

Computing made Difficult

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Chapter 1

Introduction

There are many ways in which the world tries to pretend that computing is easy. There are schemes that teach coding to children certainly starting from age 6. There are self-help books with titles along the lines of "Teach yourself programming in *so many days*". Almost every new serious programming language or software package will trumpet that it represents the next step in rendering computer use accessible to all. Finally one of the claims for Artificial Intelligence is that it means that everybody can develop computer systems by merely giving an informal explanation of what they want achieved. A rather small amount of web search (which is of course really easy!) will back up all the above. But what is hidden in all that enthusiasm is that the behaviour of computers and software and the design and construction and analysis of programs has astonishing layers of difficulty just beneath the shiny and simple-looking surface.

There are basically two reasons for investigating this difficulty. The first can obviously arise if you are trying to build a computer-based product or solve some particular problem and you come face to face with the unhappy fact that the world is messy and that naive or simplistic techniques are not good enough. If you are an optimist this may come as a nasty surprise! The second which is the one emphasised here is when you understand that clever techniques, fairly intricate details and plain weird results can be fascinating – and that coming face to face with some of them will let you build experience and understanding that lets you achieve more in the future.

So we start off here by noticing that many computing challenges had be presented in ways that do not need special skill or knowledge to appreciate. In plenty of cases there will be fairly obvious ways to start work towards resolving them. But then there are a dozen or more attitudes as well as problem categories that make it possible to unpick levels of difficulty seriously greater than were first apparent.

In some of the examples included here even the more complicated way to solve the problem will in fact be reasonably easy to grasp (once you have seen why it is necessary to go to all the trouble involved). In others it will involve somewhat messy data-structures or mathematics – but the cases we have chosen are intended to make these possible to understand and appreciate. Finally there are cases where there is no known solution or (worse) where it is known that there is no perfect solution. In such cases grasping how it is possible to demonstrate that something is impossible is of itself a challenge worth facing up to.

So here is a sort of catalogue of ways that let you start with a simple task and uncover the challenges concealed beneath its simplicity. For each idea there is a reference to a section later on here that works through an example in reasonable detail:

Look beneath the abstraction to understand how something is implemented or how it “really” works There are an amazing number of instances of this. Any time you ask a computer to sort some data, do a database lookup, compile a program, create a file, fetch a web page or encrypt a message that is supposed to be kept secret there is a great deal of technology that happily most people can just take for granted. One can view this as rather like the situation with almost all technology from digital watches to aeroplanes – almost anybody can take advantage of them. Plenty of people will be able to present an overview of how and why they work. But the details end up almost unimaginably complex. So the attitude of this point will underpin almost all of the sections here! Here an analogy might be that leaping on a motorbike or into a car and driving off probably seems pretty straightforward. Understanding in full detail how the engine and transmission works to a level where you can see how to design the next version to further optimise power or fuel efficiency or cost of manufacture is a very different business! For computers one can consider details not just of the program being written, but also the compiler used to process it, the operating system that is installed, how the instruction set of the computer is implemented in terms of gates and other logic circuitry down to the device physics behind the transistors inside. Of course for *practical* purposes it is not generally necessary to have specialist understanding of all the levels, but gaining it reveals just how difficult it is to get a proper in depth understanding of computation. A happy thought here is that the very earliest computers, while much less powerful and very much less convenient to use, were really a lot simpler and so looking back in history makes it possible to understand them from

top to bottom much more easily.

A product vs. a personal project This is perhaps best explained by an analogy with feature films. One could dream that anybody with a modern phone could capture video and a few friends could take parts in film. The status of the cameras on current phones which are inexpensive (at least in comparison with professional gear!) and can deliver remarkably high quality is analagous to the way in which these days computers are cheap, powerful and can so an amazing range of things following just a few clicks on an app or a web-resource. By choosing the plot you would not need any elaborate scenery or props, and you can avoid the need for potentially dangerous stunts. Some films pretend to be constructed by pasting together “found footage” extracted from the phones of characters (who will normally have vanished or died) further boosting the illusion that almost no serious equipment is needed. To get a clearer view of reality look at the list of credits at the end of a film and see the size of team really involved in even a seemingly simple one. Check the budget involved and let that give another indication of the complexity of the full process – even before the issues of marketing and distribution are taken into account. So the reality is that in general the finished product almost conceals the fact that it relies on the work of hundreds and sometimes thousands of people: scouting for locations, make up artists, lighting technicians, foley artists who add sound effects and many many more down to the transport catering teams who support the people more on the front line. In general while being a member of an audience you are almost unaware of all of this – save that if it was missing that would be very noticable. This situation can be seen as an “iceberg syndrome” where something that looks simple and where indeed an individual amateur could start to develop something in fact relies on such a wide range of techniques and skills that for any one person to master them all would represent a huge challenge. The message here is that making a home-movie is amazingly easy, but the simplicity there conceals the fact there is a great deal of difficulty in creating a block-buster. The same sort of things applies as between a bit of software created just for your own private use and a full-blown product.

Consideration of edge and worst cases There is a scheme called “Newton’s Method” or “The Newton-Raphson Iteration” that provides a simple to implement way of obtaining numeric solutions to equations. As a concrete and rather easy example it can be used on the equation

$x^3 - a = 0$ for some known value of a to find the cube root of a . But even with a case as easy as this there are challenges. How fast will it get an accurate result? It needs an initial guess for its answer to start from – how much does that value matter. An especially jolly issue for this case is that if one accepts that in many areas of mathematics and physics one is working with complex rather than real numbers it is necessary to accept that a will have three cube roots – and the issue of which one Newton’s method will deliver for you turns out to be a messier issue than you might have expected.

There are plenty of non-numeric instances of difficulty raised by asking about best and worst (and indeed average) cases in problems. For instance solving a Sudoku puzzle or planning how to return Rubic’s cube to a tidy state may not be totally trivial, but trying to identify the hardest possible Sudoku board or the most awkward starting point for a Rubic’s cube escalates the challenge sharply! Yet another example here come from “turtle graphics” – a computational model that has often been used in introductions of computers to the very young. In that world it can be quite easy to ask such questions as “If you continue with this pattern of movement will you ever find yourself exactly back where you began?” or “Might your turtle eventually fall off the end of the paper or indeed the world?” that make life tougher for yourself.

Insist on total correctness in every case. Pocket calculators and computers provide facilities to calculate trigonometric functions, logarithms and square roots of numbers. In normal circumstances one just trusts the computer. But there are two ways to make your own life harder. One is to express a fear that the computer will occasionally get bad results. Well in 1997 there were significant sales of computers that did not even always get division correct! How do you test things like this?

To go further note that with a computer the answers will always have been clipped to some limited precision. On a pocket calculator this may for instance be 8 decimal digits. The full perfect answer to a calculation has digits beyond that – so for instance if π is returned as 3.1415926 on a calculator (which can feel reasonable) there is an issue in that the true value is 3.14156265359... and the value quoted should have been rounded up because of the 5359... tail beyond where it ends. So the challenge is to evaluate all the elementary functions in such a way that for all possible inputs the result that is returned is correctly rounded – ie as accurate as is at all possible. And for this to be done first without intolerable extra cost and secondly with some proper

scheme that can certify that the goal of perfection has been achieved. Amazingly there are people who have been arranging that! A few years ago one could almost always just assume that the computer's results would be more than precise enough for all reasonable purposes, but now people are using rather low precision floating point arithmetic for big parallel computations in either graphics or artificial intelligence areas, and so these rather pedantic issues of precision come much closer to practical reality.

Portability For small programs that are only intended for use over a fairly short period of time and only by one person on their own computer it is not necessary to worry about portability. However larger projects raise more and more serious challenges. Windows, Macintosh, Android and Linux are amazingly not 100% compatible with each other, and it is often necessary to use techniques that apply to just one particular system. Intel and AMD processors, ARM and Risc-V are all different and are either at present in widespread use or may become so. There are areas where each of those needs custom treatment to achieve goals even as simple sounding as measuring time with the highest feasible precision. Use of a sufficiently high level language can conceal these issues by arranging that all the mess is embedded within the language and its associated library - but for our purposes that still leaves curious people with a need to know exactly what is going on. And very frequently common solutions that paper over differences carry costs that one might like to avoid.

Demand the fastest (or most compact) solution that could ever exist. The world of computer gaming is one where delivering the fastest frame-rate for the highest resolution version of the action is of serious commercial importance. And that is a matter of real directly measurable performance where there are basically few limits to what can be deployed to deliver it. That can lead to power-hungry and expensive video cards with amazingly elaborate driver software. Because much of that world is proprietary it is perhaps hard to get into, but it does illustrate that despite the fact that computers have become quite fast there are still areas where squeezing the best from them matters. This can involve both algorithm selection and coding style. For some application areas this aim to excel has the same sort of issues that mean that it will be different sets of athletes who dominate over 100 metres and over the 42195 metres of a marathon. In computing sometimes techniques that will be great for large problem instances will not show

up well on smaller inputs. There are two classical illustrations of this to be found in books on computer algorithms. One related to finding all then shortest routes from a given starting point to other locations in a maze (which is generally referred to as a graph). The other is simply multiplying numbers together. We will summarise and discuss each of these and various other cases where optimising for speed or size turns a reasonable problem into a tough one.

One nice term sometimes used is “galactic algorithms”. This is used to refer to ways of solving problems that in the long term would be better or faster than the simpler methods more commonly used, but achieve their closer approach to optimality at the cost of impressive overheads. Sometimes by way of a constant factor slowdown that would not be overcome by the fact that their cost grows a little more slowly than the standard schemes, and sometimes by such a level of implementation complication and mess that it will basically never make sense for some rather minor improvement. A few examples of this style will be covered here. The term “galactic” is to suggest something of the magnitude of problem instances where some of these would become competitive.

Challenging underpinning theory behind a simple-looking problem

I do not have nice text here, just a list of a few

1. Simplify algebraic formulae.
2. Telling if it is possible to find values for all the variables to make a boolean formula evaluate to TRUE.
3. the $3n+1$ challenge (ie Collatz).
4. Generate an unpredictable (ie random) sequence.

Constraints that rule out obvious approaches In some real sense all computer projects fall into this category because if you had a programming language or software package that aligned well enough with your task then everything would become easy. So to illustrate the point we will cover cases of extremely weak programming environments where it might not be obvious to start with that it will be possible to do anything much of interest. Each of the ones listed here and delved into in a bit more depth later can give insight into some particular aspect of computation:

1. Turing Machines and variants
2. Counter machines

3. Lambda-calculus and combinators
4. Cellular automata
5. Primitive recursive functions

Several of these start off seeming to be rather abstract and certainly not obviously practical, but for instance a highlight in one case is a report of how somebody has built a computer out of lego based on one of them such that in principle it is fully general purpose!

Seek good partial solutions to intractable problems Some – indeed many – of the challenges that arise in the real world turn out to be such that nobody knows how to solve them in general and there is serious reason to believe that that will always be the case. In a number of these it is even possible to prove that no general solution can be found. That opens the door to a lot of fun seeking either computer schemes that obtain approximate solutions or ones that sometimes or perhaps often get to the best answer, but are not guaranteed to complete the task. WQe provide several examples!

Use wilfully perverse techniques that achieve your goal. There can be great joy in exploring stupid ways of solving problems – and sometimes these emerge when a naive programmer submits their best efforts and you recognize that what they have achieved is a program that works but is magnificently slower than one might have hoped. Even experienced software engineers can end up delivering solutions that in retrospect can be seen to be pretty well absurd. We can illustrate this with cases from simple tasks the like of which are set as exercises in an elementary programming class: reversing a list, sorting some data and the like.

Challenges that make good puzzles Would any reasonable person want to invent unrealistic or pointless tasks? Why yes - those involved in recruitment of any sort might want "puzzle tasks" to set to candidates and while some questions they pose might merely test depth of knowledge, others want to look for inventiveness and the ability to think on the feet. Perhaps giving a preview of such cases undermines the joy in them, and maybe it will even be hard to tell which of the sections here have a component of this philosophy!

Future-proofing The landscape of computers has changed dramatically, and today there are still big new developments in prospect. Perhaps the two most visible are artificial intelligence and quantum computation, but one should really also note that the exploitation of the various

forms of massive parallelism that is now available represents a frontier. While some of the underpinning theoretical study of computation has remained valid for a long while, much practical software has a lifespan of a rather few decades. And segments of the theory that were central to all courses on computer science a generation back have much less clear-cut relevance than before. So we provide a few case studies that note both exciting ideas whose time seems to have passed and projects that against the odds have been kept alive.

Emphasize what must not happen as much as what will When specifying what a computer system will be for it is normal to describe what it will achieve. However the reality of things is that computer programs of any non-trivial size are liable to have flaws and will not always behave as intended. So there are cases where it becomes important to specify what must not happen as well as what should. And indeed in some cases the negatives are actually the important constraints. We will cover two areas that highlight this: security and safety-critical systems. The first can usefully be partitioned into the consideration of first encryption primitives (ie codes and the like) and then into protocols (how information is exchanged reliably between several parties). The counterpart to designing and building secure system is defeating same - and there is an amazing history of purportedly safe schemes being cracked open!

Cope with the inevitability of human error We all know that “To err is human”, and we are all used to observing that computer systems have flaws. If you were concerned with safety critical applications (think of control of anti-locking brakes on a car, of the guidance system for a missile or for a device to be implanted in somebody’s body) you may feel the need to reduce the chances of error as far as you possibly can. You may also wish to be resilient against all possible forms of hardware malfunction. There is no simple silver bullet.

Both the hardware of computers and the software they run will be constructed by humans – or these days perhaps by an artificially intelligent agent. Anybody who claims that what they deliver has been built by some special and really reliable tool should be asked about who created that tool and what basis there is for thinking it is perfect.

The inevitable result is that at least the initial version of any software must be presumed to be flawed. Testing – even very extensive testing – tends not to uncover all the defects. There are two approaches that aim toward perfection: (a) look for software building techniques

that minimise (note "minimise" rather than "eliminate") errors and (b) investigate ways to create formal proofs that the software ends up correct. We will have examples of or introductions to each of these.

