

Computational Thinking for Actuaries and Financial Professionals

With Applications in Julia

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Preface

This book is intended to enable practitioners and advanced students of financial disciplines to utilize the tools, language, and ideas of computational and related sciences in their own work.

Throughout these pages, you'll see how best to represent ideas so they remain scalable, intelligible to collaborators, and robust against changing conditions. The book covers fundamentals in programming to advanced topics like parallelization and data visualization, always returning to practical applications in insurance, asset modeling, and beyond.

The goal is straightforward: to empower financial professionals with the knowledge and confidence to manipulate computational tools in ways that reveal ever richer insights into the systems we seek to understand and influence.

A PDF of the book is available by [clicking here](#).

Warning

This book is currently being drafted. ANY AND ALL content is subject to change, including the license. To report issues or other feedback, please email Alec at firstnamelastname@gmail.com.

Part I.

Introduction

"I think one of the things that really separates us from the high primates is that we're tool builders. I read a study that measured the efficiency of locomotion for various species on the planet. The condor used the least energy to move a kilometer. And, humans came in with a rather unimpressive showing, about a third of the way down the list. It was not too proud a showing for the crown of creation. So, that didn't look so good. But, then somebody at Scientific American had the insight to test the efficiency of locomotion for a man on a bicycle. And, a man on a bicycle, a human on a bicycle, blew the condor away, completely off the top of the charts.

And that's what a computer is to me. What a computer is to me is it's the most remarkable tool that we've ever come up with, and it's the equivalent of a bicycle for our minds." - Steve Jobs (1990)

The world of financial modeling is incredibly complex and variegated. It, along with many of the sciences, is a place where practical goals harness computational tools to arrive at answers that (we hope) are meaningful in a way that tells us more about the world we live in. What this usually means specifically is that practitioners utilize computers to do the heavy work of processing data or running simulations which reveal the something about the complex systems we seek to represent. In this way, then, financial modelers must also be a craftsman who seeks not only to design new products, but must also think carefully about the tools and the process used therein.

This book seeks to aid the practitioner in developing that workmanship: we will develop new ways to look at the *process*, think about how to most clearly represent ideas, dive into details about computer hardware and bring it back up to the most abstract levels, and develop a vocabulary to more clearly express and communicate these concepts. The book contains a large number of practical examples to demonstrate that the end result is better for the journey we will take.

This book looks at programming for the applied financial professional and we will start by answering a very basic question: "*why is this relevant for financial modeling?*". The answer is simple: financial modeling is complex, data intensive, and often very abstract. Programming is the best tool humans have so far developed for rigorously transforming ideas and data into results. A builder may be the most skilled person in the world with a hammer but another with some basic training in a richer set of tools will build a better house. This book will enhance your toolkit with experience with multiple tools: a specific programming language, yes, but much more than that: a language to talk about solving problems, a deeper understanding of specific problem solving techniques, how to make decisions about what the architecture of a solution looks like, and practical advice from experienced practitioners.

The approach

The authors of the book are practicing actuaries, but we intend for the content to be applicable to nearly all practitioners in the financial industry. The discussion and examples may have an orientation towards insurance topics, but the concepts and patterns are applicable to a wide variety of related disciplines.

We will pull from examples on both sides of the balance sheet: the left (assets) and right (liabilities). We may also take the liberty to, at times, abuse traditional accounting notions: a liability is just an asset with the obligor and obligee switched. When the accounting conventions are important (such as modeling a total balance sheet) we will be mindful in explaining the accounting perspective. In practice, this means that we'll take examples that use examples of assets (fixed income, equity, derivatives) or liabilities (life insurance, annuities, long term care) and show that similar modeling techniques can be used for both.

What you will learn

It is our hope that with the help of this book, you will find it more efficient to discuss aspects of modeling with colleagues, borrow problem solving language from computer science, spot recurring structural patterns in problems that arise, and understand how best to make use of the “bicycle for your mind” in the context of financial modeling.

It is the experience of the authors that many professionals that do complex modeling as a part of their work have gotten to be very proficient *in spite of* not having substantive formal training on problem solving, algorithms, or model architecture. This book serves to fill that gap and provide the “missing semester” (or “years of practical learning”!). After reading this book, we hope that you will *appreciate* the attributes of Microsoft Excel that made it so ubiquitous, but that you *prefer* to use a programming language for the ability to more naturally express the relevant abstractions which make your models simpler, faster, or more usable by others.

Even if your direct responsibility does not entail hands-on-coding, be it management or “low-code”, the ideas and language should prove useful in guiding the work to a cleaner, more efficient solution.

The Journey Ahead

Learning a new topic, especially one that's not well trodden in a given field, can be intimidating. There are many resources available online, this book will recommend some others, and there are community support resources available - check the chat and

forums and look for the users talking about the topics that interest you. One of the wonderful things about the technology community is the degree to which content is available online for learning and reference.

Further, moving substantial parts of the financial services industry towards a digital-first, modern workflow is a monumental effort and you should seek partners on both the finance and information technology side. In general, good ideas and processes will prevail. The trick to encouraging adoption is finding the right place to plug a new idea or suggestion.

Additionally, this book provides the language and technical knowledge to partner with others (such as peers and IT) to make pragmatic decisions about the tradeoffs that will need to be made.

What to Expect

This book will guide you through:

1. Core programming concepts applied to finance
2. Modern software development practices
3. Computational approaches to common financial problems
4. Real-world examples and applications

The goal is to build both theoretical understanding and practical skills you can apply immediately in your work.

Prerequisites

Basic experience with financial modeling is not strictly required, but it will benefit the reader to be familiar so that the examples will not be attempting to teach both financial maths and computer science simultaneously.

Advanced financial maths (e.g. stochastic calculus) is *not* required. Indeed, this book is not oriented to the advanced technicalities of Wall Street “quants” and is instead directed at the multitudes of financial practitioners focused on producing results that are not measured in the microseconds of high-frequency trading.

Prior programming experience is *not* required either: Chapter 5 introduces the basic syntax and concepts while `@#sec-julia-writing` covers setting up your environment to follow along. For readers with background in programming, we recommend skimming Chapter 5.

The Contents of This Book

The book is organized into eight parts, each addressing key aspects of computational thinking and financial modeling. Part I introduces foundational concepts, explaining why programming matters for financial professionals, and why Julia is particularly well-suited for financial modeling applications.

Parts II, III, and IV establish the theoretical foundations—covering effective financial modeling practices, programming abstractions, and techniques for building performant models. These sections bridge theory with practical implementation, exploring topics like model design, functional programming, data types, and parallelization strategies.

Part V connects interdisciplinary concepts with practical applications, demonstrating how software engineering practices, computer science principles, statistical methods, and visualization techniques enhance financial modeling.

Part VI provides detailed guidance on developing in Julia, from writing and troubleshooting code to optimization. This is the section that really leans into Julia-specific ideas and workflows.

Part VII showcases applied financial modeling techniques through real-world examples, including stochastic mortality projections, scenario generation, sensitivity analysis, and portfolio optimization.

While Julia is used for the examples throughout the book, the concepts presented are largely language-agnostic. The principles of computational thinking and financial modeling remain applicable regardless of implementation language. Readers are encouraged to follow along with the examples on their own computers, with the entire book available on GitHub for reference [#TODO: determine book URL].

Notes on formatting

When a concept is defined for the first time, the term will be **bold**. Code, or references to pieces of code will be formatted in inline code style like `1+1` or in separate code blocks:

"This is a code block that doesn't show any results"

"This is a code block that does show output"

"This is a code block that does show output"

When we show inline commands are to be sent to Pkg mode in the REPL (see `?@sec-environments`), such as such as `add DataFrames`, we will try to make it clear in the context. If using Pkg mode in standalone codeblocks, it will be presented showing the full prompt, such as:

```
(@v1.10) pkg> add DataFrames
```

There will be various callout blocks which indicate tips or warnings. These should be self-evident but we wanted to point to a particular callout which is intended to convey advice that stems from practical modeling experience of the authors:

 Financial Modeling Pro-tip

This box indicates a side note that's particularly applicable to improving your financial modeling.

Colophon

The HTML and PDF book were rendered using Quarto and Quarto's open source dependencies like Pandoc and LaTeX.

The HTML version of this book uses Lato for the body font and JuliaMono for the monospace font.

The PDF version of this book uses TeX Gyre Pagella for the body font and JuliaMono for the monospace font.

The cover was designed by Alec Loudeback using Affinity Designer with the graphic used under permission by user cormullion on Github.

This book was rendered on May 14, 2025. The system used to generate the code and benchmarks was:

```
versioninfo()
```

Julia Version 1.11.5

Commit 760b2e5b739 (2025-04-14 06:53 UTC)

Build Info:

Official <https://julialang.org/> release

Platform Info:

OS: macOS (arm64-apple-darwin24.0.0)

CPU: 14 × Apple M4 Max

WORD_SIZE: 64

LLVM: libLLVM-16.0.6 (ORCJIT, apple-m1)

Threads: 1 default, 0 interactive, 1 GC (on 10 virtual cores)

Environment:

JULIA_PROJECT = @.

JULIA_LOAD_PATH = @:@stdlib

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1. Why Program?

"Humans are allergic to change. They love to say, 'We've always done it this way.' I try to fight that. That's why I have a clock on my wall that runs counterclockwise." - Grace Hopper (1987)

1.1. In this Chapter

We motivate why a financial professional should adopt programming skills which will improve their own capabilities and enjoyment of the discipline, whilst allowing themselves to better themselves and the industry we work in.

1.2. Introduction

The financial sector is undergoing a profound transformation. In an era defined by big data, (pseudo) artificial intelligence, and rapid technological advancement, the traditional boundaries of finance are expanding and blurring. From Wall Street to Main Street, from global investment banks to local credit unions, technology is reshaping how financial services are delivered, how risks are managed, and how decisions are made.

This digital revolution is not just changing the tools we use; it's fundamentally altering the skills required to succeed in finance. In the past, a strong foundation in mathematics, economics, and financial theory was sufficient for most roles in the industry. Today, these skills, while still crucial, are increasingly being augmented (and in some cases) superseded by technological proficiency.

At the forefront of this shift is the growing importance of programming skills. In the beginning of the computer era in finance, the differentiating skill was being able to utilize digital computing, data processing, and calculation engines to automate, analyze, and report on the business. These skills required low level programming and the success of many of those early programs is evident in their legacy: many of them are still around in the 2020s!

At some point, due to regulatory pressures, attempts at organization efficiencies, or management decision making, the skill of programming became highly specialized and most

1. Why Program?

financial professionals (investment analysts, actuaries, accountants, etc.) became relegated to being “business users”, utilizing either Microsoft Excel or a proprietary third-party software to accomplish their responsibilities. The reasons for this were not totally wrong, even in retrospect.

At some point, an increasingly complex stack of software sitting between the developer and the hardware, the proliferation of computer security risks it makes some sense that many financial developers were pushed out of the programming trade. Instead specialized, separate business and IT units were developed. Of course, this led to many inefficiencies and is now swinging back the other way.

What’s changed that’s enabling financial professionals to re-engage with the powerful tools that programming provides? Some reasons include:

1. **Code management tools.** Github and other version control systems provide best-in-class ways of managing codebase changes and collaboration. Tools exist to scan repositories for leaking secrets, security vulnerabilities, and dependency management.
2. **Increasingly accessible development.** Originally, very few layers of complexity existed between the written code and running it on the mainframe. Over time, drivers, operating systems, networking, dependencies, and compilers made development more complex. Today, languages, libraries, code editors, and deployment tools have smoothed many of these frictions.
3. **Competitive Pressures.** An increasingly commoditized financial product with evermore competition has led to a need to improve efficiency of manufacturing and selling financial products. Having a business developer is a lot more efficient than a business user who needs to get an IT developer to implement something. Further, pressures from outside the financial sector abound: It’s easier to teach a tech developer enough to be successful in a finance role than it is to teach a finance professional development skills.
4. **Regulatory and Risk Demands.** Pressures that previously motivated the move to proprietary software for modeling included regulatory reporting, internal risk metrics, and management performance evaluation. However, companies are realizing that their unique products, risk frameworks, preferred management measurements, and employee potential means that having a bespoke internal model is seen as a key capability. Many regulatory frameworks also encourage the use of a bespoke model, which is a particularly attractive option especially for those who view the given regulatory framework as inappropriately reflecting their own business and risk profile.

Whether you’re an investment banker modeling complex derivatives, an actuary calculating insurance risks, a financial planner optimizing client portfolios, or a risk manager stress-testing scenarios, the ability to code is becoming as fundamental as the ability to

use a spreadsheet was a generation ago. To remain competitive, adaptable, and effective in the evolving landscape of finance, professionals must embrace programming as a core skill.

Note

One subset of business analysts that *did not* start to migrate away from development as a strategic part of their value were “quants” or quantitative analysts who heavily utilized programming skills to develop unique products, trading strategies, modeling frameworks, and risk engines. This book is not really for that class of people and is instead geared towards the mass of financial professionals who want to get some of the benefits of the tools that the quants have been using for years. Quants may find value here in adapting some of their existing knowledge with the concepts and capabilities that Julia enables.

As we delve into this topic, keep in mind that learning to code is not about replacing traditional financial acumen—it’s about augmenting and enhancing it. It’s about equipping yourself with the tools to tackle the complex, data-driven challenges of modern finance. In short, it’s about future-proofing your career in an industry that is increasingly defined by its ability to innovate and adapt to technological change.

1.2.1. Market Forces

Today, there is a trend towards technological value-creation and is evident across many traditional sectors. Tesla claims that it’s a technology company; Amazon is the #1 product retailer because of its vehement focus on internal information sharing¹; Airlines are so dependent on their systems that the skies become quieter on the rare occasion that their computers give way. Companies that are so involved in *things* (cars, shopping) and *physical services* (flights) are so much more focused on improving their technological operations than insurance companies *whose very focus is ‘information-based’?* **The market has rewarded those who have prioritized their internal technological solutions.**

Commoditized investing services and challenging yield environments have reduced companies’ comparative advantage to “manage money”. Spread compression and the explosion of consumer-oriented investment services means a more competitive focus on the ability to manage the entire asset or policy’s lifecycle efficiently (digitally), perform more real-time analysis of experience and risk management, and handle the growing product and regulatory complexity.

These are problems that have technological solutions and are waiting for insurance company adoption.

¹Have you had your Bezos moment? What you can learn from Amazon.

1. Why Program?

Companies that treat data like coordinates on a grid (spreadsheets) *will get left behind*. Two main hurdles have prevented technology companies from breaking into insurance and traditional finance:

1. High regulatory barriers to entry, and
2. Difficulty in selling complex insurance products without traditional distribution.

Once those two walls are breached, traditional finance companies without a strong technology core will struggle to keep up. The key to thriving is not just adding “developers” to an organization; it’s going to be **getting domain experts like financial modelers to be an integral part of the technology transformation**.

1.3. Why Programming Matters Now

Programming is becoming as fundamental for financial professionals as spreadsheet skills were a generation ago. Here’s why:

1. **Enhanced Analysis Capabilities:** Programming allows for more complex analyses, handling of larger datasets, and application of advanced statistical and machine learning techniques.
2. **Automation and Efficiency:** Repetitive tasks can be automated, freeing up time for more value-added activities.
3. **Customization:** Bespoke solutions can be developed to address unique business needs, risk frameworks, and regulatory requirements.
4. **Data Handling:** As data volumes grow, programming provides tools to efficiently process, analyze, and derive insights from vast amounts of information.
5. **Integration:** Programming skills enable better integration across different systems and data sources, providing a more holistic view of financial operations.
6. **Competitive Edge:** In an increasingly technology-driven industry, programming skills can be a significant differentiator.

It’s now commonly accepted that to gather insights from your data, you need to know how to code. Modeling and valuation needs, too, are often better suited to customized solutions. Let’s not stop at data science when learning how to solve problems with a computer.

1.4. The Spectrum of Programming in Finance

It's important to note that becoming proficient in programming doesn't mean you need to become a full-time software developer. There's a spectrum of programming skills that can benefit financial professionals:

1. **Basic Scripting:** Automating repetitive tasks in Excel or other tools.
2. **Data Analysis:** Using languages like Python or R for statistical analysis and visualization.
3. **Model Building:** Developing financial models or risk assessment tools.
4. **Full-Scale Application Development:** Creating more complex applications for internal use or client-facing solutions.

1.5. Avoiding Red Herrings

One tantalizing path to contemplate is to avoid *really* learning how to code. Between artificial intelligence (AI) solutions being developed and low-code offerings, is there really a need to learn the fundamentals of coding? We argue that there is, for the basic reason that coding is not about mechanically typing out lines in an editor, but both a tool and a craft that is designed to enhance and apply your own creative and logical thinking.

The current generation of AI is fundamentally limited. Yann LeCun, Meta's (Facebook's) Chief AI Scientist describes the large-language model (LLM) approach as not even at the level of intelligence of a cat, and that we are decades away from true artificial general intelligence (AGI). These models have a "very limited understanding of logic . . . do not understand the physical world, do not have persistent memory, cannot reason in any reasonable definition of the term and cannot plan . . . hierarchically" (Murphy and Criddle 2024).

An important role for AI to play will be to *support* modelers in boilerplate, syntactical hurdles ("in VBA I would do it like this, but in Julia how do I do X, Y, or Z"), and basic algorithmic support. What is not likely to change in the short term is most of the value that a modeler brings to the table: creative thinking, understanding of company and market dynamics, and capability to understand broader architecture and conceptual aspects of modeling.

A similarly fraught path is low-code solutions. Low-code solutions are inherently limiting in their capabilities and lock you into a particular vendor solution. If you know enough about what you are trying to do to be able to state it in clearly in plain English, then you are most of the way to being able to program in a full coding solution (AI can actually help bridge this gap here). As soon as you hit a limitation of the system ("I'd like to use XYZ optimization algorithm at each timestep"), you are reliant on the vendor

1. Why Program?

to implement that option in the “low code” solution. Further, you are out-sourcing a lot of the important inner-workings of the model to someone else and not building that expertise yourself of in-house somewhere.

1.6. The 10x Modeler

The increasingly complex business needs will highlight a large productivity difference between a financial modeler who can code and one who can’t — simply because the former can react, create, synthesize, and model faster than the latter. From the efficiency of transforming administration extracts, summarizing and aggregating valuation output, to analyzing available data in ways that spreadsheets simply can’t handle, you can become a “10x Modeler”².

i Note

In the technology sector, a 10x developer is term for a software engineer who is an order of magnitude more productive, creative, or capable than a typical peer. Here, we extend the notion to developers of financial models.

Flipping switches in a graphical user interface versus being able to *build models* is the difference between having a surface-level familiarity and having full command over the analysis and the concepts involved — with the flexibility to do what your software can’t.

Your current software might be able to perform the first layer of analysis, but be at a loss when you want to take it a step further. Tasks like visualizations, sensitivity analysis, summary statistics, stochastic analysis, or process automation, when done programmatically, are often just a few lines of additional code over and above the primary model.

Should you drop the license for your software vendor? No, not yet anyway. But the ability to supplement and break out of the modeling box has been an increasingly important part of most professionals’ work and this trend appears to be accelerating.

Additionally, code-based solutions can leverage the entire-technology sector’s progress to solve problems that are *hard* otherwise: scalability, data workflows, integration across functional areas, version control and versioning, model change governance, reproducibility, and more.

30-40 years ago, there were no vendor-supplied modeling solutions and so you had no choice but to build models internally. This shifted with the advent of vendor-supplied modeling solutions. Today, it’s never been better for companies to leverage open and

²The 10x [Rockstar] developer is NOT a myth

1.7. Risk Governance

inner source to support their custom modeling, risk analysis/monitoring, and reporting workflows.

i Note

Open source refers to software whose source code is freely available for anyone to view, modify, and distribute. It promotes collaboration, transparency, and innovation by allowing developers worldwide to contribute to and improve the codebase. Open source projects often benefit from diverse perspectives and rapid development cycles, resulting in robust and widely-adopted solutions.

Inner source applies open source principles within a single organization. It encourages internal collaboration, code sharing, and transparency across different teams or departments. By adopting inner source practices, companies can reduce duplication of effort, improve code quality, and foster a culture of knowledge sharing. This approach can lead to more efficient development processes and better utilization of internal resources.

It is said that you cannot fully conceptualize something unless your language has a word for it. Similar to spoken language, you may find that breaking out of spreadsheet coordinates (and even a dataframe-centric view of the world) reveals different questions to ask and enables innovative ways to solve problems. In this way, you reward your intellect while building more meaningful and relevant models and analysis.

1.7. Risk Governance

Code-based workflows are highly conducive to risk governance frameworks as well. If a modern software project has all of the following benefits, then why not a modern insurance product and associated processes?

- Access control and approval processes
- Version control, version management, and reproducibility
- Continuous testing and validation of results
- Open and transparent design
- Minimization of manual overrides, intervention, and opportunity for user error
- Automated trending analysis, system metrics, and summary statistics
- Continuously updated, integrated, and self-generating documentation
- Integration with other business processes through a formal boundary (e.g. via an API)
- Tools to manage collaboration in parallel and in sequence

These aspects of business processes are what technology companies *excel* at. There is a litany of highly robust, battle-tested tools used in the information services sectors. This

1. Why Program?

book will introduce much of this to the financial professional (specifically Chapter 12, Chapter 21, and Chapter 23).

1.8. Managing and Leading the Transformation

For managers: the ability to understand the concepts, capabilities, challenges, and lingo is not a dichotomy, it's a spectrum. Most actuaries, even at fairly high levels, are still often involved in analytical work. Still above that, it's difficult to lead something that you don't understand.

Conversely, the skill and practice of coding enhances managerial capabilities. When you are really skilled at pulling apart a problem or process into its constituent parts and designing optimal solutions... that's a core attribute of leadership as well as the most essential skill in programming. This perspective also allows for a vision of where the organization *should be* instead of thinking about where it is now.

The skillset described herein is as important an aspect of career development as mathematical ability, project collaboration, or financial acumen.

1.9. Outlook

It will increasingly be essential for companies to modernize to remain competitive. That modernization isn't built with big black-box software packages; it will be with domain experts who can translate their expertise into new forms of analysis - doing it faster and more robustly than the competition.

2. Why use Julia?

We want a language that's open source, with a liberal license. We want the speed of C with the dynamism of Ruby. We want a language that's homoiconic, with true macros like Lisp, but with obvious, familiar mathematical notation like Matlab. We want something as usable for general programming as Python, as easy for statistics as R, as natural for string processing as Perl, as powerful for linear algebra as Matlab, as good at gluing programs together as the shell. Something that is dirt simple to learn, yet keeps the most serious hackers happy. We want it interactive and we want it compiled. - Bezanson, Karpinski, Shaw, Edelman (the original creators of Julia)

Julia is a relatively new¹, productive, and fast programming language. It is evident in its pragmatic, productivity-focused design choices, pleasant syntax, rich ecosystem, thriving communities, and its ability to be both very general purpose and power cutting edge computing.

With Julia: math-heavy code looks like math; it's easy to pick up, and quick-to-prototype. Packages are well-integrated, with excellent visualization libraries and pragmatic design choices.

Julia's popularity continues to grow across many fields and there's a growing body of online references and tutorials, videos, and print media to learn from.

Large financial services organizations have already started realizing gains: BlackRock's Aladdin portfolio modeling, the Federal Reserve's economic simulations, and Aviva's Solvency II-compliant modeling². The last of these has a great talk on YouTube by Aviva's Tim Thornham, which showcases an on-the-ground view of what difference the right choice of technology and programming language can make. Moving from their vendor-supplied modeling solution was **1000x faster, took 1/10 the amount of code, and was implemented 10x faster**.

The language is not just great for data science — but also modeling, ETL, visualizations, package control/version management, machine learning, string manipulation, web-backends, and many other use cases.

¹Python first appeared in 1990. R is an implementation of S, which was created in 1976, though depending on when you want to place the start of an independent R project varies (1993, 1995, and 2000 are alternate dates). The history of these languages is long and substantial changes have occurred since these dates.

²Aviva Case Study

2. Why use Julia?

Julia is well suited for financial modeling work: easy to read and write and very performant.

💡 Tip

The **two language problem** is a term describing processes and teams that separate “domain expertise” coding from “production” coding. This isn’t always a “problem”, but the “two language problem” describes the scenario where this arises not out of intention but out of the necessity of dealing with limitations of the programming languages used. The most common combination is when the domain experts utilize Python, while the quants or developers write C++. This arises because the productive, high level language hits a barrier in terms of speed, efficiency, and robustness. Then, as a necessary step to achieve the end goals of the business, the domain experts hand off the logic to be re-implemented into the lower level language. Not only does this effectively limit the architecture, it essentially defines a required staffing model which may introduce a lot of cost and redundancy in expertise. Julia solves this to a large extent, allowing for a high level, productive language to be very fast.

A similar, related dichotomy is the **two culture problem** wherein domain experts (e.g. financial analysts) exist in a different sphere from developers. This manifests in many ways, such as restricting the tools that each group is permitted to use (e.g. Excel for domain experts, codebases and Git for developers). This is less of a technical problem and more of a social one. However, Julia is also one of the better languages in this regard, because much of the associated tooling is made as straightforward as possible (e.g. packaging, distribution, workflows, etc.). See Chapter 12 and Chapter 21 through Chapter 24 for more on this.

2.1. Julia and This Book

Julia is introduced insofar that a basic understanding is necessary to illustrate certain concepts. Julia is ideal in this context, because it is generally straightforward and concise, allowing the presented idea to have the spotlight (as opposed to language boiler plate or obtuse keywords and variables). The point of structuring the book like this is to allow us to introduce a wide variety of computer science concepts to the financial professional, *not* to introduce Julia as a programming language (there are many other resources which do that just fine).

This chapter seeks to motivate to the skeptical professional why we choose Julia for teaching and for work. Then, in Chapter 5 we provide an introduction to core language concepts and syntax. After this chapter, the content is focused on illustrating a number of key concepts, with Julia taking a secondary role, serving simply as a backdrop. It’s

2.2. Expressiveness and Syntax

not until Chapter 21 where Julia regains the spotlight and we discuss particulars which generally matter only to those heavily using Julia more vigorously.

2.2. Expressiveness and Syntax

Expressiveness is the *manner in which* and *scope of* ideas and concepts that can be represented in a programming language. **Syntax** refers to how the code *looks* on the screen and its readability.

In a language with high expressiveness and pleasant syntax, you:

- Go from idea in your head to final product faster.
- Encapsulate concepts naturally and write concise functions.
- Compose functions and data naturally.
- Focus on the end-goal instead of fighting the tools.

Expressiveness can be hard to explain, but perhaps two short examples will illustrate.

2.2.1. Example 1: Retention Analysis

This is a really simple example relating `Cessions`, `Policys`, and `Lives` to do simple retention analysis. Retention is a measure of how much risk an insurance company holds on a policy after it's own reinsurance risk transfer (ceded amount of coverage are called "cessions").

First, let's define our data:

```
# Define our data structures
struct Life
    policies
end

struct Policy
    face
    cessions
end

struct Cession
    ceded
end
```

Now to calculate amounts retained. First, let's say what retention means for a Policy:

2. Why use Julia?

```
# define retention
function retained(pol::Policy)
    pol.face - sum(cession.ceded for cession in pol.cessions)
end

retained (generic function with 1 method)
```

And then what retention means for a Life:

```
function retained(l::Life)
    sum(retained(policy) for policy in life.policies)
end
```

```
retained (generic function with 2 methods)
```

It's almost exactly how you'd specify it English. No joins, no boilerplate, no fiddling with complicated syntax. You can express ideas and concepts the way that you think of them, not, for example, as a series of dataframe joins or as row/column coordinates on a spreadsheet.

We defined `retained` and adapted it to mean related, but different things depending on the specific context. That is, we didn't have to define `retained_life(...)` and `retained_pol(...)` because Julia can **dispatch** based on what you give it (this is a more powerful, generalized version of method dispatch commonly used in object-oriented programming, see Chapter 7 for more).

Let's use the above code in practice then.

```
# create two policies with two and one cessions respectively
pol_1 = Policy(1000, [Cession(100), Cession(500)])
pol_2 = Policy(2500, [Cession(1000)])

# create a life, which has the two policies
life = Life([pol_1, pol_2])

Life(Policy[Policy(1000, Cession[Cession(100), Cession(500)]), Policy(2500, Cession[Cession(1000)])], 400)

retained(pol_1)

retained(life)
```

```
1900
```

And for the last trick, something called “broadcasting”, which automatically vectorizes any function you write, no need to write loops or create if statements to handle a single vs repeated case:

```
retained.(life.policies) # retained amount for each policy
```

```
2-element Vector{Int64}:
```

```
400  
1500
```

2.2.2. Example 2: Random Sampling

As another motivating example showcasing multiple dispatch, here's random sampling in Julia, R, and Python.

We generate 100:

- Uniform random numbers
- Standard normal random numbers
- Bernoulli random number
- Random samples with a given set

Table 2.1.: A comparison of random outcome generation in Julia, R, and Python.

Julia	R	Python
using Distributions	runif(100) rnorm(100)	import scipy.stats as sps import numpy as np
rand(100)	rbern(100, 0.5)	
rand(Normal(), 100)	sample(c("Preferred","Standard"),	sps.uniform.rvs(size=100)
rand(Bernoulli(0.5), 100)	100, replace=TRUE)	sps.norm.rvs(size=100)
rand(["Preferred","Standard"], 100)		sps.bernoulli.rvs(p=0.5,size=100)
		np.random.choice(["Preferred","Standard"], size=100)

By understanding the different types of things passed to rand(), it maintains the same syntax across a variety of different scenarios. We could define rand(Cession) and have it generate a random Cession like we used above.

2. Why use Julia?

2.3. The Speed

As stated in the journal Nature, “Come for the Syntax, Stay for the Speed”.

Earlier we described Aviva’s Solvency II compliance modeling, which ran 1000x faster than the prior vendor solution mentioned earlier: what does it mean to be 1000x faster at something? It’s the difference between something taking 10 seconds instead of 3 hours — or 1 hour instead of 42 days.

With this difference in speed, you would be able to complete existing processes much faster, or extend the analysis further. This speed could allow you to do new things: a stochastic analysis of life-level claims, machine learning with your experience data, or perform much more frequent valuation.

Here’s a real example, comparing the runtime to calculate the price of a vanilla European call option using the Black-Scholes-Merton formula, as well as the associated code for each. Here’s the mathematical formula we are using:

$$\text{Call}(S_t, t) = N(d_1)S_t - N(d_2)Ke^{-r(T-t)}d_1 = \frac{1}{\sigma\sqrt{T-t}} \left[\ln\left(\frac{S_t}{K}\right) + \left(r + \frac{\sigma^2}{2}\right)(T-t) \right] d_2 = d_1 - \sigma\sqrt{T-t}$$

using Distributions

```
function d1(S,K,τ,r,σ)
    (log(S/K) + (r + σ^2/2) * τ) / (σ * √(τ))
end

function d2(S,K,τ,r,σ)
    d1(S,K,τ,r,σ) - σ * √(τ)
end

function Call(S,K,τ,r,σ)
    N(x) = cdf(Normal(),x)
    d1 = d1(S,K,τ,r,σ)
    d2 = d2(S,K,τ,r,σ)
    return N(d1)*S - N(d2) * K * exp(-r*τ)
end

S,K,τ,σ,r = 300, 250, 1, 0.15, 0.03

Call(S,K,τ,r,σ) # 58.81976813699322

from scipy import stats
import math
```

```

def d1(S,K,τ,r,σ):
    return (math.log(S/K) + (r + σ**2/2) * τ) / (σ * math.sqrt(τ))

def d2(S,K,τ,r,σ):
    return d1(S,K,τ,r,σ) - σ * math.sqrt(τ)

def Call(S,K,τ,r,σ):
    N = lambda x: stats.norm().cdf(x)
    d_1 = d1(S,K,τ,r,σ)
    d_2 = d2(S,K,τ,r,σ)
    return N(d_1)*S - N(d_2) * K * math.exp(-r*τ)

S = 300
K = 250
τ = 1
σ = 0.15
r = 0.03

Call(S,K,τ,r,σ) # 58.81976813699322

d1<- function(S,K,t,r,sig) {
  ans <- (log(S/K) + (r + sig^2/2)*t) / (sig*sqrt(t))
  return(ans)
}

d2 <- function(S,K,t,r,sig) {
  return(d1(S,K,t,r,sig) - sig*sqrt(t))
}

Call <- function(S,K,t,r,sig) {
  d_1 <- d1(S,K,t,r,sig)
  d_2 <- d2(S,K,t,r,sig)
  return(S*pnorm(d_1) - K*exp(-r*t)*pnorm(d_2))
}
S <- 300
K <- 250
t <- 1
r <- 0.03
sig <- 0.15

Call(S,K,t,r,sig) # 58.81977

```

We find in Table 2.2 that despite the syntactic similarity, Julia is much faster than the

2. Why use Julia?

other two.

Table 2.2.: Julia is nearly 20,000 times faster than Python, and two orders of magnitude faster than R.

Language	Median (<i>nanoseconds</i>)	Mean (<i>nanoseconds</i>)	Relative Mean
Python	<i>not calculated by benchmarking library</i>	817000.0	19926.0
R	3649.0	3855.2	92.7
Julia	41.0	41.6	1.0

2.3.1. Development Speed

Speed is not just great for improvement in production processes. During development, it's really helpful too. When building something, the faster feedback loop allows for more productive development. The build, test, fix, iteration cycle goes faster this way.

Admittedly, most workflows don't see a 1000x speedup, but 10x to 1000x is a very common range of speed differences vs R or Python or MATLAB.

i Note

Sometimes you will see less of a speed difference; R and Python have already circumvented this and written much core code in low-level languages. This is an example of what's called the "two-language" problem where the language productive to write in isn't very fast. For example, more than half of R packages use C/C++/Fortran and core packages in Python like Pandas, PyTorch, NumPy, SciPy, etc. do this too.

Within the bounds of the optimized R/Python libraries, you can leverage this work. Extending it can be difficult: what if you have a custom retention management system running on millions of policies every night? In technical terms, libraries like NumPy are not able to handle custom data types, and instead limit use to pre-built types within the library. In contrast, all types in Julia are effectively equal, even ones that you might create yourself.

Julia packages you are using are almost always written in pure Julia: you can see what's going on, learn from them, or even contribute a package of your own!

2.4. More of Julia's benefits

Julia is easy to write, learn, and be productive in:

2.5. Tradeoffs when Using Julia

- It's free and open-source
 - Very permissive licenses, facilitating the use in commercial environments (same with most packages)
- Large and growing set of available packages
- Write how you like because it's multi-paradigm: vector-izable (R), object-oriented (Python), functional (Lisp), or detail-oriented (C)
- Built-in package manager, documentation, and testing-library
- Jupyter Notebook support (it's in the name! **Julia-Python-R**)
- Many small, nice things that add up:
 - Unicode characters like α or β
 - Nice display of arrays
 - Simple anonymous function syntax
 - Wide range of text editor support
 - First-class support for missing values across the entire language
 - Literate programming support (like R-Markdown)
- Built-in Dates package that makes working with dates pleasant
- Ability to directly call and use R and Python code/packages with the `PythonCall.jl` and `RCall` packages
- Error messages are helpful and tell you *what line* the error came from, not just the type of error
- Debugger functionality so you can step through your code line by line

For power-users, advanced features are easily accessible: parallel programming, broadcasting, types, interfaces, metaprogramming, and more.

These are some of the things that make Julia one of the world's most loved languages on the StackOverflow Developer Survey.

In addition to the liberal licensing mentioned above, there are professional products from organizations like JuliaHub that provide hands-on support, training, IT governance solutions, behind-the-firewall package management, and deployment/scaling assistance.

2.5. Tradeoffs when Using Julia

2.5.1. Just-Ahead-of-Time Compilation

Julia is fast because it's compiled, unlike R and Python where (loosely speaking) the computer just reads one line at a time. Julia compiles code "just-in-time": right before you use a function for the first time, it will take a moment to pre-process the code section for the machine. Subsequent calls don't need to be re-compiled and are very fast.

2. Why use Julia?

A hypothetical example: running 10,000 stochastic projections where Julia needs to pre-compile but then runs each 10x faster:

- Julia runs in 2 minutes: the first projection takes 1 second to compile and run, but each 9,999 remaining projections only take 10ms.
- Python runs in 17 minutes: 100ms of a second for each computation.

Typically, the compilation is very fast (milliseconds), but in the most complicated cases it can be several seconds. One of these is the “time-to-first-plot” issue because it’s the most common one users encounter: super-flexible plotting libraries have a lot of things to pre-compile. So in the case of plotting, it can take several seconds to display the first plot after starting Julia, but then it’s remarkably quick and easy to create an animation of your model results. The time-to-first plot is a solvable problem that’s receiving a lot of attention from the core developers and will get better with future Julia releases.

For users working with a lot of data or complex calculations (like actuaries!), the runtime speedup is worth a few seconds at the start.

2.5.2. Static Binaries

Static binaries are self contained executable programs which can run very specific bits of code. Simple programs which can compile down to small (in terms of size on disk) binaries which accomplish just their pre-programmed tasks. Another use case for this is to create shared libraries which could be called from other languages (this can already be done, but again requires bundling the runtime).

Julia’s dynamic nature means that it needs to include the supporting infrastructure in order to run general code, similar to how running Python code needs to be bundled with the Python runtime.

Development is happening at the language level which would allow Julia to be compiled to a smaller, more static set of features for use in environments which are memory constrained and can’t bundle a supporting runtime.

2.6. Package Ecosystem

Using packages as dependencies in your project is assisted by Julia’ bundled package manager.

For each project, you can track the exact set of dependencies and replicate the code/process on another machine or another time. In R or Python, dependency management is notoriously difficult and it’s one of the things that the Julia creators wanted to fix from the start.

2.7. Tools in Your Toolbox

There are thousands of publicly available packages already published. It's also straightforward to share privately, such as proprietary packages hosted internally behind a firewall.

Another powerful aspect of the package ecosystem is that due to the language design, packages can be combined/extended in ways that are difficult for other common languages. This means that Julia packages often interoperable without any additional coordination.

For example, packages that operate on data tables work without issue together in Julia. In R/Python, many features tend to come bundled in a giant singular package like Python's Pandas which has Input/Output, Date manipulation, plotting, resampling, and more. There's a new Consortium for Python Data API Standards which seeks to harmonize the different packages in Python to make them more consistent (R's Tidyverse plays a similar role in coordinating their subset of the package ecosystem).

In Julia, packages tend to be more plug-and-play. For example, every time you want to load a CSV you might not want to transform the data into a dataframe (maybe you want a matrix or a plot instead). To load data into a dataframe, in Julia the practice is to use both the CSV and DataFrames packages, which help separate concerns. Some users may prefer the Python/R approach of less modular but more all-inclusive packages.

2.7. Tools in Your Toolbox

Looking at other great tools like R and Python, it can be difficult to summarize a single reason to motivate a switch to Julia, but hopefully we have sufficiently piqued your interest and we can turn to introducing important concepts.

That said, Julia shouldn't be the only tool in your tool-kit. SQL will remain an important way to interact with databases. R and Python aren't going anywhere in the short term and will always offer a different perspective on things!

Being a productive financial profession means being proficient in the language of computers so that you could build and implement great things. In a large way, the choice of tools and paradigms shape your focus. Productivity is one aspect, expressiveness is another, speed one more. There are many reasons to think about what tools you use and trying out different ones is probably the best way to find what works best for you.

Part II.

Foundations: Effective Financial Modeling

“The map is not the territory.” – Alfred Korzybski (1933)

Welcome to the heart of our exploration of financial modeling. Over the next two chapters, we’ll lay out not only the conceptual foundations that shape how and why we model financial phenomena, but also the practical approaches and data management techniques that make these models work in real-world settings. This chapter is not the technical foundations of financial modeling - instead we focus on the “how” and “why” of modeling in a professional setting.

First, we’ll focus on why financial models matter and what sets them apart from other forms of modeling. This discussion moves from the philosophical—how simplified representations of reality can still provide powerful insights—right down to the practical concerns of “small world” assumptions and “large world” complexities. You’ll discover how models can exist purely for predictive accuracy, even when they don’t fully capture the underlying causal mechanisms, and why this can be both a strength and a danger. Along the way, we’ll unpack the attributes of effective modelers and explore how curiosity and rigor create more resilient analytical tools.

Then, armed with this solid conceptual footing, we’ll turn to the practicalities: selecting the right modeling approach, handling messy data, and navigating trade-offs such as complexity vs. interpretability. You’ll learn about building processes that prepare, transform, and feed data into models in ways that ensure results remain transparent and reliable. We’ll also explore techniques for balancing simulation, machine learning, and statistical models based on your project’s unique requirements.

Think of these chapters as two sides of the same coin: if the first chapter sets the stage by clarifying ‘why’ and ‘what’ we need to model, the second chapter equips you with ‘how’. By the end, you’ll have an integrated perspective: broad conceptual principles for designing a meaningful model and the hands-on skills to bring that model to life.

3. Elements of Financial Modeling

Yun-Tien Lee and Alec Loudenback

“Truth ... is much too complicated to allow anything but approximations” -
John von Neumann

3.1. In this Chapter

We explain what constitutes a financial model and what are common uses of a model. We explore the key attributes of models, discuss different modeling approaches and their trade-offs, and some philosophy around what models are and how they capture important dynamics about the world.

3.2. What is a Model?

A **model** is a simplified representation of reality designed to help us understand, analyze, and make predictions about complex systems. In finance, models distill the intricate web of market behaviors, economic factors, and financial instruments into tractable mathematical and computational components. We may build models for a variety of reasons, as listed in Table 3.1.

Table 3.1.: The REDCAPE model use framework, from “The Model Thinker” by Scott Page.

Use	Description
Reason	To identify conditions and deduce logical implications.
Explain	To provide (testable) explanations for empirical phenomena.
Design	To choose features of institutions, policies, and rules.
Communicate	To relate knowledge and understandings.
Act	To guide policy choices and strategic actions.
Predict	To make numerical and categorical predictions of future and unknown phenomena.
Explore	To investigate possibilities and hypotheticals.

3. Elements of Financial Modeling

For example, say we want to simulate the returns for the stocks in our retirement portfolio. It would be impossible to try to build a model which would capture all of the individual people working jobs and making decisions, weather events that damage property, political machinations, etc. Instead, we try to capture certain fundamental characteristics. For example, it is common to model equity returns as cumulative pluses and minuses from random movements where those movements have certain theoretical or historical characteristics.

Whether we are using this model of equity returns to estimate available retirement income or replicate an exotic option price, a key aspect of the model is the **assumptions** used therein. For the retirement income scenario we might *assume* a healthy eight percent return on stocks and conclude that such a return will be sufficient to retire at age 53. Alternatively, we may assume that future returns will follow a stochastic path with a certain distribution of volatility and drift. These two assumption sets will produce **output** - results from our model that must be inspected, questioned, and understood in the context of the “small world” of the model’s mechanistic workings. Lastly, to be effective practitioners we must be able to contextualize the “small world” results within the “large world” that exists around us.

3.2.1. “Small world” vs “Large world”

The distinction between the modeled “small world” and real life “large world” modeling is fundamental to understanding model limitations:

- **Small World:** The constrained, well-defined environment within your model where all rules and relationships are explicitly specified.
- **Large World:** The complex, real-world environment where your model must operate, including factors not captured in your assumptions.

Say that our model is one that discounts a fixed set of future cashflows using the US Treasury rate curve. If I run my model using current rates today, and then re-run my model tomorrow with the same future cashflows and the present value of those cashflows has increased by 5%, I may ask: “why has the result has changed so much in such a short period of time?”

In the “small”, mechanistic world of the model I may be able to see that the discount rates have fallen substantially. The small world answer is that the inputs have changed which produced a mechanical change in the output. The large world answer may be that the Federal Reserve lowered the Federal Funds Rate to prevent the economy from entering a deflationary recession.

Of course, we can’t completely explain the relation between our model and the real world (otherwise we could capture that relationship in our model!). An effective practitioner will always try to look up from the immediate work and take stock of how the world at large *is* or *is not* reflected in the model.

3.3. What is a Financial Model?

3.3. What is a *Financial* Model?

Financial models are those used extensively to ascertain better understanding of complex contracts, perform scenario analysis, and inform market participants' decisions related to perceived value (and therefore price). It can't be quantified directly, but it is likely not an exaggeration that many billions of dollars is transacted each day as a result of decisions made from the output of financial models.

Most financial models can be characterized with a focus on the first or both of:

1. Attempting to project pattern of cashflows or obligations at future timepoints
2. Reducing the projected obligations into a current value

Examples of this:

- Projecting a retiree's savings through time (1), and determining how much they should be saving today for their retirement goal (2)
- Projecting the obligation of an exotic option across different potential paths (1), and determining the premium for that option (2)
- Determining a range of outcomes (1) for risk modeling purposes, and translating those outcomes into a risk or capitalization level (2)

Models are sometimes taken a step further, such as transforming the underlying **economic view** into an accounting or regulatory view (such as representing associated debits and credits, capital requirements, or associated intangible, capitalized balances). In other words, many models go further than simply modeling anticipated cashflows by building representations of entities (companies or portfolios) which are associated with the cashflows.

3.3.1. Difference from Data Science

We should also distinguish a financial model from a purely statistical model, where often the inputs and output data are known and the intention is to estimate relationships between variables (example: linear regressions). That said, a financial model may have statistical components and many aspects of modeling are shared between the two kinds.

In this book, we focus on the practice of financial modeling in the ways that it is distinct from statistics and data science. To delineate the practices:

- **Statistics, or Statistical Modeling** is the practice of applying procedures and probabilistic reasoning to data in order to determine relationships between things or to predict outcomes.

3. Elements of Financial Modeling

- **Data Science** includes statistical modeling but also incorporates the art and science of good data hygiene, data pipelines and pre-processing, and more programming than a pure statistician usually uses.

Financial Modeling is similar in the goal of modeling complex relationships or making predictions (a modeled price of an asset is simply a prediction of what its value is), however, it differs from data science in a few ways:

- Financial modeling generally incorporates a good deal of theory that assumes certain behaviors/relationships, instead of trying to infer those relationships from the data.
- Financial modeling generally contains more unique ‘objects’ than a statistical model. The latter may have derived numerical relationships between data, however a financial model would have things like the concept of a company or portfolio, or sets of individually identifiable assets, or distinct actors in a system.
- Financial modeling often involves a lot more simulation and hypothesizing, while data science is focused on drawing conclusions from what data has already been observed.

Nonetheless, there is substantial overlap in practice. For example, the assumption in a financial model (volatility, economic conditions, etc.) may be derived statistically from observed data. Given the overlap in topics, statistical content is sometimes covered in this book (including a from-the-ground-up view of modern Bayesian approaches in Chapter 14).

3.4. Key Considerations for a Model

When creating a model, whether a data model, a conceptual model, or any other type, certain key considerations are generally important to include to ensure it is effective and useful. Some essential considerations include:

Consideration	Description
Objective	Clearly define what the model aims to achieve.
Boundaries	Specify the limits and constraints of the model to avoid scope creep.
Variables	Identify and define all variables involved in the model.
Parameters	Include constants or coefficients that influence the variables.
Dependencies	Describe how variables interact with each other.
Relationships	Detail the connections between different components of the model.
Inputs	Specify the data or resources required for the model to function.
Outputs	Define what results or predictions the model produces.
Underlying Assumptions	Document any assumptions made during the model’s development to clarify its limitations and validity.

3.5. Predictive versus Explanatory Models

Consideration	Description
Validation Criteria	Outline how the model's accuracy and reliability are tested.
Performance Metrics	Define the metrics used to evaluate the model's performance.
Scalability	Ensure the model can handle increased data or complexity if needed.
Adaptability	Allow for adjustments or updates as new information or requirements arise.
Documentation	Provide comprehensive documentation explaining how the model works, including algorithms, data sources, and methods.
Transparency	Make the model's workings understandable to stakeholders or users.
User Interface	Design an intuitive interface if the model is interactive.
Ease of Use	Ensure that the model is user-friendly and accessible to its intended audience.
Ethics	Address any ethical concerns related to the model's application or impact.
Regulations	Ensure compliance with relevant laws and regulations.

Including these attributes helps create a robust, reliable, and practical model that effectively serves its intended purpose.

3.5. Predictive versus Explanatory Models

Given a set of inputs, our model will generate an output and we are generally interested in its accuracy. *The model need not have a realistic mechanism for how the world works.* That is, we may primarily be interested in accurately calculating an output value without the model having any scientific, explanatory power of how different parts of the real-world system interact.

3.5.1. A Historical Example

Consider the classic underdog story where Copernicus overthrew the status quo when he proposed (correctly) that the earth orbited the sun instead of the other way around¹.

The existing Ptolemy model used a geocentric view of the solar system in which the planets and sun orbited the Earth in perfect circles with an epicycle used to explain retrograde motion (as seen in Figure 3.1). Retrograde motion is the term used to describe

¹Prof. Richard Fitzpatrick has excellent coverage of the associated mathematics and implications in "A Modern Almagest": <https://farside.ph.utexas.edu/books/Syntaxis/Almagest.html>

3. Elements of Financial Modeling

the apparent, temporarily reversed motion of a planet as viewed from Earth when the Earth is overtaking the other planet in orbit around the sun. Despite a fundamentally flawed underlying theory, the geocentric model was accurate enough to match the observational data for the position of the planets in the sky.

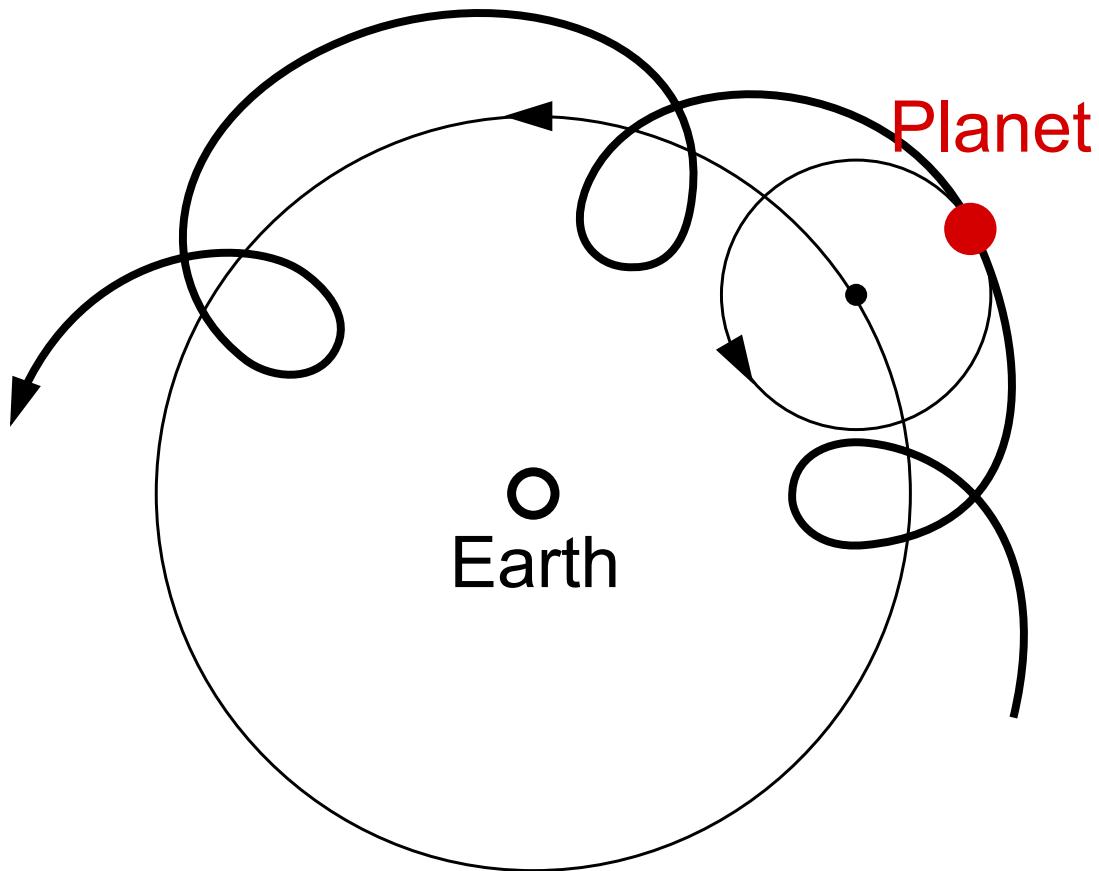


Figure 3.1.: In the Ptolemy solar model, the retrograde motion of the planets was explained by adding an epicycle to the circular orbit around the earth.

Famously, Copernicus came along and said that the sun, not the Earth, should be at the center (a heliocentric model). Earth revolves around the sun! Today, we know this to be a much better description of reality than one in which the Earth arrogantly sits at the center of the universe. However the model was actually slightly *less* accurate in predicting the apparent position of the planets (to the limits of observational precision at the time)! Why would this be?

First, the Copernican proposal still used perfectly circular orbits with an epicycle adjust-

3.5. Predictive versus Explanatory Models

ment, which we know today to be inaccurate (versus an elliptical orbit consistent with the theory of gravity). *Despite being more scientifically correct, it was still not the complete picture.*

Second, the geocentric model was already very accurate because it was essentially a Taylor-series approximation which described, to sufficient observational accuracy, the apparent position of the planet relative to the Earth. *The heliocentric model was effectively a re-parameterization of the orbital approximation.*

Third, we have considered a limited criteria for which we are evaluating the model for accuracy, namely apparent position of the planets. *It's not until we contemplate other observational data that the Copernican model would demonstrate greater modeling accuracy:* apparent brightness of the planets as they undergo retrograde motion and angular relationship of the planets to the sun.

For modelers today, this demonstrates a few things to keep in mind:

1. Predictive models need not have a scientific, causal structure to make accurate predictions.
2. It is difficult to capture the complete scientific inter-relationships of a system and much care and thought needs to be given in what aspects are included in our model.
3. We should look at, or seek out, additional data that is related to our model because we may accurately overfit to one outcome while achieving an increasingly poor fit to other related variables.

Striving to better understand the world is a *good thing to do* but trying to include more components into the model is not always going to advance a given goal.

3.5.2. Examples in the Financial Context

3.5.2.1. Home Prices

American home prices which have a strong degree of seasonality and have the strongest prices around April of each year. We may find that including a simple oscillating term in our model captures the variability in prices *better* than if we tried to imperfectly capture the true market dynamics of home sales: supply and demand curves varying by personal (job bonus payment timing, school calendars), local (new homes built, company relocation), and national (monetary policy, tax incentives for home-ownership) factors. In other words, one could likely predict a stable pattern like this with a model that contains a simple sinusoidal periodic component. One could likely spend months trying to build a more scientific model and not achieve as good of fit, *even though the latter tries to be more conceptually accurate.*

3. Elements of Financial Modeling

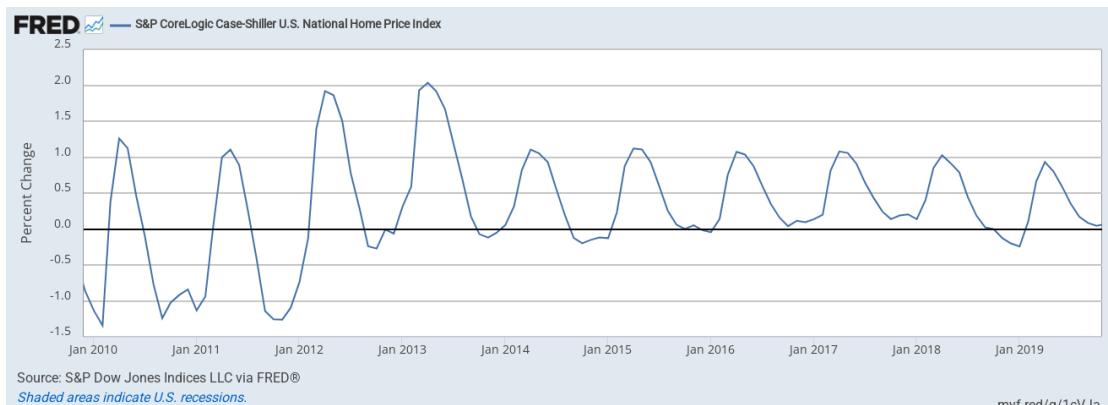


Figure 3.2.: House prices peak around April of each year.

3.5.2.2. Replicating Portfolio

Another example in the financial modeling realm: in attempting to value a portfolio of insurance contracts a **replicating portfolio** of hypothetical assets will sometimes be constructed². The point of this is to create a basket of assets that can be more quickly (minutes to hours) valued in response to changing market conditions than it would take to run the actuarial model (hours to days). This is an example where the basket of assets has no ability to explain why the projected cashflows are what they are - but retains strong predictive accuracy.

3.6. Types of Models

Different modeling approaches come with their own sets of trade-offs. Common modeling approaches, and the inherent trade-offs, may include:

Model				
Type	Description	Examples	Trade-offs	Assumptions
Statistical models	Use mathematical relationships to describe data.	Linear regression, logistic regression	Simplicity vs. Accuracy: Often simpler and more interpretable but may not capture complex relationships.	Typically rely on assumptions like linearity and normality that may not always hold true.

²See, e.g., SOA Investment Symposium March 2010. *Replicating Portfolios in the Insurance Industry* (Curt Burmeister Mike Dorsel Patricia Matson)

3.6. Types of Models

Model				
Type	Description	Examples	Trade-offs	Assumptions
Machine Learning Models	Learn patterns from data using algorithms.	Decision trees, random forests, neural networks	Complexity vs. Interpretability: Capture complex patterns but are often less interpretable. Overfitting: Risk of overfitting, requiring careful validation and tuning. Data Requirements: Require large datasets.	Performance can degrade with limited or noisy data.
Probabilistic Models	List probability distributions to model uncertainty and relationships.	Bayesian networks, probabilistic graphical models	Flexibility vs. Computational Complexity: Handle uncertainty and complex relationships but require sophisticated computations.	May require assumptions about the nature of probability distributions and dependencies.
Simulation Models	Use computational models to simulate complex systems or scenarios.	Monte Carlo simulations, agent-based models	Accuracy vs. Computational Expense: Can be computationally expensive and time-consuming. Detail vs. Generalization: High-fidelity simulations may be overkill for simpler problems.	May require simplifications or assumptions for computational feasibility.
Theoretical Models	Based on theoretical principles to explain phenomena.	Economic models, physical models	Precision vs. Practicality: Provide foundational understanding but may rely on idealizations. Applicability: Highly accurate in specific contexts but less so in broader situations.	Often rely on idealizations or simplifications.

3. Elements of Financial Modeling

Model				
Type	Description	Examples	Trade-offs	Assumptions
Empirical Models	Use historical data to predict future outcomes.	Time series forecasting, econometric models	Data Dependence vs. Predictive Power: Rely heavily on historical data and may not perform well if patterns change. Context Sensitivity: Accurate for specific data but may not generalize well.	Performance is heavily dependent on the quality and relevance of historical data.
Hybrid Models	Combine different modeling approaches to leverage their strengths.	Combining statistical and machine learning approaches	Complexity vs. Versatility: Aim to leverage strengths of different approaches but can be complex to design. Integration Challenges: May present challenges in integration and consistency.	May inherit assumptions from combined models.

Summary of Common Trade-offs:

- Complexity vs. Simplicity: More complex models can capture more nuanced details but are harder to understand and manage.
- Accuracy vs. Interpretability: High-accuracy models may be less interpretable, making it harder to understand their decision-making process.
- Data Requirements: Some models require large amounts of data or very specific types of data, which can be a limitation in practice.
- Computational Resources: More sophisticated models or simulations can require significant computational power, which may not always be feasible.

Understanding these trade-offs helps in selecting the most appropriate modeling approach based on the specific needs of the problem at hand and the resources available.

4. The Practice of Financial Modeling

Yun-Tien Lee and Alec Loudenback

“In theory there is no difference between theory and practice. In practice there is.” – Yogi Berra (often attributed)

4.1. Introduction

Having covered what models are and what they accomplish, we turn to the *craft* of modeling: what distinguishes a good model from a bad model; likewise what are attributes of an astute practitioner. Lastly, we cover some more “nuts and bolts” topics such as data handling and good governance practices.

4.2. What makes a good model?

The answer is: *it depends.*

4.2.1. Achieving original purpose

A model is built for a specific set of reasons and therefore we must evaluate a model in terms of achieving that goal. We should not critique a model if we want to use it outside of what it was intended to do. This includes: contents of output and required level of accuracy.

A model may have been created to for scenario analysis to value all assets in a portfolio to within half a percent of a more accurate, but much more computationally expensive model. If we try to add a never-before-seen asset class or use the model to order trades, then we may be extending beyond the design scope of the original model and lose predictive accuracy.

4. The Practice of Financial Modeling

4.2.2. Usability

How easy is it for someone to use? Does it require pages and pages of documentation, weeks of specialized training and an on-call help desk? *All else equal*, it is an indicator of how usable the model is by the amount of support and training required. However, one may sometimes wish to create a highly capable, complex model which is known to require a high amount of experience and expertise. An analogy here might be the cockpit of a small Cessna aircraft versus a fighter jet: the former is a lot simpler and takes less training to master but is also more limited.

Figure 4.1 illustrates this concept and shows that if your goal is very high capability that you may need to expect to develop training materials and support the more complex model. On this view, a better model is one that is able to have a shorter amount of time and experience to achieve the same level of capability.

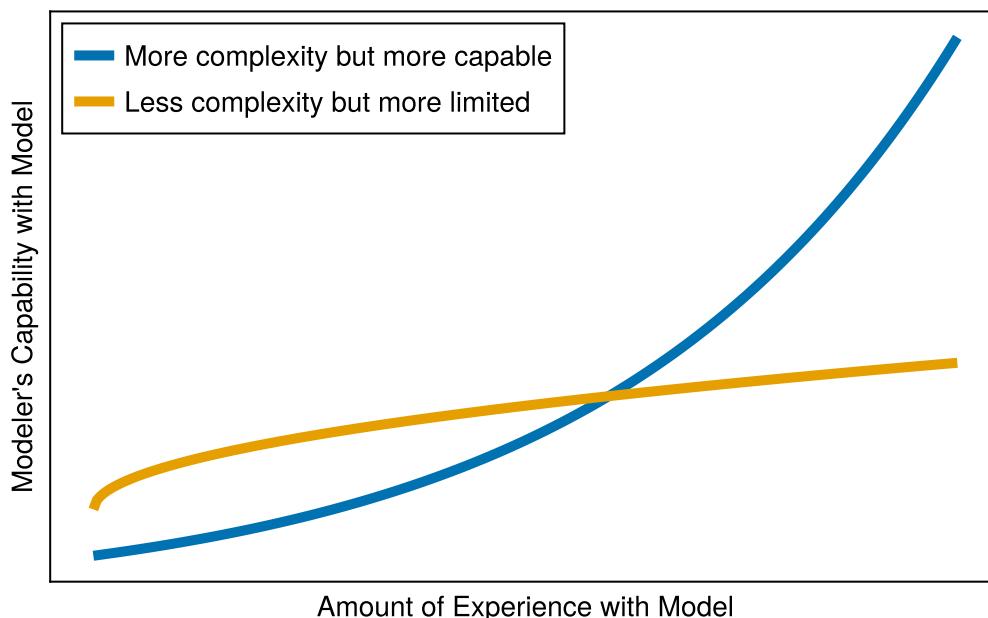


Figure 4.1.: Tradeoff between complexity and capability

4.2.3. Performance

Financial models are generally not used for their awe-inspiring beauty - users are results oriented and the faster a model returns the requested results, the better. Aside from direct computational costs such as server runtime, a shorter model runtime means

4.2. What makes a good model?

that one can iterate faster, test new ideas on the fly, and stay focused on the problem at hand.

Many readers may be familiar with the cadence of (1) try running model overnight, (2) see results failed in the morning, (3) spend day developing, (4) repeat step 1. It is preferred if this cycle can be measured in minutes instead of hours or days.

Of course, requirements must be considered here too: needs for high frequency trading, daily portfolio rebalancing, and quarterly financial reporting models have different requirements when it comes to performance.

4.2.4. Separation of Model Logic and Data

When data is intertwined with business logic it can be difficult to understand, maintain, or adapt a model. Spreadsheets are a common example of where data exists co-mingled with business logic. An alternative which separates data sources from the computations provides for better model service in the future.

4.2.5. Organization of Model Components and Architecture

If model components or data inputs are spread out in a disorganized way, it can lead to usability and maintenance issues. As an example, often times it's incredibly difficult to ascertain a model's operation if inputs are spread out across locations on many spreadsheet tabs. Or if related calculations are performed in multiple locations, or if it's not clear where the line is drawn between calculations performed in the worksheets or in macros.

If logical components or related data are broken out into discrete parts of a model, it becomes easier to reason about model behavior or make modifications. Compartmentalization is an important principle which allows a larger model to remain comprised of simpler components where whole model is greater than the sum of the pieces.

4.2.6. Abstraction of Modeled Systems

At different times we are interested in different **ladder of abstraction**: sometimes we are interested in the small details, but other times we are interested in understanding the behavior of systems at a higher level.

Say we are an insurance company with a portfolio of fixed income assets supporting long term insurance liabilities. We might delineate different levels of abstraction like so:

4. *The Practice of Financial Modeling*

Table 4.1.: An example of the different levels of abstraction when thinking about modeling an insurance company's assets and liabilities.

Item	
More Abstract	Sensitivity of an entire company's solvency position
	Sensitivity of a portfolio of assets
	Behavior over time of an individual contract
More granular	Mechanics of an individual bond or insurance policy

At different times, we are often interested in different aspects of a problem. In general, you start to be able to obtain more insights and a greater understanding of the system when you move up the ladder of abstraction.

In fact, a lot of designing a model is essentially trying to figure out where to put the right abstractions. What is the right level of detail to model this in and what is the right level of detail to expose to other systems?

Let us also distinguish between **vertical abstraction**, as described above, and **horizontal abstraction** which will refer to encapsulating different properties, or mechanics of components of model that effectively exist on the same level of vertical abstraction. For example, both asset and liability mechanics sit at the most granular level in Table 4.1, But it may make sense in our model to separate the data and behavior from each other. If we were to do that, that would be an example of creating horizontal abstraction in service of our overall modeling goals.

This book will introduce powerful, programmatic ways to handle this through things like packages, modules, namespaces, and functions.

4.3. What makes a good modeler?

A model is nothing without its operator, and a skilled practitioner is worth their weight in gold. What elements separate a good modeler from a mediocre modeler?

4.3.1. Domain Expertise

An expert who knows enough about all of the domains that are applicable is crucial. Imagine if someone said let's emulate an architect by having a construction worker and an artist work together. It's all too common for business to attempt to pair a business expert with an information technologist in the same way.

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Unfortunately, this means that there's generally no easy way out of learning enough about finance, actuarial science, computers, and/or programming in order to be an effective modeler.

Also, a word of warning for the financial analysts out there: the computer scientists may find it easier to learn applied financial modeling than the other way around since the tools, techniques, and language of problem solving is already more a more general and flexible skill-set. There's more technologists starting banks than there are financiers starting technology companies.

4.3.2. Model Theory

If it is granted that financial modeling must involve, as the essential part, a building up of modeler's knowledge, the next issue is to characterize that knowledge more explicitly. The modeler's knowledge should be regarded as a theory, in the sense of Ryle's¹ "Concept of the Mind." Very briefly: a person who has or possesses a theory in this sense knows how to do certain things and in addition can support the actual doing with explanations, justifications, and answers to queries, about the model and its results².

A financial model is rarely left in a final state. Regulatory changes, additional mechanics, sensitivity testing, market dynamics, new products, and new systems to interact with force a model to undergo change and development through its entire life. And like a living thing, it must have nurturing caregivers. This metaphor sounds extended, but Naur's point is that unless the model also lives in the heads of its developers then it cannot successfully be maintained through time:

The conclusion seems inescapable that at least with certain kinds of large programs, the continued adaption, modification, and correction of errors in them, is essentially dependent on a certain kind of knowledge possessed by a group of programmers who are closely and continuously connected with them.

Assume that we need to adapt the model to fit a new product. One possessing a high degree of model theory includes:

- the ability to describe the trade-offs between alternate approaches that would accomplish the desired change
- relate the proposed change to the design of the current system and any challenges that will arise as a result of prior design decisions

¹Ryle, G. *The Concept of Mind*. Harmondsworth, England, Penguin, 1963, first published 1949. Applying "Theory Building"

²The idea of "model theory" is adapted from Peter Naur's 1985 essay, "Programming as Theory Building". Indeed, this whole paragraph is only a slightly modified version of Naur's description of theory in the programming context.

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- provide a quantitative estimation for the impact the change will have: runtime, risk metrics, valuation changes, etc.
- Analogize how the system works to themselves or to others
- Describe key limitations that the model has and where it is most divorced from the reality it seeks to represent.

Abstractions and analogies of the system are a critical aspect of model theory, as the human mind cannot retain perfectly precise detail about how the system works in each sub-component. The ability to, at some times, collapse and compartmentalize parts of the model to limit the mental overload while at others recall important implementation details requires training - and is enhanced by learning concepts like those which will be covered in this book.

An example of how the right abstractions (and language describing those abstractions) can be helpful in simplifying the mental load:

Instead of:

The valuation process starts by reading an extract into three tabs of the spreadsheet. A macro loops through the list of policies on the first tab and in column C it gives the name of the applicable statutory valuation ruleset. The ruleset is defined as the combination of (1) the logic in the macro in the "Valuation" VBA module with, (2) the underlying rate tables from the tabs named XXX to ZZZ, along with (3) the additional policy level detail on the second tab. The valuation projection is then run with the current policy values taken from the third tab of the spreadsheet and the resulting reserve (equal to the actuarial present value of claims) is saved and recorded in column J of the first tab. Finally, a pivot table is used to sum up the reserves by different groups.

We could instead design the process so that the following could be said instead:

Policy extracts are parsed into a Policy datatype which contains a subtype ValuationKind indicating the applicable statutory ruleset to apply. From there, we map the valuation function over the set of Policies and perform an additive reduce to determine the total reserve.

There are terminologies and concepts in the second example which we will develop over the course of this section of the book - we don't want to dwell on the details bright now. However, we do want to emphasize that the process itself being able to condensed down to descriptions that are much more meaningful to the understanding of the model is a key differentiator for a code-based model instead of spreadsheets. It is not exaggerating that we could develop a handful of compartmentalized logics such that our primary valuation process described above could look like this in real code:

```
policies = parse(Policy,CSV.File("extract.csv"))
reserve = mapreduce(+,value,policies)
```

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We've abstracted the mechanistic workings of the model into concise and meaningful symbols that not only perform the desired calculations but also make it obvious to an informed but unfamiliar reader what it's doing.

`parse`, `mapreduce`, `+`, `value`, `Policy` are all imbued with meaning - the first three would be understood by any computer scientist by the nature of their training (and is training that this book covers). The last two are unique to our model and have "real world" meaning that our domain expert modeler would understand which analogizes very directly to the way we would suggest implementing the details of `value` or `Policy`. The benefit of this, again, is to provide tools and concepts which let us more easily develop model theory.

4.3.3. Curiosity

By cultivating curiosity, modelers can drive innovation, uncover new insights, and continuously improve their models and understanding of financial systems.

No model, no matter how sophisticated, ever delivers a "final" answer. If anything, a good financial model sparks as many new questions as it answers. This is where the best modelers distinguish themselves: they nurture a healthy skepticism with surface-level explanations.

Take, for instance, the gnawing feeling you get when a model's output seems "off" but you can't quite put your finger on why. The untrained eye might chalk it up to randomness or let it slide, but genuine curiosity won't settle for a hand-wavy excuse. The itch to resolve every weird edge case or apparent contradiction, to ask "what if?" and "why not?" is the spark that propels a practitioner beyond rote calculation into discovery.

In practice, carrying curiosity into modeling means:

- **Taking the time to poke holes in every story your model tells.** If two approaches give wildly different answers for the same scenario, don't sweep that under the rug. Dig until you've either found the bug or learned a new subtlety.
- **Going down rabbit holes.** Sometimes the best model improvements stem from following up on the "trivial" anomaly hiding in the the numerical 'blip' over 12 months. Ask yourself: is there a structural reason, a missing piece of data, or an assumption that needs to be made explicit?
- **Pursuing the "why" behind the numbers.** Instead of blindly running scenarios, become obsessed with the model's behavior. If changing an input slightly has points has an outsized effect elsewhere, dig into the feedback mechanism that causes it.
- **Challenging your own assumptions.** No matter how seasoned you are, ask foundational questions and reconsider the "obvious." There are no dumb questions! You'd be surprised how many "everybody knows that" ideas are actually half-remembered lore.

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- **Learning from surprises.** Whenever the model spits out something bizarre, treat it like an opportunity rather than a headache. Sometimes the oddball result teaches you more about the system than any routine validation could.
- **Trying new techniques and tools to keep your intellectual toolbox diverse.** Look for the overlap between different things, such as recognizing the similarities between different areas of practice, even if it's not your 'specialty'.

The best modelers I've worked with aren't necessarily the flashiest coders or the most fluent in finance. They're just relentless in their quest to not leave loose ends untied.

4.3.4. Rigor

If curiosity is the fuel, rigor is the steering wheel. All of that wandering through the thickets of "why?" needs a reliable process to keep from becoming noise or hand-waving. Rigor is what separates "I think it works" from "Here's why it works, and here are its limits."

When developing a model it's important to ensure that assumptions and parameters are very clear, the methodology is in line with established theory, inappropriate thought has been given to how the model will be used. Additionally one should be mindful of standards of practice. For example, professional actuarial societies have a long list of Actuarial Standards of Practice ("ASOPs"), some of which apply to modeling and the use of data that models ultimately rely on. Regardless of the applicable standards, many of them share these aspects of the best modelers:

- **Document your thinking** as you go. Write it out, whether it's in a code comment, a README, or your own notebook. If you can't explain your logic and your parameters, you probably don't understand them as well as you think.
- **Demand evidence** for your choices—don't just trust your gut or yesterday's industry standard. Check your results against reality, not just an assumed "right answer." This means obsessive test cases, sensitivity checks, and "could we break this?" scenarios.
- **Hold the model to a higher standard than tradition requires.** Don't just meet regulator norms if you can do better—set your own red-lines for quality, accuracy, and reproducibility.
- **Don't hide the warts.** Make uncertainty visible, not hypothetical. Annotate what's based on thin data versus what's on solid ground. Rigor means being honest about what you don't know—or what the model simply can't say.
- **Lean on first principles.** Often times there will be a 'simpler' way to model something, but making explicit all components of an interaction can be illuminating. For example, if you have a complex, multi-leg transaction that 'works like exotic

4.3. What makes a good modeler?

option ABC', don't always rely on that heuristic. Alternatively, model out each leg of the transaction for clarity and confirmation of your understanding.

A bad model can be worse than no model at all. Through rigorous efforts, a minimum standard of quality can be obtained.

4.3.5. Clarity

"Clarity" means never losing sight of the fact your model is only as useful as it is understandable.

- Precise language: Use well-defined terms and avoid ambiguity in communications. If a term has overloaded meanings ("reserve," "duration," "return"), either define it up front or pick a less ambiguous word.
- Spell out your assumptions, not just your output. Make the philosophy and the scaffolding explicit—what did you leave in, what did you leave out, and why?
- Visual communication: Utilize diagrams and visualizations to explain complex concepts. A simple stylized sketch rarely hurts and often helps.
- Audience-appropriate communication: Adjust your explanations depending on whether you're talking to other developers, business stakeholders, or end users.
- Regular review: Periodically update documentation to ensure ongoing clarity and accuracy. If you woke up with amnesia, would the next steps seem obvious?

Clarity is about making your future self—and your colleagues—thankful, not furious, that you were ever given keyboard access.

4.3.6. Humility

The world is complicated in ways we can sometimes describe and never fully anticipate. A humble modeler tries to understand what the model can and cannot claim. And in good faith will try to share the limitations the model has with stakeholders, such as saying something like "we have a lot of data for low-rate environments but rapidly rising environments haven't been observed in the dataset".

Irreducible & reducible (epistemic) uncertainty are critical concepts for a modeler to understand and communicate:

1. Irreducible uncertainty: Also known as aleatoric uncertainty, this refers to the inherent randomness in a system that cannot be reduced by gathering more information.
 - Examples include: future market fluctuations, individual policyholder behavior, or natural disasters.

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2. Reducible (epistemic) uncertainty: This type of uncertainty stems from a lack of knowledge and can potentially be reduced through further study or data collection.
 - Examples include: parameter estimation errors, model specification errors, or data quality issues.

Table 4.2 describes this in more detail. It's not always necessary to describe each of these types of uncertainty for every model but knowing your enemy is the first step in fighting it.

A humble modeler acknowledges these uncertainties and communicates them clearly to stakeholders. This avoids overconfidence in model predictions and keeps one open to new information and alternative perspectives. By maintaining a humble attitude, modelers can build trust with stakeholders and make more informed decisions based on model outputs.

Table 4.2.: In attempting to model an uncertain world, we can be even more granular and specific in discussing sources of that uncertainty. This table summarizes commonly noted kinds of uncertainty that arise, and whether we can reduce the uncertainty by doing better (more data, better data, better models, etc.) or not.

Type of Uncertainty	Key Characteristics	Reducibility	Example
Aleatory (Process) Uncertainty	- Inherent randomness (aka “irreducible uncertainty”) - Cannot be eliminated, even with perfect knowledge	Irreducible	Rolling dice or coin flips; outcome is inherently uncertain despite full knowledge of initial state
Epistemic (Parameter) Uncertainty	- Due to limited data/knowledge (aka “reducible uncertainty”) - Imperfect information or model parameters	Reducible (more data / improved modeling)	Uncertainty in a model’s parameters (e.g., climate sensitivity) that can be refined with more research
Model Structure Uncertainty	- Uncertainty about the correct model or framework - Often considered a special subset of epistemic uncertainty	Partially reducible (better theory/model selection)	Linear vs. nonlinear models in complex systems; risk of omitting key variables or mis-specified dynamics
Deep (Knightian) Uncertainty	- “Unknown unknowns” - Probability distributions themselves are not well-defined or are fundamentally unquantifiable	Not quantifiable (cannot assign probabilities)	Impact of radically new technology on society
Measurement Uncertainty	- Errors in data collection or instrument readings - Systematic biases or random errors in measurement	Partially reducible (improved measurement methods)	Instrument precision limits in experiments; calibration errors in sensor data

Type of Uncertainty	Key Characteristics	Reducibility	Example
Operational Uncertainty	- Uncertainty in implementation/execution - Human error, mechanical failure, or miscommunication in processes	Partially reducible (better training/processes)	Surgical errors, system failures, or incorrect handling of a financial trade order

4.3. What makes a good modeler?

4.3.7. Architecture

Any sufficiently complex project benefits from architectural thinking. Think of your model like a house: if you don't plan the plumbing, you'll have a mess down the line. Data should be separate from the logic and the model itself should not contain any substantial datum itself - instead dynamically load data from appropriate data stores and leave the "model" as the implementation of data *types* and algorithms.

- Modular design: Break complex models into reusable, independent components.
- Separation of concerns: Keep data, logic, and presentation layers distinct for better maintainability.
- Scalability: Design models to handle increasing data volumes and complexity.
- Maintainability: Use version control, stable interfaces, clear documentation, and automated tests.
- Performance optimization: Use efficient data structures and algorithms to enhance model speed.
- Security: Ensure proper data protection and regulatory compliance.

Don't underestimate the value of a well-organized model: it's how you scale from small prototypes to systems you can trust in production.

4.3.8. Planning

When tackling a large problem, it helps to have a well-structured planning process. Specific to building a financial model, one should take steps that include:

1. Clear objectives: Understand the purpose of the model and what questions it needs to answer.
2. Scope definition: Determine the boundaries of the model, including what to include and what to exclude.
3. Data assessment: Identify required data sources, assess data quality, and plan for data preparation.
4. Methodology selection: Choose appropriate modeling techniques based on the problem and available data.
5. Resource allocation: Estimate time and resources needed for model development, testing, and implementation.
6. Stakeholder engagement: Identify key stakeholders and plan for their involvement throughout the modeling process.
7. Risk assessment: Anticipate potential challenges and develop mitigation strategies.
8. Timeline development: Create a realistic timeline with key milestones and deliverables.

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9. Documentation strategy: Plan for comprehensive documentation of assumptions, methodologies, and limitations.
10. Validation and testing approach: Outline strategies for model validation and testing to ensure reliability.
11. Implementation and maintenance plan: Who will have responsibility after the model achieves its initial objectives?

Time invested at the planning stage often pays dividends through shorter model build times, fewer errors, and clarity from stakeholders at the start of the project. Additionally, it's often easier to make changes to a well-planned project halfway through since the necessary accommodations are more clearly defined.

4.3.9. Essential Tools and Skills

An experienced professional is aware of a number of approaches that can be used in solving a problem. From heuristics that are able to be calculated on a napkin to complex economic models, the ability to draw on a wide tool set allows a practitioner to find the right solution for a given problem. It is the intention of this book to enumerate a number of additional approaches that may prove useful in practice. This includes both soft and hard skills, such as those in Table 4.3

Table 4.3.: A variety of skills have their place in the proficient financial modeler's tool-belt.

Category	Examples
Diverse Modeling Techniques	<ul style="list-style-type: none">• Statistical methods (e.g. regression, time series analysis, machine learning)• Optimization techniques (e.g. linear, non-linear, black-box)• Simulation methods (e.g. monte-carlo, agent-based, seriatim)
Software Proficiency	<ul style="list-style-type: none">• Programming languages• Database and data handling• Proprietary tools (e.g. Bloomberg)
Financial Theory	<ul style="list-style-type: none">• Asset pricing• Portfolio theory• Risk Management frameworks• Numerical methods and algorithms• Bayesian inference• Stochastic calculus
Quantitative techniques	
Soft Skills	<ul style="list-style-type: none">• Verbal and written communication• Stakeholder engagement• Project Management

4.4. Feeding The Model

If the soul of the model is its logic, then the lifeblood is its data. In practice, a model's fate is often sealed not in the sophistication of its algorithms, but in the quality of the data it consumes. Even the most elegant model is helpless in the face of stale, sloppy, or misunderstood inputs.

4.4.1. “Garbage In, Garbage Out”

Every experienced modeler has a story where a subtle data quirk led to a dramatic miscalculation—a column header shifted by one, a stale price feed, or a single outlier that quietly cascaded into a million-dollar mistake. The lesson: treat the data with every bit as much skepticism (and care) as you give the model itself.

4.4.2. A Modeler’s Data Instincts

Rather than thinking of data handling as a rigid checklist, approach it as a series of habits and questions:

- **Know Your Sources.** Where did this data come from? Who collected it, and how? Is it raw, or has someone already “cleaned” it in ways you need to understand (or undo)? Data provenance is not a formality—it’s the first step in understanding what can go wrong.
- **Trust, But Verify.** Never take a dataset at face value, even if it comes from a trusted system. Run summary statistics. Plot the distributions. Check for the bizarre and the mundane: are dates reasonable, units consistent, and identifiers unique?
- **Expect Messiness.** Real-world data is rarely pristine. Missing values, odd encodings, duplicated rows, and outliers are the norm, not the exception. The best modelers are part detective, part janitor: they track down wonky values, document their triage decisions, and know when to escalate a data quality concern upstream.
- **Feature Engineering Is Judgment, Not Magic.** Choosing which fields to keep, combine, or discard is where domain expertise shines. Sometimes a new ratio or flag column, born from your unique understanding of the business, makes all the difference. Beware of “kitchen sink” modeling—too many features can obscure, rather than reveal, the truth.
- **Be Wary of Temporal Traps.** Mixing data from different time periods, or accidentally leaking future information into a model (a classic error), can invalidate results without any warning sign. When in doubt, plot your data against time and look for jumps, gaps, or trends that defy explanation.

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- **Keep Data and Logic Separate.** As harped on earlier: don't hard-code data into the model. Keep sources external, interfaces clean, and ingest paths well documented. If someone wants to rerun last year's scenario, they shouldn't have to guess which tab or variable held the original rates.

4.4.3. Data Is Never “Done”

Data handling is not a one-time hurdle to clear. Markets move, data feeds change, formats drift. Build routines to check for “data drift” and have a plan for periodic validations and refreshes.

A few practical tips:

- Maintain a simple data log or data dictionary—even if informal—so others can trace what each field means and where it came from.
- Automate the boring parts: validation scripts, input checks, and sanity tests pay off a hundredfold.
- Version your datasets, just as you do your code. Nothing is more frustrating than trying to reproduce a result only to discover “the input file changed.” See Section 12.5.3

Data is unruly, idiosyncratic, and absolutely central to every model’s fate. Treat it as a first-class concern, not an afterthought, and your models will be far sturdier for it. As a methodical guide, Table 4.4 lists key steps to follow when bringing data into the model.

Table 4.4.: Typical Steps in the Data-to-Model Process.

Step	Key Actions	Purpose / Notes
Data Collection	<ul style="list-style-type: none">• Identify sources• Acquire data (e.g., APIs, databases, scraping)	Ensures data is relevant, reliable, and timely
Data Exploration & Understanding	<ul style="list-style-type: none">• Summary statistics• Visualization	Uncovers initial insights, errors, distributions, and relationships
Data Cleaning	<ul style="list-style-type: none">• Data profiling• Handle missing values• Detect/treat outliers• Data transformation/formatting	Improves data quality, reduces noise and bias

4.4. Feeding The Model

Step	Key Actions	Purpose / Notes
Data Preprocessing	<ul style="list-style-type: none"> • Scale/normalize features • Encode categorical variables • Augment data (if needed) with other datasets 	Prepare data so it fits the format and requirements of the model
Feature Engineering	<ul style="list-style-type: none"> • Select important features • Create new features (e.g., ratios, aggregates) 	Enhance or create new variables that improve model performance
Data Splitting	<ul style="list-style-type: none"> • Divide into training, testing, (validation) sets • Apply cross-validation or static/dynamic validations 	Prevents overfitting and enables robust performance assessment
Data Storage & Management	<ul style="list-style-type: none"> • Store in databases/data lakes • Maintain version control 	Supports reproducibility, scalability, and reliable access
Ethical Considerations	<ul style="list-style-type: none"> • Evaluate bias and fairness • Ensure privacy and regulatory compliance 	Avoids perpetuating bias and protects sensitive information
Continuous Monitoring & Updating	<ul style="list-style-type: none"> • Monitor model/data performance • Detect data drift • Retrain/update as needed 	Maintains accuracy and relevance as data and conditions change

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4.5. Model Management

4.5.1. Risk Governance

An effective risk governance framework for financial modeling begins with clearly stating why such oversight is necessary—namely, to prevent costly missteps in managing complex portfolios or complying with regulations. Organizations often adopt a written policy delineating responsibilities across different levels: management or board-level committees set high-level objectives, while operational teams handle day-to-day processes.

At the heart of this framework lies a structured model inventory, a catalog of all models in use that details each model’s purpose, assumptions, and present status (for example, whether it is in a prototype phase or fully deployed in production). This inventory helps institutions understand their cumulative exposure to errors or assumptions gone awry.

In practice, many firms adopt tiered risk classifications to decide how much scrutiny a model warrants. Classification schemes may range from “low impact” for small-scale financial calculators to “mission-critical” for enterprise valuation engines. Validation and testing approaches vary according to a model’s assigned tier.

Highly critical models undergo more extensive backtesting, benchmarking, or sensitivity analyses, with results escalated to senior management. Risk governance also encompasses ongoing monitoring and scheduled reports about model health. By publicizing validation findings and model performance metrics, the organization fosters a culture where potential failures are escalated early and openly, rather than hidden away until a crisis emerges.

4.5.2. Change Management

No model remains static for long; assumptions evolve, new asset classes appear, and software libraries update. For this reason, a firm’s change management process should standardize how modifications are proposed, evaluated, and documented, ensuring continuity of both the model’s logic and the data that feeds it.

A central repository or version control system is essential: whenever the model or its associated data structures shift, the changes and their justifications must be recorded. This makes it easier to track lineage and revert to a prior version if an update proves problematic in a live environment. Later in this book, we will introduce modern version control systems and workflows that are facilitated by the code-based models that we develop.

4.5. Model Management

Equally important is assessing the ripple effects of each change. Simplifying a routine or adjusting a discount rate assumption may be minor in isolation but can have broad implications when integrated across multiple components. Projects often require up-front impact assessments to determine which historical results need recalculating and whether stakeholder training or documentation updates are needed. One strategy, that of package and model version numbering schemes, will be described in Chapter 23.

Communication around changes should be systematic, distributing concise notes on new features, potential risks, and recommended usage practices to both internal users and (where relevant) regulators. Well-handled change management fosters stability and trust, enabling prompt innovation without sacrificing the reliability of the overall modeling ecosystem.

4.5.3. Data Controls

Sound data controls are paramount in financial modeling because flawed or unverified inputs quickly undermine even the sturdiest model architecture. Organizations typically define data quality standards that address accuracy, completeness, and timeliness. These standards help detect common pitfalls, such as inconsistent formatting, delayed updates, or incorrect data mappings. Complementing formal policies, automated checks are often placed at ingestion points to spot irregularities—anything from out-of-range values that might indicate data corruption, to suspicious spikes hinting at a data input error.

Security and access protocols add another layer of protection. Role-based permission schemes or strong authentication measures minimize the risk of data tampering, accidental deletions, or unauthorized viewings of confidential information.

