



PARALLEL PROCESSING SYSTEMS

Chapter 3: Parallel Algorithm Complexity



Introduction


- Algorithms can be analyzed in two ways
 - precise
 - we count the number of operations performed in the worst or average case
 - (e.g., arithmetic, memory access, data transfer)
 - use these counts as indicators of algorithm complexity
 - is quite tedious and at sometimes impossible to perform.
 - approximate.
 - We resort to various approximate analysis methods
 - keeping in mind the error margin of the method applied
 - if such an approximate analysis indicates that Algorithm A is 1.2 times slower than Algorithm B
 - we may not be able to conclude with certainty that Algorithm B is better for the task at hand.

ASYMPTOTIC COMPLEXITY

- Suppose that
 - a parallel sorting algorithm requires $(\log_2 n)^2$ compare–exchange steps
 - another one $(\log_2 n)^2/2 + 2 \log_2 n$ steps
 - a third one $500 \log_2 n$ steps
- Ignoring lower-order terms and multiplicative constants
 - we may say that
 - the first two algorithms take on the order of $\log^2 n$ steps
 - the third one takes on the order of $\log n$ steps.
- The logic behind ignoring these details
 - when n becomes very large
 - eventually $\log n$ will exceed any constant value.
 - for such large values of n and any values of the constants c and c'
 - an algorithm with $c \log n$ is asymptotically better than an algorithm with $c' \log^2 n$



ASYMPTOTIC COMPLEXITY

- n must indeed be very large for $\log n$ to overshadow the constant 500
 - In practice
 - we do not totally ignore the constant factors
 - We take a two-step approach
 - First, we determine which algorithm is likely to be better for large problem sizes
 - through asymptotic analysis
 - If we have reason to doubt this conclusion
 - we resort to an exact analysis to determine the constant factors involved
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ASYMPTOTIC COMPLEXITY

- Some notations

- Given two functions $f(n)$ and $g(n)$ we define the relationships

- “O” (big-oh)

- $f(n) = O(g(n))$ if $\exists c, n_0$ such that $\forall n > n_0$ we have $f(n) < c g(n)$

- “ Ω ” (big-omega)

- $f(n) = \Omega(g(n))$ if $\exists c, n_0$ such that $\forall n > n_0$ we have $f(n) > c g(n)$

- “ Θ ” (theta)

- $f(n) = \Theta(g(n))$ if $\exists c, c', n_0$ such that $\forall n > n_0$ we have $c g(n) < f(n) < c' g(n)$

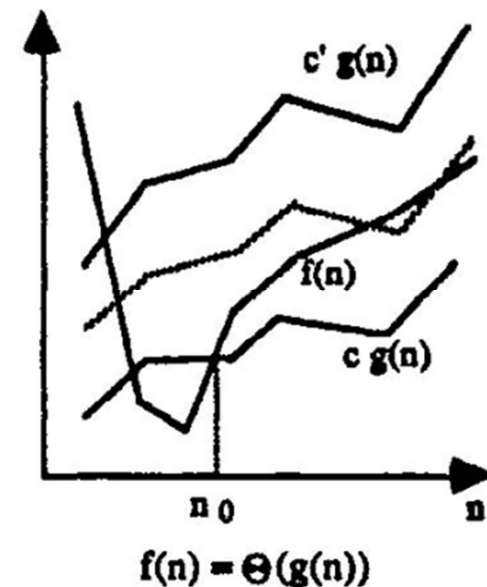
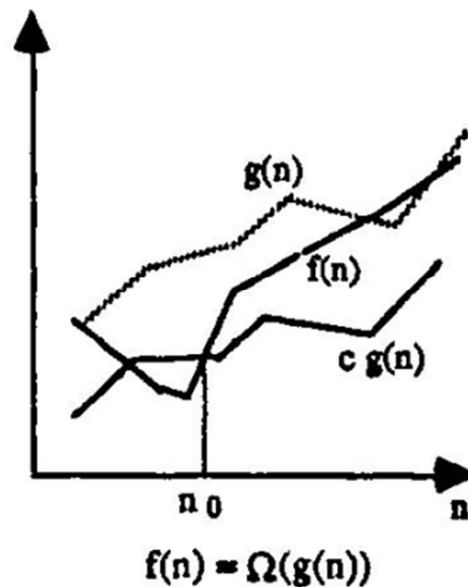
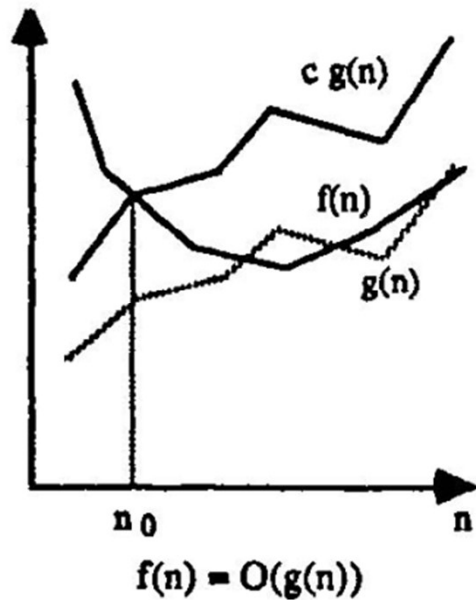
- $f(n) = \Theta(g(n))$ iff $f(n) = O(g(n))$ and $f(n) = \Omega(g(n))$