Make sense of abstract graphical presentations Various schemes generate amazingly complicated images from rather simple recipes. Many people have come across the Mandelbrot Set which is one such instance, but we will point at a number more and consider how much understanding can be gleaned from reviewing the pictures.

Obviously these schemes for making life harder overlap in places, but it is also the case that several can all apply at the same time. The result is that challenges that started off seeming easy can end up causing quite severe headaches.

One might reasonably ask whether real people apply any of these principles in their lives. The next chapter sketches a particularly extreme case, but in reality a great deal of work starts with tasks that turn out to have depths that were not at all obvious at the start.

So it is very much the case that even though computing can be seen as really simple and we can keep telling ourselves "The computer only does what you tell it to", just beneath that attractive simple veneer there are a great many wonderful, messy, complicated and difficult details.

Chapter 2

A case study

This sketches one of the most impressive cases there has been of somebody starting a project and then going to town in the way they insisted that everything should be exactly to their taste.

In the early 1960s Donald Knuth began a project to write a book. It was a time of change in that somewhat before then serious books had been prepared by armies of compositors placing metal type in trays, but it was becoming clear that computer-aided publication was going to be the way forward. Knuth could have prepared his work the old fashioned way by writing it out by hand, getting a secretary to prepare a typed version from that and then letting a publishing house and printers loose to finalise things. That would have been the easy route.

However Knuth wanted greater control and knew that the standard route would not result in a publication up to his standards. So he diverted from his main book for a while and developed his own software to lay characters out on the page. There are multiple issues that have to be faced up to in that endeavour:

- Decide where to split lines and how to stretch text so that all lines end up neatly at the right margin;
- Kern letters properly – i.e. arrange the separation between individual letters based on their shapes so that the overall visual effect is of uniformity. This includes allowing for the way that letters in *italic sloping style* abut gracefully with upright text that surrounds.
- Present displayed mathematics well. Doing so leads to a need to handle Greek letters and a huge number of special symbols, to manage subscripts and superscripts, fraction bars, nested brackets of various

styles and vertical alignment (for instance in tables or presentations of matrices);

- Schemes for the generation of an incorporation of diagrams, generation of cross references, indexes and for the formatting of a title page and chapter headers.

The resulting system, \TeX ended up a great success and since then it has been used a standard tool for scientific publication. Many would have viewed that as at least as large a task as writing the book that motivated it.

But that was not the end! Historically fonts had been designed by specialists and cut in metal. The transcription of those to computers was not in a very well developed state, and in particular Knuth felt that the computer fonts available to him were not good enough. One issue was that when you have a single style of lettering the exact shapes of small letters (for instance to use in a subscript) should not be simply scaled versions of the standard versions, and similarly huge versions for use in titles are not merely magnified copies of the original. So Knuth invented notations for describing the shaped of characters and how those shapes changed with size. And using that he developed the “Computer Modern” family of typefaces including all the symbols he would need in his book. This is another contribution that has lasted but that can be viewed as a deep diversion rendering his project of writing a book much harder than might have been expected.

It is still the case that this story is incomplete. In writing the programs that could lay out text and define fonts Knuth was very aware that reading a program written by somebody else can be really difficult because the justification behind all the choices they have made is not visible. So he developed a scheme known as “literate programming” where code and a careful textual explanation of what it was doing (and why) were woven into a single document. By running one decoder over that document he could recover code to pass to a compiler and use, but with a different decoder he extracted a \TeX document by way of explanation. Within the single combined source the concept was that each fragment of code would be positioned adjacent to a careful explanation of it, and so as and when any changes were made it would be natural to keep the code and its documentation in step. Using this he was able to turn the commentary within the \TeX source code into a book “ \TeX the program” that could accompany his book that documented how to use it.

With all that in place he could get back to “The Art of Computer Programming” which expanded from an initial expectation of being a single volume into a sequence with each focussing a particular aspects of the subject.

And these became standard issue to all aspiring and practising computer scientists and TeX became the most common way for them to prepare papers and books.

This level of starting with a task that while large might have seemed reasonably straightforward and by demanding “better” escalating the project scope amazingly is a great illustration of the thoughts we present here. But for many practical reasons we will be concentrating on cases that do not demand quite the above levels of heroic energy!

Chapter 3

Counter Machines

To fully understand computation it can be good to strip things down to absolute basics. Doing so may at first make it seem as if nothing useful can be achieved, or that it could be that setting it up would be intolerably clumsy and laborious. Happily if you are prepared to take a few liberties with notation and if you are willing to view the time that a program would take to run as a total irrelevance things turn out to be less clumsy and less laborious than one might have expected. So in this chapter the emphasis will be on one particular way of looking at programming. It is the use of flowcharts, as often used when introducing computation to real beginners.

A text that aims to get people to understand what computers do and how they are driven might start with a version of instructions to make a cup of tea along the following lines:

1. Put water in the kettle;
2. Put tea in the teapot;
3. Switch kettle on;
4. wait a bit;
5. See if kettle is boiling: if not go back to step 4;
6. Pour (boiling) water into teapot;
7. Wait 3 minutes;
8. Done!

And having introduced the laborious step by step description of actions they then draw it out as a flowgraph with actions in square boxes, tests in squashed

boxed sitting on their corner and loads of arrows that indicate the flow from box to box. One location is identified as the starting point and another and where to stop.

Well pretty well any program that does not involve defining and using functions can be rendered this way, and the chains of arrows provide a nice visual indication of what one would call “the flow of control”.

Switching a kettle on is not typically a primitive operation that a computer can perform, so in real programs the box contents will be individual statements valid in the programming language concerned. If we are trying to find a really deep understanding of computation it is natural to wonder how small a collection of different sorts of statement and different sorts of test it is possible to get away with and still have scope for interesting behavior. Counter machines¹ provide one extreme version. As considered here a counter machine has a (small) number of variables each of which can hold a non-negative whole number, i.e. $0, 1, 2, \dots$. In due course we may think of those numbers as codes for text using any of the standard ways in which characters can be coded as numbers. The program we will develop will have its input data provided in its first register (which I will call A), and all the other registers start off holding zero.

Apart from a box that is labelled “stop” the only square action boxes that can be used have as their action statement that increments one of the registers. The only lozenge-shaped test-boxes that can be present check the value of one of the registers. If that value is zero they drop through to the next part of the flowgraph, otherwise they decrease the value in that register by 1 and go somewhere else. When the machine reaches its stop state the value in register A is considered to be the result it has calculated. Although this is of course just an integer, just as was the case with the input it can be interpreted as character data. To make this point as clear as possible, and computer file containing text is stored on disc or transmitted over a network as some sequence of bits, and one frequently used scheme transmits everything in 8-bit chunks, with a scheme that means that the commonly used characters (e.g. a-z, A-Z, 0-9 and various punctuation marks are fitted into one 8-bit unit (byte) while more exotic characters such as Greek α , β , π and the rest use two bytes and specialist symbols including many geometric shapes, lots of emoticons and pictures of the pieces for a chessboard use yet more bytes. One can then interpret this potentially rather long string of bits as the denotation of a binary number. In that way every file on your

¹There are a number of different names used for these primitive models of computation and a range of different sets of operations they can perform, but all the variations can be coaxed into modelling each of the others so the key results about them are robust. The version used here follows Minsky[?]

computer as a really natural interpretation as an integer and could be passed to a counter machine! Of course from a practical point of view this is totally absurd given that the only things we can do with numbers is to increment and decrement them. It would not take a very long input string to hit a situation where the number representing it was so large that counting it down to zero would involve more steps than there are atoms in the universe, and taking those steps would take more time than most people are prepared to wait. So this is to be viewed as a theoretician's model of computation. So the previous demand that you view computing time as an utter irrelevance really has pretty sharp teeth.

The big assertion to be made here is that for any computer program that can be provided with all its input at the start and deliver all its output when it stops, and subject only to the understanding that input and output will be encoded as big numbers, that it will be possible to devise a register machine implementation of whatever that program does. Is this going to be difficult? Well at some level yes it is – but once you have grasped how to attack the translation it is going to be less horrendous than it at first seemed. So the next few paragraphs show how the various key features in “ordinary” programming languages can be supported, Once that are in place transcribing the rest of the target program will be straightforward.

It is useful to start with better arithmetic then mere adding and subtracting one.

Here I will show the bits of code in a programming-language like notation because preparing flowchart diagrams would pain me. But for the final version many need to be drawn out, perhaps especially the early ones.

For addition consider setting up something that behaves like $A = B + C$ where A , B and C are registers and where D is spare one. Well in fact it hardly deserves to be described as “difficuly”.

```
while A!=0 do A--
while D!=0 do D--
while B!=0 do B--, A++, D++
while D!=0 do D--, B++
while C!=0 do C--, A++, D++
while D!=0 do D--, C++
```

This has risked destroying B and C along the way, so it carefully preserved their values in D and then restored them. If you were willing to leave them as zero things could be simplified a little. But the overall idea is that the counter machine can do something B times by counting down in B so we sum B and C by counting up in A first B times and then C .

The big magic that makes setting up a counter machine a lot less difficult than it might first have seemed is that after having convinced yourself that the above does perform addition you can write boxes in your flowcharts with $A = B + C$ in them alongside the primitive ones that say just $A = A + 1$. You work on from there producing additional calculations that you can use in boxes but then expand out into the primitives if you are really forced to. This is not really cheating – it is just like program building in an ordinary computer language where you set up a collection of subroutines (which you sometimes call functions or procedures) and then use them freely in the higher level parts of what you do.

In what follows I will often assume that a target register starts off at zero and that there is no need to preserve anything but the trick of initially counting down in A until it is zero and of saving values in D shown above can be applied wherever it is necessary. And I will also suppose that the number of registers that my machine has, while finite, is large enough that I always have a spare one available.

Then of course multiplication can be coded as just repeated addition, so $A = B * C$ will be

```
while B!=0 do B--, A = A + C
```

and with that it will be clear that raising to a power, being merely repeated multiplication, is also straightforward. Well if expanded out fully the flowcharts concerned may start to look untidy. But if one views each operation that gets implemented as a nice block of nodes that can be packaged and thought of and presented as a unit things are not so bad.

Subtraction involves a new issue because the numbers in a counter machine may never go negative. So the statement you first thought of as being $A = B - C$ needs to be handled more like “if $B < C$ then drop through not changing anything, otherwise set $A = B - C$ and take the other exit from the lozenge”. Given that it is easy to mechanise it by decreasing B and C in turn and noticing which one hits zero first. Then as necessary things are restored (using the “ D trick”) or A can be set. It should be pretty obvious that by using this that division can be coded up in a way that leaves both a quotient and a remainder.

It is perhaps useful to note that multiplication and division by known constants is rather easier. So for instance $A = 2 * A$ needs a single workspace register D but is then

```
while A!=0 do A--, D++
while D!=0 do D--, A++, A++
\end{verbatim}
```

```

and halving  $A$  if it is even amounts to
\begin{verbatim}
  if A!=0 then
    A--;
    if A!=0 then
      A--
      D++
      go back to start
    else
      A was odd
      while D!=0 do D--, A++, A++
      exit reporting A odd and unchanged
  else
    while D!=0 do D--, A++
    exit reporting that A has been halved.

```

Given the above that multiply and divide by 2, and the fairly obvious small variations on them that multiply and divide by 3, 5, ... one can in fact with a little bit of extra encoding of data get away with a counter machine that only has two registers, say A and D . If you had really wanted say three registers A , B and C you handle that by putting the value of $2^A 3^B 5^C$ in the main register of the two-register setup. What would have been increment or decrement operations on A , B and C now expand into multiplications of (test) divisions by 2, 3 and 5. By using more primes there you can model as many registers as you feel you need. To do this properly you need to convince yourself that with one register that contains real data and one to use as temporary workspace you can manage the multiplications and divisions by 2, 3 etc., but those operations really are simple enough that that is not a severe challenge.

fundamental steps inside them this does not even make things seem much worse: you can write your code with blocks that say $A = A + 1$, $B = B + 1$ and so on for as many variables as you need and each just denotes a mess of lower level messing with the two real registers you have. This scheme is completely general save for one caveat. That is that if your machine needs input, say the number K , it will have to have that encoded into its register as 2^K , and similarly when the machine stops its result will be an encoded version of the true answer.

So all is well and you can restrict yourself to using 2-register counter machines unless your resolution falters and you consider what it means in terms of the number of steps taken to perform some calculation. But it is proper to stress again that this is a game where you have agreed not to think about that!

It is now clear that simple integer arithmetic can be handled at least if you restrict yourself to positive values. The next thing to consider will be arrays, since they are a pretty frequent component of programs. Well in the same spirit that there was an explanation of how to encode a string of text as a number, here is a recipe for dealing with an array of N integers with values a_i for i from 0 to $N - 1$. Set up a string that starts with a 1 then has a_{N-1} zeros before another 1, then a_{N-2} zeros and so on down so that the string ends in a_0 zeros. So if the array was of length 3 and the values in it were 2, 4 and 6 we would set up 100000010000100. Now view this as the representation of a number in binary and view that number as an encoding of the state of the array. Well that step is simple - but it is now necessary to verify that the key operations of accessing the j th and updating it can be performed using a counter machine.

Actually those two operations are remarkably easy to arrange. First note that counting the number of trailing zeros in the binary representation of a number just amounts to finding out how many times 2 divides into it evenly. And division by 2 is one of the things we have seen that counter machines can do. To trim off a trailing 1 from an encoding you just need the transformation $N \rightarrow (N - 1)/2$ which is also straightforward. With those two operations in place it is then easy to trim $j - 1$ entries from a packed array, leaving the item at position j as the next to inspect, and it is then easy to read its value. To replace it all that is needed is to start by deleting j items and then put back the replacement followed by everything that had been removed. To put a new value k on the “front” of an array means just extending its binary representation by a 1 followed by k zeros. And that is $N \rightarrow 2^k(N + 1)$. Again the computation there is simple enough arithmetic that the counter machine can be set up to do it. The effect of all of this is that it becomes proper to write flowcharts with actions such as $A = B[C]$ that accesses the C th element of an array B .