Although data versioning may sound like a software concept, it applies equally to financial datasets. Keeping a record of each dataset's evolution allows managers and auditors to pinpoint when and how anomalies first appeared. Where legislation like GDPR or industry-specific regulations come into play, data controls must also reflect broader requirements about personal information, consent, and retention periods. Coordinating these efforts under a unified data governance approach ensures that model outputs stand on a solid factual foundation.

4.5.4. Peer and Technical Review

Even the most experienced modelers benefit from additional eyes on their work. **Peer review**, whether informal or systematically mandated, identifies blind spots in assumptions, conceptual design, or coding. Though some organizations require independent reviewers who have not contributed to the original model, smaller teams may rely on

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a rotating schedule of internal experts sharing responsibility for checks. The key is cultivating a culture where open dialogue about potential faults is not only accepted but encouraged.

Technical review goes one step further, focusing on deeper verification of the computations themselves. Complex spreadsheets, code modules, or integrated software platforms may require structured walk-throughs in which reviewers verify arithmetic, confirm the alignment of calculation steps with business logic, or run test scenarios to ensure the model behaves as intended. This process should generate formal documentation capturing who performed the review, what methods they used, and which issues surfaced. Likewise, conceptual soundness—how well the model aligns with economic theory or domain-specific knowledge—merits discussion in a thorough review. If challenges are identified, revisions loop back into the change management system, promoting iterative refinements. By conducting peer and technical reviews in earnest, organizations reinforce consistent quality and reduce the likelihood of undetected errors slipping into production.

4.6. Conclusion

The art and science of financial modeling require a unique blend of skills, knowledge, and personal qualities. A proficient modeler combines domain expertise, theoretical understanding, and practical skills with a curious and rigorous mindset. They leverage a diverse toolset, employ sound architectural principles, and communicate with clarity. The ability to navigate the complexities of financial systems while maintaining humility in the face of irreducible uncertainties is paramount.

As the financial world continues to evolve, so too must the modeler’s approach. By cultivating these attributes and continuously refining their craft, financial modelers can create more robust, insightful, and valuable models that drive informed decision-making in an increasingly complex economic landscape. The journey of a financial modeler is one of perpetual learning and adaptation, where each challenge presents an opportunity for growth and innovation.

Part III.

**Foundations: Programming and
Abstractions**

Out of intense complexities intense simplicities emerge. - Winston Churchill
(1923)

The next several chapters build up essential concepts that enable us to create sophisticated financial models while maintaining clarity and manageability. We'll start with core programming building blocks - the vocabulary and grammar of communicating with computers. From there, we'll explore different approaches to breaking down complex problems into simpler components through abstraction: functions, types, programming paradigms, and more.

Abstraction is a form of selective ignorance - focusing attention on what matters for a particular purpose while hiding irrelevant details. Just as financial statements abstract away the details of individual transactions to reveal the bigger picture, we'll see how thoughtful abstraction in programming allows us to manage complexity by working at different levels of detail.

The goal is not to make you a computer scientist, but rather to equip you with the mental models and practical techniques needed to effectively leverage programming in your financial work. By understanding these foundations, you'll be better equipped to design clean, maintainable models that can evolve with your needs.

Think of this section as building your modeling toolkit, one concept at a time. We'll introduce ideas progressively, with plenty of concrete examples to ground the theory in practical application. The concepts build on each other, so take time to ensure your understanding before moving forward. Let's begin with the fundamental elements of programming.

5. Elements of Programming

“Programming is not about typing, it’s about thinking.” — Rich Hickey
(2011)

5.1. In this section

Start building up computer science concepts by introducing tangible programming essentials. Data types, variables, control flow, functions, and scope are introduced.

💡 On Your First Readthrough

This chapter is intended to be an introductory reference for most of the basic building blocks for which we will build abstractions on top of in chapters that follow. We want this chapter to essentially be an easy and mildly opinionated stepping-stone on your journey.

At some point, you will likely find yourself seeking more precise or thorough documentation and will begin directly searching or reading the documentation of a language or library itself. However, it may be intimidating or frustrating reading reference documentation due to the density and terminology - let this chapter (and book writ large) be a bridge for you.

If reading this book in a linear fashion and new to programming, we suggest skipping the following sections and returning when encountering the concept or term later in the book:

- Section 5.4.4 through Section 5.4.9 which covers advanced and custom data types
- After Section 5.5.3 which deals with advanced function usage and program organization via scope

🔥 Caution

This introductory chapter is intended to provide a survey of the important concepts and building blocks, not to be a complete reference. For full details on available functions, more complete definitions, and a more complete tour of all language

5. Elements of Programming

features, see the Manual at docs.julialang.org.

5.2. Computer Science, Programming, and Coding

Computer Science is the study of computing and information. As a science, it is distinct from programming languages which are merely coarse implementations of specific computer science concepts¹.

Programming (or “**coding**”) is the art and science of writing code in programming languages to have the computer perform desired tasks. While this may sound mechanistic, programming truly is one of the highest forms of abstract thinking. The design space of potential solutions is so large and potentially complex that much art and experience is needed to create a well-made program.

The language of computer science also provides a lexicon so that financial practitioners can discuss model architecture and characteristics of problems with precision and clarity. Simply having additional terminology and language to describe a concept illuminates aspects of the problem in new ways, opening one’s self up to more innovative solutions.

In this light, the financial modeling that we do can be considered a type of computer program. It takes as input abstract information (data), performs calculations (an algorithm), and returns new data as an output. We generally do not need to consider many things that a software engineer may contemplate such as a graphical user interface, networking, or access restrictions. However, the programming fundamentals are there: a good financial modeler must understand data types, algorithms, and some hardware details.

We will build up the concepts over this and the following chapter:

- This chapter will provide a survey of important concepts in computer science that will prove useful for our financial modeling. First, we will talk about data types, boolean logic, and basic expressions. We’ll build on those to discuss algorithms (functions) which perform useful work and use control flow and recursion.
- In the following chapters about abstraction, we will step back and discuss higher level concepts: the “schools of thought” around organizing the relationship between data and functions (functional versus object-oriented programming), design patterns, computational complexity, and compilation.

¹Said differently, computer science may contemplate ideas and abstractions more generally than a specific implementation, as in mathematics where a theorem may be proved ($a^2 + b^2 = c^2$) without resorting to specific numeric examples ($3^2 + 4^2 = 5^2$).

 Tip

There will be brief references to hardware considerations for completeness, but hardware knowledge is not necessary to understand most programming languages (including Julia). It's impossible to completely avoid talking about hardware when you care about the performance of your code, so feel free to gloss over the reference to hardware details on the first read and come back later after Chapter 9.

It's highly recommended that you follow along and have a Julia session open (e.g. a REPL or a notebook) when first going through this chapter. See the first part of Chapter 21 if you haven't gotten that set up yet. Follow along with the examples as we go.

 Tip

You can get some help in the REPL by typing a ? followed by the symbol you want help with, for example:

```
help?> sum
search: sum sum! summary cumsum cumsum! ...
sum(f, itr; [init])
```

Sum the results of calling function f on each element of itr.

... More text truncated...

5.3. Expressions and Control Flow

5.3.1. Assignment and Variables

One of the first things it will be convenient to understand is the concept of variables. In virtually every programming language, we can assign values to make our program more organized and meaningful to the human reader. In the following example, we assign values to intermediate symbols to benefit us humans as we convert (silly!) American distance units:

```
feet_per_yard = 3
yards_per_mile = 1760

feet = 3000
miles = feet / feet_per_yard / yards_per_mile
```

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```
0.56818181818182
```

The above is technically the same thing as just computing $3000 / 3 / 1760$, however we've given the elements names meaningful to the human user.

Beyond readability, variables are a form of **abstraction** which allows us to think beyond specific instances of data and numbers to a more general representation. For example, the last line in the prior code example is a very generic computation of a unit conversion relationship and `feet` could be any number and the expression remains a valid calculation.

Tip

We will dive a little bit deeper into variables and assignment in Section 5.3.4, distinguishing between assignment and references.

5.3.2. Expressions

Within the code examples above, we can zoom in onto small pieces of code called **expressions**. Expressions are effectively the basic block of code that gets evaluated to produce a value. Here is an expression that adds two integers together that evaluate to a new integer (3 in this case):

```
1 + 2
```

```
3
```

A bigger program is built up of many of these smaller bits of code.

5.3.2.1. Compound Expression

There's two kinds of blocks where we can ensure that sub-expressions get evaluated in order and return the last expression as the overall return value: `begin` and `let` blocks.

```
c = begin
  a = 3
  b = 4
  a + b
end

a, b, c
```

(3, 4, 7)

Alternatively, you can chain together ;s to create a compound expression:

```
z = (x = 1; y = 2; x + y)
```

3

Compound expressions allow you to group multiple operations together while still having the entire block evaluate to a single value, typically the last expression. This makes it easy to use complex logic anywhere a value is needed, like in function arguments or assignments.

5.3.2.2. Conditional Expressions

Conditionals are expressions that evaluate to a **boolean** true or false. This is the beginning of really being able to assemble complex logic to perform useful work. Here are a handful expressions that would evaluate to true:

```
1 > 0
1 == 1 # check for equality
Float64 isa Rational
(5 > 0) & (-1 < 2) # "and" expression
(5 > 0) | (-1 > 2) # "or" expression
1 != 2
```

true

i Note

In Julia, the booleans have an integer equality: true is equal to 1 (`true == 1`) and false is equal to 0 (`false == 0`). However:

- `true != 5`. Only 1 is equal to true (in some languages, any non-zero number is “truthy”).
- `true` is not equal to 1 (equal is defined later in this chapter).

Conditionals can be used to assemble different logical paths for the program to follow and the general pattern is an if block:

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```
if condition
    # do one thing
elseif condition
    # do something else
else
    # do something if none of the
    # other conditions are met
end
```

A complete example:

```
function buy_or_sell(my_value, market_price)
    if my_value > market_price
        "buy more"
    elseif my_value < market_price
        "sell"
    else
        "hold"
    end
end

buy_or_sell(10, 15), buy_or_sell(15, 10), buy_or_sell(10, 10)

("sell", "buy more", "hold")
```

5.3.3. Equality

The “Ship of Theseus²” problem is an example of how equality can be philosophically complex concept. In computer science we have the advantage that while we may not be able to resolve what’s the “right” type of equality, we can be more precise about it.

Here is an example for which we can see the difference between two types of equality:

- **Egal** equality is when a program could not distinguish between two objects at all
- **Equal** equality is when the values of two objects are the same

If two things are egal, then they are also equal.

In the following example, s and t are equal but not egal:

²The Ship of Theseus problem specifically refers to a legendary ancient Greek ship, owned by the hero Theseus. The paradox arises from the scenario where, over time, each wooden part of the ship is replaced with identical materials, leading to the question of whether the fully restored ship is still the same ship as the original. The Ship of Theseus problem is a thought experiment in philosophy that explores the nature of identity and change. It questions whether an object that has had all of its components replaced remains fundamentally the same object.

5.3. Expressions and Control Flow

```
s = [1, 2, 3]
t = [1, 2, 3]
s == t, s === t
```

```
(true, false)
```

One way to think about this is that while the values are equal, there is a way that one of the arrays could be made not equal to the other:

```
t[2] = 5
t
```

```
3-element Vector{Int64}:
1
5
3
```

Now t is no longer equal to s:

```
s == t
false
```

The reason this happens is that arrays are containers that can have their contents modified. Even though they originally had the same values, s and t are different containers, and *it just so happened* that the values they contained started out the same.

Some data can't be modified, including some kinds of collections. Immutable types like the following tuple, with the same stored values, are equal because there is no way for us to make them different:

```
(2, 4) === (2, 4)
```

```
true
```

Using this terminology, we could now interpret the "Ship of Theseus" as that his ship is "equal" but not "equal".

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5.3.4. Assignment and Variables

When we say `x = 2` we are **assigning** the integer value of 2 to the variable `x`. This is an expression that lets us bind a something to a variable so that it can be referenced more concisely or in different parts of our code. When we re-assign the variable we are not mutating the value: `x = 3` does not change the 2.

When we have a mutable object (e.g. an Array or a `mutable struct`), we can mutate the value inside the referenced container. For example:

```
x = [1, 2, 3]                                ①  
x[1] = 5                                     ②  
x
```

- ① `x` refers to the array which currently contains the elements 1, 2, and 3.
- ② We re-assign the first element of the array to be the value 5 instead of 1

```
3-element Vector{Int64}:  
5  
2  
3
```

In the above example, `x` has not been reassigned. It is possible for two variables to refer to the same object:

```
x = [1, 2, 3]  
y = x  
x[1] = 6  
y
```

- ① `y` refers to the *same* underlying array as `x`

```
3-element Vector{Int64}:  
6  
2  
3
```

Note that variables can be re-assigned unless they are marked as `const`:

```
const PHI = π * 2 # <1>
```

- ① Capitalizing constant variables is a convention in Julia.

If we tried to re-assign `PHI`, we would get an error.

 Warning

Note that if we declare a `const` variable that refers to a mutable container like an array, the container can still be mutated. It's the reference to the container that remains constant, not necessarily the elements within the container.

5.3.5. Loops

Loops are ways for the program to move through a program and repeat expressions while we want it to. There are two primary loops: `for` and `while`.

for loops are loops that iterate over a defined range or set of values. Let's assume that we have the array `v = [6,7,8]`. Here are multiple examples of using a `for` loop in order to print each value to output (`println`):

```
# use fixed indices
for i in 1:3
    println(v[i])
end

# use indices the of the array
for i in eachindex(v)
    println(v[i])
end

# use the elements of the array
for x in v
    println(x)
end

# use the elements of the array
for x ∈ v           # ∈ is typed \in<tab>
    println(x)
end
```

while loops will run repeatedly until an expression is false. Here's some examples of printing each value of `v` again:

```
# index the array
i = 1
while i <= length(v)
    println(v[i])
    global i += 1
end
```

(1)

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- ① `global` is used to increment `i` by 1. `i` is defined outside the scope of the `while` loop (see Section 5.6).

```
# index the array
i = 1
while true
    println(v[i])
    if i >= length(v)
        break
    end
    global i += 1
end
```

(1)

- ① `break` is used to terminate the loop manually, since the condition that follows the `while` will never be false.

5.3.6. Performance of loops

Loops are highly performant in Julia and often the fastest way to accomplish things. Those coming from Python or R may have developed a habit to avoid writing loops. *Fear the for loop not!*

5.4. Data Types

Data types are a way of categorizing information by intrinsic characteristics. We instinctively know that `13.24` is different than "this set of words" and types are how we will formalize this distinction. This is a key conceptual point, and mathematically it's like we have different sets of objects to perform specialized operations on. Beyond this set-like abstraction is implementation details related to computer hardware. You probably know that computers only natively "speak" in binary zeros and ones. Data types are a primary way that a computer can understand if it should interpret `01000010` as `B` or as `66`³.

Each 0 or 1 within a computer is called a **bit** and eight bits in a row form a **byte** (such as `01000010`). This is where we get terms like "gigabytes" or "kilobits per second" as a measure of the quantity or rate of bits something can handle⁴.

³This binary representations correspond to `B` and `66` with the *ASCII character set* and 8-bit integer encodings respectively, discussed later in this chapter.

⁴Some distinctions you may encounter: in short-form, "kb" means kilobits while the upper-case "B" in "kB" means kilobytes. Also confusingly, sometimes the "k" can be binary or decimal - because computers speak in binary, a binary "k" means 1024 (equal to 2^{10}) instead of the usual decimal 1000. In most computer contexts, the binary (multiples of 1024) is more common.

5.4.1. Numbers

Numbers are usually grouped into two categories: **integers** and **floating-point**⁵ numbers. Integers are like the mathematical set of integers while floating-point is a way of representing decimal numbers. Both have some limitations since computers can only natively represent a finite set of numbers due to the hardware (more on this in Chapter 9). Here are three integers that are input into the **REPL** (Read-Eval-Print-Loop)⁶ and the result is **printed** below the input:

2

2

423

423

1929234

1929234

And three floating-point numbers:

0.2

0.2

-23.3421

14e3 # the same as 14,000.0

14000.0

⁵The term floating point refers to the fact that the number's radix (decimal) point can "float" between the significant digits of the number.

⁶That is, it *reads* the code input from the user, *evaluates* what code was given to it, *prints* the result of the input to the screen, and *loops* through the process again.

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On most systems, `0.2` will be interpreted as a 64-bit floating point type called `Float64` in Julia since most architectures these days are 64-bit⁷, while on a 32-bit system `0.2` would be interpreted as a `Float32`. Given that there are a finite amount of bits attempting to represent a continuous, infinite set of numbers means that some numbers are not able to be represented with perfect precision. For example, if we ask for `0.2`, the closest representations in 64 and 32 bit are:

- `0.20000000298023223876953125` in 32-bit
- `0.200000000000000011102230246251565404236316680908203125` in 64-bit

This leads to special considerations that computers take when performing calculations on floating point maths, some of which will be covered in more detail in Chapter 9. For now, just note that floating point numbers have limited precision and even if we input `0.2`, your computations will use the above decimal representations even if it will print out a number with fewer digits shown:

`x = 0.2` (1)

`big(x)` (2)

- ① Here, we **assign** the value `0.2` to a **variable** `x`. More on variables/assignments in Section 5.3.4.
- ② `big(x)` is a arbitrary precision floating point number and by default prints the full precision that was embedded in our variable `x`, which was originally `Float64`.

`0.200000000000000011102230246251565404236316680908203125`

i Note

Note the difference in what printed between the last example and when we input `0.2` earlier in the chapter. The former had the same (not-exactly equal to `0.2`) *value*, but it printed an abbreviated set of digits as a nicety for the user, who usually doesn't want to look at floating point numbers with their full machine precision. The system has the full precision (`0.20...3125`) but is truncating the output. In the last example, we've converted the normal `Float64` to a `BigFloat` which will not truncate the output when printing.

Integers are similarly represented as 32 or 64 bits (with `Int32` and `Int64`) and are limited to exact precision:

- -32,767 to 32,767 for `Int32`
- -2,147,483,647 to 2,147,483,647 for `Int64`

⁷This means that their central processing units (CPUs) use instructions that are 64 bits long.

5.4. Data Types

Additional range in the positive direction if one chooses to use “unsigned”, non-negative numbers (`UInt32` and `UInt64`). Unlike floating point numbers, the integers have a type `Int` which will use the system bit architecture by default (that is, `Int(30)` will create a 64 bit integer on 64-bit systems and 32-bit on 32-bit systems).

💡 Floating Point and Excel

Excel’s numeric storage and routine is complex and not quite the same as most programming languages, which follow the Institute of Electrical and Electronics Engineer’s standards (such as the IEEE 754 standard for double precision floating point numbers). Excel uses IEEE for the *computations* but results (and therefore the cells that comprise many calculations interim values) are stored with 15 significant digits of information. In some ways this is the worst of both worlds: having the sometimes unusual (but well-defined) behavior of floating point arithmetic *and* having additional modifications to various steps of a calculation. In general, you can assume that the programming language result (following the IEEE 754 standard) is a better result because there are aspects to the IEEE 754 defines techniques to minimize issues that arise in floating point math. Some of the issues (round-off or truncation) can be amplified instead of minimized with Excel.

In practice, this means that it can be difficult to exactly replicate a calculation in Excel in a programming language and vice-versa. It’s best to try to validate a programming model versus Excel model using very small unit calculations (e.g. a single step or iteration of a routine) instead of an all-in result. You may need to define some tolerance threshold for comparison of a value that is the result of a long chain of calculation.

💡 Currencies and Decimals

Note that floating point numbers should **not** be used in storing transaction records! The intricacies of floating point math described above to not lend itself to accurate record-keeping, since with operations like `0.11 + 0.12` don’t precisely equal `0.23`

`BigFloat(0.11 + 0.12)`

```
0.2299999999999982236431605997495353221893310546875
```

As you can see, if we were adding US Dollar cents here, we would have destroyed a fraction of a penny. Do that for millions of transactions in a day and you have a problem!

Generally, when doing *modeling* or even creating a *valuation model*, it’s okay to use floating point math. As an example, if you are trying to determine the value of an exotic option, your model is likely just fine outputting a value like `101.987087`. If you go and sell this option, you’ll have to settle for either `101.98` or `101.99` when booking it. In most contexts this imprecision is likely okay!

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If you are implementing a transaction or trading system, ensure proper treatment of the types representing your monetary numbers. A full treatment is beyond the scope of this book, but for a good introduction to the subject, see <https://cs-syd.eu/posts/2022-08-22-how-to-deal-with-money-in-software>.

5.4.2. Type Hierarchy

We can describe a *hierarchy* of types. Both `Float64` and `Int64` are examples of `Real` numbers (here, `Real` is an **abstract** Julia type which corresponds to the mathematical set of real numbers commonly denoted with \mathbb{R}). Both `Float64` and `Int32` are `Real` numbers, so why not just define all numbers as a `Real` type? Because for performant calculations, the computer must know in advance how many bits each number is represented with.

`?@fig-julia-numeric-types` shows the type hierarchy for most built-in Julia number types.

TODO: Once Quarto Issue #10961 is resolved, render the mermaid diagram.

```
%| label: fig-julia-numeric-types
%| fig-cap: "Numeric Type Hierarchy in Julia. Leafs of the tree are concrete types."
%| fig-width: 6.5
graph TD
    Number --> Real
    Number --> Complex

    Real --> Integer
    Real --> AbstractFloat
    Real --> Rational
    Real --> Irrational

    Integer --> Signed
    Integer --> Unsigned

    Signed --> Int8
    Signed --> Int16
    Signed --> Int32
    Signed --> Int64
    Signed --> Int128
    Signed --> BigInt

    Unsigned --> UInt8
    Unsigned --> UInt16
    Unsigned --> UInt32
```

```

Unsigned --> UInt64
Unsigned --> UInt128

AbstractFloat --> Float16
AbstractFloat --> Float32
AbstractFloat --> Float64
AbstractFloat --> BigFloat

```

The integer and floating point types described in the prior section are known as **concrete** types because there are no possible sub types (child types). Further, a concrete type can be a **bit type** if the data type will always have the same number of bits in memory: a `Float32` will always be 32 bits in memory, for example. Contrast this with strings (described below) which can contain an arbitrary number of characters.

5.4.3. Collections

Collections are types that are really useful for storing data which contains many elements. This section describes some of the most common and useful types of containers.

5.4.3.1. Arrays

Arrays are the most common way to represent a collection of similar data. For example, we can represent a set of integers as follows:

`[1, 10, 300]`

```

3-element Vector{Int64}:
 1
 10
 300

```

And a floating point array:

`[0.2, 1.3, 300.0]`

```

3-element Vector{Float64}:
 0.2
 1.3
 300.0

```

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Note the above two arrays are different types of arrays. The first is `Vector{Int64}` and the second is `Vector{Float64}`. These are arrays of concrete types and so Julia will know that each element of an array is the same amount of bits which will enable more efficient computations. With the following set of mixed numbers, Julia will **promote** the integers to floating point since the integers can be accurately represented⁸ in floating point.

```
[1, 1.3, 300.0, 21]
```

4-element `Vector{Float64}`:

```
1.0  
1.3  
300.0  
21.0
```

However, if we explicitly ask Julia to use a `Real`-typed array, the type is now `Vector{Real}`. Recall that `Real` is an abstract type. Having heterogeneous types within the array is conceptually fine, but in practice limits performance. Again, this will be covered in more detail in Chapter 9.

In Julia, arrays can be multi-dimensional. Here are two three-dimensional arrays with length three in each dimension:

```
rand(3, 3, 3)
```

3×3×3 `Array{Float64, 3}`:

```
[:, :, 1] =  
0.649159 0.136041 0.152265  
0.352435 0.63539 0.77435  
0.482884 0.645682 0.470798
```

```
[:, :, 2] =  
0.704094 0.277493 0.604486  
0.590259 0.461246 0.98894  
0.353084 0.771033 0.999525
```

```
[:, :, 3] =  
0.382259 0.825086 0.455453  
0.513259 0.720049 0.206382  
0.374519 0.668256 0.590867
```

```
[x + y + z for x in 1:3, y in 11:13, z in 21:23]
```

⁸Accurate only to a limited precision, as described in Section 5.4.1.

```
3x3x3 Array{Int64, 3}:
```

```
[:, :, 1] =
```

```
33 34 35
```

```
34 35 36
```

```
35 36 37
```

```
[:, :, 2] =
```

```
34 35 36
```

```
35 36 37
```

```
36 37 38
```

```
[:, :, 3] =
```

```
35 36 37
```

```
36 37 38
```

```
37 38 39
```

The above example demonstrates **array comprehension** syntax which is a convenient way to create arrays in Julia.

A two-dimensional array has the rows by semi-colons (;):

```
x = [1 2 3; 4 5 6]
```

```
2x3 Matrix{Int64}:
```

```
1 2 3
```

```
4 5 6
```

i Note

In Julia, a `Vector{Float64}` is simply a one-dimensional array of floating points and a `Matrix{Float64}` is a two-dimensional array. More precisely, they are **type aliases** of the more generic `Array{Float64,1}` and `Array{Float64,2}` names. Arrays with three or more dimensions don't have a type alias pre-defined.

5.4.3.2. Array indexing

Array elements are accessed with the integer position, starting at 1 for the first element⁹
¹⁰:

⁹Whether an index starts at 1 or 0 is sometimes debated. Zero-based indexing is natural in the context of low-level programming which deal with bits and positional *offsets* in computer memory. For higher level programming one-based indexing is more natural: in a set of data stored in an array, it is much more natural to reference the *first* (through n^{th}) datum instead of the *zeroth* (through $(n-1)^{th}$) datum.

¹⁰Arrays in Julia can actually be indexed with an arbitrary starting point: see the package `OffsetArrays.jl`

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```
v = [10, 20, 30, 40, 50]  
v[2]
```

20

We can also access a subset of the vector's contents by passing a range:

```
v[2:4]
```

```
3-element Vector{Int64}:  
20  
30  
40
```

And we can generically reference the array's contents, such as:

```
v[begin+1:end-1]
```

```
3-element Vector{Int64}:  
20  
30  
40
```

We can assign values into the array as well, as well as combine arrays and push new elements to the end:

```
v[2] = -1  
push!(v, 5)  
vcat(v, [1, 2, 3])
```

```
9-element Vector{Int64}:  
10  
-1  
30  
40  
50  
5  
1  
2  
3
```

5.4.3.3. Array Alignment

When you have an MxN matrix (M rows, N columns), a choice must be made as to which elements are next to each other in memory. Typical math convention and fundamental computer linear algebra libraries (dating back decades!) are column major and Julia follows that legacy. **Column major** means that elements going down the rows of a column are stored next to each other in memory. This is important to know so that (1) you remember that vectors are treated like a column vector when working with arrays (that is: a N element 1D vector is like a Nx1 matrix), and (2) when iterating through an array, it will be faster for the computer to access elements next to each other column-wise. A 10x10 matrix is actually stored in memory as 100 elements coming in order, one after another in single file.

This 3x4 matrix is stored with the elements of columns next to each other, which we can see with `vec`:

```
mat = [1 2 3; 4 5 6; 7 8 9]
```

`3x3 Matrix{Int64}:`

```
1 2 3
4 5 6
7 8 9
```

```
vec(mat)
```

`9-element Vector{Int64}:`

```
1
4
7
2
5
8
3
6
9
```

5.4.3.4. Ranges

A **range** is a representation of a range of numbers. We actually used them above to index into arrays. They are expressed as `start:stop`

We don't have to actually store all of these numbers on the computer somewhere as in an `Array`. Instead, this is an object that *represents* the ordered set of numbers. So for example, we can sum up 1 through the number of atoms on the earth instantaneously:

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This is possible due to two things:

1. not needing to actually store that many numbers in memory, and
 2. Julia being smart enough to apply the triangular number formula¹¹ when `sum` is given a consecutive range.

There are more general ways to construct ranges:

Step by another number instead of the default 1:

1:2:7

1:2:7

Specify the number of values within the range, inclusive of the first number¹²:

```
range(0, 10, 21)
```

0.0:0.5:10.0

5.4.3.5. Characters, Strings, and Symbols

Characters are represented in most programming languages as letters within quotation marks. In Julia, individual characters are represented using single quotes:

'a'

'a': ASCII/Unicode U+0061 (category Ll: Letter, lowercase)

¹¹The triangular numbers (sum of integers from 1 to n) are:

$$T_n = \sum_{k=1}^n k = 1 + 2 + \dots + n = \frac{n^2 + n}{2} = \frac{n(n+1)}{2} = \binom{n+1}{2}$$

¹²Whether the last number is in the resulting range depends on if the step evenly divides the end of the range.

5.4. Data Types

Letters and other characters present more difficulties than numbers to represent within a computer (think of how many languages and alphabets exist!), and it essentially only works because the world at large has agreed to a given representation. Originally **ASCII** (American Standard Code for Information Interchange) was used to represent just 95 of the most common English characters ("a" through "z", zero through nine, etc.). Now, **UTF** (Unicode Transformation Format) can encode more than a million characters and symbols from many human languages.

Strings are a collection¹³ of characters, and can be created in Julia with double quotes:

```
"hello world"
```

```
"hello world"
```

It's easy to ascertain how 'normal' characters can be inserted into a string, but what about things like new lines or tabs? They are represented by their own characters but are normally not printed in computer output. However, those otherwise invisible characters do exist. For example, here we will use a **string literal** (indicated by the """") to tell Julia to interpret the string as given, including the invisible new line created by hitting return on the keyboard between the two words:

```
"""
hello
world
"""
```

```
"hello\nworld\n"
```

The output above shows the \n character contained within the string.

Symbols are a way of representing an identifier which cannot be seen as a collection of individual characters. :helloworld is distinct from "helloworld" - you can kind of think of the former as an un-executed bit of code - if we were to execute it (with eval(:helloworld)), we would get an error UndefVarError: 'a' not defined . Symbols can *look* like strings but do not behave like them. For now, it is best to not worry about symbols but it is an important aspect of Julia which allows the language to represent aspects of itself as data. This allows for powerful self-reference and self-modification of code but this is a more advanced topic generally out of scope of this book.

¹³Under the hood, strings are essentially a vector of characters but there are complexities with character encoding that don't allow a lossless conversion to individual characters of uniform bit length. This is for historical compatibility reasons and to avoid making most documents' file sizes larger than it needs to be.

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5.4.3.6. Tuples

Tuples are a set of values that belong together and are denoted by values inside parenthesis and separated by a comma. An example might be x-y coordinates in 2 dimensional space:

```
x = 3  
y = 4  
p1 = (x, y)
```

(3, 4)

Tuple's values can be accessed like arrays:

```
p1[1]
```

3

Tuples fill a middle ground between scalar types and arrays in more ways than one:

- Tuples have no problem having heterogeneous types in the different slots.
- Tuples are **immutable**, meaning that you cannot overwrite the value in memory (an error will be thrown if we try to do `p[1] = 5`).
- It's generally expected that within an array, you would be able to apply the same operation to all the elements (e.g. square each element) or do something like sum all of the elements together which isn't generally case for a tuple.
- Tuples are generally stack allocated instead of being heap allocated like arrays¹⁴, meaning that a lot of times they can be faster than arrays.

5.4.3.6.1. Named Tuples

Named tuples provide a way to give each field within the tuple a specific name. For example, our x-y coordinate example above could become:

```
p2 = (x=3, y=4)
```

(x = 3, y = 4)

¹⁴What this means will be explained in Chapter 9 .

5.4. Data Types

The benefit is that we can give more meaning to each field and access the values in a nicer way. Previously, we used `location[1]` to access the `x`-value, but with the new definition we can access it by name:

```
p2.x
```

3

5.4.3.7. Dictionaries

Dictionaries are a container which relates a **key** to an associated **value**. Kind of like how arrays relate an index to a value, but the difference is that a dictionary is (1) un-ordered and (2) the key doesn't have to be an integer.

Here's an example which relates a name to an age:

```
d = Dict(  
    "Joelle" => 10,  
    "Monica" => 84,  
    "Zaylee" => 39,  
)
```

```
Dict{String, Int64} with 3 entries:  
"Monica" => 84  
"Zaylee" => 39  
"Joelle" => 10
```

Then we can look up an age given a name:

```
d["Zaylee"]
```

39

Dictionaries are super flexible data structures and can be used in many situations.

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5.4.4. Parametric Types

We just saw how tuples can contain heterogeneous types of data inside a common container. **Parametric Types** are a way of allowing types themselves to be variable, with a wrapper type containing a to-be-specified inner-type.

Let's look at this a little bit closer by looking at the full type:

```
typeof(p1)
```

```
Tuple{Int64, Int64}
```

`location` is a `Tuple{Int64, Int64}` type, which means that its first and second elements are both `Int64`. Contrast this with:

```
typeof(("hello", 1.0))
```

```
Tuple{String, Float64}
```

These tuples are both of the form `Tuple{T,U}` where `T` and `U` are both types. Why does this matter? We and the compiler can distinguish between a `Tuple{Int64, Int64}` and a `Tuple{String, Float64}` which allows us to reason about things ("I can add the first element of tuple together only if both are numbers") and the compiler to optimize (sometimes it can know exactly how many bits in memory a tuple of a certain kind will need and be more efficient about memory use). Further, we will see how this can become a powerful force in writing appropriately abstracted code and more logically organize our entire program when we encounter "multiple dispatch" later on.

This is a very powerful technique - we've already seen the flexibility of having an `Array` type which can contain arbitrary inner types and dimensions. The full type signature for an `Array` looks like `Array{InnerType, NumDimensions}`.

```
let
  x = [1 2
        3 4]
  typeof(x)
end

Matrix{Int64} (alias for Array{Int64, 2})
```

Table 5.1.: Three value logic with true, missing, and false.

(a) Not logic			
NOT (!)	Value		
true	false		
missing	missing		
false	true		

(b) And logic			
AND (&)	true	missing	false
true	true	missing	false
missing	missing	missing	false
false	false	false	false

(c) Or Logic			
OR ()	true	missing	false
true	true	true	true
missing	true	missing	missing
false	true	missing	false

5.4.5. Types for things not there

nothing represents that there's nothing to be returned - for example if there's no solution to an optimization problem or if a function just doesn't have any value to return (such as in the case with input/output like `println`).

missing is to represent something *should* be there but it's not, as is all too common in real-world data. Julia natively supports missing and three-value logic, which is an extension of the two-value boolean (true/false) logic, to handle missing logical values:

 Tip

Missing and Nothing are the *types* while missing and nothing are the values here¹⁵. This is analogous to `Float64` being a type and `2.0` being a value.

¹⁵Missing and Nothing are instances of **singleton type**, which means that there is only a single value that either type can take on.

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5.4.6. Union Types

When two types may arise in a context, **union types** are a way to represent that. For example, if we have a data feed and we know that it will produce *either* a `Float64` or a `Missing` type then we can say that the value for this is `Union{Float64, Missing}`. This is much better for the compiler (and our performance!) than saying that the type of this is `Any`.

5.4.7. Creating User Defined Types

We've talked about some built-in types but so much additional capabilities come from being able to define our own types. For example, taking the x-y-coordinate example from above, we could do the following instead of defining a tuple:

```
struct BasicPoint
    x :: Int64
    y :: Int64
end

p3 = BasicPoint(3, 4)
```

```
BasicPoint(3, 4)
```

`BasicPoint` is a **composite type** because it is composed of elements of other types. Fields are accessed the same way as named tuples:

```
p3.x, p3.y
```

(1)

- ① Note that here, Julia will return a tuple instead of a single value due to the comma separated expressions.

```
(3, 4)
```

`structs` in Julia are immutable like tuples above.

But wait, didn't tuples let us mix types too via parametric types? Yes, and we can do the same with our type!

```
struct Point{T}
    x :: T
    y :: T
end
```

5.4. Data Types

Line 1 The `{T}` after the type's name allows for different Points to be created depending on what the type of the underlying `x` and `y` is.

Here's two new points which now have different types:

```
p4 = Point(1, 4)
p5 = Point(2.0, 3.0)
```

`p4, p5`

```
(Point{Int64}(1, 4), Point{Float64}(2.0, 3.0))
```

Note that the types are not equal because they have different type parameters!

```
typeof(p4), typeof(p5), typeof(p4) == typeof(p5)

(Point{Int64}, Point{Float64}, false)
```

But both are now subtypes of `PPoint2D`. The expression `X isa Y` is true when `X` is a (sub)type of `Y`:

```
p4 isa Point, p5 isa Point

(true, true)
```

Note though, that the `x` and `y` are both of the same type in each `PPoint2D` that we created. If instead we wanted to allow the coordinates to be of different types, then we could have defined `PPoint2D` as follows:

```
struct Point{T,U}
    x :: T
    y :: U
end
```

i Note

Can we define the structs above without indicating a (parametric) type? Yes!

```
struct Point
    x # no type here!
    y # no type declared here either!
end
```

But! `x` and `y` will both be allowed to be `Any`, which is the fallback type where Ju-

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lia says that it doesn't know any more about the type until runtime (the time at which our program encounters the data when running). This means that the compiler (and us!) can't reason about or optimize the code as effectively as when the types are explicit or parametric. This is an example of how Julia can provide a nice learning curve - don't worry about the types until you start to get more sophisticated about the program design or need to extract more performance from the code.

The above structs that we have defined are examples of **concrete types** types which hold data. **Abstract types** don't directly hold data themselves but are used to define a hierarchy of types which we will later exploit (Chapter 8) to implement custom behavior depending on what type our data is.

Here's an example of (1) defining a set of related types that sits above our Point2D:

```
abstract type Coordinate end
abstract type CartesianCoordinate <: Coordinate end
abstract type PolarCoordinate <: Coordinate end

struct Point2D{T} <: CartesianCoordinate
    x::T
    y::T
end

struct Point3D{T} <: CartesianCoordinate
    x::T
    y::T
    z::T
end

struct Polar2D{T} <: PolarCoordinate
    r::T
    θ::T
end
```

💡 Unicode Characters

Julia has wonderful Unicode support, meaning that it's not a problem to include characters like θ . The character can be typed in Julia editors by entering \theta and then pressing the TAB key on the keyboard.

Unicode is helpful for following conventions that you may be used to in math. For example, the math formula $\text{circumference}(r) = 2 \times r \times \pi$ can be written in Julia with `circumference(r) = 2 * r * π`.

The name for the characters follows the same for LaTeX, so you can search the internet for, e.g. “theta LaTeX” to find the appropriate name. Furhter, you can use the REPL help mode to find out how to enter a character if you can copy and paste it from somewhere:

```
help?> θ
"θ" can be typed by \theta<tab>
```

To constrain the types that could be used within our coordinates above, such as if we wanted the fields to all be Real-valued, we could modify the struct definitions with the `<:Real` annotation:

```
struct Point2D{T<:Real} <: CartesianCoordinate
    # ...
end

struct Point3D{T<:Real} <: CartesianCoordinate
    # ...
end

struct Polar2D{T<:Real} <: PolarCoordinate
    # ...
end
```

5.4.8. Mutable structs

It is possible to define structs where the data can be modified - such a data field is said to be **mutable** because it can be changed or mutated. Here's an example of what it would look like if we made Point2D mutable:

```
mutable struct Point2D{T}
    x :: T
    y :: T
end
```

You may find that this more naturally represents what you are trying to do. However, recall that an advantage of an immutable datatype is that costly memory doesn't necessarily have to be allocated for it. So you may think that you're being more efficient by re-using the same object... but it may not actually be faster. Again, more will be revealed in Chapter 9.

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💡 Financial Modeling Pro-tip

Generally you should default to using immutable types and consider only moving to mutable types in specific circumstances. You'll see some examples in the applications later in the book.

5.4.9. Constructors

Constructors are functions that return a data type (functions will be covered more generally later in the chapter). When we declare a `struct`, an implicit function is defined that takes a tuple of arguments and returns the data type that was declared. In the following example, after we define `MyType` the `struct`, Julia creates a function (also called `MyType`) which takes two arguments and will return the datatype `MyType`:

```
struct MyDate
    year::Int
    month::Int
    day::Int
end

methods(MyDate)

# 2 methods for type constructor:
[1] Main.Notebook.MyDate(year::Int64, month::Int64, day::Int64)
    @ ~/prog/julia-fin-book/foundations-of-programming.qmd:927
[2] Main.Notebook.MyDate(year, month, day)
    @ ~/prog/julia-fin-book/foundations-of-programming.qmd:927
```

Implicit constructors are nice in that you don't have to define a default method and the language does it for you. Sometimes there's reasons to want to control how an object is created, either for convenience or to enforce certain restrictions.

We can use an inner constructor (i.e. inside the `struct` block) to enforce restrictions:

```
struct MyDate
    year::Int
    month::Int
    day::Int

    function MyDate(y,m,d)
        if ~(m in 1:12)
```

```

        error("month is not between 1 and 12")
else if ~(d in 1:31)
    error("day is not between 1 and 31")
else
    return new(y,m,d)
end

end

```

And outer constructors are simply functions defined that have the same name as the data type , but are not defined inside the struct block. Extending the MyDate example, say we want to provide a default constructor for if no day is given such that the date returns the 1st of the month:

```

function MyDate(y,m)
    return MyDate(y,m,1)
end

```

5.5. Functions

Functions are a set of expressions that take inputs and return specified outputs.

5.5.1. Special Operators

Operators are the glue of expressions which combine values. We've already seen quite a few, but let's develop a little bit of terminology for these functions.

Unary operators are operators which only take a single argument. Examples include the ! which negates a boolean value or - which negates a number:

```
!true, -5
```

```
(false, -5)
```

Binary operators take two arguments and are some of the most common functions we encounter, such as + or - or >:

```
1 + 2, 1 - 2, 1 > 2
```

```
(3, -1, false)
```

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The above unary and binary operators are special kinds of functions which don't require the use of parenthesis. However, they can be written with parenthesis for greater clarity:

```
!(true), -(5), +(1, 2), -(1, 2)
```

```
(false, -5, 3, -1)
```

In Julia, we distinguish between **functions** which define behavior that maps a set of inputs to outputs. But a single function can adapt its behavior to the arguments themselves. We have just seen the function - be used in two different ways: negation and subtraction depending on whether it had one or two arguments given to it. In this way there is a conceptual hierarchy of functions that complements the hierarchy we have discussed in relation to types:

- - is the overall function
- -(x) is a unary function which negates its values, -(x,y) subtracts y from x
- Specific methods are then created for each combination of concrete types:
-(x::Float64) is a different method than -(x::Int)

Methods are specific compiled versions of the function for specific types. This is important because at a hardware level, operations for different types (e.g. integers versus floating point) differ considerably. By optimizing for the specific types Julia is able to achieve nearly ideal performance without the same sacrifices of other dynamic languages. We will develop more with respect to methods when we talk about dispatch in Chapter 8.

For example, factorial would be referred to as the *function*, while specific implementations are called *methods*. We can see all of the methods for any function with the `methods` function, like the following for factorial which has implementations taking into account the specialized needs for different types of arguments:

```
methods(factorial)
```

```
# 7 methods for generic function "factorial" from Base:
[1] factorial(n::UInt128)
    @ combinatorics.jl:26
[2] factorial(n::Int128)
    @ combinatorics.jl:25
[3] factorial(x::BigFloat)
    @ Base.MPFR mpfr.jl:769
[4] factorial(n::BigInt)
    @ Base.GMP gmp.jl:703
[5] factorial(n::Union{Int16, Int32, Int8, UInt16, UInt32, UInt8})
```

```

@ combinatorics.jl:33
[6] factorial(n::Union{Int64, UInt64})
    @ combinatorics.jl:27
[7] factorial(n::Integer)
    @ intfuncs.jl:1137

```

5.5.2. Defining Functions

Functions more generally are defined like so:

```

function functionname(arguments)
    # ... code that does things
end

```

Here's an example which returns the difference between the highest and lowest values in a collection:

```

function value_range(collection)

    hi = maximum(collection)
    lo = minimum(collection)
    return hi - lo
end

```

- ① `return` is optional but recommended to convey to readers of the program where you expect your function to terminate and return a value.

5.5.3. Defining Methods on Types

Here's another example of a function which calculates the distance between a point and the origin:

```

function distance(point)
    return sqrt(point.x^2 + point.y^2)
end

```

- ① A function block is declared with the name `distance` which takes a single argument called `point`
 ② We compute the distance formula for a point with `x` and `y` coordinates. The `return` value make explicit what value the function will output.

```
distance (generic function with 1 method)
```

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Note

An alternate, simpler function syntax for `distance` would be:

```
distance(point) = sqrt(point.x^2 + point.y^2)
```

However, we might at this point note a flaw in our function's definition if we think about the various `Coordinates` we defined earlier: our definition would currently only work for `Point2D`. For example, if we try a `Point3D` we will get the wrong answer:

```
distance(Point3D(1, 1, 1))
```

```
1.4142135623730951
```

The above value should be $\sqrt{3}$, or approximately 1.73205.

What we need to do is define a refined distance for each type, which we'll call `dist` to distinguish from the earlier definition.

```
"""
dist(point)

The euclidean distance of a point from the origin.
"""
dist(p::Point2D) = sqrt(p.x^2 + p.y^2)
dist(p::Point3D) = sqrt(p.x^2 + p.y^2 + p.z^2)
dist(p::Polar2D) = p.r

dist (generic function with 3 methods)
```

Now our result will be correct:

```
dist(Point3D(1, 1, 1))
```

```
1.7320508075688772
```

This is referred to **dispatching** on the argument types. Julia will look up to find the most specific method of a function for the given argument types, and falling back to a generic implementation if one is defined.

In Chapter 8 we will see how dispatch (single and multiple) can provide very nice abstractions to simplify the design of a model.

Docstrings (Documentation Strings)

Notice the strings preceding the definition of `dist`. In Julia, putting a string ("...") or string literal ("""...""") right above the definition will allow Julia to recognize the string as documentation and provide it to the user in help mode (`?@sec-help-mode`) and/or have a documentation tool create a webpage or PDF documentation resource.

Defining Methods for Parametric Types

We learned that `Float64 <: Real` in the type hierarchy. However, note that `Tuple{Float64}` is not a sub-type of `Tuple{Real}`. This is called being **invariant** in type theory... but for our purposes this just practically means that when we define a method we need to specify that we want it to apply to all subtypes.

For example, `myfunction(x :: Tuple{Real})` would *not* be called if `x` was a `Tuple{Float64}` because it's not a sub-type of `Tuple{Real}`. To act the way we want, would define the method with the signature of `myfunction(Tuple{<:Real})` or `myfunction{T}(Tuple{T})` where `{T<:Real}`.

5.5.4. Keyword Arguments

Keyword arguments are arguments that are passed to a function but do not use *position* to pass data to functions but instead used named arguments. In the following example, `filepath` is a **positional argument** while the two arguments after the semicolon (;) are keyword arguments.

```
function read_data(filepath; normalize_names, has_header_row)
    # ... function would be defined here
end
```

The function would need to be called and have the two keyword arguments specified:

```
read_data("results.csv"; normalize_names=true, has_header_row=false)
```

5.5.5. Default Arguments

We are able to define default arguments for both positional and keyword arguments via an assignment expression in the function signature. For example, we can make it so that the user need not specify all the options for each call. Modifying the prior example so that typical CSVs work with less customization from the user:

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```
function read_data(filepath;
    normalizeNames = true,
    hasheader = false
)
```

This is a simplified example, but if you look at the documentation for most data import packages you'll see a lot of functionality defined via keyword arguments which have sensible defaults so that most of the time you need not worry about modifying them.

5.5.6. Anonymous Functions

Anonymous functions are functions that have no name and are used in contexts where the name does not matter. The syntax is $x \rightarrow \dots$ expression with $x \dots$. As an example, say that we want to create a vector from another where each element is squared. `map` applies a function to each member of a given collection:

```
v = [4, 1, 5]
map(x → x^2, v) (1)
```

① The $x \rightarrow x^2$ is the anonymous function in this example.

```
3-element Vector{Int64}:
```

```
16
1
25
```

They are often used when constructing something from another value, or defining a function within optimization or solving routines.

5.5.7. First Class Nature

Functions in many languages including Julia are **first class** which means that functions can be assigned and moved around like data variables.

In this example, we have a general approach to calculate the error of a modeled result compared to a known truth. In this context, there are different ways to measure error of the modeled result and we can simplify the implementation of loss by keeping the different kinds of error defined separately. Then, we can assign a function to a variable and use it as an argument to another function:

```

function square_error(guess, correct)
    (correct - guess)^2
end

function abs_error(guess, correct)
    abs(correct - guess)
end

# obs meaning "observations"
function loss(modeled_obs,
    actual_obs,
    loss_function
)
    sum(
        loss_function.(modeled_obs, actual_obs)
    )
end

let
    a = loss([1, 5, 11], [1, 4, 9], square_error)          ②
    b = loss([1, 5, 11], [1, 4, 9], abs_error)            ③
    a, b
end

```

- ① `loss_function` is a variable that will refer to a function instead of data.
- ② Using a `let` block here is good practice to not have temporary variables `a` and `b` scattered around our workspace.
- ③ Using a function as an argument to another function is an example of functions being treated as “first class”.

(5, 3)

5.5.8. Broadcasting

Looking at the prior definition of `dist`, what if we wanted to compute the squared distance from the origin for a set of points? If those points are stored in an array, we can **broadcast** functions to all members of a collection at the same time. This is accomplished using the **dot-syntax** as follows:

```

points = [Point2D(1, 2), Point2D(3, 4), Point2D(6, 7)]
dist.(points) .^ 2

```

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```
3-element Vector{Float64}:
5.000000000000001
25.0
85.0
```

Let's unpack that a bit more:

1. The `.` in `dist.(points)` tells Julia to apply the function `dist` to each element in `points`.
2. The `.` in `.^` tells Julia to square each values as well

Why broadcasting is useful:

1. Without needing any redefinition of functions we were able to transform the function `dist` and exponentiation (`^`) to work on a collection of data. This means that we can keep our code simpler and easier to reason about (operating on individual things is easier than adding logic to handle collections of things).
2. When multiple broadcasted operations are joined together, Julia can **fuse** the operations so that each operation is performed at the same time instead of each step sequentially. That is, if the operation were not fused, the computer would first calculate `dist` for each point, and then apply the square on the collection of distances. When it's fused, the operations can happen at the same time without creating an interim set of values.

 Note

For readers coming from numpy-flavored Python or R, broadcasting is a way that can feel familiar to the array-oriented behavior of those two languages. Once you feel comfortable with Julia in general, you may find yourself relaxing and relying less on array-oriented design and instead picking whichever iteration paradigm feels most natural for the problem at hand: loops or broadcasting over arrays.

5.5.8.1. Broadcasting Rules

What happens if one of the collections is not the same size as the others? When broadcasting, singleton dimensions (i.e. the 1 in $1 \times N$, “1-by- N ”, dimensions) will be expanded automatically when it makes sense. For example, if you have a single element and a one dimensional array, the single element will be expanded in the function call without using any additional memory (if that dimension matches one of the dimensions of the other array).

The rules with an $M \times N$ and a $P \times Q$ array:

- either (M and P) or (N and Q) need to be the same, *and*

- one of the non-matching dimensions needs to be 1

Some examples might clarify. This 1x1 element is being combined with a 4x1, so there is a compatible dimension (N and Q match, M is 1):

```
2 .^ [0, 1, 2, 3]
```

4-element Vector{Int64}:

```
1  
2  
4  
8
```

Here, this 1x3 works with the 2x3 (N and Q match, M is 1)

```
[1 2 3] .+ [1 2 3; 4 5 6]
```

2x3 Matrix{Int64}:

```
2 4 6  
5 7 9
```

This 3x1 isn't compatible with this 2x3 array (neither M and P nor N and Q match)

```
#| error: true  
[1, 2, 3] .+ [1 2 3; 4 5 6]
```

This 2x4 isn't compatible with the 2x3 (M and P match, but N nor Q is 1):

```
#| error: true  
[1 2; 3 4] .+ [1 2 3; 4 5 6]
```

5.5.8.2. Not Broadcasting

What if you do not want the array to be used element-wise when broadcasting? Then you can wrap the array in a Ref, which is used in broadcasting to make the array be treated like a scalar. In the example below, in(needle, haystack) searches a collection (haystack) for an item (needle) and returns true or false if the item is in the collection:

```
in(4, [1 2 3; 4 5 6])
```

```
true
```

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What if we had an array of things (“needles”) that we wanted to search for? By default, broadcasting would effectively split the array up into collections of individual elements to search:

```
in.([1, 9], [1 2 3; 4 5 6])
```

```
2×3 BitMatrix:  
1 0 0  
0 0 0
```

Effectively, the result above is the result of this broadcasted result:

```
in(1, [1,2,3]) # the first row of the above result  
in(9, [4,5,6])
```

If we were expecting Julia to return `[1,0]` (that the first needle is in the haystack but the second needle is not), then we need to tell Julia not to broadcast along the second array with `Ref`:

```
in.([1, 9], Ref([1 2 3; 4 5 6]))
```

```
2-element BitVector:  
1  
0
```

5.5.9. Passing by Sharing

We often want to share data between scopes, such as between modules or by passing something into a function’s scope. Arguments to a function in Julia are **passed-by-sharing** which means that an outside variable can be mutated from within a function. We can modify the array in the outer scope (scope discussed later in this chapter) from within the function. In this example, we modify the array that is assigned to `v` by doubling each element:

```
v = [1, 2, 3]  
  
function double!(v)  
    for i in eachindex(v)  
        v[1] = 2 * v[i]  
    end  
end  
  
double!(v)
```

```
v
```

```
3-element Vector{Int64}:
6
2
3
```

 Tip

Convention in Julia is that a function that modifies it's arguments has a ! in it's name and we follow this convention in `double!` above. Another example would be the built-in function `sort!` which will sort an array in-place without allocating a new array to store the sorted values.

We won't discuss all potential ways that programming languages can behave in this regard, but an alternative that one may have seen before (e.g. in Matlab) is pass-by-value where a modification to an argument only modifies the value within the scope. Here's how to replicate that in Julia by copying the value before handing it to a function. This time, `v` is not modified because we only passed a copy of the array and not the array itself:

```
v = [1, 2, 3]
double!(copy(v))
v
```

```
3-element Vector{Int64}:
1
2
3
```

5.6. Scope

In projects of even modest complexity, it can be challenging to come up with unique identifiers for different functions or variables. **Scope** refers to the bounds for which an identifier is available. We will often talk about the **local scope** that's inside some expression that creates a narrowly defined scope (such as a function or `let` or `module` block) or the **global scope** which is the top level scope that contains everything else inside of it. Here are a few examples to demonstrate scope.

5. Elements of Programming

```
i = 1  
let  
    j = 3  
    i + j  
end
```

(1)
(2)
(3)

- ① i is defined in the global scope and would be available to other inner scopes.
- ② The let ... end block creates a local scope which inherits the defined global scope definitions.
- ③ j is only defined in the local scope created by the let block.

4

In fact, if we try to use j outside of the scope defined above we will get an error:

j

```
UndefVarError: UndefVarError(:j, Main.Notebook)  
UndefVarError: 'j' not defined in 'Main.Notebook'  
Suggestion: check for spelling errors or missing imports.
```

💡 Tip

let blocks are a great way to organize your code in bite-sized chunks or to be able to re-use common variable names without worrying about conflict. Here's an example of using let blocks to:

1. Perform intermediate calculations without fear of returning a partially modified variable
2. Re-use common variable names

```
bonds = let  
    df = CSV.read("bonds.csv", DataFrame)  
    df.issuer = lookup_issuer(df.CUSIP)  
    df  
end  
  
mortgages = let  
    df = CSV.read("bonds.csv", DataFrame)  
    df.issuer = lookup_issuer(df.CUSIP)  
    df  
end
```

If we were running this interactively (e.g. step-by step in VS Code, the REPL, or notebooks) then these two code blocks will run completely and will run separately. The short, descriptive name `df` is reused, but there's no chance of conflict. We also can't easily run the block of code (`let ... end`) and get a partially evaluated result (e.g. getting the dataframe before it has been appropriately modified to add the `issuer` column).

Here is an example with functions:

```
x = 2
base = 10
foo() = base^x
foo(x) = base^x
foo(x, base) = base^x
foo(), foo(4), foo(4, 4)
```

- ① Both `base` and `x` are inherited from the global scope.
- ② `x` is based on the local scope from the function's arguments and `base` is inherited from the global scope.
- ③ Both `base` and `x` are defined in the local scope via the function's arguments.

`(100, 10000, 256)`

In Julia, it's always best to explicitly pass arguments to functions rather than relying on them coming from an inherited scope. This is more straight-forward and easier to reason about and it also allows Julia to optimize the function to run faster because all relevant variables coming from outside the function are defined at the function's entry point (the arguments).

5.6.1. Modules and Namespaces

Modules are ways to encapsulate related functionality together. Another benefit is that the variables inside the module don't "pollute" the **namespace** of your current scope. Here's an example:

```
module Shape
    struct Triangle{T}
        base::T
        height::T
    end

```

5. Elements of Programming

```
function area(t::Triangle)          (2)
    return 1 / 2 * t.base * t.height
end
end

t = Shape.Triangle(4, 2)            (3)
area = Shape.area(t)               (4)
```

- ① module defines an encapsulated block of code which is anchored to the namespace Shape
- ② Here, area a *function* defined within the Shape module.
- ③ Outside of Shape module, we can access the definitions inside via the Module.Identifier syntax.
- ④ Here, area is a *variable* in our global scope that *does not* conflict with the area defined within the Shape module. If Shape.area were not within a module then when we said area = ... we would have reassigned area to no longer refer to the function and instead would refer to the area of our triangle.

4.0

i Note

Summarizing related terminology:

- A **module** is a block of code such as module MySimulation ... end
- A **package** is a module that has a specific set of files and associated metadata. Essentially, it's a module with a Project.toml file that has a name and unique identifier listed, and a file in a src/ directory called MySimulation.jl
 - **Library** is just another name for a package, and the most common context this comes up is when talking about the packages that are bundled with Julia itself called the **standard library** (stdlib).

6. Functional Abstractions

The purpose of abstraction is not to be vague, but to create a new semantic level in which one can be absolutely precise. - Edsger Dijkstra (1972)

6.1. In this section

Demonstrate different approaches to a problem which gradually introduce more reusable or general techniques. These techniques will allow for constructing sophisticated models while maintaining consistency and simplicity. Imperative programming, functional programming, and recursion.

6.2. Introduction

This chapter will center around a simple task: calculate the present value of a portfolio of a single fixed, risk-free, coupon-paying bonds under two different interest rate environments. The focus will be on describing different approaches to this problem, not be adding complexity to the problem (e.g. no getting into credit spreads, settlement timing, etc.).

Mathematically, the problem is to determine the Present Value, where:

$$\text{Present Value} = \sum \text{Cashflow}_t \times \text{Discount Factor}_t$$

Where

$$\text{Discount Factor}_t = \prod^t \frac{1}{1 + \text{Discount Rate}_i}$$

```
cf_bond = [10, 10, 10, 10, 110];
rate = [0.05, 0.06, 0.05, 0.04, 0.05];
```

① The rates are the one year forward rates for time 0, 1, 2, etc.

6. Functional Abstractions

We will focus on this first discount vector, and introduce more scenarios later in the chapter.

We will repeatedly solve the same problem before extending it to more examples. It may feel repetitive but the focus here is not the problem, but rather the variations between the different approaches.

6.3. Imperative Style

One of the most familiar styles of programming is called **imperative** (or **procedural**), where we provide explicit, step-by-step instructions to the computer. The programmer defines the data involved and how that data moves through the program one step at a time. It commonly uses loops to perform tasks repeatedly or across a set of data. The program's **state** (assignment and logic of the program's variables) is defined and managed by the programmer explicitly.

Here's an imperative style of calculating the present value of the bond.

```
let
    pv = 0.0
    discount = 1.0

    for i in 1:length(cf_bond)
        discount = discount / (1 + rate[i])
        pv = pv + discount * cf_bond[i]
    end
    pv
end
```

- ① Declare variables to keep track of the discount rate and running total for the present value `pv`
- ② Loop over the length of the cashflow vector.
- ③ At each step of the loop, look up (via index `i`) update the discount factor to account for the prevailing rate and add the discounted cashflow to the running total present value.

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This style is simple, digestible, and clear. If we were performing the calculation by hand, it would likely follow a pattern very similar to this. Look up the first cashflow and discount rate, compute a discount factor, and subtotal the value. Repeat for the next set of values.

6.3.1. Iterators

Note that in the prior code example we defined an index variable `i` and had to manually define the range over which it would operate (1 through the length of the bond's cashflow vector). A couple of reasons this could be sub-optimal:

1. We are making the *assumption* that the indices of the vectors start with one, when in reality Julia arrays *can* be defined to start at 0 or another arbitrary index.
2. We manually perform the lookup of the values within each iteration.

We can solve the first one (partially) by letting Julia return an iterable set of values corresponding to the indices of the `cf_bond` vector. This is an example of an **iterator** which is an object upon which we can repeatedly ask for the next value until it tells us to stop.

By using `eachindex` we can get the indices of the vector since Julia already knows what they are:

```
eachindex(cf_bond)
```

```
Base.OneTo(5)
```

Lazy Programming

The result, `Base.OneTo(5)` is a **lazy** object which represents a collection that does not get fully instantiated until asked to (which may not actually be necessary). Many (most?) iterators are lazy but we can interact with them without fully instantiating the data that they represent. **Instantiating** means fully loading the values into memory.

An analogy is that we can write the “set of all numbers from 1 to 100” without writing out each of the 100 numbers, but we are referring to the same thing.

An example of operating on a lazy iterator, is that we could find the largest index:

```
maximum(eachindex(cf_bond))
```

```
5
```

The point is if we have an object that *represents* a set, we need not actually enumerate each element of the set to interact with it.

We can fully instantiate an iterator with `collect`

```
collect(eachindex(cf_bond))
```

```
5-element Vector{Int64}:
```

```
1  
2
```

6. Functional Abstractions

```
3  
4  
5
```

Laziness is generally a good thing in programming because sometimes it can be computationally or memory expensive to fully instantiate the collection of interest (this will be discussed further in @#sec-hardware).

And when used in context:

```
let  
    pv = 0.0  
    discount = 1.0  
  
    for i in eachindex(cf_bond)  
        discount = discount / (1 + rate[i])  
        pv = pv + discount * cf_bond[i]  
    end  
    pv  
end
```

```
121.48888490821489
```

Here Julia gave us the index associated with the bond cashflows, but we are still looking up the values (why not just ask for the values instead of their index?) as well as assuming that the indices are the same for the discount rates.

We can get the value and the associated index with enumerate:

```
collect(enumerate(cf_bond))
```

```
5-element Vector{Tuple{Int64, Int64}}:  
(1, 10)  
(2, 10)  
(3, 10)  
(4, 10)  
(5, 110)
```

This would allow us to skip the step of needing to look up the bond's cashflows. However, we can go even further by just asking for value associated with both collections. With `zip` (named because it's sort of like zipping up two collections together), we get an iterator that provides the values of the underlying collections:

6.3. Imperative Style

```
collect(zip(cf_bond, rate))

5-element Vector{Tuple{Int64, Float64}}:
(10, 0.05)
(10, 0.06)
(10, 0.05)
(10, 0.04)
(110, 0.05)
```

This provides the simplest implementation of the imperative approaches:

```
let
    pv = 0.0
    discount = 1.0

    for (cf, r) in zip(cf_bond, rate)
        discount = discount / (1 + r)
        pv = pv + discount * cf
    end
    pv
end
```

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The primary downsides to iterative approaches to algorithms are:

1. Needing to keep track of state is fine in simple cases, but can quickly become difficult to reason about and error prone as the number and complexity of variables grows.
2. Program flow is explicitly stated, leaving fewer places that the compiler can automatically optimize or parallelize.

💡 Tip

In the imperative style, we mentioned needing to explicitly handle program state. In general, it's advisable to minimize as many temporary state variables as possible - more mutability tends to produce more complex, difficult to maintain code. An example of maintaining state in the examples above is keeping track of the current index as well as interim `pv` and `discount` variables.

Avoiding modifying values is often avoidable by restructuring the logic, using functional techniques, or finding the right abstractions. However, sometimes for performance reasons, clarity, or expediency you may find modifying state to be the

6. Functional Abstractions

preferred option and that's okay.

6.4. Functional Techniques and Terminology

Functional programming is a paradigm which attempts to minimize state via composing functions together.

Table 6.1 introduces a set of core functional methods to familiarize yourself with. Note that anonymous functions (Section 5.5.6) are used frequently to define intermediary steps.

Table 6.1.: Important Functional Methods.

Function	Description	Example
<code>map(f, v)</code>	Apply function <code>f</code> to each element of the collection <code>v</code> .	<code>map(</code> <code>x→x^2,</code> <code>[1,3,5]</code> <code>) # [1,9,25]</code>
<code>reduce(op, v)</code>	Apply binary <code>op</code> to pairs of values, reducing the dimension of the collection <code>v</code> . Has a couple of important, optional keyword arguments to note (which also apply to other variants of <code>reduce</code> below): <ul style="list-style-type: none">• <code>init</code> defines the identity element (e.g. the initial value of <code>+</code> and <code>*</code> is <code>0</code> and <code>1</code> respectively)• <code>dims</code> defines which dimension to reduce across (if the dimension of <code>v</code> is more than one).	<code>reduce(</code> <code>*</code> , <code>[1,3,5]</code> <code>) # 15</code>

6.4. Functional Techniques and Terminology

Function	Description	Example
<code>mapreduce(op,f,v)</code>	Maps f over collection v and returns a reduced result using op.	<code>mapreduce(*, x→x^2, [1,3,5]) # 35</code>
<code>foldl(op,v)</code>	Like reduce, but applies op from left to right (foldl) or right to left (foldr). Also has mapfoldl and mapfoldr versions.	<code>foldl(*, [1,3,5]) # 15</code>
<code>accumulate(op,v)</code>	Apply op along v , creating a vector with the cumulative result.	<code>accumulate(+, [1,3,5]) # [1, 4, 9]</code>
<code>filter(f,v)</code>	Apply f along v and return a copy of v with elements where f is true	<code>filter(>=(3), [1,3,5]) # [3, 5]</code>

This paradigm is very powerful in a few ways:

1. It provides a language for talking about what a computation is doing. Instead of “looping over a collection called `portfolio` and calling a `value` function” we can more concisely refer to this as `mapreduce(value,portfolio)`.
2. Often times you are forced to think about the design of the program more deeply, recognizing the core calculations and data used within the model.
3. The compiler is free to apply more optimizations. For example, with `reduce`, the compiler could drive the calculation in any order since the operation is associative.
4. The lack of mutable state.

Let’s build a version of the present value calculation using the functional building blocks described above. We will work up to it by discussing the core functional programming building blocks, culminating in combining `mapreduce` and `accumulate` to do the bond valuation.

6.4.1. `map`

`map` is so named for the mathematical concept of mapping an input to an output. Here, it’s effectively the same thing. We take a collection and use the given function to calculate an output. The size of the output equals the size of the input.

First, we will use `map` to compute the one-period discount factors:

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```
map(x → 1 / (1 + x), rate)
```

5-element Vector{Float64}:

```
0.9523809523809523
0.9433962264150942
0.9523809523809523
0.9615384615384615
0.9523809523809523
```

`map` transforms the `rate` collection by applying the anonymous function $x \rightarrow 1 / (1 + x)$, which is the single period discount factor. This operation is conveyed visually in Figure 6.1.



Figure 6.1.: A diagram showing that `map` creates a new collection mirroring the old one, after applying the given function to each element in the original collection.

Tip

`map` is an absolute workhorse of a function and the authors recommend using it liberally within your code. We find ourselves using `map` frequently, usually avoiding defining an explicit loop (unless we are modifying some existing collection). `map` would likely be a better tool for a loop like this:

```
output = []
```

```

for x in collection
    result = # ... do stuff ...
    push!(output,result)
end
output

```

Instead, `map` simplifies this to:

```

map(collection) do x
    # ... do stuff
end

```

Not only does this have the advantage of being clearer, more concise, and less work, it also lets Julia infer the output type of your computation so you don't have to worry about the type of `output`.

6.4.2. accumulate

`accumulate` takes an operation and a collection and returns a collection where each element is the cumulative result of applying the operation from the first element to the current one. For example, to calculate the cumulative product of the one-period discount factors:

```
accumulate(*, map(x → 1 / (1 + x), rate))
```

```

5-element Vector{Float64}:
0.9523809523809523
0.898472596585804
0.8556881872245752
0.822777103100553
0.7835972410481457

```

This results in a vector of the cumulative discount factors for each point in time corresponding to the given cashflows.

i Note

For `accumulate` and `reduce`, an important, optional value is the `init` (an optional keyword argument), which is the initial value to start the accumulation or reduction. For common operations this **identity element** is already predefined. For example, for `+` the identity is `0` while for `*` it is `1`. The identity element e is the one where for a given binary operation \odot , that $x \odot e = x$.

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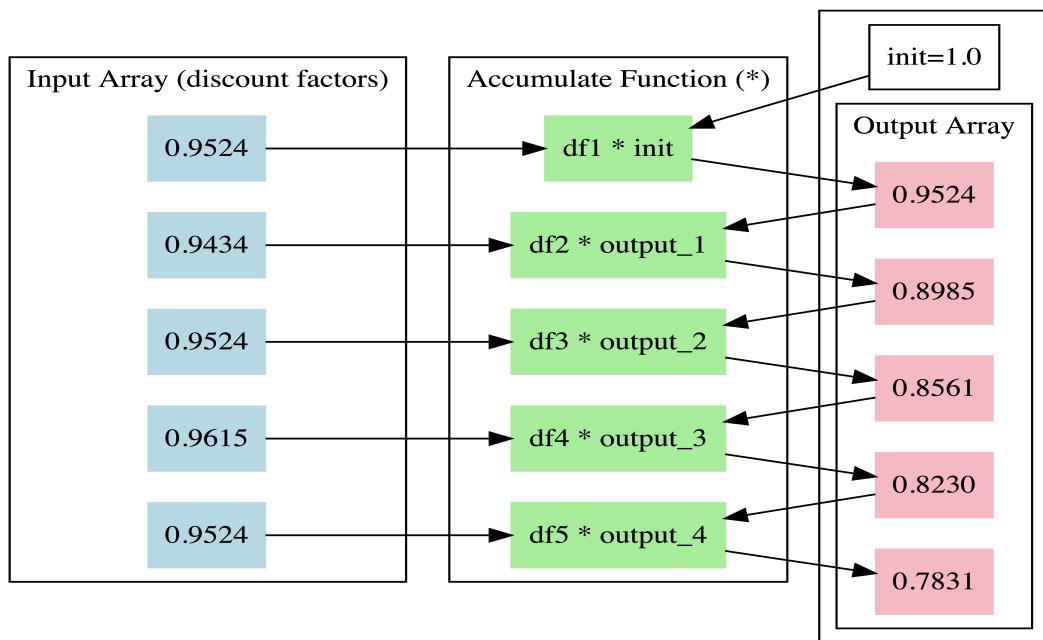


Figure 6.2.: A diagram showing that `accumulate` creates a new collection where each element is the cumulative result of applying the given operation to all previous elements.

Another example is string concatenation. In Julia, two strings are concatenated with `*` (like in mathematics, $a * b$ is also written as ab). The identity element for strings where the binary operation $\odot = *$ is `""`. For example:

```
accumulate(*, ["a", "b", "c"], init="")  
  
3-element Vector{String}:  
"a"  
"ab"  
"abc"
```

This is a taste of a branch of mathematics known as Category Theory, a very rich subject but largely beyond the immediate scope of this book. The category theoretical term for sets of things that work with the binary operator and identity elements as described above is a monoid. There will not be a quiz on this trivia.

6.4.3. reduce

`reduce` takes an operation and a collection and applies the operation repeatedly to pairs of elements until there is only a single value left.

For example, we start with the calculation of the vector of discounted cashflows

```
dfs = accumulate(*, map(x → 1 / (1 + x), rate))  
discounted_cfs = map(*, cf_bond, dfs)
```

```
5-element Vector{Float64}:  
9.523809523809524  
8.98472596585804  
8.556881872245752  
8.22777103100553  
86.19569651529602
```

Then we can sum them with `reduce`:

```
reduce(+, discounted_cfs)
```

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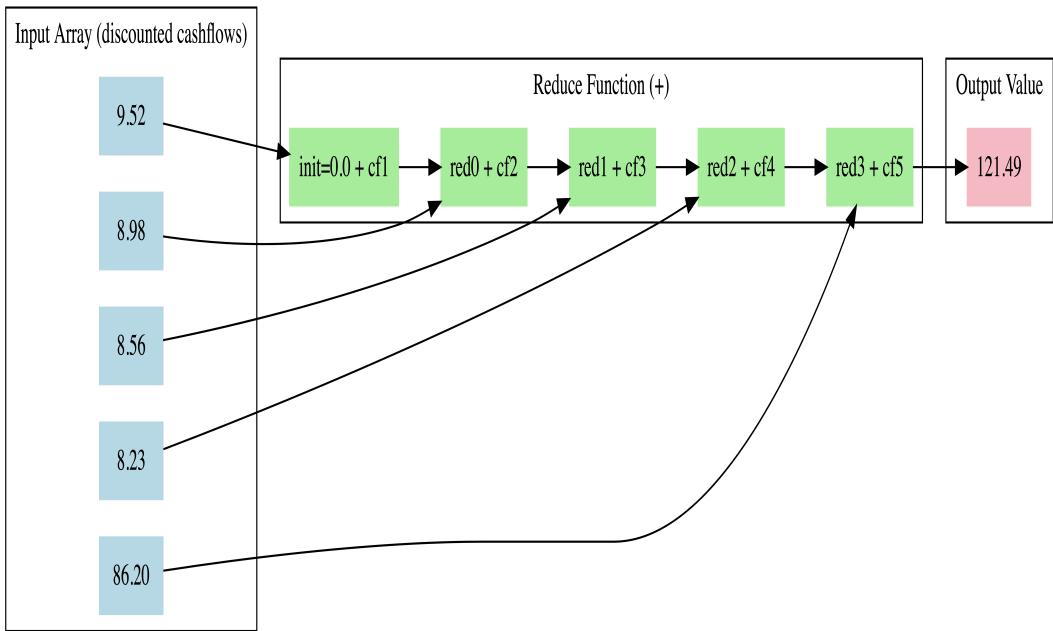


Figure 6.3.: A diagram showing how `reduce` applies the given operation to pairs of elements, ultimately reducing the collection to a single value.

6.4.4. mapreduce

We can combine `map`, `accumulate` and `reduce` to concisely calculate the present value in a functional style. This calculates the discount factors, applies them to the cashflows with `map`, and sums the result with a reduction:

```
dfs = accumulate(*, map(x → 1 / (1 + x), rate))
mapreduce(*, +, cf_bond, dfs)                                ①
                                                               ②
```

- ① Multiplicatively accumulate a discount factor derived from the rate vector.
- ② Multiply the discount factor and bond cashflows (`map` the multiplication), then sum the result (`additive reduce`).

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Contrast this example with the earlier imperative styles:

- This functional approach is more concise.
- The functions used are more descriptive and obvious (once familiar with them, of course!).
- There is no state that the user/programmer keeps track of.

6.4. Functional Techniques and Terminology

- The compiler is able to potentially optimize the code, as it can deduce that certain operations are associative.

This completes the example of using a functional approach to determine the present value of bond cashflows.

6.4.5. filter

For completeness, we will also cover `filter` even though it's not necessary for the bond cashflow example.

`filter` does what you might think - filter a collection based on some criterion that can be determined as true or false.

For example filtering out even numbers using the `isodd` function:

```
filter(isodd, 1:6)
```

```
3-element Vector{Int64}:
1
3
5
```

Or filtering out things that don't match a criteria:

```
filter(x → ~(x == 5), 1:6)
```

```
5-element Vector{Int64}:
1
2
3
4
6
```

While we didn't need `filter` to calculate a bond's present value in the example above, one can imagine how you may want to filter dates that a bond might pay a cashflow, say last day of a quarter:

```
using Dates
let d = Date(2024, 01, 01)
    filter(d → lastdayofquarter(d) == d, d:Day(1):lastdayofyear(d))
end
```

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```
4-element Vector{Date}:
2024-03-31
2024-06-30
2024-09-30
2024-12-31
```

6.4.6. More Tips on Functional Styles

6.4.6.1. do Syntax for Function Arguments

In more complex situations such as with multiple collections or multi-line logic, there is a clearer syntax that is often used. `do` is a reserved keyword in Julia that creates an anonymous function and passes its arguments to a function like `map`. For example, this (terrible) code which decides if a number is prime. The anonymous function requires a `begin` block since the logic of the function is extended into multiple lines.

```
map(x → begin
    if x == 1
        "prime"
    elseif x == 2
        "not prime"
    elseif x == 3
        "prime"
    elseif x > 4
        "probably not prime"
    end
end,
[1, 2, 3, 10]
)
```

This can be written more cleanly with the `do` syntax:

```
map([1, 2, 3, 10]) do x
if x == 1
    "prime"
elseif x == 2
    "not prime"
elseif x == 3
    "prime"
elseif x > 4
    "probably not prime"
end)
```

6.4. Functional Techniques and Terminology

6.4.6.2. Multiple Collections

`map` and the other functional operators discussed in this section can take multiple arguments. This is convenient if you have multiple arguments to a function:

```
discounts = [0.9, 0.81, 0.73]
cashflows = [10, 10, 10]

map((d, c) → d * c, discounts, cashflows)
```

```
3-element Vector{Float64}:
9.0
8.100000000000001
7.3
```

Or an example with the `do` syntax:

```
map(discounts, cashflows) do d, c
    d * c
end
```

```
3-element Vector{Float64}:
9.0
8.100000000000001
7.3
```

6.4.6.3. Using More Functions

At the risk of sounding obvious, an easy way to make the program more “functional” is to simply use more functions. Do this one thing and it will improve the model’s organization, maintainability, and reduce bugs!

Take the example from earlier:

```
pv = 0.0
discount = 1.0

for (cf, r) in zip(cf_bond, rate)
    discount = discount / (1 + r)
    pv = pv + discount * cf
end
pv
```

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We can easily turn this code into a function so that it can operate on data beyond the single pair of `cf_bond` and `rate` previously defined:

```
function pv(rates,cashflows)
    pv = 0.0
    discount = 1.0

    for (cf, r) in zip(rates, cashflows)
        discount = discount / (1 + r)
        pv = pv + discount * cf
    end
    pv
end
```

(1)

- ① Here, `cf_bond` and `rate` would refer to whatever was passed as arguments to the function instead of any globally defined values.

Now we could use this definition of `pv` on other instances of `rates` and `cashflows`.

6.4.6.4. Mixing Functional And Imperative Styles

One of the best things about Julia is how natural it can be to mix the different styles. Sometimes the best is the mix of both styles and that's one of the benefits of Julia: use the style that's most natural to the problem.

i Flexibility and the Lisp Curse

Lisp (“list processing”) is another, much older language than Julia (created in the 1950s!). One of its claims to fame is how flexible and powerful the tools are within the language to build upon. There’s a couple aspects of this curse that we wish to describe because we can learn from it while Julia is still a relatively young language. Part of the “curse” is that: because there’s so much freedom in what can be expressed in the language, there’s not an obvious “best” way of doing things. This can lead to decision paralysis where you are trying to over-analyze what’s the best way to write part of your code. Our advice: *don’t worry about it!* A working implementation of something is better than an over-optimized idea.

The other part of the “curse” is that because it’s relatively easy to implement so many things from the building blocks that Julia provides and compose them together to do what you want. This has a downside because the general approach to packages is smaller, standalone pieces that you call as needed. For example, consider Python’s Pandas library, upon which Python’s data science community was built. It came bundled with a CSV reader, Excel reader, Database reader, DataFrame type, visualization library, and statistical functions. In Julia, each of

those are separate packages that specialize for the respective topics. This is advantageous in that they can progress independently from one another, you don't have to include functionality that you don't need, and you can mix and match libraries depending on your preference.

6.5. Array-Oriented Styles

Another paradigm is **array-oriented**, which is a style that relies heavily on putting similar data into arrays and operating on the entire array at the same time (as opposed to going element-by-element).

Array-oriented programming is one that is practiced in two main contexts:

1. GPU programming
2. Python numerical computing

The former because GPUs want large blocks of similar data to operate in parallel. The latter is because native Python is too slow for many modeling problems so libraries like NumPy,SciPy, and tensor libraries utilize C++ (or similar) libraries for users to call out to.

Array-oriented programming is not always natural for financial and actuarial applications. Differences in behavior or timing of underlying cashflows can make a set of otherwise similar products difficult to capture in nicely gridded arrays. Nonetheless, certain applications (scenario generation, some valuation routines) fit very naturally into this paradigm. Furthermore, for those that work well it's often a great way to extract additional performance due to the parallelization offered via CPU or GPU array programming.

Table 6.2 shows the bond present value example in this style.

6. Functional Abstractions

Table 6.2.: The two code examples demonstrate the same logic using Julia and Numpy (Python's most popular array package). Julia's broadcasting facilitate an array-oriented style, similar to the approach that would be used with Python's NumPy.

Julia	Python (NumPy)
<pre>cf_bond = [10, 10, 10, 10, 110] rate = [0.05, 0.06, 0.05, 0.04, 0.05] discount_factors = cumprod(1 ./ (1 .+ rate)) result = sum(cf_bond .* discount_factors)</pre>	<pre>import numpy as np cf_bond = np.array([10, 10, 10, 10, 110]) discount_factors = np.array([0.05, 0.06, 0.05, 0.04, 0.05]) result = np.sum(cf_bond * discount_factors)</pre>

The downsides to this style are:

1. Sometimes it is unnatural because of non-uniformity of the data we are working with. For example if the length of the cashflows were shorter than the discount rates, we would have to perform intermediate steps to shorten or lengthen arrays in order to get them to be the same size.
2. A good bit of runtime performance is lost because the computer needs to allocate and fill many intermediate arrays (note how in Table 6.2, the `discount_factors` needs to instantiate an entirely new vector even though it's only temporarily used). See more on allocations in Chapter 9.

6.6. Recursion

A **recursive function** which is a pattern where current steps are defined in a way that depends on previous steps. Typically, an explicit starting condition is also required to be specified.

The Fibonacci sequence is a classic example of a recursive algorithm, with the starting conditions of n specified for the first two steps:

$$F(n) = \begin{cases} 0, & \text{if } n = 0 \\ 1, & \text{if } n = 1 \\ F(n - 1) + F(n - 2), & \text{if } n > 1 \end{cases}$$

In code, this translates into a function definition that refers to itself:

```
function fibonacci(n)
    if n == 0
        return 0
    elseif n == 1
        return 1
    else
        return fibonacci(n-1) + fibonacci(n-2)
    end
end
```

How could a recursive pattern be defined for valuing our bond? A possible pattern is defining the present value to be the discounted value of:

- the current period's cashflow, *plus*
- the accumulated cashflows up to that point in time

Here's how that might be defined:

```
function pv_recursive(rates,cashflows,accumulated_value=0.0,discount_factor=1.0)
    if isempty(cashflows)                                ①
        return accumulated_value
    else
        discount_factor = discount_factor / (1+first(rates))      ②
        av = first(cashflows) * discount_factor + accumulated_value ③
        remaining_rates = rates[begin+1:end]
        remaining_cfs = cashflows[begin+1:end]
        return pv_recursive(remaining_rates,remaining_cfs, av,discount_factor) ④
    end
end
```

- ① Add a terminating condition, that if we have no more cashflows then return the accumulated value.
- ② Decrement the discount factor as we step forward in time.
- ③ Take the prior accumulated value and add the first value in the given cashflows.
- ④ Pass the remaining subset of the cashflow vector, the running total, and the current discount factor to the next call of the recursive function.

`pv_recursive` (generic function with 3 methods)

And an example of its use:

```
pv_recursive(rate,cf_bond)
```

121.48888490821489

6. Functional Abstractions

The recursive pattern often works very nicely for simpler examples. However, more complex logic and conditionals can make this approach unwieldy. Nonetheless, attempting to distill the desired functionality into a single function can be a beneficial thought exercise.

7. Data and Types

I am only one, but I am one. I can't do everything, but I can do something.
The something I ought to do, I can do. And by the grace of God, I will -
Edward Everett Hale (1902)

7.1. In this section

The powerful benefits that using assigning types to data has within the model's system, some examples of utilizing types to simplify a programs logic, and comparing aspects of different type related program organization (such as object oriented design versus composition).

7.2. Using Types to Value a Portfolio

We will assemble the tools and terminology to value a portfolio of assets by leverage types (@sec-data-types). Using the constructs introduced in the prior chapter, we can describe the portfolio valuation as additively reducing the mapped value of assets in the portfolio. If `value` is our valuation function), we are trying to do the following:

```
mapreduce(value,+,portfolio)
```

The challenge is how do design an all-purpose `value` function? In `portfolio`, the assets may be heterogeneous, so we will need to define what the valuation semantics are for the different kinds of assets. To get to our end goal, we will need to:

1. Define the different kinds of assets within our portfolio
2. How the assets are to be valued.

We will accomplish this by utilizing data types.

7. Data and Types

7.3. Benefits of Using Types

As a preview of why we want to utilize types in our program, there are a number of benefits:

1. **Separate concerns.** For example, deciding how to value an option need not know how we value a bond. The code and associated logic is kept distinct which is easier to reason about and to test.
2. **Re-use code.** When a set of types within a hierarchy all share the same logic, then we can define the method at the highest relevant level and avoid writing the method for each possible type. In our simple example we won't get as much benefit here since the hierarchy is simple and the set of types small.
3. **Dispatch on type.** By defining types for our assets, we can use multiple dispatch to define specialized behavior for each type. This allows us to write generic code that works with any asset type, and the Julia compiler will automatically select the appropriate method based on the type of the asset at runtime. This is a powerful feature that enables extensibility and modularity in our code.
4. **Improve readability and clarity.** By defining types for our assets, we make our code more expressive and self-documenting. The types provide a clear indication of what kind of data we are working with, making it easier for other developers (or ourselves in the future) to understand and maintain the codebase.
5. **Enable type safety.** By specifying the expected types for function arguments and return values, we can catch type-related errors at compile time rather than at runtime. This helps prevent bugs and makes our code more robust.

With these benefits in mind, let's start by defining the types for our assets. We'll create an abstract type called `Asset` that will serve as the parent type for all our asset types. If you haven't read it already, Section 5.4.7 is a good reference for details on types at the language level (this section is focused on organization and building up the abstracted valuation process).

7.4. Defining Types for Portfolio Valuation

We will define five types of assets in this simplified universe:

- Cash
- Risk Free Bonds (coupon and zero-coupon varieties)

To do the valuation of these, we need some economic parameters as well: risk free rates for discounting.

Here's the outline of what follows to get an understanding of types, type hierarchy, and multiple dispatch.

7.4. Defining Types for Portfolio Valuation

1. Define the Cash and Bond types.
2. Define the most basic economic parameter set.
3. Define the value functions for Cash and Bonds.

```

## Data type definitions
abstract type AbstractAsset end          ①

struct Cash <: AbstractAsset
    balance::Float64
end

abstract type AbstractBond <: AbstractAsset end      ②

struct CouponBond <: AbstractBond
    par::Float64
    coupon::Float64
    tenor::Int
end

struct ZeroCouponBond <: AbstractBond
    par::Float64
    tenor::Int
end

```

- ① General convention is to name abstract types beginning with `Abstract`...
- ② There can exist an abstract type which is a subtype of another abstract type.
- ③ We define concrete data types (`structs`) with the fields necessary for valuing those assets.

Now to define the economic parameters:

```

struct EconomicAssumptions{T}
    riskfree::T
end

```

This is a parametric type because later on we will vary what objects we use for `riskfree`. For now, we will use simple scalar values, like in this potential scenario:

```
econ_baseline = EconomicAssumptions(0.05)
```

```
EconomicAssumptions{Float64}(0.05)
```

Now on to defining the valuation for `Cash` and `AbstractBonds`. `Cash` is always equal to its balance:

7. Data and Types

```
value(asset::Cash, ea::EconomicAssumptions) = asset.balance
```

value (generic function with 1 method)

Risk free bonds are the discounted present value of the riskless cashflows. We first define a method that generically operates on any fixed bond, all that's left to do is for different types of bonds to define how much cashflow occurs at the given point in time by defining `cashflow` for the associated type.