ASYMPTOTIC COMPLEXITY

- notations allow us to compare the growth rates of different functions
 - $f(n) = O(g(n))$ means that $f(n)$ grows no faster than $g(n)$
 - for n sufficiently large and a suitably chosen constant c
 - $f(n)$ always remains below $c g(n)$
 - $f(n) = \Omega(g(n))$ means that $f(n)$ grows at least as fast as $g(n)$
 - Eventually $f(n)$ will exceed $c g(n)$ for all n beyond n_0
 - $f(n) = \Theta(g(n))$ means that $f(n)$ and $g(n)$ grow at about the same rate
 - value of $f(n)$ is always bounded by $c g(n)$ and $c' g(n)$ (for $n > n_0$)

ASYMPTOTIC COMPLEXITY

- notations allow us to compare the growth rates of different functions



ASYMPTOTIC COMPLEXITY

- In other words

“The rate of growth of $f(n)$ is ___ that of $g(n)$.”

we can fill in the blank with the relational symbol ($<$, \leq , $=$, \geq , $>$) to the left of the defined relations shown below:

$<$	$f(n) = o(g(n))$	$\lim_{n \rightarrow \infty} f(n)/g(n) = 0$ {read little-oh of $g(n)$ }
\leq	$f(n) = O(g(n))$	{big-oh}
$=$	$f(n) = \Theta(g(n))$ or $\theta(g(n))$	{theta}
\geq	$f(n) = \Omega(g(n))$	{big-omega}
$>$	$f(n) = \omega(g(n))$	$\lim_{n \rightarrow \infty} f(n)/g(n) = \infty$ {little-omega}


ASYMPTOTIC COMPLEXITY

- At a very coarse level

Sublinear	$O(1)$	constant-time
	$O(\log \log n)$	double logarithmic
	$O(\log n)$	logarithmic
	$O(\log^k n)$	polylogarithmic; k is a constant
	$O(n^a)$	$a < 1$ is a constant; e.g., $O(\sqrt{n})$ for $a = 1/2$
	$O(n / \log^k n)$	k is a constant
Linear	$O(n)$	
Superlinear	$O(n \log^k n)$	
	$O(n^c)$	polynomial; $c > 1$ is a constant; e.g., $O(n\sqrt{n})$ for $c = 3/2$
	$O(2^n)$	exponential
	$O(2^{2^n})$	double exponential




Complexity classes

- Problems are divided into several complexity classes
 - Based on their running times on a single-processor
 - Problems said to belong to the P class
 - running times are upper bounded by polynomials in n
 - generally considered to be tractable.
 - Even if the polynomial is of a high degree
 - there is still hope that a reasonable running time may be obtained
 - by improvements in the algorithm or in computer performance
- 




Complexity classes

- problems said to belong to NP (nondeterministic polynomial) class
 - best known deterministic algorithm runs in exponential time
 - But the correctness of the solution can be verified in polynomial time
 - They are intractable
 - E.g., subset-sum problem
 - Given a set of n integers and a target sum s
 - determine if a subset of the integers in the given set add up to s
- 




