CS 124 Programming Assignment 2

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Caching for the Naive Algorithm

When multiplying two matrices A*B = C using the naive matrix multiplication method, the elements of C are given as follows

$$C[i][j] = \sum_k A[i][k] * B[k][j]$$

This is naturally implemented with 3 nested for loops iterating over i, j, and k. There are 3! = 6 possible permutations of their order. I tested all of them experimentally to determine which was best. Runtime measurements were taken by running the naive algorithm on $n \times n$ matrices with randomly generated entries for n = 1200.

ordering	runtime
i, j, k	5.81s
i, k, j	1.73s
j, i, k	5.54s
j, k, i	25.30s
k, i, j	2.05s
k, j, i	24.67s

The ordering i, k, j produces the best results. The caching behaviour of that ordering can be further optimized by saving the intermediate value of each A[i][k] before we iterate over j. This allows the innermost loop, with iterates over j, to only access a single row of matrix B, in order, via calls to access B[k][j]. Since the matrices are stored in row-major order, this is very cache friendly. Saving the intermediate value of A[i][k] prevents the cache from thrashing as it must repeatedly load up A[i] to recalculate A[i][k], which is wasted work since A[i][k] isn't changing during the innermost loop over j. This improves runtime of the naive algorithm by an additional factor of 2-3, which drastically affects my results for n_0 (discussed more in the experimental results section).

Strassen's Algorithm Implementation

Interaction With the Naive Algorithm

Strassen's algorithm is implemented in a method called strassen. However, the main method instead calls the multiply method, which returns strassen(parameters) if $n \geq n_0$. Otherwise, multiply returns the result of executing the naive algorithm. When strassen() recurses, it recursively calls multiply(), which then chooses whether to use strassen() or default to the naive method. strassen() never directly calls strassen(). The reason I chose to do it like this, even though it makes the recursion slightly more complicated, is that it allows the initial caller (ie the main method) to not have to worry about whether strassen's algorithm is being used on the first iteration or not. The main method simply calls multiply. Also, since strassen() must recurse in many places, strassen() would have to repeat the work of making that decision many times. This design would also make it easier if we wanted to add other matrix multiplication algorithms that were optimal under a different set of conditions.

Reducing Memory Allocation and Data Copying

Wikipedia defines Strassen's algorithm with the following variables:

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\begin{split} \mathbf{M}_1 &:= (\mathbf{A}_{1,1} + \mathbf{A}_{2,2})(\mathbf{B}_{1,1} + \mathbf{B}_{2,2}) \\ \mathbf{M}_2 &:= (\mathbf{A}_{2,1} + \mathbf{A}_{2,2})\mathbf{B}_{1,1} \\ \mathbf{M}_3 &:= \mathbf{A}_{1,1}(\mathbf{B}_{1,2} - \mathbf{B}_{2,2}) \\ \mathbf{M}_4 &:= \mathbf{A}_{2,2}(\mathbf{B}_{2,1} - \mathbf{B}_{1,1}) \\ \mathbf{M}_5 &:= (\mathbf{A}_{1,1} + \mathbf{A}_{1,2})\mathbf{B}_{2,2} \\ \mathbf{M}_6 &:= (\mathbf{A}_{2,1} - \mathbf{A}_{1,1})(\mathbf{B}_{1,1} + \mathbf{B}_{1,2}) \\ \mathbf{M}_7 &:= (\mathbf{A}_{1,2} - \mathbf{A}_{2,2})(\mathbf{B}_{2,1} + \mathbf{B}_{2,2}) \\ \text{only using 7 multiplications (one for each $M_k$) instead of 8. We may now express the $C_{i,j}$ in terms of $M_k$, like this: <math display="block"> \mathbf{C}_{1,1} &= \mathbf{M}_1 + \mathbf{M}_4 - \mathbf{M}_5 + \mathbf{M}_7 \\ \mathbf{C}_{1,2} &= \mathbf{M}_3 + \mathbf{M}_5 \\ \mathbf{C}_{2,1} &= \mathbf{M}_2 + \mathbf{M}_4 \\ \mathbf{C}_{2,2} &= \mathbf{M}_1 - \mathbf{M}_2 + \mathbf{M}_3 + \mathbf{M}_6 \end{split}
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I will use their variable naming scheme, both in my writeup and in my code. $A_{1,1}$ refers to the top right quadrant of A, and so on.