The arrays set up as above could (of course?) be nested, and one way to cope with negative integers would be to represent a positive value N as an array of length 1 $[0, N]$ and a negative value $-N$ as $[1, N]$. All the basic arithmetic operations would now need to extract and check the sign marker but that is “just a bit more programming”. And since we are interested in showing how to make things difficult that can not be seen as a problem. Those who are serious massochists could look to the standard representation of floating point numbers as arrays of length 3 with one field for sign, an exponent represented by a number in the range 0..2027 and a 52 bit mantissa. All the basic arithmetic operations on floating point values amount to integer operations on the components of these triples.

Well with simple variables mapping onto counter machine registers and

arrays packed up as explained above and strings represented as arrays of characters, with the individual characters held as integer codes (as they would be anyway in languages like C and C++) it should be clear that the flowchart for any program that does not perform input or output operations while running but is just presented with some input at the start and just generates results as it stops can be expanded into a flowchart for a counter machine. What may be more amazing will be that each textual expansion along the way when making this transformation only increases things by a constant factor, so the new flowchart based on an almost ultimately primitive model of computing will only be a constant factor larger than the natural one you started with. And when writing out your flowchart you can use essentially all the operations from the instruction set of a traditional computer in the boxes.

What about function definitions and function calls? Well even they are not a disaster. If you review the way that arrays have been introduced here you will observe that they do not have any predefined limit to the number of entries in them. In fact the representation used behaves more like flexible lists than rigid arrays and slightly different ways of looking at them allows one to perform operations that amount to pushing a new item onto the front of a list and at some later time popping it off. Those operations are just what you need to provide a stack structure that keeps track of procedure calls. And what is going on there is really a fairly close relative on how compilers work when mapping procedures and their calls onto the hardware of real machines. Working through the full details here would become tedious but anybody who has followed this far should be able to sort it out for themselves if they really wanted to. The conclusion one ends up with is that counter machines can express pretty well any computation that an ordinary computer could be programmed to perform subject mainly to the constraint that all input data must be available from the start and no output should be inspected until the program terminates. For many purposes the limitation will not be severe.

A good joke about counter machines is that one of the earliest transistorized desktop electronic computer was built at a time when hardware to do complicated things such as addition and multiplication was seriously challenging, so internally it worked by counting, and it got as far as offering a square root operation to its users. So perhaps the ideas here have more practical impact than you might have expected.

Now some of the constructions shown here are unduly general and so it would be possible to model “real computers” in a more compact and possibly more efficient way. While we are only concerned with abstraction that is again not a big issue, but it does make one ask “Given a computation that is to be done, what is the most compact counter machine that will achieve it?” or in

other words how much better that the direct modelling shown here can we do? This is not just a difficult problem. It is one that in general can not be solved in any systematic way. It is natural to feel that if one has a counter machine with M nodes that solves a problem then if resource constraints are being ignored one can enumerate all the counter machine configurations with less than M nodes (start by cataloguing all directed graphs with that size limit) and just check which if any behave in the same way as the original. By scanning from smallest upwards one would naturally come across the best. The painful reality here is that given a counter machine (or indeed any other sufficiently general model of computation) there can be no algorithm that will in general certify that example setups will have terminating patterns of computation. A consequence of that is that it can not be possible to guarantee to be able to verify that two counter-machine configurations have the same behaviour. However it will be possible to spot some cases that agree, and to identify others that do not. So for a *really* difficult problem design software that does the best you can to take the description of a counter machine and optimise it.

There is another famous challenge-problem associated with counter machines, Consider all possible machines with M nodes (and K registers). Consider their behaviours when started with all their registers zero. Among all of those which one halts and when it does has the largest possible value in its first register. Obviously there can be machines that never halt - perhaps the simplest version just increments a register and returns to its starting state, so it sits there counting up for ever. Such cases are to be discarded - it is just machines that actually terminate that are of concern.

With a truly tiny number of nodes this question is easy to answer. A machine with one node plus a stopping one can not do better than to make its first node increment its register and transfer to the halt node. So the answer there is 1. An alternative version of the challenge is to maximise the number of steps taken before the halt state is reached - both versions are remarkably tough. It is plausible that a register machine with just 2 registers and only 3 nodes (plus the stopping one) can compute for much longer than you would have expected before terminating, and I here leave that as a topic for enthusiasts to explore.

Hmmm - google AI summary for "busiest minsky machine" seems to suggest that a 2-register 3-state machine might take thousands of steps and then stop. I have not spent time working out what design behaves that way and the AI summary feels slightly incomplete and in particular does not point me at "proper" papers or reports. So I hope that somebody better at web search than me can find proper detailed documentation!

[https://codegolf.stackexchange.com/questions/279153/
long-running-section-11-4-minsky-machine](https://codegolf.stackexchange.com/questions/279153/long-running-section-11-4-minsky-machine)

‘‘Rick J. Griffiths in his dissertation in 2003 has adapted the busy beaver problem to Minsky’s “register machines”. His program is written in Java.’’

and that is a Cambridge Part II dissertation and I may be able to find a copy in the Computer Laboratory archives.

Chapter 4

Turing Machines

There have been two or three big reasons for looking for minimal schemes that are able – in some sense – to compute.

1. If you want to build a computer it is very reasonable to try to make things easy for yourself by designing the simplest possible device that you can. Even though that might make life impressively harder for those who want to write programs for it. There are modest applications of this idea that have led to very successful commercial designs the successors of which are in widespread use today – but the extremist version of it may provide a great basis for a fun practical construction project. The assertion “I have made by own computer from scratch” is obviously a good one to be able to make;
2. If you are serious about wanting to develop a robust theory that *really* lets you understand what computers can do and what they can not it is rational to start with something simple. What you will be trying to do will be difficult enough without needing to worry whether all the special capabilities built into modern computers and programming languages make big differences;
3. Those who are concerned with how long it should take to solve some problem will find there are huge and shifting complications if they consider hours, minutes and seconds on a range of desktop and laptop machines as well as mobile phones and embedded controllers. By looking at timings on truly reduced hardware their results will be less of immediate relevance but can be ones that will remain valid as this year’s computers are replaced by next years ones that are possibly from a quite different manufacturer.

The best known minimalist sort of computer is the Turing Machine. This emphasises the fact that a computer of any sort will have memory and if viewed from a distance all that it does can be seen as taking steps that each inspect and change a single item within that memory. In focussing on how data is stored it does not take any steps to make it particularly easy to coax it into solving the problem that interests you.

The storage provided by a Turing Machine is a tape which is marked out in cells. Each cell can hold one of a modest number of symbols. The number of symbols allowed in one of the parameters that describe exactly what sort of Turing machine is being considered. The tape is made long enough for whatever task you are happening to try to process. Some people would characterise that as a tape that is infinite in length, or would use the word “unbounded”, but in any particular computation that the Turing Machine takes only a finite segment of it will be used. So for practical experimentation it can be acceptable to provide a limited length “tape” and view an attempt to go beyond its end as merely reaching a limitation of the physical approximation to the abstract machine.

The Turing machine starts with whatever input data it needs ready on the tape. It then works step by step: it has a read/write head positioned over some cell of the tape and a cycle it takes will inspect the symbol there and based on an internal state (from a limited number) it will write back a possibly changed symbol and move the tape left or right. It also transfers into some internal state. One state will be special in that entering it causes the machine to halt. At that stage it is expected that it will be written its result onto the tape. The number of distinct states that the machine can be in is the second parameter alongside the size of the alphabet of symbols on the tape that characterise it.

Programming a Turing machine has to involve setting up a table that is indexed by which symbol has just been read and which state the machine is in. When that information is used to inspect a row in the table you can read off the symbol to be written, the tape movement to apply and the identify of the next state that the machine should be in. A reasonable expectation is that designing tables like that to perform even modestly elaborate computations will be a bit painful.

One way to prove that it is really worthwhile setting up this fairly clumsy looking model of computation is a proof by construction that it makes it feasible to build a mechanism that follows its behaviour pattern. In LEGO!

[https://beta.ideas.lego.com/product-ideas/
10a3239f-4562-4d23-ba8e-f4fc94eef5c7](https://beta.ideas.lego.com/product-ideas/10a3239f-4562-4d23-ba8e-f4fc94eef5c7)

is a concrete realisation of this using under 3000 LEGO parts. While that is quite a lot, it can be put in perspective by comparing with the official LEGO kit to make a model of the Star Wars Death Star, which comes with 9023 pieces – but rather fewer gearwheels. And an astounding thing is that the only limitation at all on what this Lego Turing machine could do is set by the length of its “tape” and at least in imagination it would be easy to make that really long.

When considering building a real working Turing Machine it is reasonable to ask just how much mechanism it needs to have before it can actually perform any useful calculations. The LEGO version has 8 states and handles an alphabet of 4 symbols on its tape and that feels as if it might be quite limiting, however the astounding thing is that if you had a long enough tape (and seeing how to make the tape a bit longer is surely not a terribly tricky technical challenge) this is enough to allow the machine to perform **any** computation that any other computer can. To be more specific it will be possible to set up initial contents on the tape where the first part amounts to “program” that documents what is to be done and the rest is the “input data”. A very special case of this is that the program part can explain how to behave as if the rest of the tape is being used by a Turing machine with a larger number of states and a bigger alphabet. For instance if one wanted to have 8-bit characters on the tape the machine that would be emulated would systematically use the contents of four consecutive cells on the physical tape to represent a single byte on the bigger system. Designing, understanding and using Turing machines that are really close to being minimal can be huge fun however they can sometimes demand really ugly ways of encoding the data on their tape or suffer from needing unreasonably large numbers of steps to reach results.

A variant using just 2 states and 3 symbols exists and is characterised as “weakly universal”. It will not even halt when it has completed its work and it also needs its tape initialised to a particular pattern in all regions beyond the input data, while standard machines will not read from the tape beyond the region that is explicitly data - it also tends to take a great many steps to get anywhere, while modestly larger machines are actually quite efficient.

Once one has demonstrated that a basic and fairly small Turing Machine can be used to emulate a machine with more states or a larger alphabet it can be reasonable for subsequent work to feel free to relax those constraints to make machine design simpler. and in particular at least from a theoretical perspective it is acceptable to imagine that the “state” part of the machine can do anything that an “ordinary computer” with no unbounded memory can. The tape is then just needed to provide the unbounded storage that it is good to have when analysing algorithms. To slightly simplify the proper and

general result, if the “ordinary computer” completes a run within N steps it can not possibly have touched more than N distinct memory locations. If its memory is modelled by data on the tape then the most remote bit of data ever accessed can be no more than about N cells away. Well we may aggregate raw tape cells so that some symbols are stored spanning across say K of them, but then the furthest accessed data is only KN away, and the Turing machine should only take time proportional to that to access it. So we find that each of the N steps of the ordinary machine gets emulated in time bounded by something of the form KN^2 . This quadratic overhead would of course be calamitous in practice, but is modest enough for a great many theoretical studies.

Rather simple extensions and generalisations to Turing Machines allow for yet better efficiency for many problems. One can consider a device with more than one tape or with just one tape but more than one read/write head. As a concrete example of how this makes things easier, Merge Sort was a solid solution to sorting vast amounts of data when computer memory was small and the data has to live on magnetic tapes. The treatment of those tapes and those in a multi-tape Turing machine are closely analogous to one another, so that sort of TM provides a really solid model for analysing that sort of algorithm while avoiding the need to worry about detailed characteristics of any physical computer.

The overall message here is that if you really want to make things hard for yourself try to design the most compact Turing Machine that can do the calculation you are interested in, seeing how you can make trade-offs between the number of states, the size of the alphabet you allow on the tape, the ugliness of how data has to be coded for use and just how many computational steps will be taken to obtain your solution. That is a wild and confusing space to search within. A variation on this is just to look for Turing machines that run for a seriously long time but then stop when they are started on an empty tape. For really small machines bounds are known. For instance if the tape can only hold one of two symbols in each cell and the machine has 2, 3, 4 or 5 states the number of steps it might take before termination can be 6, 21, 107 and 47176870. With more states the number of steps becomes infeasible to write using commonly-used notations! These results show rather clearly that very small systems can have extraordinarily complex behaviour and as systems as small as these can behave in such extraordinary ways the detailed understanding of larger and less artificial ones will be difficult in the extreme.

To finish this chapter it is proper to mention another variation on these machines that has enough witty consequences that it will form the basis of a whole separate chapter. Ordinary Turing machines are very much mechanical

and deterministic devices that always behave in unambiguous and predictable ways. An entertaining variation will be the class of machines where the state transition is not quite deterministic. In some cases the machine will be able to choose for itself from two different next states, with no pre-programmed or external guidance. If the choice was made by notionally tossing a coin in each such case you would have a randomised machine, and exploring the capabilities there would be entertaining. But the most important case here is where the decision between the alternate paths is made as if some good fairy waves a magic wand and the computation proceeds such that if your calculation could at all possibly succeed it now will. This is obviously a delightful fantasy, and the resulting model of computation is referred to as a “non-deterministic Turing Machine”. It seems obvious that machines like this could never exist in reality – but in fact if you abandon all concern for timings they could be emulated, and to date nobody has been able to prove that there is no way of building a rather efficient simulation of one.

Chapter 5

Galactic algorithms

Galactic algorithms are ones that have costs that grow slowly with the size of their input to an extent that for large enough cases they will do better than simpler methods, but that suffer from such large overheads that they will never become practically useful for any plausible cases. The finest ones will provably achieve the slowest possible growth rate for the problem that they solve. Some of the most impressive have overheads that can be expressed as a constant factor multiplier in their cost formula with that constant exceeding astronomical proportions. There will obviously be cases that come close to fitting into this category. Because such methods are not of a lot of practical use one could be tempted to view them as frivolous, but where they represent a better growth rate than any other known method or are provably optimal in that respect they can both be of intellectual interest and can provide ideas to guide work towards schemes that can be useful in the real world.

A number of galactic algorithms illustrate how much more complicated things get if you aim for the absolutely best possible method, but here we start off explaining one that has been known for some time and is neither terribly complicated nor terribly expensive, but that despite it giving a theoretical speedup will only achieve that in cases pushing against the size limits of the biggest available computers, and even then its advantage will be slight.