```
function value(asset::AbstractBond, r::Float64)          (2)
    discount_factor = 1.0
    value = 0.0
    for t in 1:asset.tenor
        discount_factor /= (1 + r)
        value += discount_factor * cashflow(asset, t)      (1)
    end
    return value
end

function cashflow(bond::CouponBond, time)
    if time == bond.tenor
        (1 + bond.coupon) * bond.par
    else
        bond.coupon * bond.par
    end
end

function value(bond::ZeroCouponBond, r::Float64)         (3)
    return bond.par / (1 + r)^bond.tenor
end
```

① $x /= y$, $x += y$, etc. are shorthand ways to write $x = x / y$ or $x = x + y$

② `value` is defined for `AbstractBonds` in general...

③ ... and then more specifically for `ZeroCouponBonds`. This will be explained when discussing “dispatch” below.

value (generic function with 3 methods)

7.4.1. Dispatch

When a function is called, the computer has to decide which method to use. In the example above, when we want to value a `ZeroCouponBond`, does the `value(asset::AbstractBond, r)` or `value(bond::ZeroCouponBond, r)` version get used?

Dispatch is the process of determining the right method to use and the rule is that *the most specific defined method gets used*. In this case, that means that even though our `ZeroCouponBond` is an `AbstractBond`, the routine that will be used is the most specific `value(bond::ZeroCouponBond, r)`.

Already, this is a powerful tool to simplify our code. Imagine the alternative of a long chain of conditional statements trying to find the right logic to use:

```
# don't do this!
function value(asset,r)
    if asset.type == "ZeroCouponBond"
        # special code for Zero coupon bonds
        #
    elseif asset.type == "ParBond"
        # special code for Par bonds
        #
    elseif asset.type == "AmortizingBond"
        # special code for Amortizing Bonds
        #
    else
        # here define the generic AbstractBond logic
    end
end
```

With dispatch, the compiler does this lookup for us, and more efficiently than enumerating a list of possible codepaths.

In our “don’t do this” definition of `value` above, we used a simple scalar interest rate to determine the rate to discount the cash flows. Note how in the definition of `value` for `ZeroCouponBond`, we have defined a *more specific* signature: both the first and second arguments are specific, concrete types. When we call `value(ZeroCouponBond(100.0,3),0.05)`, we avoid the loop that’s defined in the generic case and jump immediate to a more efficient definition of its value. This is dispatching on the combination of types and picking the most relevant (specific) version for what has been passed to it.

Despite the definitions above, the following will error because we haven’t defined a method for `value` which takes as its second argument a type of `EconomicAssumptions`:

7. Data and Types

```
#| error: true
value(ZeroCouponBond(100.0,5),econ_baseline)
```

Let's fix that by defining a method which takes the economic assumption type and just relays the relevant risk free rate to the `value` methods already defined (which take an `AbstractBond` and a scalar `r`).

```
value(bond::AbstractBond,econ::EconomicAssumptions) = value(bond,econ.riskfree)

value (generic function with 4 methods)
```

Now this following works:

```
value(ZeroCouponBond(100.0, 5), econ_baseline)
```

```
78.35261664684589
```

Here's an example of how this would be used:

```
portfolio = [
    Cash(50.0),
    CouponBond(100.0, 0.05, 5),
    ZeroCouponBond(100.0, 5),
]

map(asset→ value(asset,econ_baseline), portfolio)

3-element Vector{Float64}:
50.0
99.99999999999999
78.35261664684589
```

This is very close to the goal that we set out at the end of the section. We can complete it by reducing over the collection to sum up the value:

```
mapreduce(asset → value(asset,econ_baseline), +, portfolio)
```

```
228.3526166468459
```

i Note

This code:

```
mapreduce(asset-> value(asset,econ_baseline), +, portfolio)
```

is more verbose than what we set out do at the start (`mapreduce(value,+,&portfolio)`) due to the two-argument `value` function requiring a second argument for the economic variables. This works well! However, there is a way to define it which avoids the anonymous function, which in some cases will end up needing to be compiled more frequently than you want it to. Sometime we want a lightweight, okay-to-compile-on-the-fly function. Other times, we know it's something that will be passed around in compute-intensive parts of the code. A technique in this situation is to define an object which "locks in" one of the arguments but behaves like the anonymous version. There is a pair of types in the `Base` module, `Fix1` and `Fix2`, which represent partially-applied versions of the two-argument function `f`, with the first or second argument fixed to the value "x".

This is, `Base.Fix1(f, x)` behaves like `y->f(x, y)` and `Base.Fix2(f, x)` behaves like `y->f(y, x)`.

In the context of our valuation model, this would look like:

```
val = Base.Fix2(value,econ_baseline)
mapreduce(val,+,portfolio)
```

228.3526166468459

7.4.1.1. Multiple Dispatch

A more general concept is that of **multiple dispatch**, where the types of *all arguments* are used to determine which method to use. This is a very general paradigm, and in many ways is more extensible than traditional object oriented approaches, (more on that in Section 7.5). What if instead of a scalar interest rate value we wanted to instead pass an object that represented a term structure of interest rates?

Extending the example, we can use a time-varying risk free rate instead of a constant. For fun, let's say that the risk free rate has a sinusoidal pattern:

```
econ_sin = EconomicAssumptions(t -> 0.05 + sin(t) / 100)
```

```
EconomicAssumptions{var"#5#6"}(var"#5#6"())
```

7. Data and Types

Now `value` will not work, because we've only defined how `value` works on bonds if the given rate is a `Float64` type:

```
#| error: true
value(ZeroCouponBond(100.0, 5), econ_sin)
```

We can extend our methods to account for this:

```
function value(bond::ZeroCouponBond, r::T) where {T<:Function}           ①
    return bond.par / (1 + r(bond.tenor))^(bond.tenor)
end                                         ②
```

- ① The `r::T ... where {T<:Function}` says use this method if `r` is any concrete subtype of the (abstract) `Function` type.
- ② `r` is a function, where we call the time to get the zero coupon bond (a.k.a. spot) rate for the given timepoint.

```
value (generic function with 5 methods)
```

Now it works:

```
value(ZeroCouponBond(100.0, 5), econ_sin)
```

```
82.03058910862806
```

The important thing to note here is that the compiler is using the most specific method of the function (`value(bond::ZeroCouponBond, r::T) where {T<:Function}`). Both the types of the arguments are influencing the decision of which method to use. We could go on to define the appropriate method for `CouponBond` to complete the example.

7.5. Object Oriented Design

Object oriented (OO) type systems use the analogy that various parts of the system are their own objects which encapsulate both data and behavior. Object oriented design is often one of the first computer programming abstractions introduced because it is very relatable¹, however there are a number of its flaws in over-relying on OO patterns. Julia does not natively have traditional OO classes and types, but much of OO design can be emulated in Julia except for data inheritance.

¹"Many people who have no idea how a computer works find the idea of object-oriented programming quite natural. In contrast, many people who have experience with computers initially think there is something strange about object oriented systems." - David Robson, "Object Oriented Software Systems" in Byte Magazine (1981).

7.6. Assigning Behavior

We bring up object oriented design for comparison's sake, but think that ultimately choosing a data driven or functional believe is better for financial modeling. Of course, many robust, well used financial models have been built this way but in our experience the abstractions become unnatural and maintenance unwieldy beyond simple examples. We'll now discuss some of the aspects of OO design and why the overuse of OO is not preferred.

Note

For readers without background in OO programming, the main features of OO languages are:

- Hierarchical type structures, which include concrete and abstract (often called classes instead of types).
- Sub-classes inherit both behavior *and* data (in Julia, subtypes only inherit behavior, not data).
- Functions that depend on the type of the object need to be ascribed to a single class and then can dispatch more specifically on the given argument's type.

7.6. Assigning Behavior

Needing to assign methods to a single class can lead to awkward design limitations - when multiple objects are involved in a computation, why dictate that only one of them "controls" the logic?

The `value` function is a good example of this. If we had to assign `value` to one of the objects involved, should it be the economic parameters of the asset contracts? The choice is not obvious at all. Isn't it the market (economic parameters) that determines the value? But then if `value` were to be a method wholly owned by the economic parameters, how could it possibly define in advance the valuation semantics of all types of assets? What if one wanted to extend the valuation to a new asset class? Downstream users or developers would need to modify the economic types to handle new assets they wanted to value. However, because the economic types were owned by an upstream package, they can't be extended this way.

This is an issue with traditional OO designs and that resolves itself so elegantly with multiple dispatch.

7. Data and Types

7.7. Inheritance

We discussed the type hierarchy in Chapter 5 and in most OO implementations this hierarchy comes with inheriting both data *and* behavior. This is different from Julia where subtypes inherit behavior but not data from the parent type.

Inheriting the data tends to introduce a tight coupling between the parent and the child classes in OO systems. This tight coupling can lead to several issues, particularly as systems grow in complexity. For example, changes in the parent class can inadvertently affect the behavior of all its child classes, which can be problematic if these changes are not carefully managed. This is often referred to as the “fragile base class problem,” where base classes are delicate and changes to them can have widespread, unintended consequences.

Another issue with inheritance in OO design is the temptation to use it for code reuse, which can lead to inappropriate hierarchies. Developers might create deep inheritance structures just to reuse code, leading to a scenario where classes are not logically related but are forced into a hierarchy. This can make the system harder to understand and maintain.

Moreover, inheritance can sometimes lead to the duplication of code across the hierarchy, especially if the inherited behavior needs to be slightly modified in different child classes. This goes against the DRY (Don’t Repeat Yourself) principle, which is a fundamental concept in software engineering advocating for the reduction of repetition in code.

7.7.1. Composition over Inheritance

To mitigate some of the problems associated with inheritance, there’s a growing preference for *composition*. Composition involves creating objects that contain instances of other objects to achieve complex behaviors. This approach is more flexible than inheritance as it allows for the creation of more modular and reusable code. There is a general preference for “composition over inheritance” among professional developers these days.

In composition, objects are constructed from other objects, and behaviors are delegated to these contained objects. This approach allows for greater flexibility, as it’s easier to change the behavior of a system by replacing parts of it without affecting the entire hierarchy, as is often the case with inheritance.

Composition looks like this:

```
struct CUSIP
    code::String
end
```

```

struct FixedBond
    coupon::Float64
    tenor::Float64
end

struct FloatingBond
    spread::Float64
    tenor::Float64
end

struct MunicipalBond
    cusip::CUSIP
    fi::FixedBond
end

struct Swap
    float_leg::FloatingBond
    fixed_leg::FixedBond
end

struct ListedOption
    cusip::CUSIP
    #... other data fields
end

struct UnlistedBond
    fi::FixedIncome
end

# define behavior which relies on delegation to components
last_transaction(c::CUSIP) = # ...perform lookup of data
last_transaction(asset) = last_transaction(asset.cusip)

duration(f::FixedIncome) = # ... calculate duration
duration(asset) = duration(asset.fi)

```

In the above example, there are number of asset classes that have CUSIP related attributes (i.e. the 9 character code) and behavior (e.g. being able to look up transaction data). Other assets have fixed income attributes (e.g. calculating a duration). There's no clear hierarchy here.

Composition lets us bundle the data and behavior together without needing complex

7. Data and Types

chains of inheritance.

Note

A CUSIP (Committee on Uniform Security Identification Procedures) number, is a unique nine-character alphanumeric code assigned to securities, such as stocks and bonds, in the United States and Canada. This code is used to facilitate the clearing and settlement process of securities and to uniquely identify them in transactions and records.

8. Higher Levels of Abstraction

“Simple things should be simple, complex things should be possible.” —
Alan Kay (1970s)

8.1. In this section

Why we talk about abstraction as a technique in and of itself, discussion of abstraction at the level of code organization and interfaces.

8.2. Introduction

In programming and modeling, as in mathematics, abstraction permits the definition of interchangeable components and patterns that can be reused. Abstraction is a selective ignorance—focusing on the aspects of the problem that are relevant, and ignoring the others. The last two chapters described what we might call “micro” level abstractions: specific functions or types.

In this chapter, we zoom out and examine some principles that guide good model development, manifesting in architectural concerns such as how different parts of the code are organized, what parts of the program are considered ‘public’ versus ‘private’, and patterns themselves.

Chapter 5 Described a number of tools that we can utilize as interfaces within our model. We use these tools that are provided by our programming language *in service of* the conceptual abstraction described above.

- Functions let us implement behavior, where we need trouble ourselves with the low level details.
- Data types provide a hierarchical structure to provide meaning to things, and to group those things together into more meaningful structures.
- Modules allow us to combine data, and or function, into a related group of concepts which can be shared in different parts of our model

8. Higher Levels of Abstraction

8.3. Principles for Abstraction

Here is a list of some principles that arise when developing a particular abstraction. Not all abstractions serve all of these purposes but generally fit one or more of them.

Table 8.1.: Finding abstractions generally means finding patterns that fit into one of these principles.

Principle	What	Why	Example
Separation of Concerns	Divide the system into distinct parts, each addressing a separate concern	Promote modularity and reduce high degree of dependence (coupling) between components	Separating data retrieval, data processing, and output generation steps in a process
Encapsulation	Hide the internal details of a component and expose only a clean, well-defined set of functionality (interface)	Don't let other parts of the program modify internal data and make the system easier to understand and maintain	Defining a type or module with well defined behavior and responsibility
Composability	Design simple components that can be combined to create more complex behaviors, as opposed to a single component that attempts to handle all behavior.	Promote reuse and allow for the components to be combined creatively	Separate details about economic conditions into different types than contracts/instruments
Generalization	Identify common patterns and create generic components that can be specialized as needed. Often this means identifying the common behavior that arises repeatedly in a model	Avoid duplication and make the system more expressive and extensible	Defining a generic Instrument type that can be specialized for different asset classes

These principles provide guidance for creating abstractions that are modular, reusable, and maintainable. By following these principles, developers can create financial models

8.3. Principles for Abstraction

that are easier to understand, extend, and adapt to changing requirements.

8.3.1. Pragmatic Considerations for Model Design

8.3.1.1. Behavior-Oriented

This strategies is to effectively group together components with a model that behaves similarly. So, in our example of bonds and interest-rate swaps fundamentally, they share many characteristics and are used in very similar ways within a model. Therefore, it might make sense to group them together when developing a model.

8.3.1.2. Domain Expertise

It may be that components of the model require sufficient expertise that different persons or groups are involved in the development. This may warrant separating a models design, So that different groups contributing to the model can focus on any more narrow aspect, Regardless of inherent similarity of components. For example, at a higher vertical level of obstruction, financial derivatives may fall under similar grouping, but sufficient differences exist for equity credit or foreign exchange derivatives that the model should separate those three asset classes for development purposes.

8.3.1.3. Composability versus All-in-One

For some model design goals, it may be warranted to attempt to bundle together more functionality instead of allowing users to compose a functionality that comes from different packages. For example, perhaps a certain visualization of a model result is particularly useful, It is not easy to create from scratch, And virtually everyone using the model, will desire to see the model output visualized that way. Instead of relying on the user to install a separate visualization package and develop the visualization themselves, it could make sense to bundle visualization functionality with a model that is otherwise unconcerned with graphical capabilities.

In general, though it is preferred to try to loosely couple systems, you can pick and choose which components you use and that those components work well together.

8. Higher Levels of Abstraction

8.4. Interfaces

Interfaces are the boundary between different encapsulated abstractions. The user-facing interface is the set of functionality and details that the user of the package or model must consider, which is separate from the intermediate variables, logic, and complexity that may be contained within.

Example of an interface

When looking up a ticker for a market quote, one need not be mindful of the underlying realtime databases, networking, rendering text to the screen, memory management, etc. The interface is “put in symbol, get out number”. By design, there are multiple layers of interfaces and abstractions used under the hood, but the financial modeler need only be actively concerned about the points that he or she comes in contact with, not the entire chain of complexity.

For a financial model this might mean that there is an interface for bonds, or there is an interface for interest-rate swaps. There may be a different interface for calculating risk metrics or visualizing the results.

Financial model this might mean that there is an interface for bonds, or there is an interface for interest-rate swaps. There may be a different interface for calculating risk metrics or visualizing the results. A better system design will separate the concern of visualizing output from the mechanics of a fixed income contract. This is what it means to put boundaries on different parts of a models logic. One of the easiest places to see this is with the available open source packages. There are packages available for visualizations, data frames, file, storage, statistical analysis, etc. for many of these it's easy to see where the natural boundary lies.

However, it's often difficult to find where to draw lines within financial models. For example, should bonds and interest-rate swaps be in separate packages? Or both part of a broader fixed income package? This is where much of the art and domain expertise of the financial professional comes to bear in modeling. There would be no way for a pure software engineer to think about the right design for the system without understanding how underlying components share, similarities or differences and how those components interact.

8.4.1. Defining Good Interfaces

A well-designed interface should follow these principles:

1. **Be minimal and focused.** The interface should provide only the essential functionality needed, without unnecessary clutter or features. This makes the interface

easier to understand and facilitates building the necessary complexity through digestible, composable components.

2. **Be consistent and intuitive.** The interface should use consistent naming conventions, parameter orders, and behaviors. It should match the user's mental model and expectations.
3. **Hide implementation details.** The interface should abstract away the internal complexity and expose only what the user needs to know. This of details allows the implementation to change without affecting users of the interface.
4. **Be documented and contractual.** The interface should clearly specify what inputs it expects and what outputs or behaviors it provides. It forms a contract between the implementation and the users.
5. **Be testable.** A good interface allows the functionality to be easily tested through the public interface, without needing to access internal details.

8.4.2. Interfaces: A Financial Modeling Case Study

As a case study, we'll look at the `FinanceModels.jl` and related packages to discuss some of the background and design choices that went into the functionality. This suite was written by one of the authors and is publically available as set of installable Julia packages.

8.4.2.1. Background

In actuarial work, it is common to need to work with interest rate and bond yield curves to determine current forward rates, estimates of the shape of future yield curves, or discount a series of cashflows to determine a present value. Determining things like "given a par yield curve, what's the implied discount factor for a cashflow at time 10" or "what is the 10 year BBB public corporate rate implied by the current curve in five years' time" is cumbersome at best in a spreadsheet.

For example, to determine the answer to the first one ("a discount factor for time 10") actually requires quite a bit of detail and assumption to derive:

- Reference market data and a specification for how that market data should be interpreted. For example, if given the rate `0.05` for time 10, quoted as a continuous rate or annual effective? Is that a par rate, a zero-coupon bond (spot) rate, or a one-year-forward rate from time 10?
- Smoothing, interpolation, or extrapolation for noisy or sparse data. Should the rates be bootstrapped or fit to a parametrically specified curve?

This is the type of complexity that we wish to save the user from needing to keep front of mind when the primary goal is, e.g., valuation of a stream of riskless life insurance payments, which might look like this:

8. Higher Levels of Abstraction

```
risk_free_rates = [0.05,0.06,...0.06]
tenors = [1/12,3/12,...30]
yield_curve = Yields.Par(risk_free_rates,tenors)

cashflow_vector = [1e6,3e6,...,1e3]
present_value(yield_curve,cashflow_vector)
```

This is very clear from the variable and function names what the purpose and steps in the analysis are. Imagine starting with rates and cashflows in a spreadsheet, needing to perform the bootstrapping, interpolation, and discounting before getting to the simple present value sought in the analysis. What can be, with the right abstractions, distilled into five lines of code would take hundreds of cells in a spreadsheet. Providing abstractions like this at the hand of financial modelers is a productivity multiplier.

8.4.2.2. Initial Versions

There were two main abstractions to talk about from early versions of the packages.

8.4.2.2.1. Rates

Utilizing the benefit of the type system, it was decided that it would be most useful to represent rates not as simple floating point numbers (e.g. 0.05) but instead with dedicated types to distinguish between rate conventions. The abstract type `CompoundingFrequency` had two subtypes: `Continuous` and `Periodic` so that a 5% rate compounded continuously versus an effective per period rate would be distinguished via `Continuous(0.05)` versus `Periodic(0.05,1)`. The two could be converted between by extending the built-in `Base.convert` function.

This was useful because once rates were converted into `Rates` within the ecosystem, that data contained within itself characteristics that could distinguish how downstream functionality should treat the rates.

8.4.2.2.2. Yield Curves

At first, only bootstrapping was supported as a method to construct curve objects. This required that there was only one rate given per time period (no noisy data) and only supported linear, quadratic, and cubic splines.

Further, there was a specific constructor for different common types of instruments. From the old documentation:

- `Yields.Zero(rates,maturities)` using a vector of zero, or spot, rates
- `Yields.Forward(rates,maturities)` using a vector of one-period
- `Yields.Constant(rate)` takes a single constant rate for all times

- `Yields.Par(rates,maturities)` takes a series of yields for securities priced at par. Assumes that maturities ≤ 1 year do not pay coupons and that after one year, pays coupons with frequency equal to the CompoundingFrequency of the corresponding rate.
- `Yields.CMT(rates,maturities)` takes the most commonly presented rate data (e.g. Treasury.gov) and bootstraps the curve given the combination of bills and bonds.
- `Yields.OIS(rates,maturities)` takes the most commonly presented rate data for overnight swaps and bootstraps the curve.

This covered a lot of lightweight use-cases, but made a lot of implicit assumptions about how the given rates should be interpreted.

8.4.2.3. The Birth of FinanceModels

There were a multiple of insights that led to a more flexible interface in more recent versions.

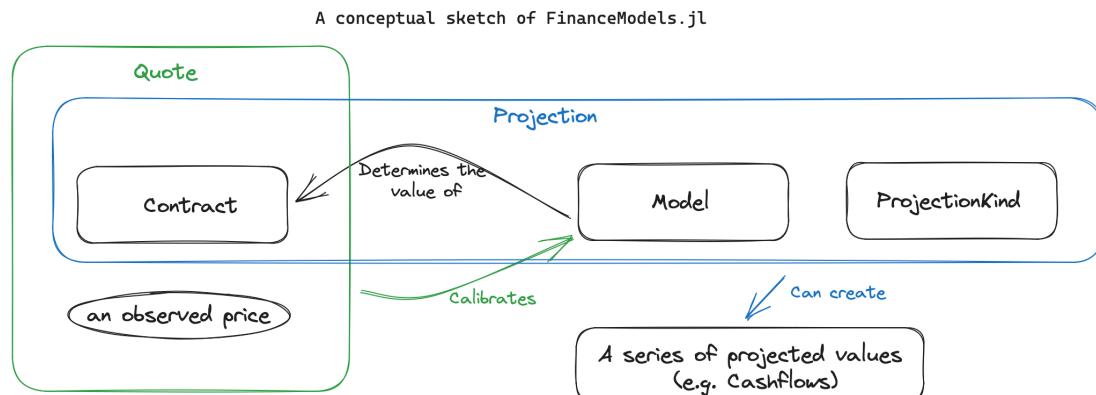


Figure 8.1.: A conceptual sketch of `FinanceModels.jl` components.

First, realizing that yield curves were just a particular kind of model - one that used interest rates to discount cashflows. But you can have different kinds of models - such as Black-Scholes option valuation or a Monte Carlo valuation approach. Likewise, the cashflows need not simply be a vector of floating point values, and instead it could be the representation of a generic financial contract. As long as the model knew how to value it, an appropriate present value could be derived.

Where previously it was:

```
present_value(yield_curve,cashflow_vector)
```

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Now, it was

```
present_value(model,contract)
```

Second, that a model was simple some generic box that had been “fit” to previously observed prices for similar types of contracts we would be trying to value in the model. The combination of a contract and a price constituted a “quote” and with multiple quotes a model could be fit using various algorithms.

With these changes, the package that was originally called Yields.jl was renamed to FinanceModels.jl. The updated code from the earlier example now would be implemented like this:

```
risk_free_rates = [0.05,0.06,...0.06]
tenors = [1/12,3/12,...30]
quotes = ParYield.(risk_free_rates,tenors)
model = fit(Spline.Cubic(),quotes,Fit.Bootstrap())

cashflow_vector = [1e6,3e6,...,1e3]
present_value(model,cashflow_vector)
```

It's slightly more verbose, but notice how much more powerful and extensible `fit(Spline.Cubic(), quotes, Fit.Bootstrap())` is than `Yields.Par(risk_free_rates, tenors)`. The end result is the same, but now the same package and interface can clearly interchange other options, such as a NelsonSiegelSvensson curve instead of a spline. And the quotes could be a combination of observed bonds of different technical parameters (though still sharing characteristics which make it relevant for the model being constructed).

The same pattern also applies for option valuation, such as this example of vanilla euro options with an assumed constant volatility assumption:

```
a = Option.EuroCall(CommonEquity(), 1.0, 1.0)                                ①
b = Option.EuroCall(CommonEquity(), 1.0, 2.0)

qs = [
    Quote(0.0541, a),
    Quote(0.072636, b),
]

model = Equity.BlackScholesMerton(0.01, 0.02, Volatility.Constant())            ③
m = fit(model, qs)                                                               ④
present_value(m,qs[1].instrument)                                                 ⑤
```

- ① The arguments to EuroCall are the underlying asset type, strike, and maturity time.
- ② A vector of observed option prices.
- ③ A BSM model with a given risk free rate, dividend yield, and a to-be-fit constant volatility component.
- ④ Fits the model and derives an approximate volatility of 0.15 .
- ⑤ Values the contract and in such a simple, noiseless model we recover the original price of 0.0541

With a consistent interface able to handle a wide variety of situations, the modeler is free to expand the model in new directions of analysis with the built in functionality allowing him or her to compose pieces together that was not possible with the less abstracted design. For example, the equity option example had no parallel when all of the available constructors were `Yields.Zero` or `Yields.Par` and would have required a completely from-scratch implementation with newly defined functions.

Further, and critically, the new design allows modelers to create their own models or contracts¹ and extend the existing methods rather than needing to create their own: the function signature `fit(model, quotes)` handles a very wide variety of cases, as does `present_value(model, contract)`.

8.5. Macros & Homoiconicity

We've talked about transforming data and restructuring logic in order to make the model more effective. We can go still deeper!(Or is it higher level?) We can actually abstract the process of writing code itself! This subject is a bit advanced, so we are simply going to introduce it because you will likely find many convenient instances of it as a *user* even if you never find a need to implement this yourself.

Homoiconicity refers to the property of a programming language where the language's code can be represented and manipulated as a data structure in the language itself. In other words, the code is data and can be treated as such. This enables powerful metaprogramming (i.e. code that writes other code) capabilities, where code can be generated or transformed during the compilation process.

Macros are a metaprogramming feature that leverage homoiconicity in Julia. They allow the programmer to write code that generates or manipulates other code at compile-time. Macros take code as input, transform it based on certain rules or patterns, and return the modified code which then gets compiled.

For example, a built-in macro is `@time` which will measure the elapsed runtime for a piece of code².

¹And projections, which is handled by defining a `ProjectionKind`, such as a cashflow or accounting basis.
This topic is covered in more detail in the `FinanceModels.jl` documentation.

²(`time?`) is a simple, built-in function. For true benchmarking purposes, see `?@sec-benchmarking`.

8. Higher Levels of Abstraction

```
@time exp(rand())
```

Will effectively expand to:

```
t0 = time_ns()
value = exp(rand())
t1 = time_ns()
println("elapsed time: ", (t1-t0)/1e9, " seconds")
value
```

Here it is when we run it:

```
@time exp(rand())
```

```
0.000003 seconds
```

```
2.219604464359098
```

8.5.1. Metaprogramming in Financial Modeling

In the context of financial modeling, macros can be used to simplify repetitive or complex code patterns, enforce certain conventions or constraints, or generate code based on data or configuration.

Here are a few potential use cases of macros in financial modeling. Again, these are more advanced use-cases but knowing that these paths exist may benefit your work in the future.

1. Defining custom DSLs (Domain-Specific Languages): Macros can be used to create expressive and concise DSLs tailored to financial modeling. For example, a macro could allow defining financial contracts using a syntax closer to the domain language, which then gets expanded into the underlying implementation code.
2. Automating boilerplate code: Macros can help reduce code duplication by generating common patterns or boilerplate code. This can include generating accessor functions³, constructors, or serialization logic based on type definitions.
3. Enforcing conventions and constraints: Macros can be used to enforce coding conventions, such as naming rules or type checks, by automatically transforming code that doesn't adhere to the conventions. They can also be used to add runtime assertions or checks based on certain conditions.

³Accessor functions are useful when working with nested data structures. For example, if you have a `struct` within a `struct` and want to conveniently access an inner `struct`'s field.

8.5. Macros & Homoiconicity

4. Optimizing performance: Macros can be used to perform code optimizations at compile-time. For example, a macro could unroll loops, inline functions, or specialize generic code based on specific types or parameters, resulting in more efficient runtime code.
5. Generating code from data: Macros can be used to generate code based on external data or configuration files. For example, a macro could read a specification file and generate the corresponding financial contract types and functions.

Part IV.

Foundations: Building Performant Models

"Premature optimization is the root of all evil (or at least most of it) in programming." - Donald Knuth

After establishing foundational programming concepts, we turn our attention to performance - a critical consideration for real-world financial models. While modern computers are remarkably powerful, thoughtlessly constructed models can still grind to a halt when faced with large portfolios or complex analyses. This section explores how to harness computational resources effectively, starting from the hardware fundamentals that both constrain and enable our work.

Understanding performance requires looking beneath the abstractions we've built. Just as a financial modeler benefits from understanding the mechanics of markets and instruments rather than treating them as black boxes, knowledge of computational infrastructure allows us to make informed decisions about model architecture and implementation. We'll examine how hardware characteristics influence algorithm design, memory usage patterns, and execution speed.

We'll introduce when optimization *does* matter, and equally important when it *doesn't*. The goal isn't to optimize prematurely or pursue performance at all costs. Rather, we aim to build models that scale gracefully as demands grow, whether through larger datasets, more sophisticated analyses, or tighter time constraints. We'll progress from single-threaded optimization techniques to parallel processing approaches, always with an eye toward practical application in financial contexts.

By the end of this section, you'll have the knowledge needed to diagnose performance bottlenecks and implement appropriate solutions, ensuring your models remain responsive and reliable as they evolve. Let's begin by examining the hardware foundation upon which all our computational work rests.

9. Hardware and Its Implications

a CPU is literally a rock that we tricked into thinking.

not to oversimplify: first you have to flatten the rock and put lightning inside it.

- Twitter user daisyowl, 2017

9.1. In this section

In this chapter, we'll explore why a basic understanding of computing hardware is essential for optimizing financial models and working efficiently with data. Understanding how data is stored and processed can help you make better design decisions, improve performance, and avoid common pitfalls. We'll cover topics like memory architecture, data storage, and the impact of hardware on computational speed.

9.2. Introduction

The quote that opens the chapter is a silly way of describing that most modern computers are made with silicone, a common mineral found in rocks. However, we will not concern ourselves with the raw materials of computers and instead will focus on the key architectural aspect.

A computer handles data at rest (in memory) or data being acted upon (processed). This chapter will try to explain both of those processes in a way that reveals key reasons why different approaches to programming can yield different results in terms of processing speed, memory usage, and compiled outputs.

9.3. Memory and Moving Data Around

The core of modeling on computers is to perform computations on data, but unfortunately the speed at which data can be *accessed* has grown much slower than the rate the actual computations can be performed. Further, the size of the available persistent data

9. Hardware and Its Implications

storage (HDDs/SSDs) has ballooned, exacerbating the problem: the overall throughput of memory is the typical workflow constraint. To try to address the bottleneck (memory throughput), solutions have been developed to create a pipeline to efficiently shuttle data to and from the processor and the persistent storage. This memory and processing architecture applies at both the single computer level as well as extending to workflows between different data stores and computers.

We will focus primarily on the architecture of a single computer, as even laptop computers today contain enough power for most modeling tasks, *if the computer is used effectively*. Further, learning how to optimize a program for a single computer/processes or is almost always a precursor step to effective parallelization as well.

9.3.1. Memory Types and Location

Memory has an inverse relationship between size and proximity to the central processor unit (CPU). The closer the data is to the processor units, the smaller the storage and the less likely the data will persist at that location for very long.

Kind	Rough Size	Lifecycle
Solid State Disk (SSD) or Hard Disk Drive (HDD)	TBs	Persistent/Permanent
Random Access Memory (RAM)	Dozens to Hundreds of GBs	Seconds to Hours (while computer is powered on)
CPU Cache - L3	8 MB to 128 MB	Microseconds to Milliseconds
CPU Cache - L2	2 MB to 16 MB	Nanoseconds to Microseconds
CPU Cache - L1	~16 KB	Nanoseconds

After requesting data from a persistent location like a Solid State Drive (SSD), the memory is read into Random Access Memory (RAM). The advantage of RAM over a persistent location is speed - typically that memory can be accessed and modified many times faster than the persistent data location. The tradeoff is that RAM is not persistent: when the computer is powered down, the RAM loses the information stored within.

When data is needed by the CPU, data is read from RAM into a small hierarchy of caches before being accessed by the CPU. The **CPU Caches** are small (physically and in capacity), but very fast. The caches are also physically colocated with the CPU for efficiency. Data is organized and funneled through the caches as an intermediary between the CPU and RAM and is fed from Level 3 (L3) cache in steps down to L1 cache as the data gets closer to the processor.

i Note

For reference, memory units are:

- 1 bit is a single binary digit (0 or 1)
- 8 bits = 1 byte
- 8 bytes = 1 word
- 1024 bytes = 1 Kilobyte (KB)
- 1024 KB = 1 Megabyte (MB)
- 1024 MB = 1 Gigabyte (GB)
- 1024 GB = 1 Terabyte (TB)
- 1024 TB = 1 Petabyte (PB)

Sometimes, you might see Kb, which is *Kilobits*, or 1024 bits. Therefore 1 KB is 8 times larger than a Kb.

The increments are 1024 and not the usual 1000, because 1024 is 2^{10} . The even binary multiple of 1024 is more convenient than 1000 when working with bits.

9.3.2. Stack vs Heap

Sitting within the RAM region of memory are two sections called Stack and Heap. These are places where variables created from our program's code can be stored. In both cases, the program will request memory space but they have some differences to be aware of.

The **stack** stores small, fixed-size (known bit length), data and program components. The stack is a last-in-first-out queue of data that is able to be written to and read from very quickly. The **heap** is a region which can be dynamically sized and has random read/write (you need not access the data in a particular order). The heap is much slower but more flexible.

9.3.2.1. Garbage Collector

The **garbage collector** is a program that gets run to free up previously requested/allocated memory. It accomplishes this by keeping track of references to data in memory by section of your code. If a section of code is no longer reachable (e.g. inside a function that will never get called again, or a loop that ran earlier in the program but is now complete), then, periodically, the garbage collector will pause execution of the primary program in order to "sweep" the memory. This step marks the space as able to be reused by your program or the operating system.

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9.4. Processor

The processor reads lines (groups of bits) from the cache into registers and then executes instructions. An example would be to take the bytes from register 10 and add the bytes from register 11 to them. This is really all a processor does at the lowest level: combining bits of data using logical circuits.

Logical circuits (**transistors**) are an arrangement of wires that output a new electrical signal that varies depending on the input. From a collection of smaller building block gates (e.g. AND, OR, NOR, XOR) more complex operations can be built up¹, into operations like addition, multiplication, division, etc. Electric impulses move the state of the program forward once time per CPU cycle (controlled by a master “clock” ticking billions of times per second). CPU cycle speed is what’s quoted for chip performance, e.g. when a CPU is advertised as 3.0 GHz (or 3 billion cycles per second).

The programmer (or compiler, if we are working in a higher level language like Julia) tells the CPU which instruction to run. The set of instructions that are valid for a given processor are called the Instruction Set Architecture, or ISA. In computer Assembly language (roughly one-level above directly manipulating the bits), the instructions are given names like ADD, SUB, MUL, DIV, and MOV. These instructions mirror the raw instruction that is part of the ISA.

All instructions are not all created equal, however. Some instructions take many CPU cycles to complete. For example, floating point DIV (division) takes 10-20 CPU cycles while ADD only takes a single CPU cycle.

Some architecture examples that may be familiar:

- Intel x86-64 (a.k.a. AMD64) are common computer processors that use registers that are 64 bits wide (the prior generation was 32 bits wide) and use the **x86** instruction set developed by Intel.
- ARM chips, including the Apple M-Series processors are characterized by the use of the **ARM** instruction set and recent processors of this kind are also 64 bit.

The ARM architecture is known as a **reduced instruction set chip** (RISC), which means that it has fewer available instructions compared to, e.g. the x86 architecture. The benefit of the reduced instruction set is that it is generally much more power efficient, but comes at the cost of sacrificing specialized instructions such as string manipulation, or in lower-end chips, even the division operation (which have to be implemented via software routines instead of CPU operations). However, for specialized workloads, the availability of a key instruction can make a program run on the CPU 10-100x faster at

¹In fact, only two logical gates are needed to reproduce all boolean logical gates: NAND (Not AND) and NOR gates can be composed to create AND, OR, NOR, etc. gates.

times. An example of this is that at the time of writing, AVX512 processors are becoming available (see Chapter 11 for a discussion of vectorization) which can benefit some workloads greatly.

 Tip

Trying to optimize your program via selecting specialized chips should be one of the *last* ways that you seek to optimize runtime, as generally a similar order of magnitude speedup can be achieved through more efficient algorithm design or general parallelization techniques. Developing programs in this way makes the performance *portable*, able to be used on other systems and not just special architectures.

When writing in Julia, you need not be concerned with the low-level instructions as the compiler will optimize the execution for you. However, should it be useful, it is easy to inspect the compiled code. For example, if we create a function to add three numbers, we can see that the ADD instruction is called twice: first adding the first and second arguments, and then adding the third argument to that intermediate sum.

```
myadd(x, y, z) = x + y + z
@code_native myadd(1, 2, 3)

.section    __TEXT,__text,regular,pure_instructions
.build_version macos, 15, 0
.globl _julia_myadd_6915          ; -- Begin function julia_myadd_6915
.p2align   2
_julia_myadd_6915:             ; @julia_myadd_6915
; Function Signature: myadd(Int64, Int64, Int64)
; | @ /Users/alecloudenback/prog/julia-fin-book/hardware.qmd:101 within `myadd'
; %bb.0:                      ; %top
; | @ /Users/alecloudenback/prog/julia-fin-book/hardware.qmd within `myadd'
; | DEBUG_VALUE: myadd:x <- $x0
; | DEBUG_VALUE: myadd:x <- $x0
; | DEBUG_VALUE: myadd:y <- $x1
; | DEBUG_VALUE: myadd:y <- $x1
; | DEBUG_VALUE: myadd:z <- $x2
; | DEBUG_VALUE: myadd:z <- $x2
; | @ /Users/alecloudenback/prog/julia-fin-book/hardware.qmd:101 within `myadd'
; | | @ operators.jl:596 within `+' @ int.jl:87
; | | add x8, x1, x0
; | | add x0, x8, x2
; | | ret
; | LL
```

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```
; -- End function
.subsections_via_symbols
```

Compilers are complex, hyper-optimized programs which turn your source code into the raw bits executed by the computer. Key steps in the process of converting Julia code you write all the way to binary machine instructions include the items in Table 9.2. Note the Julia `@code_...` macros allow the programmer to inspect the intermediate representations.

Table 9.2.: Part the key to Julia's speed is to be able to compile down to a different, specialized version of the machine code depending on the types given to a function. As described in the table above, the instructions for adding floating point together or integer numbers together are different. Julia code can reflect that distinction by compiling a different method for each combination of input types.

Step	Description	Example
Julia Source Code	The level written by the programmer in a high level language.	<code>myadd(x,y,z) = x + y + z</code>
Lowered Abstract Syntax Tree (AST)	An intermediate representation of the code after the first stage of compilation, where the high-level syntax is simplified into a more structured form that's easier for the compiler to work with.	<code>julia> @code_lowered myadd(1,2,3) CodeInfo(1 - %1 = x + y + z └ return %1)</code>
LLVM	Low-Level Virtual Machine language, which is a massively popular compiler used by languages like Julia and Rust. The core logic are the three lines with add, add, and ret. Note that the the add instruction is add i64 which means an addition operation of 64 bit integers.	<code>julia> @code_llvm myadd(1,2,3) ; @ REPL[7]:1 within `myadd` define i64 @julia_myadd_2022(i64 signe top: ; ↗ @ operators.jl:587 within `+` @ in %3 = add i64 %1, %0 %4 = add i64 %3, %2 ret i64 %4 ; }</code>

Step	Description	Example
Native	The final machine code output, specific to the target CPU architecture. This is at the same level as Assembly language. The core logic are the three lines beginning with add, add, and ret. If we used floating point addition instead, the CPU instruction would be fadd instead of add.	julia> @code_native myadd(1,2,3) .section __TEXT,__text,regular,pure_instructions .build_version macos, 14, 0 .globl _julia_myadd_1851 .p2align 2 _julia_myadd_1851: ; @ REPL[7]:1 within `myadd` ; %bb.0: ; @ operators.jl:587 within `+` @ in add x8, x1, x0 add x0, x8, x2 ret ; LL

9.4.0.1. Increasing Complexity in Search of Performance

Transistors are the building-block that creates the CPU and enables the physical process which governs the computations. For a very long time, the major source of improved computer performance was simply to make smaller transistors, allowing more of them to be packed together to create computer chips. This worked for many years and the propensity for the transistor count to double about every two years. In this way, software performance improvements came as side effect of the phenomenal scaling in hardware capability. However, raw single core performance and clock frequency (CPU cycle speed) dramatically flattened out starting a bit before the year 2010. This was due to the fact that transistor density has been starting to be limited by:

1. Pure physical constraints (transistors can be measured in width of atoms) where we have limited ability to manufacture something so small.
2. Thermodynamics, where heat can't be removed from the CPU core fast enough to avoid damaging the core and therefore operations per second are capped.

To obtain increasing performance, two main strategies have been employed in lieu of throwing more transistors into a single core:

1. Utilize multiple, separate cores and operate in an increasingly parallel way.
2. Use clever tricks to predict, schedule, and optimize the computations to make better use of the memory pipeline and otherwise idle CPU cycles.

We will cover techniques to utilize concurrent/parallel processing in Chapter 11. As for the second technique, it is capable of very impressive accelerations (on the order 2x to 100x faster than a naive implementation. However, it has sometimes caused issues. There have been some famous security vulnerabilities such as Spectre and Meltdown,

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which exploited speculative execution – a technique used to optimize CPU performance which will execute code *before being explicitly asked to* because the scheduler *anticipates* the next steps (with very good, but imperfect accuracy).

9.5. Logistics Warehouse Analogy

The problem is analogous to a logistics warehouse (persistent data) which needs to package up orders (processor instructions). There's a conveyor belt of items being constantly routed to the packaging station. In order to keep the packing station working at full capacity, the intermediate systems (RAM & CPU caches) are funneling items they *think* will be needed to the packager (data that's *expected* to be used in the processor). Most of the time, the necessary item (data) is optimally brought to the packaging station (process), or a nearby holding spot (CPU cache).

This system has grown very efficient, but sometimes the predictions miss or a never-before-ordered item needs to be picked from the far side of the warehouse and this causes significant delays to the system. Sometimes a package will start to be assembled before the packager has even gotten to that order (branch prediction) which can make the system faster most of the time, but if the predicted package isn't actually what the customer ordered, then the work is lost and has to be redone (branch mispredict).

There are a lot more optimizations along the way:

- Since the items are already mostly arranged so that related items are next to each other, the conveyor belt will bring nearby items at the same time it brings the requested item (memory blocks).
- If an item usually ordered after another one is, the conveyor system will start to bring that second item as soon as the first one is ordered (prefetching).
- Different types of packaging stations might be used for specialized items (e.g. vector processing or cryptography instructions in the CPU).

9.6. Speed of Computer Actions

In a financial model, even small delays (such as main memory references vs. L1 cache access) can accumulate quickly in high-frequency trading or risk calculation routines. Understanding these timings can guide decisions on structuring data access patterns and deciding what data to cache or load in memory for optimal performance.

Representative time is given in Table 9.3 for a variety of common actions performed on a computer. It's clear that having memory access from a local source is better for computer performance!

9.6. Speed of Computer Actions

Table 9.3.: How long different actions take on a computer. As an interesting yardstick, the distance a beam of light travels is also given to provide a sense of scale (this comparison originally comes from Admiral Grace Hopper). Source for the timings comes from: <https://cs.brown.edu/courses/csci0300/2022/assign/labs/lab4.html>

Operation	Time (ns)	Distance Light Traveled
Single CPU Cycle (e.g. one ADD or OR operation on a register)	0.3	9 centimeters
L1 cache reference	1	30 centimeters
Branch mispredict	5	150 centimeters
L2 cache reference	5	150 centimeters
Main memory reference	100	30 meters
Read 1MB sequentially from RAM	250,000	75 km (~2 marathons)
Round trip within a datacenter	500,000	150km (the thickness of Earth's atmosphere)
Read 1MB sequentially from SSD	1,000,000	300km (distance Washington D.C. to New York City)
Hard disk seek	10,000,000	3,000km (width of continental United States)
Send packet CA->Netherlands->CA	150,000,000	45,000km (circumference of earth)

As introduced in Section 9.4, calculations on the CPU vary widely. Take, for example, Table 9.4 shows runtime varying quite a bit for common mathematical operations. Despite mathematically all essentially being elementary operations, the compute intensity varies widely:

Table 9.4.: Different operations utilize different levels of required compute. On different hardware, the result may be different depending on the fundamental computational complexity of an operation and what hardware instructions are built into the CPU.

Mathematical Operation	Julia Function	Runtime in nanoseconds (Float64 arguments)
$a + b$	+	1.7
$a - b$	-	1.7
a/b	/	1.7
$a \times b$	*	1.7
a^b	^	10.8
$\exp(a)$	exp	3.9

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Mathematical Operation	Julia Function	Runtime in nanoseconds (Float64 arguments)
$\log_{10}(a)$	<code>log</code>	5.4
$\log(a)$	<code>ln</code>	3.7
$abs(a)$	<code>abs</code>	1.7

💡 Tip

One way to speed up financial models is to utilize techniques to avoid using the $^$ operation, as it's one of the most expensive basic operations, but unfortunately is also one of the most common when dealing with rates and compounding.

Some strategies:

- Instead of using interest rates which require exponentiation, utilize continuously compounded rates and use the `exp` and `log` functions instead.
- Say you are running a monthly model with inputs that are annual rates. Before running through the 'hot loop', pre-compute the annual rates into a vector of monthly rates to avoid re-computing this transformation at the cell/seriatim level.

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Perfection is achieved, not when there is nothing more to add, but when there is nothing left to take away. - Antoine de Saint-Exupéry, Airman's Odyssey

10.1. In This Chapter

Understanding single-threaded performance, strategies for efficient sequential code, recognizing when parallelization is needed, practical comparisons of single-threaded and parallel approaches, optimizing code in Julia.

10.2. Introduction

With today's hardware, the highest throughput computations utilize GPUs for massive parallelization. However, writing parallel code, let alone performant parallel code, typically relies heavily on understanding patterns of non-parallel code. Secondly, many problems are not "massively parallelizable" and a sequential model architecture is required. For these reasons, it's critical to understand sequential patterns before moving onto parallel code.

For those coming from fundamentally slower languages (such as R or Python), the common advice to speed things up is often to try to parallelize code or use array-based techniques. With many high level languages the *only* way to achieve reasonable runtime is to utilize parallelism. In contrast, fast languages like Julia can output surprisingly quick single threaded programs.

Further, it may be that a simpler, easier to maintain sequential model is preferable to a more complex parallel version if maximum performance is not required. Like the quote that opened the chapter, you may prefer a simpler sequential version of a model to a more complex parallel one.

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💡 Tip

Developer time (your time) is often more expensive than runtime, so be prepared to accept a good-enough but sub-optimal code instead of spending a lot of time optimizing every last nanosecond out of it!

ℹ Note

This section of the book continues to focus on concepts that are more general than just Julia. We will elaborate on Julia-specific techniques in the section starting with Chapter 21.

10.3. Patterns of Performant Sequential Code

Sequential code performance depends on five key patterns:

1. Efficient memory usage
2. Optimized memory access
3. Appropriate data types
4. Type stability
5. Branch prediction optimization

Understanding these patterns helps you write faster code by leveraging CPU architecture and Julia's compiler optimizations. We'll look at each of these in turn.

10.3.1. Minimize Memory Allocations

Allocating memory onto the Heap takes a lot more time than (1) not using intermediate memory storage at all, or (2) utilizing the Stack. Each allocation requires time for memory management and requires the garbage collector, which can significantly impact performance, especially in tight loops or frequently called functions.

ℹ Note

Tight loops or hot loops are the performance critical section of the code that are performed many times during a computation. They are often the “inner-most” loop of a nested loop algorithm.

A general rule of thumb is that dynamically sizable or mutable objects (arrays, mutable structs) will be heap allocated while small fixed size objects can be stack allocated. For mutable objects, a common technique is to pre-allocate an array and then re-use that

10.3. Patterns of Performant Sequential Code

array for subsequent calculations. In the following example, note how we pre-allocate the output vector instead of creating vectors for each bond *and then* summing the vectors together at the end:

```
end_time = 10
cashflow_output = zeros(end_time)

par_bonds = map(1:1000) do i
    (tenor=rand((3, 5, 10)), rate=rand() / 10)
end

for asset in par_bonds
    for t in 1:end_time
        if t == asset.tenor
            cashflow_output[t] += 1 + asset.rate
        else
            cashflow_output[t] += asset.rate
        end
    end
end
cashflow_output

10-element Vector{Float64}:
48.69352009385709
48.69352009385709
385.69352009385733
48.69352009385709
395.69352009385705
48.69352009385709
48.69352009385709
48.69352009385709
48.69352009385709
364.69352009385733
```

Julia's `@allocated` macro will display the number of bytes allocated by an expression, helping you identify and eliminate unnecessary allocations.

```
random_sum() = sum([rand() for _ in 1:10])
@allocated random_sum()
```

10. Writing Performant Single-Threaded Code

10.3.2. Optimize Memory Access Patterns

Optimizing memory access patterns is essential for leveraging the CPU's cache hierarchy effectively. Modern CPUs have multiple levels of cache (L1, L2, L3), each with different sizes and access speeds. By structuring your code to access memory in a cache-friendly manner, you can significantly reduce memory latency and improve overall performance.

What is cache-friendly memory access? Essentially it boils down to spatial and temporal locality.

10.3.2.1. Spatial Locality

Spatial locality refers to accessing data that is physically near each other in memory (e.g contiguous blocks of data in an array).

For example, it's better to access data in a linear order rather than random order. For example, if we sum up the elements of an array in order it will be significantly faster than if we do it randomly:

```
using BenchmarkTools, Random

# Create a large array of structs to emphasize memory access patterns
struct DataPoint
    value::Float64
    # Add padding to make each element 64 bytes (a typical cache line size)
    padding::NTuple{7,Float64}
end

function create_large_array(n)
    [DataPoint(rand(), tuple(rand(7)...)) for _ in 1:n]
end

# Create a large array
const N = 1_000_000
large_array = create_large_array(N)

# Function for sequential access
function sequential_sum(arr)
    sum = 0.0
    for i in eachindex(arr)
        sum += arr[i].value
    end
    sum

```

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```
end

# Function for random access
function random_sum(arr, indices)
    sum = 0.0
    for i in indices
        sum += arr[i].value
    end
    sum
end

# Create shuffled indices
shuffled_indices = shuffle(1:N)

# Benchmark
println("Sequential access:")
@btime sequential_sum($large_array)

println("\nRandom access:")
@btime random_sum($large_array, $shuffled_indices)

Sequential access:
  519.375 μs (0 allocations: 0 bytes)

Random access:
  2.227 ms (0 allocations: 0 bytes)

500005.1233954948
```

When the data is accessed in a linear order, it means that the computer can load chunks of data into the cache and it can operate on that cached data for several cycles before new data needs to be loaded into the cache. In contrast, when accessing the data randomly, then the cache frequently needs to be populated with a different set of bits from a completely different part of our array.

10.3.2.1.1. Column vs Row Major Order

All multi-dimensional arrays in computer memory are actually stored linearly. When storing the multi-dimensional array, an architectural decision needs to be made at the language-level and Julia is column-major, similar to many performance-oriented languages and libraries (e.g. LAPACK, Fortran, Matlab). Values are stored going down the columns instead of across the rows.

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For example, this 2D array would be stored as [1,2,3,...] in memory, which is made clear via `vec` (which turns a multi-dimensional array into a 1D vector):

```
let
    array = [
        1 4 7
        2 5 8
        3 6 9
    ]

    vec(array)
end

9-element Vector{Int64}:
1
2
3
4
5
6
7
8
9
```

When working with arrays, prefer accessing elements in column-major order (the default in Julia) to maximize spatial locality. This allows the CPU to prefetch data more effectively.

You can see how summing up values across the first (column) dimension is much faster than summing across rows:

```
@btime sum(arr, dims=1) setup = arr = rand(1000, 1000)
@btime sum(arr, dims=2) setup = arr = rand(1000, 1000)
```

```
71.583 μs (3 allocations: 8.08 KiB)
110.333 μs (3 allocations: 8.08 KiB)
```

```
1000×1 Matrix{Float64}:
504.9533299073725
498.82158719941634
491.85960205447134
503.0059256198045
494.85894337283185
```

```
482.38849600667015  
490.23233996867424  
510.1877685256787  
499.95789085414805  
509.0348586427372  
:  
499.7475980584672  
496.8387594577944  
492.7209711484727  
501.12224814005566  
514.5516671165186  
504.25791345119904  
512.6643742405736  
511.4758576164987  
503.78347976746375
```

i Note

In contrast to the prior example, a row-major memory layout would have the associated data stored in memory as:

```
[1,4,6,2,5,8,3,6,9]
```

The choice between row and column major reflects the historical development of scientific computing and mathematical conventions. Column-major originated in Fortran and LAPACK, cornerstones of high-performance computing and linear algebra. This aligns with mathematical notation where vectors are typically represented as matrix columns.

Row major languages include:

- C/C++
- Python (NumPy arrays)
- C#
- Java
- JavaScript
- Rust

Column major languages:

- Julia
- MATLAB/Octave
- R
- Fortran
- LAPACK/BLAS libraries

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10.3.2.2. Temporal Locality

The scheduler and branch prediction will recognize data that is accessed together closely in time and prefetch relevant blocks of data. This is an example of keeping “hot” data more readily accessible to the CPU than “cold” data.

Sometimes this happens at the operating system level - if you load a dataset that exceeds the available RAM, some of the “active” memory will actually be kept in a space on the persistent disk (e.g. your SSD) to avoid the computer crashing from being out of memory. Segments of the data that have been accessed recently will be in RAM while sections of the data not recently accessed are likely to be in the portion stored on the persistent disk.

10.3.3. Use Efficient Data Types

The right data type can lead to more compact memory representations, better cache utilization, and more efficient CPU instructions. This is another case of where having a smaller memory footprint allows for higher utilization of the CPU since computers tend to be memory-constrained in speed.

On some CPUs, you may find performance if using the smallest data type that can accurately represent your data. For example, prefer Int32 over Int64 if your values will never exceed 32-bit integer range. For floating-point numbers, use Float32 instead of Float64 if the reduced precision is acceptable for your calculations. These smaller types not only save memory but also allow for more efficient vectorized operations (see [?@sec-parallelism](#)) on modern CPUs.

Warning

When choosing data types, consider that on modern 64-bit architectures, using smaller integer or floating-point types doesn’t always improve performance. In fact, many operations are optimized for 64-bit (or even larger SIMD) registers, and using smaller data types may introduce unnecessary overhead due to conversions or misalignment.

As a rough guideline, if your data naturally fits within 64 bits (e.g. Float64 or Int64), starting there is often the best balance of performance and simplicity. You can experiment with smaller types if you know your values never exceed certain ranges and memory footprint is critical. However, always benchmark to confirm any gains—simply using a smaller type does not guarantee improved throughput on modern CPUs.

For collections, choose appropriate container types based on your use case. Arrays are efficient for calculations that loop through all or most elements, while Dictionaries are better for sparse look-ups or outside of the “hot loop” portion of a computation.

Consider using small, statically sized collections when the data is suited for it. Small, fixed-size arrays, (such as `StaticArrays.jl` in Julia) can be allocated on the stack and lead to better performance in certain scenarios than dynamically sizeable arrays. The trade-off is that the static arrays require more up-front compile time and after a certain point (length in the 50-100 element range) it usually isn't worth trying to use them.

10.3.4. Avoid Type Instabilities

Type instabilities occur when the compiler cannot infer a single concrete type for a variable or function return value. These instabilities can significantly hinder Julia's ability to generate optimized machine code, leading to performance degradation. When the compiler is not able to infer the types at compile-time (**compile time dispatch**), then while the program is running a lookup needs to be performed to find the most appropriate functions for the given type (**runtime dispatch**). When the types are known at compile-time, Julia is able to create efficient machine code that will point directly to the desired function instead of needing to perform that lookup.

To avoid type instabilities, ensure that functions have inferrable, concrete types across all code paths. For example:

```
function unstable_example(array)
    x = []
    for y in array
        push!(x, y)
    end
    sum(x)
end

function stable_example(array)
    x = eltype(array)[]
    for y in array
        push!(x, y)
    end
    sum(x)
end

data = rand(1000)
@btime unstable_example(data)
@btime stable_example(data)
```

- ① Without a type given, `[]` will create a `Vector{Any}` which can contain elements of Any type which is flexible but requires runtime dispatch to determine correct behavior.

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- ② `eltype(array)` returns the type contained within the array, which for data is `Float64`. Thus `x` is created as a `Vector{Float64}` which is more efficient code.

```
10.583 μs (2007 allocations: 53.73 KiB)
801.471 ns (9 allocations: 22.52 KiB)
```

502.83996691305686

The `unstable_example` function illustrates a common anti-pattern wherein an `Any` typed array is created and then elements are added to it. Because any type can be added to an `Any` array (we happen to just add floats to it) then Julia's not sure what types to expect inside the container and therefore has to determine it at runtime.

Note that having heterogeneous types as above is not the same thing as **type instability**, which is when Julia cannot determine in advance what the data types will be. In the example above, the return type is not unstable: the compiler recognizes that the single parametric type `Union{Float64, Int64}` will be returned., even though two different types can be returned the When the types cannot be determined by the compiler, it leads to runtime dispatch.



Tip

See Section 24.5.1 and Section 24.5.2 for tools in Julia to troubleshoot type instabilities.

10.3.5. Optimize for Branch Prediction

Modern CPUs use branch prediction to speculatively execute instructions before knowing the outcome of conditional statements. This was described in the prior chapter in the logistics warehouse analogy as trying to predict where a package will be going before you've inspected the label. CPUs execute one branch of the instructions without seeing the data (either true data, or data in the form of CPU instructions) because the CPU has higher speed/capacity than the memory throughput allows. This is another example of a technique developed for performance in a memory-throughput constrained world.

Optimizing your code for branch prediction can significantly improve performance, especially in tight loops or frequently executed code paths.

To optimize for branch prediction:

1. Structure your code to make branching patterns more predictable. For instance, in `if-else` statements, put the more likely condition first. This allows the CPU to more accurately predict the branch outcome.

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2. Use loop unrolling to reduce the number of branches. This technique involves manually repeating loop body code to reduce the number of loop iterations and associated branch instructions. See Section 11.4 for more on what this means.
3. Consider using Julia's `@inbounds` macro to eliminate bounds checking in array operations when you're certain the accesses are safe. This reduces the number of conditional checks the CPU needs to perform.
4. For performance-critical sections with unpredictable branches, consider using branch-free algorithms or bitwise operations instead of conditional statements. This can help avoid the penalties associated with branch mispredictions.
5. In some cases, it may be beneficial to replace branches with arithmetic operations (e.g., using the ternary operator or multiplication by boolean values) to allow for better vectorization and reduce the impact of branch mispredictions. An example of this would be using a statement like `y += max(x, 0)` instead of `if x > 0; y += x; end`.

Here's an example demonstrating the impact of branch prediction:

```
function process_data(data, threshold)
    sum = 0.0
    for x in data
        if x > threshold
            sum += log(x)
        else
            sum += x
        end
    end
    sum
end

# Random data = unpredictable branches
rand_data = rand(1_000_000)

# Sorted data = predictable branches
sorted_data = sort(rand(1_000_000))

@btime process_data($rand_data, 0.5)
@btime process_data($sorted_data, 0.5);

4.992 ms (0 allocations: 0 bytes)
1.605 ms (0 allocations: 0 bytes)
```

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In this example, having sorted data means that the CPU will predict which branch of the `if` statement is likely to be utilized. By then speculatively executing the code that it thinks will be used, the overall program time is faster when processing `sorted_data`.

Remember that optimizing for branch prediction often involves trade-offs. The benefits can vary depending on the specific hardware and the nature of your data. If performance critical, profile your code to ensure that your optimizations are actually improving performance in your specific use case. Over-optimizing on one set of hardware (e.g. local computer) may not translate the same on another set of hardware (e.g. server deployment).

10.3.6. Further Reading

- What scientists must know about hardware to write fast code
- Optimizing Serial Code, ScIML Book

11. Parallelization

Quantity has a quality all its own. - Attributed to Vladimir Lenin

11.1. In this section

Fundamentals of parallel workloads, different mechanisms to distribute work: vectorization, multi-threading, GPU, and multi-device workflows. Different programming models: map-reduce, arrays, and tasks.

11.2. Amdahl's Law and the Limits of Parallel Computing

An important ground-truth in computing is that there is an upper limit to how fast a workload can be sped up through distributing the workload among multiple processor units. For example, if there is a modeling workload wherein 90% of the work is independent (say policy or asset level calculations) and the remaining 10% of the workload is an aggregate (say company or portfolio level), then the theoretical maximum speedup of the process is 10x faster (1 / 90% parallelizable load). This is captured in a law known as **Amdahl's Law** and it reflects the *theoretical* maximum speedup a workload could see. In practice, the speedup is worse than this due to overhead of moving data around, scheduling the tasks, and aggregating results. This is why in many cases a good effort in sequential workloads (see Chapter 10) is often a more fruitful effort than trying to parallelize some workloads.

That said, there are still many modeling use-cases for parallelization. Modern investment and insurance portfolios can easily contain 100's of thousands or millions of seriatim holdings. In many cases, these can be evaluated independently, though on the often times there is interaction with the total portfolio (contract dividends, non-guaranteed elements, profit sharing, etc.). Further, even if the holdings are not parallelizable across the holdings dimension, we are often interested in independent evaluations across economic scenarios which is amendable to parallelization.

$$S(n) = \frac{1}{(1-p) + \frac{p}{n}}$$

Where:

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- $S(n)$ is the theoretical speedup of the execution of the whole task
- n is the number of processors
- p is the proportion of the execution time that benefits from improved resources

We can visualize this for different combinations of p and n in Figure 11.1.

```
using CairoMakie

function amdahl_speedup(p, n)
    return 1 / ((1 - p) + p / n)
end

function main()
    fig = Figure(size=(800, 600))
    ax = Axis(fig[1, 1],
              title="Amdahl's Law",
              xlabel=L"Number of processors ($n$)",
              ylabel="Speedup",
              xscale=log2,
              xticks=2 .^ (0:16),
              xtickformat=x → "2^" .* string.(Int.(log.(2, x))),
              yticks=0:2:20
    )
    n = 2 .^ (0:16)
    parallel_portions = [0.5, 0.75, 0.9, 0.95]
    linestyles = [:solid, :dash, :dashdot, :solid]
    for (i, p) in enumerate(parallel_portions)
        speedup = [amdahl_speedup(p, ni) for ni in n]
        lines!(ax, n, speedup, label="$(Int(p*100))%", linestyle=linestyles[i])
    end
    xlims!(ax, 1, 2^16)
    ylims!(ax, 0, 20)
    axislegend(ax, L"Parallel portion ($p$)", position=:lt)
    fig
end

main()
```

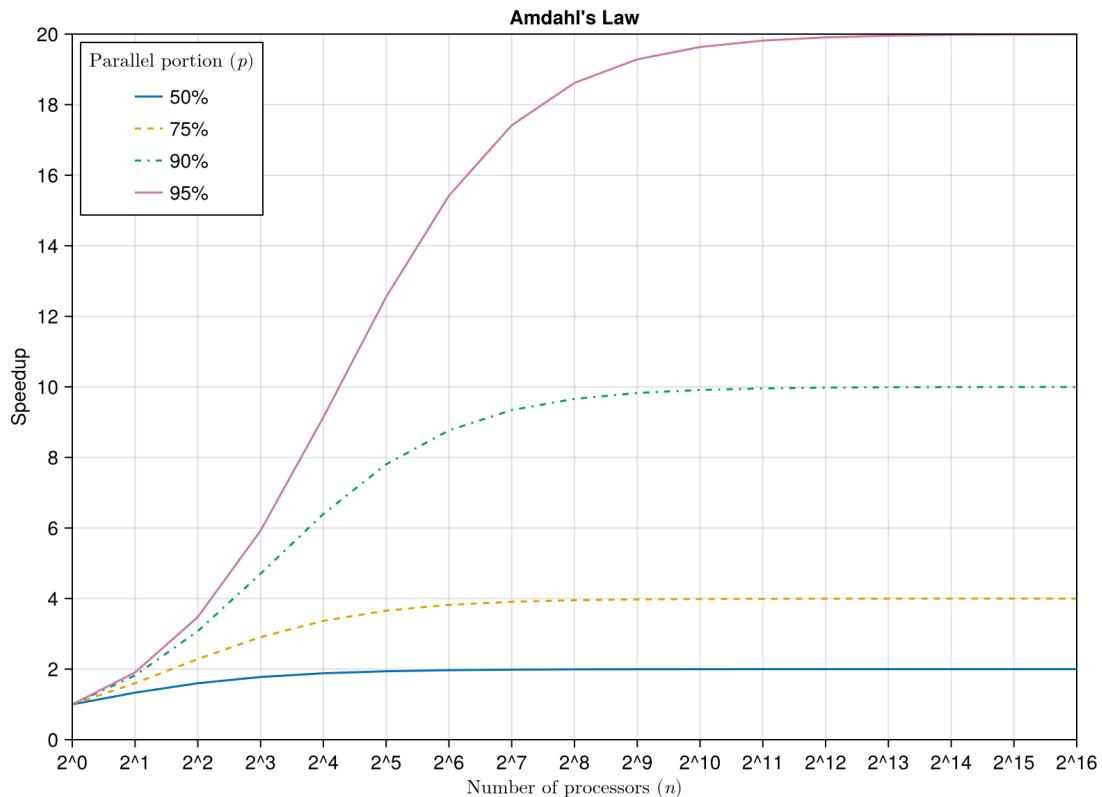


Figure 11.1.: Theoretical upper bound for speedup of a workload given the parallelizable portion p and number of processors n .

With this understanding, we will be able to set expectations and analyze the benefit of parallelization.

11.3. Types of Parallelism

Parallel processing comes in different flavors and is related to the details of hardware as discussed in Chapter 9. We will necessarily extend the discussion of hardware here, as parallelization is (mostly) inextricably tied to hardware details (we will revisit this in Section 11.8).

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Table 11.1.: Major types of computational parallelism highlighting their key characteristics, advantages, and potential drawbacks.

Type	Description	Strengths	Weaknesses
Vectorization (SIMD)	Performs same operation on multiple data points simultaneously	Efficient for data-parallel tasks, uses specialized CPU instructions	Limited to certain types of operations, data must be contiguous
Multi-Threading	Executes multiple threads concurrently on a single CPU	Good for task parallelism, utilizes multi-core processors effectively	Overhead from thread management, potential race conditions
GPU	Uses graphics processing units (GPUs) for parallel computations	Excellent for massively parallel tasks, high throughput	Specialized programming required, data transfer overhead
Multi-Device / Distributed	Spreads computation across multiple machines or devices	Scales to very large problems, can use heterogeneous hardware	Complex to implement and manage, network latency issues

11.4. Vectorization

Vectorization in the context of parallel processing refers to special circuits within the CPU wherein the CPU will load multiple data units (e.g. 4 or 8 floating point numbers) in a contiguous block and perform the same instruction on them at the same time. This is also known as **SIMD, or Single-Instruction Multiple Data**. Think of it like a `for` loop where instead of each iteration doing one operation, instead it does four or eight at a time.

The requirements for SIMD-able code are that:

- The intended section for SIMD is inside the inner-most loop.
- There are no branches (if-statements) inside the loop body.

::callout-tip Note that indexing an array is actually a branch in the code, as two cases could arise: the index is either inbounds or out-of-bounds. To avoid this, either use `for x in collection`, `for i in eachindex(collection)` or `for i in 1:n; @inbounds collection[i]` though the last of these is discouraged in favor of the prior, safer options.

:::

```

using BenchmarkTools

function prevent_simd(arr)
    sum = 0
    for x in arr
        if x > 0
            sum += x
        end
    end
    return sum
end

function allow_simd(arr)
    sum = 0
    for x in arr
        sum += max(x, 0)
    end
    return sum
end

let
    x = rand(10000)

    @btime prevent_simd($x)
    @btime allow_simd($x)
end

```

39.708 μs (0 allocations: 0 bytes)
4.815 μs (0 allocations: 0 bytes)

4956.853936109667

In testing the above code, the `allow_simd` version should be several times faster than the `prevent_simd` example. The reason is that `prevent_simd` has a branch (`if x > 0`) where the behavior of the code may change depending on the value in `arr`. Conversely, the behavior of `allow_simd` is always the same in each iteration, no matter the value of `x`. This allows the compiler to generate vectorized code automatically.

Tip

Note that Julia's the compiler is able to identify vectorizable code in many cases, though through some cases may benefit from a more manual hint to the compiler

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through macro annotations (see `?@simd` for details). See Section 24.8.4 for more.

Other types of parallelism that we will discuss in this chapter have some risk of errors or data corruption if not used correctly. SIMD isn't prone to issues like this because if the code is not SIMD-able then the compiler will not auto-vectorize the code block.

11.4.1. Hardware

Vectorization is hardware dependent. If the CPU does not support vectorization you will not see speedups from it. Many consumer and professional chips have AVX2 (Advanced Vector Extensions, with the 2 signifying second-generation 256 bit width, allowing four simultaneous 64-bit operations). The next generation is AVX512, having twice the SIMD capacity as AVX2. However, as of 2025 most consumer chips do not yet have that and commercial chips may not actually be faster than the AVX2 due to thermal restrictions (SIMD uses more power and generates more heat).

11.5. Multi-Threading

This subsection starts by introducing *tasks* — lightweight units of computation. Next, we'll see how tasks can communicate using *channels*, and then how multiple tasks (and channels) can be leveraged to achieve true parallelism through *multi-threading*. Think of it as layers building on one another: tasks define work units, channels allow them to share data, and multi-threading enables tasks to run simultaneously on multiple CPU cores. Exact details regarding tasks, channels, and multi-threading vary by language but the general ideas remain the same.

11.5.1. Tasks

To understand multi-threading examples, we first need to discuss **Tasks**, which are chunks of computation that get performed together, but after which the computer is free to switch to a new task. Technically, there are some instructions within a task that will let the computer pause and come back to that task later (such as `sleep`).

Tasks do not, by themselves, allow for multiple computations to be performed in parallel. For example, one task might be loading a data file from persistent storage into RAM. After that task is complete, the computer continues on with another task in the queue (rendering a web page, playing a song, etc.). In this way even a single processor computer could be “doing multiple things at once” (or “multi-tasking”) even though nothing is running in parallel. The scheduling of the tasks is handled automatically by the program’s compiler or the operating system.

Here's an example of a couple of tasks where we write to an array. Despite being called last, the second task should actually write to the array before the first task. This is because we asked the first task to `sleep` (pause, while allowing the computer to yield to other tasks in the queue)¹.

```
let
    shared_array = zeros(5)

    task1 = @task begin
        sleep(1)
        shared_array[1] = 1

        println("Task 1: ", shared_array)
    end

    task2 = @task begin
        shared_array[2] = 2
        println("Task 2: ", shared_array)
    end

    schedule(task1); schedule(task2)
    wait(task1)
    wait(task2)

    println("Main: ", shared_array)
end
```

```
Task 2: [0.0, 2.0, 0.0, 0.0, 0.0]
Task 1: [1.0, 2.0, 0.0, 0.0, 0.0]
Main: [1.0, 2.0, 0.0, 0.0, 0.0]
```

11.5.1.1. Channels

Channels are a way to communicate data in a managed way between tasks. You specify a type of data that the buffer (a chunk of assigned memory) will contain and how many elements it can hold. It then stores items (via `put!`) in a first-in-first-out (FIFO) queue, which can be popped off the queue (via `take!`) by other tasks.

¹Technically, it's possible that the second task doesn't write to the array first. This could happen if there's enough tasks (from our program or others on the computer) that saturate the CPU during the first task's `sleep` period such that the first task gets picked up again before the second one does.

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Here's an example of a system which generates trades in the financial markets at random time intervals, and a monitoring tasks takes the results and tabulates running statistics:

```
let

    # simulate random trades and the associated profits
    function trade_producer(channel,i)
        sleep(rand())
        profit = randn()
        put!(channel, profit)
        println("Producer: Trade Result #$i $(round(profit, digits=3))") ①
    end

    # intake trades via the communication channel
    function portfolio_monitor(channel,n)
        sum = 0.0
        for _ in 1:n
            profit = take!(channel)
            sum += profit
            println("Monitor: Received $(round(profit, digits=3)), Cumulative profit: $(round(sum, digits=3))") ②
        end
    end

    channel = Channel{Float64}(32) ④

    # Start producer and consumer tasks
    @sync begin
        for i in 1:5; @async trade_producer(channel,i); end ⑤
        @async portfolio_monitor(channel,5)
    end

    # Close the channel and wait for tasks to finish
    close(channel)
end ⑥
```

- ① Random sleep between 0 and 1 seconds to simulate real trading activity and latency.
- ② Generate a random number from standard normal distribution to simulate profit or loss from a trade.
- ③ In this teaching example, we've limited the system to produce just five "trades". In practice, this could be kept running indefinitely via, e.g., `while true`.
- ④ Create a channel with a buffer size of 32 floats (in this limited example, we could have gotten away with just 5 since that's how many the demonstration produces). In

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practice, you want this to be long enough that the consumer of the channel never gets so far behind that the channel fills up. The channel is created outside of the `@sync` block so that channel is in scope when we `close` it.

- ⑤ `@sync` waits (like `wait(task)`) for all of the scheduled tasks within the block to complete before proceeding with the rest of the program.
- ⑥ `@async` does the combination of creating a task via `@task` and `schedule-ing` in one, simpler call.

```
Producer: Trade Result #5 0.509
Monitor: Received 0.509, Cumulative profit: 0.509
Producer: Trade Result #4 -0.203
Monitor: Received -0.203, Cumulative profit: 0.307
Producer: Trade Result #2 1.324
Monitor: Received 1.324, Cumulative profit: 1.63
Producer: Trade Result #1 0.749
Monitor: Received 0.749, Cumulative profit: 2.38
Producer: Trade Result #3 1.258
Monitor: Received 1.258, Cumulative profit: 3.638
```

This is really useful for handling events that are “external” to our program. If we were just doing a modeling exercise using static data, then we could control the order of processing and not need to worry about monitoring a volatile source of data. Nonetheless, tasks can still be useful in some cases even if a model is not using “live” data: for example if one of the steps in a model is to load a very large dataset, it may be possible to perform some computations while chunked task requests are queued to load more data from the disk.

A garbage collector will usually clean up unused channels that are still open. However, it’s a good practice to explicitly `close` them to ensure proper resource management, clear signaling of completion, and to avoid potential blocking or termination issues in your programs.

🔥 Caution

If the task never finishes properly inside the `@sync`, then your program may get stuck in an infinite loop and hang. Such as if one of the tasks never has a termination condition such as an upper bound on a loop, or a clear way to break out of a `while true` loop. While not different than a normal loop, such issues become less obvious underneath the layer of task abstractions.

The key takeaway for tasks is that it’s a way to chunk work into bundles that can be run in a concurrent fashion, even if nothing is technically being processed in parallel. The multi-threading and parallel programming paradigms sections build off of tasks

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so an understanding of tasks is helpful. However, some of the higher level libraries hide the task-based building blocks from you as the user/developer and so an intricate understanding of tasks is not required to be successful in parallelizing your Julia code.

11.5.2. Multi-Threading Overview

When a program starts on your computer, a **process** is created which is where the operating system allocates some overhead items (keeping track of the the code and memory allocations and layout) and block of memory in RAM that can be utilized by that process. Different processes do not have access to each other's allocated memory.

Note

Readers may be familiar with starting Excel in different processes. When workbooks are opened within the same process (e.g. when creating a new workbook from Excel's File menu), the workbooks may seamlessly talk to each other (copy and paste from one to another). However, when Microsoft Excel is opened in different processes, then the workbooks in each respective process do not share memory and cannot create links or use full copy/paste functionality between them (this is what happens when you hold the control button and open Excel multiple times).

Within each process, a main thread is created. That thread is where the running of the code occurs. For the level of the discussion here, you can mainly think of a process as a container with shared memory for threads, which do the real work (as illustrated in Figure 11.2). Besides the main thread, other threads can be created within the process and access the same shared memory.

The advantage of threads is that within a single physical processor, there may be multiple cores. Those cores can access the shared process memory and run tasks from different threads simultaneously. This is a technique that takes advantage of modern processor architecture wherein several (sometimes as many as 32 or more) cores exist on the same chip.

Note

Technically, there are different flavors of threading. While not critical for the understanding and modeling-focused discussion here, here is a bit more detail on different thread types for completeness:

- **Multi-Tasking.** Recall that tasks are chunks of computation that get performed together, but after which the computer is free to switch to a new task. For example, one task might be loading a data file from persistent storage into RAM. After that task is complete, the computer continues on with an-

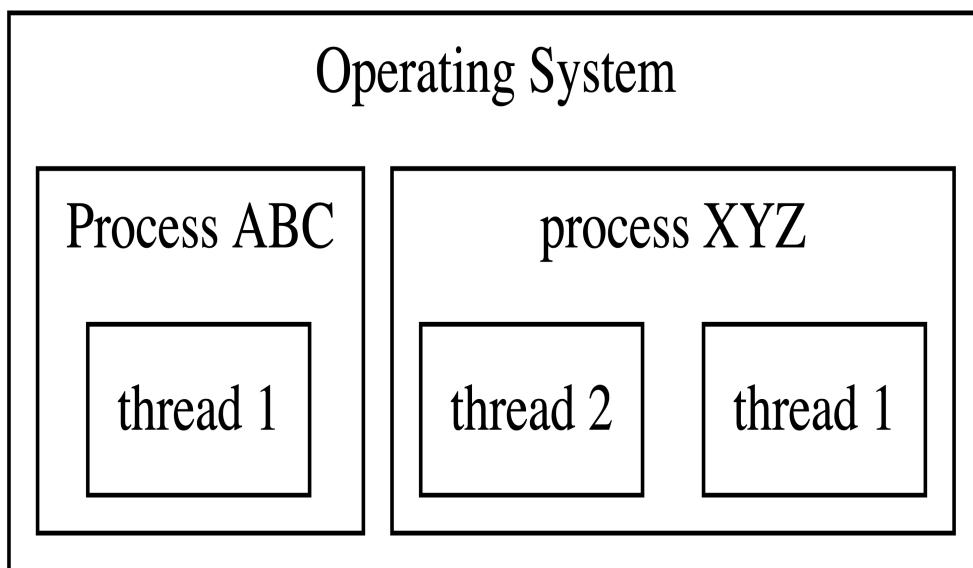


Figure 11.2.: When a program starts, the operating system creates a process for which multiple threads (a *main* thread plus optional additional threads) share memory.

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other task in the queue (rendering a web page, playing a song, etc.). In this way even with a single processor and core, a computer could be “doing multiple things at once” (or “multi-tasking”) even though nothing is running in parallel.

- **Operating System Threads** or just **Threads** are managed (as the name implies) at the operating system level. The benefit to this is that operating system level threads have more power: the operating system can pause or limit throughput on running programs if the operating system needs the resources for something it deems higher priority. It’s technically possible to use this power to force a higher priority for your own code, but Julia and many other languages do not offer creating of these types of threads in favor of the next type of threads. Operating system threads have a higher amount overhead (time and memory) involved in creating and destroying the threads.
- **Green threads, cooperative threads, fibers, or user-threads** are the type of threads that Julia provides. They are managed at the process (Julia) level and don’t have as much overhead in their creation as operating system threads. Also in Julia, a thread is implemented via Tasks

Parallelism in modern computing comes in many flavors, occurs at many different levels (hardware, OS, software, network), and has many different implementations of similar concepts. The terminology of threading in practice and online documentation is prone to confusing even seasoned developers. If you are having a discussion or asking a question, feel free to take the time to ask for clarification on the terminology being used at a given point in time.

11.5.2.1. Multi-Threading Pitfalls

Different threads being able to access the same memory is a double-edged sword. It is useful because we do not need to create multiple copies of the data in RAM or in the cache² and can improve the overall throughput of our usually memory-bandwidth-limited machines. The downside is that if we are mutating the shared data for which our program relies upon, then our program may produce unintended results if the modification occurs carelessly. There are a couple of related issues to be aware of:

11.5.2.1.1. Race Conditions

The first issue is known as a **race condition**, which occurs when a block of memory has been read from or written to in an unintended order. For example, if we have two threads

²There are some chips which do not have access to the same memory in a multi-threading context, and are known as non-uniform memory access (NUMA). These architectures work more like those in Section 11.7.

which are accumulating a sub-total, each process may read the running sub-total before the other thread has finished it's update.

In the following example, we use the `Threads.@threads` to tell Julia to automatically distribute the work across threads.

```
function sum_bad(n)
    subtotal = 0
    Threads.@threads for i in 1:n
        subtotal += i
    end
    subtotal
end

sum_bad(100_000)
```

5000050000

The result will be less than the expected sum (5000050000) due to a race condition. Here's what happens:

1. Multiple threads read the current value of `subtotal` simultaneously.
2. Each thread adds its own, local value to that reading.
3. Only one thread writes its result back to `subtotal` first.
4. A different thread then overwrites `subtotal` with its calculation based on the outdated starting point for `subtotal`.

This means some thread contributions are lost when they overwrite each other's results. The threads may not see each other's updates, leading to missing values in the final sum.

11.5.2.2. Avoiding Multi-threading Pitfalls

We will cover several ways to manage multi-threading race conditions, but it is the recommendation of the authors to primarily utilize higher level library code, which will be demonstrated after covering some of the more basic, manual techniques.

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11.5.2.2.1. Chunking up work into single-threaded work

First, let's level-set with a single-threaded result:

```
function sum_single(a)
    s = 0
    for i in a
        s += i
    end
    s
end
@btime sum_single(1:100_000)

1.500 ns (0 allocations: 0 bytes)
```

5000050000

Note that in the single-threaded case, Julia is able to identify this common pattern and use a shortcut, calculating the sum of the integers 1 through n as $\frac{n(n+1)}{2}$ through a compiler optimization and essentially avoid the loop entirely.

We can implement a correct threaded version by splitting the work into different threads, each of which is independent. Then, we can aggregate the results of each of the chunks.

```
function sum_chunker(a)

    chunks = Iterators.partition(1:a, a ÷ Threads.nthreads())           ①

    tasks = map(chunks) do chunk
        Threads.@spawn sum_single(chunk)
    end

    chunk_sums = fetch.(tasks)

    return sum_single(chunk_sums)

end

@btime sum_chunker(100_000)
```

- ① Create chunks of the integer range from 1 to a. `Iterators.partition(1:a, a ÷ Threads.nthreads())` splits the range `1:a` into contiguous subranges (chunks), each of size `a ÷ Threads.nthreads()`. For example, if `a = 100_000` and `nthreads() = 4`, you'll get four chunks of size `25_000`.

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- ② Create a set of tasks (futures) using `Threads.@spawn` that call `sum_single(chunk)` on each of the chunks. This initiates parallel computation.

3.500 µs (18 allocations: 944 bytes)

5000050000

11.5.2.2. Using Locks

Locks prevent memory from being accessed from more than one thread at a time.

```
function sum_with_lock(n)
    subtotal = 0                                ①

    lock = ReentrantLock()                      ②

    Threads.@threads for i in 1:n
        Base.@lock lock begin                  ③
            subtotal += i
        end
    end

    subtotal
end
@btime sum_with_lock(100_000)                ④
```

- ① Initialize a running total to zero.
- ② Create a reentrant lock to ensure only one thread updates `subtotal` at a time.
- ③ Parallelize the loop over 1 to n using `Threads.@threads`.
- ④ Acquire the lock before updating the shared variable `subtotal`. This ensures that only one thread updates `subtotal` at a time, preventing race conditions. The lock is automatically released at the end of the block.

3.082 ms (199469 allocations: 3.04 MiB)

5000050000

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11.5.2.2.3. Using Atomics

Atomics are certain primitive values with a reduced set of operations for which Julia and the compiler can automatically create thread-safe code. This is often significantly faster than the context-switching overhead needed with locking and unlocking memory for threaded tasks. Compared with locks, atomics are simpler to implement and easier to reason about. The downside is that atomics are limited to the available primitive atomics types and methods.

```
function sum_with_atomic(n)
    subtotal = Threads.Atomic{Int}(0)                                ①
    Threads.@threads for i in 1:n
        Threads.atomic_add!(subtotal, i)                            ②
    end
    subtotal[]
end

@btime sum_with_atomic(100_000)
```

- ① Initialize an atomic integer (`Threads.Atomic{Int}`) to store the subtotal. An atomic variable ensures that increments are performed atomically, preventing race conditions without needing explicit locks.
- ② Atomically add `i` to `subtotal`. The `Threads.atomic_add!` function ensures that the addition and update of `subtotal` happens as one atomic step, preventing multiple threads from interfering with each other's updates.

```
160.917 μs (8 allocations: 608 bytes)
```

```
5000050000
```

11.6. GPU and TPUs

11.6.1. Hardware

Graphics Processing Units (GPUs) and Tensor Processing Units (TPUs) are hardware accelerators for massively parallel computations. A TPU is very similar to a GPU but have special ability to handle data types and instructions that are more specialized for linear algebra operations; going forward we will simply refer to these types of accelerators as GPUs.

GPUs have similar components as the CPU as discussed in Chapter 9. They have RAM, caches for the cores, and cores that run the coded instructions on the data. The differences from a CPU are primarily:

- A GPU typically has thousands of cores while a CPU generally has single or double digit cores.
 - The cores typically operate at a *slower* clock speed than CPUs, relying on the sheer number of cores to perform computations faster.
- The GPU cores essentially have to be running the same set of instructions on all of the data, not unlike vectorization (Section 11.4).
 - GPU code is not suited for code with branching conditions (e.g. if statements) and so is more limited in the kinds of computations it can handle compared to the CPU.
- The RAM is typically much more constrained, typically less than a quarter of what primary RAM might be.
 - As a result, GPUs may need strategies to move chunks of data to and from the GPU memory for moderately large datasets. Further, it's actually fairly common to use a lower-precision datatype (e.g. Float16 or Float32) to improve overall program throughput at the cost of some precision.
- The caches are similar in concept to CPU, but unlike most CPU caches, there is relative locality to data (wherein core #1 will have much quicker access to a different subset of data than, say, core #1024).
- A GPU is usually a secondary device of sorts: it physically and in device architecture is separate from the CPU. The CPU remains in charge of overall computer execution.
 - The implication of this (as with any movement of memory) is that there is overhead to moving data to and from the GPU. Your calculations will need to be in the single milliseconds range of time in order to start to see benefit from utilizing a GPU.
 - To some extent, separable CPU, RAM, and GPU is changing with some of the latest computer hardware. For example, the M-series of Apple processors have the CPU, GPU, and RAM in a single tightly integrated package for efficiency and computational power.

11.6.1.1. Notable Vendors and Libraries

Like the difference between x86 and ARM architectures, GPU also have specific architectures which vary by the vendor. To make full use of the hardware, the vendors need to (1) provide device drivers which allow the CPU to talk to the GPU, and (2) provide the libraries (lower level application programming interfaces, or APIs) which allow developers to utilize different hardware features without needing to write machine code.

As of the mid 2020s, here are the most important GPU vendors and the associated programming library for utilizing their specific hardware:

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Table 11.2.: Important GPU and TPU vendors and the associated library/interface.

Vendor	Hardware	API Library/Package
NVIDIA	Geforce, GTX/RTX, various Data Center focused hardware	CUDA
AMD	Radeon, various Data Center focused hardware	ROCM
Intel	Core, Xeon, Arc processors	OneAPI
Apple	M Series processors	Metal
Google	Tensor processors	TensorFlow

11.6.2. Utilizing a GPU

With some of the key conceptual differences between CPUs and GPUs explained, let's explore how to incorporate these powerful hardware accelerators. We will use Julia libraries to illustrate GPU programming, though the concepts are generally applicable to other high-level languages that offer GPU interface libraries.

11.6.2.1. Julia GPU Libraries

There's essentially two types of GPU programming we will discuss here:

1. **Array-based programming**, where arrays are stored and operated on directly on the GPU memory. This approach abstracts away the low-level details, allowing you to work with familiar array operations that are automatically executed on the GPU.
2. **Custom kernels**, which are specialized functions that define exactly how each GPU thread should process data in parallel. A kernel explicitly specifies the computation that each GPU thread will perform on its portion of the data.

i Note

Kernels in this context are specialized functions that run directly on the GPU. Rather than relying solely on high-level array operations, kernels explicitly define the sequence of low-level, parallel instructions executed across many GPU threads. In other words, a kernel directly expresses the computation you want to perform on the data, enabling fine-grained control over GPU execution.

A third approach would be to implement GPU code in a low-level vendor toolkit (such as C++ and associate CUDA libraries), but this approach will not be illustrated here.

Julia has wonderful support for several of the primary vendors (at the time of writing, CUDA, Metal, OneAPI, and ROCm) via the JuliaGPU organization. Installation of the required dependencies is also very straightforward and the interfaces at the array and generated kernel levels are very similar. The differences are obvious at the lower level vendor-API wrappers (which is the lower-level technique that will not be covered here).

The benefit of the consistency of the higher level libraries we will use here is that examples written for one of the types of accelerators will be largely directly translatable to another. This is especially true for array programming, though less so for the kernel style as architecture-specific considerations often creep in³.

i Note

This book will be rendered on a Mac and therefore the examples will use Metal in order to run computational cells, however we'll show a CUDA translation for some of the examples in order to show the straight-forward nature of translating higher level GPU code in Julia is.

GPU API	GPU Array Type	Kernel Macro
CUDA	CuArray	@cuda
Metal	MtlArray	@metal
oneAPI	oneArray	@oneAPI
ROCM	ROCArray	@roc

11.6.2.2. Array Programming on the GPU

First described in Section 6.5, array programming eschews writing loops and instead favors initializing blocks of heap-allocated memory and filling it with data to be operated on at a single point in time. While this is often not the most efficient way to utilize CPUs, it's essentially the required style of code to utilize GPUs.

For the example below, we will calculate the present value of a series of cashflows across a number of different scenarios. An explanation of the code is given below the example.

