Partitioning and Divide-and-Conquer Strategies Partitioning Strategies

Partitioning simply divides the problem into parts

Example - Adding a sequence of numbers

We might consider dividing the sequence into *m* parts of *n/m* numbers each, $(x_0 \dots x_{(n/m)})$. 1), $(x_{n/m} \dots x_{(2n/m)-1}), \dots, (x_{(m-1)n/m} \dots x_{n-1})$, at which point *m* processors (or processes) can each add one sequence independently to create partial sums.



Using separate send() has and recv()s

Master

	0110[3
	{
)* accumulate partial ammu */	imus_treq + mus = mus
	recv(∂_sum, P _{ANY});
* wait for results from slaves */	for (i = 0; i < m; i++; fm > i (0 = i) rot
	:0 = uns
)* send a numbers to slave */	send(&numbers[x], s, P _i);
(s + x = x	for $(\dot{t} = 0, x = 0; \dot{t} < m; \dot{t} < m$
<pre>*savets tor stampers for slaves'</pre>	:u/u = s

Slave

tsum, P_master ', 's send sum to master ',	люд») bnэг
m = part_sum + munbers[i];	าร_วรชช
0; i < s; i < sdd numbers */	= ए) उ०उ
:0 =	wns_jzed
ers, s, P _{master}); /* receive s numbers from master	quunu) voər

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Using Broadcast/multicast Routine

!u/u = s Master

* accumulate partial sums */ :wns-jzed + wns = wns i(%part_sum, PANY); * savels mort stiuser rot tisw */ }(++i ;m > i ;0 = i) rol :0 = uns bcast(numbers, s, P_{slave_group}); * send all numbers to slaves */

 $^{\prime *}$ number of numbers for slaves $^{\prime \prime}$

Slave

{

;(%part_sum, P_{master}); /* retarm of mus bras */ ![i]sredmun + mus_tred = mus_tred /* add numbers */ (++i ibne > i itaste = i) tot :0 = wns_jied end = start + s; $^{\prime \star}$ slave number obtained earlier $^{\star \prime}$ is * redmun_evels = frsts; $^{\rm \star}$ receive all numbers from matter $^{\rm \star}$ bcast(numbers, s, P_{master});

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Using scatter and reduce routines

Master

reduce_add(&så, p_{group}, root=master); * results from slaves */ ;(retter(numbers,&s,P_{group},root=master); * send numbers to slaves */ $^{\prime\,\star}$ areadmun to redmun $^{\star\,\prime}$!u/u = s

Slave

/* aredmun bbs */ scatter(numbers,&s,Pgroup,root=master); /* receive s numbers */

'* reduce_add(&part_su,mus_terve,root_matter);/* så,mus_tred%)bbs_ebucer */

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Divide and Conquer

Characterized by dividing a problem into subproblems that are of the same form as the larger problem. Further divisions into still smaller sub-problems are usually done by recursion

A sequential recursive definition for adding a list of numbers is

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Requires n - 1 additions with a time complexity of O(n).

Parallel

Using individual send and receive routines D and D and D are continues D and D are continued on D and D are contained on D are contained on D and D are contained on D and D are contained on D are contained on

 $(_{\text{startup}}) = m(t_{\text{startup}}) + q_{\text{startup}}) = m_{\text{startup}}$

Phase 2 — Computation

I - m/n = Iqmod

 $\label{eq:theta} Returning \ partial results using individual send and receive routines$

 $t_{\text{comm2}} = m(t_{\text{startup}} + t_{\text{data}})$

Phase 4 — Computation

Final accumulation

I - m = 2 dmost

[[byotheral]]

 $l_p = (l_{\text{comm}1} + l_{\text{comm}2}) + (l_{\text{comm}1} + l_{\text{comm}2}) = 2m l_{\text{startup}} + (m + m) + (m + m + m) + 2m l_{\text{startup}1} + 2m l_{\text{startup}1}$

OL

(u + u)O = di

We see that the parallel time complexity is worse than the sequential time complexity.

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Parallel Code

Suppose we statically create eight processors (or processes) to add a list of numbers.

Frocess P_0

/* divide al into two, sl and s2 */
divide (sl, sl, s2);
f* divide sl into two, sl and s2 */
divide (sl, sl, s2);
send(s2, P₂);
send(s2, P₂);
send(s2, P₁);
send(s2

The code for process P_4 might take the form

i (% fart_suml, P4);

:(%part_suml, P2);

iuns_jred + mus_jred = mus_jred

bart_sum = part_sum + part_suml;

Process P_4

recv(al, P₀); /* division phase */ divide(al, sl, s2); send(s2, P₆); for the send(s2, P₆); part_sum = *sl; recv(&part_sum1, P₆); part_sum = part_sum + part_sum1; send(&part_sum, P₆); part_sum = part_sum + part_sum1; send(&part_sum, P₀); send(&part_sum, P₀);

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The following phases are needed:

- Partition numbers.
- 2. Sort into small buckets.
- Send to large buckets.
- Sort large buckets.

Phase 1 — Computation and Communication

 $u = I^{duoo} \mu$

 $n_{\text{startup}}t + q_{\text{utrats}}t = 1 \text{mmos}t$

Phase 2 — Computation

 $d/u = z^{duios} t$

Phase 3 — Communication.