5.1 Matrix Multiplication

The straightforward scheme for multiplying a pair of square matrices of size $n \times n$ has cost that grows proportional to n^3 . Our first galactic algorithm represented a breakthrough discovery that this growth rate was not the best possible, and that a tolerably straightforward scheme could achieve a cost that grows as $n^{\log_2 7}$ which is about $n^{2.80735}$. To show how this is done it is

sufficient to consider the very simple case of forming the product of a pair of 2×2 matrices A and B . The four elements of their product as evaluated in the classical manner are just $A_{11}B_{11} + A_{12}B_{21}$ and three rather similar looking values and a total of 8 multiplications and 4 additions are used. A paper by Strassen[?] instead managed to show that using just 7 multiplications that worked on carefully chosen sums and differences (e.g. $(A_{11} + A_{22})(B_{11} + B_{22})$) it was possible to express all the values needed in the product matrices as sums and differences involving these 7 values. Here we leave finding exactly how to do that as either a puzzle to work on or a topic to research! A variety of different schemes have been found, all ending up with just 7 multiplications but some with more and some with fewer additions.

Once the base case of 2×2 matrices has been handled large case can be treated by partitioning each matrix into four blocks (padding out with zeros if necessary to make those blocks all square and all the same size) and coping there by using just 7 multiplications (which are now matrix ones) on those half size blocks. An analysis of the growth rate in time spent using this scheme fairly clearly suggests that doubling the side of the matrices multiplies the amount of work by a factor of 7, hence the $n^{\log_2 7}$ result. This is clever and quite elegant, and not even too hard to program. But as against the classical method it struggles even to break even until matrices are larger than almost all applications, and it certainly does not deliver truly valuable speedup for some way beyond that. A further concern with this approach is its impact on how rounding errors propagate through the calculation.

But this is not the end. Since the Strassen work others have sought to find schemes that are at least asymptotically faster, with a long sequence of reports of gradual reduction of the exponent in the cost function. As of 2026 the best growth rate that has been reported is around $n^{2.37134}$ which looks like a distinct improvement over Strassen. So far nobody has been able to show what the ultimate limit for solving this problem will be. Most of these improved schemes work by splitting large matrices into more than 4 blocks and all of them are even further from practicality than the Strassen method.

5.2 Integer multiplication

Multiplying integers is a pretty fundamental operation, but here to make things harder the emphasis will be on large integers – typically ones with from hundreds to millions of decimal digits. Things start off in a style almost parallel to the matrix multiplication one, with first the simple “schoolbook” method that will form the product of a pair of N digit numbers using around

N^2 operations. The first improvement on this is due to Karatsuba[?] and works by observing that for 2-digit numbers one can form the partial products a_1b_1 , $a_1b_2 + a_2b_1$ and a_2b_2 using 3 rather than the obvious 4 multiplications by calculating a_1b_1 , $(a_1 + a_2) \times (b_1 + b_2)$ and a_2b_2 and doing a couple of subtractions. By taking large numbers are splitting each in half this can be applied recursively leading to a method where costs grow like $N^{\log_3 2}$. This scheme can become proper to use when numbers exceed say 10^{300} , but the exact break even point will be quite sensitive to details of the computer used and just how everything was coded. So this not a galactic algorithm.

The next scheme of note may start to be competitive when multiplying numbers of magnitude around 10^{5000} . The exact point where it becomes realistic can vary quite a lot and 5000 digits is perhaps better characterised as stellar rather than galactic, so this is a method that is at least relevant in some slightly specialist cases. It is however much harder to explain in elementary terms, and so what will be presented here is an overview perhaps uses enough technical language to show that it starts to count as “difficult” or at the very least “reasonably advanced”. So for those who do not recognise the words used here this will just show how much messy technology a simple-seeming task such as integer multiplication can involve. For those who then want to follow through by reading textbooks or scouring the web it may give a broad overview so they know how the things they read up on fit together. And those who already understand this well can just smugly skip to the next section.

Schönhage-Strassen works by first padding its integer inputs so that each have a suitable power of two bits. At least as an abstract algorithm it expresses its operations in terms of bits not digits so it is kept fully honest and does not hide costs within arithmetic on digit-sized units. Then if the numbers are of length N it clumps the bits into clumps of size around \sqrt{N} and views each clump as a digit, and so there are around \sqrt{N} digits. Well there can be some pedantry about the exact size of digits that is needed but the decomposition as described here is close enough for use in this overview.

The next issue is that it views each collection of digits as a vector and it takes a Fourier Transform of it. So what is a Fourier Transform? Well it amounts to multiplying a vector of length K by a special form of matrix where the entries in the matrix are all powers of a value ω where $\omega^K = 1$ but no lower power of ω has that value. In the normal world one would say that ω is a complex K th root of unity, and that it is primitive, so at first it looks as if complex numbers are involved. However to avoid them (and any risk of creeping rounding error that could hurt if floating point arithmetic was used), Schönhage and Strassen used modular arithmetic. In other words they have some number P so that after any operation on integers they just retain

the remainder when the natural result is divided by P . A big part of their cleverness is that by choosing $P = 2^M + 1$ for a suitable M two wonderful things can be achieved. First it becomes possible to use 2 as the value for ω and that means that multiplication by powers of ω can be mechanised as simple cheap shift operations. Secondly the remaindering operation where a double length integer is to be reduced modulo P can be performed by just subtracting the top half of the number from its bottom half.

One interpretation of what a Fourier Transform does is that it views its input vector as the coefficients in a polynomial and evaluates that polynomial at all the powers of ω . If that is done to two polynomials and the resulting values are pairwise multiplied the result will be the values of the product polynomial at all those points. An inverse Fourier Transformation will act as interpolation and recovers the coefficients of the product polynomial. This can be re-interpreted as an integer following some simple carry operations. A further key feature of Fourier Transforms is that on using vectors of size K can be completed using around $K \log K$ arithmetic operations. Of course here each of those arithmetic operations is using modular arithmetic modulo P where P is a pretty huge multi-digit number – but because of all the powers of 2 involved that is not in fact seriously painful. In fact the worst part of all of this is the pairwise multiplication of elements of the transformed vectors and that generally needs to recurse into a further layer of use of the whole procedure – at least until things get small enough that classical methods will win.

The “polynomial product” produced by the Fourier Transform scheme has been computed modulo P but we want exact and correct integer results. Well provided P was greater than any value that could legitimately arise in that product all is well and the calculated values will be the perfect integer results as needed.

The above explanation know it is a sketch. It has omitted the need to pad various things up to powers of 2 and the proper care about splitting the inputs up so that the value of P used is big enough to ensure that results are correct. Those details of course matter in a proper formal explanation of the algorithm and in any implementation of it, but perhaps they represent details best deferred to a second reading about the method.

The Schönhage-Strassen scheme for multiplying N bit numbers has a cost that grows proportional to $N \log N \log \log N$. Because it is concerned with truly gigantic input numbers (i.e. values of N) it becomes proper to worry that if digits making of numbers are stored in computer arrays of length N then index arithmetic used to work out which digit to touch next can become beyond the scope of single computer operations. For instance a modern 64-bit computer will not cope trivially with arrays with more than

2^{64} elements! So the proper analysis of costs here needs to be done in terms of some suitable abstract machine that does not have any limits at all on the bulk of data it can store and manipulate. Variations on Turing machines are commonly used, and amazingly some of them can deliver the cost growth rate noted above.

This scheme, however messy, is still used in some practical applications. But it is not the end of the road. A succession of authors have described yet more elaborate schemes that reduce the $\log \log N$ term in the cost function, culminating one[?] where the cost is precisely proportional to $B \log N$. This is a great theoretical result in that it is conjectured (but not yet proved) that this is the best growth rate that can be reached. The technology applied is significantly higher powered than the scheme covered above and so there is not even an attempt to explain it here, but those who really enjoy difficulty can read the 45-page paper that introduced it! One of the authors say that for this method to be cost competitive the numbers involved would need to be rather large: “Even if each digit was written on a hydrogen atom, there would not be nearly enough room available in the observable universe to write them down.” That surely qualifies it to count as galactic.

5.3 Regular expressions

The separate chapter here on pattern matching also presents a problem that might have seemed tame but is in fact galactic.

5.4 Primality checking

It is easy to check if a small number N is prime – just check each potential factor up to \sqrt{N} . For larger numbers there are fine and practically very sensible methods some of which rely on a supply of (genuine) random numbers to deliver a result with arbitrarily small probability of error, or that rely on currently unproven results in number theory. But a scheme introduced by Agrawal, Kayal and Saxena[?] runs in time bounded by a polynomial in the number of digits it takes to express the number and guarantees a correct response without needing to make any questionable assumptions. This is difficult in two senses! The mathematics behind explaining and justifying it are somewhat tough, and if it was implemented the suggestion is that its cost would be significantly slower than using probabilistic methods for all numbers with no more than 10^{1000} digits. It is so laborious to use that even for numbers of a few thousand bits (as used in many cryptographic contexts)

is is not really feasible at all as well as it being hugely slower than other methods. So here we can report that this exists, that it is a fine example of the levels of difficulty you get when trying to get the very best solution to a problem, but that those who want full details may first need to attend courses on number theory!

5.5 Minimum Spanning Trees

A graph here is a collection of vertices with edges that join some of them. A spanning sub-tree is a subset of the edges such that it is possible to get from any vertex to any other one only traversing edges in that subgraph. A graph may have weights (or costs) associated with each edge, and then a minimum spanning tree is a spanning sub-tree such that the sum of the weights on all its edges is as small as possible. Often in university courses on datastructures and algorithms procedures by Prim and Kruskal[?] that find such minimum trees are presented. In the more advanced versions of such courses improvements on the basic versions of those using forms of priority queues known as Fibonacci Heaps[?] get discussed. If the graph has m edges and n vertices these schemes can guarantee to find a minimum spanning tree within time proportional to $m + n \log n$ and that is typically good enough for practical purposes. However a fancier method by Chazelle[?] has a growth rate that scales as $m\alpha(m, n)$ where α is an inverse Ackerman function and grows exceptionally slowly such that for any conceivably feasible practical graph its value will be no greater than 4. So until we are really at or beyond galactic scale the costs of this method grow linearly with the number of edges regardless of the number of vertices. As with the previous examples in this chapter this represents a breakthrough such that for large enough graphs the method will beat all the ones that can be developed before. As with other examples in this chapter the details of the extreme method defy compact explanation, but the overview is that they replace the Fibonacci heaps (which are themselves somewhat messy!) with data structures known as “soft heaps” that mostly provide priority queue operations in a really cheap way but that do not always yield the correct result. However they do offer bounds on the frequency with which errors can be introduced, and the Chazelle method as a while allows for that by making checks as it goes and where necessary back-tracking to correct for mistakes. The presence of the inverse Ackermann function in the eventual cost bound gives clear warning that this is all far from straightforward even though right here we are not documenting just what that function is or how it arises!

5.6 Perfect data compression

When data of any sort is to be transmitted over a slow network or stored in an archive it may well be good to produce a reduced size version of it such that when the original data is eventually required it can be reconstructed. The case considered here is where the original data must be recovered perfectly – that is in contrast to data compression often used with images (.jpg) and sound (.mp3) where ending up with a version that is in some sense “close enough” to the original will suffice. There are many commonly used schemes for compressing data and a number of them use the word **zip** in their names. But none of those even pretend to achieve the best possible reduction in size.

There is a scheme that is in a sense optimal, but the big problem with it is that it is even worse than the previous galactic algorithms just considered, in that it can be shown that it is and will always be impossible to write a program to implement it!

This scheme is known as Kolmogorov compression and the idea behind it is really simple. A body of data is represented by the most concise program that, when run, will regenerate it. Well that definition raises an obvious issue: “what notation should this program be expressed in?”. From a theoretician’s perspective the response is that any notation at all may be used. If then one wanted a version of the data to be decoded on a different computer or one that provided implementations of different languages all you do is construct code in the notation that will suit this different computer that will emulate the one the data was compressed by. This simulation code will be of some definite finite size and so sticking it on the front of your compressed data will only alter its simply-measured size by some additive constant.

The unsolvable problem that makes this scheme one only to dream of is that there is no way to guarantee to find a shortest program to generate some given output. This is a consequence of the fact that given a program there is no algorithm that can guarantee even to tell if it will terminate in finite time, far less give information about its output. But this is only a practical limitation! One might expect that something that was impossible to implement would not have many uses, but amazingly that is not the case. Here are two that illustrate the power of this concept:

5.6.1 Testing for randomness

A fundamental result is that if data is truly random then the most concise way to express it is basically to put it in a simple print statement. For instance this shows that a string of digits such as those from 10000 to 1000000 within the decimal expansion of π are not really random, because

this very sentence can be viewed as a recipe from which they could be reconstructed. And if desired the sentence could be preceded by a program for calculating digits of π . So the fact that the string of digits that starts off 85667227966198867... which may have no pattern discernable to the casual eye could be spotted by Kolmogorov compression as being extracted from the list of digits of π .

So suppose you have a sequence of values that are supposed to be truly random – such as numbers drawn each week in some lottery – the thought here is that if somehow you could find a scheme that could generate exactly that sequence where the description of your scheme is notably shorter than merely listing the numbers themselves then you have good reason to believe that they were not random after all. At least in some sense all the techniques used to try to assess procedures that generate pseudo-random numbers are modelled on this in that they look for patterns that represent traces left by the mechanisms actually used to generate the numbers. But Kolmogorov compression would uncover the implementation of the generating program and in so doing show that the sequences were pseudo-random rather than genuine.

5.6.2 How common are primes?

The result sketched here uses the fact that almost all data must be impossible to compress. This result emerges from counting how many different possible byte-sequences of length N exist and concluding that any loss-free compression can only possibly map that number of longer strings onto them. So all the other longer strings can not be squashed. This line of reasoning does not indicate exactly just which strings will end up being mapped onto shorter ones and which will not! But good practical compressions methods try to arrange that the ones that are allocated short representations are ones more liable to arise in practise.