One way to implement this code would be to define every matrix named in the above description, then carry out every computation explicitely. My algorithm sacrifices that readibility in favor of having to define fewer variables (and therefore allocate more space) and copy data from one matrix to another fewer times. I define only 3 intermediate $\frac{n}{2} \times \frac{n}{2}$ matrices, and additionally an $n \times n$ matrix C to store my results. C is ultimately returned. The four submatrices of C, and the need to copy their contents into the final C, can be entirely eliminated by indexing correctly into C when adding on the various M's. The other 3 matrices I allocated I called M_{temp} , A_{temp} , and B_{temp} . There is no need to declare 7 matrices to store $M_1...M_7$ when one could instead reuse a single matrix M_{temp} , calculate its value, add it to the appropriate places in C, then reuse it for the next M. A_{temp} and B_{temp} contain the appropriate combination of

submatrices of A and B that correspond to the M we are currently calculating and saving in M_{temp} . It is unneccessary to ever define $A_{1,1}$, etc. We instead index appropriately into A and B when calculating A_{temp} and B_{temp} respectively. Not defining the 4 submatrices of A and B explicitly saves memory and removes the need to copy all of that data an additional time.

Other Optimizations

I/O is expensive, and was taking more time than the algorithm itself in my initial implementation (which read and wrote characters one at a time). I fixed this by using Java's built-in BufferedReader to read in the input, and by building then printing a single string with many linebreaks (meaning there's only a single call to file output). In java, the most efficient way to build a long string is to use the StringBuilder class, since Java replaces strings instead of modifying them by default because it treats strings as immutable objects.

Experimental Results for n_0

Testing Correctness

I used the unused first parameter as input for n_0 to determine what the experimental value for the optimal n-0 is. The program will default to whatever I determine that value to be if 0 is input for the first parameter (as the problem set description said it would be for grading). I wrote a bash script called check to help me test correctness, that would run the program first on an inputted $n_0 < n$, then on $n_0 > n$, and use the shell diff command to check that their respective outputs are identical. When $n_0 < n$, Strassen's algorithm will be used at least once. When $n_0 > n$, only the traditional matrix multiplication algorithm will be used. Therefore, this process checks my implementation of Strassen's algorithm against the traditional matrix multiplication algorithm.

Runtime for Different n_0

I wrote another bash script called testn0 to experiment efficiently with timing the runtime of different values of n and n_0 . I was curious if the value of the optimal crossover point might vary with n as well as with n_0 . It seemed natural to test powers of 2 for n_0 , and I experimented with several different values of n. I averaged over 5 trials to reduce statistical variation. The first entry in the table, with $n_0 > n$, represents using the traditional multiplication algorithm.

n	n_0	runtime
1000	> 1000	0.82s
1000	64	1.38s
1000	128	0.83s
1000	256	0.82s
1000	512	0.67s

 $n_0 \approx 512$ looks promising. We experiment with a few more values of n.

n	n_0	runtime
2000	> 2000	4.43s
2000	64	5.36s
2000	128	3.47s
2000	256	2.76s
2000	512	2.77s
2000	1024	3.83s

n	n_0	runtime
1500	> 1500	2.11s
1500	64	2.72 s
1500	128	2.10s
1500	256	1.68s
1500	512	1.45s
1500	1024	1.78s

n	n_0	runtime
2500	> 2500	$8.80 \mathrm{\ s}$
2500	64	13.59 s
2500	128	$8.60 \mathrm{\ s}$
2500	256	$6.85 \mathrm{\ s}$
2500	512	$6.33 \mathrm{\ s}$
2500	1024	$5.40 \; s$

 $n_0=512$ seems to be producing the best result, ie to use Strassen's algorithm for $n\geq 512$ and not use Strassen's algorithm for n<512. This was much higher than the values for n_0 that my classmates seemed to be getting. I am convinced that the majority of the difference comes not from slowness in my Strassen's algorithm, but from the optimizations I did for my naive multiplication algorithm. Not only did I rearrange the matrix order as suggested, but I also saved some intermediate values (discussed above). Saving intermediate values speeds up the naive multiplication algorithm by an additional factor of 2-3 in my tests, which has the consequence of pushing n_0 much higher than it would be if I disabled that. In fact, when I commented out that cacheing optimization, I got values of n_0 as low as 64.

This table gives results for n_0 with the correct i, j, k ordering but without the additional caching optimization on the naive algorithm.

n	n_0	runtime
1024	> 1024	1.9 s
1024	32	1.8 s
1024	64	1.4 s
1024	128	1.4 s
1024	256	$1.5 \mathrm{\ s}$
1024	512	$1.5 \mathrm{\ s}$
1024	1024	1.6 s

In this implementation (since n is a power of 2 and the program executes Strassen's algorithm if $n = n_0$), the last time Strassen's algorithm is used should be at either n = 64 or n = 128, with the naive algorithm beginning at n = 32 or n = 64.

Theoretical Results for n_0