```
using Metal

function calculate_present_values!(present_values,cashflows, discount_matrices) ①
```

³The KernelAbstractions.jl library actually allows you to write generic kernels which then get compiled into different code depending on the backend you are using.

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```
# Perform element-wise multiplication and sum along the time dimension
present_values .= sum(cashflows .* discount_matrices, dims=1)      ②
end

# Example usage using 100 time periods, 100k scenarios
num_scenarios = 10^5
pvs = zeros(Float32, 1, num_scenarios)
cashflows = rand(Float32, 100)                                         ③
discount_matrices = rand(Float32, 100, num_scenarios)                  ④

# copy the data to the GPU
pvs_GPU = MtlArray(pvs)
cashflows_GPU = MtlArray(cashflows)                                     ⑤
discount_matrices_GPU = MtlArray(discount_matrices)

@btime calculate_present_values!($pvs,$cashflows, $discount_matrices)
@btime calculate_present_values!($pvs_GPU,$cashflows_GPU, $discount_matrices_GPU) ⑥
```

- ① The function `calculate_present_values!` is written the same way as if we were just writing CPU code. Note that we are also passing a pre-allocated vector, `present_values` to store the result. This will allow us to isolate the performance of the computation, rather than including any overhead of allocating the array for the result.
- ② The code is broadcasted across the first dimension so that the single set of cashflows is discounted for each scenario's discount vector.
- ③ Metal only supports 32 bit floating point (some CUDA hardware will support 64 bit floating point)
- ④ Using 100 thousand scenarios for this example.
- ⑤ `MtlArray(array)` will copy the array values to the GPU.
- ⑥ Note that the data still lives on the GPU and is of the `MtlMatrix` (a type alias for a 2-D `MtlArray`).

```
2.561 ms (6 allocations: 38.53 MiB)
72.541 μs (515 allocations: 13.93 KiB)
```

```
1×100000 MtlMatrix{Float32, Metal.PrivateStorage}:
22.3762 22.771 23.9453 25.5413 ... 23.3689 22.318 25.1744 26.9168
```

The testing suggests approximately 200 times faster computation when performed on the GPU. Note however, that does not include the overhead of (1) moving the data to the GPU (in the initial `MtlArray(cashflows)` call), or (2) returning the data to the CPU (since the return type for the GPU version is `MtlArray`). We can measure this overhead by wrapping the data transfer inside another function and benchmarking it:

```

function GPU_overhead_test(present_values, cashflows, discount_matrices)
    pvs_GPU = MtlArray(present_values)
    cashflows_GPU = MtlArray(cashflows)
    discount_matrices_GPU = MtlArray(discount_matrices) (5)
    calculate_present_values!(pvs_GPU, cashflows_GPU, discount_matrices_GPU)

    Array(pvs_GPU) # convert to CPU array
end

@btime GPU_overhead_test($pvs,$cashflows,$discount_matrices)

```

6.404 ms (987 allocations: 418.45 KiB)

```

1×100000 Matrix{Float32}:
 22.3762  22.771  23.9453  25.5413 ... 23.3689  22.318  25.1744  26.9168

```

With the additional overhead, the computation on the GPU takes more total time than if the work were done just on the CPU. This is a very simple example, and the balance tips heavily in favor of the GPU when:

1. The computational demands are significantly higher (e.g. we were to do more calculations than just a simple multiply/divide/sum).
2. The data size grows bigger.

Note

The previous example can be translated to CUDA by simply exchanging `MtlArray` for `CuArray`.

Warning

This example again underscores that hardware parallelization is not an automatic “win” for performance. A lot of uninformed discussion around modeling performance is to simply try to get things to run on the GPU and it is often *not* the case that the models will run faster. Further, as the modeling logic gets more complex, it does require greater care to keep in mind GPU constraints (acceptable data types, memory limitations, avoiding scalar operations, data transfer between CPU and GPU, etc.). A best practice is to contemplate sequential performance and memory usage before leveraging GPU accelerators.

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11.6.2.3. Kernel Programming on the GPU

Another approach to GPU programming is often referred to as kernel programming, or being much more explicit about *how* a computation is performed. This is as opposed to the declarative approach in the array-oriented style (Section 11.6.2.2) wherein we specified *what* we wanted the computation to be.

The key ideas here are that we need to manually specify several aspects which came ‘free’ in the array-oriented style. The tradeoff is that we can be more fine-tuned about how the computation leverages our hardware, potentially increasing performance.

The GPU libraries in Julia abstract much of the low level programming typically necessary for this style of programming, but we still need to explicitly look at:

1. How the GPU will iterate across different cores/threads threads.
 2. How many threads to utilize, the optimal number depends on the shape of the computation (long vectors, multi-dimensional arrays), memory constraints, and hardware specifics.
- GPU threads: Individual units of execution within a kernel. Each thread runs the same kernel code but operates on a different portion of the data.
3. How to chunk (group) the data to distribute the data to the different GPU threads

Our strategy for the present values example will be to distribute the work such that different GPU threads are working on different scenarios. Within a scenario, the loop is a very familiar approach: initialize a subtotal to zero and then accumulate the calculated present values.

```
function calculate_present_values_kernel!(present_values,cashflows, discount_matrices)
    idx = thread_position_in_grid_1d()                                ①
    pv = 0.0f0
    for t in 1:size(cashflows, 1)
        pv += cashflows[t] * discount_matrices[t, idx]
    end
    present_values[idx] = pv                                         ④
    return nothing
end                                                               ⑤
```

- ① As the work is distributed across threads, `thread_position_in_grid_1d()` will give the index of the current thread so that we can index data appropriately for the work as we decide to split it up (we’ve split up the work by scenario in this example).
- ② Recall that we are working with `Float32` on the GPU here, so the zero value is set via the `f0` notation indicating a 32-bit floating point number.

- ③ The loop is across timesteps within each thread, while the thread index is tracked with `idx`.
- ④ The result is written to the pre-allocated array of present values, and we avoid race conditions because the different threads are working on different scenarios.
- ⑤ We don't explicitly have to `return nothing` here, but it makes it extra clear that the intention of the function is to mutate the `present_values` array given to it. This mutation intention is also signaled by the `!` convention in the function name.

```
calculate_present_values_kernel! (generic function with 1 method)
```

The kernel above was fairly similar to how we might write code for CPU-threaded approaches, but we now need to specify the technicals of launching this on the GPU. The `threads` defines how many independent calculations to run at a given time, and the maximum will be dependent on the hardware used. The `groups` argument defines the number of threads that share memory and synchronize results together (meaning that group will wait for all threads to finish before moving onto the next chunk of data). The push-pull here is that threads that can share data avoid needing to create duplicate copies of that data in memory. However, if there is variability in how long each calculation will take, then the waiting time for synchronizing results may slow the overall computation down.

Our task utilizes shared memory of the cashflows for each thread, so through some experimentation in advance, we find that a relatively large group size of ~512 is optimal.

We bring this all together through the use of the kernel macro `@metal`:

```
threads = 1024
groups = cld(num_scenarios, 512)

@btime @metal threads=$threads groups=$groups calculate_present_values_kernel!(
    $pvs_GPU,
    $cashflows_GPU,
    $discount_matrices_GPU
)
```

16.750 µs (149 allocations: 3.33 KiB)

```
Metal.HostKernel{typeof(calculate_present_values_kernel!)}, Tuple{MtlDeviceMatrix{Float32, 1}, M
```

This is approximately seven times faster than the array-oriented style above, meaning that the GPU kernel version's computation is over 1000 times faster than the CPU version. However, we saw previously that the cost of moving the data to the GPU memory and then back to the CPU memory was the biggest time sink of all - again we'd need to have more scale in the problem to make offloading to the GPU beneficial overall.

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i Note

The Metal GPUs are able to iterate threads across three different dimensions. In the prior example, we only used one dimension and thus used `thread_position_in_grid_1d()`. If we were distributing the threads across, say, three dimensions then we would use `thread_position_in_grid_2d()`.

How do you determine how many dimensions to use? A good approach is to mimic the data you are trying to parallelize. In the example of calculating a vector of present values across 100k scenarios, that was the primary ‘axis of parallelization’. If instead of a one-dimensional set of cashflows (e.g. a single asset with fixed cashflows), we had a two-dimensional set of cashflows (e.g. a portfolio of many assets), then we may find the best balance of code simplicity and performance to iterate across two dimensions of threads (but we are still limited by the same number of total available threads).

i Note

The above example would be translated to CUDA by changing just a few things:

- The thread indexing would be `idx = threadIdx().x` instead of `i = thread_position_in_grid_1d()`
- The GPU arrays should be created with `CuArray` instead of `MtlArray`.
- The kernel macro would be `@cuda threads=1024 calculate_present_values_kernel!(...)` instead of `@metal threads=threads groups=groups calculate_present_values_kernel(...)`. The memory sharing and synchronizing between threads is more manual than Metal.jl, but this is not strictly necessary for our example.

```
function calculate_present_values_kernel!(present_values,cashflows, discount_matrices)
    idx = threadIdx().x

    pv = 0.0f0
    for t in 1:size(cashflows, 1)
        pv += cashflows[t] * discount_matrices[t, idx]
    end

    present_values[idx] = pv
    return nothing
end

groups = cld(num_scenarios, 512)
```

```
@cuda threads=threads calculate_present_values_kernel!(
    pvs_GPU,
    cashflows_GPU,
    discount_matrices_GPU
)
```

11.7. Multi-Processing / Multi-Device

Multiple device, or **multi-device** computer refers to using separate groups of memory/processor combinations to accomplish tasks in parallel. This can be as simple as multiple distinct cores on within a single desktop computer, or many separate computers networked across the internet, or many processors within a high performance cluster or a computing-as-a-service provider like Amazon Web Services or JuliaHub.

Everything discussed previously related to hardware (Chapter 9, Section 11.5, Section 11.6) continues to apply. The additional complexity is attempting to synchronize the computation across multiple sets of the same (homogeneous) or different (heterogeneous) hardware.

As you might imagine, approaches to multi-device computing can vary widely. Julia's approach tries to strike the balance between capability and user-friendliness and uses a primary/worker model wherein one of the processors is the main coordinator while other processors are "workers". If only one processor is started, then the main processor is also a worker processor. This main/worker approach uses a "one-sided" approach to coordination. The main worker utilizes high level calls and the workers respond, with some of the communication and hand-off handled by Julia transparently from the user's perspective.

A useful mental model is the asynchronous task-based concepts discussed in Section 11.5.1, as the main worker will effectively queue work with the worker processors. Because there may be a delay associated with the computation or the communication between the processors, the worker runs asynchronously.

Description	Task API	Distributed Analogue
Create a new task	Task()	@spawnat
Run task asynchronously	@async	@spawnat
Retrieve task result	fetch	fetch
Wait for task completion	@sync	sync
Communication between tasks	Channel	RemoteChannel

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Adapting the trade producer and monitor example from above to run on multiple processors (see #sec-channels to review the base model and algorithm), we make a few key changes:

- using `Distributed` loads the `Distributed` standard Julia library, providing the interface for multi-processing across different hardware.
- `addprocs(n)` will add n worker processors (the main processor is already counted as one worker). When adding local machine processors, the processors are part of the local machine. This starts new Julia processes (you can see this in the task manager of the machine) which inherit the package environment (i.e. `Project.toml` and environment variables) from the main process; this does not occur automatically if not part of the same local machine.
 - To add processors from other machines, see the Distributed Computing section of the Julia docs.
- `myid()` is the identification number of the given processor that's been spun up.
- We use a `RemoteChannel` instead of a `Channel` to facilitate communication across processors.
- Instead of `@async`, we use `@spawnat n` to create a task for processor number n (or `:any` will automatically assign a processor).

See

```
using Distributed
let

    # Add worker processes if not already added
    if nworkers() == 1
        addprocs(4) # Add 4 worker processes
    end

    @everywhere function trade_producer(channel, i)
        sleep(rand())
        profit = randn()
        put!(channel, profit)
        println("Producer $(myid()): Trade Result #$i $(round(profit, digits=3))")
    end

    @everywhere function portfolio_monitor(channel, n)
        sum = 0.0
        for _ in 1:n
            profit = take!(channel)
            sum += profit
            println("Monitor $(myid()): Received $(round(profit, digits=3)), Cumulative profit $sum")
        end
    end
end
```

```

        end
    end

function run_distributed_simulation()
    channel = RemoteChannel(() -> Channel{Float64}(32))

    # Start producer and consumer tasks
    @sync begin
        for i in 1:5
            @spawnat :any trade_producer(channel, i)
        end
        @spawnat :any portfolio_monitor(channel, 5)
    end

    # Close the channel and wait for tasks to finish
    close(channel)
end

# Run the simulation
run_distributed_simulation()
end

```

```

From worker 5: Producer 5: Trade Result #4 -0.817
From worker 3: Producer 3: Trade Result #2 -0.164
From worker 3: Monitor 3: Received -0.164, Cumulative profit: -0.164
From worker 3: Monitor 3: Received -0.817, Cumulative profit: -0.981
From worker 3: Monitor 3: Received 0.757, Cumulative profit: -0.224
From worker 2: Producer 2: Trade Result #1 0.757
From worker 2: Producer 2: Trade Result #5 1.845
From worker 3: Monitor 3: Received 1.845, Cumulative profit: 1.621
From worker 3: Monitor 3: Received -0.414, Cumulative profit: 1.207
From worker 4: Producer 4: Trade Result #3 -0.414

```

Given the similarity to the single-process task-based version above, what's the motivation for this bothering with a distributed approach? A few differences:

- In this simplified example, we are simply starting additional Julia processes on the same machine. Like a threaded approach, the work will be split across the same multi-core processor. In this context, the main difference is that the processes do not share memory.
 - Communicating across processes generally has a little bit more overhead than communicating across threads.

11. Parallelization

- If distributing across machines, avoiding memory sharing is advantageous if using the different machines that have their own memory stores, which need not compete with the main process (such as distributed chunks of large datasets). This essentially helps with memory constrained problems since you are no longer limited by the memory size or throughput of a single machine.
- The worker processors don't need to be the same architecture as the main processor, allowing usage of different machines or cloud computing that is communicating with a local main process.

11.8. Parallel Programming Models

The previous sections have explained the different parallel programming models and how to directly utilize them to harness additional computing power. Each approach (multi-threading, GPU, distributed processing, etc.) has unique considerations and trade-offs. These approaches in Julia are generally much more accessible to beginning and intermediate users than other languages, but admittedly still requires a decent amount of thought and care.

It is possible, if you are willing to give up some fine-grained control, to utilize some higher level approaches which look to abstract away some of the particularities of the implementation.

11.8.1. Map-Reduce

Map-Reduce (Section 6.4.4) operations are inherently parallelizable and various libraries provide parallelized versions of the base `mapreduce`. This is the workhorse function of many ‘big data’ workloads and many statistical operations are versions of `mapreduce`.

11.8.1.1. Multi-Threading

11.8.1.1.1. OhMyThreads

`ThreadsX.jl` provides the threaded versions of essential functions such as `tmap`, `tmapreduce`, `tcollect`, and `tforeach` (see Table 6.1). In most cases, the chunking and data sharing is handled automatically for you.

```
import OhMyThreads
@btime OhMyThreads.tmapreduce(x → x, +, 1:100_000)

4.416 μs (17 allocations: 896 bytes)
```

```
5000050000
```

11.8.1.1.2. ThreadsX

ThreadsX.jl is built off of the wonderful Transducers.jl package, though the latter is a bit more advanced (more abstract, but as a result more composable and powerful). ThreadsX provides threaded versions of many popular base functions. It offers a wider set of ready-made threaded functions, but has a much more complex codebase. For the vast majority of threading needs, OhMyThreads.jl should be sufficient and performant. See the documentation for all of the implemented functions, but for our illustrative example:

```
import ThreadsX
@btime ThreadsX.mapreduce(x -> x, +, 1:100_000)
```

```
75.500 µs (50 allocations: 3.73 KiB)
```

```
5000050000
```

11.8.1.2. Multi-Processing

`reduce(op, pmap(f, collection))` will use a distributed map and reduce the resulting map on the main thread. This pattern works well if each application of `f` to elements of `collection` is costly.

`@distributed (op) for x in collection; f(x); end` is a way to write the loop with the reduction `op` for which the `f` need not be costly.

The difference between the two approaches is that with `pmap`, `collection` is made available to all workers. In the `@distributed` approach, the collection is partitioned and only a subset is sent to the designated workers.

Here's an example of both of these, calculating a simple example of counting coin flips:

```
# this is a example of poor utilization of pmap, since the operation is
# fast and the overhead of moving the whole collection dominates
@btime reduce(+, pmap(x -> rand((0,1)), 1:10^3))
```

```
74.303 ms (68104 allocations: 3.28 MiB)
```

```
500
```

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```
function dist_demo()
    @distributed (+) for _ in 1:10^5
        rand((0,1))
    end
end

@btime dist_demo()
```

```
182.834 μs (313 allocations: 13.92 KiB)
```

```
50071
```

11.8.2. Array-Based

Array based approaches will often utilize the parallelism of SIMD on the CPU or many cores on the GPU. It's as simple as using generic library calls which will often be optimized at the compiler level. Examples:

```
let
    x = rand(Float32, 10^8)
    x_GPU = MtlArray(x)
    @btime sum($x)
    @btime sum($x_GPU)
end
```

```
4.638 ms (0 allocations: 0 bytes)
1.279 ms (794 allocations: 19.90 KiB)
```

```
5.0001724f7
```

`sum(x)` compiles to SIMD instructions on the CPU, while using the GPU array type in `sum(x_GPU)` is enough to let the compiler dispatch on the GPU type and emit efficient, parallelized code for the GPU.

Distributed array types allow for large datasets to effectively be partitioned across multiple processors, and have implementations in the `DistributedArrays.jl` and `Dagger.jl` libraries.

11.8.3. Loop-Based

Loops which don't have race conditions can easily become multi-threaded. Here, we have three versions of updating a collection to square the contained values:

```
let v = collect(1:10000)

for i in eachindex(v)
    v[i] = v[i]^2
end
v[1:3]
end
```

3-element Vector{Int64}:

```
1
4
9
```

Using multi-threading

```
let v = collect(1:10000)
Threads.@threads for i in eachindex(v)
    v[i] = v[i]^2
end
v[1:3]
end
```

3-element Vector{Int64}:

```
1
4
9
```

Using multi-processing:

```
let v = collect(1:10000)
Distributed.@distributed for i in eachindex(v)
    v[i] = v[i]^2
end
v[1:3]
end
```

11. Parallelization

```
3-element Vector{Int64}:
1
2
3
```

For more advanced usage, including handling shared memory see Section 11.7 and Section 11.5.

11.8.4. Task-Based

Task based approaches attempt to abstract the scheduling and distribution of work from the user. Instead of saying how the computation should be done, the user specifies the intended operations and allows the library to handle the workflow. The main library for this in Julia is Dagger.jl.

Effectively, the library establishes a network topology (a map of how different processor nodes can communicate) and models the work as a directed, acyclic graph (a DAG, which is like a map of how the work is related). The library is then able to assign the work in the appropriate order to the available computation devices. The benefit of this is most apparent with complex workflows or network topologies where it would be difficult to manually assign, communicate, and schedule the workflow.

Here's a very simple example which demonstrates Dagger waiting for the two tasks which work in parallel (we already added multiple processors to this environment in Section 11.7):

```
import Dagger

# This runs first:
a = Dagger.@spawn rand(100, 100)

# These run in parallel:
b = Dagger.@spawn sum(a)
c = Dagger.@spawn prod(a)

# Finally, this runs:
wait(Dagger.@spawn println("b: ", b, ", c: ", c))

From worker 3:    Unhandled Task ERROR: On worker 1:
From worker 3:    On worker 4:
From worker 3:    peer 3 has not connected to 4
From worker 3:    Stacktrace:
From worker 3:        [1] error
```

```

From worker 3:      @ ./error.jl:35
From worker 3:      [2] wait_for_conn
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [3] #27
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [4] exec_conn_func
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [5] exec_conn_func
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [6] #invokelatest#2
From worker 3:      @ ./essentials.jl:1055
From worker 3:      [7] invokelatest
From worker 3:      @ ./essentials.jl:1052
From worker 3:      [8] #110
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [9] run_work_thunk
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [10] #109
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      Stacktrace:
    From worker 3:      [1] #remotecall_fetch#159
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [2] remotecall_fetch
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [3] remotecall_fetch
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [4] #23
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [5] #invokelatest#2
From worker 3:      @ ./essentials.jl:1055
From worker 3:      [6] invokelatest
From worker 3:      @ ./essentials.jl:1052
From worker 3:      [7] #110
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [8] run_work_thunk
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      [9] #109
From worker 3:      @ ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
    From worker 3:      Stacktrace:
From worker 3:      [1] remotecall_fetch(::Function, ::Distributed.Worker, ::Int64, ::Vararg{Ir
From worker 3:      @ Distributed ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/sha
From worker 3:      [2] remotecall_fetch(::Function, ::Distributed.Worker, ::Int64, ::Vararg{Ir
From worker 3:      @ Distributed ~/julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/sha

```

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```
From worker 3:      [3] remotecall_fetch
From worker 3:      @ ~/.julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
From worker 3:      [4] (::Distributed.var"#22#25"{Distributed.Worker})()
From worker 3:      @ Distributed ~/.julia/juliaup/julia-1.11.4+0.aarch64.apple.darwin14/share/julia/std
From worker 4:      b: 4993.310552760367, c: 0.0
```

11.9. Choosing a Parallelization Strategy

There is no one-size-fits-all strategy to parallelization. Here are some general guides to thinking about what parallelization technique to try given the circumstances:

- **CPU-bound workloads with manageable memory demands:** If your entire dataset fits comfortably in RAM and your operations are primarily arithmetic or straightforward loops, start by optimizing your single-threaded performance and consider **vectorization (SIMD)** for inner loops and **multi-threading** for parallelizable tasks. This approach leverages your CPU cores efficiently without introducing significant complexity.
- **Large-scale linear algebra or highly data-parallel computations:** If your problem involves large matrix operations, linear algebra routines, or embarrassingly parallel computations that can be batched over many independent elements, a GPU or other specialized accelerators may be beneficial. GPUs excel at uniform computations over large datasets and can provide substantial speedups—assuming data transfer overhead and memory constraints are managed effectively. Note that standard linear algebra libraries are likely to already parallelize on the CPU without any explicit parallelization needed to be coded on your part.
- **Distributing work across multiple machines or heterogeneous resources:** If you need to scale beyond a single machine’s CPU and GPU capabilities—whether due to extremely large datasets, the need for concurrent access to geographically distributed resources, or leveraging specialized hardware—then consider **distributed computing**. Spreading tasks across multiple processes, servers, or clusters can scale performance horizontally. Just keep in mind the overhead of communication, potential data partitioning strategies, and the complexity of managing a distributed environment.

In practice, you may find that a combination of these approaches is ideal: start simple, measure performance, and iterate. By understanding your workload and hardware constraints, you can make informed decisions that balance complexity, cost, and the performance gains of parallel computing.

11.10. References

- https://book.sciml.ai/notes/06-The_Different_Flavors_of_Parallelism/
- <https://docs.julialang.org/en/v1/manual/parallel-computing/>
- <https://enccs.github.io/julia-for-hpc/>

Part V.

**Interdisciplinary Concepts and
Applications**

This section explores concepts from related fields that enhance financial modeling. We examine how ideas from computer science, statistics, and other disciplines intersect with and improve modeling practices. Through examples, we'll uncover the theoretical foundations supporting advanced techniques. This interdisciplinary approach aims to broaden your perspective and equip you with diverse tools for tackling complex financial problems.

12. Applying Software Engineering Practices

"Programs must be written for people to read, and only incidentally for machines to execute." — Harold Abelson and Gerald Jay Sussman (1984)

12.1. In this section

Modern software engineering practices such as version control, testing, documentation, and pipelines which makes modeling more robust and automated. Data practices and workflow advice.

12.2. Introduction

In addition to the core concepts from computer science described so far, there's also a similar set of ideas about the *practice* and *experience* of working with a code-based workflow that makes the end-to-end approach more powerful than the code itself.

It's likely that the majority of a professional financial modeler's time is often spent *doing things other than building models*, such as testing the model's results, writing documentation, collaborating with others on the design, and figuring out how to share the model with others inside the company. This chapter covers how a code-first workflow makes each one of those responsibilities easier or more effective.

12.2.1. Regulatory Compliance and Software Practices

Financial models often face regulatory requirements around model validation, change management, and reproducibility. Software engineering practices directly support these requirements - version control provides a complete audit trail of all model changes, automated testing helps validate model behavior and demonstrates ongoing quality control, and comprehensive documentation meets regulatory demands for model transparency. For example, the European Central Bank's Targeted Review of

12. Applying Software Engineering Practices

Internal Models (TRIM) Guide explicitly requires financial institutions to maintain documentation of model changes and validation procedures, which is naturally supported by Git commit history and continuous integration test reports which will be discussed in this chapter.

12.2.2. Chapter Structure

There are three essential topics covered in this chapter:

- **Testing** is the practice of implementing automated checks for desired behavior and outputs in the system.
- **Documentation** is the practice of writing plain English (or your local language) to compliment the computer code for better human understanding.
- **Version Control** is the systematic practice of tracking changes and facilitating collaborative workflows on projects.

As the chapter progresses, some highly related topics are covered, bridging some of the conceptual ideas into practical implementations for your own code and models.

As a reminder, this chapter is heavily oriented to concepts that are applicable in any language, though the examples are illustrated using Julia for consistency and its clarity. The code examples would have direct analogues in other languages. In Chapter 21 many of these concepts will be reinforced and expanded upon with Julia-specific workflows and tips.

12.3. Testing

Testing is a crucial aspect of software engineering that ensures the reliability and correctness of code. In financial modeling, where accuracy is paramount, implementing robust testing practices is essential, and in many cases now legally required by the regulatory body or financial reporting authority. It's good practice regardless of requirements.

A test is implemented by writing a boolean expression after a `@test` macro:

```
@test model_output = desired_output
```

If the expression evaluates to `true`, then the test passes. If the expression is anything else (`false`, or produces an error, or nothing, etc.) then the test will fail.

Here is an example of modeled behavior being tested. We have a function which will discount the given cashflows at a given annual effective interest rate. The cashflows are assumed to be equally spaced at the end of each period:

12.3. Testing

```
function present_value(discount_rate, cashflows)
    v = 1.0
    pv = 0.0
    for cf in cashflows
        v = v / (1 + discount_rate)
        pv = pv + v * cf
    end
    return pv
end

present_value (generic function with 1 method)
```

We might test the implementation like so:

```
using Test

@test present_value(0.05, [10]) ≈ 10 / 1.05
@test present_value(0.05, [10, 20]) ≈ 10 / 1.05 + 20 / 1.05^2
```

Test Passed

The above test passes because the expression is true. However, the following will fail because we have defined the function assuming the given `discount_rate` is compounded once per period. This test will fail because the test target presumes a continuous compounding convention. The failing test will show the stacktrace of where the error occurred.

```
@test present_value(0.05, [10]) ≈ 10 * exp(-0.05 * 1)
```

💡 Tip

When testing results of floating point math, it's a good idea to use the approximate comparison (`≈`, typed in a Julia editor with by entering `\approx<TAB>`). Recall that floating point math is a discrete, computer representation of continuous real numbers. As perfect precision is not efficient, very small differences can arise depending on the specific numbers involved or the order in which the operations are applied.

In tests (as in the `isapprox/≈` function), you can also further specify a relative tolerance and an absolute tolerance:

```
@test 1.02 ≈ 1.025 atol = 0.01
@test 1.02 ≈ 1.025 rtol = 0.005
```

Test Passed

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The testing described in this section is sort of a ‘sampling’ based approach, wherein the modeler decides on some pre-determined set of outputs to test and determines the desired outcomes for that chosen set of inputs. That is, testing that $2 + 2 == 4$ versus testing that a positive number plus a positive number will always equal another positive number. There are some more advanced techniques that cover the latter approach in Section 13.6.1.

💡 Tip

More Julia-specific testing workflows are covered in Section 21.14.

12.3.1. Test Driven Design

Test Driven Design (TDD) is a software development approach where tests are written before the actual code. The process typically follows these steps:

1. Write a test that defines a desired function or improvement.
2. Run the test, which should fail since the feature hasn’t been implemented.
3. Write the minimum amount of code necessary to pass the test.
4. Run the test again. If it passes, the code meets the test criteria.
5. Refactor the code for optimization while ensuring the test still passes.

TDD can be particularly useful in financial modeling as it helps clarify (1) intended behavior of the system and (2) how you think the system should work.

For example, if we want to create a new function which calculates an interpolated value between two numbers, we might first define the test like this:

```
# interpolate between two points (0,5) and (2,10)
@test interp_linear(0,2,5,10,1) ≈ (10-5)/(2-0) * (1-0) + 5
```

We’ve defined how it should work for a value inside the bounding x values, but writing the test has us wondering... should the function error if x is outside of the left and right x bounds? Or should the function extrapolate outside the observed interval? The answer to that depends on exactly how we want our system to work. Sometimes the point of such a scheme is to extrapolate, other times extrapolating beyond known values can be dangerous. For now, let’s say we would like to have the function extrapolate, so we can define our test targets to work like that:

```
@test interp_linear(0,2,5,10,-1) ≈ (10-5)/(2-0) * (-1 - 0) + 5
@test interp_linear(0,2,5,10,3) ≈ (10-5)/(2-0) * (3 - 0) + 5
```

By thinking through what the correct result for those different functions is, we have forced ourselves to think how the function should work generically:

12.3. Testing

```
function interp_linear(x1,x2,y1,y2,x)
    # slope times difference from x1 + y1
    return (y2 - y1) / (x2 - x1) * (x - x1) + y1
end
```

```
interp_linear (generic function with 1 method)
```

And we can see that our tests defined above would pass after writing the function.

```
@testset "Linear Interpolation" begin
    @test interp_linear(0,2,5,10,1) ≈ (10-5)/(2-0) * (1-0) + 5
    @test interp_linear(0,2,5,10,-1) ≈ (10-5)/(2-0) * (-1 - 0) + 5
    @test interp_linear(0,2,5,10,3) ≈ (10-5)/(2-0) * (3 - 0) + 5
end;
```

```
Test Summary: | Pass Total Time
Linear Interpolation | 3 3 0.0s
```

12.3.2. Test Coverage

Testing is great, but what if some things aren't tested? For example, we might have a function that has a branching `if/else` condition and only ever test one branch. Then when the other branch is encountered in practice it is more vulnerable to having bugs because its behavior was never double checked. Wouldn't it be great to tell whether or not we have tested all of our code?

The good news is that there is! **Test coverage** is a measurement related to how much of the codebase is covered by at least one associated test case. In the following example, code coverage would flag that the ... other logic is not covered by tests and therefore encourage the developer to write tests covering that case:

```
function asset_value(strike,current_price)
    if current_price > strike
        # ... some logic
    else
        # ... other logic
    end
end

@test asset_value(10,11) ≈ 1.5
```

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From this, it's possible to determine a score for how well a given set of code is tested. 100% coverage means every line of code has at least one test that double checked its behavior.

⚠ Warning

Testing is only as good as the tests that are written. You could have 100% code coverage for a codebase with only a single rudimentary test covering each line. Or the test itself could be wrong! Testing is not a cure-all, but does encourage best practices.

Test coverage is also a great addition when making modification to code. It can be set up such that you receive reports on how the test coverage changes if you were to make a certain modification to a codebase. An example might look like this for a proposed change which added 13 lines of code, of which only 11 of those lines were tested ("hit"). The total coverage percent has therefore gone down (-0.49%) because the proportion of new lines covered by tests is $11/13 = 84\%$, which is lower than the original coverage rate of 90%.

```
@@          Coverage Diff          @@
##      original    modif    +/-    ##
=====
- Coverage   90.00%   89.51%   -0.49%
=====
Files        2        2
Lines       130      143      +13
=====
+ Hits       117      128      +11
- Misses     13       15       +2
```

12.3.3. Types of Tests

Different tests can emphasize different aspects of model behavior. You could be testing a small bit of logic, or test that the whole model runs if hooked up to a database. The variety of this kind of testing has given rise to various named types of testing, but it's somewhat arbitrary and the boundaries between the types can be fuzzy.

Test Type	Description
Unit Testing	Verifies the functionality of individual components or functions in isolation. It ensures that each unit of code works as expected.
Integration Testing	Checks if different modules or services work together correctly. It verifies the interaction between various components of the system.

Test Type	Description
End-to-End Testing	Simulates real user scenarios to test the entire application flow from start to finish. It ensures the system works as a whole.
Functional Testing	Validates that the software meets specified functional requirements and behaves as expected from a user's perspective.
Regression Testing	Ensures that new changes or updates to the code haven't broken existing functionality. It involves re-running previously completed tests.

There are other types of testing that can be performed on a model, such as performance testing, security testing, acceptance testing, etc., but these types of tests are outside of the scope of what we would evaluate with an `@test` check. It is possible to create more advanced, mathematical-type checks and tests, which is introduced in [?@sec-related-topics-formal-verification](#).

💡 Financial Modeling Pro-tip

Test reports and test coverage are a wonderful way to demonstrate regular and robust testing for compliance. It is important to read and understand any limitations related to testing in your choice of language and associated libraries.

12.4. Documentation

The most important part of code for maintenance purposes is plainly written notes for humans, not the compiler. This includes in-line comments, docstrings, reference materials, and how-to pages, etc. Even as a single model developer, writing comments for your future self is critical for model maintenance and ongoing productivity.

12.4.1. Comments

Comments are meant for the developer to aid in understanding a certain bit of code. A bit of time-tested wisdom is that after several weeks, months, or years away from a piece of code, something that seemed 'obvious' at the time tends to become perplexing at a later time. Writing comments is as much for yourself as it is your colleagues or successors.

Here's an example of documentation with single-line comments (indicated with the preceding `#`) and multi-line comments (enclosed by `#=` and `=#`):

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```
function calculate_bond_price(face_value, coupon_rate, years_to_maturity, market_rate)
    # Convert annual rates to semi-annual
    semi_annual_coupon = (coupon_rate / 2) * face_value
    semi_annual_market_rate = market_rate / 2
    periods = years_to_maturity * 2

    # Calculate the present value of coupon payments
    pv_coupons = 0
    for t in 1:periods
        pv_coupons += semi_annual_coupon / (1 + semi_annual_market_rate)^t
    end

    # Calculate the present value of the face value
    pv_face_value = face_value / (1 + semi_annual_market_rate)^periods

    #=
    Sum up the components for total bond price
    1. Present value of all coupon payments
    2. Present value of the face value at maturity
    =#
    bond_price = pv_coupons + pv_face_value

    return bond_price
end
```

12.4.2. Docstrings

Docstrings (documentation strings) are intended to be a user-facing reference and help text. In Julia, docstrings are just strings placed in front of definitions. Markdown¹ is available (and encouraged) to add formatting within the docstring.

Here's an example with some various features of documentation shown:

```
"""
    calculate_bond_price(face_value, coupon_rate, years_to_maturity, market_rate) # <1>

Calculate the price of a bond using discounted cash flow method.
```

Parameters:

- `face_value`: The bond's par value

¹Markdown is a type of plain text which can be styled for better communication and aesthetics. E.g. ****some text**** would render as boldface: **some text**. This book was written in Markdown and all of the styling arose as a result of plain text files written with certain key elements.

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- `coupon_rate`: Annual coupon rate as a decimal
- `years_to_maturity`: Number of years until the bond matures
- `market_rate`: Current market interest rate as a decimal

Returns:

- The calculated bond price

Examples: # <2>

```
```julia-repl
julia> calculate_bond_price(1000, 0.05, 10, 0.06)
925.6126256977221
````
```

Extended help: # <3>

This function uses the following steps to calculate the bond price:

1. Convert annual rates to semi-annual rates
2. Calculate the present value of all future coupon payments
3. Calculate the present value of the face value at maturity
4. Sum these components to get the total bond price

The calculation assumes semi-annual coupon payments, which is standard in many markets.

```
"""
function calculate_bond_price(face_value, coupon_rate, years_to_maturity, market_rate)
    # ... function defintion as above
end
```

- ① The typical docstring on a method includes the signature, indented so it's treated like code in the Markdown docstring.
- ② It's good practice to include a section that includes examples of appropriate usage of a function and the expected outcomes.
- ③ The *Extended help* section is a place to put additional detail that's available on generated docsites and in help tools like the REPL help mode.

The last point, the *Extended help* section is shown when using help mode in the REPL and including an extra ?. For example, in the REPL, typing ?calculate_bond_price will show the docstring up through the examples. Typing ??calculate_bond_price will show the docstring in its entirety.

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12.4.3. Docsites

Docsites, or documentation sites, are websites that are generated to host documentation related to a project. With modern tooling around programming projects, a really rich set of interactive documentation can be created while the developer/modeler focuses on simple documentation artifacts.

Specifically, modern docsite tooling generally takes in markdown text pages along with the in-code docstrings and generates a multi-page site that has navigation and search.

Typical contents of a docsite include:

- A *Quickstart* guide that introduces the project and provides an essential or common use case that the user can immediately run in their own environment. This helps to convey the scope and capability of a project and showcase how to model is intended to be used.
- *Tutorials* are typically worked examples that introduce basic aspects of a concept or package usage and work up to more complex use cases.
- *Developer documentation* is intended to be read by those who are interested in understanding, contributing, or modifying the codebase.
- *Reference documentation* describes concepts and available functionality. A subset of this is API, or **Application Programming Interface**, documentation which is the detailed specification of the available functionality of a package, often consisting largely of docstrings like the previous example.

Docsite generators are generally able to look at the codebase and from just the docstrings create a searchable, hierarchical page with all of the content from the docstrings in a project. This basic docsite feature is incredibly beneficial for current and potential users of a project.

To go beyond creating a searchable index of docstrings requires additional effort (time well invested!). Creating the other types of documentation (quick start, tutorials, etc.) is mechanically as simple as creating a new markdown file. The hard part is learning how to write quality documentation. Good technical writing is a skill developed over time - but at least the technical and workflow aspects have been made as easy as possible!

12.5. Version Control

Version control systems (VCS) refer to the tracking of changes to a codebase in a verifiable and coordinated way across project contributors. VCS underpins many aspects of automating the mundane parts of a modeler's job. Benefits of VCS include (either directly, or contribute significantly to):

- Access control and approval processes
- Versioning of releases
- Reproducibility across space and time of a model's logic
- Continuous testing and validation of results
- Minimization of manual overrides, intervention, and opportunity for user error
- Coordinating collaboration in parallel and in sequence between one or many contributors

These features are massively beneficial to a financial modeler! A lot of the overhead in a modeler's job becomes much easier or automated through this tooling.

Among several competing options (CVS, Mercurial, Subversion, etc.), Git is the predominant choice as of this book's writing and therefore we will focus on Git concepts and workflows.

12.5.1. Git Overview

This section will introduce Git related concepts at a level deeper than "here's the five commands to run" but not at a reference-level of detail. The point of this is to reveal some of the underlying technology and approaches so that you can recognize where similar ideas appear in other areas and understand some of the limitations of the technology.

Git is a free and open source software that tracks changes in files using a series of snapshots. Git itself is the underlying software tool and is a command-line tool at its core. However, you will often interact with it through various interfaces (such as a graphical user interface in VS Code, GitKraken, or other tool).

Each snapshot, or **commit**, stores references to the files that have changed². All of this is stored in a `.git` subfolder of a project, which is automatically created upon initializing a repository. This folder may be hidden by default by your filesystem. You generally never need to modify anything inside the `.git` folder yourself as Git handles this for you. Git tracks files that are contained in sibling directories to the `.git` subfolder.

i Note

A directory structure demonstrating where git data is stored and what content is tracked after initializing a repository in the `tracked-project` directory. Note how `untracked` folder is not the parent directory of the `.git` directory so it does not get tracked by Git.

²Git uses a content-addressable filesystem, meaning it stores data as key-value pairs. The key is a hash of the content, ensuring data integrity and allowing efficient storage of identical content. For more on hashes, jump to ([note-content-hash?](#)).

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```
/home/username
└── untracked-folder
    ├── random-file.txt
    └── ...
└── tracked-project
    ├── .git
    │   ├── config
    │   ├── HEAD
    │   └── objects
    │       └── ...
    ├── .gitignore
    ├── README.md
    └── src
        ├── MainProject.jl
        ├── module1.jl
        └── module2.jl
```

Under the hood, the Git data stored inside the `.git/` include things such as binary blobs, trees, commits, and tags. Blobs store file contents, trees represent directories, commits point to trees and parent commits, and tags provide human-readable names for specific commits. This structure allows Git to efficiently handle branching, merging, and maintaining project history.

Hashes indicate a verifiable snapshot of a codebase. For example, a full commit ID (a hash) `40f141303cec3d58879c493988b71c4e56d79b90` will *always* refer to a certain snapshot of the code, and if there is a mismatch in the git history between the contents of the repository and the commit's hash then the git repository is corrupted and it will not function. A corrupted repository usually doesn't happen in practice, of course! You might see hashes shortened to the last several characters (e.g. `6d79b90`) and in the following examples we'll shorten the hypothetical hashes to just three characters (e.g. `b90`).

A codebase can be **branched**, meaning that two different version of the same codebase can be tracked and switched between. Git lends itself to many different workflows, but a common one is to designate a primary branch (call it `main`) and make modifications in a new, separate branch. This allows for non-linear and controlled changes to occur without potentially tainting the `main` branch. It's common practice to always have the `main` branch be a 'working' version of the code that can always 'run' for users, while branches contain works-in-progress or piecemeal changes that temporarily make the project unable to run.

The commit history forms a directed acyclic graph (DAG) representing the project's

12.5. Version Control

history. That is, there is an order to the history: the second commit is dependent on the first commit. From the second commit, two child commits may be created which both have the second commit as their parent. Each one of this commits represents a stored snapshot of the project at the time of the commit.

Table 12.2 shows an example workflow for trying to fix an erroneous function `present_value` which was written as part of a hypothetical `FileXYZ.jl` file. This example is trivial, but in a larger project where a ‘fix’ or ‘enhancement’ may span many files and take several days or weeks to implement this type of workflow becomes like a superpower compared to traditional, manual version control approaches. It sure beats an approach where you end up with filenames like `FileXYZ v23 July 2022 Final.jl`!

Table 12.2.: A workflow demonstrating branching, staging, committing, and merging in order to fix an incorrect function definition for a present value (pv) function. The branch name/commit ID shows which version you would see on your own computer/filesystem. The inactive branches are tracked in Git but do not manifest themselves on your filesystem unless you checkout that branch.

| Branch main, FileXYZ.jl file | Branch fix_function, FileXYZ.jl file | Action | Active Branch (Commit ID) |
|--|---|---|---------------------------|
| <code>function pv(rate,amount,time)
 amount / (1+rate)
end</code> | Does not yet exist | Write original function which forgets to take into account the time. Stage and commit it to the main branch. | main
(...58b) |
| " | <code>function pv(rate,amount,time)
 amount / (1+rate)
end</code> | Git commands:
<code>git add FileXYZ.jl</code>
<code>git commit -m 'add pv function'</code>
Create a new branch fix_function. | main
(...58b) |
| " | " | Git command:
<code>git branch fix_function</code>
Checkout the new branch and make it active for editing. The branch is different but is starting from the existing commit. | fix_function
(...58b) |
| " | <code>function pv(rate,amount,time)
 amount / (1+rate)^time
end</code> | Git command: <code>git checkout fix_function</code>
Edit and save the changed file. No git actions taken. | fix_function
(...58b) |
| " | " | "Stage" the modified file, telling git that you are ready to record ("commit") a new snapshot of the project. | fix_function
(...58b) |
| " | " | Git command: <code>git add FileXYZ.jl</code>
Commit a new snapshot of the project by committing with a note to collaborators saying fix: present value logic
Git command: <code>git commit -m 'fix: present value logic'</code> | fix_function
(...6ac) |

| Branch main, FileXYZ.jl file | Branch fix_function, FileXYZ.jl file | Action | Active Branch
(Commit ID) |
|---|--------------------------------------|---|------------------------------|
| " | " | Switch back to the primary branch | main
(...58b) |
| <pre>function pv(rate,amount,time) amount / (1+rate)^time end</pre> | " | Git command: git checkout main
Merge changes from other branch into the main branch, incorporating the corrected version of the code. | main
(...b90) |
| | | Git command: git merge fix_function | |

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A visual representation of the git repository and commits for the actions described in Table 12.2 might be as follows, where the ...XXX is the shortened version of the hash associated with that commit.

```
main branch      : ...58b      → ...b90
                  ↓           ↗
fix_function branch : ...58b → ...6ac
```

The “staging” aspect will be explained next.

12.5.1.1. Git Staging

Staging (sometimes called the “staging area” or “index”) is an intermediate step between making changes to files and recording those changes in a commit. Think of staging as preparing a snapshot - you’re choosing which modified files (or even which specific changes within files) should be included in your next commit. When you modify files in your Git repository, those changes are initially “unstaged.” Using the git add command (or your Git GUI’s staging interface) moves changes into the staging area. This two-step process - staging followed by committing - gives you precise control over which changes get recorded together. Here’s a typical workflow:

1. You modify several files in your modeling project
2. You review the changes and decide which ones are ready to commit
3. You stage only the changes that belong together as a logical unit
4. You create a commit with just those staged changes
5. Repeat as needed with other modified files

This is particularly useful when you’re working on multiple features or fixes simultaneously. For example, imagine you’re updating a financial model and you:

- Fix a bug in your present value calculation
- Add comments to improve documentation
- Start working on a new feature

You could stage and commit the bug fix and documentation separately, keeping the work-in-progress feature changes unstaged until they’re complete. This creates a cleaner, more logical project history where each commit represents one coherent change.

Think of staging as preparing a shipment: you first gather and organize the items (staging), then seal and send the package (committing). This extra step helps maintain a well-organized project history where each commit represents a logical, self-contained change.

12.5.1.2. Git Tooling

Git is traditionally a command-line based tool, however we will not focus on the command line usage as more beginner friendly and intuitive interfaces are available from different available software. The branching and nodes are well suited to be represented visually.

Note

Some recommend Git tools with a **graphical user interface (GUI)** :

- *Github Desktop* interfaces nicely with Github and provides a GUI for common operations.
- *Visual Studio Code* has an integrated Git pane, providing a powerful GUI for common operations.
- *GitKraken* is free for open source repositories but requires payment for usage in enterprise or private repository environments. GitKraken provides intuitive interfaces for handling more advanced operations, conflicts, or issues that might arise when using Git.

12.5.2. Collaborative Workflows

Git is a distributed VCS, meaning that copies of the repository and all its history and content can live in multiple locations, such as on two colleagues' computer as well as a server. In this distributed model, what happens if you make a change locally?

Git maintains a local repository on each user's machine, containing the full project history. This local repository includes the working directory (current state of files), staging area (changes prepared for commit), and the .git directory (metadata and object database). When collaborating, users push and pull changes to and from remote repositories. Git uses a branching model, allowing multiple lines of development to coexist and merge when ready.

Note

By convention, the primary branch of a project is named the `main` branch. Historically, this was called the `master` branch (after the Master-Slave technology terms). Due to connotations of human slavery most software will now default to the `main` name. However, many projects continue to use or have never renamed their original `master` branch.

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12.5.2.1. Pull Requests

Layered onto core Git functionality, services like Github provide interfaces which enhance the utility of VCS. A major example of this is **Pull Requests** (or PRs), which is a process of merging git branches in a way that allows for additional automation and governance.

The following is an example of a PR on a repository adding a small bit of documentation to help future users. We'll walk through several elements to describe what's going on:

Referencing Figure 12.1, several elements are worth highlighting. In this pull request the author of this change, alecloudenback, is proposing to modify the QuartoNotebookRunner repository, which is not a repository for which the user alecloudenback has any direct rights to modify. After having made a copy of the repository (a "fork"), creating a new branch, making modifications, and committing those changes... a pull request has been made to modify the primary repository.

- All changes are recorded using Git, keeping track of authorship, timestamps, and history.
- At the top of the screenshot, the title "Note that Quarto..." allows the author to summarize what is changed in the branch to be merged.
- The "Conversation" tab allows for additional details about the change to be discussed.
- In the top right, a **+1 -1**  is an indication of what's changed. In this case, a single line of code (documentation, really) was removed (-1) and replaced with another (+1).
- Not shown in the Figure 12.1, the "Files Changed" tab shows a file-by-file and line-by-line comparison of what has changed (see Figure 12.2).
- The reviewer (user MichaelHatherly) is a collaborator with rights to modify the destination repository has been assigned to review this change before merging it into the `main` branch.
- "CodeCoverage" was discussed above in testing, and in this case tests were automatically run when the PR was created, and the coverage indicates that there was no added bit of code that was untested.
- A section of "9 checks" that pass, which validated that tests within the repository still passed, on different combinations of operating systems and Julia versions. Additionally, the repository was checked to ensure that formatting conventions were followed.
- Additionally, there are a number of project management features not really showcased here. A few to note:
 - The cross-referencing to another related issue in a different repository (the reference to issue #156).
 - Assignees (assigned doers), labels, milestones, and projects are all functionality to help keep track of codebase development.

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Note that Quarto doesn't follow project-wide engine yet #157

The screenshot shows a GitHub pull request interface. At the top, a purple banner indicates the pull request has been merged. Below this, the main header includes 'Conversation 2', 'Commits 1', 'Checks 9', 'Files changed 1', and a diff summary '+1 -1'. Navigation buttons for 'Edit' and 'Code' are also present.

The pull request details show it was merged by MichaelHatherly from alecloudenback:patch-1 into PumasAI:main on Jun 20. The author, alecloudenback, commented on Jun 19, noting that Quarto doesn't follow a project-wide engine yet. This comment is marked as 'Verified' and has a green checkmark icon. MichaelHatherly approved the changes on Jun 19, leaving a thank you message. A bot named codecov commented on Jun 20, reporting a 'Codecov Report' where all modified and coverable lines are covered by tests (indicated by a green checkmark).

On the right side of the pull request page, there are sections for 'Reviewers' (MichaelHatherly), 'Assignees' (None assigned), 'Labels' (None yet), 'Projects' (None yet), 'Milestone' (No milestone), and 'Development' (Successfully merging may close issues). There is also a 'Notifications' section with an 'Unsubscribe' button and a note about receiving notifications because the user authored the thread. A checkbox for 'Allow edits by maintainers' is also present.

At the bottom of the pull request view, a summary shows 9 checks passed, including various test and format checks across different operating systems and versions.

Figure 12.1.: A Pull Request on Github, demonstrating several utilities which enhance change management, automation, and governance.

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The features described here can take modelers many manhours of time - testing, review of changes, sign-off tracking, etc. In this Git-based workflow, the workflow above happens frictionlessly and much of it is automated. This is such a powerful paradigm that should be adopted within the financial industry and especially amongst modelers.

Note that Quarto doesn't follow project-wide engine yet #157

The screenshot shows a GitHub pull request interface. At the top, a purple button says 'Merged' and indicates 'MichaelHatherly merged 1 commit into PumasAI:main from alecloudenback:patch-1'. Below this, there are tabs for 'Conversation' (2), 'Commits' (1), 'Checks' (9), and 'Files changed' (1). The 'Files changed' tab is selected, showing a diff for 'README.md'. The diff highlights changes in lines 11 through 17. Lines 11-13 show text being added. Line 14 shows text being removed (indicated by a red background) and replaced by new text (indicated by green background). Lines 15-17 show unchanged text. The GitHub interface includes standard navigation and search tools at the bottom.

Figure 12.2.: A diff shows a file-by-file and line-by-line comparison of what has changed between commits in a codebase. Red indicates something was removed or changed, and green shows what replaced it. Note that even within a line, there's extra green highlighting to show the newly added text while the unchanged text remains a lighter shade.

12.5.3. Data Version Control

Git is not well suited for large files (images, videos, datasets) that change regularly. Instead, the approach is to combine git with a **data version control** tool. These tools essentially replace the data (often called binary data or a blob) with a content hash in the git snapshots. The actual content referred to by the hash is then located elsewhere (e.g. a hosted server).

```
::: {#note-content-hash callout-note} ## Content Hashes
```

Content hashes are the output of a function that transforms arbitrary data into a string of data. Hashes are very common in cryptography and in areas where security/certainty is important. Eliding the details of how they work exactly, what's important to understand for our purposes is that a content hash function will take arbitrary bits and output the same set of data each time the function is called on the original bits.

For example, it might look something like this:

```
data = "A data file contents 123"
Base.hash(data)
```

0x7a1f7a6e6f61fa64

If it's run again on the same inputs, you get the same output:

```
Base.hash(data)
```

0x7a1f7a6e6f61fa64

And if the data is changed even slightly, then the output is markedly different:

```
data2 = "A data file contents 124"
Base.hash(data2)
```

0xb0b70b13a7d93dcb

This is used in content addressed systems like Data Version Control (Section 12.5.3) and Artifacts (sec ([artifacts?](#))) to ask for, and confirm the accuracy of, data instead of trying to address the data by its location. That is, instead of trying to ask to get data from a given URL (e.g. <http://adatasource.com/data.csv>) you can set up a system which keeps track of available locations for the data that matches the content hash. Something like (in Julia-ish psuedocode):

```
storage_locations = Dict(
    0x4d7e8e449af1c48 => [
        "http://datasets.com/1231234",
        "http://companyintranet.com/admin.csv",
        "C:/Users/your_name/Documents/Data/admin.csv"
    ]
)
function get_data(hash, locations)
```

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```
for location in locations[hash]
    if is_available(location)
        return get(location)
    end
end

# if loop didn't return data
return nothing
end
```

⋮

12.6. Distributing the Package

Once you have created something, the next best feeling after having it working is having someone else also use the tool. Julia has a robust way to distribute and manage dependencies. This section will cover essential and related topics to distributing your project publicly or with an internal team.

12.6.1. Registries

Registries are a way to keep track of packages that are available to install, which is more complex than it might seem at first. The registry needs to keep track of:

- What dependencies your package has.
- What versions of the dependencies your package is compatible with.
- Metadata about the package (name, unique identifier, authors).
- Versions of your package that you have made available and the Git hash associated with that version for provenance and tracking
- The location where your package's repository lives so that the user can grab the code from there.

The Julia General Registry (“General”) is the default registry that comes loaded as the default registry when installing Julia. From a capability standpoint, there's nothing that separates General from other registries, including ones that you can create yourself. At its core, the registry is seen as a Git repository where each new commit just adds information associated with the newly registered package or version of a package.

For distributing packages in a private, or smaller public group see Section 23.9.2.1.

12.6.1.1. General Registry and other Hosted Registries

At its core, General is essentially the same as a local registry described in the prior section. However, there's some additional infrastructure supporting General. Registered packages get backed up, cached for speed, and multiple servers across the globe are set up to respond to Pkg requests for redundancy and latency. Nothing would stop you from doing the same for your own hosted registry if it got popular enough!

💡 Tip

A local registry is a great way to set up internal sharing of packages within an organization. Services do exist for “managed” package sharing, adding enterprise features like whitelisted dependencies, documentation hosting, a ‘hub’ of searchable packages.

12.6.2. Versioning

Versioning is an important part of managing a complex web of dependencies. Versioning is used to let both users and the computer (e.g. Pkg) understand which bits of code are compatible with others. For this, consider your model/program’s **Application Programming Interface** (API). The API is essentially defined by the outputs produced by your code given the same inputs. If the same inputs are provided, then for the same API version the same output should be provided. However, this isn’t the only thing that matters. Another is the documentation associated with the functionality. If the documentation said that a function would work a certain way, then your program should follow that (or fix the documentation)!

Another case to consider is what if new functionality was *added*? Old code need not have changed, but if there’s new functionality, how to communicate that as an API change? This is where **Semantic Versioning** (SemVer) comes in: *semantic* means that your version something is intended to convey some sort of meaning over and above simply incrementing from v1 to v2 to v3, etc.

12.6.2.1. Semantic Versioning

Semantic Versioning (SemVer) is one of the most popular approaches to software versioning. It’s not perfect but has emerged as one of the most practical ways since it gets a lot about version numbering right. Here’s how SemVer is defined³:

³You can read more at SemVer.org.

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Given a version number MAJOR.MINOR.PATCH, increment the:

- MAJOR version when you make incompatible API changes
- MINOR version when you add functionality in a backward compatible manner
- PATCH version when you make backward compatible bug fixes

So here are some examples of SemVer, if our package's functionality for v1.0.0 is like this:

```
""" my_add(x,y)

Add the numbers x and y together
"""

my_add(x,y) = x - y
```

Patch change (v1.0.1): Fix the bug in the implementation:

```
""" my_add(x,y)
Add the numbers x and y together
"""

my_add(x,y) = x + y
```

This is a patch change because it fixes a bug without changing the API. The function still takes two arguments and returns their sum, as originally documented.

Minor change (v1.1.0): Add new functionality in a backward-compatible manner:

```
""" my_add(x,y)
Add the numbers x and y together

my_add(x,y,z)
Add the numbers x, y, and z together
"""

my_add(x,y) = x + y
my_add(x,y,z) = x + y + z
```

This is a minor change because it adds new functionality (the ability to add three numbers) while maintaining backward compatibility with the existing two-argument version.

Major change (v2.0.0): Make incompatible API changes:

```
""" add_numbers(numbers...)
Add any number of input arguments together. This
function replaces 'my_add' from prior versions.
"""

add_numbers(numbers...) = sum(numbers)
```

12.6. Distributing the Package

This is a major change because it fundamentally alters the API. The function name has changed, and it now accepts any number of arguments instead of specifically two or three. This change is not backward compatible with code using the previous versions, so a major version increment is necessary.

Note

Numbers need not roll over to the next digit when they hit 10. That is, it's perfectly valid to go from `v1.09.0` to `v1.10.0` in SemVer.

Tip

Sometimes you'll see a package with a version that starts with zero, such as `v0.23.1`. **We recommend that as soon as you register a package, to make it a v1.** `v1` need not indicate the package is "complete" (what software is?), so don't hold back on calling it `v1`. You're letting users install it easily, so you might as well call it the first version and move on!

According to SemVer's rules, there are no patch versions when the major version is zero. This means that you have one less meaningful digit to communicate to users what's going on with a change. Most packages put an upper bound on compatibility so that major or minor changes in upstream packages are less likely to cause issues in their own packages. This can be somewhat painful to depend on a package which has a `v0` and is iterating through 'fixes' but is incrementing the minor version. You have to assume it's making backward incompatible changes and should have skepticism of just upgrading to the new version of the dependency. It takes work on the downstream dependencies to decide if they should upgrade, adding mental and time loads to other authors and users.

12.6.3. Artifacts

Artifacts are a way to distribute content-addressed ((**note-content-hash?**)) data and other dependencies. An example use case is if you want to distribute some demonstration datasets with a package. When a package is added or updated, the associated data is pulled and un-archived by Pkg instead of the author of the package needing to manually handle data dependencies. Aside from this convenience, it means that different packages could load the same data without duplicating the data download or storage (since the data is content-addressed). The use-case is not real-time data, as the content-hash can only be updated per package version.

For example, the `MortalityTable.jl` package redistributes various publicly available, industry mortality tables. Inside the repository, there's an `Artifacts.toml` file specified like:

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```
#/Artifacts.toml  
  
["mort.soa.org"]  
git-tree-sha1 = "6164a6026f108fe95828b689fc3b992acb7c3"  
①  
  
[["mort.soa.org".download]]  
sha256 = "6f5eb4909564b55a3397ccf4f4c74290e002f7e2e2474cebeb224bb23a9a2606" ②  
url = "https://github.com/JuliaActuary/Artifacts/raw/v2020-02-15/mort.soa.org/2020-02-15.t
```

- ① The sha1 hash of the un-archived data once downloaded, used to verify that extraction was successful.
- ② The sha256 hash of the archived (compressed) data to ensure that the data downloaded was as intended.

Then, within the package the data artifact can be referenced and handled by the artifact system rather than needing to manually handle it. That is, the data is reference-able like this:

```
table_dir = artifact"mort.soa.org" ①
```

- ① artifact"..." is a string macro, which is a special syntax for macros that interact with strings. md"..." is another example, specifying that the content of the string is Markdown content.

As opposed to something like this:

```
# psuedo Julia code  
table_dir = if is_first_package_run  
  
    data_path = download("url_of_data.tar.gz") # download to a temp location  
    mv(data_path, "/somewhere/to/keep/data.tar.gz") # move to a 'permanent' location  
    extract("/somewhere/to/keep/data.tar.gz") # extract contents  
    "/somewhere/to/keep/data/" # return data path  
else  
    "/somewhere/to/keep/data/" # return pre-processed path  
  
end
```

Note

Utility Packages such as `ArtifactUtils.jl` can assist in creating correct entries for `Artifact.toml` files.

i Note

Artifacts support .tar (uncompressed) and .tar.gz because that compression format enjoys more universal support and features than the .zip format most common on Windows systems.

12.7. Example Repository

The is a good example of a repository which shows the organization of files and code, setting up testing, and documentation.

12.8. Example Repository Structure

A well-structured Julia package demonstrates key software engineering principles in action. The JuliaTemplateRepo repository demonstrates best practices for:

- Logical file organization and code structure
- Comprehensive test coverage and continuous integration
- Clear, accessible documentation with examples
- Standard tooling configuration for package development

This open-source template serves as a reference implementation.

The PkgTemplates.jl package will allow you to create an empty repository with all of the testing, documentation, Git, and continuos integration scaffolding already in place.

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“Fundamentally, computer science is a science of abstraction—creating the right model for a problem and devising the appropriate mechanizable techniques to solve it. Confronted with a problem, we must create an abstraction of that problem that can be represented and manipulated inside a computer. Through these manipulations, we try to find a solution to the original problem.” - Al Aho and Jeff Ullman (1992)

13.1. In this section

Adapting computer science concepts to work for financial professionals. Computability, computational complexity, and the language of algorithms and problem solving. A survey of important data structures. More advanced testing and verification topics.

13.2. Computer Science for Financial Professionals

Computer science as a term can be a bit misleading because of the overwhelming association with the physical desktop or laptop machines that we call “computers”. The discipline of computer science is much richer than consumer electronics: at its core, computer science concerns itself with areas of research and answering tough questions:

- **Algorithms and Optimization.** How can a problem be solved efficiently? How can that problem be solved *at all*? Given constraints, how can one find an optimal solution?
- **Theory of Computation.** What sorts of questions are even answerable? Is an answer easy to compute or will resolving it require more resources than the entire known universe? Will a computation ever stop calculating?
- **Data Structures.** How to encode, store, and use data? How does that data relate to each other and what are the trade-offs between different representations of that data?
- **Information Theory**¹. Given limited data, what *can* be known or inferred from it?

¹This topic will be covered in Chapter 14.

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For a reader in the twenty-first century we hope that it is patently obvious how impactful the *applied* computer science has been as an end-user of the internet, artificial intelligence, computational photography, safety control systems, etc. have been to our lives. It is a testament to the utility of being able to harness computer science ideas for practical use.

It's common for beneficial advances in knowledge and application to occur at the boundary between two disciplines. It's here in this chapter that we desire to bring together the financial discipline together with computer science and to provide the financial practitioner with the language and concepts to leverage some of computer science's most relevant ideas.

13.3. Algorithms

An **Algorithm** is a general term for a process that transforms an input to an output. It's the set of instructions dictating how to carry out a process. That process needs to be specified in sufficient detail to be able to call itself an algorithm (versus a *heuristic* which lacks that specificity).

An algorithm might be the directions to accomplish the following task: summing a series of consecutive integers integers from 1 to n . There are multiple ways that this might be accomplished, each one considered a distinct algorithm:

- iterating over each number and summing them up (starting with the smallest number)
- iterating over each number and summing them up (starting with the largest number)
- summing up the evens and then the odds, then adding the two subtotals
- and many more distinct algorithms...

We will look at some specific examples of these alternate algorithms as we introduce the next topic, computational complexity.

13.4. Complexity

13.4.1. Computational Complexity

We can characterize the computational complexity of a problem by looking at how long an algorithm takes to complete a task when given an input of size n . We can then compare two approaches to see which is computationally less complex for a given n . This is

13.4. Complexity

a way of systematically evaluating an algorithm to determine its efficiency when being computed.

Note that computational complexity isn't quite the same as how fast an algorithm will run on your computer, but it's a very good guide. Modern computer architectures can sometimes execute multiple instructions in a single cycle of the CPU making an algorithm that is, on paper slower than another, actually run faster in practice. Additionally, sometimes algorithms are able to substantially limit the number of *computations* to be performed, at the expense of using a lot more *memory* and thereby trading CPU usage with RAM usage.

You can think of computational complexity as a measure of how much work is to be performed. Sometimes the computer is able to perform certain kinds of work more efficiently.

Further, when we analyze an algorithm recall that ultimately our code gets translated into instructions for the computer hardware. Some instructions are implemented in a way that for any type of number (e.g. floating point), it doesn't matter if the number is 1.0 or 0.41582574300044717, the operation will take the exact same time and number of instructions to execute (e.g. for the addition operation).

Sometimes a higher level operation is implemented in a way that takes many machine instructions. For example, division instructions may require many CPU cycles when compared to multiplication or division. Sometimes this is an important distinction and sometimes not. For this book we will ignore this granularity of analysis.

13.4.1.1. Example: Sum of Consecutive Integers

Take for example the problem of determining the sum of integers from 1 to n . We will explore three different algorithms and the associated computational complexity for them.

13.4.1.1.1. Constant Time

A mathematical proof can show a simple formula for the result. This allows us to compute the answer in **constant time**, which means that for any n , our algorithm is essentially the same amount of work.

```
nsum_constant(n) = n * (n + 1) / 2
```

```
nsum_constant (generic function with 1 method)
```

In this we see that we perform three operations: a multiplication, a sum, and a division, no matter what n is. If n is 10_000_000 we'd expect this to complete in about a single unit of time.

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13.4.1.1.2. Linear Time

This algorithm performs a number of operations which grows in proportion with n by individually summing up each element in 1 through n :

```
function nsum_linear(n)
    result = 0
    for i in 1:n
        result += i
    end

    result
end

nsum_linear (generic function with 1 method)
```

If n were 10_000_000, we'd expect it to run with roughly 10 million operations, or about 3 million times as many operations as the constant time version. We can say that this version of the algorithm will take approximately n steps to complete.

13.4.1.1.3. Quadratic Time

What if we were less efficient, and instead we were only ever able to increment our subtotal by one. That is, instead of adding up 1+3, we had to instead do four operations: 1 + 1 + 1 + 1 . We can add a second loop which increments our result by a unit instead of simply adding the current i to the running total $result$. This makes our algorithm work much harder since it has to add numbers so many more times (Recall that to a computer adding two numbers is the same computational effort regardless of what the numbers are).

```
function nsum_quadratic(n)
    result = 0
    for i in 1:n
        for j in 1:i
            result += 1
        end
    end

    result
end
```

- ① The outer loop with iterator i . loops over the integers 1 to n .
- ② The inner loop with iterator j does the busy work of adding 1 to our subtotal i times.

```
nsum_quadratic (generic function with 1 method)
```

Breaking down the steps:

- When i is 1 there is 1 addition in the inner loop
- When i is 2 there are 2 additions in the inner loop
- ...
- When i is n there are n additions in the inner loop

Therefore, this computation takes $1 + \dots + (n - 2) + (n - 1) + n$ steps to complete. Algebraically, this simplifies down to our constant time formula $n * (n + 1) \div 2$ or $n^2 + n \div 2$ steps to complete.

13.4.1.2. Comparison

13.4.1.2.1. Big-O Notation

We can categorize the above implementations using a convention called **Big-O Notation**² which is a way of distilling and classifying computational complexity. We characterize the algorithms by the most significant term in the total number of operations. Table 13.1 shows for the examples constructed above what the description, order, and order of magnitude complexity is.

Table 13.1.: Complexity comparison for the three sample cases of summing integers from 1 to n .

| Function | Computational Cost | Complexity Description | Big-O Order | Steps ($n = 10,000$) |
|----------------|--------------------|------------------------|-------------|------------------------|
| nsum_constant | fixed | Constant | $O(1)$ | ~1 |
| nsum_linear | n | Linear | $O(n)$ | ~10,000 |
| nsum_quadratic | $n^2 + n \div 2$ | Quadratic | $O(n^2)$ | ~100,000,000 |

Table 13.2 shows a comparison of a more extended set of complexity levels. For the most complex categories of problems, the cost to compute grows so fast that it boggles the mind. What sorts of problems fall into the most complex categories? $O(2^n)$, or exponential complexity, examples include the traveling salesman problem³ if solved with

²“Big-O”, so named because of the “O” in used in $O(1)$. $O(n)$, etc. The “O” refers to the “order” of growth for the function.

³The Traveling Salesperson Problem is a classic computer science problem where you need to find the shortest possible route that visits each city in a given set exactly once and returns to the starting city. It’s a seemingly simple problem that becomes computationally intensive very quickly as the number of cities increases.

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dynamic programming or the recursive approach to calculating the n th Fibonacci number. The beastly $O(n!)$ algorithms include brute force solving the traveling salesman problem or enumerating all partitions of a set. In financial modeling, we may encounter these sorts of problems in portfolio optimization (using the brute-force approach of testing every potential combination assets to optimize a portfolio).

Table 13.2.: Different Big-O Orders of Complexity

| Big-O Order | Description | $n = 10$ | $n = 1,000$ | $n = 1,000,000$ |
|-----------------------|-------------------|-----------|------------------|---------------------|
| $O(1)$ | Constant Time | 1 | 1 | 1 |
| $O(n)$ | Linear Time | 10 | 1,000 | 1,000,000 |
| $O(n^2)$ | Quadratic Time | 100 | 1,000,000 | 10^{12} |
| $O(\log(n))$ | Logarithmic Time | 3 | 7 | 14 |
| $O(n \times \log(n))$ | Linearithmic Time | 30 | 7,000 | 14,000,000 |
| $O(2^n)$ | Exponential Time | 1,024 | $\sim 10^{300}$ | $\sim 10^{301029}$ |
| $O(n!)$ | Factorial Time | 3,628,800 | $\sim 10^{2567}$ | $\sim 10^{5565708}$ |

i Note

We care only about the most significant term because when n is large, the most significant term tends to dominate. For example, in our quadratic time example which has $n^2 + n \div 2$ steps, if n is a large number like 10^6 , then we see that it will result in:

$$\begin{aligned} n^2 + \frac{n}{2} &= (10^6)^2 + \frac{10^6}{2} \\ &= 10^{12} + \frac{10^6}{2} \end{aligned}$$

10^{12} is significantly more important than $\frac{10^6}{2}$ (sixty-four million times as important, to be precise). This is why Big-O notation reduces the problem down to only the most significant complexity cost term.

If n is small then we don't really care about computational complexity in general. This is a lesson for our efforts as developers: focus on the most intensive parts of calculations when looking to optimize, and don't worry about seldom used portions of the code.

13.4.1.2.2. Empirical Results

The preceding examples of constant, linear, and exponential times are *conceptually* correct but if we try to run them in practice we see that the description doesn't seem to hold at all for the linear time version, as it runs as quickly as the constant time version.

```
using BenchmarkTools
@btime nsum_constant(10_000)

0.833 ns (0 allocations: 0 bytes)

50005000

@btime nsum_linear(10_000)

1.500 ns (0 allocations: 0 bytes)

50005000

@btime nsum_quadratic(10_000)

1.104 μs (0 allocations: 0 bytes)
```

What happened was that the compiler was able to understand and optimize the linear version such that it effectively transformed it into the constant time version and avoid the iterative summation that we had written. For examples that are simple enough to use as a teaching problem, the compiler can often optimize different written code down to the same efficient machine code (this is the same Triangular Number optimization we saw in Section 5.4.3.4).

13.4.1.3. Expected versus worst-case complexity

Another consideration is that there may be one approach which performs better in the majority of cases, at the expense of having very poor performance in specific cases. Sometimes we may risk those high cost cases if we expect the benefit to be worthwhile on the rest of the problem set.

This often happens when the data we are working with has some concept of “distance”. For example, in multi-stop route planning we can use the idea that it's likely to be more efficient to visit nearby destinations first. Generally this works, but sometimes the nearest distance actually has a high cost (such as needing to avoid real-world obstacles in the way which force you to drive past other further away locations to get there).

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13.4.2. Space Complexity

So far we have focused on computational complexity, however similar analysis could be performed for **space complexity**, which is how much computer memory is required to solve a problem. Sometimes, an algorithm will trade computational complexity for space complexity. That is, we might be able to solve a problem much faster if we have more memory available.

For example, there has been research to improve the computational efficiency of matrix multiplication which do indeed run faster than traditional techniques. However, those algorithms don't get implemented on computers because they require way more memory than is available!

13.4.3. Complexity: Takeaways

The idea of algorithmic complexity is important because it grounds us in the harsh truth that some problems are *very* difficult to compute. It's in these cases that a lot of the creativity and domain specific heuristics can become the foremost consideration.

We must remember to be thoughtful about the design of our models. When searching for additional performance to look for the “loops-within-loops” which is where combinatorial explosions tend to happen. Focusing on the places that have large n or poor Big-O order that you can transform the performance of the overall model. Sometimes though, the fundamental complexity of the problem at hand forbids greater efficiency.

13.5. Data Structures

The science of **data structures** is about how data is represented conceptually and in computer applications.

For example, how should the string “abcdef” be represented and analyzed?

There are many common data structures and many specialized subtypes. We will describe some of the most common ones here. Julia has many data structures available in the Base library, but an extensive collection of other data structures can be found in the DataStructures.jl package.

13.5.1. Arrays

An **array** is a contiguous block of memory containing elements of the same type, accessed via integer indices. Arrays have fast random access and are the fastest data structure for linear/iterated access of data.

In Julia, an array is a very common data structure and is implemented with a simple declaration, such as:

```
x = [1,2,3]
```

In memory, the integers are stored as consecutive bits representing the integer values of 1, 2, and 3, and would look like this (with the different integers shown on new lines for clarity):

This is great for accessing the values one-by-one or in consecutive groups, but it's not efficient if values need to be inserted in between. For example, if we wanted to insert 0 between the 1 and 2 in x, then we'd need to overwrite the second position in the array, ask the operating system to allocate more memory⁴, and re-write the bytes that come after our new value. Inserting values at the end (`push!(array, value)`) is usually fast unless more memory needs to be allocated.

13.5.2. Linked Lists

A **linked list** is a chain of nodes where each node contains a value and a pointer to the next node. Linked lists allow for efficient insertion and deletion but slower random access compared to arrays.

In Julia, a simple linked list node could be implemented as:

```
mutable struct Node
    value::Any
    next::Union{Node, Nothing}
end
```

```
z = Node(3, Nothing)
```

⁴In practice, the operating system may have already allocated space for an array that's larger than what the program is actually using so far, so this step may be 'quick' at times, while other times the operating system may actually need to extend the block of memory allocated to the array.

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```
y = Node(2,z)
x = Node(1,y)
```

- ① Nothing represents the end of the linked list.

Inserting a new node between existing nodes is efficient - if we wanted to insert a new node between the ones with value 2 and 3, we could do this:

```
a = Node(0,z)           ①
y.next = a               ②
```

Accessing the nth element requires traversing the list from the beginning to check each Node's next value. This iterative approach makes it $O(n)$ time complexity for random access. This is in contrast to an array where you know right away where the n th item will be in the data structure.

Also, the linked list is one-directional. Items further down the chain don't know what node points to them, so it's impossible to traverse the list backwards.

There are many related implementations which make random access or traversal in reverse order more efficient such as doubly-linked lists or "trees" (Section 13.5.6) which organize the data not as a chain but as a tree with branching nodes.

13.5.3. Records/Structs

An aggregate of named fields, typically of fixed size and sequence. Records group related data together. We've encountered structs in Section 5.4.7, but here we'll add that simple structs with primitive fields can themselves be represented without creating pointers to the data stored:

```
struct SimpleBond
    id::Int
    par::Float64
end

struct LessSimpleBond
    id::String
    par::Float64
end

a = SimpleBond(1, 100.0)
b = LessSimpleBond("1", 100.0)
isbits(a), isbits(b)

(true, false)
```

Because `a` is comprised of simple elements, it can be represented as a contiguous set of bits in memory. It would look something like this in memory:

- ① The bits of 1
 - ② The bits of 100.0

In contrast, the `LessSimpleBond` uses a `String` to represent the ID of the bond. `Strings` are essentially arrays of containers, and the arrays themselves are mutable containers which is by definition not a constant set of bits. In memory, `b` would look like:

- ① a pointer/reference to the array of characters that comprise the string ID
 - ② The bits of 100.0

In performance critical code, having data that is represented with simple bits instead of references/pointers can be much faster (see Chapter 25 for an example).

i Note

For many mutable types, there are immutable, bits-types alternatives. For example:

- Arrays have a `StaticArray` counterpart (from the `StaticArrays.jl` package).
 - Strings have `InlineStrings` (from the `InlineStrings.jl` package) which use fixed-width representations of strings.

The downsides to the immutable alternatives (other than the loss of potentially desired flexibility that mutability provides) are that they can be harder on the compiler (more upfront compilation cost) to handle the specialized cases involved.

13.5.4. Dictionaries (Hash Tables)

13.5.4.1. Hashes and Hash Functions.

Hashes are the result of a **hash function** that maps arbitrary data to a fixed size value. It's sort of a "one way" mapping to a simpler value which has the benefits of:

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1. One way so that if someone knows the hashed value, it's *very* difficult to guess what the original value was. This is most useful in cryptographic and security applications.
2. Creating (probabilistically) unique IDs for a given set of data.

For example, we can calculate a type of hash called an SHA hash on any data:

```
import SHA
let
    a = SHA.sha256("hello world") ▷ bytes2hex
    b = SHA.sha256(rand(UInt8, 10^6)) ▷ bytes2hex
    println(a)
    println(b)
end
```

```
b94d27b9934d3e08a52e52d7da7dabfac484efe37a5380ee9088f7ace2efcde9
e85bbae82c82d096bb8181fd05f1539ddf4bfde22286ac918cde2e4df7de5265
```

We can easily verify that the sha256 hash of "hello world" is the same each time, but it's virtually impossible to guess "hello world" if we are just given the resulting hash. This is the premise of trying to "crack" a password when the stored password hash is stolen.

One way to check if two sets of data are the same is to compute the hash and see if the resulting hashes are equal. For example, maybe you want to see if two data files with different names contain the same data - comparing the hashes is a sure way to determine if they contain the same data.

13.5.4.2. Dictionaries

Dictionaries map a *key* to a *value*. More specifically, they use the *hash of a key* to store a reference to the *value*.

Dictionaries offer constant-time average case access but must handle potential collisions of keys (generally, the more robust the collision handling means higher fixed cost for access).

Here's an illustrative portfolio of assets indexed by CUSIPs:

```
assets = Dict(
    # CUSIP ⇒ Asset
    "037833AH4" ⇒ Bond("General Electric", ...),
    "912828M80" ⇒ Equity("AAPL", ...),
    "594918BQ1" ⇒ Bond("ENRON", ...),
)
```

Then, lookup is performed by indexing the dictionary by the desired key:

```
assets["037833AH4"] # gives the General Electric Bond
```

13.5.5. Graphs

A **graph** is a collection of nodes (also called vertices) connected by *edges* to represent relationships or connections between entities. Graphs are versatile data structures that can model various real-world scenarios such as social networks, transportation systems, or computer networks.

In Julia, a simple graph could be implemented using a dictionary where keys are nodes and values are lists of connected nodes:

```
struct Graph
    nodes::Dict{Any, Vector{Any}}
end

function add_edge!(graph::Graph, node1, node2)
    push!(get!(graph.nodes, node1, []), node2)
    push!(get!(graph.nodes, node2, []), node1)
end

g = Graph(Dict())
add_edge!(g, 1, 2)
add_edge!(g, 2, 3)
add_edge!(g, 1, 3)
```

This implementation represents an undirected graph. For a directed graph, you would only add the edge in one direction.

Graphs can be traversed using various algorithms such as depth-first search (DFS) or breadth-first search (BFS). These traversals are useful for finding paths, detecting cycles, or exploring connected components.

For more advanced graph operations, the `Graphs.jl` package provides a comprehensive set of tools for working with graphs in Julia.

13.5.6. Trees

A tree is a hierarchical data structure with a root node and child subtrees. Each node in a tree can have zero or more child nodes, and every node (except the root) has exactly one parent node. Trees are widely used for representing hierarchical relationships, organizing data for efficient searching and sorting, and in various algorithms.

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A simple binary tree node in Julia could be implemented as:

```
mutable struct TreeNode
    value::Any
    left::Union{TreeNode, Nothing}
    right::Union{TreeNode, Nothing}
end

# Creating a simple binary tree
root = TreeNode(1,
    TreeNode(2,
        TreeNode(4, nothing, nothing),
        TreeNode(5, nothing, nothing)
    ),
    TreeNode(3,
        nothing,
        TreeNode(6, nothing, nothing)
    )
)
```

Trees have various specialized forms, each with its own properties and use cases:

- Binary Search Trees (BST): Each node has at most two children, with all left descendants less than the current node, and all right descendants greater.
- AVL Trees: Self-balancing binary search trees, ensuring that the heights of the two child subtrees of any node differ by at most one.
- B-trees: Generalization of binary search trees, allowing nodes to have more than two children. Commonly used in databases and file systems.
- Trie (Prefix Tree): Used for efficient retrieval of keys in a dataset of strings. Each node represents a common prefix of some keys.

Trees support efficient operations like insertion, deletion, and searching, often with $O(\log n)$ time complexity for balanced trees. They are fundamental in many algorithms and data structures, including heaps, syntax trees in compilers, and decision trees in machine learning.

13.5.7. Data Structures Conclusion

Data structures have strengths and weakness depending on whether you want to prioritize computational efficiency, memory (space) efficiency, code simplicity, and/or mutability. Due to the complexity of real world modeling needs, it can be the case that different representations of the data are more natural or more efficient for the use case at hand.

13.6. Verification and Advanced Testing

13.6.1. Formal Verification

Formal verification is a technique used to prove or disprove the correctness of algorithms with respect to a certain formal specification or property. In essence, it's a mathematical approach to ensuring that a system behaves exactly as intended under all possible conditions.

In formal verification, we use mathematical methods to:

1. Create a formal model of the system
2. Specify the desired properties or behaviors
3. Prove that the model satisfies these properties

This process can be automated using specialized software tools called theorem provers or model checkers.

13.6.1.1. Formal Verification in Practice

It sounds like the perfect risk management and regulatory technique: prove that the system works exactly as intended. However, there has been very limited deployment of formal verification in industry. This is for several reasons:

1. Incomplete Coverage: It's often impractical to formally verify entire large-scale financial systems. Verification, if at all, is typically limited to critical components.
2. Incomplete Specification: Actually reasoning through how the system should behave in all scenarios requires actually contemplating mathematically complete and rigorous possibilities that could occur.
3. Model-Reality Gap: The formal model may not perfectly represent the real-world system, especially in finance where market behavior can be unpredictable.
4. Changing Requirements: Financial regulations and market conditions change rapidly, potentially outdated formal verifications.
5. Performance Trade-offs: Systems designed for easy formal verification might sacrifice performance or flexibility.
6. Cost: The process can be expensive in terms of time and specialized labor.

13.6.2. Property Based Testing

Testing will be discussed in more detail in Chapter 12, but an intermediate concept between Formal Verification and typical software testing is **property-based** testing, which tests for general rules instead of specific examples.

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For example, a function which is associative ($(a + b) + c = a + (b + c)$) or commutative ($a + b = b + a$) can be tested with simple examples like:

```
using Test

myadd(a,b) = a + b

@test myadd(1,2) == myadd(2,1)
@test myadd(myadd(1,2),3) == myadd(1,myadd(2,3))
```

However, we really haven't proven the associative and commutative properties in general. There are techniques to do this, which is a more comprehensive alternative to testing specific examples above. Packages like Supposition.jl provide functionality for this. Note that like Formal Verification, property-based testing is a more advanced topic.

13.6.3. Fuzzing

Fuzzing is kind of like property based testing, but instead of testing general rules, we generalize the simple examples using randomness. For example, we could test the commutative property using random numbers instead, therefore statistically checking that the property holds:

```
@testset for i in 1:10000
    a = rand()
    b = rand()

    @test myadd(a,b) == myadd(b,a)
end
```

This is a good advancement over the simple `@test myadd(1,2) == myadd(2,1)`, in terms of checking the correctness of `myadd`, but it comes at the cost of more computational time and non-deterministic tests.

14. Statistical Inference and Information Theory

"My greatest concern was what to call [the amount of unpredictability in a random outcome]. I thought of calling it 'information,' but the word was overly used, so I decided to call it 'uncertainty.'

When I discussed it with John von Neumann, he had a better idea. Von Neumann told me, 'You should call it entropy, for two reasons. In the first place, your uncertainty function has been used in statistical mechanics under that name, so it already has a name. In the second place, and more important, no one really knows what entropy really is, so in a debate you will always have the advantage.'"- Claude Shannon (1971)

14.1. In This Chapter

A brief introduction to information theory and its foundational role in statistics. Entropy and probability distributions. Bayes' rule and model selection comparison via likelihoods. A brief tour of modern Bayesian statistics.

14.2. Introduction

Statistics has an invaluable role in any data-driven modeling enterprise. As financial professionals dealing inherently with risk and uncertainty - we use probability and statistics to understand, model, and communicate these aspects.

Statistics curricula and practice is undergoing a significant transformation, with a larger focus on information theory and Bayesian methods (as opposed to the common Frequentist methods that have dominated the statistics field for more than a century). Why the change? In short, these methods work better in a wider range of situations and convey more meaningful information about model performance and uncertainty. Until now, computational challenges have limited Bayesian methods to simpler problems, but newer algorithms and better hardware are overcoming this limitation.

14. Statistical Inference and Information Theory

In our experience, it is rare that a financial professional has had exposure to non-Frequentist theory and methods. Given how central probability and statistics is to this endeavor, we have drafted this chapter as a introduction - a proper treatment is beyond the scope of this book. Armed with this knowledge, some of the terminology and tools should be more accessible, and possibly included in new financial models.

14.3. Information Theory

Probability, statistics, machine learning, signal processing, and even physics have a foundational link in **information theory** which is the description and analysis of how much useful data is contained within something. We will work through this with concrete examples.

14.3.1. Example: The Missing Digit

Let's consider the following situation: we are studying a poorly made copy of a financial statement. Amongst many associated exhibits, we are interested in the par value of a particular asset class. Unfortunately, for one reason or another one of the digits is completely indecipherable. Here's what you can read, with the _ indicating that one of the digit is missing from the scanned copy:

32,000,_00

It is likely that you quickly formed an opinion on what the missing number is, but let us make that intuition more formal and quantitative.

Given that we know that par values of assets tend to be nice round numbers, our **prior assumption** for what the probability of the missing digit is may be something like the $p(x_i)$ row of Table 14.1. This prior distribution assumes that the missing digit is most likely a 0. We shall call the individual outcomes x_i and the overall set of probabilities $\{x_0, x_1, \dots, x_9\}$ is called X .

The **information content** of an outcome, $h(x)$ is measured in bits and defined as¹:

$$h(x_i) = \log_2 \frac{1}{p(x_i)} \quad (14.1)$$

¹Log base two turns out to be the most natural representation of information content as it mimics the fundamental 0 or 1 value bit. A more complete introduction is available in "Information Theory, Inference, and Learning Algorithms" by David MacKay.

14.3. Information Theory

Looking at Table 14.1, we can see that the information content of an outcome is *lower* when that outcome has a higher probability than the other potential outcomes. Specifically, If the digit was indeed 0, we have gained less information relative to our expectation than if the missing digit were anything other than 0 .

 Tip

The information content is sometimes referred to as a measure of *surprise* that one would have when observing a realized outcome. In our missing digit example (@tbl-digit-information-human), we would not be surprised at all to find out that the missing digit were 0. In contrast, we would be more surprised to find out the digit were an 8.

We can characterize the entire distribution X via the **entropy**, $H(X)$, of a probability set is the ensemble's average information content:

$$H(X) = \sum p(x_i) \log_2 \frac{1}{p(x_i)} \quad (14.2)$$

The entropy $H(X)$ of the presumed outcomes in Table 14.1 distribution of outcomes is 0.722bits. In just

Table 14.1.: Probability distribution of missing digit, knowing the human inclination to prefer round numbers for par values of assets.

| x_i | 0 | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 |
|----------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| $p(x_i)$ | .91 | .01 | .01 | .01 | .01 | .01 | .01 | .01 | .01 | .01 |
| $h(x_i)$ | 0.136 | 6.644 | 6.644 | 6.644 | 6.644 | 6.644 | 6.644 | 6.644 | 6.644 | 6.644 |

To be clear, we have take a non-uniform view on the probability distribution for the missing digit, and we'll refer to this as the **prior assumption** (or just **prior**). This is unashamedly an opinionated assumption, just like your intuition when you encountered 32,000,-00! All we are doing is giving a quantitative basis for describing this assumption. Taking a view on a prior distribution is a quantitatively incorporating previously encountered data and professional judgment. Having a prior assumption like this is completely compatible with information theory.

Our professional judgment notwithstanding: what if we had another colleague who believed humans are completely rational and without bias for certain numbers? They think an asset's par value need not be rounded at all. They argue for a prior distribution consistent with Table 14.2.

With the uniform prior assumption, $H(X) = 3.322\text{bits}$ and $h(x_i)$ is also uniform. Note that H is higher for the uniform prior than the prior in Table 14.1. We will not prove it

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here, but a uniform probability over a set of outcomes is the highest entropy distribution that can be assumed for this problem. A higher entropy prior distribution can typically be viewed as a less biased prior assumption than a lower entropy prior.

Table 14.2.: Probability distribution of missing digit with uniform, maximal entropy for the assumed probability distribution.

| x_i | 0 | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 |
|----------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| $p(x_i)$ | .10 | .10 | .10 | .10 | .10 | .10 | .10 | .10 | .10 | .10 |
| $h(x_i)$ | 3.322 | 3.322 | 3.322 | 3.322 | 3.322 | 3.322 | 3.322 | 3.322 | 3.322 | 3.322 |

The choice of prior assumption can significantly impact the interpretation and analysis of the missing information. If we have strong reasons to believe that the human bias prior is more appropriate given the context (e.g., knowing that the number is likely a round number), then we would expect the missing digit to be '0' with high probability. However, if we have no specific knowledge about the nature of the number and prefer to make a more conservative assumption, the uniform prior may be more suitable.

In real-world scenarios, the choice of prior assumptions often depends on domain knowledge, available data, and the specific problem at hand. It is important to carefully consider and justify the prior assumptions used in information-theoretic and statistical analyses.

14.3.2. Example: Classification

In this example, we will determine the optimal splits for a decision tree² based on the information gained at each node in the tree.