Using this principle that most data is not compressible it is possible to derive a result that shows that primes must be reasonably common. This result is not going to be anything like as good as the ones that “proper number theory” comes up with, but the fact that it uses data compression as its main tool is perhaps witty!

So consider a really huge number that is written using N digits. The principle that most data can not be compressed says that in general it will not be possible to find a way to describe it using significantly fewer than N digits. Well one alternate way of specifying the number would be to list its prime factors, so for instance the integer 856672 would be presented as $2 \times 2 \times 2 \times 2 \times 2 \times 19 \times 1409$. and now we imagine numbering the primes, so

2 is the first, 3 is the second and so on, Then this representation of 856672 might be written as (1, 1, 1, 1, 1, 8, 223) because 19 is the 8th prime and 1409 turns out to be the 223rd.

Now if you imagine that primes are really very uncommon then the index of a prime (eg 233) will be very much less than the value of that prime (eg 1409) and so writing out the index will use less space. That could potentially give a way of finding a way to create a compressed representation of the number, and while that can happen occasionally that has to be rare. This establishes that primes can not be hugely rare.

Rather than attempting to get the sharpest possible result out of this, pretend for a moment that the density of primes was such that the n th prime had value comparable with n^2 which would mean that its simply written value used about twice as many digits as were needed for writing out its index. Then at least a simplistic reading is that the compressed representation of a number a list of the indices of its factors will use half the number of digits plus a number of commas based on how many factors are present. for big enough numbers this will pretty well always be significantly shorter than the representation that merely writes out the number simply, and this being impossible establishes that primes must be more common than had been assumed.

This illustrates that even without being able to exhibit perfect compression a consideration of it can be used as a basis for reasoning about probability distributions and the like. But it remains the fact that optimal compression is not merely galactic it is uncomputable.

Chapter 6

Intractable problems

The chapter about galactic algorithms looked at extraordinarily laborious ways of solving problems, but various of the problems considered could be resolved in sensible amounts of time in practical-sized cases using an alternative method which would lose out to the galactic algorithm in the long run, but only the very very very long run. This chapter considers problems where in some cases all we can say is that the current best solution known is obnoxiously time consuming and in others where it is possible to prove that an algorithm, including all the ones not yet invented, will be infeasibly slow.

Chapter 7

Practical Performance

Galactical algorithms is all about pushing theory and asymptotic performance estimates to the limit regardless of whether the result would be practical. So a scheme that ran in $10^{100}N$ when faced with a problem of size N would be preferred to one that needed $0.001N^{1.0001}$ because for large enough N it would be the winner. This chapter considers the other end of a spectrum and considers cases where absolute timings measured in minutes and seconds or space use in bytes is to be optimized, but in the pursuit of perfection all other constraints and limitations are ignored. This can certainly involve use of significantly messy and elaborate code to achieve even small improvements over straightforward implementation.

7.1 Abandoning portability

7.2 Special hardware support

7.3 Composite methods

7.4 Dynamic code generation

Chapter 8

Lambda calculus and Combinators

Turing machines provide a model for computation that (perhaps unexpectedly) can give really good insight into costs. But setting up programs for them is not really anything like conventional programming and they do not adapt well to day to day use. Minsky or Counter machines have their behaviour expressed by flowcharts and come much closer to being amenable to casual use, but they exhibit totally absurd run-times that render them quite unsuitable for anything other than theoretical purposes. That leaves a gap in the market to be filled by a formal model of computation that can be reasonably easily adapted for use as a practical programming system. A key candidate for this is called “lambda calculus”.

The fundamental idea here is that functions should be seen as the basic building block for everything else. When first introduced this was accepted as a neat theoretical idea and it led to what I will impolitely describe as niche programming languages, but over recent years it has been recognised as such a good idea that basically all modern programming languages incorporate features that are directly derived from it. A *lambda expression* denotes a function and is seen as a freestanding entity. For instance in earlier usage one could only introduce a function by a notation that gave it a name, as in

$$f(x) = 3*x + 1$$

as a definition of a function called **f**. With lambda calculus the name **f** is not needed and we can write just this function as

$$\lambda x . 3*x + 1$$

where the name just after the λ indicates what the formal argument is and the section after the dot is the body of the function. About the only action that can apply to a lambda expression is that it can be applied to an argument,

and the consequence of that is just about the same as when the previous function called f is applied to an argument. It is almost as simple as that!

Well there are two issues that need discussion that arise when several lambda expressions arise together. The first rule is just to give a pedantic confirmation that things happen the way you would naturally hope, and it insists that when the body of a function has a value for the function's formal parameter substituted in nothing improper happens. One version of it starts by noting that in any lambda expression one could rename the variable that is bound to anything else and so the example shown above is to be viewed as entirely equivalent to

$$\lambda y . 3*y + 1$$

and the mapping from one text string to the other is known as alpha conversion. Now if you applied out first lambda expression to an argument that was the literal symbol x rather than a number there would be a potential source of confusion about substituting for x in the body. All possibility of uncertainty can be overcome if one used alpha conversion on the function just before applying it to make its formal parameter a new symbol that could not conflict with anything. Another part of what is essentially the same issue arises when the body of a lambda expression includes further nested lambdas and variable names are reused. It is proper to apply scoping rules so that inner lambda bindings take priority over outer ones, but even here use of enthusiastic alpha conversion can keep all names separate and avoid risk.

The second issue is that if one has a reasonably large nest of lambdas (and some examples will arise later) it could be that there are several places in the whole where a lambda could be applied to an argument. When there are several options for the order to do these operations does it matter which is selected and if so what is the proper strategy? On this there is some divergence of opinion! From a theoretical point of view it is considered proper that where there is a choice one should select the leftmost outermost application. This is known as normal order reduction and if there is any chance at all of continued application of lambda expressions to arguments ever terminating with a form where there are not more available than this guarantees to reach that state. The alternative is that innermost lambdas are processed first. This can sometimes risk unnecessary exploding or looping sequences of operations, but is generally seen as easier and cheaper to implement. So many but not all programming languages with lambda support follow the second path.

It is perfectly feasible to write out everything that is to be discussed in this chapter using loads of λ symbols, But for many people it will be easier to work with a notation that feels more familiar (but that can be expanded directly into the raw lambdas). A paper by P. J. Landin[?] has as its title

“The Next 700 Programming Languages” and introduced ISWIM (If You See What I Mean) as a sketch, and so rather than latching onto any particular real or modern notation that exposition used here will be based on that.

ISWIM starts by using a keyword **where** to give names to things. To be a little pedantic here are a couple of examples and how that would expand onto raw lambda expressions:

```
3*x+2  where x = a + b      : (λ x . 3*x+2) (a + b)
f(a+b)  where f(x) = 3*x + 2 : (λ f . f(a + b)) (λ x . 3*x+2)
```

The hope here is obviously that the ISWIM versions are un-threatening and so will make things easy to understand. As well as **where** that puts a definition after the use of something it will also be possible to use **let** to put a definition first, and here we will often use that with the intent that the definition provided should be available to all subsequent input. What is more likely to cause some concern to start with is that in this world functions are first class citizens with all the rights and privileges entailed. In particular they can be passed as arguments and returned as values. A major lesson that lambda-calculus teaches is that this is liberating and powerful and so it is used a very great deal.

The ISWIM examples above use function definitions and **where** clauses to provide the structure of a program, and then have numbers and arithmetic operations. As a prototype for 700 variant languages ISWIM can imagine a wide variety of built-in data types and functions that can apply to them, so that it can be tuned for particular application areas. I will allow for the creation of user-specified data types. But it tends to de-emphasise sequential programming, with its documentation noting that in Algol 60 it was already possible to avoid that quite a lot by virtue of allowing conditionals as expressions as well as statements. It can view a sequence of several values as in **A**, **B**, **C** to be viewed as a single item, and that item could be returned as the value of a conditional expression or passed to provide three arguments for a function that needed that many. The variant or variants of it used here will be ones tuned to be good for describing things that can be done with lambda expressions!

From the tiny examples shown above it is possible to push explanations either up or down in terms of the level of abstraction involved. Here we will first show that lambda calculus using nothing more than function definitions and even restricting those to the one-argument case is in fact capable of modelling all computation. And that amazingly the steps it needs to take to do this are all remarkably concise and may even not be terribly inefficient. After that it will be explained how lambda expressions can act as a convenient and well defined intermediate stage to be used by compilers while mapping

from source code down to some machine-specific level. And When that has been done it is easy to observe that any programs in the languages that those compilers support can be rendered as lambda expressions. A final section will comment on a somewhat archaic but thoroughly practical language that was explicitly built to illustrate this but that has been used to build large and complicated programs.

To build up an understanding of computation from lambda calculus we start by noting that a lambda expression only has a single argument. In much of what follows it is going to be convenient to think in terms of functions with two or more arguments. These can be rendered using a technique called “currying” after H. B. Curry[?] who popularised it. We will write function definitions as if they used several arguments as in

```
let f a b c = a+b+c+2
```

using spaces between the multiple parameters rather than enclosing the collection of them in parentheses. We write in parentheses to show how we intend or expect to view this as grouped. In this case both the description of the pattern for the function (i.e. the bit to the left of the =) and the additions in the body associate to the left. Then there is a bit of rather dubious renaming and a use of (where) and the name F that is really not quite proper but that turns the main definition into one with a single argument which is then easy to transform into a lambda.

```
let ((f a) b) c = ((a+b)+c)+2
⇒
let F c = a+b+c+2
  where F = (f a) b
⇒
let (f a) b = λ c . a+b+c+2
⇒
let f = λ a . λ b . λ c . a+c+c+2
```

Even if the last line might look slightly awkward to start with it shows that the ISWIM-style notation that defines a function that in a certain sense has three arguments does correspond to respectable lambda calculus.

Given that the first thing to be done is to show how easy it is to provide truth values (i.e. *true* and *false*) and a conditional that tests them. And instantly from that **and**, **or** and **not**:

```
let true a b = a
let false a b = b
let if p X Y = p X Y
```

```

let and a b = a b false
let or a b = a true b
let not a = a false true

```

These can also look a bit strange to start with, but observe how very concise they all are and then just verify them by writing down combinations of them and using the function definitions as rules to check how things behave. However there is one big caveat in all this – it only works nicely provided you use normal order reduction, because that manages to arrange that for instance when you write `if p X Y` that the evaluation of `X` and `Y` gets deferred until it can be determined that only one actually needs to be processed. With the evaluation strategy that deals with innermost expressions first both `X` and `Y` would get expanded there and often that would lead to a loop or other disaster.

Note that for the above to behave the value `p` tested by `if` must evaluate to either `true` or `false`. In a bit there will be a discussion about how that might be enforced.

Given truth values and conditionals the next feature of “real programming” to consider will be data structures. A useful building block from which pretty well anything else can be constructed will be the “ordered pair”, where we will have a function `pair` that creates one and then `left` and `right` that retrieve one or other of the two included items. Again the lambda forms that implement this are amazingly concise and you just have to try following through what happens when you use them to convince yourself that for instance `left (pair A B)` will recover `A`:

```

let pair A B f = f A B
let left P = P true
let right P = P false

```

There are two things that can be done instantly once you have pairs. The first is to use clusters of them to build triples, 4-tuples and larger structures with any fixed number of components. The second is to pair an object with a tag (which might be a truth value) so that as a while it can have two (or more) possible shapes. As an illustration of this consider a way of handling lists that can have any number of elements (including zero):

```

let emptylist = pair false <anything>
let node A B = pair true (pair A B)
let isempty L = not left P
let first L = if (isempty L) <error case> (left (right L))
let rest L = if (isempty L) <error case> (right (right L))

```

which exposes the fact that dynamic structures like this have to carry extra information so that it is possible to detect when the end of a list is reached. A similar scheme could describe trees. The sections in the above that are written in angle brackets are parts that ought never to be used so anybody confident that their code would not contain any errors could use absolutely anything to fill those gaps.

Next comes a representation of numbers and arithmetic. Well one scheme that could handle positive integers tolerably well would be to express them all in binary and use a list of the bits, with `true` used for 1 and `false` for 0. With that representation arithmetic operations would have costs driven by the bit-length of the numbers, and really that is a proper cost even though we tend to be used to thinking of arithmetic as having unit cost per operation. This modelling comes much closer to revealing how many electronic components will be activated in performing an operation.

But coding up binary arithmetic is a bit tedious, so it is fun to show that it is seriously easy to make pure lambda calculus explain the behaviour of integer arithmetic¹ really neatly. Rather than thinking in terms of zero, one, two, three and so on it is good here to think in terms of never, once, twice, thrice and the like, and a “number” can then be a function that applies a second function multiple times:

```
let never f a = a
let once f a = f a
let twice f a = f (f a)
let thrice f a = f (f (f a))
```

and this pattern allows you to have a function corresponding to any number of copies of `f` you could want to set up. With this representation of numbers a scheme to do basic arithmetic is almost frighteningly simple!

```
let increment n f a = n f (f a)
let add m n f a = m f (n f a)
let multiply m n f a = m (n f) a
let power m n f a = n m f a
```

and also a test for zero is remarkably easy to arrange

```
let iszero n = n f true where f x = false
```

Again one can appreciate how compact all the above are even without (at first) understanding how they work. And to get insight into the working it is

¹To be more pedantic natural number arithmetic, i.e. only using positive values

possible to start with simple expressions such as `power twice thrice f x` and follow through the definitions to observe it expanding until it ends up as (in this case) `f(f(f(f(f(f(f(f(a))))))))` which used `f` $8 = 2^3$ times. Or `iszero twice = twice f true = f(f true) = false`

There is one thing to be done here that is slightly more tricky, and one issue that has so far been swept under the carpet. The tricky issue is a function `decrement` to find the predecessor of a `n`. It should be expected that this will not be quite so straightforward because `never` does not have a proper predecessor. There are however several ways of managing this. The one shown here is not the standard one but is perhaps the most compact – but it relies on the precise way in which `increment` has been coded and it achieves its result by handing that an argument that is rather improperly not a number but that has been chosen so that the result returned is `never`. See how that can be done, starting with a letter `X` standing for what we will invent:

```
increment X f a = never
increment X f a = X f (f a) = never
let X a b = never
```

and the three lines above that introduce a function `X` that in just this one context serves as `-1` we end up with an implementation

```
let decrement n f a = n f X a where X a b = never
```

Concise it possibly confusing²

Any reasonable programmer will see that given `decrement` it will now be pretty trivial to code up subtraction and then division. If negative numbers are needed that can be handled using a sign-and magnitude representation by using `pair` to glue an explicit sign onto every number.

