

Lesson 4: Partitioning and Divide-and-Conquer strategies

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Partitioning strategies

Partitioning strategy is used as a generic term for any procedure based on the division of a problem into parts.

Example (adding numbers): The goal is to add a sequence of n numbers x_0, \ldots, x_{n-1} . The algorithm divides the sequence into p parts

$$x_0, \dots, x_{(n/p)-1}$$

 $x_{(n/p)}, \dots, x_{2(n/p)-1}$
 \vdots
 $x_{(p-1)(n/p)}, \dots, x_{p(n/p)-1} (= x_{n-1})$

The numbers in each subsequence are added independently, then the resulting partial sums are added to get the final result.



..Partitioning strategies

1st implementation (using separate send() and recv()):

Slave:

```
recv(numbers, r, P_{master});

partSum = 0;

for (i=0; i<r; i++)

    partSum = partSum + numbers[i];

send(partSum, P_{master});
```



..Partitioning strategies

Analysis:

Sequential: Computation requires n-1 additions.

Hence time complexity is O(n).

Parallel (1st implementation):

- —1st communication: $t_{comm1} = p(t_{startup} + (n/p)t_{data})$
- —computation (in slaves): $t_{comp1} = n/p 1$
- —2nd communication (return partial sum):

$$t_{comm1} = p(t_{startup} + t_{data})$$

- —final computation: $t_{comp1} = p 1$
- —overall:

$$t_p = 2pt_{startup} + (n+p)t_{data} + n/p + p - 2$$



.. Partitioning strategies

Conclusion:

- The computation part is *decreasing* from n-1 to n/p+p-2.
- The communication part is *linear* on both the size of data and the number of processes.
- The overall parallel time complexity is *worse* than sequential time complexity.
- To be useful, the slaves should have heavier computations (say, t(n) such that the $t(n) t(n/p) > 2pt_{startup} + (n+p)t_{data}$)

..Partitioning strategies

2nd implementation (using broadcast/multicast):

Master:

```
r = n/p
bcast (numbers, r, P_{master}, P_{group});
sum = 0;
for (i=0; i< p; i++)
       recv (partSum, P_{any});
       sum = sum + partSum;
Slave:
```

```
bcast (numbers, r, P_{master}, P_{group});
start = slaveNumber * r;
end = start + r;
partSum = 0;
for (i=start; i<end; i++)</pre>
       partSum = partSum + numbers[i];
send (partSum, P_{master});
```



..Partitioning strategies

3rd implementation (using scatter/reduce):

Master:

```
r = n/p

scatter(numbers, r, P_{master}, P_{group});

reduceAdd(sum, P_{master}, P_{group});
```

Slave:

```
scatter (numbers, r, P_{master}, P_{group});

start = slaveNumber * r;

end = start + r;

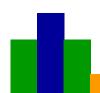
partSum = 0;

for (i=start; i<end; i++)

    partSum = partSum + numbers[i];

reduceAdd (partSum, P_{master}, P_{group});
```

Notice: The analysis of the 2nd or 3rd implementation is similar. (The real values depend on the network type and the particular implementations of broadcasting procedures.)



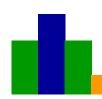
Divide-and-conquer is a particular partitioning strategy where the subproblems are of the same form as the larger problem.

Hence one can recursively apply the partition procedure to get smaller and smaller problems.

Add a list of numbers:

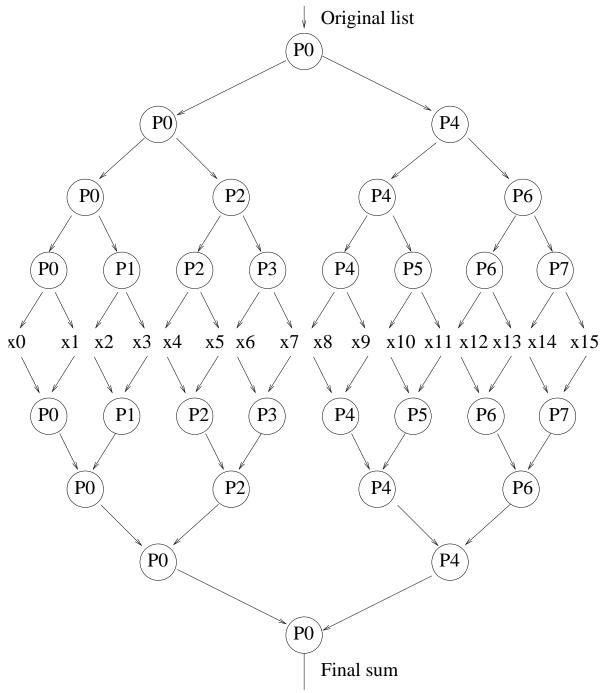
```
int add(list) {
  if (numbersOfElements(list) =< 2) return theirSum;
  else{
         divide(list, list1, list2);
         return(add(list1) + add(list2));
     }
}</pre>
```

Notice: This is a general procedure which may be implemented either in a sequential or a parallel way.