```
using DataFrames

employed = [true, false, true, true, true, false, true]
good_credit = [true, true, false, true, false, false, true]
default = [true, false, true, true, true, false, true]
default_data = DataFrame(; employed, good_credit, default)
```

The entropy of the default rate data is, per Equation 14.2:

```
H0 = let
    p_default = sum(default_data.default) / nrow(default_data)
    p_good = 1 - p_default
```

²A decision tree is a classification algorithm which attempts to optimally classify an output based on if/else type branches on the input variables.

Table 14.3.: Fictional data regarding loan attributes and whether or not a loan defaulted before its maturity.

| | employed | good_credit | default |
|---|----------|-------------|---------|
| | Bool | Bool | Bool |
| 1 | 1 | 1 | 1 |
| 2 | 0 | 1 | 0 |
| 3 | 1 | 0 | 1 |
| 4 | 1 | 1 | 1 |
| 5 | 1 | 0 | 1 |
| 6 | 0 | 0 | 1 |
| 7 | 0 | 0 | 0 |
| 8 | 1 | 1 | 1 |

```
p_default * log2(1 / p_default) + p_good * log(1 / p_good)
end
```

0.6578517147391054

Our goal is to determine which attribute (`employed` or `good_credit`) to use as the first split in the decision tree. Intuitively, we are looking for the most important factor in predicting default rates. We will quantitatively evaluate this by calculating the information gain, which is the difference in entropy between the prior node and the candidate node. Whichever criteria gains us the most information is the preferred attribute to create a decision split.

In our case we start with H_0 as calculated above for the output variable `default` and calculate the difference in entropy between it and the average entropy of the data if we split on that node. The name for this is the **information gain**, $IG(inputs, attributes)$:

$$IG(T, a) = H(T) - H(T|a)$$

In words, the information gain is simply the difference in entropy before and after learning the value of an outcome a . We will illustrate that by determining the first branch in the decision tree.

Let's first consider splitting the tree based on the `employed` status. We will calculate the entropy of each subset: with employment and without employment.

If we split the data based on being employed, we'd get two sub-datasets:

```
df_employed = filter(:employed => ==(true), default_data)
```

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| | employed | good_credit | default |
|---|----------|-------------|---------|
| | Bool | Bool | Bool |
| 1 | 1 | 1 | 1 |
| 2 | 1 | 0 | 1 |
| 3 | 1 | 1 | 1 |
| 4 | 1 | 0 | 1 |
| 5 | 1 | 1 | 1 |

and

```
df_unemployed = filter(:employed => ==(false), default_data)
```

| | employed | good_credit | default |
|---|----------|-------------|---------|
| | Bool | Bool | Bool |
| 1 | 0 | 1 | 0 |
| 2 | 0 | 0 | 1 |
| 3 | 0 | 0 | 0 |

Let's call it's entropy H_{employed} , which should be zero because there is no variability in the `default` outcome for this subset.

```
H_employed = let
    p_default = sum(df_employed.default) / nrow(df_employed)
    p_good = 1 - p_default
    # p_default * log2(1 / p_default) + p_good * log(1 / p_good)
    p_default * log2(1 / p_default) + 0
end
```

①

- ① In the case of $p_i = 0$ the value of h (the second term in the sum above) is taken to be 0, which is consistent with the $\lim_{p \rightarrow 0^+} p \log(p) = 0$.

0.0

And the corresponding candidate leaf is $H_{\text{unemployed}}$:

```
H_unemployed = let
    p_default = sum(df_unemployed.default) / nrow(df_unemployed)
    p_good = 1 - p_default
    p_default * log2(1 / p_default) + p_good * log(1 / p_good)
end
```

0.7986309056458281

To balance these two results, we weight them according to the amount of data (number of observations) that would fall into each leaf:

```
H1_employment = let
    p_emp = nrow(df_employed) / nrow(default_data)
    p_unemp = 1 - p_emp

    p_emp * H_employed + p_unemp * H_unemployed
end
```

0.29948658961718555

The information gain for splitting the tree using employment status is the difference between the root entropy and the entropy of the employment split:

```
IG_employment = H0 - H1_employment
```

0.35836512512191987

We could repeat the analysis to determine the information gain if we were to split the tree based on having good credit. However, given that there are only two attributes we can already conclude that employed is a better attribute to split the data on. This is because the information gain of `IG_employment` (0.358) is the majority of the overall entropy `H0` (0.658). Entropy is always additive and you cannot have negative entropy, therefore no other other attribute could have greater information gain. This also matches our intuition when looking at Table 14.3 as the eye can spot a higher correlation between `employed` and `default` than `good_credit` and `default`.

The above example demonstrates how we can use information theory to create more optimal inferences on data.

14.3.3. Maximum Entropy Distributions

Why is information theory a useful concept? Many financial models are statistical in nature and concepts of randomness and entropy are foundational. For example, when trying to estimate parameter distributions or assume a distribution for a random process you can lean on information theory to use the most conservative choice: the distribution with the highest entropy given known constraints. These distributions are referred to as **maximum entropy (maxent) distributions**.

In many real-world problems, we seek a distribution that is “least biased” or “most conservative” given certain known information (such as a mean or a range). Maximum entropy distributions accomplish this by spreading out probability as widely as possible under the constraints we know to be true (e.g., average value, bounded domain). They make no additional assumptions beyond those constraints, thereby avoiding unwarranted specificity.

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By using a maxent distribution, we effectively acknowledge that everything else about the system's behavior is unknown and should remain as "random" (unconstrained) as possible. This is a powerful principle because it aligns well with real-world modeling scenarios where we might know just a few key facts—like a process's average rate or finite variance—but have no strong reason to assume anything else about its structure.

Many of the most common probability distributions (Normal, Exponential, Gamma, etc.) can be derived by applying the maximum entropy principle under simple, natural constraints:

- **Normal distribution** when the mean and variance are finite but otherwise unconstrained.

- **Exponential distribution** when we know only that the mean is positive, with outcomes over $([0, \infty))$.

- **Uniform distribution** when outcomes are bounded within a certain interval, and we have no further information about how likely each point is.

Additional maxent distributions and associated constraints are listed in Table 14.4. Those distributions arise again and again in nature because of the second law of thermodynamics - nature likes to have constantly increasing entropy and therefore it should be no surprise (random) processes that maximize entropy pop up all over the place. The **second law of thermodynamics** in physics is an analogy: it states that a closed system tends to move toward higher entropy states. Similarly, in purely probabilistic settings, when few constraints are imposed, the system's "natural" distribution tends to be the one that maximizes entropy.

Maxent distributions have practical modeling use:

- **Conservative Assumptions:** Using a maxent distribution guards against overfitting or adding hidden assumptions. It essentially says: "Given only these constraints, let the data spread out in the most uniform (least structured) way consistent with what I know."
- **Simplicity and Clarity:** It's often easier to justify a maxent model to stakeholders or regulators. If you only know a mean and a variance, a Normal distribution may be the least-biased fit. If you only know a mean and posit that values must be positive, the Exponential distribution is your maxent choice.
- **Built-In Neutrality:** In financial or actuarial contexts, adopting a maxent framework can prevent overly optimistic or pessimistic models. By sticking to the distribution with the fewest assumptions, the risk analysis remains transparent and more robust to model misspecification.

Some discussion of maximum entropy distributions in the context of risk assessment is available in an article by Duracz³.

³https://www.researchgate.net/publication/239752412_Derivation_of_Probability_Distributions_for_Risk_Assessment

Table 14.4.: Maximum Entropy Distributions and the conditions under which they are applicable. For example, if you know that a probability must be continuous and have a positive mean (and can't be normalized), then the MED is the Exponential Distribution.

| Constraint | Discrete Distribution | Continuous Distribution |
|--|-----------------------|-------------------------|
| Bounded range | Uniform (discrete) | Uniform (continuous) |
| Bounded range (0 to 1) with information about the mean or variance | | Beta |
| Mean is finite, two possible values | Binomial | |
| Mean is finite and positive | Geometric | Exponential |
| Mean is finite and range is > zero | | Gamma |
| Mean and Variance is finite | | Gaussian (Normal) |
| Positive and equal mean and variance | Poisson | |

As an example, let's look at processes that behave like the Gaussian (Normal) distribution.

14.3.3.1. Processes that give rise to certain distributions

A random walk can be viewed as the cumulative impact of nudges pushing in opposite directions. This behavior culminates in the random, terminal position being able to be described by a Gaussian distribution. The center of a Gaussian distribution is "thick" because there are many more ways for the cumulative total nudges to mostly cancel out, while it's increasingly rare to end up further and further from the starting point (mean). The distribution then spreads out as flat (randomly) as it can while still maintaining the constraint of having a given, finite variance. Any other continuous distribution that has the same mean and variance has lower entropy than the Gaussian.

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Table 14.5.: Underlying processes create typical probability distributions. That there is significant overlap with the distributions in Section 14.3.3 is not a coincidence.

| Process | Distribution of Data | Examples |
|--|----------------------|--|
| Many <i>additive</i> pluses and minus that move an outcome in one dimension | Normal | Sum of many dice rolls, errors in measurements, sample means (Central Limit Theorem) |
| Many <i>multiplicative</i> pluses and minus that move an outcome in one dimension | Log-normal | Incomes, sizes of cities, stock prices |
| Waiting times between independent events occurring at a constant average rate | Exponential | Time between radioactive decay events, customer arrivals |
| Discrete trials each with the same probability of success, counting the number of successes | Binomial | Coin flips, defective items in a batch |
| Discrete trials each with the same probability of success, counting the number of trials until the first success | Geometric | Number of job applications until getting hired |
| Continuous trials each with the same probability of success, measuring the time until the first success | Exponential | Time until a component fails, time until a sales call results in a sale |
| Waiting time until the r-th event occurs in a Poisson process | Gamma | Time until the 3rd customer arrives, time until the 5th defect occurs |

💡 Probability Distributions

There are a *lot* of specialized distributions. There are lists of distributions you can find online or in references such as Leemis and McQueston (2008) which has a full-page network diagram of the relationships.

The information-theoretic and Bayesian perspective on it is to eschew memorization of a bunch of special cases and statistical tests. If you pull up the aforementioned diagram in Leemis and McQueston (2008), you can see just a handful of

distributions that have the most central roles in the universe of distributions. Many distributions are simply transformations, limiting instances, or otherwise special cases of a more fundamental distribution. Instead of trying to memorize a bunch of probability distributions, it's better to think critically about:

1. The fundamental processes that give rise to the randomness we are interested in modeling.
2. Transformations of the data to make it nicer to work with, such as translations, scaling, or other non-destructive changes.

Then when you encounter an unusual dataset, you don't need to comb the depths of Wikipedia to find the perfect probability distribution for that situation.

14.3.3.2. Additive and Multiplicative Processes

Table 14.5 describes some examples, let us discuss further what it means to have a process that arises via an additive vs multiplicative effect⁴. Additive processes result in a normal distribution while multiplicative processes give rise to a log-normal distribution⁵.

An outcome is additive and results in a normal distribution if it's the sum or difference of multiple independent processes. Examples of this include:

- Rolling multiple dice and taking their sum.
- A random walk along the natural numbers wherein with equal probability you take a step left or right.
- Calculating the arithmetic mean of samples (the Central Limit Theorem).

However, many processes are multiplicative in nature. For example the population density of cities is distributed in a log-normal fashion. If we think about the factors that contribute to choice of place to live, we can see how these factors multiply: an attractive city might make someone 10% more likely to move, a city with water features 15% more likely, high crime 30% less likely, etc. These forces combine in a multiplicative way in the generative process of deciding where to move. In finance, many price processes are considered multiplicative.

⁴Multiplicative process are often referred to as “geometric”, as in “geometric Brownian motion” or “geometric mean”. Additive processes are sometimes referred to as “arithmetic”. This root of this confusing terminology appears to be due to the fact that series involving repeated multiplication were solved via geometric (triangles, angles, etc.) methods while those using sums and differences were solved via arithmetic.

⁵See Tip 1.

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💡 Tip 1: Logarithms

The logarithm of a geometric process transforms the outcomes into “log-space”. The information is the same, but is often a more convenient form for the analysis. That is, if:

$$Y = x_1 \times x_2 \times \dots \times x_i$$

Then,

$$\log(Y) = \log(x_1) + \log(x_2) + \dots + \log(x_i)$$

This is effectively the transformation that gives rise to the Normal versus Log-Normal distribution.

In the context of computational thinking:

First, we should think about how to transform data or modeling outcomes into a more convenient format. The log transform doesn't eliminate any information but may map the information into a shape that is easier for an optimizer or Monte Carlo simulation to explore.

Second, per Chapter 5, floating point math is a *lossy* transformation of real numbers into a digital computer representation. Some information (in the literal Shannon information sense) is lost when computing and this tends to be worst with very small real numbers, such as those we encounter frequently in probabilities and likelihoods. Logarithms map very small numbers into negative numbers that don't encounter the same degree of truncation error that tiny numbers do.

Third, modern CPUs are generally much faster at adding or subtracting numbers than multiplying or dividing. Therefore working with the logarithm of processes may be computationally faster than the direct process itself.

14.4. Bayes' Rule

With some of the foundational concepts laid down, we now turn to the perpetual challenge of attempting to make inferences and predictions given a set of data. We covered basic information theory, probability distributions, log transformations, and random processes because modern statistical analysis relies heavily on those concepts and techniques. We'll introduce Bayes' Rule, but then analysis beyond trivial applications one will typically quickly encounter challenges posed by those ideas.

The remainder of this chapter will re-introduce Bayes' Rule and then build up modeling applications that illustrate some core concepts of applying Bayes Rule to complex, data-intensive problems.

14.4.1. Bayes' Rule Formula

The minister and statistician Thomas Bayes derived a relationship of conditional probabilities that we today know as **Bayes' Rule**. Laplace⁶ furthered the notion, and developed the modern formulation, commonly written as:

$$P(H|D) = \frac{P(D|H) \times P(H)}{P(D)}$$

The components of this are:

- $P(H | D)$ is the conditional probability of event H occurring given that D is true.
- $P(D | H)$ is the conditional probability of event D occurring given that H is true.
- $P(H)$ is the prior probability of event H .
- $P(D)$ is the prior probability of event D .

If we take the following:

- D is the available data
- H is our hypothesis

Then we can draw conclusions about the probability of a hypothesis being true given the observed data. When thought about this way, Bayes' rule is often described as:

$$\text{posterior} = \frac{\text{likelihood} \times \text{prior}}{\text{evidence}}$$

This is a very useful framework, which we'll return to more completely in Section 14.5. First, let's look at combining information theory and Bayes' rule in an applied example.

14.4.2. Model Selection via Likelihoods

Let's say that we have competing hypothesis about a data generating process, such as: "given a set of data representing risk outcomes, what distribution best fits the data"? We will not be able to determine an *absolute* probability of a model given the data, but amazingly we can determine *relative* probability of models given the data. This is powerful because often times one of the most difficult modeling tasks is to select a model formulation - and Bayes gives us a powerful tool to help choose.

⁶Laplace actually deserves most of the credit, as it was he who formalized the modern notion of Bayes' rule and cemented the mathematical formulation. Bayes just described it first, in a way that actually had almost no direct impact on math or science. See "The Theory That Would Not Die".

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We can compare these models using Bayes' rules by observing the following: Suppose we have two models, H_1 and H_2 , and we want to compare their likelihoods given the observed data, D . We can use Bayes' rule to calculate the posterior probability of each model:

$$P(H_1|D) = (P(D|H_1) \times P(H_1))/P(D)$$

$$P(H_2|D) = (P(D|H_2) \times P(H_2))/P(D)$$

Where:

- $P(H_1|D)$ and $P(H_2|D)$ are the posterior probabilities of models H_1 and H_2 , respectively, given the data D .
- $P(D|H_1)$ and $P(D|H_2)$ are the likelihoods of the data D under models H_1 and H_2 , respectively.
- $P(H_1)$ and $P(H_2)$ are the prior probabilities of models H_1 and H_2 , respectively.
- $P(D)$ is the marginal likelihood of the data, which serves as a normalizing constant.

To compare the likelihoods of the two models, we can calculate the ratio of their posterior probabilities, known as the **Bayes factor**, BF :

$$BF = \frac{P(H_1|D)}{P(H_2|D)}$$

Substituting the expressions for the posterior probabilities from Bayes' rule, we get:

$$BF = \frac{P(D|H_1) \times P(H_1)}{P(D|H_2) \times P(H_2)}$$

The marginal likelihood $P(D)$ cancels out since it appears in both the numerator and denominator. If we assume equal prior probabilities for the models, i.e., $P(H_1) = P(H_2)$, then the Bayes factor simplifies to the likelihood ratio:

$$BF = \frac{P(D|H_1)}{P(D|H_2)}$$

The Bayes factor then is a statement about the relative probability of two competing models for the given data. We can interpret the results as:

- If $BF > 1$, the data favors model H_1 over model H_2 .
- If $BF < 1$, the data favors model H_2 over model H_1 .

- If $BF = 1$, the data do not provide evidence in favor of either model.

In practice, the likelihoods $P(D|H_1)$ and $P(D|H_2)$ are often calculated using the probability density or mass functions of the models, evaluated at the observed data points. The prior probabilities $P(H_1)$ and $P(H_2)$ can be assigned based on prior knowledge or assumptions about the models. By comparing the likelihoods of the models using the Bayes factor, we can quantify the relative support for each model given the observed data, while taking into account the prior probabilities of the models.

Another way of interpreting this is the evaluation of which model has the higher likelihood given the data.

Null Hypothesis Statistical Test

Null Hypothesis Statistical Tests (NHST) is the idea of trying to statistically support an alternative hypothesis over a null hypothesis. The support in favor of alternative versus the null is reported via some statistical power, such as the **p-value** (the probability that the test result is as, or more extreme, than the value computed). The idea is that there's some objective way to push science towards greater truths and NHST was seen as a methodology that avoided the subjectivity of the Bayesian approach. However, while pure in intention, the NHST choices of both null hypothesis and model contain significant amounts of subjectivity! There is subjectivity in the null hypothesis, data collection methodologies, study design, handling of missing data, choice of data *not* to include, which statistical tests to perform, and interpretation of relationships.

We might as well call the null hypothesis a prior and stop trying to disprove it absolutely. Instead: focus on model comparison, model structure, and posterior probabilities of the competing theories.

Over 100 statistical tests have been developed in service of NHST Lewis (2013), but it's widely viewed now that a focus on NHST has led to *worse* science due to a multitude of factors, such as:

- “P-hacking” or trying to find subsets of data which can (often only by chance) support rejecting some null.
- Cognitive anchoring to the importance of a p-value of 0.05 or less – why choose that number versus 0.01 or 0.001 or 0.49?
- Bias in research processes where one may stop data collection or experimentation after achieving a favorable test result.
- Inappropriate application of the myriad of statistical tests.
- Focus on p-values rather than effects that simply matter more or have greater effect.
 - For example, which is of more interest to doctors? A study indicating a 1 in a billion chance of serious side effect , with p-value of 0.0001 or

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a study indicating a 1 in 3 chance with p-value 0.06? Many journals would only publish the former study, even though the latter study intuitively suggests a potentially more risky drug.

- Difficulty to determine *causal* relationships.

The authors of this book recommend against basic NHST and memorization of statistical tests in favor of principled Bayesian approaches. For the actuarial readers, NHST is analogous to traditional credibility methods (of which the authors also prefer more modern statistical approaches).

14.4.2.1. Example: Rainfall Risk Model Comparison

The example we'll look at relates to the annual rainfall totals for a specific location in California⁷, which could be useful for insuring flood risk or determining the value of a catastrophe bond. Acknowledging that we are attempting to create a geocentric model⁸ instead of a scientifically accurate weather model, we narrow the problem to finding a probability distribution that matches the historical rainfall totals.

Our goal is to recommend a model that best fits the data and justify that recommendation quantitatively. Before even looking at the data, Table 14.6 shows three competing models based on thinking about the real-world outcome we are trying to model. These three are chosen for the increasingly sophisticated thought process that might lead the modeler to recommend them - but which is supportable by the statistics?

Table 14.6.: Three alternative hypothesis about the distribution of annual rainfall totals.

| Hypothesis | Process | Possible Rationale |
|------------|--|---|
| H_1 | A Normal
(Gaussian)
distribution | The sum of independent rainstorms creates annual rainfall totals that are normally distributed |
| H_2 | A LogNormal
distribution | Since it's normal-ish, but skewed and can't be negative |
| H_3 | A Gamma
Distribution | Since rainfall totals would be the sum of exponentially-distributed independent rainfall events |

⁷<https://data.ca.gov/dataset/annual-precipitation-data-for-northern-california-1944-current>

⁸See @sec-predictive-vs-explanatory.

i Note

In the literature for rainfall modeling, H_3 (the Gamma distribution) is known as the “Log-Pearson Type III distribution”. It’s actually recommended by the US Corps of Army Engineers as the recommended way to model rainfall totals. We are able to avoid learning and memorizing specialty distributions and statistical tests, which are so common in Frequentist approaches. First-principles reasoning on the probabilistic processes can get one to a reasonable hypothesis, comparable to ‘specialist’ knowledge one would encounter in the literature for a particular applied field.

Here’s the data:

```
rain = [
    39.51, 42.65, 44.09, 41.92, 28.42, 58.65, 30.18, 64.4, 29.02,
    37.00, 32.17, 36.37, 47.55, 27.71, 58.26, 36.55, 49.57, 39.84,
    82.22, 47.58, 51.18, 32.28, 52.48, 65.24, 51.12, 25.03, 23.27,
    26.11, 47.3, 31.8, 61.45, 94.95, 34.8, 49.53, 28.65, 35.3, 34.8,
    27.45, 20.7, 36.99, 60.54, 22.5, 64.85, 43.1, 37.55, 82.05, 27.9,
    36.55, 28.7, 29.25, 42.32, 31.93, 41.8, 55.9, 20.65, 29.28, 18.4,
    39.31, 20.36, 22.73, 12.75, 23.35, 29.59, 44.47, 20.06, 46.48,
    13.46, 9.34, 16.51, 48.24
];
```

Plotted, we see some of the characteristics that align with our prior assumptions and knowledge about the system itself, such as: the data being constrained to positive values and a skew towards having some extreme weather years with lots of rainfall.

```
using CairoMakie
hist(rain)
```

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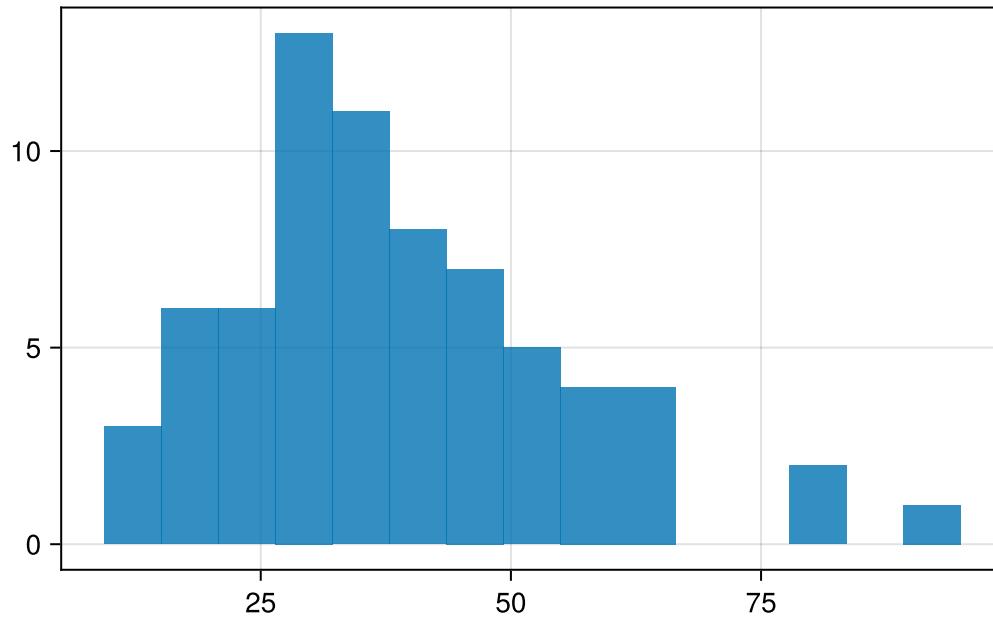


Figure 14.1.: Annual rainfall totals for a specific location in California.

We will show the likelihood of the three models after deriving the **maximum likelihood (MLE)**, which is simply finding the parameters that maximize the calculated likelihood. In general, this can be accomplished by an optimization routine, but here we will just use the functions built into Distributions.jl:

```
using StatsBase
using Distributions

n = fit_mle(Normal, rain)
ln = fit_mle(Normal, log.(rain))
lg = fit_mle(Gamma, log.(rain))
@show n
@show ln
@show lg;

n = Distributions.Normal{Float64}(\mu=38.91442857142857, σ=16.643603630714306)
ln = Distributions.Normal{Float64}(\mu=3.5690550009062663, σ=0.44148379736539156)
lg = Distributions.Gamma{Float64}(α=61.58531301458412, θ=0.05795302201453571)
```

Let's look at the likelihoods by applying the maximum likelihood distribution to the observed data. For the practical reasons described in Tip 1, we will compare the the log-

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likelihoods to maintain convention with what you'd likely see or deal with in practice. Taking the log of the likelihood does not change the ranking of the likelihoods.

```
let
    n_lik = sum(log.(pdf.(n, rain)))
    ln_lik = sum(log.(pdf.(ln, log.(rain))))
    lg_lik = sum(log.(pdf.(lg, log.(rain))))

    @show n_lik
    @show ln_lik
    @show lg_lik
end;

n_lik = -296.16751566478115
ln_lik = -42.09272021737913
lg_lik = -43.79151806348801
```

The results indicate that the LogNormal and the Gamma model for rainfall distribution are very superior to the Normal model, consistent with the visual inspection of the quantiles in Figure 14.2. We reach that conclusion by noting how much more likely the latter two are, as the likelihoods of -42 and -44 is much greater than -296 ⁹.

```
let x = rain

    range = 1:0.1:100
    fig, ax, _ = lines(range, cdf.(n, range), label="Normal", axis=(xgridvisible=false, ygridvisible=false))
    lines!(ax, range, cdf.(ln, log.(range)), label="LogNormal")
    lines!(range, cdf.(lg, log.(range)), label="LogGamma")
    lines!(quantile.(Ref(x), 0.01:0.01:0.99), 0.01:0.01:0.99, label="Data", color=:black, 0.6)
    fig[1, 2] = Legend(fig, ax, "Model", framevisible=false)
    fig
end
```

⁹The values are negative because we are taking the logarithm of a number less than 1. The likelihoods are less than 1 because the likelihood is the joint (multiplicative) probability of observing each of the individual outcomes.

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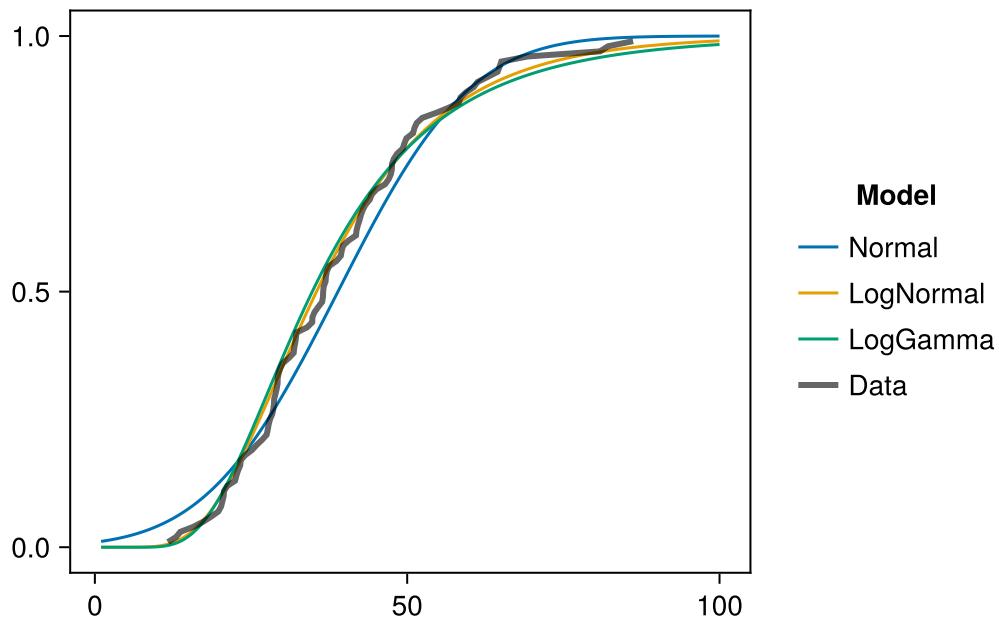


Figure 14.2.

We evaluated the likelihood at a single point estimate of the parameters, but a true posterior probability of the parameters of the distributions will be represented by a *distribution* rather than a point. The rest of chapter will describe how to express the posterior probabilities of the the parameters for H_1 , H_2 , and H_3 using Bayesian statistical methods.

14.5. Modern Bayesian Statistics

14.5.1. Background

Bayesian statistics is generally *not* taught in undergraduate statistics. Bayes' rule is introduced, basic probability exercises are assigned, and then statistics moves on to a curriculum of regression and NHSTs (of the Frequentist school). Why is the applied practice of statistics currently gravitating towards Bayesian approaches? There are both philosophical and practical reasons why.

14.5.1.1. Philosophical Motivations

Philosophically, one of the main reasons why Bayesian thinking is appealing is its ability to provide straightforward interpretations of statistical conclusions.

For example, when estimating an unknown quantity, a Bayesian probability interval can be directly understood as having a high probability of containing that quantity. In contrast, a Frequentist confidence interval is typically interpreted only in the context of a series of similar inferences that could be made in repeated practice. In recent years, there has been a growing emphasis on interval estimation rather than hypothesis testing in applied statistics. This shift has strengthened the Bayesian perspective since it is likely that many users of standard confidence intervals intuitively interpret them in a manner consistent with Bayesian thinking.

i “The fallacy of placing confidence in confidence intervals”

“The fallacy of placing confidence in confidence intervals” is the title of an article ([Morey_Hoekstra_Rouder_Lee_Wagenmakers_2016?](#)) describing the issues with confidence intervals, with one of the primary issues being what they call the “Fundamental Confidence Fallacy,” or FCF.

The FCF is the belief that if you have a $X\%$ confidence interval, that the probability that the true value of interest lies within that interval is $X\%$. This is false, and an abbreviated explanation of why is as follows:

- An $X\%$ confidence interval refers to a procedure that results in a range R wherein $X\%$ of the time the true value of interest lies within R .
- However, once the data is observed we can know with higher probability than X whether the true value lies within that range.

Here’s an example taken from the article referenced above: we wish to estimate the mean of a continuous random variable. We observe two data points y_1 and y_2 . If $y_1 < y_2$ then we say our interval is $(-\infty, \infty)$ otherwise the interval is empty. Our credibility procedure creates an interval for which 50% of the time it contains the true value. However, once we’ve observed the data and created the interval, then “post-data” we can tell with certainty whether our interval contains the variable of interest. We can only say “pre-data”, or pre-observations, that the credibility interval is really $X\%$ probable to contain the right value. Similar (and other) issues arises with more “real world” examples of confidence intervals.

In contrast, the Bayesian procedure of estimating an interval using the posterior distribution of the data *can* be interpreted as an interval for which you can say that you believe the true value is contained within the interval. Typically, this is referred to as a *credibility interval* to distinguish from a *confidence interval*.

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Another meaningful way to understand the contrast between Bayesian and Frequentist approaches is through the lens of decision theory, specifically how each view treats the concept of randomness. This perspective pertains to whether you regard the data being random or the parameters being random.

Frequentist statistics treats parameters as fixed and unknown, and the data as random — that data you collect is but one realization of an infinitely repeatable random process. Consequently, Frequentist procedures, like hypothesis testing or confidence intervals, are generally based on the idea of long-run frequency or repeatable sampling.

Conversely, Bayesian statistics turns this on its head by treating the data as fixed — after all, once you've collected your data, it's no longer random but a fixed observed quantity. Parameters, which are unknown, are treated as random variables. The Bayesian approach then allows us to use probability to quantify our uncertainty about these parameters.

The Bayesian approach tends to align more closely with our intuitive way of reasoning about problems. Often, you are given specific data and you want to understand what that particular set of data tells you about the world. You're likely less interested in what might happen if you had infinite data, but rather in drawing the best conclusions you can from the data you do have.

14.5.1.2. Practical Motivations

Practically, recent advances in computational power, algorithm development, and open-source libraries have enabled practitioners to adapt the Bayesian workflow.

For most real-world problems, deriving the posterior distribution is analytically intractable and computational methods must be used. Advances in raw computing power only in the 1990's made non-trivial Bayesian analysis possible, and recent advances in algorithms have made the computations more efficient. For example, one of the most popular algorithms, NUTS, was only published in the 2010's.

Many problems require the use of compute clusters to manage runtime, but if there is any place to invest in understanding posterior probability distributions, it's financial companies trying to manage risk!

The availability of open-source libraries, such as Turing.jl, PyMC3, and Stan provide access to the core routines in an accessible interface. To get the most out of these tools requires the mindset of computational thinking described in this book - understanding model complexity, model transformations and structure, data types and program organization, etc.

14.5.1.3. Advantages of the Bayesian Approach

The main advantages of this approach over traditional actuarial techniques are:

1. **Focus on distributions rather than point estimates of the posterior's mean or mode.** We are often interested in the distribution of the parameters and a focus on a single parameter estimate will understate the risk distribution.
2. **Model flexibility.** A Bayesian model can be as simple as an ordinary linear regression, but as complex as modeling a full insurance mechanics.
3. **Simpler mental model.** Fundamentally, Bayes' theorem could be distilled down to an approach where you count the ways that things could occur and update the probabilities accordingly.
4. **Explicit Assumptions.**: Enumerating the random variables in your model and explicitly parameterizing prior assumptions avoids ambiguity of the assumptions inside the statistical model.

14.5.1.4. Challenges with the Bayesian Approach

With the Bayesian approach, there are a handful of things that are challenging. Many of the listed items are not unique to the Bayesian approach, but there are different facets of the issues that arise.

1. **Model Construction.** One must be thoughtful about the model and how variables interact. However, with the flexibility of modeling, you can apply (actuarial) science to make better models!
2. **Model Diagnostics.** Instead of R^2 values, there are unique diagnostics that one must monitor to ensure that the posterior sampling worked as intended.
3. **Model Complexity and Size of Data.** The sampling algorithms are computationally intensive - as the amount of data grows and model complexity grows, the runtime demands cluster computing.
4. **Model Representation.** The statistical derivation of the posterior can only reflect the complexity of the world as defined by your model. A Bayesian model won't automatically infer all possible real-world relationships and constraints.

i Subjectivity of the Priors?

There are two ways one might react to subjectivity in a Bayesian context: It's a feature that should be embraced or it's a flaw that should be avoided.

14.5.1.5. Subjectivity as a Feature

A Bayesian approach to defining a statistical model is an approach that allows for explicitly incorporating professional judgment. Encoding assumptions into

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a Bayesian model forces the actuary to be explicit about otherwise fuzzy predilections. The explicit assumption is also more amenable to productive debate about its merits and biases than an implicit judgmental override.

14.5.1.6. Subjectivity as a Flaw

Subjectivity is inherent in all useful statistical methods. Subjectivity in traditional approaches include how the data was collected, which hypothesis to test, what significant levels to use, and assumptions about the data-generating processes. In fact, the “objective” approach to null hypothesis testing is so prone to abuse and misinterpretation that in 2016, the American Statistical Association issued a statement intended to steer statistical analysis into a “post $p < 0.05$ era.” That “ $p < 0.05$ ” approach is embedded in most traditional approaches to actuarial credibility¹⁰ and therefore should be similarly reconsidered.

14.5.2. Implications for Financial Modeling

Like Bayes’ Formula itself, another aspect of financial literature that is taught but often glossed over in practice is the difference between process risk (volatility), parameter risk, and model formulation risk. When performing analysis that relies on stochastic results, in practice typically only process/volatility risk is assessed.

Bayesian statistics provides the tools to help financial modelers address parameter risk and model formulation. The posterior distribution of parameters derived is consistent with the observed data and modeled relationships. This posterior distribution of parameters can then be run as an additional dimension to the risk analysis.

Additionally, best practices include skepticism of the model construction itself, and testing different formulation of the modeled relationships and variable combinations to identify models which are best fit for purpose. Tools such as Information Criterion, posterior predictive checks, Bayes factors, and other statistical diagnostics can inform the actuary about trade-offs between different choices of model.

i Bayesian Versus Machine Learning

Machine learning (ML) is *fully compatible* with Bayesian analysis - one can derive posterior distributions for the ML parameters like any other statistical model and the combination of approaches may be fruitful in practice.

However, to the extent that actuaries have leaned on ML approaches due to the shortcomings of traditional actuarial approaches, Bayesian modeling may provide

¹⁰Note that the approach discussed here is much more encompassing than the Bühlmann-Straub Bayesian approach described in the actuarial literature.

an attractive alternative without resorting to notoriously finicky and difficult-to-explain ML models. The Bayesian framework provides an explainable model and offers several analytic extensions beyond the scope of this introductory chapter:

- Causal Modeling: Identifying not just correlated relationships, but causal ones, in contexts where a traditional designed experiment is unavailable.
- Bayes Action: Optimizing a parameter for, e.g., a CTE95 level instead of a parameter mean.
- Information Criterion: Principled techniques to compare model fit and complexity.
- Missing data: Mechanisms to handle the different kinds of missing data.
- Model averaging: Posteriors can be combined from different models to synthesize different approaches.
- Credibility Intervals: A posterior representation around the likely range of values for parameters of interest.

14.5.3. Basics of Bayesian Modeling

A Bayesian statistical model has four main components to focus on:

1. **Prior** encoding assumptions about the random variables related to the problem at hand, before conditioning on the data.
2. A **Model** that defines how the random variables give rise to the observed outcome.
3. **Data** that we use to update our prior assumptions.
4. **Posterior** distributions of our random variables, conditioned on the observed data and our model

While this is simply stating Bayes' formula in words, it's also the blueprint for a workflow to implement more advanced Bayesian methods.

Having defined a prior assumption, selected a model, and collected our data, the computation of the posterior can be the most challenging. The workflow involves computationally sampling the posterior distribution, often using a technique called **Markov Chain Monte-Carlo** (MCMC). The result is a series of values that are sampled statistically from the posterior distribution. Introducing this process is the focus of the rest of this chapter.

14.5.4. Markov-Chain Monte Carlo

Computing the posterior distribution for most model parameters is analytically intractable. However, we can probabilistically sample from the posterior distribution and achieve an approximation of the posterior distribution. MCMC samplers, as they are

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called, do this by moving through the parameter space (the set of possible values for the parameters) in a special way. The statistical marvel is that they travel to different points *in proportion* to the posterior probability. It is a “Markov-Chain” because the probability of the next point’s location is influenced by the prior sampling point’s location.

14.5.4.1. Example: MCMC from Scratch

Here is a simple example demonstrated with one of the oldest MCMC algorithm, called Metropolis-Hastings. The general idea is this:

1. Start at an arbitrary point and make that the `current_state`.
2. Propose a new point which is the `current_state` plus some movement that comes from a random distribution, `proposal_dist`.
3. Calculate the likelihood ratio of the proposed versus current point (`acceptance_ratio` below).
4. Draw a random number - if that random number is less than the `acceptance_ratio`, then move to that new point. Otherwise do not move.
5. Repeat steps 2-4 until the distribution of points converges to a stable posterior distribution.

This gets us what we desire because the resulting distribution of samples has frequency that’s proportional to the posterior distribution.

We will try to find the posterior of an arbitrary set of normally distributed asset returns. We set the true, (unobserved in reality) values for μ and σ and then draw 250 generated observations:

```
# In reality, we don't observe the parameters
# we are interested in determining the values for.
σ = 0.15
μ = 0.1

n_observations = 250
return_dist = Normal(μ,σ)
returns = rand(return_dist,n_observations)

# plot the distribution of returns
μ_range = LinRange(-0.5, 0.5, 400)
σ_range = LinRange(0.0, 3.0, 400)

f = Figure()
ax1 = Axis(f[1,1],title="True Distribution of Returns")
```

```

ax2 = Axis(f[2,1],title="Simulated Outcomes", xlabel="Return")
plot!(ax1,return_dist)
vlines!(ax1,[μ],color=(:black,0.7))
text!(ax1,μ,0;text="mean ($μ)",rotation=pi/2)
hist!(ax2,returns)
vlines!(ax2,[mean(returns)],color=(:black,0.7))
text!(ax2,mean(returns),0;text="mean ($round(mean(returns);digits=3))",rotation=pi/2)

linkxaxes!(ax1,ax2)

```

f

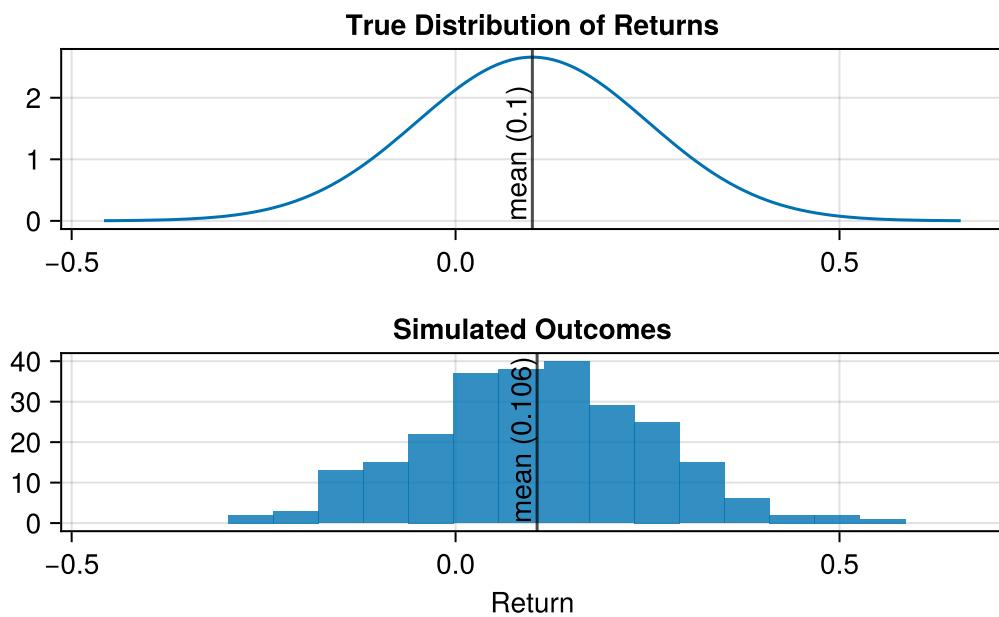


Figure 14.3.: The target probability densities which we will attempt to infer via MCMC.

Having generated sample data, we will next define a probability distribution for the random step that we take from the `current_point` on the Markov chain. We choose a 2D Gaussian for this, since the parameter space to explore is two-dimensional (μ and θ). The `proposal_std` controls how big of a movement is taken at each step.

```

# Define the proposal step distribution
proposal_std = 0.05
proposal_dist = Normal(0,proposal_std)

Distributions.Normal{Float64}({μ=0.0, σ=0.05})

```

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We next define how many steps we want the chain to sample for, and implement the algorithm's main loop containing the logic in the steps above.

```
# MCMC parameters
num_samples = 5000
burn_in = 500

# Define priors
μ_prior = Normal(0, 0.25)
σ_prior = Gamma(0.5)

# Initialize the Markov chain
μ_current, σ_current = 0.0, 0.25
current_prob = sum(logpdf(Normal(μ_current, σ_current), r) for r in returns) +
    logpdf(μ_prior, μ_current) +
    logpdf(σ_prior, σ_current)

chain = zeros(num_samples, 2)

count = 0

# MCMC sampling loop
while count < num_samples

    # Generate a new proposal
    ḡ, ḡ = μ_current + rand(proposal_dist), σ_current + rand(proposal_dist)
    if ḡ > 0

        # Calculate the acceptance ratio

        proposal_prob = sum(logpdf(Normal(ḡ, ḡ), r) for r in returns) +
            logpdf(μ_prior, ḡ) + logpdf(σ_prior, ḡ)
        log_acceptance_ratio = proposal_prob - current_prob

        # Accept or reject the proposal
        if log(rand()) < log_acceptance_ratio
            μ_current, σ_current = ḡ, ḡ
            current_prob = proposal_prob
        end

        # Store the current state as a sample
        count += 1
        chain[count, :] .= μ_current, σ_current
    else

```

```

# skip because σ can't be negative
end
end

chain

5000×2 Matrix{Float64}:
0.0      0.25
0.0      0.25
0.0477587  0.240734
0.0399676  0.183676
0.0399676  0.183676
0.0594153  0.14767
0.0594153  0.14767
0.0594153  0.14767
0.0594153  0.14767
0.0594153  0.14767
⋮
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774
0.108293  0.142774

```

The resulting chain contains a list of points that the algorithm has moved along during the sampling process. Note that there is a burn-in parameter. This is because we want the chain iterate long enough to be effectively independent of both (1) the starting point for the sample, and (2) that different chains are effectively independent.

After having performed the sampling, we can now visualize the chain versus the target_distribution. A few things to note:

1. The red line indicates the “warm up” or “burn-in” phase and we do not consider that as part of the sampled chain because those values are too correlated with the arbitrary starting point.
2. The blue line indicates the path traveled by the Metropolis-Hastings algorithm. Long asides into low-probability regions are possible, but in general the path will traverse areas in proportion to the probability of interest.

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```
# Plot the chain
let
    f = Figure()

    # μ lines
    ax1 = Axis(f[1, 1], ylabel="σ", xticklabelsvisible=false)

    # burn in lines
    scatterlines!(ax1, chain[1:burn_in, 1],
                  chain[1:burn_in, 2],
                  color=(:red, 0.1),
                  markercolor=(:red, 0.1))

    # sampled lines
    scatterlines!(ax1, chain[burn_in+1:end, 1],
                  chain[burn_in+1:end, 2],
                  color=(:blue, 0.1),
                  markercolor=(:blue, 0.1))

    # μ histogram
    ax2 = Axis(f[2, 1], xlabel="μ", yticklabelsvisible=false)
    hist!(ax2, chain[burn_in+1:end, 1], color=:blue)
    linkxaxes!(ax1, ax2)

    # σ histogram
    ax3 = Axis(f[1, 2], xticklabelsvisible=false)
    hist!(ax3, chain[burn_in+1:end, 2], color=:blue, direction=:x)
    linkyaxes!(ax1, ax3)

f
end
```

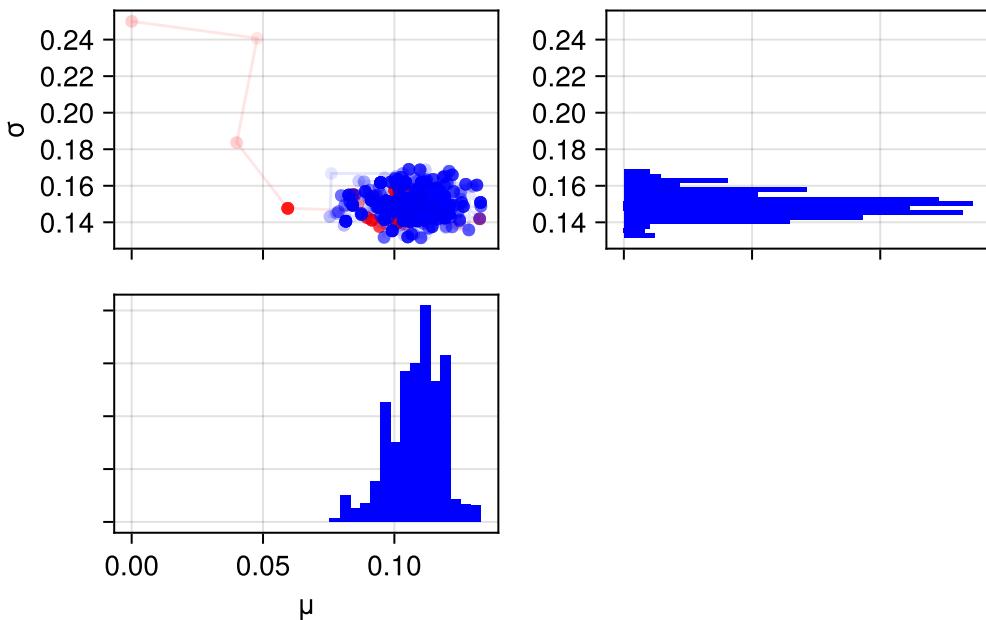


Figure 14.4.: The blue lines of the MCMC chain explore the posterior density of interest (after discarding the burn-in samples in red). Note that locations where the sampler remained longer (rejected more proposals) show up as darker points.

In this example, μ and σ are independent, but if there were a correlation (such as when μ were higher, σ were also higher) then the sampler would pick up on this, and we would see a skew in the plotted chain.

The point of this short, ground-up introduction to MCMC is that the technique is not magic by demonstrating that we could do-it-from-scratch with small amounts of code. The challenge is that it's just computationally intensive. Modern libraries perform the sampling for you with more advanced algorithms than Metropolis-Hastings.

14.5.5. MCMC Algorithms

The Metropolis-Hastings algorithm is simple, but somewhat inefficient. Some challenges with MCMC sampling are both mathematical and computational:

1. Often times the algorithm will back-track (take a “U-Turn”), wasting steps in regions already explored.

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2. The algorithm can have a very high rate of rejecting proposals if the proposal mechanism generates steps that would move the current state into a low-probability regions.
3. The choice of proposal distribution and parameters can greatly influence the speed of convergence. Too large of movement and key regions can be entirely skipped over, while small movements can take much longer than necessary to explore the space.
4. As the number of parameters grows, the dimensionality of the parameter space to explore also grows making posterior exploration much harder.
5. The shape of the posterior space can be more or less difficult to explore. Complex models may have regions of density that are not nicely “round” - regions may be curved, donut shaped, or disjointed.

The issues above mean that MCMC sampling is very computationally expensive for more complex examples. Compared with Metropolis-Hastings, modern algorithms (such as the No-U-Turn (NUTS)) algorithm explore the posterior distribution more efficiently by avoiding back-tracking to already explored regions and dynamically adjusting the proposals to adaptively fit the posterior. Many of them take direct influence from particle physics, with the algorithm keeping track of the energy of the current state as it explores the posterior space.

Algorithms have only brought so much relief to the modeler with finite resources and compute. There is still a lot of responsibility for modeler to design models that are computationally efficient, transformed to eliminate oddly-shaped density regions, or find the right simplifications to the analysis in order to make the problem tractable.

i Note

What does it mean to transform the parameter space?

An example will be shown in Chapter 30 where we want to ensure that a binomial variable is constrained to the region $[0, 1]$ but the underlying factors are allowed to vary across the entire real numbers. We use a logit (or inverse logit, a.k.a. logistic) to transform the parameters to the required probability range for the binomial outcome.

Another common transform is “Normalizing” the data to center the data around zero and to scale the outcomes such that the sample standard deviation is equal to one.

14.5.6. Rainfall Example (Continued)

We will construct a Bayesian model using the Turing.jl library. Using a battle-tested library allows us to step back from the intricacies of defining our own sampler and routine and focus on the models and analysis. The goal is to fit the parameters of one of the

competing models from above in order to demonstrate an MCMC analysis workflow and essential concepts.

The first thing that we will do is use Turing's `@model` macro to define a model. This has a few components:

1. The “model” is really just a Julia function that takes in data and relates the data to the statistical outcomes modeled.
2. The `~` is the syntax to either relate a parameter to a prior assumptions.
3. A loop (or broadcasted `.~`) that ties specific data observations to the random process.

Think of the `@model` block really as a model *constructor*. It isn't until we pass data to the model that you get a fully instantiated `Model` type¹¹.

Here's what defining the LogNormal model looks like in Turing. We have to specify prior distributions for LogNormal parameters.

```
using Turing
```

```
@model function rainLogNormal(logdata) (1)
    # Prior Assumptions for the (Log) Normal Parameters
    μ ~ Normal(4,1) (2)
    σ ~ Exponential(0.5) (3)
    # Link observations to the random process
    for i in 1:length(logdata)
        logdata[i] ~ Normal(μ, σ)
    end
end

m = rainLogNormal(log.(rain));
```

- (1) Defining the model uses the `@model` macro from Turing.
- (2) We presume that there will be positive rainfall and 96% of mean annual rainfall will be somewhere between $\exp(2)$ and $\exp(6)$, or 7 and 403 inches.
- (3) In a LogNormal model, 0.5 deviations covers a lot of variation in outcomes.

14.5.6.1. Setting Priors

In the example above, we used “weakly informative” priors. We constrained the prior probability to plausible ranges, knowing enough about the system of study (rainfall)

¹¹Specifically: a `DynamicPPL.Model` type (PPL = Probabilistic Programming Language).

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that it would be completely implausible for there to be a `Uniform(0, Inf)` distribution of mean log-rainfall total, knowing that rain can't fall in infinite quantities.

Admittedly, we haven't confirmed with a meteorologist that `exp(20)` (485 million) inches of rain per year is impossible. But such is the beauty of the transparency of Bayesian analysis that the prior assumption is right there! Front and center and ready to be debated by other modelers! If you think that 485 million inches of rain is possible next year than you can challenge this assumption and propose another explicit alternative.

"Strongly informative" priors would be something where we want to encode a stronger assumption about the plausible range of outcomes, such as if we knew enough about the problem domain that we could tell given the location of the rainfall, we'd expect 95% of the rainfall to be between, say, 10 and 30 inches per year. Then we could constrain the prior even more than we did above.

"Uninformative" priors use only maximum entropy or uniform priors to avoid encoding other assumptions into the model.

14.5.6.2. Sampling

Analysis should begin by evaluating the prior assumptions for reasonability and coverage over possible outcomes of the process we are trying to model. The top plot in Figure 14.8 shows the modeled rainfall outcomes taking on a wide range of possible outcomes. If we had more knowledge of the system we could enforce a stronger (narrower) prior assumption to constrain the model to a smaller set of values.

The object returned is an MCMCChains structure containing the samples as well as diagnostic information. Summary information gets printed below.

```
chain_prior = sample(m, Prior(), 1000)
```

Assessment of samples from the prior should include:

- Confirming that the model's behavior is reasonable that
- Confirming that the model covers the range of possible data that might be observed.

The sample outcomes from the modeled prior are shown in Figure 14.8.

Next, we sample the posterior by using the No-U-Turns (NUTS) algorithm and drawing 1000 samples (not including the warm-up phase). This is the primary result we will analyze further.

```
chain_posterior = sample(m, NUTS(), 1000)
```

Chains MCMC chain (1000×3×1 Array{Float64, 3}):

```
Iterations      = 1:1:1000
Number of chains = 1
Samples per chain = 1000
Wall duration    = 0.21 seconds
Compute duration = 0.21 seconds
parameters       = μ, σ
internals        = lp
```

Summary Statistics

| parameters | mean | std | mcse | ess_bulk | ess_tail | rhat | ... |
|------------|---------|---------|---------|-----------|----------|---------|-----|
| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 | Float64 | ... |
| μ | 4.0245 | 0.9933 | 0.0310 | 1025.0582 | 983.2831 | 0.9994 | ... |
| σ | 0.4996 | 0.4886 | 0.0154 | 1037.9963 | 974.4703 | 0.9992 | ... |

1 column omitted

Quantiles

| parameters | 2.5% | 25.0% | 50.0% | 75.0% | 97.5% |
|------------|---------|---------|---------|---------|---------|
| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 |
| μ | 2.2068 | 3.3300 | 3.9961 | 4.6884 | 5.9541 |
| σ | 0.0126 | 0.1449 | 0.3481 | 0.7167 | 1.7801 |

Figure 14.5.: Model output for the sampled prior. This isn't running an MCMC algorithm, it's simply taking draws from the defined prior assumptions.

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Chains MCMC chain (1000×14×1 Array{Float64, 3}):

```
Iterations      = 501:1:1500
Number of chains = 1
Samples per chain = 1000
Wall duration    = 2.17 seconds
Compute duration = 2.17 seconds
parameters       = μ, σ
internals        = lp, n_steps, is_accept, acceptance_rate, log_density, hamiltonian_energy, hamil
```

Summary Statistics

| parameters | mean | std | mcse | ess_bulk | ess_tail | rhat | e ... |
|------------|---------|---------|---------|----------|----------|---------|-------|
| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 | Float64 | ... |
| μ | 3.5707 | 0.0532 | 0.0018 | 891.4371 | 660.1217 | 0.9992 | ... |
| σ | 0.4518 | 0.0410 | 0.0014 | 914.6074 | 589.6091 | 0.9991 | ... |

1 column omitted

Quantiles

| parameters | 2.5% | 25.0% | 50.0% | 75.0% | 97.5% |
|------------|---------|---------|---------|---------|---------|
| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 |
| μ | 3.4658 | 3.5366 | 3.5710 | 3.6046 | 3.6681 |
| σ | 0.3825 | 0.4229 | 0.4484 | 0.4774 | 0.5358 |

Figure 14.6.: Model output for the sampled posterior.

14.5.6.3. Diagnostics

Before analyzing the result itself, we should check a few things to ensure the model and sampler were well behaved. MCMC techniques are fundamentally stochastic and randomness can cause an errant sampling path. Or a model may be mis-specified such that the parameter space to explore is incompatible with the current algorithm (or any known so far).

A few things we can check:

First, the `ess` or **effective sample size** which adjusts the number of samples for the degree of autocorrelation in the chain. Ideally, we would be able to draw independent samples from the posterior, but due to the Markov-Chain approach the samples can have autocorrelation between neighboring samples. We collect less information about the posterior in the presence of positive autocorrelation.

An `ess` greater than our sample indicates that there was less (negative) autocorrelation than we would have expected for the chain. An `ess` much less than the number of samples indicates that the chain isn't sampling very efficiently but, aside from needing to run more samples, isn't necessarily a problem.

```
ess(chain_posterior)
```

| ESS | | |
|------------|----------|-------------|
| parameters | ess | ess_per_sec |
| Symbol | Float64 | Float64 |
| μ | 891.4371 | 410.9899 |
| σ | 914.6074 | 421.6724 |

Second, the `rhat` (\hat{R}) is the Gelman-Rubin convergence diagnostic and it's value should be very close to 1.0 for a chain that has converged properly. Even a value of 1.01 may indicate an issue and quickly gets worse for higher values.

```
rhat(chain_posterior)
```

| R-hat | |
|------------|---------|
| parameters | rhat |
| Symbol | Float64 |
| μ | 0.9992 |
| σ | 0.9991 |

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Next, we can look at the “trace” plots for the parameters being sampled (Figure 14.7). These are sometimes called “hairy caterpillar” plots because in a healthy chain sample, we should see a series without autocorrelation and that the values bounce around randomly between individual samples.

```
let
    f = Figure()
    ax1 = Axis(f[1,1],ylabel="μ")
    lines!(ax1,vec(get(chain_posterior,:μ).μ.data))
    ax2 = Axis(f[2,1],ylabel="σ")
    lines!(ax2,vec(get(chain_posterior,:σ).σ.data))
    f
end
```

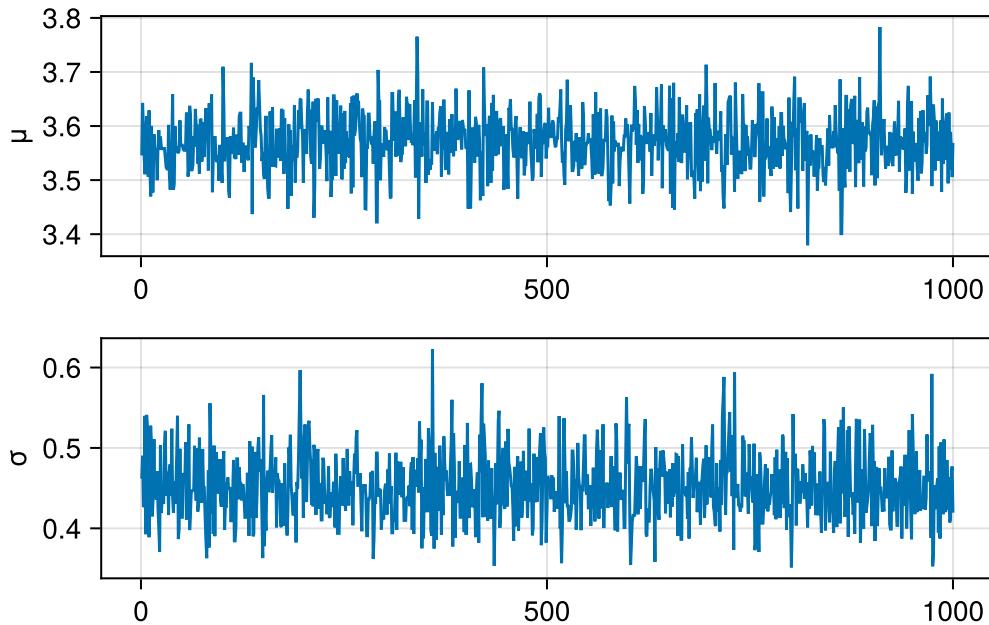


Figure 14.7.: The trace plots indicate low autocorrelation which is desirable for an MCMC sample.

The ess, rhat, and trace plots all look good for our sampled chain so we can next we will analyze the results in the context of our rainfall problem.

14.5.6.4. Analysis

Let's see how it looks compared to the data first. Figure 14.8 shows 200 samples from the prior and posterior. The prior (top) shows how wide the range of possible rainfall outcomes could be using our weakly informative prior assumptions. The bottom shows that after having learned from the data, the posterior probability of rainfall has narrowed considerably.

```

function chn_cdf!(axis,chain,rain)
    n = 200
    s = sample(chain, n)
    vals = get(s, [:μ, :σ])
    ds = Normal.(vals.μ, vals.σ) # ,get(s,:σ)[i])
    rg = 1:200
    for (i, d) in enumerate(ds)
        lines!(axis, rg,cdf.(d,log.(rg)),color=(:gray,0.3))
    end

    # plot the actual data
    percentiles= 0.01:0.01:0.99
    lines!(axis,quantile.(Ref(rain),percentiles),percentiles,linewidth=3)
end

let
    f = Figure()
    ax1 = Axis(f[1,1],title="Prior", xgridvisible=false, ygridvisible=false,ylabel="Quantile")
    chn_cdf!(ax1,chain_prior,rain)

    ax2 = Axis(f[2,1],title="Posterior", xgridvisible=false, ygridvisible=false,xlabel="Annual
    chn_cdf!(ax2,chain_posterior,rain)

    linkxaxes!(ax1, ax2)

    f
end

```

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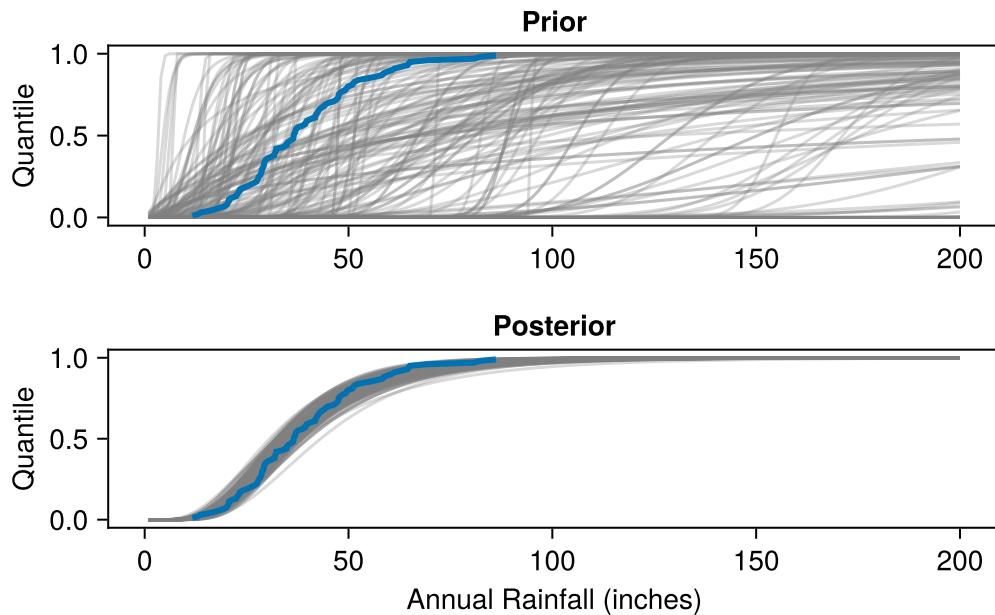


Figure 14.8.: The prior model show a wide range of possible outcomes, and the shape of the distribution is reasonable: there's a nice 'S' shape to the CDF, indicating a dense region where most outcomes would fall in the PDF. The fitted posterior model (bottom) has good coverage of the observed data (shown in blue).

Comparing to the maximum likelihood analysis from before by plotting the MLE point estimate onto the marginal densities in Figure 14.9. The peak of the the posterior is referred to as the **maximum a posteriori** (MAP) and would be the point estimate proposed by this Bayesian analysis. However, the Bayesian way of thinking about distributions of outcomes rather than point estimates is one of the main aspects we encourage for financial modelers. Using the posterior distribution of the parameters, we can assess parameter uncertainty directly instead of ignoring it as we tend to do with point estimates.

```
let
    # get the parameters from the earlier MLE approach
    p = params(ln)

    f = Figure()

    # plot μ posterior
    ax1 = Axis(f[1,1],title="μ posterior",xgridvisible=false)
    hideydecorations!(ax1)
```

```

d = density!(ax1,vec(get(chain_posterior,:μ).μ.estimate))
l = vlines!(ax1,[p[1]],color=:red)

# plot σ posterior
ax2 = Axis(f[2,1],title="σ posterior", xgridvisible=false)
hideydecorations!(ax2)
density!(ax2,vec(get(chain_posterior,:σ).σ.estimate))
vlines!(ax2,[p[2]],color=:red)

Legend(f[1,2],[d,l],["Posterior Density", "MLE Estimate"])

f
end

```

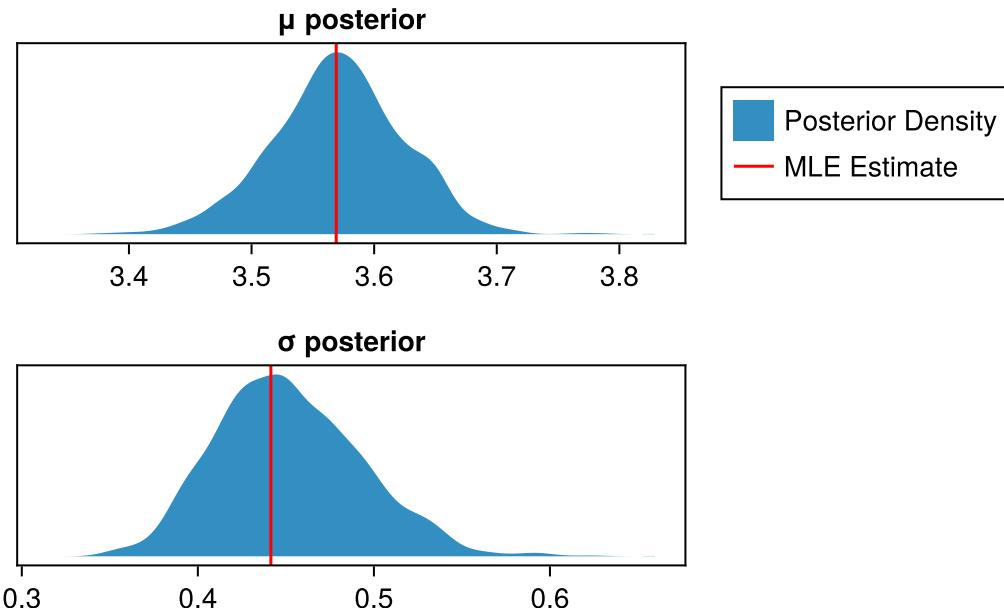


Figure 14.9.: The MLE point estimate need not necessarily align with the peak or center of posterior densities (e.g. in the case of a bimodal distribution).

14.5.6.5. Model Limitations

We have built and assessed a simple statistical model that could be used in the estimation of risk for a particular location. Nowhere in our model did we define a mechanism to

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capture a more sophisticated view of the world. There is no parameter for changes over time due to climate change, or inter-annual seasonality for El Niño or La Niña cycles, or any of a multitude of other real-world factors that can influence the forecasting. All we've defined is that there is a LogNormal process generating rainfall in a particular location. This may or may not be sufficient to capture the dynamics of the problem at hand.

Part of the benefits of the Bayesian approach is that it allows us to extend the statistical model to be arbitrarily complex in order to capture our intended dynamics. We are limited by the availability of data, computational power and time, and our own expertise in the modeling. Regardless of the complexity of the model, the same fundamental techniques and idea apply in the Bayesian approach.

14.5.6.6. Continuing the Analysis

Like any good model, you can often continue the analysis in any number of directions, such as: collecting more data, evaluating different models, creating different visualizations, making predictions about future events, creating a multi-level model that predicts rainfall for multiple related locations simultaneously, among many other threads of analysis.

Earlier we discussed model comparison. To compute a real Bayes Factor in comparing the different models, we would take the average likelihood across the posterior samples instead of just comparing the maximum likelihood points as we did earlier. There are more sophisticated tools for estimating out of sample performance of the model, or measures that evaluate a model for over-fitting by penalizing the diagnostic statistic for the model having too many free parameters. See LOO (leave-one-out) cross-validation and various "information criteria" in the resources listed in Section 14.5.8.

14.5.7. Conclusion

This chapter has attempted to make accessible the foundations of statistical inference and the modern tools and approaches available. Underlying this approach to thinking about statistical problems are informational theoretic and mathematical concepts that can be challenging to learn. This is especially true when traditional finance and actuarial curricula is not centered on the necessary computational foundations associated with modern statistical analysis. Most importantly, moving beyond single 'best estimate' values to embrace full probability distributions leads to richer financial analysis and more comprehensive risk assessment.

14.5.8. Further Reading

Bayesian approaches to statistical problems are rapidly changing the professional statistical field. To the extent that the actuarial profession incorporates statistical procedures, financial professionals should consider adopting the same practices. The benefits of this are a better understanding of the distribution of risk and return, results that are more interpretable and explainable, and techniques that can be applied to a wider range of problems. The combination of these things would serve to enhance best practices related to understanding and communicating about financial quantities.

Textbooks recommended by the author are:

- Statistical Rethinking (McElreath)
- Bayes Rules! (Johnson, Ott, Dogucu)
- Bayesian Data Analysis (Gelman, et. al.)

Chi Feng has an interactive demonstration of different MCMC samplers available at:
<https://chi-feng.github.io/mcmc-demo/>.

15. Stochastic Modeling

Yun-Tien Lee and Alec Loudenback

The Monte Carlo Method: (i) A last resort when doing numerical integration, and (ii) a way of wastefully using computer time. - Malvin H. Kalos¹ (c. 1960)

15.1. In This Chapter

Brief overview of stochastic modeling concepts. Random inputs vs. embedded randomness in stepwise models. Key components: distributions, time horizons, state space. Evaluating outputs with expectation, variance, and risk measures. Special properties like Markov, stationarity, and SDEs. Practical finance examples with Julia code for scenario generation and macroeconomic simulations.

15.2. Introduction

Stochastic modeling is the technique of running many similar scenarios through a model in order to evaluate different aspects of a given model. Some contexts in which stochastic modeling arises in the financial context include:

- **Valuation.** Market consistent pricing often involves calibrating a set of stochastic economic scenarios to market-observed prices. Those calibrated scenarios are then used to determine the value of another set of contracts for which market prices are not directly observable.
- **Scenario Analysis.** Many financial models have a high degree of complexity and interactions between components not well understood *a priori*. Using a set of calibrated or otherwise arbitrary scenarios can illuminate unexpected or interrelated behavior that arises from the model. Sometimes financial assets and liabilities have a path-dependency, meaning that the overall behavior or value is a function of the entire path taken (e.g. an asian option).

¹Kalos was a pioneer in Monte Carlo techniques, quoted via https://doi.org/10.1007/978-3-540-74686-7_3

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- **Risk and Capital Analysis.** Whether model parameters are point estimates (“an asset has a 1% probability of default”) or distributions (“an asset has a probability of default distributed as $Beta(1, 99)$ ”), stochastic modeling can be utilized to determine projected capital levels for a company or line of business at certain security levels.

The modeled outcomes are only as good as the scenarios and model themselves. For example, for many years the predominant economic scenario generator for US regulated insurance liabilities (the American Academy of Actuaries’ Economic Scenario Generator) would produce interest rate paths that were substantially higher than then-current market-implied paths. It’s erroneous in those circumstances to compute the 99th-percentile scenario and characterize the scenario as “extreme.” Similarly, there may be known or unknown deficiencies with the modeled behavior and this should be considered by the modeler when discussing the results.

Recall the kinds of uncertainties in Table 4.2 - stochastic modeling implementations often mechanically only address the aleatory (process) volatility component of outcome uncertainty. Through techniques like developing a posterior distribution for model parameters (@sec-bayes-rule) we can also explicitly attempt to model epistemic (parameter) uncertainty as well.

15.3. Fundamentals of Stochastic Modeling

Stochastic modeling represents a significant departure from deterministic approaches by explicitly incorporating randomness and uncertainty. Before diving into specific modeling techniques, it’s essential to understand the fundamental concepts that form the foundation of any stochastic model.

15.3.1. Random Variables and Probability

At the core of stochastic modeling is the concept of random variables—quantities whose values depend on random outcomes. In financial contexts, random variables might represent:

- Future interest rates
- Stock price movements
- Insurance claim frequencies
- Default probabilities
- Mortality rates

15.3. Fundamentals of Stochastic Modeling

These random variables are characterized by probability distributions that describe the likelihood of different outcomes. For example, stock returns might follow a normal distribution, while insurance claims might follow a Poisson distribution for frequency and a LogNormal distribution for severity.

15.3.2. The Role of Monte Carlo Simulation

Despite the opening quote's humorous skepticism, **Monte Carlo methods** have become indispensable in modern financial modeling. Named after the famous casino, these techniques use repeated random sampling to obtain numerical results when analytical solutions are impractical or impossible.

The basic procedure involves:

1. Generating random inputs based on specified probability distributions
2. Running the model with these inputs
3. Collecting and analyzing the outputs
4. Repeating until sufficient samples are obtained for statistical significance

As computational power has increased, Monte Carlo methods have moved from "last resort" to standard practice in many financial applications. In addition to this conceptual chapter, Chapter 26 demonstrates this process to generate stochastic economic scenarios.

15.3.3. Time Dimensions in Stochastic Models

Stochastic models can be categorized based on how they treat time:

- Discrete-time models: The system evolves in distinct time steps (daily, monthly, annually)
- Continuous-time models: The system evolves continuously, often described by stochastic differential equations

Financial applications frequently use both approaches. For example, option pricing might use continuous-time models (Black-Scholes), while asset-liability management might use discrete-time models with monthly or quarterly steps.

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15.3.4. State Spaces and Transitions

A stochastic model's **state space** represents all possible states the system can occupy. This might be:

- Discrete: Limited number of possible states (e.g., credit ratings: AAA, AA, A, etc.)
- Continuous: Infinite possible states (e.g., stock prices can take any positive real value)

The transitions between states are governed by probability distributions or transition matrices, depending on the model type. Understanding the state space is crucial for properly formulating and solving stochastic problems in finance.

15.3.5. Aleatory vs. Epistemic Uncertainty

Stochastic models address two fundamental types of uncertainty:
- Aleatory uncertainty: The inherent randomness in the system (e.g., market volatility)
- Epistemic uncertainty: Uncertainty due to limited knowledge about model parameters

Most basic stochastic models focus on aleatory uncertainty, assuming model parameters are known with certainty. More sophisticated approaches incorporate epistemic uncertainty by treating parameters themselves as random variables, often using Bayesian methods.

15.3.6. Calibration and Validation

Before a stochastic model can be applied, it must be properly calibrated to match observed data or market prices. This typically involves:

1. Establishing initial parameters via professional judgment or statistical means.
2. Adjusting parameters to match desired calibration data: e.g. minimizing a loss function with respect to the predicted outcomes versus the calibration data (see @sec-optimization).
3. Validating the model against out-of-sample data.

For financial applications, models are often calibrated to be "market-consistent," meaning they reproduce the prices of traded instruments before being used to value more complex products.

15.3.7. Computational Considerations

Stochastic modeling is computationally intensive, particularly when:

- Many scenarios are required for convergence
- The model has high dimensionality
- Path dependencies require storing entire trajectories

Modern implementations leverage parallel processing, variance reduction techniques, and sometimes machine learning approximations to make complex stochastic models computationally feasible.

With these fundamental concepts in mind, we can now explore specific types of stochastic models and their applications in finance and actuarial science.

15.3.8. Kinds of Stochastic Models

15.3.8.1. Input Ensemble

In this approach, you first generate many sets of random inputs (scenarios) and then run the same model repeatedly—one run per scenario²:

- **Interest Rate Scenarios for Bond Portfolios**

Generate thousands of yield-curve scenarios (e.g., parallel shifts, twists) to test how a bond portfolio might perform over time. By averaging the results over all simulations, you can estimate the portfolio's expected return or risk metrics.

- **Commodity Price Paths for Hedging**

Create multiple possible paths for commodity prices under different macro conditions. Use these to evaluate various hedging strategies and see which one consistently mitigates risk across a wide range of market conditions.

Generally, models programmed in code can easily map across a large number of stochastic inputs.

15.3.8.2. Random State

Here, the model itself “rolls the dice” at each step. Instead of pre-generating input scenarios, the randomness is embedded in the logic that evolves the state over time³:

²An extended example of this approach is discussed in Chapter 26.

³A worked example of this approach is illustrated in Chapter 25.

15. Stochastic Modeling

- **Property & Casualty Claims Simulation**

In each simulated month, randomly draw the number of claims (from a Poisson or negative binomial distribution), then randomly assign the severity of each claim (often a lognormal or gamma distribution). The insurer's financial state (reserves, surplus) evolves based on these simulated losses.

- **Call-Center (Queueing) Model**

Model the random arrival of customer calls following a Poisson process, and at each arrival decide if the call is answered immediately or queued. Service times for each call can be drawn from an exponential distribution. By simulating many days, you can estimate average waiting times or the probability that a caller has to wait more than a certain threshold.

i Note

Many financial models project outcomes using *expected values* of random variables. Many actuarial models utilize this approach to determine expected outcomes given an assumption (such as expected present value of life insurance claims). This is different than evaluating many possible discrete realized scenarios wherein each insured life either dies or survives each period, and models designed for projected expected payments need to be adapted to handle a random state approach.

15.3.8.3. Closed-Form

Some stochastic processes admit analytical solutions or formulas for key quantities, so you do not need to run a simulation for each scenario. For example, here are a couple of common models that have closed form solutions to the underlying stochastic differential equation.

- **Black Scholes Option Price**

The Black-Scholes European call and put-call price parity are results for analytically pricing options despite the underlying theory being based on a stochastic evolution of asset prices.

- **Ornstein-Uhlenbeck Interest Rate Model**

An Ornstein-Uhlenbeck process is often used to model mean-reverting interest rates. Despite being fundamentally stochastic, it has closed-form expressions for zero-coupon bond prices and other interest-rate derivatives, saving you from running thousands of random simulations for valuation.

The focus of this section will be on the other kinds of computational-driven, stochastic models.

15.3.8.4. Special Cases or Properties

Stochastic processes often combine several structural or theoretical properties to reflect real-world phenomena in financial and actuarial models. Before diving into each property—Markov, stationarity, martingales, and SDEs—it can be useful to see how they serve as building blocks. For example, Markov processes allow today's decisions or valuations to depend only on the current state, whereas stationarity implies that the underlying statistical characteristics (mean or variance) do not change over time. Martingales capture a notion of “fairness” in gambling and risk-neutral pricing, and SDEs inject random shocks in a continuous-time setting. By understanding these properties, one can better see which assumptions are embedded in derivatives pricing, credit migration models, or mortality forecasts.

15.3.8.4.1. Markov Models

A stochastic process has the Markov property if the future state depends only on the current state and not on the past states (memorylessness). Markov chains and hidden Markov models contain this property. The Markov property simplifies modeling and computation by reducing dependencies on past states. For instance, a credit rating model might assume that the probability of going from BBB to BB depends only on today's rating, without direct reference to prior years' ratings.

15.3.8.4.2. Stationary Models

A stochastic process is stationary if its statistical properties, such as mean and variance, do not change over time. Strict stationarity means all moments of the process are invariant over time, while weak Stationarity means only the mean and variance are invariant over time. Stationarity simplifies the analysis and modeling of stochastic processes, especially for time series data. Sometimes a non-stationary model can be made stationary by transformation (such as removing a constant trend component).

15.3.8.4.3. Martingales

A martingale is a stochastic process where the expected future value, given all prior information, is equal to the current value. Examples include fair gambling games or financial asset prices in efficient markets. Martingales are important in financial modeling and risk-neutral pricing, especially for derivatives.

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15.3.8.4.4. Stochastic Differential Equations (SDEs)

These are differential equations that incorporate stochastic terms (typically driven by Brownian motion or other noise sources). The well-known Black-Scholes equation for option pricing is one example of how SDE can be applied in real world. SDEs are essential for modeling systems where both deterministic and random factors drive behavior over time.

A special kind of SDE is **Brownian Motion**. This is a continuous-time stochastic process where changes over time are independent and normally distributed. This is extensively used in financial models for asset price movements (e.g., the Black-Scholes model). In addition to financial markets, Brownian motion is also widely used in modeling physical systems (diffusion).

In practice, these properties often occur together in financial models.

15.3.9. Components of Stochastic Models

Stochastic modeling has additional terminology to introduce.

15.3.9.1. Random variables

A random variable represents a quantity whose value is determined by the outcome of a random event. It can be discrete or continuous. It is essentially a mapping from event space to numerical values. Examples include stock prices, waiting time in queues and number of claims in insurance, etc. Random variables form the basis of stochastic models by introducing uncertainty into the model.

15.3.9.2. Probability distributions

A probability distribution describes the likelihood of different outcomes for a random variable. Examples include normal distribution, Poisson distribution and exponential distribution, etc. Probability distributions help in modeling the behavior of random variables and in defining how likely different events or outcomes are.

15.3.9.3. State space

The state space represents all possible states or values that a stochastic process can take. For example, in a Markov process, the state space might be the set of all possible values that the system can occupy at any given time. The state space helps in defining the scope of the model by specifying possible outcomes.

15.3. Fundamentals of Stochastic Modeling

15.3.9.4. Stochastic processes

A stochastic process is a collection of random variables indexed by time (or some other variable) that evolve in a probabilistic manner. Examples include Brownian motion, Markov chains and Poisson processes. Stochastic processes model how random variables change over time, which is essential for understanding dynamic systems influenced by randomness.

15.3.9.5. Transition probabilities

These represent the probabilities of transitioning from one state to another in a stochastic process. For example, in a Markov chain, the transition matrix contains the probabilities of moving from one state to another. Transition probabilities determine how the system evolves from one time step to the next, reflecting the underlying randomness.

15.3.9.6. Time horizon

The time horizon refers to the period over which the stochastic process is observed. It can be discrete (e.g., steps in a Markov chain) or continuous (e.g., continuous-time models like Brownian motion). The time horizon helps in determining how the process behaves over short or long periods.

15.3.9.7. Initial conditions

These are the starting points or initial values of the random variables or the system at time zero. They may be initial price of a stock, initial number of customers in a queue, etc. The starting condition influences the future evolution of the process, and different initial conditions can lead to different outcomes.

15.3.9.8. Noise (random shocks)

Random noise represents unpredictable random fluctuations that can affect the outcome of a stochastic process. This can be market volatility, measurement errors or environmental variations. Noise is a critical element in stochastic models as it introduces randomness and uncertainty into otherwise deterministic systems.

15.4. Evaluating and Applying Stochastic Models

15.4.1. Evaluating Stochastic Results

These types of models are evaluated simply by running them many times until the measure of interest converges on a stable result: for example, we might run a model until the mean of the results no longer varies materially as we add more scenarios.

15.4.1.1. Expectation and variance

The expected value (mean) represents the average or mean outcome of a random variable over many trials or realizations. The variance measures the spread or variability of outcomes around the expected value. These statistical measures provide insights into the central tendency and the uncertainty or risk in a stochastic model.

15.4.1.2. Covariance and correlation

Covariance measures how two random variables change together. Positive covariance indicates that the variables tend to increase together. On the other hand correlation is the standardized version of covariance that measures the strength of the linear relationship between two variables. Understanding how different random variables interact helps in building more complex models, especially in multivariate stochastic processes.

15.4.1.3. Risk Measures

Risk measures encompass the set of functions that map a set of outcomes to an output value characterizing the associated riskiness of those outcomes. As is usual when attempting to compress information (e.g. condensing information into a single value), there are multiple ways we can characterize this riskiness.

15.4.1.3.1. Coherence & Other Desirable Properties

Further, it is desirable that a risk measure has certain properties, and risk measures that meet the first four criteria are called “Coherent” in the literature. From “An Introduction to Risk Measures for Actuarial Applications” ((Hardy 2006)), she describes four properties. Using H as a risk measure and X as the associated risk distribution:

15.4. Evaluating and Applying Stochastic Models

15.4.1.3.1.1. 1. Translation Invariance

For any non-random c

$H(X+c) = H(X)+c$ This means that adding a constant amount (positive or negative) to a risk adds the same amount to the risk measure. It also implies that the risk measure for a non-random loss, with known value c , say, is just the amount of the loss c .

15.4.1.3.1.2. 2. Positive Homogeneity

For any non-random $\lambda > 0$:

$$H(\lambda X) = \lambda H(X)$$

This axiom implies that changing the units of loss does not change the risk measure.

15.4.1.3.1.3. 3. Subadditivity

For any two random losses X and Y ,

$$H(X + Y) \leq H(X) + H(Y)$$

It should not be possible to reduce the economic capital required (or the appropriate premium) for a risk by splitting it into constituent parts. Or, in other words, diversification (ie consolidating risks) cannot make the risk greater, but it might make the risk smaller if the risks are less than perfectly correlated.

15.4.1.3.1.4. 4. Monotonicity

If $Pr(X \leq Y) = 1$ then $H(X) \leq H(Y)$.

If one risk is always bigger than another, the risk measures should be similarly ordered.

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15.4.1.3.1.5. Other Properties

In “Properties of Distortion Risk Measures” (Balbás, Garrido, and Mayoral 2009) also note other properties of interest:

Complete

Completeness is the property that the distortion function associated with the risk measure produces a unique mapping between the original risk's survival function $S(x)$ and the distorted $S^*(x)$ for each x . See Distortion Risk Measures for more detail on this.

In practice, this means that a non-complete risk measure ignores some part of the risk distribution (e.g. CTE and VaR don't use the full distribution and have the same)

Exhaustive

A risk measure is “exhaustive” if it is coherent and complete.

Adaptable

A risk measure is “adapted” or “adaptable” if its distortion function (see Distortion Risk Measures).

1. g is strictly concave, that is g is strictly decreasing.
2. $\lim_{u \rightarrow 0^+} g'(u) = \inf$ and $\lim_{u \rightarrow 1^-} g'(u) = 0$.

