequence of $2^{31} 2$ different numbers



Parallel Random Number Generation

It turns out that

$$x_{i+1} = (ax_i + c) \bmod m$$

$$x_{i+k} = (Ax_i + C) \mod m$$

where $A = a^k \mod m$, $C = c(a^{k-1} + a^{n-2} + ... + a^1 + a^0) \mod m$, and k is a selected "jump" constant.

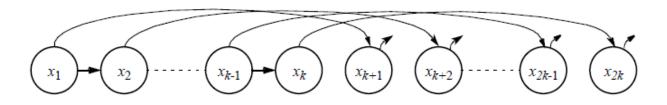


Figure 3.10 Parallel computation of a sequence.