If all the communications could overlap:

 $t_{\text{comm3}} = (1 - 1)(t_{\text{startup}} + (n/p^2))$

Phase 4 — Computation

(d/u) Sol(d/u) = pqmos

Uverall

 $(q/n) \operatorname{gol}(q/n) + (_{\operatorname{stab}} \mathfrak{l}(^2 q/n) + _{\operatorname{qutual}} \mathfrak{l})(1-q) + q/n + n_{\operatorname{stab}} \mathfrak{l} + _{\operatorname{qutual}} \mathfrak{l} = q^{\mathfrak{l}}$

It is assumed that the numbers are uniformly distributed to obtain these formulas. The worst-case scenario would occur when all the numbers fell into one bucket!

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Further Parallelization

By partitioning the sequence into m regions, one region for each processor.

Each processor maintains p "small" buckets and separates the numbers in its region into its own small buckets.

These small buckets are then "emptied" into the p final buckets for sorting, which requires each processors to send one small bucket to each of the other processors (bucket i to processor). sor i).



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Numerical Integration

A general divide-and-conquer technique divides the region continually into parts and lets some optimization function decide when certain regions are sufficiently divided.

Example: numerical integration:

$$xp(x)f_q^p \int = I$$

Can divide the area into separate parts, each of which can be calculated by a separate process. Each region could be calculated using an approximation given by rectangles:











Three-Dimensional Space

Since the bodies are in a three-dimensional space, all values are vectors and have to be resolved into three directions, x, y, and z.

In a three-dimensional space having a coordinate system (x, y, z), the distance between the bodies at (x_0, y_0, z_0) and (x_b, y_b, z_b) is given by

$$\binom{p_2 - q_2}{7} + \binom{p_3 - q_3}{7} + \binom{p_3 - q_3}{7} + \frac{q_3}{7} + \frac{q_3}{7}$$

The forces are resolved in the three directions, using, for example,

$$E^{z} = \frac{b}{Quu^{n}u^{p}} \frac{5}{Quu^{n}u^{p}} \left(\frac{b}{2^{p}-z^{n}}\right)$$
$$E^{\lambda} = \frac{b}{Quu^{n}u^{p}} \left(\frac{b}{\lambda^{p}-\lambda^{n}}\right)$$
$$E^{\lambda} = \frac{b}{Quu^{n}u^{p}} \frac{5}{2^{p}} \frac{b}{2^{p}} \frac{b}{2^{p}}$$

where the particles are of mass m_a and m_b and have the coordinates (x_a, y_a, z_a) and (x_b, y_b, z_b) .

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Gravitational N-Body Problem

The objective is to find the positions and movements of the bodies in space (say planets) that are subject to gravitational forces from other bodies using Newtonian laws of physics.

The gravitational force between two bodies of masses m_a and m_b is given by

$$\frac{z'}{q u^{p} u g} = d$$

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

Subject to forces, a body will accelerate according to Newton's second law:

where m is the mass of the body, F is the force it experiences, and a is the resultant acceleration.

Let the time interval be Δt . Then, for a particular body of mass m, the force is given by

$$\frac{i\nabla}{\left(\frac{\Lambda}{1}-\frac{1}{1+1}\Lambda\right)u}=A$$

and a new velocity

$$\frac{u}{u} + v = v$$

where v^{t+1} is the velocity of the body at time t + 1 and v^{t} is the velocity of the body at time t.

If a body is moving at a velocity v over the time interval Δt , its position changes by

$$i \nabla v = x - \frac{1}{1+i} x$$

where x^{1} is its position at time t.

Once bodies move to new positions, the forces change and the computation has to be repeated.

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Distant cluster of bodies , or ו•• Center of mass ter of mass of the cluster: can be approximated as a single distant body of the total mass of the cluster sited at the cen-The time complexity can be reduced using the observation that a cluster of distant bodies esting N-body problems where N is very large. by each of the other N - 1 bodies. It is not feasible to use this direct algorithm for most inter-The algorithm is an $O(N^2)$ algorithm (for one iteration) as each of the N bodies is influenced Parallel Code



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$;_{W \in \Pi}[\dot{L}] V = [\dot{L}] V$ $i_{W \ominus \Pi}[\dot{L}] X = [\dot{L}] X$ *noitized pns yitoolev elaber */ for (i = 0; i < mmx; i++) { /* for each body */ ł /* (to respect to the position of the position (leap-frog) * (t) * (t) * (t) * (t) * im / Jb * F + [i]v = wen[i]v * compute new velocity and * compute force on ith body */ Force_routine(i); for (i = 0; i < N > i :0 = i) /* for each body */ * for each time period */ (++J ;xsmJ > J ;0 = J) rol

The overall gravitational N-body computation can be described by the algorithm

Sequential Code

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Barnes-Hut Algorithm

Starts with the whole space in which one cube contains the bodies (or particles).

First, this cube is divided into eight subcubes.

If a subcube contains no particles, the subcube is deleted from further consideration.

If a subcube contains more than one body, it is recursively divided until every subcube contains one body.

This process creates an octrree; that is, a tree with up to eight edges from each node. The leaves represent cells each containing one body.

After the tree has been constructed, the total mass and center of mass of the subcube is stored at each node.

The force on each body can then be obtained by traversing the tree starting at the root, stopping at a node when the clustering approximation can be used, e.g. when:

$\frac{\theta}{p} \ge J$

where θ is a constant typically 1.0 or less (θ is called the opening angle).

Constructing the tree requires a time of $O(n \log n)$, and so does computing all the forces, so that the overall time complexity of the method is $O(n \log n)$.

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Orthogonal Recursive Bisection

Example for a two-dimensional square area.

First, a vertical line is found that divides the area into two areas each with an equal number of bodies. For each area, a horizontal line is found that divides it into two areas each with an equal number of bodies. This is repeated until there are as many areas as processors, and then one processor is assigned to each area.