Now for the issue that had been avoided that far (but when you look back at all the fragments of code to date it does not impact them). In raw lambda calculus the functions do not have names. That means there is no obvious way for the definition of a function to refer back to the function itself. In ISWIM you could have felt tempted to write things rather like

```
let factorial n = if n=1 then 1 else n*factorial(n-1)
```

²A scheme that does not cheat by passing things to the increment function that are not numbers works by having a function that accepts an ordered pair as its argument: `let inc (a,b) = (a+1,a)`. If this is applied n times to a starting value $(0,0)$ the result is $(n,n-1)$ so the second component is the predecessor of n

as a simple recursive definition. On expanding that to lambda calculus the use of the word **factorial** within the body of the function will be improper and it certainly has no simple way to relate to the name you happen to be using for the function. So it looks as if recursion is going to be a problem! Happily this can be resolved. Often the way of sorting it is thought of as a bit strange. The presentation here is based on thinking of recursive definitions as just a notation for programs of infinite length that unwind it:

```
let f1 n = if n=1 then 1 else n*f2(n-1)
  where f2 n = if n=1 then 1 else n*f3(n-1)
    where f3 n = if n=1 then 1 else n*f4(n-1)
      where f4 n = if n=1 then 1 else n*f5(n-1)
        ...
```

and provided that chain of nested definitions is continued to a depth that is at least as great as the value of an argument n we are about to use this will do just what is needed. This sort of expansion can be applied to any recursively defined function³.

The magic that makes it possible to create unboundedly nested things in raw lambda calculus is a function traditionally called **Y**. Its behaviour can be characterised with the identity

$$Y\ f = f\ (Y\ f) = f(f(f(\dots$$

but because this is an explicitly recursive presentation it can not be used directly. However there are many ways in which ISWIM or lambda calculus can provide a **Y** function without needing recursion, and the simplest is

```
let Y f = g g where g h = f (h h)
```

In ISWIM notation this is really concise and rather easy to understand! Just simply following the steps we have $Y\ f = g\ g = f\ (g\ g) = f\ (Y\ f)$. and the ISWIM definition did not mention **Y** within the right hand side of its definition so it translated into 100% valid lambda calculus.

The general use of **Y** is that recursive function definitions can be expanded using the following rule:

```
let f x = ... f x' ...
 $\Rightarrow$ 
let f = Y G
  where G f =
    F where F x = ... f x' ...
```

³ISWIM insists on the use of **let rec** when a recursive function is being defined to highlight that something special is going on.

and the key thing here is that the recursive reference to `f` now has a meaning because it is talking about the argument to `G`, and if you trace through expanding the definitions out a few times you can see that this ends up being exactly what `f` was to end up as.

We now have boolean logic, data structures, arithmetic and recursion and with those it is pretty easy to code up anything else that can be imagined⁴. And the mappings from them down onto the extraordinarily simple basic rules of lambda calculus are concise and in general not too unpleasant to work with, so if anybody *really* wants to reason about the behaviour of code starting from first principles it can be reasonable to select lambda expressions and their behaviour as those fundamental principles. But now numbers and all the rest can (when necessary) be explained in this rather gory manner it is perhaps reasonable to consider a version of ISWIM that has all such features as built in primitives that can be implemented using traditional computer hardware and reasoned about at their own level of abstraction. The fundamentalist view of lambda calculus then just sits in the background providing assurance that the foundations for all that is built are really secure.

⁴Well interactive input and output, file-system access and the like are not parts of the world of concern just here!

Chapter 9

Primitive Recursion

Chapter 10

Solving a quadratic equation

There are introductions to programming that give as one of their earliest examples the challenge of creating an application that reads in three numbers, a , b and c and then prints out the two solutions to the equation $ax^2 + bx + c = 0$. The clear expectation is that this will be done using the well-known formula $\frac{-b \pm \sqrt{b^2 - 4ac}}{2a}$. When done by just writing out use of that formula the task is indeed easy enough to be in an introductory section about use of computers.

However if one looks at the task more carefully and start to insist on getting as correct a result as possible for all legitimate inputs things become rather different. I will suppose that since this is a task set for beginners it is to be solved using the default way in which your computer handles numbers. In almost all cases this will follow an international standard called IEEE 754 and the program that is written will use a representation they call “binary64” which is commonly referred to as “double precision”. The task in hand here highlights several ways in which the computer arithmetic that is thereby provided can do things that a naive user might not have thought about. So let’s take these in turn and see how the equation solver will need to be made more elaborate to avoid them hurting.

The (IEEE) floating point representation in a computer can not handle arbitrarily large numbers. Specifically it runs out of steam when values exceed around 1.8×10^{308} . Beyond that it stores values that represent “infinity”. Now of course sensible people will not tend to work with problems that lead to numbers so absurdly large, and so it is common not to think much about it! However if one seeks perfection then the quadratic solver should deliver accurate results for any inputs where the results are sensible, and it should report failure in exactly the cases where the inputs do not lead to answers that can be represented in the floating point format that the computer uses.

So now consider the case when some user provides the input $a = 1, b = 1, c = -1$. Substituting these into the formula leads to a pair of very sensible roots with values about 0.618034 and -1.618034 . Everything seems good. But now a rather less helpful user enthusiastically offers $a = b = 2 \times 10^{154}$ and $c = -2 \times 10^{154}$. This fairly obviously has exactly the same two roots. However the with this and with various fairly closely related cases involving huge numbers it is possible to arrange that in the computation of $b^2 - 4ac$ that either b^2 overflow or $4ac$ overflow or both, or that neither overflow but their difference does. It is furthermore possible for one or both of the terms to exceed the 1.8×10^{308} maximum and hence overflow even though when they are combined the result is in range. This is all a mess! However this particular case can be tidied up by starting the calculation by dividing all of a, b and c by some large value so as to reduce them to sensible size numbers. At which point an additional issue comes into play. If you divide a floating point number by a general scale factor doing so can introduce rounding errors. Consider for instance the calculation of just $1/3$ and the computer will produce something in the style of 0.3333333333333333^1 which is not precisely the same since it terminates after a finite number of digits. Continuing the calculation with this corrupted value will naturally lead to a (small) error in the final result. Happily with IEEE floating point it is legitimate to divide by any power of 2 and in that case no precision will be lost. So it will be necessary to identify a power of 2 that is about the right size so that it can be used to rescale the input to avoid overflow.

The situation with very very small numbers is in fact even more curious although the way to sort things out is basically the same. The smallest floating point value that can exist (with smaller values being flushed to zero) is about 4.94×10^{-3242} but once values get lower than 2.23×10^{-308} the representation starts to work with them at lower than normal precision. So to preserve as much accuracy as possible the scaling has to keep things away from not the point where underflow maps values to zero, but from some rather earlier-encountered threshold.

For now and to prevent things getting quite out of hand the issues of rounding errors that might arise when multiplying b by itself and so on up to and including those that could be introduced by the square root calculation are going to be ignored, but anybody who really wants to push for perfection regardless of the cost in difficulty can explore those paths. Also it would be proper to review cases where the true results are going to be very close to or beyond the overflow or underflow points so that good answers are produced

¹Since it is working in binary internally it will actually be more like $0.01010101\dots$
 _{$2^{2^{-1917}}$}

in every case where that is possible. Part of that might involve finding a scale factor σ and replacing a with σa and c with c/σ^3 . Then a final result can be generated by multiplying or dividing the solutions to the scaled version by σ , and any overflow or underflow will then be captured in and limited to that final re-scaling operation.

However there is a further and distinct way in which the naive use of the standard formula can generate seriously inaccurate results even after overflow is avoided and regardless of minor rounding errors. Suppose that the value of $-b$ and $\sqrt{b^2 - 4ac}$ are rather similar in absolute value. This easily happens when b is rather larger than either a or c . Then one of their sum and their difference will add two similar values and give a good result, while the other can lead to massive loss of precision as leading digits match and cancel out. To illustrate this consider the use of 8-digit decimal floating point on a pocket calculation and look at the equation $x^2 - 4003x - 1 = 0$. Here $-b$ is obviously just 4003.0000 when written so that the 8 digits of precision are made explicit. A careful calculation of $\sqrt{b^2 - 4ac}$ gives its value as 4003.000499625249859... and to 8 digits that is 4003.0005. When the computer subtracts these it has no access to any of the lower digits so the only result it can offer is -0.0005 while the ideal result would have been -0.00049962524 . So although the arithmetic had been being performed keeping 8 significant digits throughout the cancellation of leading digits in the subtraction means that our final result has at best only 4 of its leading digits correct. Using a number larger than 4003 would make this precision loss worse to an extent that even with full 64-bit IEEE arithmetic results can be unsatisfactory.

While we can all accept that the example shown here might have been chosen with awkward numbers deliberately picked to give this severe cancellation of leading digits, there is no fundamental reason why such cases might not arise in real life and in fact it can occur so easily that it should be considered a common risk.

Happily (for those who like to deliver accurate answers) or painfully (for those who see this adding yet an extra layer of complication and difficulty to the task) it is possible to avoid this calamity, because it is known that the product of the two roots of a quadratic will always be equal to c/a and because whenever calculating one of the two roots does a subtraction that can lead to leading digit cancellation the other will be found by doing an addition that gives full precision results. So in effect one should calculate the more delicate root as $2c/(-b - \sqrt{b^2 - 4ac})$ which will now be very respectable. Of course it will be necessary to judge which of the plus or minus cases is the one to use and which the one to avoid.

³with σ a power of 2 to avoid introducing corruption through rounding

So overall the program that solves a simple quadratic and that does not even worry about cases where there is no (real) solution is dramatically messier and calls for much more understanding that those novice programmers will have been ready to deploy. But if you are writing a library or application that is to be used by others you have a responsibility to cope with all cases, not just the ones you think about first!

Chapter 11

Turtle Graphics

One of the schemes often proposed for getting the very young into computation involved letting them draw pictures using a “turtle”. This moves around the world leaving a trail that shows where it has been (so why is it not described as a snail?). It is possible to instruct it to move directly forward by some number of steps or to turn left or right by an angle that is usually specified in degrees. Combinations of these two operations can be repeated. So two very easy initial examples of what can be done are

```
repeat 4 times
  move forward by 10
  turn left by 90 degrees
end
```

```
repeat 5 times
  move forward by 10
  turn left by 90
  move forward by 10
  turn right by 90
```

where one of these draws a square and the other a zig-zag. By giving instructions that are less repetitive it will be possible to draw a house or other interesting outlines. It can be useful to be able to say “pen up” and “pen down” so that the drawing being produced does not have to use a single continuous line and then perhaps the turtle can be used to create any drawing then could be made using a pencil. That sets one path towards difficulty: set out the instructions for a turtle to approximate some of the pencil work from Leonardo da Vinci or Albrecht Durer! That would probably just ends up as a hugely long list of movements and although the output would be spectacular

the text of the sequence of instructions to the turtle would not be very interesting or informative! So here we will concentrate on examples where the sequence of instructions is reasonably tidy. It was easy to understand what the square and zigzag scripts would lead to, but the point of this chapter is that with fairly harmless-looking extensions to the set of operations that a turtle can be asked to perform it gets remarkably harder to predict what will emerge or reason about it in detail. So what we provide here are a selection of more or less difficult questions and challenges regarding turtle behaviour and we will not spoil them all by giving all the answers!

1. Start with something that is not too hard. For exactly what angles of turn will a sequence rather like the one that draws a square return to its starting point, and how many steps will that take? Angles do not need to be whole numbers of degrees.
2. If the turtle position is computed using computer arithmetic that is only precise to say around 16 or 17 significant figures, for a pattern that would close up in an ideal world how far from joining up can it be in reality? What are the consequences if the turtle keeps following the same pattern of activity for a really long time? In other words why might it be that when I use a turtle to draw pentagons and try tracing all the way round many many times in some cases I keep following exactly the same path on each cycle, while in others I drift very very slowly away from where I started?
3. What rule will lead to the turtle following a nice spiral path, and how does it behave beyond the time it reaches the centre (of it ever does)?
4. Suppose that at each step the turtle moves forward by unit distance and then spins so that the next direction is utterly at random. That could include it keeping on in its original direction, totally backtracking or anything in between. After N steps about how far from its starting point is it likely to be?
5. As above, but the new random direction is limited, say to correspond to turning right by an angle uniformly chosen between 0 and 180 degrees? How much does this impact things as against the fully random turn?
6. Suppose the turtle has a home and it makes a random turn that is almost fully random but that has a rather small bias toward pointing it homewards. How big does this bias be to give it a good change of getting within a reasonable range of its island? This case can be interpreted as a reasonable first attempt to model real bird or animal

long range migration skills by exploring just how much navigational precision they actually need. And the traces of movement of several turtles trying this out can make nice pictures!

7. For parameters x and N consider the turtle instructions:

```
a = b = c = 0
repeat N times
  a = a + x
  b = b + a
  c = c + b
  move forward by 1
  turn left by c
```

This has introduced some arithmetic and is a generalisation of the challenge to understand what drawing will emerge from “move 1;turn 1;move 1;turn 2; move 1; turn 3;...” where the angle turned at each step grows. Note that turning by angles over 360° is perfectly respectable in that you just spin all the way around once (not having any overall effect!) and the turn by the specified angle less 360. The challenge here is to understand what values of x lead to closed paths, how long the paths are before they join up (i.e. how large should N be to make this neat), and what symmetries there will be in the picture created. For some values of x one gets a 3-fold symmetry. Just what values of x lead to that? Can one get a 5-fold symmetry ever? And why are the pictures so decorative?