Parallel implementation:

- (a) In a first stage, the list is recursively divided till each process receives a two-element list.
- (b) In the next stage the partial results are collected using an opposite tree structure.



Parallel code (Process P_i): Let

```
k(i) = \begin{cases} \bullet \ r - \text{if } i = 0 \text{ (and there are } 2^r \text{ processes)} \\ \bullet \text{ the greatest power of 2 which divides } i - \text{otherwise} \\ \text{[i.e., } 2^{k(i)} \text{ divides } i, \text{ but } 2^{k(i)+1} \text{ does not divide } i \end{cases}
               if (!(myRank == 0)) recv(list,P_{myRank-2^{k(i)}});
               for (i=k-1; i>=0; i--)
                        divide (list, list, list2);
                         send(list2,P_{myRank+2^i});
               partSum = sumOf(list);
               for(i=0; i<k; i++){
                        recv(partSum2,P_{mvRank+2^i});
                        partSum = partSum + partSum2;
               if (!(myRank == 0)) send(partSum,P_{myRank-2^{k(i)}});
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```



Examples:

Process P0

```
divide(list, list, list2);
send(list2, P4);
divide(list, list, list2);
send(list2, P2);
divide(list, list, list2);
send(list2, P1);
partSum = sumOf(List);
recv(partSum2, P1);
partSum = partSum + partSum2;
recv(partSum2, P2);
partSum = partSum + partSum2;
recv(partSum2, P4);
partSum = partSum + partSum2;
```

Process P4 (k(4) = 2):

```
recv(list,P0);
divide(list,list,list2);
send(list2,P6);
divide(list,list,list2);
send(list2,P5);
partSum = sumOf(List);
recv(partSum2,P5);
partSum = partSum + partSum2;
recv(partSum2,P6);
partSum = partSum + partSum2;
send(partSum,P0);
```



Analysis: Assume $n = 2^k$ and $p = 2^r$. Then:

—division:
$$t_{comm1} = t_{startup} \log_2 p + (n/2 + n/4 + \dots + n/(2^r)) t_{data}$$

= $t_{startup} \log_2 p + [n(p-1)/p] t_{data}$

- —collecting results: $t_{comm2} = t_{startup} \log_2 p + t_{data} \log_2 p$
- —computation: $t_{comp} = n/p 1 + \log_2 p$
- —overall:

$$t_p = 2t_{startup}\log_2 p + [\log_2 p + n(p-1)/p]t_{data} + n/p - 1 + \log_2 p$$



Conclusion:

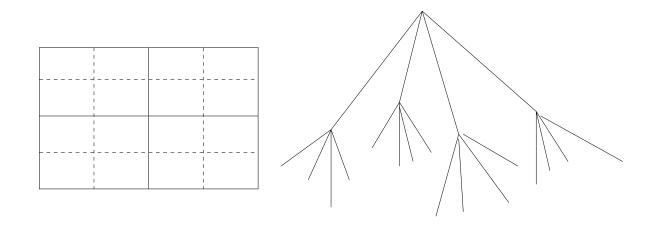
- The computation part is *decreasing*.
- The communication is still large: *linear* on the size of data, but now *logarithmic* on the number of processes.
- The algorithm becomes efficient when the computation part is large enough.

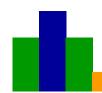
Notice that the work is not equally loaded on processes.



M-ary divide and conquer: A similar approach may be used when the problem is divided into more than two parts at each stage [if possible]. Now, m - ary trees are to be used, where m is the branching degree used at each stage.

Example: Such a technique may be applied to image processing by dividing each dimension into two parts at each stage, hence 4-ary trees are to be used [trees in which each node has four children.]





Case studies: 1. Sorting

Sorting using bucket sort:

1. The range of the numbers, say the interval [min, max), is divided into p equal regions:

```
[min, min + (max - min)/p) (Range 1)

[min + (max - min)/p, min + 2(max - min)/p) (Range 2)