Adaptive risk measures give more weight to larger losses. They are exhaustive but the converse is not true.

15.4.1.4. Summary of Risk Measure Properties

| Measure | Coherent | Complete | Exhaustive | Adaptable |
|--|----------|----------|------------|-----------|
| VaR | No | No | No | No |
| CTE | Yes | No | No | No |
| DualPower
($y > 1$) | Yes | Yes | Yes | No |
| ProportionalHazard
($\gamma > 1$) | Yes | Yes | Yes | No |
| WangTransform | Yes | Yes | Yes | Yes |

i Distortion Risk Measures

Distortion Risk Measures are a way of remapping the probabilities of a risk distribution in order to compute a risk measure H on the risk distribution X .

Adapting (Wang 2002), there are two key components:

15.4.1.5. Distortion Function $g(u)$

This remaps values in the $[0,1]$ range to another value in the $[0,1]$ range, and in H below, operates on the survival function S and $F = 1 - S$.

Let $g : [0, 1] \rightarrow [0, 1]$ be an increasing function with $g(0) = 0$ and $g(1) = 1$. The transform $F^*(x) = g(F(x))$ defines a distorted probability distribution, where “ g ” is called a distortion function.

Note that F^* and F are equivalent probability measures if and only if $g : [0, 1] \rightarrow [0, 1]$ is continuous and one-to-one. Definition 4.2. We define a family of distortion risk-measures using the mean-value under the distorted probability $F^*(x) = g(F(x))$:

15.4.1.6. Risk Measure Integration

To calculate a risk measure H , we integrate the distorted F across all possible values in the risk distribution (i.e. $x \in X$):

$$H(X) = E^*(X) = - \int_{-\infty}^0 g(F(x))dx + \int_0^{+\infty} [1 - g(F(x))]dx$$

That is, the risk measure (H) is equal to the expected value of the distortion of the risk distribution ($E^*(X)$).

15.4.1.7. Risk Measures: Examples

15.4.1.7.1. Basic Comparison

We won’t re-implement the logic here, but here’s a very simple example demonstrating the relative values of VaR, CTE, and a Wang Transform at the 90% level. These functions come from the public library ActuaryUtilities.jl.

```
using ActuaryUtilities

let outcomes = rand(1000)

@show VaR(0.90)(outcomes)
@show CTE(0.90)(outcomes)
@show WangTransform(0.90)(outcomes)
end

(VaR(0.9))(outcomes) = 0.918416152077607
(CTE(0.9))(outcomes) = 0.9574360390555113
(WangTransform(0.9))(outcomes) = 0.8280642174963452
```

15. Stochastic Modeling

0.8280642174963452

15.4.1.8. Plotting a Range of Values

We will generate a random outcome and show how the risk measures behave:

```
using Distributions
using ActuaryUtilities
using CairoMakie

# the assumed distribution of outcomes
outcomes = Weibull(1, 5)

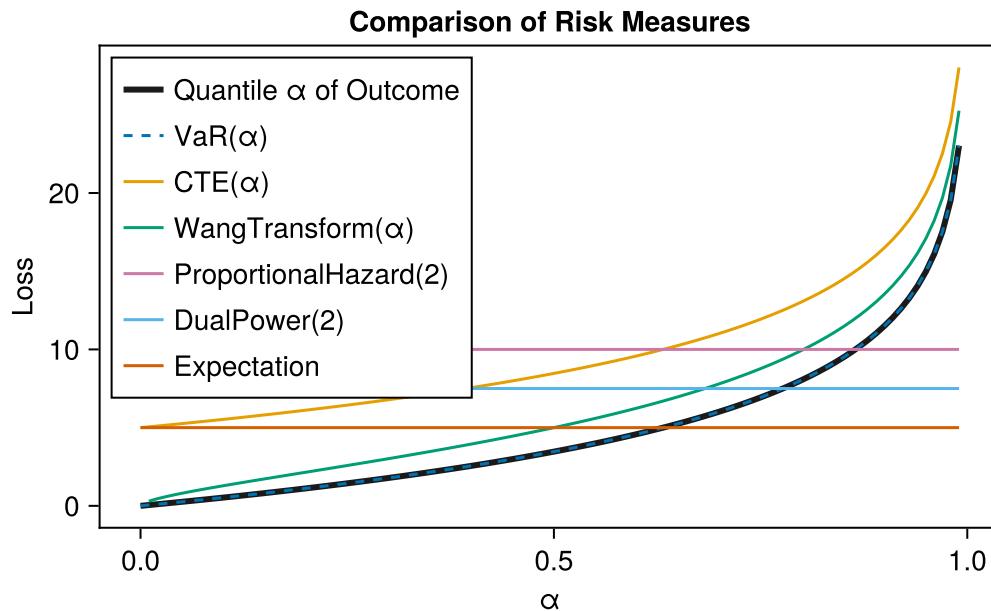
αs = range(0.00, 0.99; length=100)

let
    f = Figure()
    ax = Axis(f[1, 1],
              xlabel="α",
              ylabel="Loss",
              title="Comparison of Risk Measures",
              xgridvisible=false,
              ygridvisible=false,
    )
    lines!(ax,
           αs,
           [quantile(outcomes, α) for α in αs],
           label="Quantile α of Outcome",
           color=:grey10,
           linewidth=3,
    )
    lines!(ax,
           αs,
           [VaR(α)(outcomes) for α in αs],
           label="VaR(α)",
           linestyle=:dash
    )
    lines!(ax,
           αs,
           [CTE(α)(outcomes) for α in αs],
```

15.4. Evaluating and Applying Stochastic Models

```
    label="CTE(α)",  
)  
lines!(ax,  
    αs[2:end],  
    [WangTransform(α)(outcomes) for α in αs[2:end]],  
    label="WangTransform(α)",  
)  
lines!(ax,  
    αs,  
    [ProportionalHazard(2)(outcomes) for α in αs],  
    label="ProportionalHazard(2)",  
)  
  
lines!(ax,  
    αs,  
    [DualPower(2)(outcomes) for α in αs],  
    label="DualPower(2)",  
)  
lines!(ax,  
    αs,  
    [RiskMeasures.Expectation()(outcomes) for α in αs],  
    label="Expectation",  
)  
axislegend(ax, position=:lt)  
  
f  
end
```

15. Stochastic Modeling



15.4.2. Variance Reduction Techniques

While Monte Carlo simulation is a powerful approach for modeling complex financial or actuarial systems, it can require a very large number of random draws for the results to converge. When there have been a limited number of runs, the variance of the sampled results can be high, and vary materially between different sets of stochastic runs. This can be addressed by simply running more scenarios, but can be computationally expensive, particularly when path-dependent processes—or high-dimensional models—are involved. For this reason, several “variance reduction techniques” have been established to improve efficiency without reducing accuracy. They work by designing the way we draw samples (or process them) so that desired statistics converge more quickly.

15.4.2.1. Control Variates

A control variate is a random variable whose expected value is already known (or can be estimated with high precision). By simulating the main process alongside the control variate and leveraging the difference in outcomes, the overall variance of the estimator is reduced. In finance, for instance, one might use a simpler instrument whose price is analytically known as a control variate when simulating a more complex derivative.

15.4.2.2. Antithetic Variates

In this technique, each random draw (X) is paired with a corresponding draw ($-X$) or another symmetric transformation. The idea is that positive and negative deviations tend to cancel out, thereby reducing overall variability. For example, if simulating geometric Brownian motion paths, each path with a particular series of random shocks could be matched with a path using the same shocks negated. Updating estimates with the average result of these pairs generally yields a lower variance for the same number of total draws.

15.4.2.3. Stratified Sampling

Instead of sampling purely at random across the entire probability distribution, the distribution is divided into “strata” (or sub-intervals), and draws are forced to fall into each of these strata with a pre-specified frequency. By covering the distribution more systematically, the resulting estimates often converge faster. This is sometimes called **quasi-Monte Carlo** and involves techniques to span the sample space that is less random but still unbiased. Julia has a package for this called QuasiMonteCarlo.jl

Together, these techniques reflect a crucial principle of computational thinking: by leveraging structure or known properties of the simulation, you can achieve more accurate estimates with fewer overall scenarios. That translates into faster runs, less strain on computational resources, and better performance when embedding these methods in large-scale models or iterative processes.

15.4.3. Stochastic Modeling: Examples

15.4.3.1. Macroeconomic Analysis

Here we show still another stochastic process in macroeconomic analysis. Stochastic macroeconomic analysis often involves modeling random shocks and their effects on macroeconomic variables such as output, consumption, inflation, and employment. One common approach is through Dynamic Stochastic General Equilibrium (DSGE) models, which are widely used in macroeconomic analysis. These models incorporate randomness (stochastic elements) to capture real-world uncertainty in economic systems.

```
using Random, CairoMakie, Distributions

# Parameters
α = 0.33                      # Capital share of output
δ = 0.05                        # Depreciation rate
```

15. Stochastic Modeling

```
s = 0.2                      # Savings rate
n = 0.01                      # Population growth rate
g = 0.02                      # Technology growth rate
σ = 0.01                      # Standard deviation of productivity shocks
T = 100                        # Number of periods to simulate
K0 = 1.0                       # Initial capital stock
A0 = 1.0                       # Initial productivity

# Shock distribution (normal distribution for productivity shocks)
shock_distribution = Normal(0, σ)

# Random number generator (Xoshiro as default with a seed to ensure reproducability)
rng_gen = Xoshiro(1234)

# Function to simulate the model
function simulate_stochastic_solow(T, α, δ, s, n, g, σ, K0, A0)
    K = zeros(T)                # Capital over time
    Y = zeros(T)                # Output over time
    A = zeros(T)                # Productivity shocks over time
    A[1] = A0                   # Initial productivity
    K[1] = K0                   # Initial capital

    for t in 1:T-1
        # Apply random productivity shock
        ε_t = rand(rng_gen, shock_distribution)
        A[t+1] = A[t] * exp(ε_t) # Productivity evolves stochastically

        # Output based on Cobb-Douglas production function
        Y[t] = A[t] * K[t]^α

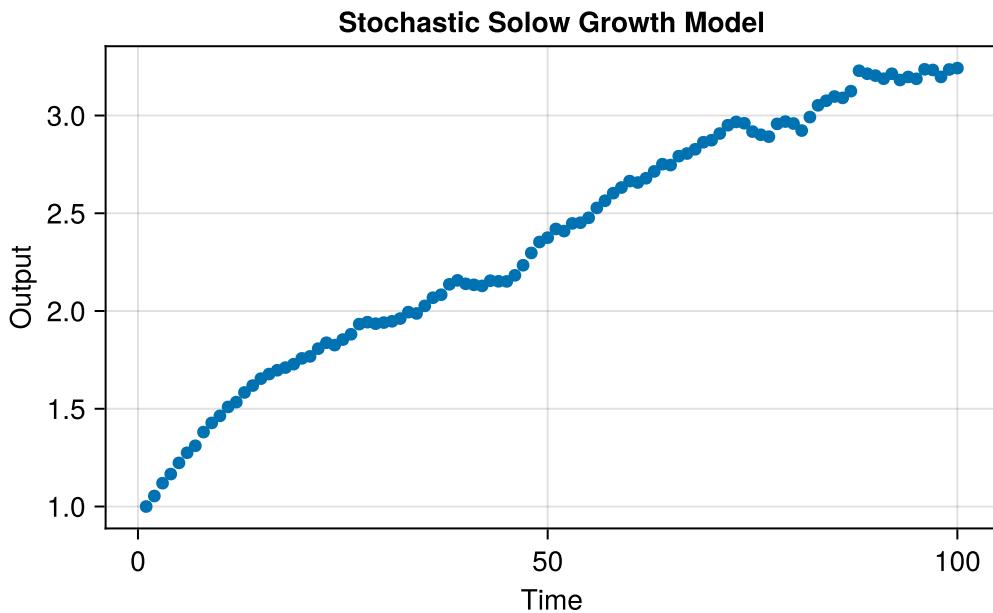
        # Capital accumulation equation
        K[t+1] = s * Y[t] + (1 - δ) * K[t]

        # Population and technology growth
        K[t+1] *= (1 + n) * (1 + g)
    end

    Y[T] = A[T] * K[T]^α # Final output
    return K, Y, A
end

# Simulate the model
K, Y, A = simulate_stochastic_solow(T, α, δ, s, n, g, σ, K0, A0)
```

```
# Plot the results
f = Figure()
axis = Axis(f[1, 1], xlabel="Time", ylabel="Output", title="Stochastic Solow Growth Model")
scatter!(1:T, Y, label="Output (Y)")
f
```



15.4.3.2. Other Examples

Stochastic examples can be found in many other sections of this book, such as:

- Portfolio optimization applications (Chapter 29).
- Genetic algorithms for certain optimization problems (Chapter 17).
- Stochastic state-based mortality projections (Chapter 25).

15.5. Conclusion

Stochastic modeling involves the study of systems influenced by random factors and uncertainty. By combining random variables, probability distributions, and processes like Markov chains or Brownian motion, stochastic models provide insights into systems that cannot be described purely deterministically.

16. Automatic Differentiation

“It is unworthy of excellent men to lose hours like slaves in the labour of calculation which could safely be relegated to anyone else if machines were used. — Gottfried Wilhelm Leibniz (Describing, in 1685, the value of his hand-cranked calculating machine invention.)”

16.1. In This Chapter

Harnessing the chain rule to compute derivatives not just of simple functions, but of complex programs.

16.2. Motivation for (Automatic) Derivatives

Derivatives are one of the most useful analytical tools we have. Determining the rate of change with respect to an input is effectively sensitivity testing. Knowing the derivative lets you optimize things faster (see Chapter 17). You can test properties and implications (monotonicity, maxima/minima). These applications make derivatives foundational for machine learning, generative models, scientific computing, and more.

We will work up concepts on computing derivatives, from the most basic (finite differentiation) to automatic differentiation (forward mode and then reverse mode). These tools can be used in many modeling applications.

16.3. Finite Differentiation

Finite differentiation is evaluating a function $f(x)$ at a value x and then at a nearby value $x + \epsilon$. The line drawn through these two points effectively estimates the line that is tangent to the function f at x - effectively the derivative has been found by approximation. That is, we are looking to approximate the derivative using the property:

$$f'(x) = \lim_{\epsilon \rightarrow 0} \frac{f(x_0 + \epsilon) - f(x_0)}{\epsilon}$$

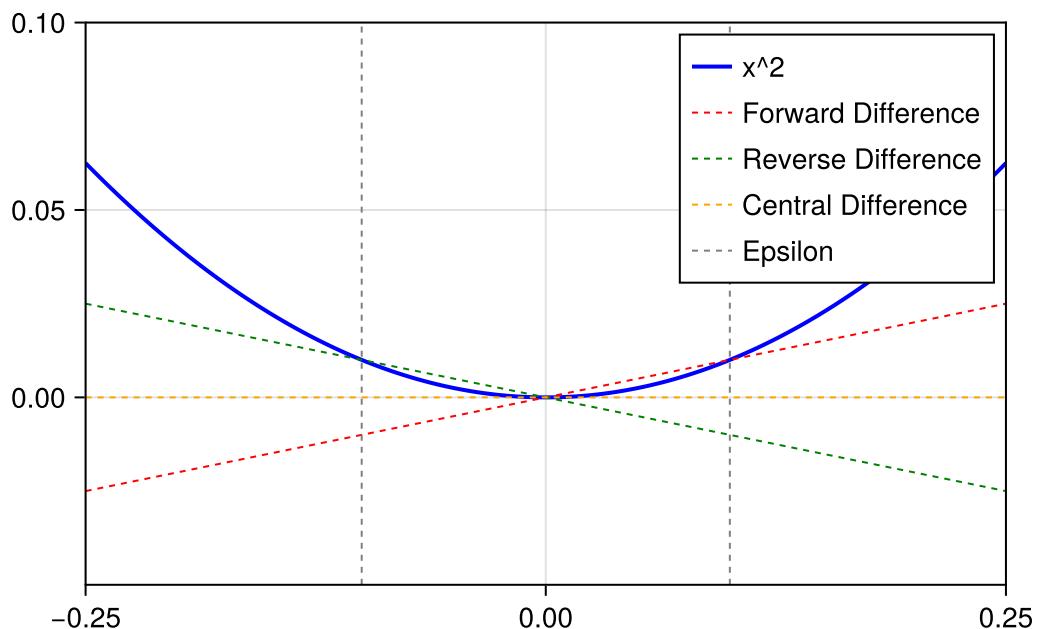
16. Automatic Differentiation

We can approximate the result by simply choosing a small ϵ .

There's also flavors of finite differentiation to approximate derivatives to be aware of:

- forward difference is as defined in the above equation, where ϵ is *added* to x_0
- reverse difference is as defined in the above equation, where ϵ is *subtracted* from x_0
- central difference is where we evaluate at $x_0 \pm \epsilon$ and then divide by 2ϵ

The benefit of the central difference is that it limits issues around minima and maxima where the trough or peak respectively would seem much steeper if using forward or reverse. Here's a picture of this:



One benefit of the central difference method is that it is often more accurate than forward or reverse differences. However, it comes at the cost of needing an additional function evaluation/computation in many circumstances. Take, for example, the process of optimizing a function to find a maxima or minima.

Maxima-finding algorithms usually involve guessing an initial point, evaluating the function at that point, and determining what the derivative of the function is at that point. Both items are used to update the guess to one that's closer to the solution. This approach is used in many optimization algorithms such as Newton's Method. At each step you need to evaluate the function three times: for x , $x + \epsilon$, and $x - \epsilon$. With forward or reverse finite differences, you can reuse the prior function evaluation of the prior guess x . As one of the components in the estimation of the derivative, thereby saving an evaluation of the function for each iteration.

16.3. Finite Differentiation

There are additional challenges with the finite differentiation method. In practice, we are often interested in much more complex functions than x^2 . For example, we may actually be interested in the sum of a series that is many elements long or contains more complex operations than basic algebra. In the prior example, the ϵ is set unusually wide for demonstration purposes. As ϵ grow smaller generally, the accuracy of all three finite different methods increases. However, that's not always the case due to both the complexity of the function that you may be trying to differentiate or due to numerical inaccuracies of floating point math.

To demonstrate, here is a more complex example using an arbitrary function

$$f(x) = \exp(x)$$

for this example we'll show the results of the three methods calculated at different values of ϵ :

```
using DataFrames
```

```
f(x) = exp(x)
ε = 10 .^ (range(-16, stop=0, length=100))
x₀ = 1
estimate = @. (f(x₀ + ε) - f(x₀ - ε)) / 2ε
actual = f(x₀)                                ①

fig = Figure()
ax = Axis(fig[1, 1], xscale=log10,yscale=log10, xlabel="ε", ylabel="absolute error")
scatter!(ax, ε, abs.(estimate .- actual))
fig
```

- ① The derivative of $f(x) = \exp(x)$ is itself. That is $f'(x) = f(x)$ in this special case.

16. Automatic Differentiation

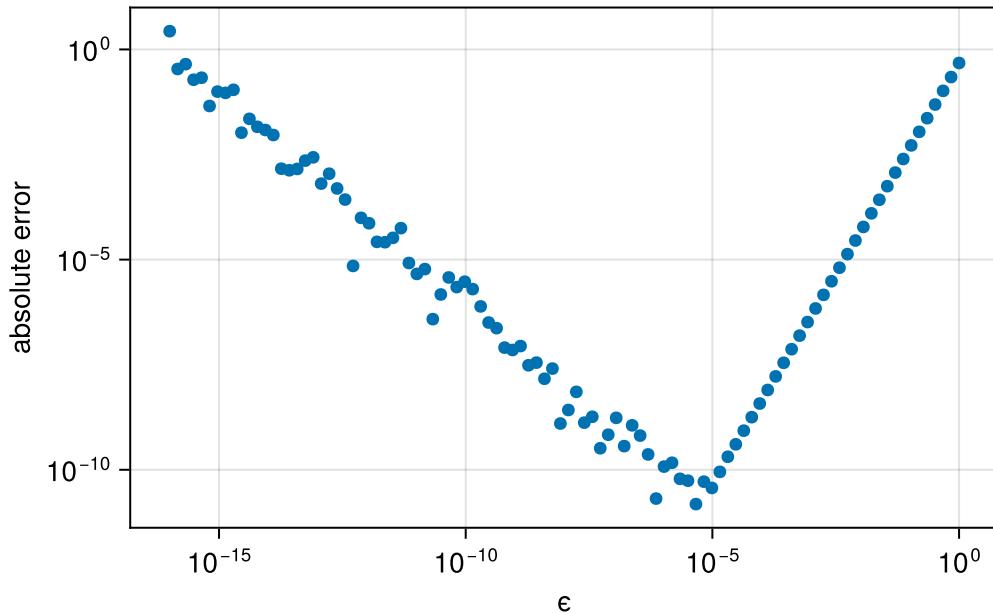


Figure 16.1.: A log-log plot showing the absolute error of the finite differences. Further to the left, roundoff error dominates while further to the right, truncation error dominates.

i Note

The `@.` in the code example above is a macro that applies broadcasting each function to its right. `@. (f(x0 + ε) - f(x0 - ε)) / 2ε` is the same as `(f.(x0 .+ ε) .- f.(x0 .- ε)) ./ (2 .* ε)`

A few observations:

1. At virtually every value of ϵ we observe some error from the true derivative.
2. That error is the sum of two parts:
 1. **Truncation error** is inherent in that we are using a given non-zero value for ϵ and not determining the limiting analytic value as $\epsilon \rightarrow 0$. The larger ϵ is, the larger the truncation error.
 2. **Roundoff error** which arises due to the limited precision of floating point math.

The implications of this are that we need to often be careful about the choice of ϵ , as the optimal choice will vary depending on the function and the point we are attempting to evaluate. This presents a number of practical difficulties in various algorithms.

16.4. Automatic Differentiation

Additionally, when computing the finite difference we must evaluate the function multiple times to determine a single estimate of the derivative. When performing something like optimization, the process typically involves iteratively making many guesses requiring many evaluations of the approximate derivative. Further, the efficiency of the algorithm usually depends on the accuracy of computing the derivative!

Despite the accuracy and computational overhead, finite differences can be very useful in many circumstances. However, a more appealing alternative approach will be covered next.

16.4. Automatic Differentiation

Automatic differentiation (“autodiff” or “AD” for short) is essentially the practice of defining algorithmically what the derivatives of function should be. We are able to do this through a creative application of the chain rule. Recall that the **chain rule** allows us to compute the derivative of a composite function using the derivatives of the component functions:

$$\begin{aligned} h(x) &= f(g(x)) \\ h'(x) &= f'(g(x))g'(x) \end{aligned}$$

Using this rule, we can define how elementary operations act when differentiated. Combined with the fact that most computer code is building up from a bunch of elementary operations, we can get a very long way in differentiating complex functions.

16.4.1. Dual Numbers

To understand where we are going, let’s remind ourselves about complex numbers. Complex numbers are of the form which has an real part (r) and an imaginary part (iq):

$$r + iq$$

By definition we say that $i^2 = -1$. This is useful because it allows us to perform certain types of operations (e.g. finding a square root of a negative number) that is otherwise unsolvable with just the real numbers¹. After defining how the normal algebraic operations (addition, multiplication, etc.) work for the imaginary number, we are able to utilize the imaginary numbers for a variety of practical mathematical tasks.

¹Richard Feynman has a wonderful, short lecture on algebra here:
https://www.feynmanlectures.caltech.edu/I_22.html

16. Automatic Differentiation

What is meant by extending the algebraic operations for imaginary numbers? For example, stating how addition should work for imaginary numbers:

$$(r + iq) + (s + iu) = (r + s) + i(q + u)$$

In a similar fashion as extending the Real (\mathbb{R}) numbers with an *imaginary* part, for automatic differentiation we will extend them with a *dual* part. A **dual number** is one of the form:

$$a + \epsilon b$$

Where $\epsilon^2 = 0$ and $\epsilon \neq 0$ by definition. While a represents the function value, b carries its derivative. An example should make this clearer. First let's define a DualNumber:

```
struct DualNumber{T,U}                                ①
    a :: T
    b :: U
    function DualNumber(a :: T, b :: U=zero(a)) where {T,U}
        return new{T,U}(a, b)
    end
end
```

- ① We define this type parametrically to handle all sorts of `<:Real` types and allow `a` and `b` to vary types in case a mathematical operation causes a type change (e.g. as in the case of integers becoming a floating point number like `10/4 == 2.5`)
- ② In the constructor, we set the default value of `b` to be `zero(a)`. `zero(a)` is a generic way to create a value equal to zero with the same type of the argument `a`. E.g. `zero(12.0) == 0.0` and `zero(12) == 0`.

Now let's define how dual numbers work under addition. The mathematical rule is:

$$(a + \epsilon b) + (c + \epsilon d) = (a + c) + (b + d)\epsilon$$

We then need to define how it works for the combinations of numbers that we might receive as arguments to our function (this is an example where multiple dispatch greatly simplifies the code compared to object oriented single dispatch!):

```
Base.:+(d :: DualNumber, e :: DualNumber) = DualNumber(d.a + e.a, d.b + e.b)
Base.:+(d :: DualNumber, x) = DualNumber(d.a + x, d.b)
Base.:+(x, d :: DualNumber) = d + x
```

And here's how we would get the derivative of a very simple function:

16.4. Automatic Differentiation

```
f1(x) = 5 + x  
f1(DualNumber(10, 1))  
  
DualNumber{Int64, Int64}(15, 1)
```

That's not super interesting though - the derivative of f_1 is just 1 and we supplied that in the construction of `DualNumber`. We did at least prove that we can add the 10 and 5!

Let's make this more interesting by also defining the multiplication operation on dual numbers. We'll follow the product rule:

$$(u \times v)' = u' \times v + u \times v'$$

```
Base.*(d::DualNumber, e::DualNumber) = DualNumber(d.a * e.a, d.b * e.a + d.a * e.b)  
Base.(x, d::DualNumber) = DualNumber(d.a * x, d.b * x)  
Base.(d::DualNumber, x) = x * d
```

Now what if we evaluate this function:

```
f2(x) = 5 + 3x  
f2(DualNumber(10, 1))  
  
DualNumber{Int64, Int64}(35, 3)
```

We have found that the second component is 3, which is indeed the derivative of $5 + 3x$ with respect to x . And in the first part we have the value of f_2 evaluated at 10.

i Note

When calculating the derivative, why do we start with 1 in the dual part of the number? Because the derivative of a variable with respect to itself is 1. From this unitary starting point, the various operations applied accumulate the derivative of the various operations in the b part of $a + \epsilon b$.

We can also define this for things like transcendental functions:

```
Base.exp(d::DualNumber) = DualNumber(exp(d.a), exp(d.a) * d.b)  
Base.sin(d::DualNumber) = DualNumber(sin(d.a), cos(d.a) * d.b)  
Base.cos(d::DualNumber) = DualNumber(cos(d.a), -sin(d.a) * d.b)  
exp(DualNumber(1, 1))
```

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```
DualNumber{Float64, Float64}(2.718281828459045, 2.718281828459045)
```

```
sin(DualNumber(0, 1))
```

```
DualNumber{Float64, Float64}(0.0, 1.0)
```

```
cos(DualNumber(0, 1))
```

```
DualNumber{Float64, Float64}(1.0, -0.0)
```

And finally, to put it all together in a more usable wrapper, we can define a function which will calculate the derivative of another function at a certain point. This function applies `f` to an initialized `DualNumber` and then returns the `b` component from the result:

```
derivative(f, x) = f(DualNumber(x, one(x))).b
```

```
derivative (generic function with 1 method)
```

And then evaluating it on a more complex function like $f(x) = 5e^{\sin(x)} + 3x$ at $x = 0$, we would analytically derive 8, which matches what we calculate next:

```
f3(x) = 5 * exp(sin(x)) + 3x
derivative(f3, 0)
```

```
8.0
```

We have demonstrated that through the clever use of dual numbers and the chain rule that complex expressions can be automatically differentiated by a computer to an exact level, limited only by the same machine precision that applies to our primary function of interest as well.

Libraries exist (such as `ChainRules.jl`) which define large numbers of predefined rules for many more operations, even beyond basic algebraic functions. This allows complex programs to be differentiated automatically.

16.5. Performance of Automatic Differentiation

Recall that in the finite difference method, we generally had to evaluate the function two or three times to *approximate* the derivative. Here we have a single function call that provides both the value and the derivative at that value. How does this compare performance-wise to simply evaluating the function a single time? Let's check how long it takes to compute a `Float64` versus a `DualNumber`:

```
using BenchmarkTools
@btime f3(rand())
```

7.042 ns (0 allocations: 0 bytes)

5.635169096968871

```
@btime f3(DualNumber(rand(), 1))
```

11.177 ns (0 allocations: 0 bytes)

`DualNumber{Float64, Float64}(11.396382856542257, 10.29150397468039)`

In performing this computation, the compiler has been able to optimize it such that we effectively are able to compute the function and its derivative at less than two times the cost of the base function evaluation. As the function gets more complex, the overhead does increase but is still a generally preferred versus finite differentiation. This advantage becomes more pronounced as we contemplate derivatives with respect to many variables at once or for higher-order derivatives.

i Note

In fact, it's largely due to the advances in applications of automatic differentiation that has led to the explosion of machine learning and artificial intelligence techniques in the 2010s/2020s. The "learning" process relies on solving parameter weights and would be too computationally expensive if using finite differences. These applications of AD in specialized C++ libraries underpin the libraries like PyTorch, Tensorflow, and Keras. These libraries specialize in allowing for AD on a limited subset of operations. Julia's available AD libraries are more general and can be applied to many more scenarios.

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16.6. Automatic Differentiation in Practice

We have, of course, not defined an exhaustive list of operations, covering only `+`, `*`, `exp`, `sin`, and `cos`. There are only a few more arithmetic (`-`, `/`) and transcendental (`log`, more trigonometric functions, etc.) before we would have a very robust set of algebraic operations defined for our `DualNumber`. In fact, it's possible to go even further and to define the behavior through conditional expressions and iterations to differentiate fairly complex functions or to extend the mechanism to partial derivatives and higher-order derivatives as well.

```
import Distributions
import ForwardDiff

N(x) = Distributions.cdf(Distributions.Normal(), x)

function d1(S, K, τ, r, σ, q)
    return (log(S / K) + (r - q + σ^2 / 2) * τ) / (σ * √(τ))
end

function d2(S, K, τ, r, σ, q)
    return d1(S, K, τ, r, σ, q) - σ * √(τ)
end

"""
eurocall(parameters)
```

Calculate the Black-Scholes implied option price for a european call where 'parameters' is a vector containing [S, K, τ, r, σ, q].

- '`S`' is the current asset price
- '`K`' is the strike or exercise price
- '`τ`' is the time remaining to maturity (can be typed with `\tau[tab]`)
- '`r`' is the continuously compounded risk free rate
- '`σ`' is the (implied) volatility (can be typed with `\sigma[tab]`)
- '`q`' is the continuously paid dividend rate

```
"""
function eurocall(parameters)
    S, K, τ, r, σ, q = parameters
    iszero(τ) && return max(zero(S), S - K)
    d₁ = d1(S, K, τ, r, σ, q)
    d₂ = d2(S, K, τ, r, σ, q)
    return (N(d₁) * S * exp(τ * (r - q)) - N(d₂) * K) * exp(-r * τ)
end
```

- ① We put the various variables inside a single `parameters` vector to allow calling a single gradient call instead of multiple derivative calls for each parameter.

```
Main.Notebook.eurocall
```

```
S = 1.0
K = 1.0
tau = 30 / 365
r = 0.05
sigma = 0.2
q = 0.0
params = [S, K, tau, r, sigma, q]
eurocall(params)
```

```
0.02493376819403728
```

💡 Tip

Some terminology in differentiation:

- **Scalar-valued function:** A function whose output is a single scalar.
- **Vector-valued (array-valued) function:** A function whose output is a vector or array.
- **Derivative (or partial derivative):** The (instantaneous) rate of change of the output with respect to one input variable. In a multivariate context, these are partial derivatives, e.g., $\frac{\partial f}{\partial x} f(x, y, z)$.
- **Gradient:** For a scalar-valued function of several variables, the gradient is the vector of all first partial derivatives, e.g. $\nabla f(x, y, z) = \left[\frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}, \frac{\partial f}{\partial z} \right]^T$.
- **Jacobian:** For a vector-valued function, the Jacobian is the matrix of first partial derivatives. If $f: \mathbb{R}^n \rightarrow \mathbb{R}^m$, its Jacobian is an $m \times n$ matrix.
- **Hessian:** For a scalar-valued function of several variables, the Hessian is the matrix of all second partial derivatives (i.e., it's an $n \times n$ matrix).

With the above code, now we can get the partial derivatives with respect to each parameter. The first, third, fourth, fifth, and sixth correspond to the common “greeks” *delta*, *theta*, *rho*, *vega*, and *epsilon* respectively. The second term is the partial derivative with respect to the strike price:

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```
ForwardDiff.gradient(eurocall, params)
```

```
6-element Vector{Float64}:
 0.5399635456230838
 -0.5150297774290467
 0.16420676980838977
 0.042331214583209334
 0.11379886104405816
 -0.04438056539367815
```

We can also get the second order greeks with another call. This includes many uncommon second order partial derivatives, but the popular *gamma* is in the [1,1] position for example:

```
ForwardDiff.hessian(eurocall, params)
```

```
6x6 Matrix{Float64}:
 6.92276   -6.92276    0.242297   0.568994   -0.0853491   -0.613375
 -6.92276    6.92276   -0.07809   -0.526663    0.199148    0.568994
 0.242297   -0.07809   -0.846846   0.521448    0.685306   -0.559878
 0.568994   -0.526663    0.521448   0.0432874   -0.0163683   -0.0467667
 -0.0853491   0.199148   0.685306   -0.0163683    0.00245525   0.007015
 -0.613375    0.568994   -0.559878   -0.0467667    0.007015   0.0504144
```

16.6.1. Performance

Earlier we assessed the impact on performance for the derivatives using `DualNumber` on a very basic function. What about if we take a more realistic example like `eurocall`? We can observe approximately a 9x slowdown when computing all of the first order derivatives which isn't bad considering we are computing 6x of the outputs!

```
@btime eurocall($params)
```

```
17.952 ns (0 allocations: 0 bytes)
```

```
0.02493376819403728
```

```
let
  g = similar(params)                                ①
  @btime ForwardDiff.gradient!($g, eurocall, $params)
end
```

16.7. Forward Mode and Reverse Mode

- ① To avoid benchmarking memory allocation for the new array, we pre-allocate the array in memory to store the result and then call `gradient!` to fill in `g` for each result.

```
259.328 ns (3 allocations: 704 bytes)
```

```
6-element Vector{Float64}:
 0.5399635456230838
 -0.5150297774290467
 0.16420676980838977
 0.042331214583209334
 0.11379886104405816
 -0.04438056539367815
```

16.7. Forward Mode and Reverse Mode

The approach of AD outlined about is called **forward mode** AD where the derivative is brought forward through the computation and accumulated through each step. The alternative to this is to first evaluate the function and then work backwards by accumulating the partial derivatives in what's called **reverse mode** AD.

Reverse mode requires more book-keeping because unlike the forward mode the derivative needs to be carried backwards, unlike the `DualNumber` approach of forward mode.

16.8. Practical tips for Automatic Differentiation

Here are a few practical tips to keep in mind.

16.8.1. Choosing between Reverse Mode and Forward Mode

Forward mode is more efficient when the number of outputs is much larger than the number inputs. When the number of inputs is much larger than the number of outputs, then reverse mode will generally be more efficient. Examples of the number of inputs being larger than the outputs might be in a statistical analysis where many features are used to predict a limited number of outcome variables or a complex model with a lot of parameters.

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16.8.2. Mutation

Auto-differentiation works through most code, but a particularly tricky part to get right is when values within arrays are mutated (changed). It's possible to do so but may require a little bit more boilerplate to setup. As of 2024, Enzyme.jl has the best support for functions with mutation inside of them.

16.8.3. Custom Rules

Custom rules for new or unusual functions can be defined, but this is an area that should be explored equipped with a bit of calculus and a deeper understanding of both forward-mode and reverse-mode. ChainRules.jl provides an interface for defining additional rules that hook into the AD infrastructure in Julia as well as provide a good set of documentation on how to extend the rules for your custom function.

16.8.4. Available Libraries in Julia

- **ForwardDiff.jl** provides robust forward-mode AD.
- **Zygote.jl** is a reverse-mode package with the innovations of being able to differentiate structs in addition to arrays and scalars.
- **Enzyme.jl** is a newer package which allows for both forward and reverse mode, but has the advantage of supporting array mutation. Additionally, Enzyme works at the level of LLVM code (an intermediate level between high level Julia code and machine code) which allows for different, sometimes better, optimizations.
- **DifferentiationInterface.jl** is a wrapper library providing a consistent API while being able exchange different backends.

In the authors' experience, they would probably recommend DifferentiationInterface.jl as a starting point, and diving into specific libraries if certain features are needed.

16.9. References

- [https://book.sciml.ai/notes/08-Forward-Mode_Automatic_Differentiation_\(AD\)_via_High_Dimension.html](https://book.sciml.ai/notes/08-Forward-Mode_Automatic_Differentiation_(AD)_via_High_Dimension.html)
- <https://blog.esciencecenter.nl/automatic-differentiation-from-scratch-23d50c699555>

17. Optimization

Yun-Tien Lee and Alec Loudenback

Far better an approximate answer to the right question, which is often vague, than an exact answer to the wrong question, which can always be made precise. - John Tukey, 1962

17.1. In This Chapter

Optimization as root finding or minimization/maximization of defined objectives. Differentiable programming and the benefits to optimization problems. Other non-gradient based optimization approaches. Model fitting as an optimization problem.

17.2. Introduction

Local and global optimization: A **local optimum** value refers to a solution where the objective function (or cost function) has the best possible value in a neighborhood surrounding that solution. A **global optimum** value, on the other hand, is the best possible value of the objective function across the entire feasible domain. For smooth and convex functions, the gradient points towards the global minimum (or maximum), making it extremely efficient for finding the optimal solution. Even for non-convex functions, the gradient provides valuable information about the direction to move towards improving the objective function value locally.

```
using CairoMakie
let

    f(x) = x^8 - 3x^4 + x

    # Generate a range of x and y values
    xs = range(-1.5, 1.5, length=300)
    ys = f.(xs)

    fig = Figure(resolution=(600, 400))
```

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```
ax = Axis(fig[1, 1],
          xlabel="x",
          ylabel="f(x)",
          title="Function with Local and Global Maxima",
          limits=(-1.5, 1.5, -5, 3)
        )

lines!(ax, xs, ys, color=:blue)

x_global_max = -1.1226

scatter!(ax,
        [x_global_max],
        [f(x_global_max)],
        color=:red,
        markersize=10,
        label="Global Minimum"
      )

x_local_min = [1.08835]

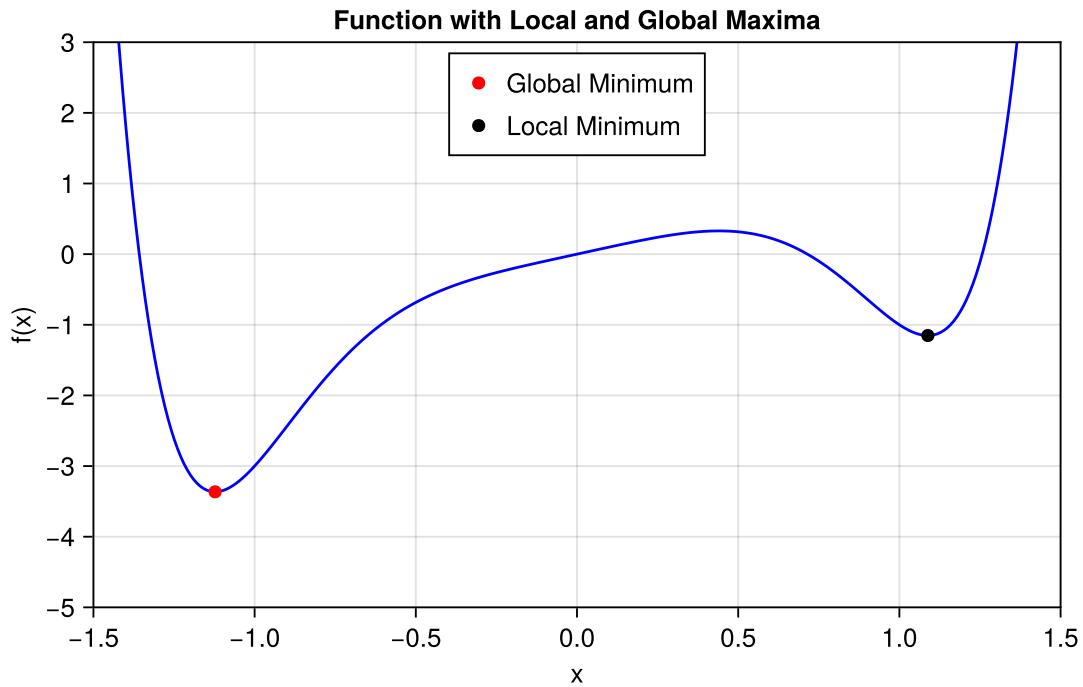
scatter!(ax,
        x_local_min,
        f.(x_local_min),
        color=:black,
        markersize=10,
        label="Local Minimum"
      )

# Add a legend
axislegend(ax, position=:ct)

fig
end
```

```
┌ Warning: Found 'resolution' in the theme when creating a 'Scene'. The 'resolution' keyword for 'S
```

```
└ @ Makie ~/julia/packages/Makie/ux0Te/src/scenes.jl:238
```



Optimization techniques are incredibly important and find uses in many areas:

- Machine learning relies on “training”, which is essentially optimizing parameters to match target data.
- Determining parameters for market consistent models where the modeled price is fit to market-observed prices.
- Minimizing a certain risk outcome by optimizing asset allocations
- Maximizing risk-adjusted yield in a portfolio.

We will introduce some basic concepts and categories of optimization techniques in this chapter:

- Objective and Loss Functions
- Optimization Techniques
 - Nonlinear Optimization
 - * Gradient Based Techniques
 - * Bracketing Methods
 - * Other Techniques
 - Linear Optimization

Below is an expanded version of the section that aims to be more thorough while keeping the writing concise and direct.

17. Optimization

17.3. Objective and Loss Functions

Any optimization algorithm needs to know what is being optimized. We call this function the **objective function**. Usually, the objective function is something we want to minimize—though maximizing a function is mathematically equivalent to minimizing its negative. Concretely, if you want to maximize $f(x)$, you can minimize $-f(x)$. The **optimal point** (also known as the $\arg \min$) is simply the input value that produces the smallest objective value.

In many practical situations—especially those with noisy or uncertain data—the function we want to optimize is called a **loss function**. This term is common in statistical modeling and machine learning, where the aim is to measure how far a model’s predictions deviate from real observations. For instance, the following squared loss function computes the sum of squared errors between a model’s predictions and the targets:

```
loss(f, guess, target) = sum((f.(guess) .- target).^2)
```

Here, `guess` might represent model parameters (or input values), and `target` is the observed data. Minimizing this loss aligns the model’s predictions with reality as closely as possible.

17.4. Optimization Techniques

17.4.1. Non-Linear Optimization

Non-linear optimization refers to problems where the relationship between the inputs and outputs is not constrained to be a linear relationship, which is incredibly common in financial modeling.

In many financial settings, variables must obey certain bounds or relationships (e.g., portfolios cannot exceed a total budget, a particular risk measure must remain below a threshold). Handling constraints often involves methods like Lagrange multipliers, projected gradient, or analytic penalty functions. We won’t cover these in depth here, but be aware that most general optimization libraries can handle constraints in non-linear problems as well.

17.4.1.1. Gradient-Based Optimization

The most efficient non-linear optimization algorithms tend to utilize the gradient (i.e. multi-variable derivative) in order to make the optimization substantially more efficient. The gradient is useful because it can tell you if you are at a maxima or minima (when the derivative is zero, the function is at a maxima or minima), but also

17.4. Optimization Techniques

because algorithms can be ‘smarter’ about searching for a solution using the additional information. The gradient provides the direction of the steepest ascent of a function. Optimization algorithms often iteratively update parameters in the direction opposite to the gradient (for minimization problems), which tends to converge towards a local minimum (or maximum for maximization problems). Besides, computing the gradient is often computationally feasible and relatively inexpensive compared to other methods for determining function behavior, such as higher-order derivatives or finite-difference methods. Beyond just the direction, the magnitude (or norm) of the gradient also indicates how steep the function change is in that direction. This information is used to adjust step sizes in optimization algorithms, balancing between convergence speed and stability.

Calculating gradients in the context of computer algorithms is discussed at length in Chapter 16, but a quick recap of the available approaches:

- Finite differences: evaluate the function at nearby points and determine the rates of change associated with each change in direction. This
- Analytic derivatives: A human-derived or computer tool (such as Mathematica) is able to analytically determine a derivative that is coded into the optimization algorithm
- Automatic differentiation (AD): elementary rules are applied to decompose elementary code operations into derivatives, allowing for very efficient computation of complex algorithms.

To compare the approaches, here is an example of determining the derivative of a simple function at a certain point:

```
using Zygote

# Define a differentiable function
f(x) = 3x^2 + 2x + 1
# Define an input value
x = 2.0

finite_diff = (f(x + 0.001) - f(x)) / 0.001

println("Value of f(x) at x=", f(x))
println("Gradient of f(x) at x=", finite_diff)                                ①
println("Gradient of f(x) at x=", gradient(f, x))                            ②
println("Gradient of f(x) at x=", 6 * x + 2)                                ③
```

- ① A finite-difference approach to determining a gradient. Note that this method required two function calls of `f` to calculate an approximation.
- ② An automatic, exact derivative determined by the rules in the Zygote library.
- ③ An analytic derivative using a user-coded derivative.

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```
Value of f(x) at x=17.0
Gradient of f(x) at x=14.002999999998877
Gradient of f(x) at x=(14.0,)
Gradient of f(x) at x=14.0
```

17.4.1.1. Root finding

Root finding, also known as root approximation or root isolation, is the process of finding the values of the independent variable (usually denoted as x) for which a given function equals zero. In mathematical terms, if we have a function $f(x)$, root finding involves finding values of x such that $f(x) = 0$.

There are various algorithms for root finding, each with its own advantages and disadvantages depending on the characteristics of the function and the requirements of the problem. One notable approach is Newton's method, an iterative method that uses the derivative or gradient of the function to approximate the root with increasing accuracy in each iteration.

We will again use a simple function to illustrate the process:

```
using Zygote

# Define a differentiable function
f(x) = 2x^2 - 3x + 1
# Define an initial value
x = 0.0
# tolerance of difference in value
tol = 1e-6
# maximum number of iteration of the algorithm
max_iter = 1000
iter = 0
while abs(f(x)) > tol && iter < max_iter
    x -= f(x) / gradient(f, x)[1]
    iter += 1
end
if iter == max_iter
    println("Warning: Maximum number of iterations reached.")
else
    println("Root found after ", iter, " iterations.")
end
print("Approximate root: ", x)
```

```
Root found after 5 iterations.
Approximate root: 0.499999998835846
```

Although it might look different, **root-finding** (solving $(g(x)=0)$) can be cast as a minimization problem by defining an objective function such as $((g(x))^2)$. In this view, driving $((g(x))^2)$ to zero compels $(g(x))$ itself to be zero, so the methods and algorithmic ideas from minimization apply naturally to root-finding scenarios as well.

17.4.1.1.2. BFGS

BFGS—named for Broyden, Fletcher, Goldfarb, and Shanno—is a popular member of the quasi-Newton family of optimization algorithms. Although it does require first-order gradient information, BFGS does not need the exact Hessian (i.e., second derivatives). Instead, it updates an approximation to the inverse Hessian at each step using the gradients from previous iterations. This extra curvature information allows BFGS to converge much faster than steepest descent in practice. Moreover, because it never explicitly forms the full Hessian matrix, it remains efficient for moderately sized problems. In finance and actuarial settings, BFGS can be especially useful for model calibration or parameter estimation tasks where one needs to handle nonlinear functions relatively quickly but cannot afford the computational overhead of second derivatives.

The following example uses the `Optim` package available in Julia to do the BFGS optimization. We don't illustrate the complete algorithm as it is a bit longer, but wanted to ensure that this workhorse of an algorithm was mentioned.

```
using Optim

# Define the objective function to minimize
function objective_function(x)
    return sum(x .* x)
end

# Initial guess for the minimization
initial_x = [1.0]
# Perform optimization using BFGS method
result = optimize(objective_function, initial_x, BFGS())
# Extract the optimized solution
solution = result.minimizer
minimum_value = result.minimum

# Print the result
println("Optimized solution: x = ", solution)
println("Minimum value found: ", minimum_value)

Optimized solution: x = [-6.359357485052897e-13]
Minimum value found: 4.0441427622698303e-25
```

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17.4.1.2. Gradient-Free Optimization

This category includes algorithms that do not rely on gradients or derivative information. They often explore the objective function using heuristics or other types of probes to guide the search.

17.4.1.2.1. Bracketing Methods

A bracketed search algorithm is a technique used in optimization and numerical methods to confine or “bracket” a minimum or maximum of a function within a specified interval. The primary goal is to reduce the search space systematically until a satisfactory solution or range containing the optimal value is found.

```
function bisection_method(f, a, b; tol=1e-6, max_iter=100)
    """
    Bisection method to find a root of the function f(x) within the interval [a, b].

    Parameters:
    - f: Function to find the root of.
    - a, b: Initial interval [a, b] where the root is expected to be.
    - tol: Tolerance for the root (default is 1e-6).
    - max_iter: Maximum number of iterations allowed (default is 100).

    Returns:
    - root: Approximate root found within the tolerance.
    - iterations: Number of iterations taken to converge.
    """
    fa = f(a)
    fb = f(b)
    if fa * fb > 0
        error("The function values at the endpoints must have opposite signs.")
    end
    iterations = 0
    while (b - a) / 2 > tol && iterations < max_iter
        c = (a + b) / 2
        fc = f(c)
        if fc == 0
            return c, iterations
        end
        if fa * fc < 0
            b = c
            fb = fc
        else
            a = c
            fa = fc
        end
        iterations += 1
    end
    return b, iterations
```

```

        a = c
        fa = fc
    end
    iterations += 1
end
root = (a + b) / 2
return root, iterations
end

# Define the function we want to find the root of
function f(x)
    return x^3 - 6x^2 + 11x - 6.1
end

# Initial interval [a, b] and tolerance
a = 0.5
b = 10
tolerance = 1e-6
# Apply the bisection method
root, iterations = bisection_method(f, a, b, tol=tolerance)

# Print results
println("Approximate root: ", root)
println("Iterations taken: ", iterations)
println("Function value at root: ", f(root))

Approximate root: 3.046680122613907
Iterations taken: 23
Function value at root: -9.356632642010254e-7

```

A popular practical algorithm is called Brent's Method, which uses additional heuristics to accelerate the optimization routine in most cases.

17.4.1.3. Other Non-Gradient Based Optimization Techniques

17.4.1.3.1. Nelder-Mead simplex method

The Nelder-Mead simplex method is a popular optimization algorithm used for minimizing (or maximizing) nonlinear functions that are not necessarily differentiable. It's particularly useful when gradient-based methods cannot be applied. It is often used in low-dimensional problems due to its simplicity and robustness. We will use the Rosenbrock function which can be useful in certain portfolio optimization problem to illustrate the process.

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```
using Optim

# Define the Rosenbrock function
function rosenbrock(v)
    x, y = v[1], v[2]
    return (1 - x)^2 + 100 * (y - x^2)^2
end

# Initial guess for (x, y)
initial_guess = [-1.5, 2.0]

# Perform optimization using the Nelder-Mead method
result = optimize(rosenbrock, initial_guess, NelderMead())

# Extract results
optimal_point = Optim.minimizer(result)
minimum_value = Optim.minimum(result)

println("Optimal Point: ", optimal_point)
println("Minimum Value: ", minimum_value)
```

```
Optimal Point: [0.9999913430783984, 0.9999847519696222]
Minimum Value: 5.016695917763382e-10
```

17.4.1.3.2. Simulated annealing

Simulated Annealing (SA) is a probabilistic optimization technique inspired by the annealing process in metallurgy. It is used to find near-optimal solutions to optimization problems, particularly in cases where traditional gradient-based methods may get stuck in local minima/maxima. SA accepts worse solutions with a certain probability, allowing it to explore the search space more broadly initially and then gradually narrow down towards better solutions as it progresses. In this section the Rastrigin function is used to illustrate the process. The function can be useful for asset modeling.

```
using Random

Random.seed!(1234)

# Parameters
max_iterations = 1000          # Number of iterations
initial_temperature = 100.0     # Starting temperature
cooling_rate = 0.99             # Cooling rate (temperature multiplier)
bounds = (-5.12, 5.12)          # Bounds for the search space
```

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```
dimension = 5 # Number of dimensions in the search space

# Objective function: Rastrigin function
function rastrigin(x)
    A = 10
    n = length(x)
    return A * n + sum(xi^2 - A * cos(2 * π * xi) for xi in x)
end

# Random initialization within bounds
function initialize_solution()
    return rand(bounds[1]:0.01:bounds[2], dimension)
end

# Random perturbation within bounds
function perturb_solution(solution)
    perturbed = copy(solution)
    index = rand(1:dimension)
    perturb_amount = rand(-0.1:0.01:0.1) # Small random change
    perturbed[index] += perturb_amount
    # Ensure perturbed solution is within bounds
    perturbed[index] = clamp(perturbed[index], bounds[1], bounds[2])
    return perturbed
end

# Simulated Annealing main function
function simulated_annealing()
    current_solution = initialize_solution()
    current_value = rastrigin(current_solution)
    best_solution = copy(current_solution)
    best_value = current_value
    temperature = initial_temperature

    for iteration in 1:max_iterations
        # Generate new candidate solution by perturbation
        candidate_solution = perturb_solution(current_solution)
        candidate_value = rastrigin(candidate_solution)

        # Acceptance probability (Metropolis criterion)
        ΔE = candidate_value - current_value
        if ΔE < 0 || rand() < exp(-ΔE / temperature)
            current_solution = candidate_solution
            current_value = candidate_value
    end
end
```

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```
end

# Update best solution found so far
if current_value < best_value
    best_solution = copy(current_solution)
    best_value = current_value
end

# Decrease temperature
temperature *= cooling_rate
if iteration % 100 == 0
    println("Iteration $iteration: Best Value = $best_value, Temperature = $temperature")
end

return best_solution, best_value
end

# Run the simulated annealing algorithm
best_solution, best_value = simulated_annealing()
println("Best Solution: ", best_solution)
println("Best Value (Minimum): ", best_value)

Iteration 100: Best Value = 79.6456061966745, Temperature = 36.60323412732294
Iteration 200: Best Value = 68.74807121999402, Temperature = 13.397967485796167
Iteration 300: Best Value = 35.612137805380314, Temperature = 4.9040894071285726
Iteration 400: Best Value = 22.630909076712427, Temperature = 1.7950553275045138
Iteration 500: Best Value = 22.116366182140844, Temperature = 0.6570483042414603
Iteration 600: Best Value = 22.11468680700349, Temperature = 0.24050092913110663
Iteration 700: Best Value = 21.995266535865557, Temperature = 0.08803111816824594
Iteration 800: Best Value = 21.93833868942773, Temperature = 0.03222223628802339
Iteration 900: Best Value = 21.918338689427735, Temperature = 0.011794380589564411
Iteration 1000: Best Value = 21.918338689427735, Temperature = 0.004317124741065788
Best Solution: [-0.9999999999999993, -1.3357370765021415e-16, -1.9799999999999993, 3.979999999999999
Best Value (Minimum): 21.918338689427735
```

17.4.1.3.3. Particle swarm optimization (PSO)

Particle swarm optimization is a metaheuristic optimization algorithm inspired by the social behavior of birds flocking or fish schooling. It is used to solve optimization problems by iteratively improving a candidate solution based on the velocity and position of particles (potential solutions) in the search space. The PSO algorithm differs from other methods in a key way, that instead of updating a single candidate solution at each

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iteration, we update a population (set) of candidate solutions, called a swarm. Each candidate solution in the swarm is called a particle. We think of a swarm as an apparently disorganized population of moving individuals that tend to cluster together while each individual seems to be moving in a random direction. The POS algorithm aims to mimic the social behavior of animals and insects.

```
using Random, CairoMakie

# Define the Rosenbrock function
function rosenbrock(x)
    return (1 - x[1])^2 + 100 * (x[2] - x[1]^2)^2
end

# PSO Implementation
function particle_swarm_optimization(objective, n_particles, n_iterations, bounds, dim)
    # Initialize particles
    positions = [rand(bounds[1]:0.1:bounds[2], dim) for _ in 1:n_particles]
    velocities = [rand(-1.0:0.1:1.0, dim) for _ in 1:n_particles]
    personal_best_positions = deepcopy(positions)
    personal_best_scores = [objective(p) for p in positions]
    global_best_position = personal_best_positions[argmin(personal_best_scores)]
    global_best_score = minimum(personal_best_scores)

    # PSO parameters
    ω = 0.5           # Inertia weight
    c1, c2 = 2.0, 2.0 # Cognitive and social learning factors

    # Optimization loop
    for iter in 1:n_iterations
        for i in 1:n_particles
            # Update velocity
            r1, r2 = rand(), rand()
            velocities[i] .= ω .* velocities[i] +
                c1 * r1 .* (personal_best_positions[i] - positions[i]) +
                c2 * r2 .* (global_best_position - positions[i])

            # Update position
            positions[i] .= positions[i] .+ velocities[i]

            # Clamp positions within bounds
            positions[i] .= clamp.(positions[i], bounds[1], bounds[2])

            # Evaluate fitness
            score = objective(positions[i])
        end
    end
end
```

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```
if score < personal_best_scores[i]
    personal_best_positions[i] = deepcopy(positions[i])
    personal_best_scores[i] = score
end

if score < global_best_score
    global_best_position = deepcopy(positions[i])
    global_best_score = score
end
end
if iter % 100 == 0
    println("Iteration $iter: Best Score = $global_best_score")
end
end
return global_best_position, global_best_score
end

# Parameters
n_particles = 30
n_iterations = 100
bounds = (-2.0, 2.0)
dim = 2

# Run PSO
best_position, best_score = particle_swarm_optimization(rosenbrock, n_particles, n_iterations,
    println("Best Position: $best_position")
    println("Best Score: $best_score")

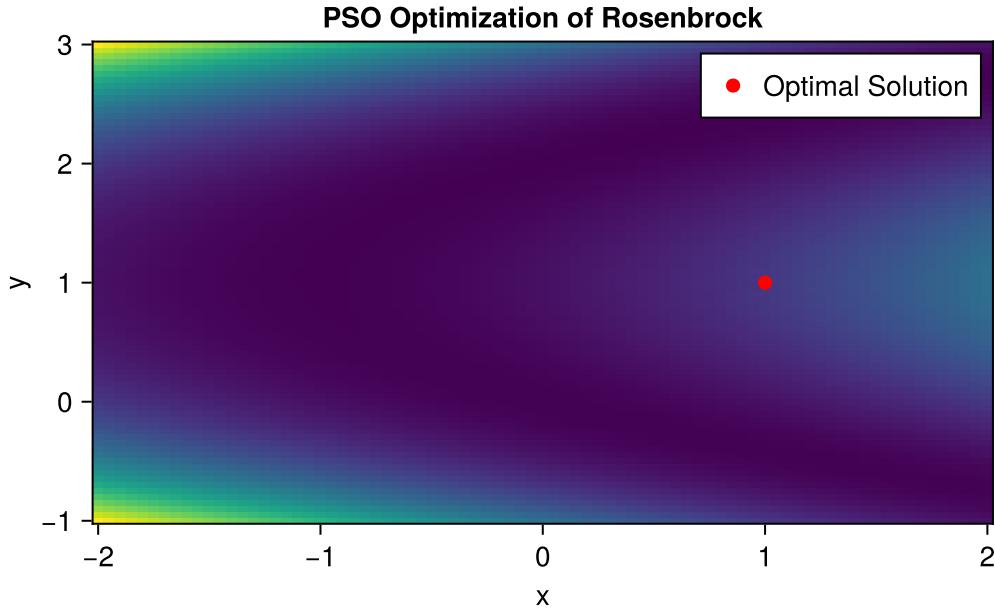
# Visualization
x = -2.0:0.05:2.0
y = -1.0:0.05:3.0
Z = [(1 - xi)^2 + 100 * (yi - xi^2)^2 for yi in y, xi in x]

# Heatmap and Scatter Plot
fig = Figure()
ax = Axis(fig[1, 1], title="PSO Optimization of Rosenbrock", xlabel="x", ylabel="y")
heatmap!(ax, x, y, Z, colormap=:viridis)
scatter!(ax, [best_position[1]], [best_position[2]], color=:red, markersize=10, label="Optimal")
axislegend(ax)
fig

Iteration 100: Best Score = 5.829480589621494e-19
```

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```
Best Position: [1.0000000001510196, 1.0000000003768819]
Best Score: 5.829480589621494e-19
```



17.4.1.3.4. Evolutionary Algorithm

An evolutionary algorithm (EA) is a family of optimization algorithms inspired by the principles of biological evolution. They are particularly useful for solving complex optimization problems where traditional gradient-based methods may struggle due to non-linearity, multimodality, or high dimensionality of the search space.

The following shows an example to maximize population fitness in terms of an objective function, with common crossover and mutation processes throughout all generations.

```
using Random

Random.seed!(1234)

# Parameters
population_size = 50          # Number of individuals in the population
chromosome_length = 5          # Number of genes in each individual (dimensionality)
generations = 100              # Number of generations
mutation_rate = 0.1             # Probability of mutation
crossover_rate = 0.7            # Probability of crossover
bounds = (-5.12, 5.12)          # Boundaries for each gene
```

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```
# Target function: Rastrigin function
function rastrigin(x)
    A = 10
    n = length(x)
    return A * n + sum(xi^2 - A * cos(2 * π * xi) for xi in x)
end

# Initialize population randomly within bounds
function initialize_population()
    return [rand(bounds[1]:0.01:bounds[2], chromosome_length) for _ in 1:population_size]
end

# Fitness function (negative because we are minimizing)
function fitness(individual)
    return -rastrigin(individual)
end

# Selection: Tournament selection
function tournament_selection(population, fitnesses)
    candidates = rand(1:population_size, 2)
    return ifelse(fitnesses[candidates[1]] > fitnesses[candidates[2]],
        population[candidates[1]], population[candidates[2]])
end

# Crossover: Single-point crossover
function crossover(parent1, parent2)
    if rand() < crossover_rate
        point = rand(1:chromosome_length)
        child1 = vcat(parent1[1:point], parent2[point+1:end])
        child2 = vcat(parent2[1:point], parent1[point+1:end])
        return child1, child2
    else
        return parent1, parent2
    end
end

# Mutation: Randomly change genes with some probability
function mutate(individual)
    for i in 1:chromosome_length
        if rand() < mutation_rate
            individual[i] = rand(bounds[1]:0.01:bounds[2])
        end
    end
end
```

```

end
return individual
end

# Main Genetic Algorithm loop
function genetic_algorithm()
    population = initialize_population()
    best_individual = nothing
    best_fitness = -Inf

    for gen in 1:generations
        # Evaluate fitness
        fitnesses = [fitness(ind) for ind in population]

        # Find best individual in current population
        current_best = argmax(fitnesses)
        if fitnesses[current_best] > best_fitness
            best_fitness = fitnesses[current_best]
            best_individual = population[current_best]
        end

        # Generate new population
        new_population = []
        while length(new_population) < population_size
            # Selection
            parent1 = tournament_selection(population, fitnesses)
            parent2 = tournament_selection(population, fitnesses)

            # Crossover
            child1, child2 = crossover(parent1, parent2)

            # Mutation
            child1 = mutate(child1)
            child2 = mutate(child2)

            # Add children to new population
            push!(new_population, child1, child2)
        end
        population = new_population[1:population_size]

        if gen % 100 == 0
            println("Generation $gen: Best Fitness = ", best_fitness)
        end
    end
end

```

17. Optimization

```
    end

    return best_individual, -best_fitness
end

# Run the genetic algorithm
best_solution, best_value = genetic_algorithm()
println("Best Solution: ", best_solution)
println("Best Value (Minimum): ", best_value)

Generation 100: Best Fitness = -2.9892957379930465
Best Solution: [-0.02, -1.02, -0.04, 0.02, -1.04]
Best Value (Minimum): 2.9892957379930465
```

17.4.1.3.5. Bayesian optimization

Bayesian Optimization (BO) is a powerful technique for global optimization of expensive-to-evaluate black-box functions. It leverages probabilistic models to predict the objective function's behavior across the search space and uses these models to make informed decisions about where to evaluate the function next. This approach efficiently balances exploration (searching for promising regions) and exploitation (exploiting regions likely to yield optimal values), making it particularly suitable for optimization problems where function evaluations are costly, such as tuning hyperparameters of machine learning models or optimizing parameters of complex simulations.

```
using Random

# Define your objective function to be optimized
function objective(x::Float64)
    return -(x^2 + 0.1 * sin(5 * x)) # Example objective function (negative because we seek minima)
end

# Bayesian optimization function
function bayesian_optimization(objective, bounds::Tuple{Float64,Float64}, num_iterations::Int)
    Random.seed!(1234) # Setting a seed for reproducibility
    X = Float64[] # List to store evaluated points
    Y = Float64[] # List to store objective values
    # Initial random point (you can choose other initial points as well)
    x_init = rand() * (bounds[2] - bounds[1]) + bounds[1]
    push!(X, x_init)
    push!(Y, objective(x_init))
    # Main loop
    for i in 1:num_iterations
        # Fit a model to the observed data (Gaussian Process in this case)
```

17.4. Optimization Techniques

```
# For simplicity, let's just use the current best observed value
x_next = rand() * (bounds[2] - bounds[1]) + bounds[1] # Random sampling
# Evaluate the objective function at the chosen point
y_next = objective(x_next)
# Update the data with the new observation
push!(X, x_next)
push!(Y, y_next)
# Here, we will just print the current best observed value
println("Iteration $i: Best value = $(maximum(Y))")

end
# Return the best observed value and corresponding parameter
best_idx = argmax(Y)
return X[best_idx], Y[best_idx]
end

best_x, best_value = bayesian_optimization(objective, (-5.0, 5.0), 10)
println("Best x found: $best_x, Best value: $best_value")

Iteration 1: Best value = -0.3041807254074535
Iteration 2: Best value = -0.3041807254074535
Iteration 3: Best value = -0.3041807254074535
Iteration 4: Best value = -0.3041807254074535
Iteration 5: Best value = -0.3041807254074535
Iteration 6: Best value = -0.3041807254074535
Iteration 7: Best value = -0.3041807254074535
Iteration 8: Best value = 0.02504980758369785
Iteration 9: Best value = 0.02504980758369785
Iteration 10: Best value = 0.02504980758369785
Best x found: -0.057501331095793695, Best value: 0.02504980758369785
```

While gradient-free methods help optimize complex non-linear or noisy functions, some financial and operational problems are best modeled with linear relationships. The next section shows how to set up such problems, including constraints, using linear and integer programming.

17.4.2. Linear optimization

Linear optimization, also known as linear programming (LP), is a mathematical method for finding the best outcome in a mathematical model with linear relationships. It involves optimizing a linear objective function subject to a set of linear equality and inequality constraints. Linear programming has a wide range of applications across various fields, including operations research, economics, engineering, and logistics.

17. Optimization

We will use linear optimization to solve the following problem, with n the number of elements in b :

$$\begin{aligned} \max_x \quad & c \cdot x \\ \text{Subject to} \quad & x \geq 0 \\ & A_i \cdot x \leq b_i \quad \forall i \in n \end{aligned}$$

```
using JuMP, GLPK, LinearAlgebra

# Define the objective coefficients
c = [1.0, 2.0, 3.0]
# Define the constraint matrix (A) and right-hand side (b)
A = [1.0 1.0 0.0;
      0.0 1.0 1.0]
b = [10.0, 20.0]
# Create a JuMP model
linear_model = Model(GLPK.Optimizer)
# Define decision variables
@variable(linear_model, x[1:3] >= 0)
# Define objective function
@objective(linear_model, Max, dot(c, x))
# Add constraints
@constraint(linear_model, constr[i=1:2], dot(A[i, :], x) <= b[i])
# Solve the optimization problem
optimize!(linear_model)

# Print results
println("Objective value: ", objective_value(linear_model))
println("Optimal solution:")
for i in 1:3
    println("\tx[$i] = ", value(x[i]))
end
```

```
WARNING: using JuMP.objective_function in module Notebook conflicts with an existing identifier.
Objective value: 70.0
Optimal solution:
x[1] = 10.0
x[2] = 0.0
x[3] = 20.0
```

17.4.2.1. Integer programming

Integer Programming (IP) is a type of optimization problem where some or all of the variables are restricted to be integers. Although the problem definition seems similar to an LP, the complexity of solving an IP hugely increases as the solution space is not continuous but discrete.

Let us use IP to solve this problem. A factory produces two types of products x_1 and x_2 with the following details:

$$\begin{aligned} \max_x & 40 \cdot x_1 + 50 \cdot x_2 \text{ Subject to } x_1 \text{ and } x_2 \text{ are integers} \\ & 4 \cdot x_1 + 3 \cdot x_2 \leq 200 \text{ (labor)} \\ & x_1 + 2 \cdot x_2 \leq 40 \text{ (material)} \end{aligned}$$

```
# Import necessary packages
using JuMP, GLPK

# Create a model with the GLPK solver
model = Model(GLPK.Optimizer)

# Define decision variables (x1 and x2 are integers)
@variable(model, x1 >= 0, Int)
@variable(model, x2 >= 0, Int)

# Define the objective function (maximize profit)
@objective(model, Max, 40 * x1 + 50 * x2)

# Add constraints
@constraint(model, 4x1 + 3x2 <= 200) # Labor constraint
@constraint(model, x1 + 4x2 <= 40) # Material constraint

# Solve the model
optimize!(model)

# Check the solution status
if termination_status(model) == MOI.OPTIMAL
    println("Optimal solution found!")
    println("x1 (Product x1 units): ", value(x1))
    println("x2 (Product x2 units): ", value(x2))
    println("Maximum Profit: ", objective_value(model))
else
    println("No optimal solution found.")
end
```

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```
Optimal solution found!
x1 (Product x1 units): 40.0
x2 (Product x2 units): 0.0
Maximum Profit: 1600.0
```

17.5. Example: Modeling fitting

In model fitting, the “best fitting curve” refers to the curve or function that best describes the relationship between the independent and dependent variables in the data. The goal of model fitting is to find the parameters of the chosen curve or function that minimize the difference between the observed data points and the values predicted by the model.

The process of finding the best fitting curve typically involves:

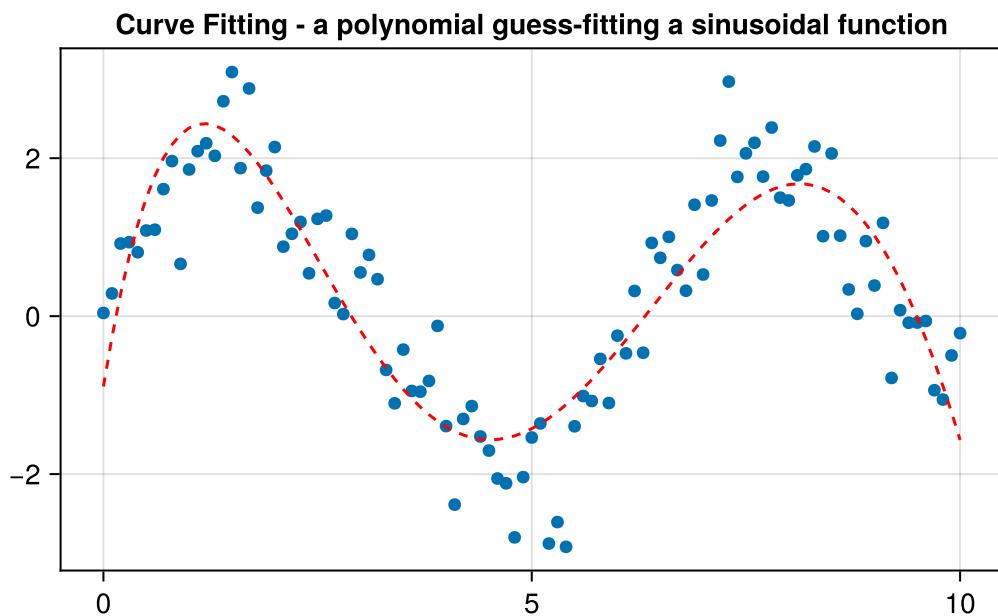
- Choosing a model: Based on the nature of the data and the underlying relationship between the variables, a suitable model or family of models are selected.
- Estimating parameters: Using the chosen model, one estimates the parameters that best describe the relationship between the variables. This is often done using optimization techniques such as least squares regression, maximum likelihood estimation, or Bayesian inference.
- Evaluating the fit: Once the parameters are estimated, one evaluates the goodness of fit of the model by comparing the predicted values to the observed data. Common metrics for evaluating fit, or error functions, include the residual sum of squares, the coefficient of determination (R-squared), and visual inspection of the residuals.
- Iterating if necessary: If the fit is not satisfactory, one may need to iterate on the model or consider alternative models until you find a satisfactory fit to the data.

While we have shown general approaches like BFGS or gradient-free schemes, libraries such as LsqFit wrap these concepts into convenient functions. Under the hood, these packages may employ gradient-based methods (including automatic differentiation) to refine parameters. Below is a demonstration:

```
using LsqFit, CairoMakie

x_data = 0:0.1:10
y_data = 2 .* sin.(x_data) .+ 0.5 .* randn(length(x_data))
# Define the model function
curve_model(x, p) = p[1] * x .^ 5 + p[2] * x .^ 4 .+ p[3] * x .^ 3 .+ p[4] * x .^ 2 .+ p[5] *
# Initial parameter guess
p₀ = [1.0, 1.0, 1.0, 1.0, 1.0]
```

```
# Fit the model to the data
fit_result = curve_fit(curve_model, x_data, y_data, p₀)
# Extract the fitted parameters
params = coef(fit_result)
# Evaluate the model with the fitted parameters
y_fit = curve_model(x_data, params)
# Plot the data and the fitted curve
fig = Figure()
Axis(fig[1, 1], title="Curve Fitting - a polynomial guess-fitting a sinusoidal function")
scatter!(x_data, y_data, label="Data")
lines!(x_data, y_fit, label="Fitted Curve", linestyle=:dash, color=:red)
fig
```



17.6. More Resources

The textbook “Algorithms For Optimization” (by Kochenderfer And Wheeler) is a comprehensive introduction to optimization and uses Julia for it’s examples.

In Julia, Optimization.jl provides a unified front-end for all kinds of general optimization problems. JuMP.jl provides a unified front-end in a specialized optimization mini-domain specific language.

18. Visualizations

Yun-Tien Lee and Alec Loudenback

Graphical excellence is that which gives to the viewer the greatest number of ideas in the shortest time with the least ink in the smallest space. - Edward Tufte, 2001

18.1. In This Chapter

The evolved brain and pattern recognition, a general guide for creating and iterating on visualizations, and principles for creating good visualizations while avoiding common mistakes.

18.2. Introduction

Visualization is a cornerstone of data analysis, statistical modeling, and decision-making. It transforms raw data into something we can see and understand, making it easier to uncover patterns, communicate ideas, and make informed decisions.

The human brain can only parse a relatively small number of textual datapoints at a single time. We are incredibly visual creatures, with our brains able to process visually many orders of magnitude more information per second than through text.

Consider the following example of tabular data, with four sets of paired x and y coordinates.

```
using DataFrames
```

```
# Define the Anscombe Quartet data
anscombe_data = DataFrame(
    x1=[10.0, 8.0, 13.0, 9.0, 11.0, 14.0, 6.0, 4.0, 12.0, 7.0, 5.0],
    y1=[8.04, 6.95, 7.58, 8.81, 8.33, 9.96, 7.24, 4.26, 10.84, 4.82, 5.68],
    x2=[10.0, 8.0, 13.0, 9.0, 11.0, 14.0, 6.0, 4.0, 12.0, 7.0, 5.0],
    y2=[9.14, 8.14, 8.74, 8.77, 9.26, 8.10, 6.13, 3.10, 9.13, 7.26, 4.74],
    x3=[10.0, 8.0, 13.0, 9.0, 11.0, 14.0, 6.0, 4.0, 12.0, 7.0, 5.0],
```

18. Visualizations

```

y3=[7.46, 6.77, 12.74, 7.11, 7.81, 8.84, 6.08, 5.39, 8.15, 6.42, 5.73],
x4=[8.0, 8.0, 8.0, 8.0, 8.0, 8.0, 19.0, 8.0, 8.0, 8.0],
y4=[6.58, 5.76, 7.71, 8.84, 8.47, 7.04, 5.25, 12.50, 5.56, 7.91, 6.89]
)

```

| | x1 | y1 | x2 | y2 | x3 | y3 | x4 | y4 |
|----|---------|---------|---------|---------|---------|---------|---------|---------|
| | Float64 |
| 1 | 10.0 | 8.04 | 10.0 | 9.14 | 10.0 | 7.46 | 8.0 | 6.58 |
| 2 | 8.0 | 6.95 | 8.0 | 8.14 | 8.0 | 6.77 | 8.0 | 5.76 |
| 3 | 13.0 | 7.58 | 13.0 | 8.74 | 13.0 | 12.74 | 8.0 | 7.71 |
| 4 | 9.0 | 8.81 | 9.0 | 8.77 | 9.0 | 7.11 | 8.0 | 8.84 |
| 5 | 11.0 | 8.33 | 11.0 | 9.26 | 11.0 | 7.81 | 8.0 | 8.47 |
| 6 | 14.0 | 9.96 | 14.0 | 8.1 | 14.0 | 8.84 | 8.0 | 7.04 |
| 7 | 6.0 | 7.24 | 6.0 | 6.13 | 6.0 | 6.08 | 8.0 | 5.25 |
| 8 | 4.0 | 4.26 | 4.0 | 3.1 | 4.0 | 5.39 | 19.0 | 12.5 |
| 9 | 12.0 | 10.84 | 12.0 | 9.13 | 12.0 | 8.15 | 8.0 | 5.56 |
| 10 | 7.0 | 4.82 | 7.0 | 7.26 | 7.0 | 6.42 | 8.0 | 7.91 |
| 11 | 5.0 | 5.68 | 5.0 | 4.74 | 5.0 | 5.73 | 8.0 | 6.89 |

Something not obvious by looking at the tabular data above is that each set of data has the same summary statistics. That is, the four sets of data are all described by the same linear features.

```

using Statistics, Printf
let d = anscombe_data
    map([[:x1, :y1], [:x2, :y2], [:x3, :y3], [:x4, :y4]]) do pair
        x, y = eachcol(d[:, pair])

        # calculate summary statistics
        mean_x, mean_y = mean(x), mean(y)
        intercept, slope = ([ones(size(y)) x] \ y)
        correlation = cor(x, y)

        (; mean_x, mean_y, intercept, slope, correlation)
    end ▷ DataFrame
end

```

| | mean_x | mean_y | intercept | slope | correlation |
|---|---------|---------|-----------|----------|-------------|
| | Float64 | Float64 | Float64 | Float64 | Float64 |
| 1 | 9.0 | 7.50091 | 3.00009 | 0.500091 | 0.816421 |
| 2 | 9.0 | 7.50091 | 3.00091 | 0.5 | 0.816237 |
| 3 | 9.0 | 7.5 | 3.00245 | 0.499727 | 0.816287 |
| 4 | 9.0 | 7.50091 | 3.00173 | 0.499909 | 0.816521 |

Analytical summarization alone is not enough to understand the data. We need to visualize the data to see the patterns emerge, wherein each of the four datasets tells a very different story:

```
using CairoMakie

# Create the plots
fig = Figure(resolution=(800, 800))

ax1 = Axis(fig[1, 1], title="Dataset 1")
scatter!(ax1, anscombe_data.x1, anscombe_data.y1)
lines!(ax1, 2:14, x → 3 + 0.5x, color=:red)

ax2 = Axis(fig[1, 2], title="Dataset 2")
scatter!(ax2, anscombe_data.x2, anscombe_data.y2)
lines!(ax2, 2:14, x → 3 + 0.5x, color=:red)

ax3 = Axis(fig[2, 1], title="Dataset 3")
scatter!(ax3, anscombe_data.x3, anscombe_data.y3)
lines!(ax3, 2:14, x → 3 + 0.5x, color=:red)

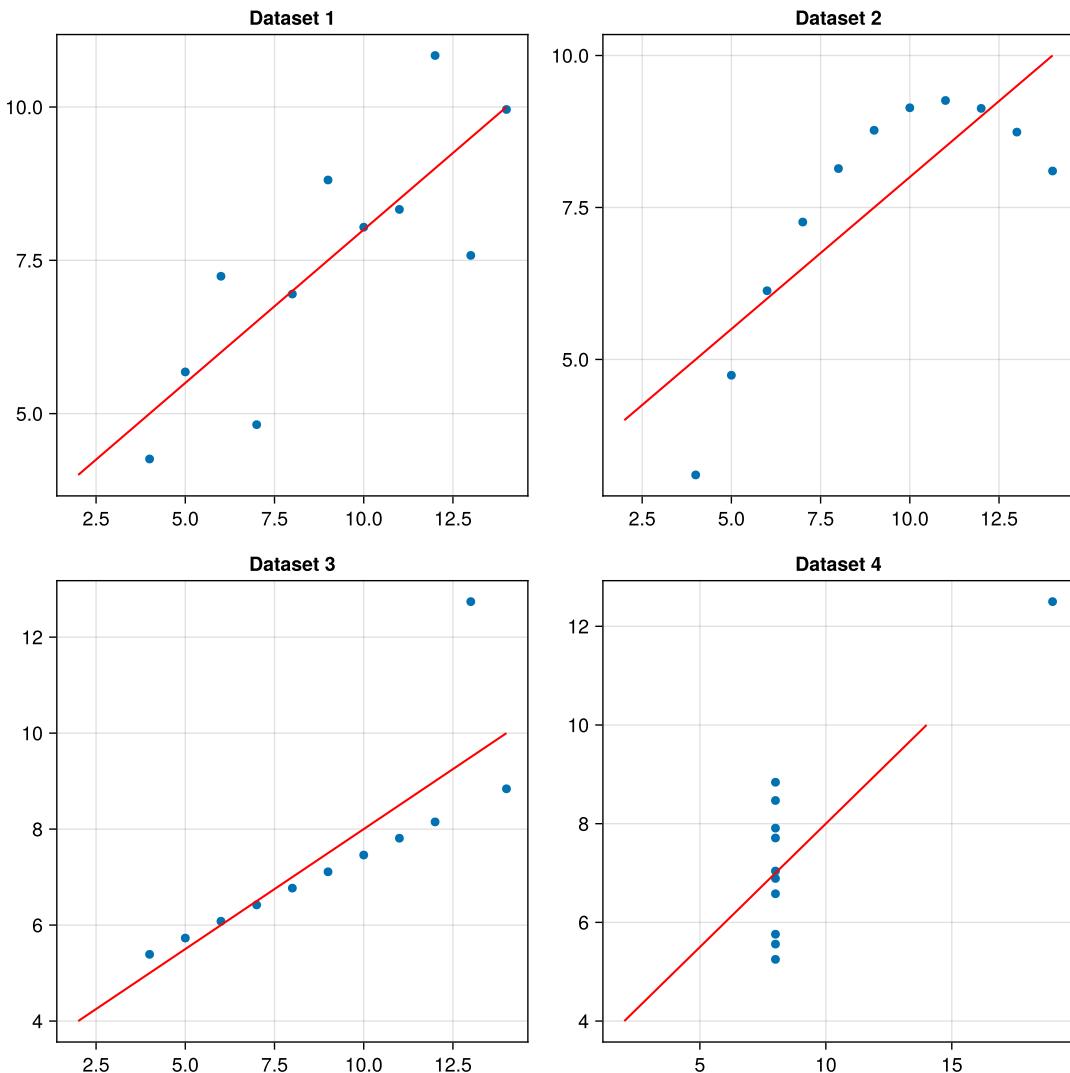
ax4 = Axis(fig[2, 2], title="Dataset 4")
scatter!(ax4, anscombe_data.x4, anscombe_data.y4)
lines!(ax4, 2:14, x → 3 + 0.5x, color=:red)

fig
```

↳ Warning: Found 'resolution' in the theme when creating a 'Scene'. The 'resolution' keyword for 'Scene'

↳ @ Makie ~/.julia/packages/Makie/ux0Te/src/scenes.jl:238

18. Visualizations



This dataset is known as Anscombe's Quartet and is a famous statistical example which is used here to underscore the importance of visualization when seeking to understand or communicate data. However, there are more reasons to refine your experience in the art and science of data visualization which we list in Table 18.1.

Table 18.1.: A list of reasons to practice the art and science of data visualization.

| Purpose | Description |
|---------------------|---|
| Simplify Complexity | Raw data can be overwhelming, especially with large datasets or many variables. A single visualization can convey thousands of points of data into a clear picture. |

| Purpose | Description |
|-----------------------------------|---|
| Reveal Patterns and Relationships | Some insights are hidden in plain sight until you visualize them, such as the relationships in the Anscombe's Quartet example. |
| Support Better Decisions | Understanding patterns and relationships can then translate into better decision making, such as highlighting trends or risks at a glance. |
| Communicate Effectively | Conveying information to others in a visual manner is one of the most effective ways at aiding in understanding. The best visualizations don't just inform—they tell a story that's useful for understanding and decision making. |
| Encourage Exploration | Visual exploration is at the heart of understanding data, uncovering distributions, relationships, or unusual patterns before diving into formal models. |

Visualization isn't just about making data look pretty—it's about making it useful. Whether you're exploring data for the first time, presenting findings to stakeholders, or refining a model, good visualizations are essential tools for turning information into insight.

18.3. Developing Visualizations

How does one develop effective visualizations? While a specialty all its own, we present the following considerations when creating quantitative displays of visual information. Consider this a guide of 'how' to create visualizations of data.

18.3.1. Define Your Message

- **Clarify the Objective:** Start with a clear purpose. What is the key insight you want to communicate? Whether it's forecasting trends, assessing risk, analyzing variances, or comparing financial scenarios, your visualization should be laser-focused on delivering that message, stripping out unnecessary details.
- **Know Your Audience:** Tailor every aspect of your visualization to the needs and expertise of your audience. For financial professionals or actuaries, this means ensuring that the visual elements align with their analytical requirements and technical proficiency.

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18.3.2. Emphasize Accuracy and Integrity

- **Maintain Consistent Scales:** Ensure axes are uniformly scaled and proportional to avoid misleading interpretations. For instance, when showing growth rates or volatility, avoid truncating or exaggerating axes. Consider using logarithmic axes when plotting growth or exponential relationships.
- Think about human perception of the shapes. For example:
 - We have a hard time comparing arc distances compared to linear distances, so pie charts are almost always a bad idea.
 - When using area to convey data (such as the size of a circle), note that the area scales quadratically, so small changes in diameter can lead to large perceptual differences in area.
- **Data-Driven Design:** Strip away unnecessary decorative elements. Every visual component should serve the core purpose of guiding interpretation based on the data.

18.3.3. Prioritize Clarity Over Complexity

- **Simplify Graphics:** Use straightforward charts, clean lines, and precise labels. Avoid embellishments that detract from the data's message such as color variation for aesthetics' sake.
- **Eliminate "Chartjunk":** Remove distracting elements like excessive gridlines, complex legends, or overly varied colors. Each element should have a clear role in supporting the narrative.
- **Leverage White Space:** Thoughtful use of white space can help separate key elements and make comparisons more intuitive.

18.3.4. Organize Data Thoughtfully

- **Decompose Complexity:** When dealing with multi-variable or time-series data, consider breaking it into small multiples or related charts for side-by-side comparison.
- **Layered Information:** Combine related datasets (e.g., financial performance vs. risk exposure), but ensure each layer is visually distinct and does not obscure others.
- **Provide Multiple Views:** Offer both high-level summaries for quick insights and detailed views for deeper analysis.

18.3.5. Enhance Readability

- **Clear Annotations:** Label axes, data points, and key takeaways explicitly. Use annotations to highlight critical insights, such as shifts in trends or activation of risk triggers.
- **Consistent Design Elements:** Stick to legible fonts and cohesive color schemes. For example, use consistent colors across charts to represent comparable data points for easier pattern recognition.

18.3.6. Validate and Iterate

- **Test for Clarity:** Share your visuals with peers or stakeholders to ensure they are interpreted as intended. Feedback can help identify areas of confusion or misrepresentation.
- **Iterate Continuously:** Treat visualization design as an evolving process. Refine layouts, scales, and annotations based on feedback and changing analytical needs.

Effective financial visualizations are built on clarity, accuracy, and thoughtful organization. By adhering to these principles—streamlined design, data integrity, and iterative refinement—you can transform complex datasets into actionable insights that empower decision-making in financial modeling and actuarial work.

 Tip

Most financial modelers are familiar with putting together plots in Excel....

18.3.7. Example: Improving a Disease Funding Visualization

We will take a visualization ([?@fig-vox-plot](#)) which has a number of issues and apply some of the principles above to improve the communication.

This example was found via (Schwarz 2016), which also identifies several of the following issues with the graphic:

- **Misleading Circle Areas:** Using circle diameters to represent values distorts perception because people intuit a comparison of area, not diameter. This results in misleading comparisons, such as the area for Breast Cancer funds appearing four times larger than Prostate Cancer, despite being only twice the value.
- **Data Dimensionality:** There are two dimensions of data conveyed: deaths and funding, while it's presented with four degrees of variation: (1) color, (2) vertical ranking, (3) horizontal categorization, and (4) bubble size.

18. Visualizations

- **Labeling Issue:** Disease names should be placed directly on circles to improve readability and assist color-blind individuals.
- **Comparison Clarity:** To effectively compare funds raised to deaths caused, these metrics should be displayed side-by-side or connected with lines.
- **Precision Overload:** Excessive precision in numerical data, such as using eight digits for funds raised, is unnecessary and can confuse interpretation.
- **Missing Data Label:** The graph omits a label for the last dollar amount, likely related to Diabetes, which could be avoided by labeling directly on the graph.

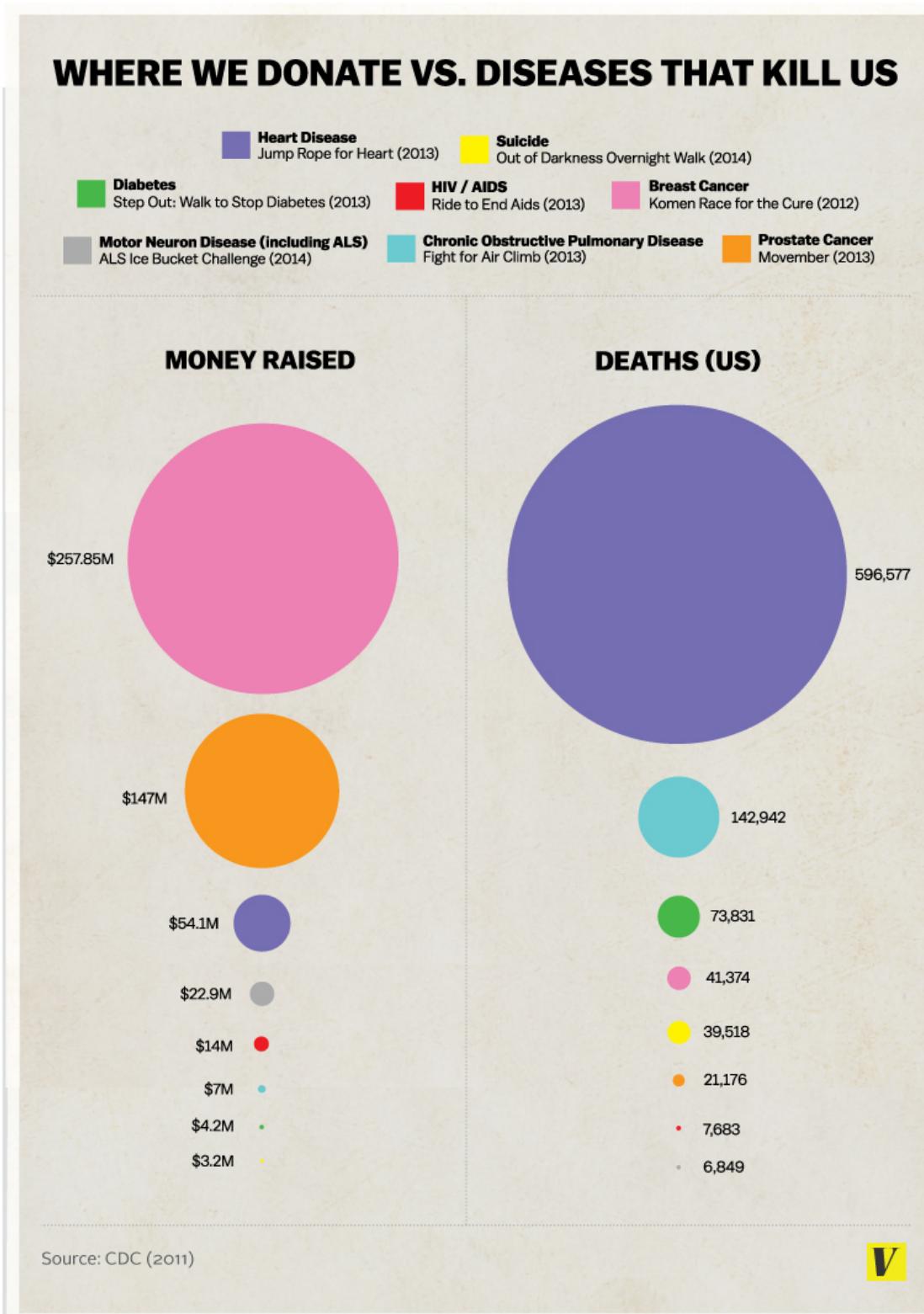


Figure 18.1.: Vox Media Infographic that inappropriately and ineffectively conveys data.
{#fig-vox-plot}

18. Visualizations

In this revised version, we take the data as accurate and simply recast the visualization of the data. The revised plot takes the following steps:

- Use a simpler 2D scatterplot mirroring the two dimensional data.
- Eliminate unnecessary color and let the labels themselves sit within the plot to avoid the eye needing to jump between the legend and the datapoints.
- Remove precision in the axis ticks, since decimal level precision is not necessary to tell the story.
- Remove unnecessary plot elements including gridlines and axes without tick labels.