Complexity classes

- problems said to belong to NP-complete class
 - any problem in NP can be transformed to any one of these problems
 - by a computationally efficient process
 - The subset-sum problem is known to be NP-complete
 - if one ever finds an efficient solution one of these problems
 - this proves $P = NP$
 - are the “hardest” problems in the NP class
 - proving that a problem is NP-complete removes any hope of finding an efficient algorithm
- 



Complexity classes

- problems said to belong to NP-hard class
 - problems that are not even in NP
 - verifying that a claimed solution to such a problem is correct is currently intractable
 - we do not know to be in NP
 - but do know that any NP problem can be reduced to it by a polynomial-time algorithm
 - are at least as hard as any NP problem
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Complexity classes

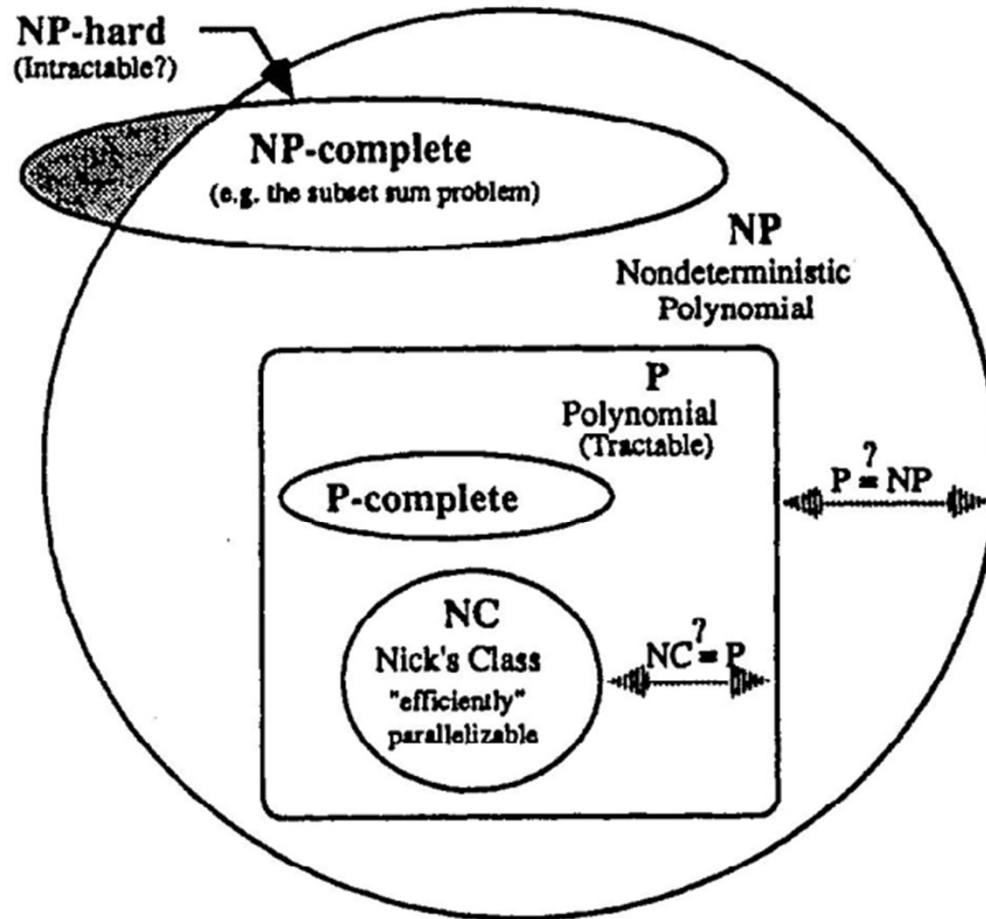





Figure 3.4. A conceptual view of complexity classes and their relationships.