\vdots

[min + (p-1)(max - min)/p, max) (Range p)
```

- 2. For each region a *bucket* is assigned to hold the numbers that fall within that region.
- 3. The numbers are placed into the appropriate buckets.
- 4. Finally, the numbers from each bucket are sorted using a sequential algorithm.

Notice: Works well if the numbers are uniformly distributed in [min, max).



Sequential algorithm: Suppose there are n numbers uniformly distributed in [min, max). Then each bucket has an average of n/p numbers. Hence

$$t_s = n + p(n/p)\log_2(n/p)$$

where: \bullet *n* — time to parse the numbers and place into buckets

- *p* number of buckets
- $(n/p)\log_2(n/p)$ sequential time to sort n/p numbers of each bucket



Parallel time (1st version): Sort the buckets in parallel. Hence

$$t1_p = n + (n/p)\log_2(n/p) + t_{startup} + nt_{data}$$

where, comparing with the former equation,

- p in the 2nd term disappears (work in parallel)
- $t_{startup} + nt_{data}$ is added (send data, via broadcast)

Parallel time (2nd version): A further parallelization is possible by *placing numbers into buckets in parallel*, as well. Suppose we have *n* numbers and *p* regions.

- 1. Separate the unsorted numbers a_0, \ldots, a_{n-1} into groups of n/p elements: a_0 to $a_{n/p-1}, a_{n/p}$ to $a_{2n/p-1}, \ldots$, one for each processor $P_i (i = 1, \ldots, p)$.
- 2. Each processor P_i parses its group of n/p numbers and generates *small buckets* b_{i1}, \ldots, b_{ip} (b_{ik} contains the numbers that fall into Range k).
- 3. The small buckets are sent to appropriate processors (P_i sends b_{ik} to P_k)
- 4. Each processor sorts its bucket of numbers.

Analysis: Assume we have *n* numbers, *p* processes and the numbers are uniformly distributed into buckets. Then:

- 1. send numbers to processes: $t_{comm1} = t_{startup} + nt_{data}$ (broadcast)
- 2. generate small buckets: $t_{comp1} = n/p$
- 3. send small buckets: $t_{comm2} = (p-1)(t_{startup} + (n/p^2)t_{data})$ (if communications could overlap; otherwise multiply by p)
- 4. sort large buckets: $t_{comp} = (n/p)\log_2(n/p)$

Overall:

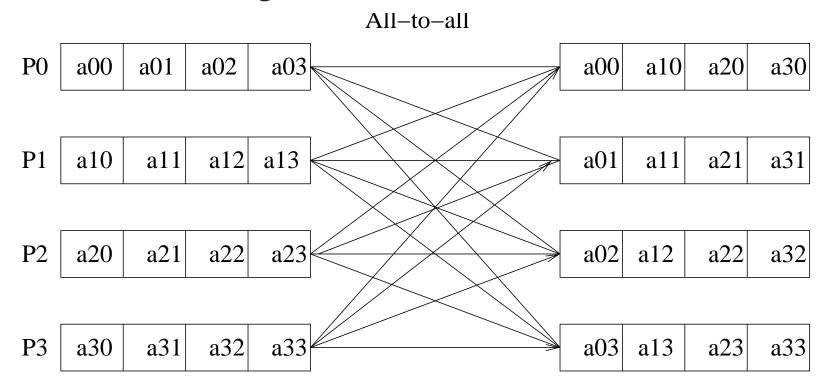
$$t_p = n/p + (n/p)\log_2(n/p) + pt_{startup} + (n+n/p^2)t_{data}$$

Conclusion: The computation time to generate buckets is decreasing from n to n/p, but the communication time is increasing.



All-to-all routines:

- A point where further improvement is possible is Step 3 where each process has to send specific data to all other processes.
- Such sort of communication may be made more efficient using standard routines, e.g., MPI_Alltoall().





Case studies: 2. Numerical integration

Numerical integration:

- A general divide-and-conquer technique divides the region continuously into parts.
- The procedure stops when some optimization function decides that the regions are sufficiently divided.

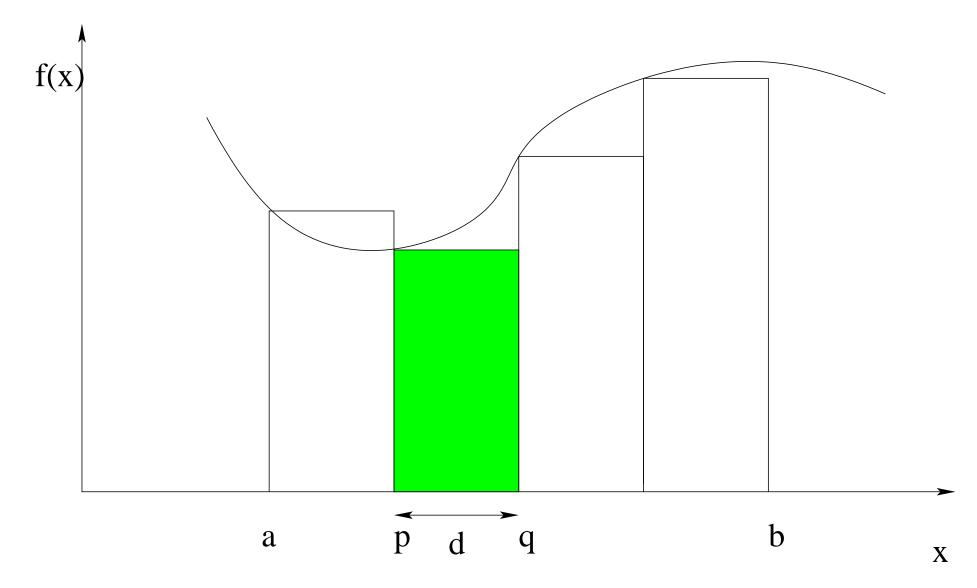
Example: For

$$I = \int_{a}^{b} f(x)dx$$

divide the range [a,b] into separate parts and perform the computation in each part using a separate process.



Approximation using rectangles:

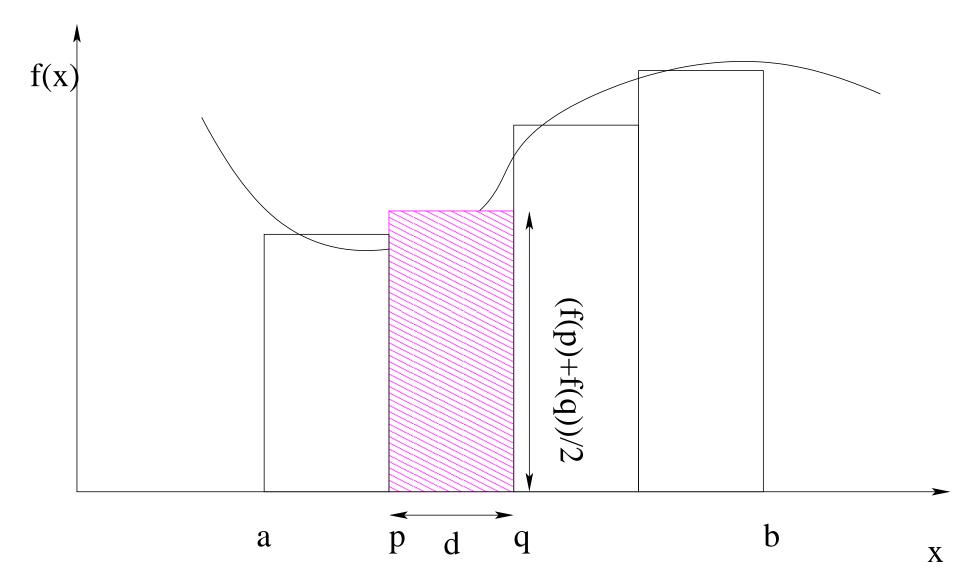


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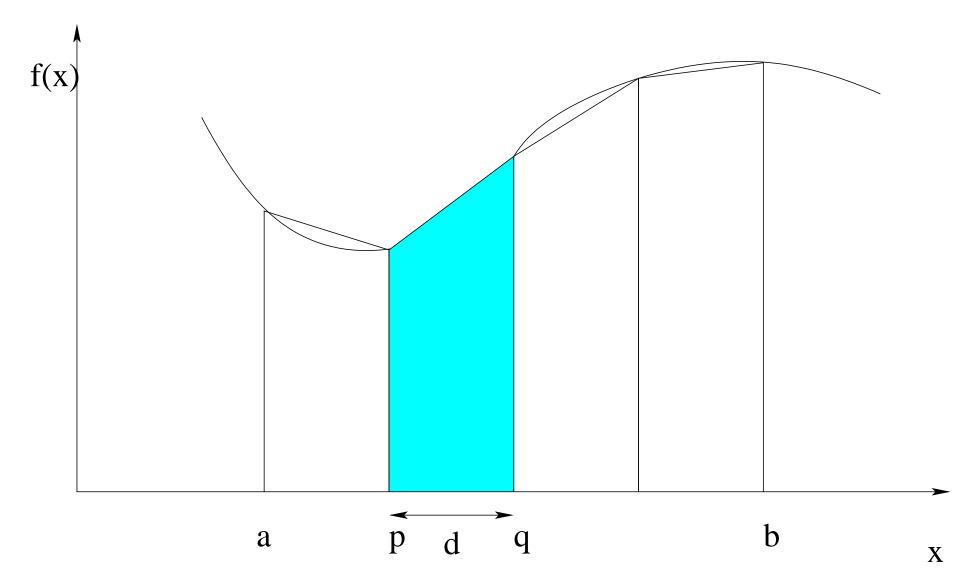


Better approximation using rectangles (average value):





Using trapezoidal approximations:



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Sample pseudocode:

(Static assignment, SPMD program, 2nd rectangular method)

Process P_i :