```
using CairoMakie

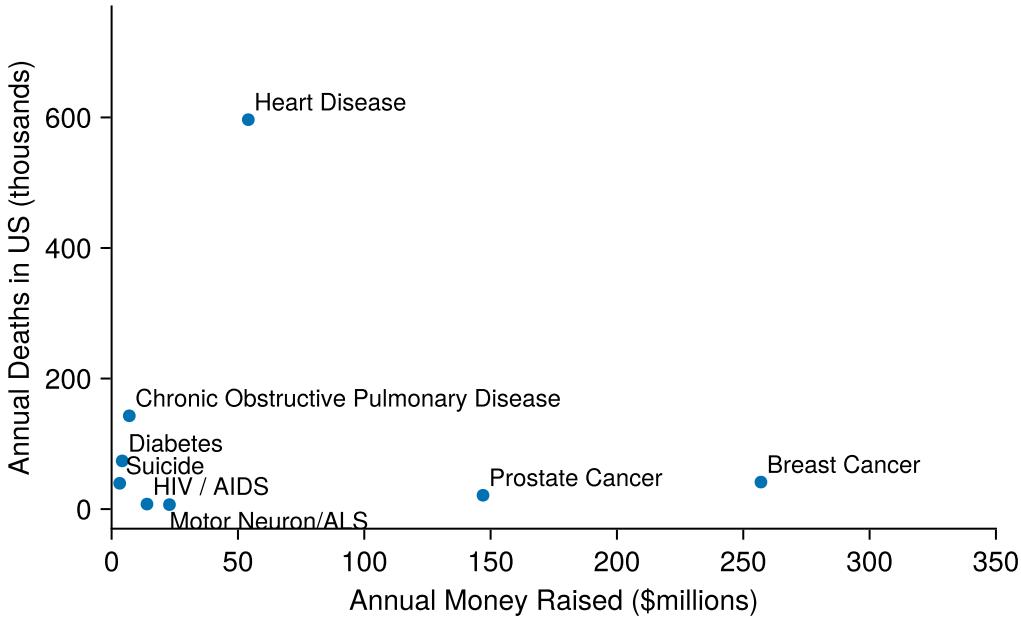
# Data
diseases = ["Breast Cancer", "Prostate Cancer", "Heart Disease", "Motor Neuron/ALS",
            "HIV / AIDS", "Chronic Obstructive Pulmonary Disease", "Diabetes", "Suicide"]
money_raised = [257, 147, 54.1, 22.9, 14, 7, 4.2, 3.2]
deaths_us = [41.374, 21.176, 596.577, 6.849, 7.683, 142.942, 73.831, 39.518]

# Create the scatter plot
fig = Figure()
ax = Axis(
    fig[1, 1],
    xlabel="Annual Money Raised (\$millions)",
    ylabel="Annual Deaths in US (thousands)",
    limits=(0, 350, -30, 770),
    xgridvisible=false,
    ygridvisible=false,
)
hidespines!(ax, :t, :r)
scatter!(ax, money_raised, deaths_us)

# Annotate each point with the disease name
for (i, disease) in enumerate(diseases)
    # avoid overlapping labels
    offset = if disease == "Motor Neuron/ALS"
              (0, -15)
            else
              (3, 2)
            end
    text!(ax, money_raised[i], deaths_us[i], text=disease, fontsize=12, offset=offset)
end
```

18.3. Developing Visualizations

```
# Display the plot  
fig
```



From the revised plot, a few key insights emerge naturally and immediately:

- The cancers receive outsized funding relative to the deaths caused.
- Heart disease remains an outsized killer compared to all other causes of death present.

And it raises some interesting questions:

- Is there an inverse relationship between the perceived “control” one has over a disease and how much funding people are willing to allocate to a cause?
- Given the wide dispersion in funding, does it matter? Ho does funding correlate with progress? E.g. has there been faster progress in extending lifespan from avoiding cancer deaths than other diseases?

The new visualization is easier to understand and draw comparisons. From the clarity, relationships are revealed and the visualizations itself reveals interesting followup questions and suggests follow-on analysis to be performed.

18.4. Principles of Good Visualization

Extending the above “how”, we now present the “what”; principles of good visualization (some elements taken from (Tufte 2001)):

- Clearly represent the data without distortions of size or space.
 - Refrain from clipping axes.
 - Do not rely on features such as shape area unless you have fully considered how viewers perceive them.
- Utilize variations of features to represent data dimensionality with purpose.
 - If colors vary in a plot, the different colors should have meaning.
 - Don’t jump to a 3D plot - use variations in marker/line styles or small multiples to convey higher dimensions.
- Encourage the eye to compare different pieces of data.
- Reveal the data at several levels of detail, from a broad overview to the fine structure.
 - Instead of summary statistics, try plotting all of the data with reduced transparency and let the viewer draw summary conclusions.
- Maintain consistency throughout the exhibit.
 - Any change in font, color, size, weight, etc. can be interpreted as an intentional choice that the viewer will try to interpret - don’t overburden the viewer.
- Serve a reasonably clear purpose: description, exploration, tabulation, or decoration and cut out what’s not purposeful.
 - Maximize the data to ink ratio.

18.5. Types of visualization tools

While not an exhaustive list by any means, we take a brief tour through some very common plots and the associated Julia code.

Note

In several of the examples, we could go further to abide by the previously listed principles of good visualizations, such as removing unnecessary gridlines or chart elements. Here, the intention is to provide a sense of how these types of plots might be constructed programmatically. Therefore, we seek not just to streamline

the plots themselves, but to also ensure that the code examples are simple and understandable.

- Basic Charts and Graphs: Bar charts, line graphs, scatter plots, histograms, pie charts.
 - Bar charts are best for comparing categorical data or discrete values across different categories. Sometimes categories can be grouped for a stacked bar chart to show for example how each category changes over time.
 - Line graphs are best for showing trends over time or when we want to highlight the rate of change. It is very intuitive to use line graphs to track trends or patterns.
 - Scatter plots are best for showing relationships or correlations between two variables. It is used a lot when one looks for patterns, clusters or outliers, or would like to explore the distribution of data points across different dimensions.
 - Histograms are best for showing the distribution of a single continuous variable, or visualizing the distribution of data points across different ranges or intervals.

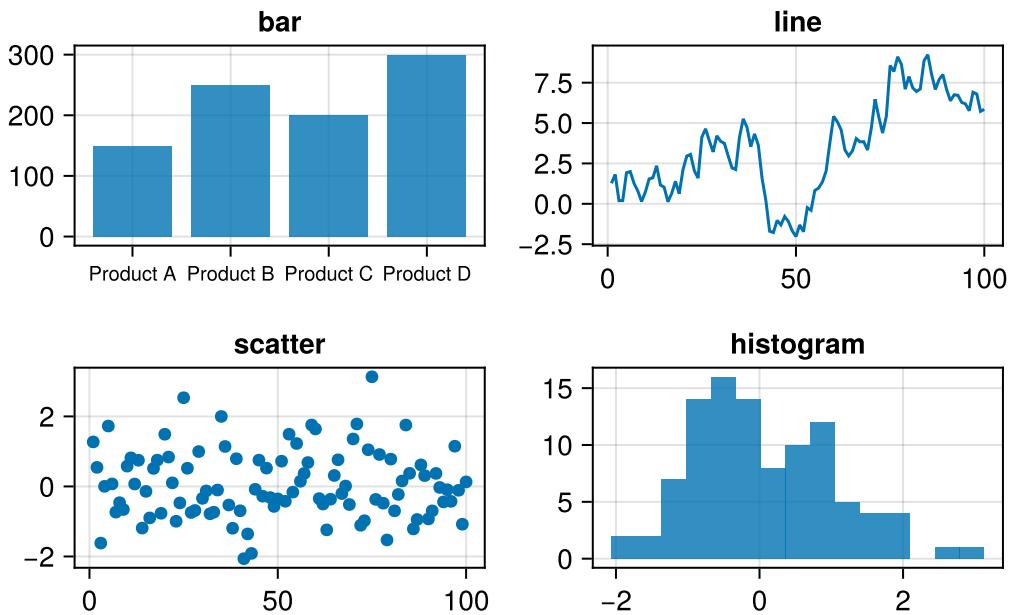
```
using Random, CairoMakie

# Data for the plots
categories = ["Product A", "Product B", "Product C", "Product D"]
sales = [150, 250, 200, 300] # For bar chart

x = randn(100) # For scatter plot

# Combine individual plots into a 2x2 layout
f = Figure()
barplot(f[1, 1], 1:4, sales, axis=(xticks=(1:4, categories), title="bar", xticklabelsize=10))
axis = Axis(f[1, 2], title="line")
lines!(f[1, 2], cumsum(x))
axis = Axis(f[2, 1], title="scatter")
scatter!(axis, x)
axis = Axis(f[2, 2], title="histogram")
hist!(axis, x)
f
```

18. Visualizations



- Multivariate Visualizations: Heatmaps, parallel coordinates plots, radar charts, bubble charts.
 - Heatmaps are best for visualizing the intensity, interactions or relationships of values across two dimensions.
 - Bubble charts which are variants of scatter plots are best for showing relationships between three variables. One can easily highlight relative importance or magnitude using the size of bubbles (e.g., revenue, population).
 - Parallel coordinates plots are best for comparing multiple variables across different observations. They are often used for detecting patterns, correlations or relationships across multiple dimensions.
 - Radar charts are best for comparing multiple variables for a single or few observations, especially when one needs to show comparisons of several quantitative variables for one or more items, with each variable represented on an axis.

```
using Random, CairoMakie

Random.seed!(1234)

# Data for plots
# For heatmap
xs = range(0, π, length=10)
ys = range(0, π, length=10)
zs = [sin(x * y) for x in xs, y in ys]
```

```

bubble_x = rand(10) * 10
bubble_y = rand(10) * 10
bubble_size = rand(10) * 100

# Dummy data for radar chart
radar_data = [0.7, 0.9, 0.4, 0.6, 0.8]

# Dummy data for parallel coordinates plot
parallel_data = rand(10, 5)

f = Figure()

# Heatmap (1,1)
ax1 = Axis(f[1, 1], title="Heatmap")
heatmap!(ax1, xs, ys, zs)

# Parallel coordinates (1,2)
ax2 = Axis(f[1, 2], title="Parallel Coordinates")
for i in 1:size(parallel_data, 2)
    lines!(ax2, 1:size(parallel_data, 2), parallel_data[i, :])
end

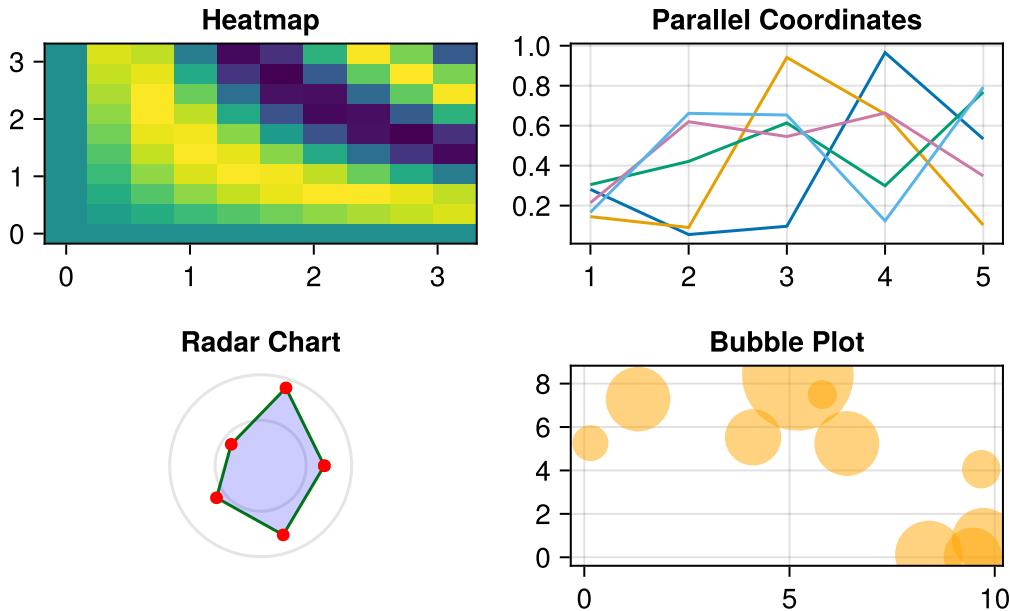
# Radar Chart (2,1)
ax3 = Axis(f[2, 1], title="Radar Chart", aspect=1)
angles = range(0, 2π, length=length(radar_data) + 1)
r = [radar_data; radar_data[1]]
arc!(ax3, Point2f(0), 0.5, -π, π, color=:grey90)
arc!(ax3, Point2f(0), 1, -π, π, color=:grey90)
lines!(ax3, cos.(angles) .* r, sin.(angles) .* r, color=:green)
poly!(ax3, cos.(angles) .* r, sin.(angles) .* r, color=(:blue, 0.2))
scatter!(ax3, cos.(angles) .* r, sin.(angles) .* r, color=:red)
hidedecorations!(ax3)
hidespines!(ax3)

# Bubble Plot (2,2)
ax4 = Axis(f[2, 2], title="Bubble Plot")
scatter!(ax4, bubble_x, bubble_y, markersize=bubble_size, color=:orange, alpha=0.5)

f

```

18. Visualizations



- Dimensionality Reduction: PCA plots, t-SNE, and UMAP for visualizing high-dimensional data. Here we show an example how high-dimensional data can be shown on a t-SNE plot.

Here we show an example to cluster synthetic stocks based on financial indicators like:

– Volatility – Momentum (6-month return) – Market Cap – P/E Ratio – Dividend Yield

t-SNE will reduce the dimensions and help us visualize clusters of stocks with similar characteristics.

```
using TSne, DataFrames, Random, Distributions, CairoMakie, StatsBase

# Generate synthetic financial dataset
Random.seed!(42)
num_stocks = 100

df = DataFrame(
    Stock=["Stock_$(i)" for i in 1:num_stocks],
    Volatility=rand(Uniform(10, 50), num_stocks), # % annualized
    Momentum=rand(Uniform(-10, 30), num_stocks), # 6-month return
    Market_Cap=1:num_stocks, # in billion USD
    P_E_Ratio=rand(Uniform(5, 50), num_stocks),
    Dividend_Yield=rand(Uniform(0, 5), num_stocks) # in %
)
```

```

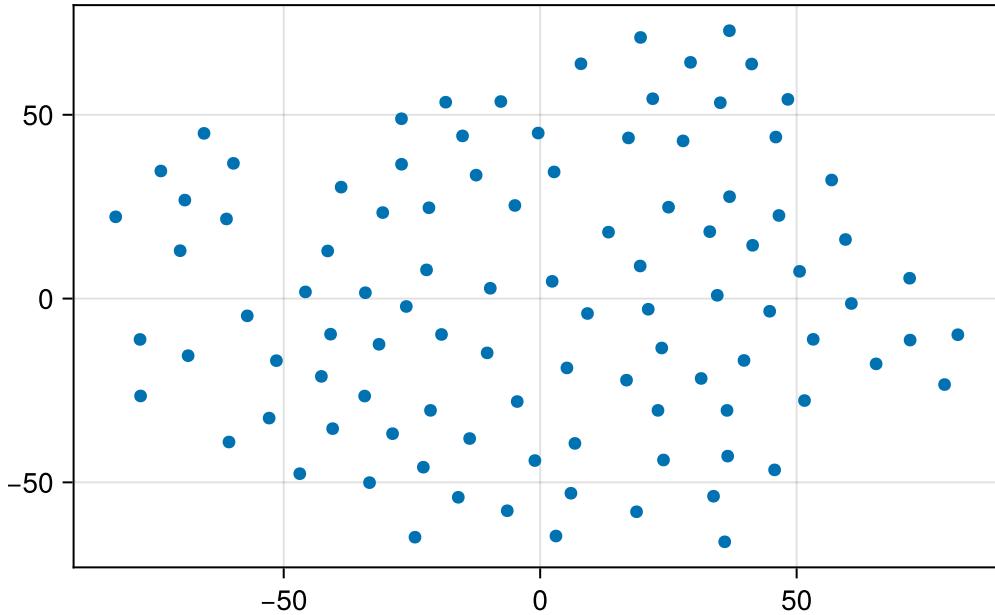
# Normalize features
features = [:Volatility, :Momentum, :Market_Cap, :P_E_Ratio, :Dividend_Yield]
X = Matrix(df[:, features])
X = StatsBase.standardize(ZScoreTransform, X, dims=1) # Standardize data

# Apply t-SNE
tsne_result = tsne(X)

# Add t-SNE components to DataFrame
df.TSNE_1 = tsne_result[:, 1]
df.TSNE_2 = tsne_result[:, 2]

# A graph of somewhat randomly distributed but also small patterns of linearity.
Makie.scatter(df.TSNE_1, df.TSNE_2)

```



- Time Series Visualization: Line charts, area charts, time-series decomposition plots. Refer to Chapter 20 for a time series plot.
- Geospatial Visualization: Maps, choropleth maps for visualizing spatial data. Here we show how spatial data can be visualized using a choropleth map.
- Interactive Dashboards: Tools like Tableau, Power BI, Pluto for interactive and dynamic data exploration.

18. Visualizations

18.5.1. Additional Examples

For more kinds of visualizations, see:

- <https://datavizproject.com>
- A Tour Through the Visualization Zoo ((Heer, Bostock, and Ogievetsky 2010))

18.6. Julia Plotting Packages

Julia has several powerful packages for data visualization, each with different strengths depending on your needs (e.g., interactive vs. static plots, ease of use vs. customization). Each package has unique strengths depending on the use case, so the best choice depends on our specific needs, the type of data, and whether we need interactive or static visualizations. Below are some of the most common visualization packages in Julia:

18.6.1. CairoMakie.jl and GLMakie.jl

Makie is designed for high-performance, interactive, and 3D visualization. It supports real-time interaction and is highly customizable. It supports 2D and 3D plotting and real-time interactivity. It is also extremely fast with GPU acceleration for certain operations. CairoMakie is suitable for print-quality vector output, while GLMakie utilizes GPU acceleration for high quality 2D and 3D plots. CairoMakie was chosen as the tool for this book because it offers very sensible default behavior and aesthetics, is easy to customize, and generally straightforward code.

18.6.2. Plots.jl

Plots is one of the most versatile and popular Julia plotting libraries. It provides a high-level interface for different plotting backends (e.g., GR, Plotly, PyPlot, PGFPlotsX, etc.). It uses a high-level syntax that is easy to use. It supports multiple backends for both static and interactive plots, but can be more limited in its customization options.

18.6.2.1. StatsPlots.jl

StatsPlots extends Plots by adding statistical plot types such as boxplots, violin plots, histograms, and density plots. It's ideal for users who frequently work with statistical data. It is specialized for statistical visualizations. It allows easy integration with Julia's statistical packages like DataFrames and StatsBase.

18.6.3. GraphPlot.jl

This package is used to plot graphs (networks), such as social network visualizations or other graph-related problems. It supports integration with the LightGraphs package for graph analytics.

18.6.4. UnicodePlots.jl

UnicodePlots provides simple plotting capabilities in the terminal using Unicode characters, making it lightweight and fast. There are no external dependencies. It is great for quick plotting within the terminal.

18.7. References

Much of the principles and some of the examples are inspired by (Tufte 2001).

19. Matrices and Their Uses

Yun-Tien Lee

“The essence of mathematics is not to make simple things complicated, but to make complicated things simple. — Stan Gudder”

19.1. In This Chapter

Matrices and their myriad uses: reframing problems through the eyes of linear algebra, an intuitive refreshing on applicable maths, and recurring patterns of matrix operations in financial modeling.

19.2. Matrix manipulation

We first review basic matrix manipulation routines before going into more advanced topics.

19.2.1. Addition and subtraction

Think of each matrix as a data grid (like a spreadsheet). Adding or subtracting values element by element is analogous to combining two sets of financial figures—such as merging cash inflows and outflows.

Example: Combining variations in A and B where each element represents a specific scenario’s impact when doing cash flow projections. We would like to see combined variations of A and B, and we also would like to know the difference in variations between A and B.

```
# Define two matrices
A = [1 2 3;
     4 5 6;
     7 8 9]
B = [9 8 7;
     6 5 4;
```

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```
3 2 1]
# Perform element-wise matrix addition and subtraction
C = A .+ B
D = A .- B
# Display the result
println("Result of matrix addition:")
println(C)
println("Result of matrix subtraction:")
println(D)

Result of matrix addition:
[10 10 10; 10 10 10; 10 10 10]
Result of matrix subtraction:
[-8 -6 -4; -2 0 2; 4 6 8]
```

19.2.2. Transpose

Transposing a matrix is akin to flipping a dataset over its diagonal—turning rows into columns. This operation is useful when aligning data for regression or matching dimensions in financial models.

Example: Converting a time-series (rows as time points) in A into a format suitable for cross-sectional analysis (columns as different variables) in B.

```
# Define a matrix
A = [1 2 3;
      4 5 6;
      7 8 9]
# Perform matrix transpose
B = A'
# Display the result
println("Result of matrix transpose:")
println(B)

Result of matrix transpose:
[1 4 7; 2 5 8; 3 6 9]
```

19.2.3. Determinant

The determinant acts as a “volume-scaling” factor. It indicates how much a linear transformation stretches or compresses space. A zero determinant signals that the transformation collapses the space into a lower dimension, implying that the matrix cannot be inverted.

Given a matrix A

$$\mathbf{A} = \begin{pmatrix} a_{11} & a_{12} & \cdots & a_{1n} \\ a_{21} & a_{22} & \cdots & a_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{n1} & a_{n2} & \cdots & a_{nn} \end{pmatrix}$$

the determinant of matrix A can be calculated as

$$\det(\mathbf{A}) = \sum_{j=1}^n (-1)^{1+j} a_{1j} \det(\mathbf{A}_{1j})$$

Example: In portfolio theory, a near-zero determinant of a covariance matrix might indicate multicollinearity among assets.

```
using LinearAlgebra
```

```
# Define a matrix
A = [1 2 3;
      5 10 20;
      7 8 9]
# Perform matrix determinant calculation
B = det(A)
# Display the result
println("Result of matrix determinant:")
println(B)
```

```
Result of matrix determinant:
30.000000000000007
```

19.2.4. Trace

The trace, being the sum of the diagonal elements, offers a quick summary that can reflect the total variance or influence of a matrix.

Example: In risk analysis, the trace of a covariance matrix may provide insights into the overall market volatility captured by the diagonal elements.

```
using LinearAlgebra
```

```
# Define a matrix
A = [1 2 3;
      5 10 20;
      7 8 9]
```

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```
# Perform matrix determinant calculation
B = tr(A)
# Display the result
println("Result of matrix trace:")
println(B)
```

```
Result of matrix trace:
20
```

19.2.5. Norm

A matrix norm measures the “size” or “energy” of the matrix. It generalizes the concept of vector length to matrices, quantifying the overall magnitude.

The Frobenius norm of a matrix is defined as:

$$\|\mathbf{A}\|_F = \sqrt{\sum_{i=1}^m \sum_{j=1}^n |A_{ij}|^2}$$

Example: A common usage of norms include error analysis, where the norm of the difference between two matrices measures how far an approximation has deviated from the true values. Another common usage in machine learning is during regularization where it allows the training process to know how large an error would be to guide the direction of updating parameters.

```
using LinearAlgebra

# Define a matrix
A = [1 2 3;
      4 5 6;
      7 8 9]
# Perform matrix norm calculation
B = norm(A)
# Display the result
println("Result of matrix norm:")
println(B)
```

```
Result of matrix norm:
16.881943016134134
```

19.2.6. Multiplication

Matrix multiplication (non-element-wise) represents the composition of linear transformations. It's like applying a sequence of financial adjustments—first transforming the data with one factor and then modifying it with another. Other applications include:

- Transforming asset returns by a matrix representing factor loadings to obtain risk contributions.
- Neural network construction. Matrix multiplication is fundamental for training and using neural networks.
- Systems of linear equations. Many real-world problems reduce to solving systems of linear equations.

```
# Define two matrices
A = [1 2 3;
      4 5 6;
      7 8 9]
B = [9 8 7;
      6 5 4;
      3 2 1]
# Perform non-element-wise matrix multiplication
C = A * B
# Display the result
println("Result of non-element-wise matrix multiplication:")
println(C)
```

Result of non-element-wise matrix multiplication:
[30 24 18; 84 69 54; 138 114 90]

On the other hand, element-wise multiplication multiplies corresponding elements directly (like applying a weight matrix).

Example: Adjusting individual cash flow items by their respective risk weights in stress testing.

```
# Define two matrices
A = [1 2 3;
      4 5 6;
      7 8 9]
B = [9 8 7;
      6 5 4;
      3 2 1]
# Perform element-wise matrix multiplication
```

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```
C = A .* B
# Display the result
println("Result of element-wise matrix multiplication:")
println(C)
```

```
Result of element-wise matrix multiplication:
[9 16 21; 24 25 24; 21 16 9]
```

19.2.7. Inversion

Matrix inversion “reverses” a transformation. If a matrix transforms one set of financial assets into another state, its inverse would bring them back.

Example: In solving linear systems for equilibrium pricing, obtaining the inverse of the coefficient matrix allows you to revert to the original asset prices.

```
# Define a matrix
A = [1 2; 3 4]
# Compute the inverse of the matrix
A_inv = inv(A)
# Display the result
println("Inverse of matrix A:")
println(A_inv)
```

```
Inverse of matrix A:
[-1.999999999999996 0.999999999999998; 1.499999999999998 -0.499999999999999]
```

i Note

For a matrix to be inverted, it must meet several important criteria.

- Square Matrix.
- The matrix must be square, meaning it has the same number of rows and columns.
- The determinant of the matrix must be non-zero.

19.3. Matrix decomposition

19.3.1. Eigenvalues

Eigenvalue decomposition, also known as eigendecomposition, is a matrix factorization that decomposes a matrix into its eigenvectors and eigenvalues. This technique uncovers the intrinsic “modes” or principal directions in a dataset. The eigenvalues indicate the

19.3. Matrix decomposition

strength of each mode, while eigenvectors show the direction or pattern associated with that strength. Eigenvalues and eigenvectors are fundamental concepts in linear algebra and play key roles include:

- Eigenvalues help in analyzing how linear transformations affect vectors in a vector space.
- Eigenvalues facilitate the diagonalization of matrices and simplify the calculations.
- In systems of differential equations, eigenvalues help determine the stability of equilibrium points.
- Identifying the main factors that cause variance in a set of asset returns, which is critical for risk management or stress testing portfolios.
- In graph theory, eigenvalues of the adjacency matrix provide insights into the properties of the graph, such as connectivity, stability, and clustering.
- Many algorithms in data science, like clustering and factorization methods, rely on eigenvalues to identify patterns and reduce dimensionality, which enhances computational efficiency and interpretability.

```
using LinearAlgebra

# Create a square matrix
A = [1 2 3;
     4 5 6;
     7 8 9]
# Perform eigenvalue decomposition
eigen_A = eigen(A)
# Extract eigenvalues and eigenvectors
λ = eigen_A.values
V = eigen_A.vectors

# Display the results
println("Original Matrix:")
println(A)
println("\nEigenvalues:")
println(λ)
println("\nEigenvectors:")
println(V)

Original Matrix:
[1 2 3; 4 5 6; 7 8 9]

Eigenvalues:
[-1.1168439698070434, -8.582743335036247e-16, 16.11684396980703]
```

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Eigenvectors:

```
[-0.7858302387420671 0.4082482904638635 -0.2319706872462857; -0.0867513392566285 -0.81649658092
```

i Note

For a matrix to get eigenvalues, it must be square, meaning it has the same number of rows and columns.

19.3.2. Singular values

Singular value decomposition (SVD) breaks a matrix into three matrices U, Σ , and V, representing the left singular vectors (analogous to the primary features), the singular values (diagonal matrix capturing the importance), and the right singular vectors (detail on how features interact), respectively. Singular values are key to:

- Matrix factorization, which simplifies many matrix operations, making it easier to analyze and manipulate data.
- Dimensionality reduction. This is particularly useful in high-dimensional data scenarios, where reducing dimensions helps eliminate noise and improve computational efficiency.
- SVD can be used for data compression, particularly in image processing.
- SVD helps filter out noise in data analysis.
- SVD provides a robust method for solving linear equations, particularly when the matrix is ill-conditioned or singular.
- In machine learning, SVD helps extract important features from datasets.
- SVD provides insights into the relationships within data. The singular values indicate the strength of the relationship, while the singular vectors offer a way to visualize and interpret those relationships.

```
using LinearAlgebra

# Create a random matrix
A = rand(4, 3)
# Perform Singular Value Decomposition (SVD)
U, Σ, V = svd(A)
# U: Left singular vectors
# Σ: Singular values (diagonal matrix)
# V: Right singular vectors (transpose)
# Reconstruct original matrix
A_reconstructed = U * Diagonal(Σ) * V'

# Display the results
```

```

println("Original Matrix:")
println(A)
println("\nLeft Singular Vectors:")
println(U)
println("\nSingular Values:")
println(Sigma)
println("\nRight Singular Vectors:")
println(V)
println("\nReconstructed Matrix:")
println(A_reconstructed)

```

Original Matrix:

[0.009965247091482832 0.6791834875425091 0.6957196130045858; 0.8609011723735872 0.5186428053028]

Left Singular Vectors:

[-0.3688251962643321 0.7049590485750195 0.6026073314440659; -0.3907787202692736 -0.708347985765]

Singular Values:

[2.0580089207591965, 0.7632647924407799, 0.46861733702594904]

Right Singular Vectors:

[-0.621674675150867 -0.7608919447091893 -0.18591408433126203; -0.4563876271441306 0.15897601468]

Reconstructed Matrix:

[0.009965247091482853 0.6791834875425096 0.6957196130045863; 0.8609011723735871 0.5186428053028]

19.3.3. Matrix factorization and fatorization machines

Matrix factorization is a popular technique in recommendation systems for modeling user-item interactions and making personalized recommendations. The core idea behind matrix factorization is to decompose the user-item interaction matrix into two lower-dimensional matrices, capturing latent factors that represent user preferences and item characteristics. By learning these latent factors, the recommendation system can make predictions for unseen user-item pairs.

Factorization Machines (FM) are a type of supervised machine learning model designed for tasks such as regression and classification, especially in the context of recommendation systems and predictive modeling with sparse data. FM models extend traditional linear models by incorporating interactions between features, allowing them to capture complex relationships within the data.

Example: In credit scoring or recommendation systems for financial products, these techniques reveal latent factors that influence customer behavior.

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```
using Recommendation, SparseArrays, MLDataUtils

# Generate synthetic user-item interaction data
num_users = 100
num_items = 50
num_ratings = 500
user_ids = rand(1:num_users, num_ratings)
item_ids = rand(1:num_items, num_ratings)
ratings = rand(1:5, num_ratings)
# Create a sparse user-item matrix
user_item_matrix = sparse(user_ids, item_ids, ratings)
# Split data into training and testing sets
train_data, test_data = splitobs(user_item_matrix, 0.8)
# Set parameters for matrix factorization
num_factors = 10
num_iterations = 10
# Train matrix factorization model
data = DataAccessor(user_item_matrix)
recommender = MF(data) # FactorizationMachines(data) alternatively
fit!(recommender)
# Predict ratings for the test set
rec = Dict()
for user in 1:num_users
    rec[user] = recommend(recommender, user, num_items, collect(1:num_items))
end
# Evaluate model performance
predictions = []
for (i, j, v) in zip(findnz(test_data.data)[1], findnz(test_data.data)[2], findnz(test_data.da
    for p in rec[i]
        if p[1] == j
            push!(predictions, p[2])
            break
        end
    end
end
end
rmse = measure(RMSE(), predictions, nonzeros(test_data.data))
println("Root Mean Squared Error (RMSE): ", rmse)
```

Root Mean Squared Error (RMSE): 1.2742717388336229

19.3.4. Principal component analysis

Principal Component Analysis (PCA) is a widely used technique in various fields for dimensionality reduction, data visualization, feature extraction, and noise reduction. PCA can also be applied to detect anomalies or outliers in the data by identifying data points that deviate significantly from the normal patterns captured by the principal components. Anomalies may appear as data points with large reconstruction errors or as outliers in the low-dimensional space spanned by the principal components.

Example: Compressing various economic indicators into a handful of principal components to illustrate predominant trends in market dynamics or risk factors.

```
using MultivariateStats
```

```
# Generate some synthetic data
data = randn(100, 5) # 100 samples, 5 features
# Perform PCA
pca_model = fit(PCA, data; maxoutdim=2) # Project to 2 principal components
# Transform the data
transformed_data = transform(pca_model, data)
# Access principal components and explained variance ratio
principal_components = pca_model.prinvars
explained_variance_ratio = pca_model.prinvars / sum(pca_model.prinvars)

# Print results
println("Principal Components:")
println(principal_components)
println("Explained Variance Ratio:")
println(explained_variance_ratio)

Principal Components:
[30.669854346985133, 24.42364238063687]
Explained Variance Ratio:
[0.5566873799754357, 0.4433126200245643]
```


20. Learning from Data

Yun-Tien Lee

“In God we trust; all others bring data.” — W. Edwards Deming

“The goal is to turn data into information, and information into insight.” — Carly Fiorina

20.1. In this chapter

We will touch on how to use data to inform a model: fitting parameters, forecasting, and fundamental limitations on prediction.

20.2. How to learn from data

20.2.1. Understand the problem and define goals

- Clarify objectives: What we want to achieve with the data (e.g., prediction, classification, clustering, or insight extraction).
- Identify key metrics: Determine how success will be measured (accuracy, RMSE, precision, etc.).
- Know the context: Understand the domain and business problem one is addressing to shape the data analysis process.

20.2.2. Collect data

Various data may be available in different formats. - Ensure data relevance: The data should be relevant to the problem. - Consider data quality: Collect data with high accuracy, completeness, and consistency.

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20.2.3. Explore and preprocess the data

This involves data cleaning and preparation to ensure the dataset is suitable for analysis.

- Handle missing data: We could impute missing values (mean, median, or KNN imputation), or drop rows/columns with excessive missing data.
- Deal with outliers: Use statistical techniques (e.g., z-scores) to detect and remove or cap extreme values.
- Feature scaling: Apply normalization or standardization to ensure features are on comparable scales (important for algorithms like SVM, K-means, etc.).
- Encode categorical data: Use techniques such as: one-hot encoding for nominal data, or label encoding or ordinal encoding for ordered categories.
- Data visualization: Use tools like `Makie.jl` to visualize distributions, correlations, and missing values.

20.2.4. Exploratory data analysis (EDA)

EDA helps discover patterns, relationships, and insights within the data. One can do the following, but not limited, to these analyses:

- Summary statistics: Check mean, variance, skewness, and correlations between variables.
- Visualize relationships: Use histograms, scatter plots, box plots, and heatmaps to identify trends and correlations.
- Detect multicollinearity: Check correlations between independent variables (e.g., Pearson's correlation matrix).

20.2.5. Select and engineer features

Feature selection and engineering help improve model performance by focusing on the most relevant information.

20.2.6. Choose the right algorithm or model

Depending on our problem type, choose appropriate algorithms for learning from the data:

- Supervised Learning (with labeled data):
 - Classification: Logistic regression, SVM, decision trees, random forests, or neural networks.

- Regression: Linear regression, ridge regression, or gradient boosting.
- Unsupervised Learning (without labeled data):
 - Clustering: K-means, DBSCAN, hierarchical clustering.
 - Dimensionality Reduction: PCA, t-SNE, or UMAP.
- Reinforcement Learning: Learn from interactions with an environment (e.g., Q-learning, Deep Q-Networks).

20.2.7. Train and evaluate the model

- Split the data: Use either a train-test split (e.g., 80/20 or 70/30 split) or a cross-validation (e.g., k-fold cross-validation).
- Fit the model: Train the model on the training set.
- Evaluate the model: Use evaluation metrics appropriate to the task.

20.2.8. Tune hyperparameters

Hyperparameters control how models learn. One can use techniques like the following to tune hyperparameters:

- Grid search: Test a range of hyperparameter values.
- Random search: Randomly explore combinations of hyperparameters.
- Bayesian optimization: Use probabilistic models to guide hyperparameter search.

20.2.9. Deploy and Monitor the Model

Once the model performs well, deploy it to make predictions on new data.

- Model deployment platforms: Use tools like Flask, FastAPI, or MLOps platforms.
- Monitor performance: Continuously monitor metrics to detect concept drift or performance degradation.

20.2.10. Draw Insights and Make Decisions

Finally, interpret the results and use insights to make decisions or recommendations. Effective communication of findings is essential, especially for stakeholders.

- Visualization: Use dashboards or reports to communicate findings.
- Interpretability: Use explainable AI (e.g., SHAP values) to make model predictions transparent.

20. Learning from Data

20.2.11. Limitations

However, there are certain fundamental limitations:

- There may often be inherent uncertainty and noise in the data itself.
- Every model has its own assumptions and simplifications.
- There may be non-stationarity in the data, especially in financial data. Non-stationary processes change over time, meaning that patterns learned from past data may no longer be valid in the future.
- Models may be overfitting or underfitting. Overfitting occurs when a model is too complex and captures noise instead of the underlying pattern, leading to poor generalization to new data. Underfitting occurs when the model is too simple to capture the relevant structure in the data.
- Sometimes in high-dimensional spaces, data becomes sparse, and meaningful patterns are harder to identify.
- Some predictions may be limited by ethical concerns (e.g., predicting criminal behavior) or legal restrictions (e.g., privacy laws that limit data collection).

20.3. Applications

20.3.1. Parameter fitting

Refer to Chapter 17 on Optimization for more details.

20.3.2. Forecasting

Forecasting is the process of making predictions about future events or outcomes based on historical data, patterns, and trends. It involves the use of statistical methods, machine learning models, or expert judgment to estimate future values in a time series or predict the likelihood of specific events. Forecasting is widely used in fields like economics, finance, meteorology, supply chain management, and business planning.

Here is an example how to do time series forecasting in Julia, where point sizes show covariance of predictions:

```
using CSV, DataFrames, CairoMakie, StateSpaceModels

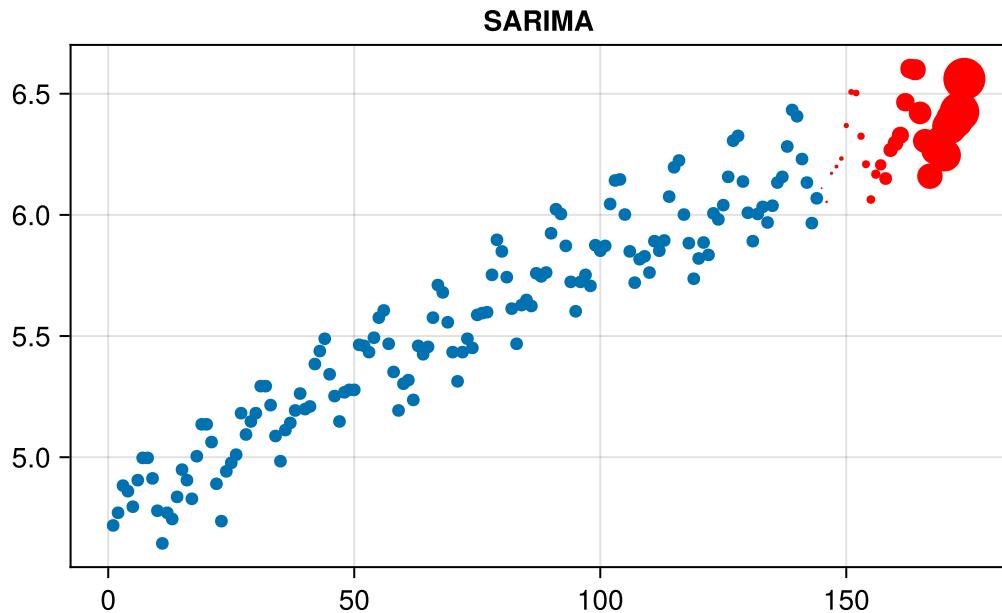
airp = CSV.read(StateSpaceModels.AIR_PASSENGERS, DataFrame)
log_air_passengers = log.(airp.passengers)
steps_existing = length(log_air_passengers)
steps_ahead = 30
```

20.3. Applications

```
# SARIMA
model_sarima = SARIMA(log_air_passengers; order=(0, 1, 1), seasonal_order=(0, 1, 1, 12))
fit!(model_sarima)
forec_sarima = forecast(model_sarima, steps_ahead)

f = Figure()
axis = Axis(f[1, 1], title="SARIMA")
scatter!(1:steps_existing, log_air_passengers)
scatter!(steps_existing+1:steps_existing+steps_ahead, map(x → x[1], forec_sarima.expected_val
f

Warning: f_tol is deprecated. Use f_abstol or f_reltol instead. The provided value (1.0e-6) will
@ Optim ~/.julia/packages/Optim/8dE7C/src/types.jl:120
```



Part VI.

Developing In Julia

Aside from the essentials necessary to convey the topics, the prior chapters have provided relatively limited insight into the *workflow* of developing in Julia. An analogy is that we have, up to this point, talked about using hammers and saws to create widgets. And some advanced topics like how to make a widget run faster and smoother in different machines. This section of the book talks about the *process* of building and designing widgets, helping you to make your widget making factory run smoothly.

This chapter is heavy with Julia-specific tips and advice. We have deliberately delayed this content until well into the book, focusing on the concepts instead of getting bogged down on language-specific details. In this section, we dive into the ‘messy’ business of building bigger, integrated things and tools to make that easier. Readers taking the concepts to other languages need not burden themselves with the details of Julia workflows and therefore can jump to the section beginning with Chapter 25.

Note

The chapters in this section are adapted from Modern Julia Workflows, originally written by G. Dalle, J. Smit, A. Hill. These chapters are a derivative of that work, and is also licensed CC BY-SA 4.0.

The content was modified to fit the other content of this book (adding cross-references, removing duplicated content) and to add/subtract elements that the authors of this book deemed more appropriate for the financial modeler.

21. Writing Julia Code

21.1. In this chapter

Installing and setting up your Julia environment. Text editor and REPL editing environments. Setting up your global environment for development. Creating packages. Logging and debugging code.

21.2. Getting help

Before you write any line of code, it's good to know where to find help. The official help page is a good place to start. In particular, the Julia community is always happy to guide beginners.

As a rule of thumb, the Discourse forum is where you should ask your questions to make the answers discoverable for future users. If you just want to chat with someone, you have a choice between the open source Zulip and the closed source Slack.

21.3. Installation

The most natural starting point to install Julia onto your system is the Julia downloads page, which will tell you to use `juliaup`.

1. Windows users can download Julia and `juliaup` together from the Windows Store.
2. OSX or Linux users can execute the following terminal command:

```
curl -fsSL https://install.julialang.org | sh
```

In both cases, this will make the `juliaup` and `julia` commands accessible from the terminal (or Windows Powershell). On Windows this will also create an application launcher. All users can start Julia by running

```
julia
```

21. Writing Julia Code

Meanwhile, `juliaup` provides various utilities to download, update, organize and switch between different Julia versions. As a bonus, you no longer have to manually specify the path to your executable. This all works thanks to adaptive shortcuts called “channels”, which allow you to access specific Julia versions without giving their exact number.

For instance, the `release` channel will always point to the current stable version, and the `lts` channel will always point to the long-term support version. Upon installation of `juliaup`, the current stable version of Julia is downloaded and selected as the default.

💡 Tip

To use other channels, add them to `juliaup` and put a `+` in front of the channel name when you start Julia:

```
juliaup add lts  
julia +lts
```

You can get an overview of the channels installed on your computer with

```
juliaup status
```

When new versions are tagged, the version associated with a given channel can change, which means a new executable needs to be downloaded. If you want to catch up with the latest developments, just do

```
juliaup update
```

21.4. REPL

The Read-Eval-Print Loop (or REPL) is the most basic way to interact with Julia, check out its documentation for details. You can start a REPL by typing `julia` into a terminal, or by clicking on the Julia application in your computer. It will allow you to play around with arbitrary Julia code:

```
julia> a, b = 1, 2;  
julia> a + b  
3
```

This is the standard (Julia) mode of the REPL, but there are three other modes you need to know. Each mode is entered by typing a specific character after the `julia>` prompt.

Once you're in a non-Julia mode, you stay there for every command you run. To exit it, hit backspace after the prompt and you'll get the `julia>` prompt back.

21.4.1. Help mode (?)

By pressing `?` you can obtain information and metadata about Julia objects (functions, types, etc.) or unicode symbols. The query fetches the docstring of the object, which explains how to use it.

```
help?> println
search: println print sprint pointer printstyled

println([io::IO], xs...)
Print (using print) xs to io followed by a newline. If io is not supplied, prints to the def

See also printstyled to add colors etc.
```

Examples

```
=====
```

```
julia> println("Hello, world")
Hello, world

julia> io = IOBuffer();

julia> println(io, "Hello", ',', " world.")
julia> String(take!(io))
"Hello, world.\n"
```

If you don't know the exact name you are looking for, type a word surrounded by quotes to see in which docstrings it pops up.

21.4.2. Package mode ()

By pressing `[]` you access `Pkg.jl`, Julia's integrated package manager, whose documentation is an absolute must-read. `Pkg.jl` allows you to:

- `]activate` different local, shared or temporary environments;
- `]instantiate` them by downloading the necessary packages;
- `]add`, `]update` (or `]up`) and `]remove` (or `]rm`) packages;

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- get the `]status` (or `]st`) of your current environment.

As an illustration, we download the package `Example.jl` inside a new environment we call `demo` (which will create an associated folder if it does not exist):

```
(demo) pkg> activate demo
Activating new project at `~/demo`

(demo) pkg> add Example
Resolving package versions...
Updating `~/demo/Project.toml'
[7876af07] + Example v0.5.5
Updating `~/demo/Manifest.toml'
[7876af07] + Example v0.5.5

(demo) pkg> status
Status `~/demo/Project.toml'
[7876af07] Example v0.5.5
```

Note that the same keywords are also available in Julia mode:

```
julia> using Pkg

julia> Pkg.rm("Example")
Updating `~/demo/Project.toml'
[7876af07] - Example v0.5.5
Updating `~/demo/Manifest.toml'
[7876af07] - Example v0.5.5
```

The package mode itself also has a help mode, accessed with `?`, in case you're lost among all these new keywords.

21.4.3. Shell mode (`;`)

By pressing `;` you enter a terminal, where you can execute any command you want, such as changing the working directory to the folder we just created:

```
shell> cd demo
/Users/myself/demo
```

21.5. Editor

In theory, any text editor suffices to write and modify Julia code. In practice, an Integrated Development Environment (or IDE) makes the experience much more pleasant, thanks to code-related utilities and language-specific plugins.

The best IDE for Julia is Visual Studio Code, or VSCode, developed by Microsoft. The Julia VSCode extension is the most feature-rich of all Julia IDE plugins. You can download it from the VSCode Marketplace and read its documentation.

 Tip

In what follows, we will sometimes mention commands and keyboard shortcuts provided by this extension. But the only shortcut you need to remember is **Ctrl + Shift + P** (or **Cmd + Shift + P** on Mac): this opens the VSCode command palette, in which you can search for any command. Type `julia` in the command palette to see what you can do.

21.6. Running code

You can execute a Julia script from your terminal, but in most cases that is not what you want to do.

```
julia myfile.jl # avoid this
```

Julia has a rather high startup and compilation latency. If you only use scripts, you will pay this cost every time you run a slightly modified version of your code. That is why many Julia developers fire up a REPL at the beginning of the day and run all of their code there, chunk by chunk, in an interactive way. Full files can be run interactively from the REPL with the `include` function.

```
julia> include("myfile.jl")
```

Alternatively, `includet` from the `Revise.jl` package can be used to “include and track” a file. This will automatically update changes to function definitions in the file in the running REPL session.

 Tip

Running code is made much easier by the following commands:

- **Julia:** Restart REPL (shortcut **Alt + J** then **Alt + R**) - this will open or restart the integrated Julia REPL. It is different from opening a plain VSCode

21. Writing Julia Code

- terminal and launching Julia manually from there.
- **Julia:** Execute Code in REPL and Move (shortcut Shift + Enter) - this will execute the selected code in the integrated Julia REPL, like a notebook.

When keeping the same REPL open for a long time, it's common to end up with a "polluted" workspace where the definitions of certain variables or functions have been overwritten in unexpected ways. This, along with other events like `struct` redefinitions, might force you to restart your REPL now and again, and that's okay.

21.7. Notebooks

Notebooks are a popular alternative to IDEs when it comes to short and self-contained code, typically in data science. They are also a good fit for literate programming, where lines of code are interspersed with comments and explanations.

The most well-known notebook ecosystem is Jupyter, which supports **Julia**, **Python** and **R** as its three core languages. To use it with Julia, you will need to install the `IJulia.jl` backend. Then, if you have also installed Jupyter with `pip install jupyterlab`, you can run this command to launch the server:

```
jupyter lab
```

If you only have `IJulia.jl` on your system, you can run this snippet instead:

```
julia> using IJulia  
julia> IJulia.notebook()
```

💡 Tip

Jupyter notebooks can be opened, modified and run directly from VSCode. Thanks to the Julia extension, you don't even need to install `IJulia.jl` or Jupyter first.

`Pluto.jl` is a newer, pure-Julia tool, adding reactivity and interactivity. It is also more amenable to version control than Jupyter notebooks because notebooks are saved as plain Julia scripts. Pluto is unique to Julia because of the language's ability to introspect and analyze dependencies in its own code. Pluto also has built-in package/environment management, meaning that Pluto notebooks contains all the code needed to reproduce results (as long as Julia and Pluto are installed).

To try out Pluto, install the package and then run

```
julia> using Pluto  
julia> Pluto.run()
```

21.8. Markdown

Markdown is a markup language used to add formatting elements to plain text content, such as Julia docstrings. Additionally, other tools such as Quarto (described below) are built using Markdown notation as the basis for their formatting, so it's useful to know about Markdown and the most essential

21.8.1. Plain Text Markdown

Plain text markdown files, which have the .md extension, are not used for interactive programming, meaning one cannot run code written in the file. As a result, plain text markdown files are usually rendered into a final product by other software.

This is an example of a plain text markdown file, including a code example contained within the "~~~~~" block:

```
# Title

## Section Header

This is example text.

```julia
println("hello world")
```
```

21.8.2. Quarto

Quarto “is an open-source scientific and technical publishing system.” Quarto makes a plain text markdown file (.md) alternative called Quarto markdown file (.qmd).

Quarto markdown files like plain text markdown files also integrate with editors, such as VSCode.



Install the Quarto extension for a streamlined experience.

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Unlike plain text markdown files, Quarto markdown files have executable code chunks. These code chunks provide a functionality similar to notebooks, thus Quarto markdown files are an alternative to notebooks. Additionally, Quarto markdown files give users additional control over output and styling via the YAML header at the top of the .qmd file.

As of Quarto version 1.5, users can choose from two Julia engines to execute code - a native Julia engine and IJulia.jl. The primary difference between the native Julia engine and IJulia.jl is that the native Julia engine does not depend on Python and can utilize local environments. For this reason it's recommended to start with the native Julia engine. Learn more about the native Julia engine in Quarto's documentation.

This book is built using Quarto documents to create the associated typeset book and website.

21.9. Environments and Dependencies

Julia comes bundled with Pkg.jl, an environment and package manager. It enables installation of packages from registries, pinning versions for compatibility, and analyzing your dependencies. Environment is meant to mean, in general, the computer you use and software installed in it. When we speak about **environments** in the Julia context, this means the Julia version and packages available to the current Julia code. For example, from the current code is a given package installed and usable?

If you open a Julia REPL, by default you will be in the *global* environment. If you hit] to enter Pkg mode, you should see:

```
(@v1.10) pkg>
```

The (@v1.10) indicates that you are using the global environment for the current Julia version (there is no global environment which applies across all Julia versions installed). You can activate a new environment with `activate [environment name]`.

```
(@v1.10) pkg> activate MyNewEnv  
Activating new project at `~/MyNewEnv`
```

This will... not do anything. Yet! When we add a package to this environment, *then* it will create a `Project.toml` and `Manifest.toml` file in that directory. Now that directory is a full fledged Julia project!

Tip

Activate a temporary environment with `activate --temp`. This will give you a temporary environment with a random name, which is very useful for testing out

things in a clean, simplified environment (the global environment, like @1.10 still applies.)

21.9.1. Project.toml

A `Project.toml` file defines attributes about the current project and its dependencies. Julia uses this to understand how to reference your current project and what dependencies it should look for from registries when instantiating the project.

i Note

TOML (Tom's Obvious Markup Language) is a modern configuration file format used to store settings and data in a human-readable, plaintext format.

This is a bit abstract, so here is a quick, annotated tour of an example `Project.toml` file:

```
name = "FinanceCore"                                     (1)
uuid = "b9b1ffdd-6612-4b69-8227-7663be06e089"          (2)
authors = ["alecloudenback <alecloudenback@users.noreply.github.com> and contributors"]
version = "2.1.0"                                         (3)

[deps]
Dates = "ade2ca70-3891-5945-98fb-dc099432e06a"
LoopVectorization = "bdcacae8-1622-11e9-2a5c-532679323890"
Roots = "f2b01f46-fcfa-551c-844a-d8ac1e96c665"

[compat]
Dates = "1"
LoopVectorization = "^0.12"
Roots = "^1.0, 2"
julia = "1.6"
```

- ① The `name` is the name of your current project which only matters if you turn your project into a package.
- ② A **UUID** is a unique identifier and can be created with Julia's UUIDs standard library.
- ③ The version follows Semantic Versioning ("SemVer") to convey to Pkg (and users!) information that ties a specific version to a specific code commit¹.
- ④ The `deps` section records the name of direct dependencies and their UUIDs so that Julia can know which packages to grab in order to make your project run.

¹When registering a package to a repository, the repository will record the version indicated in the `Project.toml` file to the git commit id of the package when it is registered.

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- ⑤ The `compat` section defines compatibility with packages can be enforced (via SemVer) to clarify which versions are allowed to be installed in case incompatibilities arise.

When you instantiate a project (see `?@sec-environment-details` for more), Julia will essentially add the packages listed under `deps`, and will **resolve** the compatible versions, generally picking the highest version number for the packages so long as the `compat` section rule are note broken.

When adding the dependencies, those packages themselves likely specify their own set of dependencies and Julia must resolve the entire **dependency graph** or **dependency tree** to allow your current project to work.

Semantic Versioning

Semantic Versioning (“SemVer”) is a scheme which uses the three-component version code to convey meaning about different versions of a package to both users and computer systems. With the version scheme `vMAJOR.MINOR.PATCH`, the meaning is roughly as follows:

1. MAJOR increments denote changes to the code which make it incompatible with prior versions.
2. MINOR increments denote changes which add features that are compatible with the prior versions.
3. PATCH increments denote changes which fix issues in prior versions and code written against the prior version is still compatible.

As an example, say we are currently using `v2.10.4` of a package, and the following theoretical options are available for us to upgrade to:

- `v2.10.5` - The 4 to 5 indicates that something may have been broken in the prior release and so we should upgrade without fear that we need to make changes to our code (unless we relied on the previously broken code!).
- `v2.11.0` - The 10 to 11 bump suggests that the new release contains some features which should not require us to change any of our previously written code.
- `v3.0.0` - The 2 to 3 indicates that we will potentially have to modify code that we have written that interfaces with this dependency.

SemVer cannot distill all possible compatibility and upgrade information about a set of packages (e.g. an author may release an update with a MINOR version which also includes fixes).

21.9.2. Manifest.toml

The `Manifest.toml` file includes a record of all external dependencies used by the project at hand. Unlike `Project.toml`, this file gets machine generated when Julia instantiates or updates the environment. The contents are basically a long list of your direct dependencies and the dependencies of those direct dependencies and looks something like this:

```
julia_version = "1.10.0"
manifest_format = "2.0"
project_hash = "5fea00df4808d89f9c977d15b8ee992bd408081b"

[[deps.AbstractFFTs]]
deps = ["LinearAlgebra"]
git-tree-sha1 = "d92ad398961a3ed262d8bf04a1a2b8340f915fef"
uuid = "621f4979-c628-5d54-868e-fcf4e3e8185c"
version = "1.5.0"
weakdeps = ["ChainRulesCore", "Test"]

[deps.AbstractFFTs.extensions]
AbstractFFTsChainRulesCoreExt = "ChainRulesCore"
AbstractFFTsTestExt = "Test"

... many more lines
```

 Note

Starting in Julia 1.11, Manifest files will include a version indication, making it nicer to work with multiple Julia versions at one time on a single system.

21.9.3. Reproducibility

Reproducibility fulfills both practical and principled goals. *Practical* in that we can record the complex chain of dependencies that is used in modern computing in order to potentially re-create a result or demonstrate an audit trail of the tools used. *Principled* in that there are circumstances (like science research) in which we want to be able to replicate results. The combination of `Project.toml` and `Manifest.toml` go a long way towards accomplishing this, as you can share both and with the same hardware and Julia version should be able to get the exact same set of dependencies and therefore run the same code. In practice, this level of reproducibility isn't *usually* needed, as most time a set of code can be run accurately without requiring the exact same set of dependencies.

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Since dependencies can have variation between systems (Windows/Mac) and architectures (x86 vs x64), you may not be able to recreate the Manifest exactly. Nevertheless, it's a fairly low bar if you are trying to maintain the utmost level of rigor around the toolchain and Julia is one of the most robust languages regarding tools to support open replication of results.

💡 Artifacts

Julia has a system called **artifacts** which allows specification of a location and hash (a cryptographic key) for data and binaries. The artifact system used to download and verify the contents of a file match the hash. This is designed for more permanent data and less end-user workflows, but we call it out here as another example where Julia takes steps to promote consistency and reproducibility.

For more on data workflows for the end-user, see Chapter 12.

💡 Tip

You can configure the environment in which a VSCode Julia REPL opens. Just click the `Julia env: ...` button at the bottom. Note however that the Julia version itself will always be the default one from `juliaup`.

21.10. Creating Local packages

Once your code base grows beyond a few scripts, you will want to create a package of your own. The first advantage is that you don't need to specify the path of every file: using `MyPackage` is enough to get access to the names you define and export (or using `MyPackage: myfunc1, myfunc2` to use bring non-exported functions into your environment). Furthermore, by structuring your project as a Pacakge, you can specify versions for your package and its dependencies, making your code easier and safer to reuse.

To create a new package locally, the easy way is to use `]generate` (we will discuss a more sophisticated workflow in the next blog post).

```
Pkg.generate(sitepath("MyPackage")); # ignore sitepath
```

This command initializes a simple folder with a `Project.toml` and a `src` subfolder. As we have seen, the `Project.toml` specifies the dependencies. Meanwhile, the `src` subfolder contains a file `MyPackage.jl`, where a module called `MyPackage` is defined. It is the heart of your package, and will typically look like this when you're done:

```
module MyPackage
```

```

# imported dependencies
using OtherPackage1
using OtherPackage2

# files defining functions, types, etc.
include("file1.jl")
include("subfolder/file2.jl")

# names you want to make public
export myfunc # e.g. defined in `file1.jl`
export MyType

end

```

21.10.1. PkgTemplates.jl

PkgTemplates.jl is like `]generate` from Pkg.jl but provides a number of options to pre-configure the repository for things such as continuous integration, testing, and compatibility. If you are not yet making use of that more advanced functionality, the `]generate` method will work just fine for you.

This will walk you through an interactive prompt to create a package in the desired folder. `~/.julia/dev` is a suggested location, but technically any folder will make do:

```

using PkgTemplates
cd("~/.julia/dev")
Template(interactive=true)("MyPkg")

```

21.11. Development workflow

Once you have created a package, your development routine might look like this:

1. Open a REPL in which you import `MyPackage`
2. Run some functions interactively, either by writing them directly in the REPL or from a Julia file that you use as a notebook
3. Modify some files in `MyPackage`
4. Go back to step 2

For that to work well, you need code modifications to be taken into account automatically. That is why Revise.jl exists. If you start every REPL session by explicitly `Revise.jl` (using `Revise`), then all the other packages you import after that will have their code

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tracked. Whenever you edit a source file and hit save, the REPL will update its state accordingly. To automatically do this for every session, see Section 21.12.

i Note

The Julia extension imports Revise.jl by default when it starts a REPL.

This is how you get started using your own package once it's set up:

```
using Revise, Pkg
Pkg.activate("./MyPackage")
using MyPackage
myfunc() # defined and exported in MyPackage
MyPackage.myfunc2() # defined and *not* exported in MyPackage
```

i Note

If you are working on a set of interrelated packages, you may need to tell those packages to use the *development* version of the package which you are modifying, instead of using the latest available from a registry. For example, say you are working on revisions to PkgA in the following dependency tree:

```
PkgB -- depends on -- > PkgA
```

If you are modifying PkgA, then you might need to tell PkgB to use the development version. For this, then you would need to:

1. Create an outer environment where you want to run the packages for interactive use while developing (say `activate @mydevenv`).
2. `]dev PkgB` which will download the associated repository into `~/.julia/dev/PkgB`
3. Go into the environment `~/.julia/dev/PkgB` and tell that environment to use the development version of PkgA with `]dev PkgB` (assuming you are modifying PkgA also in the `~/.julia/dev/` folder)

Now, in the `@mydevenv` environment, when you load `PkgB` it will load the version of PkgA

21.12. Configuration

Julia accepts startup flags to handle settings such as the number of threads available or the environment in which it launches. In addition, most Julia developers also have a

startup file which is run automatically every time the language is started. It is located at `.julia/config/startup.jl`.

The basic component that everyone puts in the startup file is `Revise.jl`. Users also commonly import packages that affect the REPL experience, as well as esthetic, benchmarking or profiling utilities. A typical example is `OhMyREPL.jl` which is widely used for syntax highlighting in the REPL. While other packages are often used, we suggest the following as a minimum:

```
# save as a file in /.julia/config/startup.jl
try
    using Revise
    using OhMyREPL
catch e
    @warn "Error with startup packages"
end
```

More generally, the startup file allows you to define your own favorite helper functions and have them immediately available in every Julia session. `StartupCustomizer.jl` can help you set up your startup file.

Tip

Here are a few more startup packages that can make your life easier once you know the language better:

- `AbbreviatedStackTraces.jl` allows you to shorten error stacktraces, which can sometimes get pretty long (beware of its interactions with VSCode)
- `Term.jl` offers a completely new way to display things like types and errors (see the advanced configuration to enable it by default).

21.13. Interactivity

The Julia REPL comes bundled with `InteractiveUtils.jl`, a bunch of very useful functions for interacting with source code.

Here are a few examples:

```
using InteractiveUtils # not necessary in a REPL session
supertypes(Int64)

(Int64, Signed, Integer, Real, Number, Any)
```

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```
subtypes(Integer)

3-element Vector{Any}:
Bool
Signed
Unsigned

methodswith(Integer)[1:5] # first five methods that take an integer argument

[1] Array(s::LinearAlgebra.UniformScaling, m::Integer, n::Integer) @ LinearAlgebra ~/.julia/juli...
[2] Float16(x::Integer) @ Base float.jl:234
[3] GenericMemoryRef(mem::GenericMemoryRef, i::Integer) @ Core boot.jl:527
[4] GenericMemoryRef(mem::GenericMemory, i::Integer) @ Core boot.jl:526
[5] Integer(x::Integer) @ Core boot.jl:925

@which exp(1) # where the currently used function is defined

exp(x::Real)
    @ Base.Math math.jl:1528

apropos("matrix exponential") # search docstrings

Base.^
Base.exp
```

When you ask for help on a Julia forum, you might want to include your local Julia information:

```
versioninfo()

Julia Version 1.11.5
Commit 760b2e5b739 (2025-04-14 06:53 UTC)
Build Info:
    Official https://julialang.org/ release
Platform Info:
    OS: macOS (arm64-apple-darwin24.0.0)
    CPU: 14 × Apple M4 Max
    WORD_SIZE: 64
    LLVM: libLLVM-16.0.6 (ORCJIT, apple-m1)
Threads: 1 default, 0 interactive, 1 GC (on 10 virtual cores)
Environment:
    JULIA_PROJECT = @.
    JULIA_LOAD_PATH = @:@stdlib
```

 Tip

The following packages can give you even more interactive power:

- InteractiveCodeSearch.jl to look for a precise implementation of a function.
- InteractiveErrors.jl to navigate through stacktraces.
- CodeTracking.jl to extend InteractiveUtils.jl

21.14. Testing

Testing in Julia primarily revolves around the built-in `Test` package, which provides a straightforward way to write and run tests using the `@test` macro. The basic syntax is simple - you write `@test expression` where the expression should evaluate to `true` for the test to pass.

To run tests in Julia, navigate to the package directory and run `Pkg.test()`, or use the `test YourPackageName` command in the Julia REPL. This will run the file in the package directory contained in `test/runtests.jl`. The test infrastructure automatically handles setting up the correct environment and dependencies for testing. Test coverage reports can be generated to see which lines of code are exercised by your tests.

`runtests.jl` is just a normal Julia file and can include other files to help organize your tests. This structure integrates naturally with Julia's package manager and testing tools.

Test organization is handled through `@testset` blocks, which group related tests together and provide summary statistics when tests are run. For example, extending the introduction to testing from Section 12.3:

```
using Test
function present_value(discount_rate, cashflows)
    v = 1.0
    pv = 0.0
    for cf in cashflows
        v = v / (1 + discount_rate)
        pv = pv + v * cf
    end
    return pv
end

@testset "Scalar Discount" begin
    @test present_value(0.05, 10) ≈ 10 / 1.05
    @test present_value(0.05, 20) ≈ 20 / 1.05
```

21. Writing Julia Code

```
end
@testset "Vector Discount" begin
    @test present_value(0.05, [10]) ≈ 10 / 1.05
    @test present_value(0.05, [10, 20]) ≈ 10 / 1.05 + 20 / 1.05^2
end;

Test Summary: | Pass  Total  Time
Scalar Discount | 2      2  0.1s
Test Summary: | Pass  Total  Time
Vector Discount | 2      2  0.0s
```

There are many more related testing facilities described in the Julia Docs, such as combining for loops with test sets.

💡 Tip

For floating-point comparisons, you'll often want to use `isapprox` (as a shorter symbol: `≈`, typed as `~`) instead of `==` to handle small numerical differences. Here's some examples:

```
@test 1/3 ≈ 0.3333333333333333  
①
@test 1/3 ≈ 0.333 atol = 1e-3  
②
@test 1/3 ≈ 0.333 rtol = 1e-33  
③
```

- ① Passes because the values are sufficiently close.
- ② Passes because the absolute difference between the values is less than 1/1000.
- ③ Passes because the difference between values is less than 1/1000 times the larger of the two values.

Here's the default behavior for `isapprox`, excerpted from its docstring:

For real or complex floating-point values, if an `atol > 0` is not specified, `rtol` defaults to the square root of `eps` of the type of `x` or `y`, whichever is bigger (least precise). This corresponds to requiring equality of about half of the significant digits. Otherwise, e.g. for integer arguments or if an `atol > 0` is supplied, `rtol` defaults to zero.

The testing workflow in Julia supports both test-driven development and continuous integration seamlessly. Tests can be run locally during development, and services like GitHub Actions can automatically run your test suite on multiple Julia versions and operating systems when you push changes.

Good testing practices in Julia involve testing edge cases, using appropriate numerical tolerances, organizing tests logically, and ensuring adequate coverage of your code's

21.14. Testing

functionality. It's also important to write tests that are clear and maintainable - each test should have a specific purpose and test one thing well.

22. Troubleshooting Julia Code

“Debugging is twice as hard as writing the code in the first place. Therefore, if you write the code as cleverly as possible, you are, by definition, not smart enough to debug it.” - Brian Kernighan

22.1. In this Chapter

Debugging in Julia involves a mix of strategies, including using print statements, the Debugger package for step-by-step inspection, logging with the Logging module, and interactive debugging with Infiltrator. These tools and techniques can help you identify and fix issues in our code efficiently.

22.2. Error Messages and Stack Traces

Julia’s error messages and stack traces can be quite informative. When an error occurs, Julia provides a traceback that shows the function call stack leading to the error, which helps in identifying where things went wrong.

```
#| error: true
function mysqrt(x)
    return sqrt(x)
end

mysqrt(-1) # This will raise a `DomainError`
```

The **stacktrace** will show us the sequence of function calls that led to the error. The print out will show the list of functions that were called (the **callstack**) which led to the code that errored. Additionally, help text is often printed, potentially offering some advice for resolving the issue. When you encounter errors in an interactive session, you can click on different parts of the stacktrace and be taken to the associated code in your editor.

22. Troubleshooting Julia Code

22.2.1. Error Types

Notice that errors are given specific types and not just result in a generic `Error`. This aids in understanding for the user: if a `DomainError` then you know that you passed the right type (e.g. a `Float64` to a function that takes a number), just that the value was not acceptable (as in the example above). Contrast that with a `MethodError` which will tell you that you've passed an invalid kind of thing to the function, not just that its value was off:

```
#| error: true
mysqrt("a string isn't OK")
```

22.3. Logging

When you encounter a problem in your code or want to track progress, a common reflex is to add `print` statements everywhere.

```
function printing_func(n)
    for i in 1:n
        println(i^2)
    end
end

printing_func (generic function with 1 method)

printing_func(3)

1
4
9
```

A slight improvement is given by the `@show` macro, which displays the variable name:

```
function showing_func(n)
    for i in 1:n
        @show i^2
    end
end

showing_func (generic function with 1 method)

showing_func(3)
```

22.3. Logging

```
i ^ 2 = 1  
i ^ 2 = 4  
i ^ 2 = 9
```

But you can go even further with the macros `@debug`, `@info`, `@warn` and `@error`. They have several advantages over printing:

- They display variable names and a custom message
- They show the line number they were called from
- They can be disabled and filtered according to source module and severity level
- They work well in multithreaded code
- They can write their output to a file

```
function warning_func(n)  
    for i in 1:n  
        @warn "This is bad" i^2  
    end  
end
```

```
warning_func (generic function with 1 method)
```

```
warning_func(3)
```

```
[ Warning: This is bad  
  i ^ 2 = 1  
  @ Main.Notebook ~/prog/julia-fin-book/julia-debugging.qmd:75  
[ Warning: This is bad  
  i ^ 2 = 4  
  @ Main.Notebook ~/prog/julia-fin-book/julia-debugging.qmd:75  
[ Warning: This is bad  
  i ^ 2 = 9  
  @ Main.Notebook ~/prog/julia-fin-book/julia-debugging.qmd:75
```

Refer to the logging documentation for more information.

i Note

In particular, note that `@debug` messages are suppressed by default. You can enable them through the `JULIA_DEBUG` environment variable if you specify the source module name, typically `Main` or your package module.

22. Troubleshooting Julia Code

Beyond the built-in logging utilities, ProgressLogging.jl has a macro `@progress`, which interfaces nicely with VSCode and Pluto to display progress bars. And Suppressor.jl can sometimes be handy when you need to suppress warnings or other bothersome messages (use at your own risk).

22.4. Commonly Encountered Macros

Aside from those mentioned in the context of Logging, there are a number of different useful macros, many of which are highlighted in the following table:

Table 22.1.: Useful macros for modeling work. There are others related to parallelism which will be covered in Chapter 11.

| Macro | Description |
|--|--|
| <code>BenchmarkTools.@benchmark</code> | Runs the given expression multiple times, collecting timing and memory allocation statistics. Useful for benchmarking and performance analysis. |
| <code>BenchmarkTools.@btime</code> | Similar to <code>@benchmark</code> , but focuses on the minimum execution time and provides a more concise output. |
| <code>@edit</code> | Opens the source code of a function or module in an editor for inspection or modification. |
| <code>@which</code> | Displays the method that would be called for a given function call, helping to understand method dispatch. |
| <code>@code_warntype</code> | Shows the type inference results for a given function call, highlighting any type instabilities or performance issues. |
| <code>@info, @warn, @error</code> | Used for logging messages at different severity levels (info, warning, error) during program execution. |
| <code>@assert</code> | Asserts that a given condition is true, throwing an error if the condition is false. Useful for runtime checks and debugging. |
| <code>@view, @views</code> | Access a subset of an array without copying the data in that slice. <code>@views</code> applies to all array slicing operations within the expressions that follow it. |
| <code>Test.@test,</code>
<code>Test.@testset</code> | Used for defining unit tests. <code>@test</code> checks that a condition is true, while <code>@testset</code> groups related tests together. |
| <code>@raw</code> | Encloses a string literal, disabling string interpolation and escape sequences. Useful for writing raw string data. This is especially helpful when working with filepaths where the \ in Windows paths otherwise needs to be escaped with a leading slash (e.g. \\). |

| Macro | Description |
|-----------|---|
| @fastmath | Enables aggressive floating-point optimizations within a block, potentially sacrificing strict IEEE compliance for performance. |
| @inbounds | Disables bounds checking for array accesses within a block, improving performance but removing safety checks. |
| @inline | Suggests to the compiler that a function should be inlined at its call sites, potentially improving performance by reducing function call overhead. |

22.5. Debugging

The limitation of printing or logging is that you cannot interact with local variables or save them for further analysis. The following two packages solve this issue (consider adding to your default environment @v1.X, like Revise.jl).

22.5.1. Setting

Assume you want to debug a function checking whether the n -th Fermat number $F_n = 2^{2^n} + 1$ is prime:

```
function fermat_prime(n)
    k = 2^n
    F = 2^k + 1
    for d in 2:isqrt(F) # integer square root
        if F % d == 0
            return false
        end
    end
    return true
end

fermat_prime (generic function with 1 method)

fermat_prime(4), fermat_prime(6)

(true, true)
```

Unfortunately, $F_4 = 65537$ is the largest known Fermat prime, which means F_6 is incorrectly classified. Let's investigate why this happens!

22. Troubleshooting Julia Code

22.5.2. Infiltrator.jl

Infiltrator.jl is a lightweight inspection package, which will not slow down your code at all. Its `@infiltrate` macro allows you to directly set breakpoints in your code. Calling a function which hits a breakpoint will activate the Infiltrator REPL-mode and change the prompt to `infil>`. Typing `?` in this mode will summarize available commands. For example, typing `@locals` in Infiltrator-mode will print local variables:

```
using Infiltrator

function fermat_prime_infil(n)
    k = 2^n
    F = 2^k + 1
    @infiltrate
    for d in 2:isqrt(F)
        if F % d == 0
            return false
        end
    end
    return true
end
```

What makes Infiltrator.jl even more powerful is the `@exfiltrate` macro, which allows you to move local variables into a global storage called the `safehouse`.

```
julia> fermat_prime_infil(6)
Infiltrating fermat_prime_infil(n::Int64)
  at REPL[2]:4

infil> @exfiltrate k F
Exfiltrating 2 local variables into the safehouse.

infil> @continue

true

julia> safehouse.k
64

julia> safehouse.F
1
```

The diagnosis is a classic one: integer overflow. Indeed, 2^{64} is larger than the maximum integer value in Julia:

```
typemax(Int)
2^63-1
```

And the solution is to call our function on “big” integers with an arbitrary number of bits:

```
fermat_prime(big(6))
```

22.5.3. Debugger.jl

Debugger.jl allows us to interrupt code execution anywhere we want, even in functions we did not write. Using its `@enter` macro, we can enter a function call and walk through the call stack, at the cost of reduced performance.

The REPL prompt changes to `1|debug>`, allowing you to use custom navigation commands to step into and out of function calls, show local variables and set breakpoints. Typing a backtick ` will change the prompt to `1|julia>`, indicating evaluation mode. Any expression typed in this mode will be evaluated in the local context. This is useful to show local variables, as demonstrated in the following example:

```
julia> using Debugger

julia> @enter fermat_prime(6)
In fermat_prime(n) at REPL[7]:1
 1  function fermat_prime(n)
>2      k = 2^n
 3      F = 2^k + 1
 4      for d in 2:isqrt(F)  # integer square root
 5          if F % d == 0
 6              return false

About to run: (^)(2, 6)
1|debug> n
In fermat_prime(n) at REPL[7]:1
 1  function fermat_prime(n)
 2      k = 2^n
>3      F = 2^k + 1
 4      for d in 2:isqrt(F)  # integer square root
 5          if F % d == 0
 6              return false
 7      end

About to run: (^)(2, 64)
```

22. Troubleshooting Julia Code

```
1| julia> k  
64
```

💡 Tip

VSCode offers a nice graphical interface for debugging. Click left of a line number in an editor pane to add a *breakpoint*, which is represented by a red circle. In the debugging pane of the Julia extension, click Run and Debug to start the debugger. The program will automatically halt when it hits a breakpoint. Using the toolbar at the top of the editor, you can then *continue*, *step over*, *step into* and *step out* of your code. The debugger will open a pane showing information about the code such as local variables inside of the current function, their current values and the full call stack.

The debugger can be sped up by selectively compiling modules that you will not need to step into via the + symbol at the bottom of the debugging pane. It is often easiest to start by adding ALL_MODULES_EXCEPT_MAIN to the compiled list, and then selectively remove the modules you need to have interpreted.

23. Distributing and Sharing Julia Code

23.1. In this Chapter

Applying software engineering best practices (Chapter 12) in Julia, including testing, documentation, and coverage metrics. Collaborating on code. Publishing packages for others to use.

23.2. Setup

A vast majority of Julia packages are hosted on GitHub (although less common, other options like GitLab are also possible). GitHub is a platform for collaborative software development, based on the version control system Git (see Chapter 12 for an introduction).

The first step is therefore creating an empty GitHub repository on GitHub (don't add a README License, etc. at this step).

💡 Tip

You should try to follow package naming guidelines and add a ".jl" extension at the end, like so: "MyAwesomePackage.jl".

Locally, use PkgTemplates.jl (see Section 21.10.1) to then create the package's folder locally on your computer, which will create a package with several subfolders (these will be described as the chapter progresses).

To sync this up with the newly created GitHub repository, you git push this new folder to the remote repository <https://github.com/myuser/MyAwesomePackage.jl> (GitHub should show you how to do this on the associated repository page).

23.3. GitHub Actions

The most useful aspect of PkgTemplates.jl is that it automatically generates workflows for GitHub Actions. These are stored as YAML files in .github/workflows, with a

23. Distributing and Sharing Julia Code

slightly convoluted syntax that you don't need to fully understand. For instance, the file `CI.yml` contains instructions that execute the tests of your package (see below) for each pull request, tag or push to the `main` branch. This is done on a GitHub server and should theoretically cost you money, but if your GitHub repository is public, you get an unlimited workflow budget for free.

A variety of workflows and functionalities are available through optional plugins. The interactive setting `Template(..., interactive=true)` allows you to select the ones you want for a given package. Otherwise, you will get the default selection, which you are encouraged to look at.

23.4. Testing

The purpose of the `test` subfolder in your package is unit testing: automatically checking that your code behaves the way you want it to. For instance, if you write your own square root function, you may want to test that it gives the correct results for positive numbers, and errors for negative numbers.

```
using Test

@test sqrt(4) ≈ 2

@testset "Invalid inputs" begin
    @test_throws DomainError sqrt(-1)
    @test_throws MethodError sqrt("abc")
end;
```

Such tests belong in `test/runtests.jl`, and they are executed with the `]test` command (in the REPL's Pkg mode). Unit testing may seem rather naive, or even superfluous, but as your code grows more complex, it becomes easier to break something without noticing. Testing each part separately will increase the reliability of the software you write.

Tip

To test the arguments provided to the functions within your code (for instance their sign or value), avoid `@assert` (which can be deactivated) and use `ArgCheck.jl` instead.

That is, avoid this:

```
function mysqrt(x)
    @assert x >= 0
```

...

And do this instead:

```
function mysqrt(x)
    @argcheck x < 0 DomainError
    ...
end
```

At some point, your package may require test-specific dependencies. In essence, you give the test subfolder its own environment and `Project.toml` file. This often happens when you need to test compatibility with another package, on which you do not depend for the source code itself. Or it may simply be due to testing-specific packages like the ones we will encounter below. For interactive testing work, use `TestEnv.jl` to activate the full test environment (faster than running `]test` repeatedly).

💡 Tip

The Julia extension also offers a more advanced own testing framework, which relies on defining “test items” the code. The benefit of this is that the tests will integrate more directly with the VS Code interface and specific subgroups of tests can be run independently, on-demand. See `TestItemRunner.jl` for more.

💡 Tip

If you want to have more control over your tests, you can try

- `ReferenceTests.jl` to compare function outputs with reference files.
- `ReTest.jl` to define tests next to the source code and control their execution.
- `TestSetExtensions.jl` to make test set outputs more readable.
- `TestReadme.jl` to test whatever code samples are in your `README`.
- `ReTestItems.jl` for an alternative take on VSCode’s test item framework.

23.4.1. Code Coverage

Code coverage refers to the fraction of lines in your source code that are covered by tests (described in more detail in Section 12.3.2). `Codecov` is a website that provides easy visualization of this coverage, and many Julia packages use it. It is available as a `PkgTemplates.jl` plugin, but you have to perform an additional configuration step on the repo for `Codecov` to communicate with it.

23. Distributing and Sharing Julia Code

23.5. Code Style

To make your code easy to read, it is recommended to follow a consistent set of guidelines. The official style guide is very short, so most people use third party style guides like BlueStyle or SciMLStyle.

JuliaFormatter.jl is an automated formatter for Julia files which can help you enforce the style guide of your choice. Just add a file `.JuliaFormatter.toml` at the root of your repository, containing a single line like

```
style = "blue"
```

Then, the package directory will be formatted in the BlueStyle whenever you call

```
using JuliaFormatter  
JuliaFormatter.format(MyAwesomePackage)
```

Note

The default formatter uses JuliaFormatter.jl.

Tip

You can format code automatically in GitHub pull requests with the `julia-format` action, or add the formatting check directly to your test suite.

23.6. Code quality

Of course, there is more to code quality than just formatting. Aqua.jl (Auto QUality Assurance) provides a set of routines that examine other aspects of your package, from ensuring that there are no unused dependencies to catching ambiguous methods statically.

Include the following in your tests to have Aqua.jl run various checks against your code each time tests get run.:

```
using Aqua, MyAwesomePackage  
Aqua.test_all(MyAwesomePackage)
```

JET.jl is tool that is similar to a static linter in other languages. This means that it can inspect your code and ‘understand’ it well enough to catch many types of errors before runtime. It does this by running type inference and figuring out how a given type will flow through the call stack of methods.

You can either use it in report mode (with a nice VSCode display) or in test mode as follows:

```
using JET, MyAwesomePackage
JET.report_package(MyAwesomePackage)
JET.test_package(MyAwesomePackage)
```

Note that both Aqua.jl and JET.jl might pick up false positives: refer to their respective documentations for ways to make them less sensitive.

Finally, ExplicitImports.jl can help you get rid of generic imports to specify where each of the variables in your package comes from. As a project gets more complex, using SomePackage can bring many, sometimes conflicting symbols into your current namespace. ExplicitImports forces you to either qualify the usage (e.g. SomePackage.somefunction(...)) or explicitly opt into importing certain variables.

23.7. Documentation

Refer to for more detail on documentation and its importance in Section 12.4.2. Here are some additional workflow tips for setting up documentation for your package.

DocStringExtensions.jl provides a few shortcuts that can speed up docstring creation by taking care of the obvious parts.

In addition to docstrings, Documenter.jl allows you to design a website for all of this, based on Markdown files contained in the docs subfolder of your package. Unsurprisingly, its own documentation is excellent and will teach you a lot. To build the documentation locally, just run

```
julia> using Pkg

julia> Pkg.activate("docs")

julia> include("docs/make.jl")
```

Then, use LiveServer.jl from your package folder to visualize and automatically update the website as the code changes (similar to Revise.jl, but for your docpages instead of your code):

```
julia> using LiveServer

julia> servedocs()
```

23. Distributing and Sharing Julia Code

To host the documentation online easily, just select the Documenter plugin from PkgTemplates.jl during creation. Not only will this fill the docs subfolder with the appropriate starting files: it will also initialize a GitHub Actions workflow to build and deploy your website on GitHub pages. Lastly, select the gh-pages branch as source in the Github settings for your repository.