Parallelizable tasks and the NC class

- parallel processing
 - is generally of no avail for solving NP problems
 - A problem that takes 400 billion centuries on a uniprocessor
 - if it can be perfectly parallelized over 1 billion processors
 - It would still take 400 centuries
 - is useful for speeding up the execution time of the problems in P
 - even a factor of 1000 speed-up can mean the difference between practicality and impracticality
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Parallelizable tasks and the NC class

- The NC class
 - efficiently parallelizable problems in P
 - defined as
 - problems that can be solved
 - By at most polylogarithmic in the problem size n , i.e., $T(p) = O(\log^k n)$ for some constant k
 - using no more than a polynomial number $p = O(n^l)$ of processors
 - Sorting is a good example
- 




Parallel programming paradigms

- Divide and Conquer
 - Decompose the problem of size n into two or more “smaller” subproblems.
 - takes $T_d(n)$ time when done in parallel
 - Solve the subproblems independently and obtain the corresponding results
 - the time T_s to solve them will likely be less than $T(n)$
 - Finally, combine the results of the subproblems to compute the answer to the original problem.
 - If the combining can be done in time $T_c(n)$,
 - the total computation time is given by
$$T(n) = T_d(n) + T_s + T_c(n).$$
- 



Parallel programming paradigms

- Divide and Conquer
 - is perhaps the most important tool for devising parallel algorithms
 - E.g., sorting a list of n keys
 - decompose the list into two halves
 - sort the two sublists independently in parallel
 - merge the two sorted sublists into a single sorted list
- 



Parallel programming paradigms

- Randomization
 - Often balanced divide and conquer is impossible, or computationally difficult
 - Obstacles for achievable effective speed-up
 - Large decomposition and combining overheads
 - wide variations in the solution times of the subproblems
 - it might be possible to use random decisions
 - lead to good results with very high probability.
 - Has led to the solution of many otherwise unsolvable problems.




Parallel programming paradigms

- Randomization
 - Again, sorting provides a good example
 - each of p processors begins with a sublist of size n/p
 - each processor selects a random sample of size k from its local sublist
 - The kp samples from all processors form a smaller list that can be readily sorted
 - on a single processor
 - or using an efficient parallel algorithm for small lists.
 - this sorted list of samples is now divided into p equal segments
 - the beginning values in the p segments used as thresholds to divide the original into p sublists
 - the lengths of these sublists will be approximately balanced with high probability





Parallel programming paradigms

- Randomization
 - Again, sorting provides a good example
 - The sorting problem has thus been transformed into
 - an initial random sampling
 - a small sorting problem for the k_p samples
 - broadcasting of the p threshold values to all processors
 - permutation of the elements among the processors according to the p threshold values
 - p independent sorting problems of approximate size n/p
 - The average case running time can be quite good
 - However, there is no useful worst-case guarantee on its running time.
- 



Parallel programming paradigms

- Randomization
 - can be applied in several other ways
 - Input randomization
 - used to avoid bad data patterns that a particular algorithm
 - is efficient on the average
 - might have close to worst-case performance
 - Random search
 - Control randomization
 - Symmetry breaking



Parallel programming paradigms

- Randomization
 - can be applied in several other ways
 - Input randomization
 - Random search
 - Searching a large space for an abundant element
 - random search can lead to very good average-case performance
 - A deterministic linear search can lead to poor performance
 - if all of the desired elements are clustered together
 - Control randomization
 - Symmetry breaking



Parallel programming paradigms

- Randomization
 - can be applied in several other ways
 - Input randomization
 - Random search
 - Control randomization
 - Randomly choosing the algorithm or an algorithm parameter
 - avoid consistently experiencing close to worst-case performance
 - with one algorithm
 - For some unfortunate distribution of inputs
 - Symmetry breaking



Parallel programming paradigms

- Randomization
 - can be applied in several other ways
 - Input randomization
 - Random search
 - Control randomization
 - Symmetry breaking
 - deterministic processes may exhibit a cyclic behavior that leads to deadlock
 - Randomization can be used to break the symmetry and thus the deadlock



Parallel programming paradigms

- Approximation
 - Iterative numerical methods often use approximation
 - begin with some rough estimates for the answers
 - successively refine these estimates using numerical calculations
 - Advantage
 - fairly precise results can be obtained rather quickly
 - additional iterations may be used to increase the precision if desired
 - It is a powerful method for time/cost/accuracy trade-offs because
 - the computations for each iteration can be easily parallelized over any number p of processors
 - the computation can still be performed at lower precision in case of deadline limitations
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