Chapter 12

Reduce to range 0 to π

When needing to compute a trigonometric function such as sine or cosine of some number (and when you are working in radians rather than degrees) it will normally be right to start by reducing the argument to by subtracting off multiples of 2π . That is because these functions are periodic so altering the argument by any multiple of 2π does not change the value to be returned. It might in fact be better to reduce the range to $0 \dots \pi/2$ or possibly $0\pi/4 \dots \pi/4$ but the principles explained here apply in all such variations.

A naive way of doing the reduction would be something along the lines of

```
n = integer_part_of(x/(2*pi))
x = x - 2*n*pi
```

and if computer arithmetic was in perfect agreement with mathematical ideas of numbers this would be enough to do the job. However floating point numbers on a computer have limited precision. An immediate consequence of that is that the mathematical value of π can not be represented exactly there. So computationally (as distinct from mathematically) the subtraction of $2*n*\pi$ is going to subtract something that is not exactly what was ideally needed. A fine illustration of this arises when you simply ask the computer to show you the value of `sin(2*pi)`. To illustrate this imagine that the computer works to a precision of 4 digits in decimal (while double precision would be 52 binary digits which is messier to present here). Then the computer's idea of `pi` will be 3.142 precisely, and its model for $2*\pi$ will be 6.283. Note for a start that the second of these is not just twice the first! The computer can not do any better within the 4-digit precision we are working with. If you calculate the true value of `sin` of that exact value and keep just 4 significant figures you do not get zero. You obtain -0.0004073 . If the computer had

Some would try to claim that the point being made here is aa bit pedantic and a result of zero is what the user “expected” so would be better. Here the view is that a good implementation of elementary functions should treat input values as standing for exactly what they say (which will always be values with only a limited number of digits) and that results should match the mathematical result based on those values. And that anything else is sub-standard.

Looking at the above case it appears that to calculate accurate values of trig functions for large arguments (perhaps stupidly large arguments that sensible people do not use a lot – but perfectionists want to cope well with even awkward cases) is going to need serious multiple-precision arithmetic, as suggested by the need for 50 digits here. And for proper computer floating point the largest valid double precision float has value around $1.8 \cdot 10^{308}$ and that starts to suggest that it may be necessary to work with precision equivalent to over 300 decimals. That would start to be expensive – a.k.a. “difficult”.

0.1591549430918953357688837633725143620344596457404564...

6283000

¹With binary computer floating point one would multiply by a power of 2 here

interested in its value. So when you are doing the multiplication there is no point at all about forming partial products that will only add into that part. Also any partial products that contribute much further down than the four digits immediately after the decimal point are not relevant unless they lead to carries up into the important four digits. The result of that is that you can get the required digits of the fractional part of your input divided by 2π doing little more than a single length multiplication! The only cost is that you need to have stored rather a lot of digits of the value of $2/\pi$.

The above shows that by doing something just slightly tricky it is possible to turn something that looked at first as if it would either be impossible, or at best that it would be grimly expensive, into a fairly short calculation.

But for perfectionists this is not the end of the difficulty! Note that in the imaginary 4-decimal precision arithmetic being used here to illustrate what is done² every floating point number bigger than 999.9 in fact describes a whole number. This is in the spirit that for instance 6.233e3 has the integer value 6283 and the next floating point value up, i.e. 6.284e3 is then 6284. That is a simple observation to make. The challenging question is “How close to an exact multiple of 2π can a non-zero whole number less than 10^{308} be?” The reason this is important is that when we find, for some x , the fractional part of $x/(2\pi)$ some number of digits immediately after the decimal point might be zeros. So in fact it will not be good enough as suggested by the previous paragraph to collect just four digits there. It is necessary to collect enough that there will be four *significant* digits following any leading zeros³.

A useful observation is that π is a transcendental number and that means that the remainder can never be zero, so we always have a finite number of leading zeros there, and if we limit the values of interest to the range up to 10^{308} (a value picked as an approximation to the range supported by standard computer arithmetic) there must be a worst case. For almost all values of x there are no or very few leading zeros, but the concern here has to be to get every single case correct and that involves seeking the most extreme behaviour.

²Using real 52 or 53-bit machine floating point does not add any significant extra issues

³... and equivalent cases when the fraction has many leading 9 digits

Chapter 13

Ingenious Data Structures

Textbook of algorithms are often full of explanations of clever ways to do things that would not at first have been at all obvious. These have typically been invented because the more straightforward ways of solving the problems concerned can be improved on, often by huge amounts once you get to big enough test cases. So in this chapter a few of these schemes are presented as an illustration of the way in which seeking the most efficient scheme can escalate difficulty quite a lot.

13.1 Binomial Heaps

A problem that can have plenty of real-world application is the maintenance of a priority queue. With such a queue anybody joining the queue has an associated weight or priority. The queue is processed by insisting that whenever the server finishes with one customer they next look after whichever one in the queue has highest priority. Clearly this scheme arranges that a new customer but very important customer can arrive and jump much of the queue. There are two main operations whose cost needs to be considered: inserting a new customer into the queue and identifying and removing the next to be served.

The most naive scheme will be to keep all customers in the queue in a list arranged such that adding a new one has unit cost – and not at that stage worrying about the priorities. Then to pick the next one to serve it may be necessary to scan the whole of that list to identify the most worthy member of it, and excise them from the list regardless of whether they are its head, tail or somewhere in the middle. That may have optimised adding new customers to the queue but it makes selecting the next one to be served have a cost proportionate to the queue length. If this was in fact the best that

could be achieved all would be simple, but a datastructure called a “heap” improves on it sharply making the cost of both queue activities proportional to the logarithm of the queue length. For long queues this gives a very good saving.

The explanation of heaps given here is not going to go into all the details and tricks that proper textbooks do because the main payoff of this section is something that builds on the general idea. So for now a heap is a structure arranged as a (binary) tree. Each node of the tree holds a customer and references to two sub-trees. Two key rules are applied: the customer in each node is one who will have priority over every customer in either of the two sub-trees. And things are arranged so that the number of customers in each sub-tree are close to the same. The first of these means that server has immediate access to the most important customer – they are the one in the top node of the tree. The second ensures that the tree is nicely balanced and that if there are N customers in all the the height of the tree is only $\log_2(N)$ ¹. The management of such a heap is based on needing to be able to add new items to it while preserving its properties in time proportionate to its height, and equally being able to repair it when its top item is removed in a similarly fast way. Go and read the books to discover how that can be achieved, because the interest here is in making the problem slightly harder yet in a way that calls for further ingenuity.

Imagine you now have priority queues all sorted, and your setup now happens to have two servers each with their own separate queue. You are not allowed to make any assumption about which arriving customers joined which queue. Now one server finishes their shift, and the remaining one is left to handle all the work. This immediately calls for the two separate priority queues to be consolidated. One could do that by asking each customer who has been abandoned by the server they were waiting for to join the other queue one at a time, but since the queues we are now thinking about represent the queues as trees it is likely that this has a significant cost. The scheme sketched next reduces that almost as much as is possible!

Priority queues that may need to be merged can usefully be represented by a data structure known as a “Binomial Heap”. All operations on such heaps have costs no worse than the logarithm of the number of items stored. The explanation here will start from the top level so that it can motivate the data structure used by showing why it is good.

The key clever idea here is that if we have any number N we can look at how it would be written in binary notation and express it as the sum of a

¹Well that logarithm usually has a value that is not a whole number, so we need to round it up to get one!

bunch of powers of 2. So for instance 39 is 100111 in binary and that means $39 = 2^0 + 2^2 + 2^2 + 2^5$. If the number is N then we can say that there are $\log N$ bits in its binary representation.

If a Binomial Heap (which is going to act as a priority queue) has N items stored in it then they are arranged in a bunch of sub-heaps each of which has size that is a power of 2. It is not yet clear just how these will be represented, but it will be arranged that the highest priority item in each sub-heap is instantly accessible. That means that the top item in the whole heap can be found by checking each of the (up to) $\log N$ sub-heaps.

The clever part now emerges when you wish to consolidate two such heaps into one. The steps taken follow exactly the pattern used in performing addition on binary values. It can consider each possible power of 2 in turn. If neither input has a sub-heap that size then the output will not. If just one has then that will appear in the result. But if both do then those two sub-heaps get consolidated in a way that will be described soon into a single one that corresponds to the next power of 2 up. In terms of the binary addition this is a “carry”. Provided that the sub-heaps are kept with the 2^0 one first and provided consolidating a pair of sub-heaps on size 2^j into a single one of size 2^{j+1} is cheap this manages to add (or perhaps we say form the union) of the two binomial heaps in logarithmic time.

Now what about that consolidation step? Well a good way to represent a sub-heap of size 2^j is to have the highest priority item in it picked out and sitting at the top, and the remaining $2^j - 1$ items kept in a list of smaller heaps of size $2^{j-1}, 2^{j-2}, \dots, 4, 2, 1$. Happily this satisfies our hope that the top item in the sub-heap would be easy to find, and it means that the double size sub-tree can be formed very easily by comparing the top items in the two trees to be merged and just pushing the smaller one onto the list held by the larger.

The task of removing the top item from a heap is equally straightforward. We already know we can identify which sub-heap had the desired element at its head. Remove that whole sub-heap from the top-level list. Now if you lop the top item from the bit of structure you have just retrieved you have a nice list of sub-heaps each of whose size is a power of two. Gosh that is just the shape of a general Binomial Heap and you can re-insert all its data into the main one using just the binary addition process already described.

Those who are properly pedantic will observe that in each of the various sub-heaps you will want to have stored not just the top element and not just a reference to the list of sub-sub heaps, but something to explain how many items are present (i.e. which power of 2 is involved) and probably also a reference to the tail end of the chain of sub-sub-heaps so that tagging a new item on the end is really cheap. Doing all that carries some overhead

but for cases with enough customers the savings by having costs that grow only logarithmically with the size of the queues is so much more valuable that it is not a big issue.

The tricks and the elaboration of data representation here may seem extreme enough that it would be natural to expect it was the best that could be achieved. But massochists can look in the next chapter of their Big Book of Algorithms to learn about “Fibonacci Heaps” that are yet more bizarre – and which may only rather rarely be useful in practise because although from a theoretical analysis its costs grow slowly in practise the overheads mean that simpler schemes test to win. But for anybody keen to see how difficult computing can be made looking at the them, and at the Brodal Queue and all other options for priority queue implementation can provide a fine collection of rabbit holes to dive into.

13.2 What next?

I have not decided yet!

Chapter 14

Simple Pattern Matching

Sometimes you might want to search within some text but what you want to find is not just a fixed string. Perhaps it can allow options or repetition of sub-parts. Perhaps you want to put some sort of wild-cards into the pattern that is your target. There is a very well established scheme for setting up patterns for use in cases like this, and variations on it. Very many programming languages and even dialog-boxes in user interfaces use at least subsets of it. So for instance the pattern `*.jpeg` may be used to let you look for all files with the “jpeg” suffix, while at least in a Linux shell the pattern `*.\{cpp,h\}` will match names that end in either `.cpp` or `.h`. A fuller scheme used for pattern matching as part of the language PERL and available through libraries in almost all other programming languages as a bit more formal. A pattern is built up starting with the very simplest: patterns that consist of and match just one letter¹. These simple patterns are combined using three constructions, If P and Q are existing patterns then one can write

- PQ – this is a pattern that matches anything that can start with a sequence that matches P and follows that with one that matches Q . Obviously the very easiest use of this is that it means that you can write a sequence of individual letters and they form a word to be spotted;
- $P|Q$ – here we accept anything that matches either P or Q . So for instance `cat|dog|rabbit` matches strings that name creatures suitable as pets, and `p(e|a)t` illustrates that it is sometimes useful to have parentheses to group things. This pattern will match either `pet` or how you might treat one, i.e. `pat`. If you need one of the characters (,) or | in the alphabet you are matching over some way of distinguishing

¹It can also in fact be useful to have a basic pattern that matches an empty string.

raw characters from punctuation used to build up the pattern will be called for.

- P^* – This is the big one, It indicates an arbitrary repetition of the pattern P . So it is in effect equivalent to $(|P|PP|PPP|\dots)$. Note there the initial option of no instances of P , i.e. of this matching the empty string. A really simple instance of this would be $B(\text{an})^*a$ which matches Ba , $Bana$, $Banana$, $Bananana$ and so on.

There are two viewpoints that can be taken about this. One is a practical one that adds a number of shorthand notations for things one might frequently want to do. A particular instance of this arises because these patterns (which are referred to as “regular expressions”) provide an excellent way of characterising the ways in which tokens or symbols can be written in programming languages, and there are software tools that take a list of patterns and create a program that splits textual input up based on the. A first extension to notation that is used there is being able to give a name to a pattern fragment and then use it later. In the programs `lex` and `flex` one can name a fragment and then to refer to it you put the name in braces. You also enclose literal text in your pattern in double quotes. So for instance:

```
digit    "0"|"1"|"2"|"3"|"4"|"5"|"6"|"7"|"8"|"9"
number   {digit}{digit}*
```

gives a pattern for any (non-empty) string of digits and calls it “number”. This example motivates two further expansions which clearly do not alter the range of patterns that can be expressed but that can make the presentation of the regular expressions concerned much more compact. Enclosing a collection of characters in square brackets and allowing character ranges is a help. If the opening square bracket is followed by \wedge then the expression is treated as if was a square bracket form enclosing all letters in your character set except the ones actually shown. With this the tabulation of digits becomes just `[0-9]`. The second expansion allows for the fact that $*$ can indicate zero or more uses of the pattern that precedes it and sometimes as here you want at least one. Replacing the $*$ with $+$ does that. Hence you can now write

```
digit    [0-9]
number   {digit}+
```

Note that this could be textually expanded to the slightly clunky but basic for regular expressions, so the extended notations are in general a matter of convenience rather than things that bring genuine new capabilities. And in that spirit here are a few more useful extensions that similarly do not alter

the range of patterns that can be expressed but that may make it easier to specify them.

P & Q
 ~ P
 .

The idea is that the first of these will match every pattern provided that both P and Q do, while the second matches any input that P would fail to accept. The single dot (.) will match any single character, and so of course $.^*$ matches either anything at all (including nothing).