```
if (i == master) {
        printf('Enter number of interval'');
        scanf(%d, &n);
}
bcast(&n, P<sub>master</sub>, P<sub>group</sub>);
region = (b-a)/p;
start = a + region * i;
end = start + region;
area = 0.0;
for(x=start; x<end; x=x+d)
        area = area + f(x) + f(x+d);
area = 0.5 * area * d;
reduceAdd(&integral, &area, P<sub>group</sub>);
```



Final comments:

- This program is using a simple partitioning strategy. An approach using divide-and-conquer strategy may be used as well.
- The above program does not consider any termination condition. It may be (1) repeated for larger and larger *n* till a good approximation is obtained, or (2) the convergence criteria may be included into the program.
- Sometimes the function is not so smooth and a nonuniform division of the interval may fit better. E.g., using more processes closer to 0 gives a better approximation for $\int_{0.01}^{1} (x + \sin(1/x)) dx$.



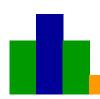
Case studies: 3. Gravitational N-body problem

Gravitational *N***-body problem:**

- This is a (computationally) difficult, practical problem which was wildly studied.
- Goal: find positions and movements of bodies in space subject to gravitational forces using Newtonian laws of physics.
- The gravitational force between two bodies of masses m_a and m_b is

 $F = \frac{Gm_am_b}{r^2}$

where *G* is the gravitational constant and *r* is the distance between the bodies.



• Subject to the (resulting) force *F*, a body will accelerate according to Newton's 2nd law

$$F = ma$$

where *m* is the mass of the body and *a* is the resulting acceleration.

• Let the time interval be Δt and v^{t+1} (resp. v^t) be the velocity at time t+1 (resp. t). Using the approximation

$$F = \frac{m(v^{t+1} - v^t)}{\Delta t}$$

one gets

$$v^{t+1} = v^t + \frac{F\Delta t}{m}$$

• The position of the body is changed from x^t to x^{t+1} by

$$x^{t+1} = x^t + v\Delta t$$

• After all these computations a new position is obtained and the procedure is repeated:

 $position \mapsto force \mapsto acceleration \mapsto velocity \mapsto position...$

• In a 3-dim space, the distance between bodies $a = (x_a, y_a, z_a)$ and $b = (x_b, y_b, z_b)$ is

$$r = \sqrt{(x_b - x_a)^2 + (y_b - y_a)^2 + (z_b - z_a)^2}$$

and the general force gives actions along each direction

$$F_x = k(x_b - x_a)/r$$
, $F_y = k(y_b - y_a)/r$, $F_z = k(z_b - z_a)/r$
where $k = \frac{Gm_am_b}{r^2}$



Sequential (pseudo)code: (sketched, one iteration step)



Parallel approach: (sketched)

- The algorithm uses $O(N^2)$ operations for one iteration (each body acts on each other body), hence is not practical for usual *N*-body problems where *N* is large.
- The time complexity is reduced using the observation that a cluster of distant bodies can be approximated as a single body (containing the total mass of the bodies in the cluster)

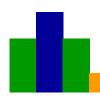


Barnes-Hut algorithm:

- Start with a cube containing all bodies. The cube is divided into eight subcubes.
- If a subcube contains no particles, then the subcube is deleted from further considerations
- If a subcube contains more than one body, then it is recursively divided until every subcube contains one body.

Notice: This way octtrees are obtained (each node has up to 8 children).

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- The total mass and the center of mass is stored in each node of the tree.
- The force on each body can be obtained by traversion the tree from the root and stopping when a clustering approximation may be used, usually,

$$r \ge d/\theta$$

where r - distance to the mass center, d - cube dimension, and θ constant, typically 1.0.

Both, the construction of the tree and the computation of all forces require $O(n \log n)$ time, hence the algorithm is $O(n \log n)$ (rather than the direct approach requiring $O(N^2)$ time).



The resulting tree may be not well balanced, hence a different approach will be to divide the space into subcubes having (closed to) an equal number of bodies. An illustration is below (in 2-dim case).