While it is fairly straightforward to show that adding these constructs does not add any ability to match new sorts of pattern – all they do is make it easier and more flexible to specify them – the notes here are not going to explain the details there. Head for a suitable textbook if you need to know exactly how it can be done! But the typical places on your computer that provides support for regular expressions will typically not support these last two because in fact they unlock levels of practical difficulty that are hard to comprehend.

Thus far the use of regular expressions to provide patterns that you can try to match against input text seems really rather easy despite the pessimistic statement above. So here to give some insight into just why adding those two last capabilities is so bad let's state what problem it turns from tolerable into being solvable in principle and theory but utterly dreadful in practise.

The sections here will next work through showing that it is possible to set up a regular expression of reasonable size that such that the only strings it can match will be absurdly long. Once that is done it becomes possible to argue that while answering various questions about them and their relatives is theoretically possible, the cost of doing so will be at least as great as the length of the shortest string they match, and hence is beyond all feasibility now and for ever.

The way of doing this will be by using the term “ruler” to characterise an expression that only matches strings of some given length L , and showing that if you are given a ruler for L you can derive one for a length of the order of 2^L such that the new expression is only some constant factor larger than the original one. By iterating this process you can obtain rulers of length $2^{2^{\dots^L}}$ for any height tower of powers of 2 that you want, and that leads to truly huge values rather rapidly.

To illustrate the process we will start with a ruler of length 3, and the simple regular expression $x\ x\ x$ consisting of just 3 characters does the job. The general concept that will be set up is based on counting in binary with

(in this case) 3 bit numbers, which may be presented to start with on two lines as

```
#000#001#010#011#100#101#110#111#
...#xxx#xxx#xxx#xxx#xxx#xxx#xxx#xxx#...
```

Here the upper line can be seen to be counting and it uses the hashes to keep the separate binary values neatly apart. The lower line uses our ruler and will help to enforce the regular pattern of where the hashes appear. It can be rendered as the simple expression `#(xxx#)*` which just takes our ruler and sets up repeats of it separated by `#` characters. Well this looks like two strings, one for counting and one for ruler. So then we can use the language of regular expressions we interleave the two so we have a single string where odd characters are from the top line and even from the bottom:

```
##0x0x0x##0x0x1x##0x1x0x##0x1x1x##1x0x0x##1x0x1x##1x1x0x##1x1x1x##
```

This may be less easy for a human reader to decode and so in the following presentation we will sometimes use the 2-line format, but what it will always mean will be the flattened out version. Note that with our ruler of length 3 the string we have here is of length $2 \times (3 + 1) \times 2^3 + 2$ where the $3 + 1$ and the final 2 come from counting the hashes and the extra factor of 2 is because there were two lines or text merged. Calling this just 2^3 is something of an understatement but that does not alter the main thrust of the presentation, which is that this will allow us to build extraordinarily long rulers.

Initially it might seem that setting up regular expressions that constrain text strings to count in binary is going to be hard. The way of doing so is to consider what strings are *not* counting sequences. So we will establish a set of regular expressions exact of which characterises something that we do not want to see, and using negation and the “and” operator ensure that none of them apply. What remains will only be able to be our counting sequence. This of course is the just what Sherlock Holmes taught us: “When you have eliminated all which is impossible, then whatever remains, however improbable, must be the truth.”

To start with it makes sense to enforce the block format of the string, which amounts to demanding that hashes in the top line can only be in places where they lie above hashes in the bottom line. Well any string that violated that rule will have either a hash above a non-hash or a non-hash above a hash somewhere. If just deal with the first of those it can be described using a pattern that carefully uses pairs of characters everywhere

```
(...)* (# (~#)) (...)*
```

This little expression and its friend that matches the bottom to top will form part of the expression we are building to describe our long string. As written

above it is only 17 tokens long. If you needed to use more primitive regular expressions each wildcard character `.` might expand into something like `a|b|c|d...0|1|#` and have a size proportional to the number of characters in your alphabet, but that again is just a constant.

The next constraint can be expressed positively and is that the string we are going to match must start at 000 and stop at 111. Again this is easy:

```
#. (0.)* #. (...)* #. (1.)* #.
```

where this version demands that our initial ruler is of length at least 2 so there are at least two hashes between the first and last segments. This works because `# 0* #` insures that there are only zeros in that region, and similarly for 1s at the end. And the previous rule has made certain that the `#` marks all line up and that forces those strings of zeros and ones to be the desired length. Note that all the messy things like `(0.)*` and indeed all the dots here are just arranging that we only pay attention to top-line (eg odd position) characters. because that is really just a bit of technicality in future we will present things in the form as if they were only applied to the top line and the extra mess to ignore the interleaved bottom line will be assumed to be applied:

```
# 0* # .* # 1* #
```

Looking ahead it will make sense to ensure that we only count through the binary sequence once, so that it is not possible to have a second section consisting all of zeros. This is also rather easy if we make it a pattern that we will say must not apply that puts two separating hash marks ahead of a block of all zeros. It is written here ignoring the issue of the lower ruler line...

```
.* # .* # 0* # .*
```

which is anything then `#`, more anything and then a block like `#000#`. By disallowing that we will count just once.

That leaves what may feel like the hard part which is to ensure that sequential blocks of digits count. It feels easiest to first explain what must happen, but the fragments of pattern we end up using will all need to be set up to disallow anything else. To increment a binary number one looks for the rightmost 0 and of course that means any digits to the right of it must all be 1s. The next number up must preserve all bits to the left of that key zero, flip the zero into a one, and turn all the trailing ones into zeros. A way of handling this is to imagine that corresponding bits in the number and its successor (which must be a distance apart set by our ruler) are underlined

and displayed in bold, then four cases apply which are shown a bit informally here:

```

...0 ... 0 ... # ... 0 ...
...1 ... 0 ... # ... 1 ...
...0      1*    # ... 1 ...
...1      1*    # ... 0 ...

```

The first two cases have at least one zero after the particular character and insist that the character remains unchanged, while the second two are when there are only 1s to the right in which case the bit is flipped. These cases are mutually exclusive but between them cover all possibilities.

The key trick is clearly to get the effect of the underlining set up in our petterns, and that can be explained in a “two row” presentation again. where for that first the above four cases what must be avoided is

```

.* 0 .* 0 .* # .* 1 .*
.* #      xxx.      # .*

```

which can be read out as the lower line using the ruler **xxx.** to insist that the two # marks are the length of our ruler plus one (to allow for the # in the main pattern) apart. The top line looks for zero which is followed by at last one zero before a #, and that after the # there is a one that has been forced to be just the right distance away. Matching a 1 there captures the case that is wrong, so saying that this pattern does not match leaves only the good cases. The ruler line used here is not the same as the one used before in that it puts in only two markers that are properly separated, and the leading and trailing .* parts say that it can match and indicate that separation anywhere. Ah well – adding this in will mean interleaving these two patters again, and the interleaving with the previous repeating ruler will have to remain in place. The result is that characters are going to end up being worked on in blocks of 3 or perhaps 4 rather than individually. This just makes the big ruler that we are creating with our counting scheme a bit longer yet and is not a big cause for worry.

Each of the fragments of regular expression introduced here are rather small even when all the interleaving stuff is allowed for, and in the end we have shown how to create a pattern whose size is some modest multiple of the size of the original ruler, but that will only match strings that are exponentially larger.

So why might this be a worthwhile exercise? Apart of course from it being a demonstration that pattern matching can do more than was first apparent. Well it can be part of the process of showing that understanding particular patterns can be hard. Begin with a fundamental result about these patterns, which is that any regular expression will have a corresponding deterministic

finite automaton that can recognise exactly the strings that the pattern will match. Well to rephrase that in less technical language for any pattern there is a really simple program that can be used to apply it to input strings. The program is basically of the form:

```
int state = 0;
for each character in the input
    state = transitionTable[state, character];
return isAccepting[state];
```

where the transition table is used to adjust the state as each fresh character is presented and the `isAccepting` vector reports at the end whether the string matched. It is clear that the work done for each input character is rather small – just accessing the array, and so this is liable to be really efficient way of performing pattern matching. There are readily available programs² that can build the tables and they will tend to be a bit cleverer than using a simple 2 dimensional array for the transitions would be because many entries in the full rectangular block would never be used.

Both for practical and theoretical purposes it is worthwhile to consider how many distinct states will be required to build a program of the above form that will correspond to a given pattern. One bound that is pretty clear cut is that there must be at least as many states as the shortest string that matches the pattern. To see that imagine the sequence of states encountered in the process of a successful match of such a shortest string. If there were to be fewer states than the length of the string then one of the along the way would have to be repeated, as in

ABCDEFDGH I J

where state D has been entered twice. Well if that happened it would be possible to provide input that ran up to that situation and then continued on a home run, omitting all the steps between the first and second encounter with the repeating state.

ABCDGH I J

This is a shorter string that has just been accepted, contradicting the assertion that we had started by looking at the shortest matching input. What we can read off from this observation is that provided we allow extended regular expressions with a negation operator there are tolerably concise patterns that would necessarily lead to super-galactically large numbers of states. This

²notably `lex` and `flex`

is a pretty convincing reason not to allow negation in practical applications of these patterns!

So next consider a challenge that if it was readily solvable would be of great use: given some pattern expressed as a regular expression is there another more concise pattern that matches exactly the same set of inputs? From the perspective of a theoretician this can be addressed by emulating all possible regular expressions with all the ones using k characters before the ones with $k + 1$ characters, and with ones that are the same length in alphabetic order. Over any fixed alphabet of the characters you are working with this is clear-cut. Now for each such you want to test if it happens to match the same inputs that your original one did. Well if I have two regular expressions P and Q then the language $(P \& Q | Q \& P)$ will match any things that P does but Q does not and vice versa. So if the two expressions do the same thing this new one I have just set out will not match anything at all. But that is something that there is a clear-cut way of checking for. Make the transition table for this new composite expression and supposing it turns out that there are N states then if there is any string that it accepts at all there must be one of length at most N . This is just the flip-side of the previous remark about lengths of matched strings and the number of states. If your alphabet is of size k then there are only(!) k^N strings of length N to check and if none of those match then nothing will. The explanation here is a bit needlessly inefficient but it demonstrates that in a finite amount of work one can tell if two patterns match the same set of strings and that then by exhaustive search in a finite amount of time it would be possible to exhibit the most concise pattern for any particular behaviour. But although this shows that the challenge is decidable in the sense that one can have a computer program that would guarantee to address it and eventually complete its work with a correct result, the resources needed to achieve that might well be truly ridiculous.

The explanations given here show how huge rulers and hence regular expressions that necessarily lead to transition tables with huge numbers of states can arise. They do not complete the proof that there can not be some clever way of reasoning about e.g. equivalence between patterns that manages to sidestep generating the associated transition table. And details of that will be left as a further literature search or research project for the reader. But a key step that it may by now be easy to recognise is that the sorts of techniques explained here that block a string into segments each representing the next binary number on from the previous can be modified so that the string is split into blocks such that each represents the next state of a fairly arbitrary computer one step on from the previous. It would be normal to set this up describing a Turing machine with a tape whose length is limited

to the size of our ruler. Given that and the large amount of theory known about the problem of telling if a Turing machine will halt you can imagine that formalising a statement that there can be no amazing short cuts and that something very much like building the full state transition table for the description is going to be necessary to tell if certain messy patterns will match anything at all.

In writing this the authors here want to leave the reader understanding that there is always more to be looked into!

Chapter 15

Parallelism and Concurrency

Chapter 16

Puzzles

Chapter 17

Floating Point

Chapter 18

Formal Proofs

Chapter 19

Iteration

Chapter 20

Hardware

Chapter 21

Networks and other interconnects

Chapter 22

AI

Chapter 23

Unusual Computer Language

Chapter 24

Undecidable problems

Chapter 25

Lessons that have been learned

The comments here are just ACN rambling a bit more

I have just picked up a copy of Wolfram's New Kind of Science for not a huge amount on eBay. That is a much bulkier volume than I had expected, not having checked it before. But what is maybe more terrifying is that Wolfram now has a "20 years on" book that he explains basically as "When I write ANKoS I thought it was a real breakthrough for the world, but 20 years on I see it is way better than even the extreme level of importance I saw in it back then".

I have only just started reading it and the main think that comes across is how utterly Wolfram wants to make a point that everybody from before the ancient Greeks has just skipped past the motherload of overwhelmingly important stuff that he and he alone has discovered. He is asserting that what he has discerned upends every scientific and many other disciplines totally and provides a way to discern the true nature of everything. Gosh it is amazing in that way. Wow – what a guy.

However for the purposes of what this book wants to do it is perhaps rather nice in that I think a major point he is wanting to hammer on is that something that follows very simple rules can have astonishingly complicated behaviour, and that this applies not to just one sort of "simple thing" but rather generally.

It is less clear to me (as yet) whether he can then do anything interesting with the complexity apart from show it off in loads of pictures. And some of us sort of believed that the not-too-bad equations of fluid dynamics could lead to very messy turbulence and not just smooth flow, that looking at multiplication and division dumped one into the quasi regularity of the distribution of prime numbers and great depth, and that the investigation simple problems like "boolean satisfiability" could tell you about all the other NP-complete problems. But that some of these are discrete and some continuous so he will

have to do a really merry dance to convince me that they are all the same even if all show complicated behaviours. But still reviewing all he has to say is liable to reveal a range of very fine examples of things where the starting point is simple and the end-point really is not.

A lot of what he talks about is cellular automata.

A different think to consider is “solve your problem by first designing and building your computer, then the software stack...”. This is of course just what people had to do in the 1940s. And indeed Babbage/Lovelace had a go there, and there was Konrad Zuse and his relay-based computer where one can even at least imagine constructing the relays...

In yet a different direction I think of steam engines. The pistons must slide nicely into their cylinders so they have to move in straight lines, but if you support them with sliding supports that might introduce friction you do not like. So how do you make a pin-jointed linkage so that the end-point moves in exactly a straight line? One answer is Hart’s Inversor. Now having invented that if you are a masochist you set up all the simultaneous equations that characterise the way that the other end of a rod that has a fixed pivot at one end lies on a circle, and so on. You then see if you can simplify and solve all those equations to prove that the key endpoint lies on a straight line. This is a horrid thing to try Groebner Bases on.