23.8. Literate programming

Literate programming is so-called for combining written documents with the output of programs (literature + code = literate programming). These tools allow you to interleave code with texts, formulas, images and so on.

In addition to the Pluto.jl and Jupyter notebooks, take a look at Literate.jl to enrich your code with comments and translate it to various formats. Books.jl is relevant to draft long documents in a pure Julia way.

Quarto is an open-source scientific and technical publishing system that supports Python, R and Julia. Quarto can render markdown files (.md), Quarto markdown files (.qmd), and Jupyter Notebooks (.ipynb) into documents (Word, PDF, presentations), web pages, blog posts, books, and more. Additionally, Quarto makes it easy to share or publish rendered content to various online hosts.

PPTX.jl will create Microsoft PowerPoint files.

23.9. Versions and registration

23.9.1. Versions and Compatibility

The Julia community has adopted semantic versioning, which means every package must have a version, and the version numbering follows strict rules (the concept of versioning was covered in Section 12.6.2).

To comply with the versioning requirements in Pkg's resolver, you need to specify compatibility bounds for your dependencies: this happens in the [compat] section of your Project.toml. To initialize these bounds with current dependency versions, use the]compat command in the Pkg mode of the REPL, or the package PackageCompatUI.jl.

Over time, new versions of your dependencies will be released. The CompatHelper.jl GitHub Action will help you monitor upstream Julia dependencies and suggest changes to your Project.toml's [compat] section accordingly. In addition, Dependabot can monitor the dependencies... of your GitHub actions themselves. Both of these can be included in the default PkgTemplates setup.

 Tip

It may also happen that you incorrectly promise compatibility with an old version of a package and not realize it (since Pkg prefers newer versions within the compatibility bounds, not all combinations get tested). To prevent that, the julia-downgrade-compat GitHub action tests your package with the oldest possible version of every dependency, and verifies that everything still works.

23.9.2. Registration

If your package is useful to others in the community, it may be a good idea to register it, that is, make it part of the pool of packages that can be installed with

```
pkg> add MyAwesomePackage # made possible by registration
```

Note that unregistered packages can also be installed by anyone from the GitHub URL, but this is a less reproducible solution:

```
pkg> add https://github.com/myuser/MyAwesomePackage # not ideal
```

To register your package, check out the general registry guidelines. The Registrator.jl bot can help you automate the process. Another handy bot, provided by default with PkgTemplates.jl, is TagBot: it automatically tags new versions of your package following each registry release. If you have performed the necessary SSH configuration, TagBot will also trigger documentation website builds following each release.

23.9.2.1. Local Registry

For distributing privately (or publicly if you make the repository public), LocalRegistry.jl provides convenience functions for creating a new registry, adding new packages, and updating versions for the packages. If you want to share packages internally, create and register packages to a repository that's hosted somewhere you and your team can access. If you wanted to make the repository public, you can publish the registry repository somewhere publicly accessible (such as a public GitHub repository).

Once established, other users can add a repository as easily as entering package mode and running `registry add`. Say that we have already put a registry we called `FinancePackages` in a repository on the company intranet:

```
pkg> registry add http://company-intranet.com/git/FinancePackages.git
```

23. Distributing and Sharing Julia Code

23.9.2.2. Hosted Registries

Alternatively to a self-hosted local registry, third party services such as JuliaHub provide managed registries well suited for corporate environments.

23.10. Reproducibility

Obtaining consistent and reproducible results is an essential part of model auditing and compliance. One tool to consider is DrWatson.jl. It is a general toolbox for running and re-running models in an orderly fashion.

Some specific issues come up in attempting to ensure reproducibility

A first hurdle is random number generation, which is not guaranteed to remain stable across Julia versions. To ensure that the random streams remain exactly the same, you need to use StableRNGs.jl. The downside to this is that the random number generation will be considerably slower than the usual generator.

Another aspect is dataset download and management. The packages DataDeps.jl, DataToolkit.jl and ArtifactUtils.jl can help you bundle non-code elements with your package (some of these rely on artifacts - discussed in Section 12.6.3).

23.11. Interoperability

To ensure compatibility with earlier Julia versions, Compat.jl is your best ally.

Making packages play nice with one another is a key goal of the Julia ecosystem. Since Julia 1.9, this can be done with package extensions, which override specific behaviors based on the presence of a given package in the environment. For example, if you want to provide pre-configured plotting, but don't in general need to include a plotting library as part of your package for all users and use cases. PackageExtensionTools.jl eases of setting up extensions for your package.

Furthermore, the Julia ecosystem as a whole plays nice with other programming languages too. C and Fortran are natively supported. Python can be easily interfaced with the combination of CondaPkg.jl and PythonCall.jl. Other language compatibility packages can be found in the JuliaInterop organization, like RCall.jl.

23.12. Customization

Part of interoperability is also flexibility and customization: the Preferences.jl package gives a nice way to specify various options in TOML files. These customizable preferences persist across sessions and provide the preferences at both compile and runtime. For example, say different parts of a company had different preferred data sources but otherwise used the same code. This could be set in a way via Preferences.jl so that each team can share the logic while seamlessly defaulting to different data sources.

23.13. Collaboration

Once your package grows big enough, you might need to bring in some help. Working together on a software project has its own set of challenges, which are partially addressed by a good set of ground rules like SciML ColPrac. Of course, collaboration goes both ways: if you find a Julia package you really like, you are more than welcome to contribute as well, for example by opening issues or submitting pull requests.

24. Optimizing Julia Code

The two fundamental principles for writing fast Julia code:

1. Ensure that **the compiler can infer the type** of every variable.
2. Avoid **unnecessary (heap) allocations**.

24.1. Type Inference

The compiler's job is to optimize and translate Julia code it into runnable machine code. If a variable's type cannot be deduced before the code is run, then the compiler won't generate efficient code to handle that variable. This phenomenon is called "type instability". Enabling type inference means making sure that every variable's type in every function can be deduced from the types of the function inputs alone.

That is, the compiler will be able to create more optimized code if it can analyze the function you've written and determine what type will be returned. In the following function, the compiler can analyze the expressions and determine with certainty that if `mysum` is given two `Ints` as arguments, the return value will also be an `Int`.

```
function mysum(a,b)
    a + b
end
```

24.2. Avoiding Heap Allocations

A "heap allocation" (or simply "allocation") occurs when we create a new variable without knowing how much space it will require (like a `Vector` with flexible length). This has two implications:

1. Allocating memory on the heap takes substantially more time than stack allocated memory to be created.
2. Periodically, a **garbage collector** (GC), needs to run to de-allocate (free up) memory on the heap which is no longer used by the program

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Execution of code is stopped while the garbage collector runs, so minimising its usage is important.

The vast majority of performance tips come down to these two fundamental ideas.

Typically, the most common beginner pitfall is the use of untyped global variables without passing them as arguments. Why is it bad? Because the type of a global variable can change outside of the body of a function, so it causes type instabilities wherever it is used. Those type instabilities in turn lead to more heap allocations.

Much more detail on performance considerations is covered in Chapter 10.

24.3. Measuring performance

The simplest way to measure how fast a piece of code runs is to use the `@time` macro, which returns the result of the code and prints the measured runtime and allocations. Because code needs to be compiled before it can be run, you should first run a function without timing it so it can be compiled, and then time it:

```
sum_abs(vec) = sum(abs(x) for x in vec);
v = rand(100);

using BenchmarkTools
@time sum_abs(v); # Inaccurate, note the >99% compilation time
@time sum_abs(v); # Accurate

0.008580 seconds (70.38 k allocations: 3.393 MiB, 99.85% compilation time)
0.000002 seconds (1 allocation: 16 bytes)
```

Using `@time` is quick but it has flaws, because your function is only measured once. That measurement might have been influenced by other things going on in your computer at the same time. In general, running the same block of code multiple times is a safer measurement method, because it diminishes the probability of only observing an outlier.

24.3.1. BenchmarkTools

`BenchmarkTools.jl` is the most popular package for repeated measurements on function executions. Similarly to `@time`, `BenchmarkTools` offers `@btime` which can be used in exactly the same way but will run the code multiple times and provide an average. Additionally, by using `$` to interpolate external values, you remove the overhead caused by global variables.

```
using BenchmarkTools
@btime sum_abs(v);
@btime sum_abs($v);

22.381 ns (1 allocation: 16 bytes)
17.410 ns (0 allocations: 0 bytes)
```

In more complex settings, you might need to construct variables in a setup phase that is run before each sample. This can be useful to generate a new random input every time, instead of always using the same input.

```
my_matmul(A, b) = A * b;
@btime my_matmul(A, b) setup =
    A = rand(1000, 1000); # use semi-colons between setup lines
    b = rand(1000)
);

137.417 μs (3 allocations: 7.94 KiB)
```

For better visualization, the `@benchmark` macro shows performance histograms:

i Note

Certain computations may be optimized away by the compiler before the benchmark takes place. If you observe suspiciously fast performance, especially below the nanosecond scale, this is very likely to have happened.

24.3.2. Other tools

`BenchmarkTools.jl` works fine for relatively short and simple blocks of code (microbenchmarking). To find bottlenecks in a larger program, you should rather use a profiler (see next section) or the package `TimerOutputs.jl`. It allows you to label different sections of your code, then time them and display a table of grouped by label.

Finally, if you know a loop is slow and you'll need to wait for it to be done, you can use `ProgressMeter.jl` or `ProgressLogging.jl` to track its progress. This will display a progress bar in VS Code or in a notebook, indicating how far along a loop has progressed.

24.4. Profiling

Profiling can identify performance bottlenecks at function level, and graphical tools such as `ProfileView.jl` are the best way to use it.

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24.4.1. Sampling

Whereas a benchmark measures the overall performance of some code, a profiler breaks it down function by function to identify bottlenecks. Sampling-based profilers periodically ask the program which line it is currently executing, and aggregate results by line or by function. Julia offers two kinds: one for runtime (in the module `Profile`) and one for memory (in the submodule `Profile.Allocs`).

These built-in profilers print textual outputs, but the result of profiling is best visualized as a flame graph. In a flame graph, each horizontal layer corresponds to a specific level in the call stack, and the width of a tile shows how much time was spent in the corresponding function. Here's an example:

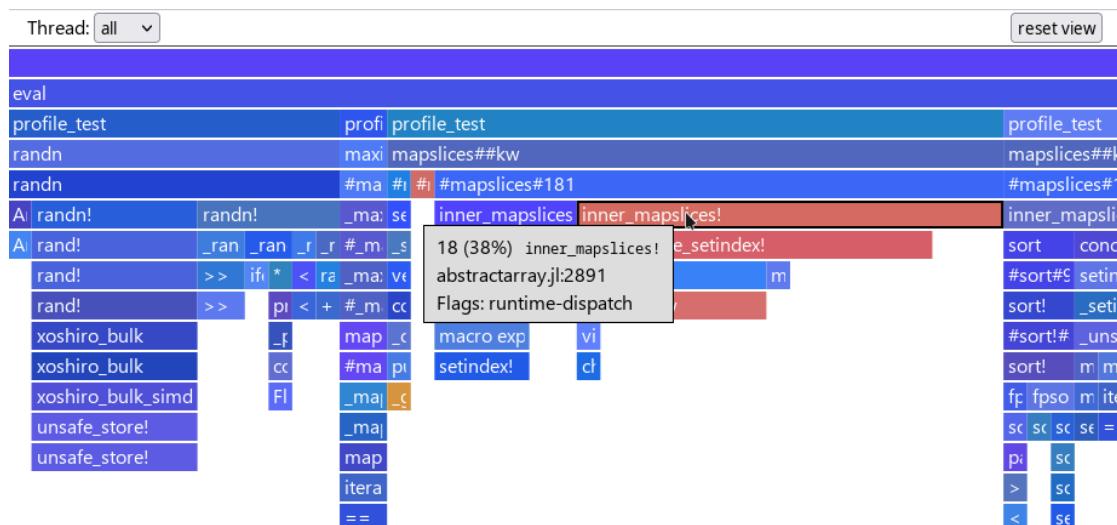


Figure 24.1.: flamegraph

24.4.2. Visualization Profile Results

The packages `ProfileView.jl` and `PProf.jl` both allow users to record and interact with flame graphs. `ProfileView.jl` is simpler to use, but `PProf` is more featureful and is based on `pprof`, an external tool maintained by Google which applies to more than just Julia code. Here we only demonstrate the former:

```
julia profileview-example using ProfileView @profview do_work(some_input)
```

 Tip

Calling `@profview do_work(some_input)` in the integrated Julia REPL will open an interactive flame graph, similar to `ProfileView.jl` but without requiring a separate package.

To integrate profile visualisations into environments like Jupyter and Pluto, use `ProfileSVG.jl` or `ProfileCanvas.jl`, whose outputs can be embedded into a notebook.

No matter which tool you use, if your code is too fast to collect samples, you may need to run it multiple times in a loop.

 Tip

To visualize memory allocation profiles, use `PProf.jl` or `VSCode's @profview_allocs`. A known issue with the allocation profiler is that it is not able to determine the type of every object allocated, instead `Profile.Allocs.UnknownType` is shown instead. Inspecting the call graph can help identify which types are responsible for the allocations.

24.4.3. External profilers

Apart from the built-in `Profile` standard library, there are a few external profilers that you can use including Intel VTune (in combination with `IntelITT.jl`), NVIDIA Nsight Systems (in combination with `NVTX.jl`), and Tracy.

24.5. Type stability

For a section of code to be considered type stable, the type inferred by the compiler must be “concrete”, which means that the size of memory that needs to be allocated to store its value is known at compile time. Types declared abstract with `abstract type` are not concrete and neither are parametric types whose parameters are not specified:

```
@show isconcretetype(Any)
@show isconcretetype(AbstractVector)
@show isconcretetype(Vector) # Shorthand for `Vector{T}` where T
@show isconcretetype(Vector{Real})
@show isconcretetype(eltype(Vector{Real}))
@show isconcretetype(Vector{Int64})
```

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```
isconcretetype(Any) = false
isconcretetype(AbstractVector) = false
isconcretetype(Vector) = false
isconcretetype(Vector{Real}) = true
isconcretetype(eltype(Vector{Real})) = false
isconcretetype(Vector{Int64}) = true

true
```

Note

`Vector{Real}` is concrete despite `Real` being abstract for subtle typing reasons but it will still be slow in practice because the type of its elements is abstract.

While type-stable function calls compile down to fast GOTO statements, type-unstable function calls generate code that must read the list of all methods for a given operation and find the one that matches. This phenomenon called “dynamic dispatch” prevents further optimizations.

Type-stability is a contagious thing: if a variable’s type cannot be inferred, then the types of variables that depend on it may not be inferrable either. Most code should be type-stable unless it has a good reason not to be.

24.5.1. Detecting Instabilities

The simplest way to detect an instability is with the builtin macro `@code_warntype`: The output of `@code_warntype` is difficult to parse, but the key takeaway is the return type of the function’s Body: if it is an abstract type, like `Any`, something is wrong. In a normal Julia REPL, such cases would show up colored in red as a warning.

```
using InteractiveUtils # loaded automatically if using a notebook or REPL
function put_in_vec_and_sum(x)
    v = []
    push!(v, x)
    return sum(v)
end;

@code_warntype put_in_vec_and_sum(1)

MethodInstance for put_in_vec_and_sum(::Int64)
  from put_in_vec_and_sum(x) @ Main In[7]:2
Arguments
```

```
#self#::Core.Const(Main.put_in_vec_and_sum)
x::Int64
Locals
  v::Vector{Any}
Body::Any
1 -      (v = Base.vect())
    %2 = Main.push!::Core.Const(push!)
    %3 = v::Vector{Any}
          (%2)(%3, x)
    %5 = v::Vector{Any}
    %6 = Main.sum(%5)::Any
    return %6
```

Unfortunately, `@code_warntype` is limited to one function body: calls to other functions are not expanded, which makes deeper type instabilities easy to miss. That is where JET.jl can help: it provides optimization analysis aimed primarily at finding type instabilities. While test integrations are also provided, the interactive entry point of JET is the `@report_opt` macro.

```
using JET
@report_opt put_in_vec_and_sum(1)
```

[Info: tracking Base

```
===== 4 possible errors found =====
put_in_vec_and_sum(x::Int64) @ Main ./In[7]:5
  sum(a::Vector{Any}) @ Base ./reducedim.jl:982
    sum(a::Vector{Any}; dims::Colon, kw::Kwargs{}) @ Base ./reducedim.jl:982
      _sum(a::Vector{Any}, ::Colon) @ Base ./reducedim.jl:986
        _sum(a::Vector{Any}, ::Colon; kw::Kwargs{}) @ Base ./reducedim.jl:986
          _sum(f::typeof(identity), a::Vector{Any}, ::Colon) @ Base ./reducedim.jl:987
            _sum(f::typeof(identity), a::Vector{Any}, ::Colon; kw::Kwargs{}) @ Base ./reducedim.jl:987
              mapreduce(f::typeof(identity), op::typeof(Base.add_sum), A::Vector{Any}) @ Base ./reducedim.jl:987
                mapreduce(f::typeof(identity), op::typeof(Base.add_sum), A::Vector{Any}; dims::Colon,
                  _mapreduce_dim(f::typeof(identity), op::typeof(Base.add_sum), ::Base._InitialValue,
                    _mapreduce(f::typeof(identity), op::typeof(Base.add_sum), ::IndexLinear, A::Vector{Any}),
                      mapreduce_impl(f::typeof(identity), op::typeof(Base.add_sum), A::Vector{Any}, ifirst,
                        mapreduce_impl(f::typeof(identity), op::typeof(Base.add_sum), A::Vector{...}, ifirst,
```

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```
||| runtime dispatch detected: op::typeof(Base.add_sum)(%91::Any, %110::Any)::Any
|||   mapreduce_impl(f::typeof(identity), op::typeof(Base.add_sum), A::Vector{...}, ifirst
|||     runtime dispatch detected: op::typeof(Base.add_sum)(%187::Any, %189::Any)::Any
|||       _mapreduce(f::typeof(identity), op::typeof(Base.add_sum), ::IndexLinear, A::Vector{...
|||         runtime dispatch detected: op::typeof(Base.add_sum)(%97::Any, %115::Any)::Any
|||           _mapreduce(f::typeof(identity), op::typeof(Base.add_sum), ::IndexLinear, A::Vector{...
|||             runtime dispatch detected: op::typeof(Base.add_sum)(%118::Any, %139::Any)::Any
```

💡 Tip

The Julia extension features a static linter, and runtime diagnostics with JET can be automated to run periodically on your codebase and show any problems detected.

Cthulhu.jl¹ exposes the `@descend` macro which can be used to interactively “step through” lines of the corresponding typed code, and “descend” into a particular line if needed. This is akin to repeatedly calling `@code_warntype` deeper and deeper into your functions (“slowly succumbing to the madness...” of type instability).

24.5.2. Fixing Instabilities

The Julia manual has a collection of tips to improve type inference.

💡 Tip

To be more forceful about ensuring type stability in your code, one approach is to error whenever a type instability occurs: the macro `@stable` from DispatchDoctor.jl allows exactly that.

When working with parametric types, look to avoid usage of generic type parameters (e.g. `Array{Any}`) whenever possible. For custom types, make use of parametric types to create type-stable abstractions. For example, the latter `struct Bond2` or `struct Bond3` will allow Julia to create distinct concrete types and methods as opposed to needing generic runtime dispatch due to the unpredictable potential types of the `struct` fields.

¹So-named for the “slow descent into madness” when descending into functions to follow the Julia compiler’s type inference across many layers of function calls.

The difference between Bond2 and Bond3 is that the fields in Bond3 could be any type (such as `Float16` or `String`) as long as both of them equaled the type variable `T`.

```
struct Bond1
    par
    coupon
end

struct Bond2
    par::Float64
    coupon::Float64
end

struct Bond3{T}
    par::T
    coupon::T
end
```

24.6. Memory management

After ensuring type stability, one should try to reduce the number of heap allocations a program makes. Again, the Julia manual has a series of tricks related to arrays and allocations which you should take a look at. In particular, try to modify existing arrays instead of allocating new objects and try to access arrays in the right order for Julia, i.e. accessing data down columns instead of across rows.

Alternatively, to ensure that non-allocating functions never regress in future versions of your code, you can write a test set to check allocations by providing the function and a concrete type-signature.

```
julia AllocCheck @testset "non-allocating" begin      @test isempty(AllocCheck.check_allocs(my_func(Float64, Float64))) end
```

24.7. Compilation

A number of tools allow you to reduce Julia's latency, also referred to as TTFX (time to first X, where X was historically plotting a graph).

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24.7.1. Precompilation

PrecompileTools.jl reduces the amount of time taken to run functions loaded from a package or local module that you wrote. It allows module authors to specify methods to precompile when a module is loaded for the first time. The methods chosen should represent those which would be used during typical user workflows. These methods then have the same latency as if they had already been run by the end user. This adds upfront pre-compilation time when installing a package version for the first time, but then subsequent uses will be much quicker.

Here's an example of precompilation, adapted from the package's documentation:

```
module MyPackage

using PrecompileTools: @compile_workload

struct MyType
    x::Int
end

myfunction(a::Vector) = a[1].x

@compile_workload begin
    a = [MyType(1)]
    myfunction(a)
end

end
```

Note that every method that is called will be compiled, no matter how far down the call stack or which module it comes from. To see if the intended calls were compiled correctly or diagnose other problems related to precompilation, use SnoopCompile.jl. This is especially important for writers of registered Julia packages, as it allows you to diagnose recompilation that happens due to invalidation.

24.7.2. Package compilation

To reduce the time that packages take to load, you can use PackageCompiler.jl to generate a custom version of Julia, called a **sysimage** (system image), with its own custom standard library. As packages in the standard library are already compiled, any `using` or `import` statement involving them is almost instant.

Once PackageCompiler.jl is added to your global environment, activate a local environment for which you want to generate a sysimage, ensure all of the packages you want

to compile are in its `Project.toml`, and run `create_sysimage` as in the example below. The filetype of `sysimage_path` differs by operating system: Linux has `.so`, MacOS has `.dylib`, and Windows has `.dll`.

```
using PackageCompiler # installed in global environment
packages_to_compile = ["Makie", "DifferentialEquations"]
create_sysimage(packages_to_compile; sysimage_path="MySysimage.so")
```

Once a sysimage is generated, it can be used with the command line flag: `julia --sysimage=path/to/sysimage`.

💡 Tip

The generation and loading of sysimages can be streamlined with VSCode. By default, the command sequence Task: Run Build Task followed by Julia: Build custom sysimage for current environment will compile a sysimage containing all packages in the current environment, but additional details can be specified in a `/.vscode/JuliaSysimage.toml` file. To automatically detect and use a custom sysimage, set `useCustomSysimage` to `true` in the application settings.

24.7.3. Static compilation

`PackageCompiler.jl` also facilitates the creation of apps and libraries that can be shared to and run on machines that don't have Julia installed.

At a basic level, all that's required to turn a Julia module `MyModule` into an app is a function `julia_main()::Cint` that returns `0` upon successful completion. Then, with `PackageCompiler.jl` loaded, run `create_app("MyModule", "MyAppCompiled")`. Command line arguments to the resulting app are assigned to the global variable `ARGS::Array{ASCIIString}`, the handling of which can be made easier by `ArgParse.jl`.

In Julia, a library is just a sysimage with some extras that enable external programs to interact with it. Any functions in a module marked with `Base.@ccallable`, and whose type signature involves C-conforming types e.g. `Cint`, `Cstring`, and `Cvoid`, can be compiled into an externally callable library with `create_library`, similarly to `create_app`. Unfortunately, the process of compiling and sharing a standalone executable or callable library must take relocability into account, which is beyond the scope of this blog.

ℹ Note

An alternative way to compile a shareable app or library that doesn't need to compile a sysimage, and therefore results in smaller binaries, is to use `StaticCom-`

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piler.jl and its sister package StaticTools.jl. The biggest tradeoff of not compiling a sysimage, is that Julia's garbage collector is no longer included, so all heap allocations must be managed manually, and all code compiled *must* be type-stable. To get around this limitation, you can use static equivalents of dynamic types, such as a `StaticArray` (`StaticArrays.jl`) instead of an `Array` or a `StaticString` (`StaticTools.jl`), use `malloc` and `free` from `StaticTools.jl` directly, or use arena allocators with `Bumper.jl`. The README of `StaticCompiler.jl` contains a more detailed guide on how to prepare code to be compiled.

Note

Starting with Julia 1.12, it is anticipated that there will be a new way to compile Julia to small, static binaries with a tool called `juliac`.

24.8. Parallelism

Code can be made to run faster through parallel execution with multithreading (shared-memory parallelism) or multiprocessing / distributed computing. Parallelism was covered in a whole chapter (Chapter 11), but this section provides insight into recommended packages and patterns specific to Julia.

Many common operations such as maps and reductions can be trivially parallelised through either method by using their respective Julia packages (e.g `pmap` from `Distributed.jl` and `tmap` from `OhMyThreads.jl`). Multithreading is available on almost all modern hardware, whereas distributed computing is most useful to users of high-performance computing clusters.

24.8.1. Multithreading

To use multi-threading, Julia needs to be started with more than one thread. This can be done by either setting the environment variable `JULIA_NUM_THREADS` to either `auto` or specify a number like `4`. You can also specify how many threads to start julia with if given the `-t` command line argument (such as running `julia -t 4` to start Julia with four threads from the command line). Once Julia is running, you can check if this was successful by calling `Threads.threads()`.

 Tip

In VS Code, the default number of threads can be edited by adding "julia.NumThreads": auto, to your settings. This will be applied to the integrated terminal.

Why doesn't Julia automatically start with more than one thread? Between "hyper-threading" (synthetic additional thread capacity), multi-core architectures, and the different types of threads it's actually difficult to predict how many threads will be optimal for a given system. Julia's current default is to take the more conservative approach and start single-threaded unless otherwise specified. The "auto" option is a best-guess but can, on certain systems and configurations, be very bad for performance. The authors recommend for most common systems to just use "auto".

 Tip

Linear algebra code calls the low-level libraries BLAS and LAPACK. These libraries manage their own pool of threads, so single-threaded Julia processes can still make use of multiple threads. Multi-threaded Julia processes that call these libraries may run into performance issues due to the limited number of threads available in a single core.

In this case, once `LinearAlgebra` is loaded, BLAS can be set to use only one thread by calling `BLAS.set_num_threads(1)`. For more information see the docs on multithreading and linear algebra.

Regardless of the number of threads, you can parallelise a for loop with the macro `Threads.@threads`. The macros `@spawn` and `@async` function similarly, but require more manual management of tasks and their results. For this reason `@threads` is recommended for those who do not wish to use third-party packages.

When designing multithreaded code, you should generally try to write to shared memory as rarely as possible. Where it cannot be avoided, you need to be careful to avoid "race conditions", i.e. situations when competing threads try to write different things to the same memory location. It is usually a good idea to separate memory accesses with loop indices, as in the example below:

```
results = zeros(Int, 4)
Threads.@threads for i in 1:4
    results[i] = i^2
end
```

Almost always, it is **not** a good idea to use `threadid()`.

Even if you manage to avoid any race conditions in your multithreaded code, it is very easy to run into subtle performance issues (like false sharing). For these reasons, you

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might want to consider using a high-level package like OhMyThreads.jl, which provides a user-friendly alternative to Threads and makes managing threads and their memory use much easier. The helpful translation guide demonstrates how Base multi-threading loops can be translated into the OhMyThreads API.

If the latency of spinning up new threads becomes a bottleneck, check out Polyester.jl for very lightweight threads that are quicker to start.

If you're on Linux, you should consider using ThreadPinning.jl to pin your Julia threads to CPU cores to obtain stable and optimal performance. The package can also be used to visualize where the Julia threads are running on your system (see `threadinfo()`).

24.8.2. Distributed computing

Julia's multiprocessing and distributed relies on the standard library `Distributed`. The main difference with multi-threading is that data isn't shared between worker processes. Once Julia is started, processes can be added with `addprocs`, and their can be queried with `nworkers`.

The macro `Distributed.@distributed` is a *syntactic* equivalent for `Threads.@threads`. Hence, we can use `@distributed` to parallelise a for loop as before, but we have to additionally deal with sharing and recombining the results array. We can delegate this responsibility to the standard library `SharedArrays`. However, in order for all workers to know about a function or module, we have to load it `@everywhere`:

```
using Distributed

# Add additional workers then load code on the workers
addprocs(3)
@everywhere using SharedArrays
@everywhere f(x) = 3x^2

results = SharedArray{Int}(4)
@sync @distributed for i in 1:4
    results[i] = f(i)
end
results

4-element SharedVector{Int64}:
 3
 12
 27
 48
```

Note that `@distributed` does not force the main process to wait for other workers, so we must use `@sync` to block execution until all computations are done.

One feature `@distributed` has over `@threads` is the possibility to specify a reduction function (an associative binary operator) which combines the results of each worker. In this case `@sync` is implied, as the reduction cannot happen unless all of the workers have finished.

```
using Distributed # hide

@distributed (+) for i in 1:4
    i^2
end
```

30

Alternately, the convenience function `pmap` can be used to easily parallelise a `map`, both in a distributed and multi-threaded way.

```
results = pmap(f, 1:100; distributed=true, batch_size=25, on_error=ex → 0)

100-element Vector{Int64}:
 3
 12
 27
 48
 75
 108
 147
 192
 243
 300
 363
 432
 507
  :
23763
24300
24843
25392
25947
26508
27075
27648
```

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28227
28812
29403
30000

For more functionalities related to higher-order functions, `Transducers.jl` and `Folds.jl` are the way to go.

 Tip

`MPI.jl` implements the Message Passing Interface standard, which is heavily used in high-performance computing beyond Julia. The C library that `MPI.jl` wraps is *highly* optimized, so Julia code that needs to be scaled up to a large number of cores, such as an HPC cluster, will typically run faster with MPI than with plain Distributed.

`Elemental.jl` is a package for distributed dense and sparse linear algebra which wraps the `Elemental` library written in C++, itself using MPI under the hood.

24.8.3. GPU programming

GPUs are specialized in executing instructions in parallel over a large number of threads. While they were originally designed for accelerating graphics rendering, more recently they have been used to train and evaluate machine learning models.

Julia's GPU ecosystem is managed by the JuliaGPU organization, which provides individual packages for directly working with each GPU vendor's instruction set. The most popular one is `CUDA.jl`, which also simplifies installation of CUDA drivers for NVIDIA GPUs. Through `KernelAbstractions.jl`, you can easily write code that is agnostic to the type of GPU where it will run.

24.8.4. SIMD instructions

In the Single Instruction, Multiple Data (SIMD) paradigm, several processing units perform the same instruction at the same time, differing only in their inputs. The range of operations that can be parallelized (or “vectorized”) like this is more limited than in the previous sections, and slightly harder to control. Julia can automatically vectorize repeated numerical operations (such as those found in loops) provided a few conditions are met:

1. Reordering operations must not change the result of the computation.
2. There must be no control flow or branches in the core computation.
3. All array accesses must follow some linear pattern.

While this may seem straightforward, there are a number of important caveats which prevent code from being vectorized. Performance annotations like `@simd` or `@inbounds` help enable vectorization in some cases, as does replacing control flow with `if else`.

If this isn't enough, SIMD.jl allows users to force the use of SIMD instructions and bypass the check for whether this is possible. One particular use-case for this is for vectorizing non-contiguous memory reads and writes through `SIMD.vgather` and `SIMD.vscatter` respectively.

 Tip

You can detect whether the optimizations have occurred by inspecting the output of `@code_llvm` or `@code_native` and looking for vectorized registers, types, instructions. Note that the exact things you're looking for will vary between code and CPU instruction set, an example of what to look for can be seen in this blog post by Kristoffer Carlsson.

24.8.5. Additional Packages

Some additional packages to be aware of include:

- LoopVectorization.jl which can enhance the vectorized loops even further, such as handling the “tail” of a vectorized loops more efficiently than the base compiler. The “tail” refers to situations like where you have a vector width of 8, but don’t have a collection that’s a nice multiple of 8 (say 1001 elements).
- Octavian.jl implements a linear algebra-like library, utilizing parallelism via vectorization to generate efficient code for the system it’s running on.
- Tulio.jl is an einsum library, a domain-specific language for tensor (a specific subset of vectors) operations, common in machine learning and linear algebra.

24.9. Efficient types

Using an efficient data structure is a tried and true way of improving the performance. While users can write their own efficient implementations through officially documented interfaces, a number of packages containing common use cases are more tightly integrated into the Julia ecosystem.

24.9.1. Static arrays

Using StaticArrays.jl, you can construct arrays which contain size information in their type. Through multiple dispatch, statically sized arrays give rise to specialised, efficient

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methods for certain algorithms like linear algebra. In addition, the `SArray`, `SMatrix` and `SVector` types are immutable, so the array does not need to be garbage collected as it can be stack-allocated. Creating a new `SArrays` comes at almost no extra cost, compared to directly editing the data of a mutable object. With `MArray`, `MMatrix`, and `MVector`, data remains mutable as in normal arrays.

To handle mutable and immutable data structures with the same syntax, you can use `Accessors.jl`:

```
using StaticArrays, Accessors

sx = SA[1, 2, 3] # SA constructs an SArray
@set sx[1] = 3 # Returns a copy, does not update the variable
@reset sx[1] = 4 # Replaces the original

3-element SVector{3, Int64} with indices SOneTo(3):
4
2
3
```

24.9.2. Classic data structures

All but the most obscure data structures can be found in the packages from the Julia-Collections organization, especially `DataStructures.jl` which has all the standards from the computer science courses (stacks, queues, heaps, trees and the like). Iteration and memoization utilities are also provided by packages like `IterTools.jl` and `Memoize.jl`.

24.9.3. Bits types

When you create custom structs, keeping the fields as simple, concrete types makes it more likely that the compiler will be able to allocate these objects on the stack instead of the heap. An example of this was shown in Section 13.5.3.

Part VII.

Applied Financial Modeling Techniques

Here we focus on practical implementation of financial models. This section provides concrete examples and strategies for building effective models across various applications. We'll discuss model design, optimization, and validation, with an emphasis on real-world usage. The goal is to bridge theory and practice, giving you hands-on knowledge to apply advanced techniques to actual financial challenges.

25. Stochastic Mortality Projections

25.1. In This Chapter

A term life insurance policy is used to illustrate: selecting key model features, design tradeoffs between a few different approaches, and a discussion of the performance impacts of the different approaches to parallelism.

25.2. Introduction

Monte Carlo simulation is common in risk and valuation contexts. This worked example will create a population of term life insurance policies and simulate the associated claims stochastically. For this chapter, the focus is not so much on the outcomes of the model, but instead *how* and *why* the model was chosen to be setup in the way that it is.

The general structure of the example is:

1. Define the datatypes and sample data
2. Define the core logic that governs the projected outcomes for the modeled policies
3. Evaluate a few ways to structure the simulation, including:
 4. allocating and non-allocating approaches
 5. single threaded and multi-threaded approaches

As will be shown, the number of simulations able to be completed on modern CPU hardware is really remarkable!

```
using CSV, DataFrames
using MortalityTables, FinanceCore
using Dates
using ThreadsX
using BenchmarkTools
using Random
using CairoMakie
```

25. Stochastic Mortality Projections

25.3. Data and Types

25.3.1. @enums and the Policy Type

Our core unit of simulation will be a single life insurance Policy. Important characteristics include: the age a policy was issued at, the sex of the insured, and risk class to determine the assumed mortality rate. To make the example more realistic and demonstrate how it might look for a real block of inforce policies, additional fields have been included such as ID (not really used, but a common identifier) and COLA which is a cost-of-living-adjustment, used to modify the policy benefit through time. To be clear: the Policy type has more fields than will actually be used in the calculations, with the purpose to show how typical fields used in practice might be defined.

Before we define the core Policy type, there's a couple of types we might consider defining that would subsequently get used by Policy: types representing sex and risk class. A typical approach might be to simply define associated types, like this:

```
abstract type Sex end

struct Male <: Sex end
struct Female <: Sex end
```

Then, if we were to include a sex field in Policy, we could write it like this:

```
struct Policy
    # ...
    sex::Sex
    # ...
end
```

This would be a totally valid and logical approach! However, high performance is a top priority for this simulation, and therefore this approach would sacrifice being able to keep Policy data on the stack instead of the heap. This is because Sex is of unknown concrete type, which could be as simple as our definition above (with no fields in Male and Female) or someone could add a new Alien subtype of Sex with a number of associated data fields! This is why Julia cannot assume that subtypes of Sex will always be a simple Singletons with no associated data.

Instead, we can utilize Enums, which are a sort of lightweight type where the only thing that matters with it is distinguishing between categories. Enums in Base Julia are basically a set of constants grouped together that reference an associated integer.

```
@enum Sex Female = 1 Male = 2
@enum Risk Standard = 1 Preferred = 2
@enum PaymentMode Annual = 1 Quarterly = 4 Monthly = 12
```

Enums are convenient because it lets us use human-meaningful names for integer-based categories. Julia will also keep track of valid options: we cannot now use anything other than `Female` or `Male` where we have said a `Sex` must be specified.

Note

There exist Julia packages which are more powerful versions of Enums, essentially leaning into the ability to use the type system instead of just nice names for categorical variables.

Moving on to the definition of the policy itself, here's what that looks like. Note that every field has a type annotation associated with it.

```
struct Policy
    id::Int
    sex::Sex
    benefit_base::Float64
    COLA::Float64
    mode::PaymentMode
    issue_date::Date
    issue_age::Int
    risk::Risk
end
```

The benefit of the way we have defined it here, using simple bits-types for each field is that our new composite `Policy` type is also a bitstype:

```
let
    p = Policy(1, Male, 1e6, 0.02, Annual, today(), 50, Standard)

    isbits(p)
end

true
```

Note

Type annotations are optional, but providing them is able to coerce the values to be all plain bits (i.e. simple, non-referenced values like arrays are) when the type is constructed. We could have defined `Policy` without the types specified:

```
struct Policy
    id
    sex
```

25. Stochastic Mortality Projections

```
...  
risk  
end
```

Leaving out the annotations here forces Julia to assume that `Any` type could be used for the given field. Having the field be of type `Any` makes the instantiated struct data be stored in the heap, since Julia can't know the size of `Policy` in bits in advance.

We would also find that the un-annotated type is about 50 times slower than the one with annotation due to the need to utilize runtime lookup and reference memory on the heap instead of the stack.

25.3.2. The Data

To partially illustrate a common workflow, we'll pretend that the data we are interested in comes from a CSV file, which will be defined inline using an `IOBuffer` so that the structure of the source data is clear to the reader. Only two policies will be listed for brevity, but we will duplicate them for simulation purposes later on.

```
sample_csv_data =  
    IOBuffer(  
        raw"id,sex,benefit_base,COLA,mode,issue_date,issue_age,risk  
        1,M,100000.0,0.03,12,1999-12-05,30,Std  
        2,F,200000.0,0.03,12,1999-12-05,30,Pref"  
    )
```

```
IOBuffer(data=UInt8[...], readable=true, writable=false, seekable=true, append=false, size=152,
```

We will not load the sample data using a common pattern:

1. Load the source file into a `DataFrame`
2. `map` over each row of the dataframe, and return an instantiated `Policy` object
3. Within the map, apply basic data parsing and translation logic as needed.

```
policies = let  
  
    # read CSV directly into a dataframe  
    df = CSV.read(sample_csv_data, DataFrame) (1)  
  
    # map over each row and construct an array of Policy objects  
    map(eachrow(df)) do row  
        Policy(
```

```

    row.id,
    row.sex == "M" ? Male : Female,
    row.benefit_base,
    row.COLA,
    PaymentMode(row.mode),
    row.issue_date,
    row.issue_age,
    row.risk == "Std" ? Standard : Preferred,
)
end

end

```

- ① CSV.read("sample_inforce.csv",DataFrame) could be used if the data really was in a CSV file named sample_inforce.csv instead of our demonstration IOBuffer.

```

2-element Vector{Policy}:
Policy(1, Male, 100000.0, 0.03, Monthly, Date("1999-12-05"), 30, Standard)
Policy(2, Female, 200000.0, 0.03, Monthly, Date("1999-12-05"), 30, Preferred)

```

25.4. Model Assumptions

25.4.1. Mortality Assumption

MortalityTables.jl provides common life insurance industry tables, and we will use two tables: one each for male and female policies respectively.

```

mort = Dict(
    Male => MortalityTables.table(988).ultimate,                               ①
    Female => MortalityTables.table(992).ultimate,
)

function mortality(pol::Policy, params)
    return params.mortality[pol.sex]
end

```

- ① ultimate refers to not differentiating the mortality by a 'select' underwriting period, which is common but unnecessary for the deomstration in this chapter.

```
mortality (generic function with 1 method)
```

25.5. Model Structure and Mechanics

25.5.1. Core Model Behavior

The overall flow of the model loop will be as follows:

1. Determine some initialized values for each policy at the start of the projection.
2. Step through annual timesteps and simulate whether a death has occurred.
 1. If a death has occurred, log the benefit paid out.
 2. If a death has not occurred keep track of the remaining lives inforce and increment policy values.

The code is shown first and then discussion will follow it:

```
function pol_project!(out, policy, params)
    # some starting values for the given policy
    dur = length(policy.issue_date:Year(1):params.val_date) + 1
    start_age = policy.issue_age + dur - 1
    COLA_factor = (1 + policy.COLA)
    cur_benefit = policy.benefit_base * COLA_factor^(dur - 1)

    # get the right mortality vector
    qs = mortality(policy, params)

    # grab the current thread's id to write to results container
    # without conflicting with other threads
    tid = Threads.threadid()

    ω = lastindex(qs)

    @inbounds for t in 1:min(params.proj_length, ω - start_age)           ①
        q = qs[start_age+t] # get current mortality

        if (rand() < q)
            # if dead then just return and don't increment the results anymore
            out.benefits[t, tid] += cur_benefit                           ②
            return
        else
            # pay benefit, add a life to the output count, and increment the benefit for next
            out.lives[t, tid] += 1
            cur_benefit *= COLA_factor
        end
    end
```

```
    end
end
```

- ① `inbounds` turns off bounds-checking, which makes hot loops faster but first write loop without it to ensure you don't create an error (will crash if you have the error without bounds checking)
- ② Note that the loop is iterating down a column (i.e. across rows) for efficiency (since Julia is column-major).

```
pol_project! (generic function with 1 method)
```

25.5.2. Inputs and Outputs

25.5.2.1. Inputs

The general approach for non-allocating model runs is to provide a previously instantiated container for the function to write the results to. Here, the incoming argument `out(put)` will be a named tuple with associated matrices as the `lives` and `benefits` fields. We know how many periods the model will simulate for and can therefore size the array appropriately at creation.

Other inputs include: `params` which defines some global-like parameters and `policy` which is a single `Policy` object.

i Note

Note that the unit of the core model logic is a single policy. This simplifies the logic and reduces the chance for error due to needing to code for entire arrays of policies at a single time (as would be the case for array oriented programming style, as described in Section 6.5).

25.5.3. Threading

The simulations is using a threaded parallelism approach where it could be operating on any of the computer's available threads. Multi-processor (multi-machine) or GPU-based computation would require some modifications see ([?@sec-parallelism](#)). For the scale and complexity of this example, thread-based parallelism on a single CPU is all one should need for compute.

The threads are handled by distributed the work across threads (this is done by `ThreadsX` in the `foreach` loop below), but we need to write the appropriate place in the matrix so that threads do not compete for the same column in the output data.

25. Stochastic Mortality Projections

Therefore, when we create the `lives` and `benefits` matrices we need to have them sized so that the number of rows is the number of projection periods and the number of columns is the number of threads.

25.5.4. Simulation Control

Parameters for our projection:

```
params = (
    val_date=Date(2021, 12, 31),
    proj_length=100,
    mortality=mort,
)
```

```
(val_date = Date("2021-12-31"), proj_length = 100, mortality = Dict{Sex, OffsetArrays.OffsetVector}
```

Having defined the model behavior at the unit of the policy above, we now need to define how the model should iterate over the entire population of interest. Given a vector of `Policies` in the `policies` argument, the `project` function will:

1. Create output containers, keeping in mind the projection length and number of threads being used.
2. Loop over each policy, letting `ThreadsX.foreach` distribute the work across different threads.
3. Sum up the results across threads via `reduce`.

```
function project(policies, params)
    threads = Threads.nthreads()
    benefits = zeros(params.proj_length, threads)
    lives = zeros(Int, params.proj_length, threads)
    out = (; benefits, lives)
    ThreadsX.foreach(policies) do pol
        pol_project!(out, pol, params)
    end
    map(x → vec(reduce(+, x, dims=2)), out) ①
end
```

① `vec` turns the result into a 1D vector instead of a 1D matrix for later convenience.

```
project (generic function with 1 method)
```

25.6. Running the projection

Example of a single projection:

```
project(repeat(policies, 100_000), params) # <!>
```

```
(benefits = [1.2778972659544663e9, 1.297735757564328e9, 1.426909659059295e9, 1.531826777336355e9,
```

- repeat creates a vector that repeats the two demonstration policies many times.

25.6.1. Stochastic Projection

This defines a loop to calculate the results n times (this is only running the two policies in the sample data n times). This is emulating running our population of policies through n stochastic scenarios, similar to what might be done for a risk or pricing exercise.

```
function stochastic_proj(policies, params, n)

    ThreadsX.map(1:n) do i
        project(policies, params)
    end
end

stochastic_proj (generic function with 1 method)
```

25.6.1.1. Demonstration

We'll simulate the two policies' outcomes 1,000 times and visualize the resulting distribution of claims value:

```
stoch = stochastic_proj(policies, params, 1000)
```

```
1000-element Vector{@NamedTuple{benefits::Vector{Float64}, lives::Vector{Int64}}}:  

(benefits = [0.0, 0.0, 0.0, 215659.12675438428, 0.0, 0.0, 0.0, 485452.4942379325, 0.0, 0.0 ... 0.0  

(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

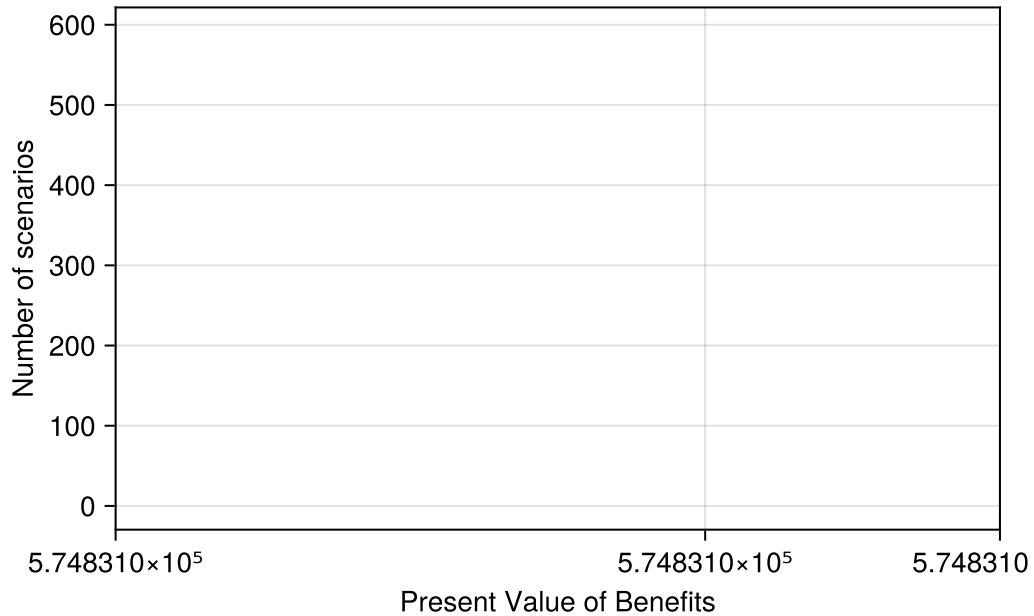
(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

(benefits = [0.0, 203279.4106460404, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0,  

(benefits = [0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0 ... 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0, 0.0
```

25. Stochastic Mortality Projections



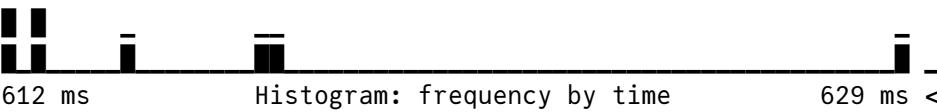
25.7. Benchmarking

Using a 2024 Macbook Air M3 laptop, about 45 million policies able to be stochastically projected per second:

```
policies_to_benchmark = 4_500_000
# adjust the `repeat` depending on how many policies are already in the array
# to match the target number for the benchmark
n = policies_to_benchmark ÷ length(policies)

@benchmark project(p, r) setup = (p = repeat($policies, $n); r = $params)

BenchmarkTools.Trial: 8 samples with 1 evaluation per sample.
Range (min ... max): 611.597 ms ... 628.634 ms | GC (min ... max): 0.00% ... 0.00%
Time (median):      613.131 ms           | GC (median):      0.00%
Time (mean ± σ):   615.495 ms ±  5.711 ms | GC (mean ± σ):  0.00% ± 0.00%

Histogram: frequency by time



Memory estimate: 9.09 KiB, allocs estimate: 68.
```

25.8. Conclusion

This example has worked through a recommended pattern of setting up and running a stochastic simulation using a threaded approach to parallelism. The results show that quite a bit of simulation power is available using even consumer laptop hardware!

26. Scenario Generation

Yun-Tien Lee and Alec Loudenback

“Scenarios are the rehearsal of possible futures. They are not predictions, but pathways that help us prepare for the unknown.” — Peter Schwartz

26.1. In This Chapter

How to generate synthetic data for your model using sub-models, with applications to economic scenario generation and portfolio composition.

26.2. Pseudo-Random Number Generators

Modern computers utilize pseudo-random number generators (PRNGs) to generate random-like numbers. PRNGs are algorithms used to generate sequences of numbers that appear to be random but are actually determined by an initial value, known as the seed. These generators are called “pseudo-random” because the sequences they produce are deterministic; if you provide the same seed, you’ll get the same sequence of numbers. In addition, they have a finite period, which means that after a certain number of generated values, the sequence will repeat. It’s important to choose or design PRNGs with a long enough period for practical applications.

Financial modelers should understand how PRNGs work because many financial models rely on Monte Carlo simulations, risk analysis, and other stochastic modeling techniques that require random sampling. A good PRNG is essential for robust financial modeling. Choosing the right PRNG ensures accurate, reproducible, and unbiased results with efficiency, which is critical when making financial decisions.

26. Scenario Generation

26.2.1. Common PRNGs

26.2.1.1. Mersenne Twister

One of the strengths of the Mersenne Twister is its exceptionally long period. The period is $2^{19937} - 1$, which means it can generate $2^{19937} - 1$ pseudo-random numbers before repeating. This long period is crucial for applications requiring a large number of independent random numbers. It is also known for its good statistical properties. It passes many standard tests for randomness and provides a relatively uniform distribution of random numbers. Moreover, it is designed to allow multiple independent instances to be used concurrently without interfering with each other. This makes it suitable for parallel computing. Although there are faster generators for specific use cases, the Mersenne Twister is still often favored for its balance between speed and quality, and has been one of the recommended PRNGs for financial modeling purposes.

26.2.1.2. Xorshift

Xorshift is a family of PRNGs known for their simplicity and relatively fast operation. The name “xorshift” comes from the bitwise XOR (exclusive or) and bit-shifting operations that are the core of the algorithm. Xorshift generators are often used in applications where speed is a priority and cryptographic-strength randomness is not a strict requirement. Xorshift PRNGs use bitwise XOR, left shifts, and right shifts to update the internal state and generate pseudo-random numbers. The basic idea is to repeatedly apply these operations to the state to produce a sequence of numbers. The period of a typical xorshift generator is relatively short compared to some other PRNGs like the Mersenne Twister. However, there are variations of xorshift algorithms that can have longer periods. One of the main advantages of xorshift is its simplicity and speed. The bitwise XOR and bit-shifting operations can be efficiently implemented in hardware, making xorshift generators suitable for applications where fast random number generation is crucial.

26.2.1.3. Xoshiro

Xoshiro is a family of PRNGs known for their high performance and good statistical properties. The name “Xoshiro” is derived from the Japanese word “xoroshiro,” meaning “random.” Xoshiro algorithms, including Xoshiro128 and others, use a combination of bitwise XOR, bit-shifting, and addition operations. They often have more complex update rules than basic Xorshift algorithms. In addition, they typically have longer periods, making them suitable for applications that require more pseudo-random numbers before repetition.

26.2.2. Consistent Interface

Julia offers a consistent interface for random numbers due to its design and multiple dispatch principles. Consider the following random numbers in different data types.

```
using Random
```

```
rng = MersenneTwister(1234)
rand(Int, (2, 3))
```

2×3 Matrix{Int64}:

| | | |
|----------------------|----------------------|----------------------|
| -2988873214447539226 | -1606048053734002818 | 2884491630289756814 |
| 8541285957272188319 | -2756444082977533404 | -4506841968268069559 |

```
using Random
```

```
rng = MersenneTwister(1234)
rand(Float64, (2, 3))
```

2×3 Matrix{Float64}:

| | | |
|----------|----------|----------|
| 0.338631 | 0.711375 | 0.372296 |
| 0.927057 | 0.323612 | 0.241972 |

```
using Random
```

```
rng = Xoshiro(1234)
rand(Bool, (2, 3))
```

2×3 Matrix{Bool}:

| | | |
|---|---|---|
| 0 | 1 | 1 |
| 1 | 0 | 1 |

26.3. Common Use Cases of Scenario Generators

Scenario generators are widely used in risk management, investment analysis, and regulatory compliance to model potential future outcomes. If the goal is forecasting actual market behavior, real world scenarios (RW) are commonly used. If, on the other hand, pricing financial instruments is needed, risk neutral (RN) scenarios are often used.

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26.3.1. Risk management, especially Value at Risk (VaR) & Expected Shortfall (ES).

RW scenario generators are used to simulate market movements to estimate potential portfolio losses. Basel III regulatory capital requirements have adopted these approaches.

26.3.2. Stress Testing & Regulatory Compliance

RW scenario generators can also be used to generate extreme but plausible market conditions to assess resilience, which is required by central banks and financial regulators (e.g., Federal Reserve and ECB).

26.3.3. Portfolio Optimization & Asset Allocation

RW scenario generators are used to simulate thousands of market conditions to determine optimal portfolio allocations which is commonly used in modern portfolio theory (MPT) and Black-Litterman models.

26.3.4. Pension & Insurance Risk Modeling

RW scenario generators can be used to simulate longevity risk, policyholder behavior, and interest rate movements. They are also used for economic capital estimation under uncertain economic scenarios.

26.3.5. Economic & Macro-Financial Forecasting

Central banks and institutions (e.g., IMF, World Bank) use RW scenario generators to predict macroeconomic trends.

26.3.6. Asset Pricing & Hedging

RN scenario generators help value options using stochastic models (e.g., Black-Scholes, Heston model). They can help simulate future stock price movements under different volatility conditions. They can also be used for hedging purposes to test how a portfolio performs under different inflation, interest rate, or commodity price scenarios.

26.4. Common Economic Scenario Generation Approaches

26.3.7. Fixed Income & Interest Rate Modeling

Yield curve modeling uses RN scenarios to value bonds and interest rate derivatives. Swaps, swaptions, and credit default swaps (CDS) also rely on RN pricing. RN scenario generators can also simulate yield curves for bond and fixed-income pricing. Models like Cox-Ingersoll-Ross (CIR) or Hull-White generate future interest rate paths.

26.3.8. Regulatory & Accounting Valuations

IFRS 13 & fair value accounting uses RN models to determine the market-consistent value of liabilities. Solvency II for insurers asks valuation of policyholder guarantees using RN scenarios.

26.4. Common Economic Scenario Generation Approaches

Economic scenario generation involves the development of plausible future economic scenarios to assess the potential impact on financial portfolios, investments, or decision-making processes. Various approaches are used to generate economic scenarios, such as adapting underlying stochastic differential equations (SDEs) for Monte Carlo scenario generation techniques.

26.4.1. Interest Rate Models

26.4.1.1. Vasicek and Cox Ingersoll Ross (CIR)

The Vasicek model is a one-factor model commonly used for simulating interest rate scenarios. It describes the dynamics of short-term interest rates using a stochastic differential equation (SDE). In a Monte Carlo simulation, we can use the Vasicek model to generate multiple interest rate paths. The CIR model is an extension of the Vasicek model with non-constant volatility. It addresses the issue of negative interest rates by ensuring that interest rates remain positive. Vasicek is defined as

$$dr(t) = \kappa(\theta - r(t)) dt + \sigma dW(t)$$

where

- $r(t)$ is the short-term interest rate at time t .
- κ is the speed of mean reversion, representing how quickly the interest rate reverts to its long-term mean.
- θ is the long-term mean or equilibrium level of the interest rate.

26. Scenario Generation

- σ is the volatility of the interest rate.
- $dW(t)$ is a Wiener process or Brownian motion, representing a random shock.

And CIR is defined as

$$dr(t) = \kappa(\theta - r(t)) dt + \sigma\sqrt{r(t)} dW(t)$$

where

- $r(t)$ is the short-term interest rate at time t .
- κ is the speed of mean reversion, representing how quickly the interest rate reverts to its long-term mean.
- θ is the long-term mean or equilibrium level of the interest rate.
- σ is the volatility of the interest rate.
- $dW(t)$ is a Wiener process or Brownian motion, representing a random shock.

The following code shows a simplified implementation of a CIR model. The specification of dr can be changed to make it a Vasicek model.

```
using Random, CairoMakie

# Set seed for reproducibility
Random.seed!(1234)

# CIR model parameters
κ = 0.2      # Speed of mean reversion
θ = 0.05     # Long-term mean
σ = 0.1      # Volatility

# Initial short-term interest rate
r₀ = 0.03

# Number of time steps and simulations
num_steps = 252
num_simulations = 1_000

# Time increment
Δt = 1 / 252

# Function to simulate CIR process
function cir_simulation(κ, θ, σ, r₀, Δt, num_steps, num_simulations)
    interest_rate_paths = zeros(num_steps, num_simulations)
    for j in 1:num_simulations
        interest_rate_paths[1, j] = r₀
```

26.4. Common Economic Scenario Generation Approaches

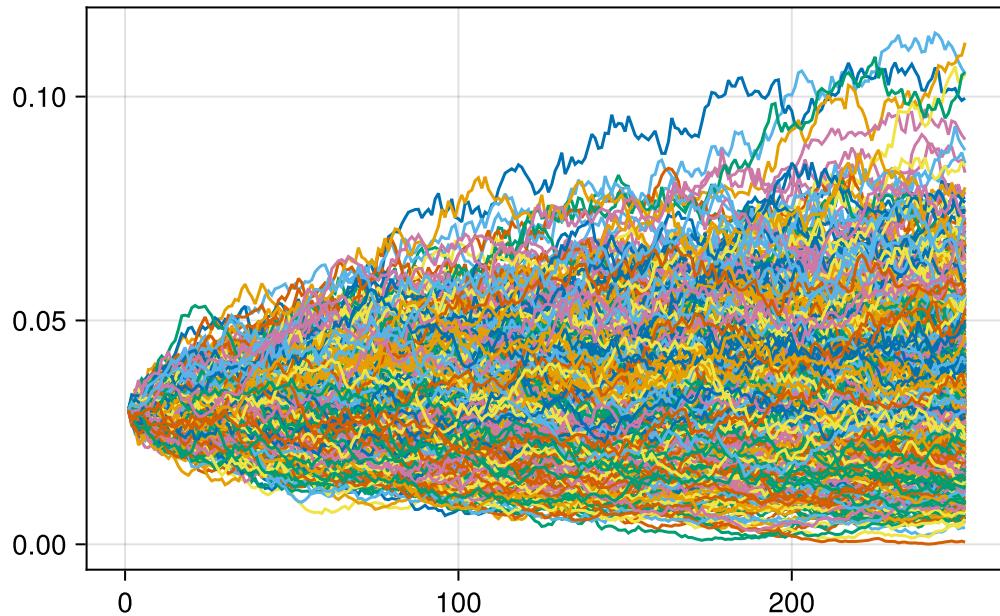
```

for i in 2:num_steps
    dW = randn() * sqrt(Δt)
    # for Vasicek
    # dr = κ * (θ - interest_rate_paths[i-1, j]) * Δt + σ * dW
    dr = κ * (θ - interest_rate_paths[i-1, j]) * Δt + σ * sqrt(interest_rate_paths[i-1, j])
    interest_rate_paths[i, j] = max(interest_rate_paths[i-1, j] + dr, 0) # Ensure non-negative
end
end
return interest_rate_paths
end

# Run CIR simulation
cir_paths = cir_simulation(κ, θ, σ, r₀, Δt, num_steps, num_simulations)

# Plot the simulated interest rate paths
f = Figure()
Axis(f[1, 1])
for i in 1:num_simulations
    lines!(1:num_steps, cir_paths[:, i])
end
f

```



26. Scenario Generation

26.4.1.2. Hull White

The Hull-White model is a one-factor model that extends the Vasicek model by allowing the mean reversion and volatility parameters to be time-dependent. It is commonly used for pricing interest rate derivatives. Brace-Gatarek-Musiela (BGM) Model extends the Hull-White model to incorporate more factors. It is one of the Libor Market Model (LMM) that describes the evolution of forward rates. It allows for the modeling of both the short-rate and the entire yield curve. It is defined as

$$dr(t) = (\theta(t) - ar(t)) dt + \sigma(t) dW(t)$$

where

- $r(t)$ is the short-term interest rate at time t .
- θ is the long-term mean or equilibrium level of the interest rate.
- a is the speed of mean reversion.
- $\sigma(t)$ is the time-dependent volatility of the interest rate.
- $dW(t)$ is a Wiener process or Brownian motion, representing a random shock.

```
using Random, CairoMakie

# Set seed for reproducibility
Random.seed!(1234)

# Hull-White model parameters
α = 0.1      # Mean reversion speed
σ = 0.02     # Volatility
r₀ = 0.03    # Initial short-term interest rate

# Number of time steps and simulations
num_steps = 252
num_simulations = 1_000

# Time increment
Δt = 1 / 252

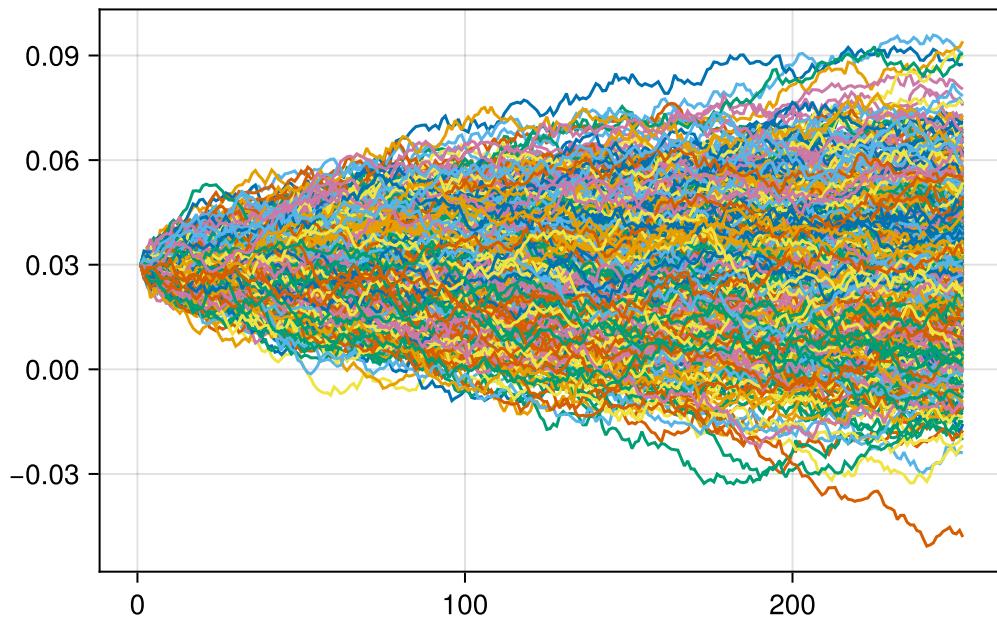
# Function to simulate Hull-White process
function hull_white_simulation(α, σ, r₀, Δt, num_steps, num_simulations)
    interest_rate_paths = zeros(num_steps, num_simulations)
    for j in 1:num_simulations
        interest_rate_paths[1, j] = r₀
        for i in 2:num_steps
            dW = randn() * sqrt(Δt)
```

26.4. Common Economic Scenario Generation Approaches

```
dr = α * (σ - interest_rate_paths[i-1, j]) * Δt + σ * dW
interest_rate_paths[i, j] = interest_rate_paths[i-1, j] + dr
end
end
return interest_rate_paths
end

# Run Hull-White simulation
hull_white_paths = hull_white_simulation(α, σ, r₀, Δt, num_steps, num_simulations)

# Plot the simulated interest rate paths
f = Figure()
Axis(f[1, 1])
for i in 1:num_simulations
    lines!(1:num_steps, hull_white_paths[:, i])
end
f
```



26. Scenario Generation

26.4.2. Equity Models

26.4.2.1. Geometric Brownian Motion (GBM)

GBM is a stochastic process commonly used to model the price movement of financial instruments, including stocks. It assumes constant volatility and is characterized by a log-normal distribution. It is defined as

$$dS(t) = \mu S(t) dt + \sigma S(t) dW(t)$$

where

- $S(t)$ is the stock price at time t .
- μ is the drift coefficient (expected return).
- σ is the volatility coefficient.
- $dW(t)$ is a Wiener process or Brownian motion, representing a random shock.

```
using Random, CairoMakie

# Set seed for reproducibility
Random.seed!(1234)

# GBM parameters
μ = 0.05      # Drift (expected return)
σ = 0.2       # Volatility

# Initial stock price
S₀ = 100

# Number of time steps and simulations
num_steps = 252
num_simulations = 1_000

# Time increment
Δt = 1 / 252

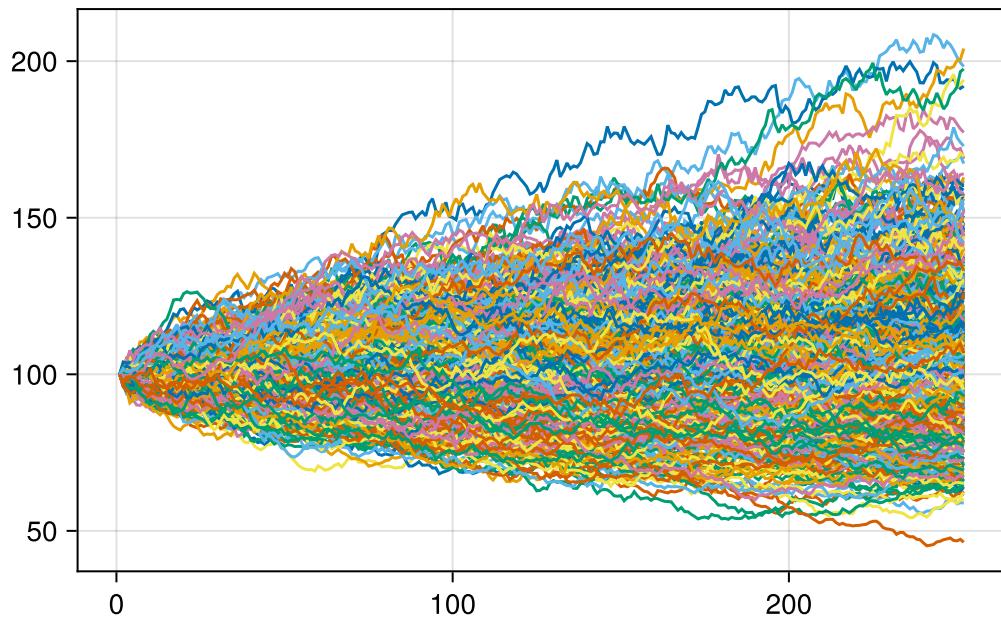
# Function to simulate GBM
function gbm_simulation(μ, σ, S₀, Δt, num_steps, num_simulations)
    stock_price_paths = zeros(num_steps, num_simulations)
    for j in 1:num_simulations
        stock_price_paths[1, j] = S₀
        for i in 2:num_steps
            dW = randn() * sqrt(Δt)
```

26.4. Common Economic Scenario Generation Approaches

```
S = stock_price_paths[i-1, j]
dS = μ * S * Δt + σ * S * dW
stock_price_paths[i, j] = S + dS
end
end
return stock_price_paths
end

# Run GBM simulation
gbm_paths = gbm_simulation(μ, σ, S₀, Δt, num_steps, num_simulations)

# Plot the simulated stock price paths
f = Figure()
Axis(f[1, 1])
for i in 1:num_simulations
    lines!(1:num_steps, gbm_paths[:, i])
end
f
```



26.4.2.2. Generalized Autoregressive Conditional Heteroskedasticity (GARCH)

GARCH models capture time-varying volatility. They are often used in conjunction with other models to forecast volatility. It is defined as

26. Scenario Generation

$$\sigma_t^2 = \omega + \alpha_1 r_{t-1}^2 + \beta_1 \sigma_{t-1}^2$$

$$r_t = \varepsilon_t \sqrt{\sigma_t^2}$$

- σ_t^2 is the conditional variance at time t
- r_t is the return at time t
- ε_t is a white noise or innovation process
- $\omega, \alpha_1, \beta_1$ are model parameters

```
using Random, CairoMakie

# Set seed for reproducibility
Random.seed!(1234)

# GARCH(1,1) parameters
α₀ = 0.01      # Constant term
α₁ = 0.1       # Coefficient for lagged squared returns
β₁ = 0.8       # Coefficient for lagged conditional volatility

# Number of time steps and simulations
num_steps = 252
num_simulations = 1_000

# Time increment
Δt = 1 / 252

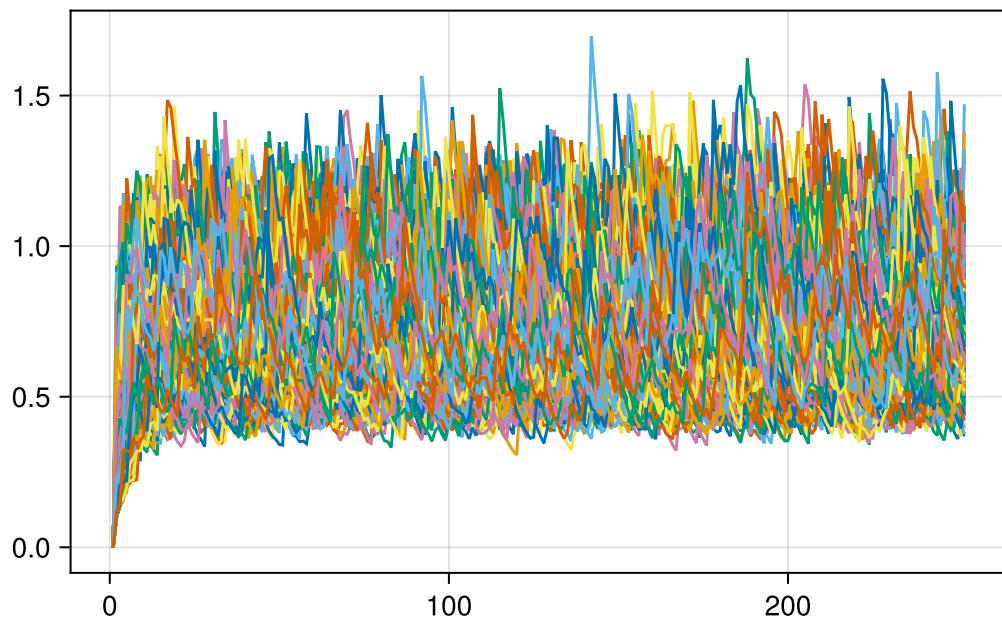
# Function to simulate GARCH(1,1) volatility
function garch_simulation(α₀, α₁, β₁, num_steps, num_simulations)
    volatility_paths = zeros(num_steps, num_simulations)
    for j in 1:num_simulations
        ε = randn(num_steps)
        squared_returns = zeros(num_steps)
        for i in 2:num_steps
            squared_returns[i] = α₀ + α₁ * ε[i-1]^2 + β₁ * squared_returns[i-1]
            volatility_paths[i, j] = sqrt(squared_returns[i])
        end
    end
    return volatility_paths
end

# Run GARCH simulation
```

26.4. Common Economic Scenario Generation Approaches

```
garch_paths = garch_simulation(α₀, α₁, β₁, num_steps, num_simulations)

# Plot the simulated volatility paths
f = Figure()
Axis(f[1, 1])
for i in 1:num_simulations
    lines!(1:num_steps, garch_paths[:, i])
end
f
```



26.4.3. Copulas

Simulating data using copulas involves generating multivariate samples with specified marginal distributions and a copula structure.

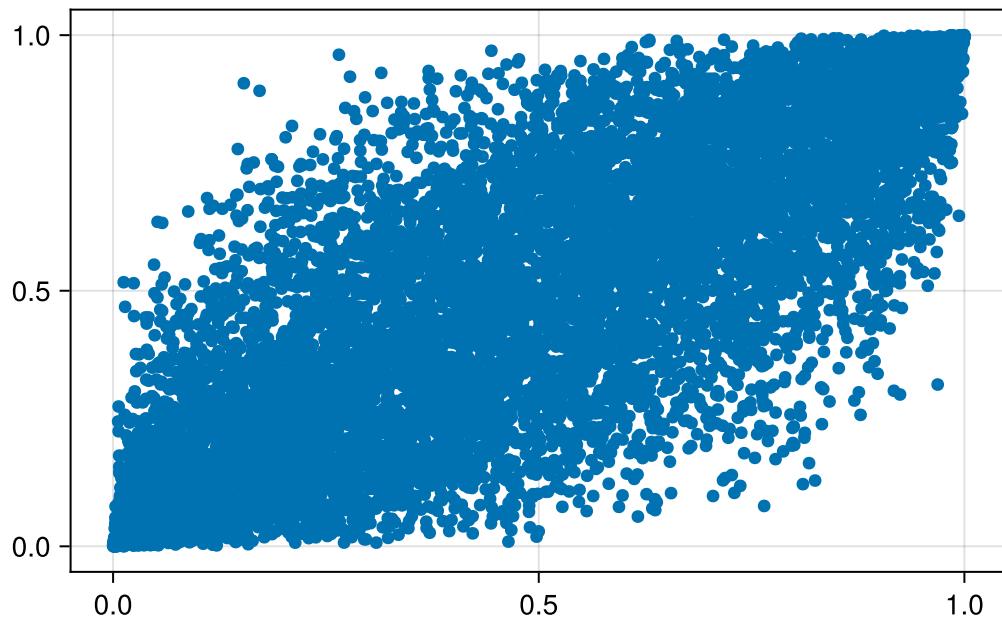
```
using Random, CairoMakie, BivariateCopulas

# Set seed for reproducibility
Random.seed!(1234)

# Generate a Gaussian copula
gaussian_copula = Gaussian(0.8)
```

26. Scenario Generation

```
# Show simulated copula
f = scatter(rand(gaussian_copula, 10^4))
f
```



Copulas can also be used to infer combined distributions from data samples.

```
using Copulas, Distributions, Random

X₁ = Gamma(2, 3)
X₂ = Pareto()
X₃ = LogNormal(0, 1)
C = ClaytonCopula(3, 0.7) # A 3-variate Clayton Copula with θ = 0.7
D = SklarDist(C, (X₁, X₂, X₃)) # The final distribution

# Generate a dataset
simu = rand(D, 1000)
# We may estimate a copula, or get parameters of underlying distributions, using the 'fit' function
D̂ = fit(SklarDist{ClaytonCopula,Tuple{Gamma,Normal,LogNormal}}, simu)

Copulas.SklarDist{Copulas.ClaytonCopula{3, Float64}, Tuple{Distributions.Gamma{Float64}, Distr...}
C: Copulas.ClaytonCopula{3, Float64}(
G: Copulas.ClaytonGenerator{Float64}(0.7255762179151387)
)
```

26.4. Common Economic Scenario Generation Approaches

```
m: (Distributions.Gamma{Float64}(α=1.9509359315325794, θ=3.0668504198367565), Distributions.Nor
```


27. Similarity Analysis

Yun-Tien Lee

“By understanding the similarities, we unlock the potential for new insights, as patterns across different contexts often reveal hidden truths.” — Unknown

27.1. In This Chapter

Given a set of interest, understanding the relative similarity (or not) of features of interest is useful in classification and data compression techniques.

27.2. The Data

Stored data can generally be categorized into two formats: tabular (structured) and non-tabular (unstructured). Structured data format is a structured way of organizing and presenting data in rows and columns, resembling a table. This format is widely used for storing and representing structured datasets, making it easy to read, analyze, and manipulate data. The most common example of structured data is a spreadsheet, where data is organized into rows and columns. Structured data can also be stored in relational databases for easier lookups and matching. On the other hand, unstructured data refers to data that lacks a predefined data model or structure. Unlike structured data, which fits neatly into tables or databases, unstructured data does not have a predefined schema. It can include text documents, images, audio files, video files, social media posts, and more.

Structured data can be further categorized into numerical and categorical data based on the types of values they represent. The following data tables will be referenced throughout the chapter. Real numerical data can easily be converted or normalized to a series of floating points, and real categorical data to a series of binary literals through one-hot encoding procedures.

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```
sample_csv_data =
    IOBuffer(
        raw"id,sex,benefit_base,education,occupation,issue_age
1,M,100000.0,college,1,30.0
2,F,200000.0,master,3,20.0
3,M,150000.0,high_school,4,40.0
4,F,50000.0,college,2,60.0
5,M,250000.0,college,1,40.0
6,F,200000.0,high_school,2,30.0"
    )

IOBuffer(data=UInt8[...], readable=true, writable=false, seekable=true, append=false, size=278,
using CSV, DataFrames, TableTransforms

df = CSV.read(sample_csv_data, DataFrame)
df_num = apply(MinMax(), df[:, [:benefit_base, :issue_age]])[1]



	benefit_base	issue_age
	Float64	Float64
1	0.25	0.25
2	0.75	0.0
3	0.5	0.5
4	0.0	1.0
5	1.0	0.5
6	0.75	0.25



using StatsBase

arr_cat = hcat(indicatormat(df.sex)', indicatormat(df.education)', indicatormat(df.occupation)

6x9 Matrix{Bool}:
 0  1  1  0  0  1  0  0  0
 1  0  0  0  1  0  0  1  0
 0  1  0  1  0  0  0  0  1
 1  0  1  0  0  0  1  0  0
 0  1  1  0  0  1  0  0  0
 1  0  0  1  0  0  1  0  0
```

For unstructured data, due to the nature of their variety, the choice of representation depends on the type of data and the specific task at hand. For text data, a Word2Vec embedding is commonly used, while Convolutional Neural Networks (CNNs) are for image data and wave transforms are for audio data. No matter which transformation is applied, unstructured data can generally be converted to a series of floating points, just like numerical structured data.

27.3. Common Similarity Measures

The following measures are commonly used to calculate similarities.

27.3.1. Euclidean Distance (L2 norm)

Euclidean distance, also known as the L2 norm, is defined as

$$d = \sqrt{\sum_{i=1}^n (w_i - v_i)^2}$$

The distance is usually meaningful when applied to numerical data. The following Julia code shows the Euclidean distance for the first two rows in df_num.

```
using LinearAlgebra

#d12 = √(Σ((Array(df_num[1, :]) .- Array(df_num[2, :])) .* (Array(df_num[1, :]) .- Array(df_num[2, :])))
d12 = LinearAlgebra.norm(Array(df_num[1, :]) .- Array(df_num[2, :]))

0.5590169943749475
```

27.3.2. Manhattan Distance (L1 Norm)

Manhattan distance, also known as the L1 norm, is defined as

$$d = \sum_{i=1}^n |w_i - v_i|$$

The distance is also usually meaningful when applied to numerical data. The following Julia code shows the Euclidean distance for the first two rows in df_num.

```
using LinearAlgebra

#d12 = Σ(abs.(Array(df_num[1, :]) .- Array(df_num[2, :])))
d12 = LinearAlgebra.norm1(Array(df_num[1, :]) .- Array(df_num[2, :]))

0.75
```

27. Similarity Analysis

27.3.3. Cosine Similarity

Cosine similarity is defined as

$$d = \frac{\sum_{i=1}^n w_i \cdot v_i}{\sqrt{\sum_{i=1}^n w_i^2} \cdot \sqrt{\sum_{i=1}^n v_i^2}}$$

The distance would be meaningful when applied to both numerical and categorical data.

The following Julia code shows the cosine similarity for the first two rows in df_num.

```
using LinearAlgebra

d12 = (Array(df_num[1, :]) · Array(df_num[2, :])) / norm(df_num[1, :]) / norm(df_num[2, :])

0.7071067811865475
```

The following Julia code shows the cosine similarity for the first and the third rows in arr_cat.

```
using LinearAlgebra

d13 = (arr_cat[1, :] · arr_cat[3, :]) / norm(arr_cat[1, :]) / norm(arr_cat[3, :])

0.3333333333333337
```

Note how similar the syntax of processing for numerical or categorical data is. Multiple dispatch allows Julia to identify most efficient underlying procedure for different types of data. For categorical data, the *dot* operation on binary vectors is essentially count of 1's, while for numerical data it is the *dot* operation for most numerical processing libraries.

27.3.4. Jaccard Similarity

Jaccard similarity is defined as

$$d = \frac{|W \cap V|}{|W \cup V|}$$

The distance is usually meaningful when applied to categorical data. The following Julia code shows the Jaccard similarity for the first and the third rows in arr_cat.

```
d13 = (arr_cat[1, :] ∙ arr_cat[3, :]) / sum(arr_cat[1, :] .| arr_cat[3, :])

0.2
```

27.4. *k*-Nearest Neighbor (*kNN*) Clustering

27.3.5. Hamming Distance

Hamming distance is defined as $d = \text{Number of positions at which } w \text{ and } v \text{ differ}$. The distance is usually meaningful when applied to categorical data. The following Julia code shows the Hamming distance for the first and the third rows in arr_cat.

```
d13 = sum(arr_cat[1, :] .⊤ arr_cat[3, :])
```

4

27.4. *k*-Nearest Neighbor (*kNN*) Clustering

kNN is primarily known as a classification algorithm, but it can also be used for clustering, particularly in the context of density-based clustering. Density-based clustering identifies regions in the data space where the density of data points is higher, and it groups points in these high-density regions. The core idea of *kNN* clustering is to assign each data point to a cluster based on the density of its neighbors. A data point becomes a core point if it has at least a specified number of neighbors within a certain distance.

```
using Random, NearestNeighbors, CairoMakie

# Step 1: Generate synthetic data
Random.seed!(1234) # For reproducibility
data = rand(10, 2) # 10 points with 2 dimensions
println("Dataset:\n", data)

# Step 2: Create a KD-tree for efficient nearest neighbor search
kdtree = KDTree(data)

# Step 3: Define a query point (for which we want to find nearest neighbors)
query_point = [0.5, 0.5]

# Step 4: Find the nearest neighbors
# Specify how many neighbors to find
k = 1
indices, distances = knn(kdtree, query_point, k)

# Step 5: Display the results
println("\nQuery Point: ", query_point)
println("Indices of Nearest Neighbors: ", indices)
println("Distances to Nearest Neighbors: ", distances)
```

27. Similarity Analysis

```
# Step 6: Visualize the points and the query
f = Figure()
axis = Axis(f[1, 1], title="Nearest Neighbors Search")
scatter!(data[:, 1], data[:, 2], label="Data Points", color=:blue)
scatter!([query_point[1]], [query_point[2]], label="Query Point", color=:red, marker=:cross, m=100)
# Highlight nearest neighbors
for idx in indices
    lines!([query_point[1], data[idx, 1]], [query_point[2], data[idx, 2]], color=:black)
end
f
```

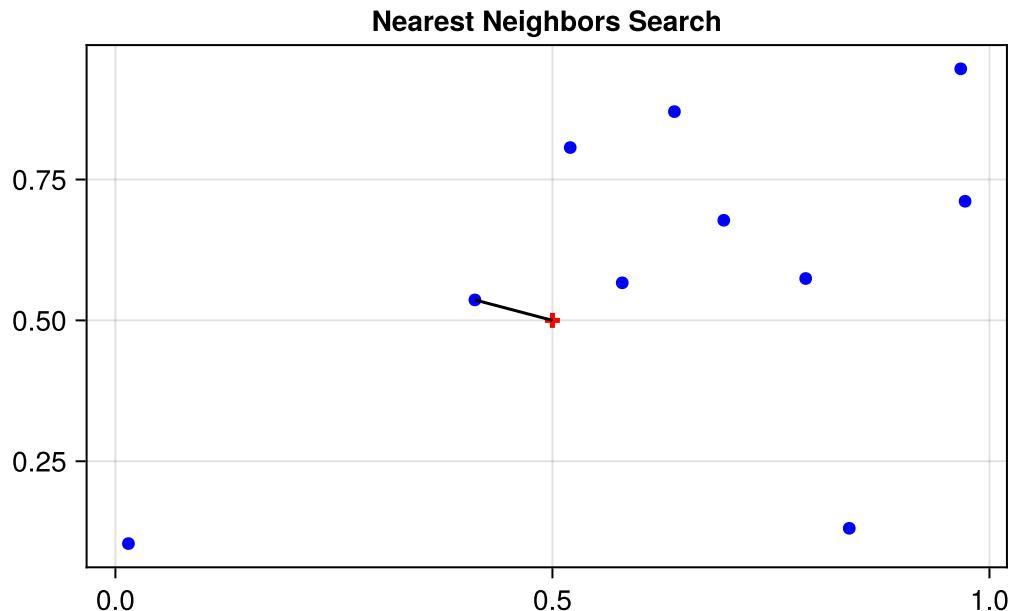
Dataset:

```
[0.5798621201341324 0.5667043025437501; 0.4112941179498505 0.5363685687293304; ... ; 0.78976440953
```

Query Point: [0.5, 0.5]

Indices of Nearest Neighbors: [2]

Distances to Nearest Neighbors: [0.07597457975710152]



28. Sensitivity Analysis

Yun-Tien Lee

“Sensitivity analysis is the art of understanding how small changes in assumptions can lead to big changes in outcomes, helping us navigate uncertainty with clarity.” — Unknown

28.1. In This Chapter

Different approaches to understanding the sensitivity of a model to changes in its inputs: derivatives, finite differences, global sensitivity analysis approaches, and statistical approaches.

28.2. Setup

Let's assume there are certain insurance policies throughout the chapter.

```
using Dates

@enum Sex Female = 1 Male = 2
@enum Risk Standard = 1 Preferred = 2

mutable struct Policy
    id::Int
    sex::Sex
    benefit_base::Float64
    COLA::Float64
    mode::Int
    prem::Float64
    pp::Int
    issue_date::Date
    issue_age::Int
    risk::Risk
end
```

28. Sensitivity Analysis

28.3. The Data

```
using MortalityTables
sample_csv_data =
    IOBuffer(
        raw"id,sex,benefit_base,COLA,mode,prem,pp,issue_date,issue_age,risk
        1,M,100000.0,0.03,1,1000.0,3,1999-12-05,30,Std"
    )

mort = Dict(
    Male => MortalityTables.table(988).ultimate,
    Female => MortalityTables.table(992).ultimate,
)

Dict{Sex, OffsetArrays.OffsetVector{Float64, Vector{Float64}}} with 2 entries:
Male => [0.022571, 0.022571, 0.022571, 0.022571, 0.022571, 0.0225...
Female => [0.00745, 0.00745, 0.00745, 0.00745, 0.00745, 0.00745, 0.0...

using CSV, DataFrames

policies = let

    # read CSV directly into a dataframe
    # df = CSV.read("sample_inforce.csv",DataFrame) # use local string for notebook
    df = CSV.read(sample_csv_data, DataFrame)

    # map over each row and construct an array of Policy objects
    map(eachrow(df)) do row
        Policy(
            row.id,
            row.sex == "M" ? Male : Female,
            row.benefit_base,
            row.COLA,
            row.mode,
            row.prem,
            row.pp,
            row.issue_date,
            row.issue_age,
            row.risk == "Std" ? Standard : Preferred,
        )
    end
end
```

28.3. The Data

1-element Vector{Policy}:

```
Policy(1, Male, 100000.0, 0.03, 1, 1000.0, 3, Date("1999-12-05"), 30, Standard)
```

Given a basic insurance product, a pure whole of life (WOL) policy with level benefits and level premiums payable within the first 10 years, the reserve at the end of the y^{th} policy year is defined by

$$res(y) = \sum_{t=age+y}^{120} (sur_{t-age-y} * mort_t * B_y * \sqrt{1+r}) - (P_y * sur_{t-age-y})$$

where

- $mort_t$ is the mortality at age t
- p_y is the survival probability adjusted with COLA, with values of
 - $p_{y-1} = 1$,
 - $p_x = p_{x-1} * (1 - mort_{age+y}) / (1 + COLA)$ for $x \geq y$, and
 - 0 for $x < y - 1$ or $age + x \geq 120$, or ultimate age of the current mortality table
- B_y is the level benefit throughout the policy
- P_y is the level premium within the first 10 policy years which is 0 for policy years after 10
- r is the level interest rate throughout the policy

```
function sur(y::Int, pol::Policy)
    if y == 0
        1
    elseif y < 0 || 120 - y <= pol.issue_age
        0
    else
        sur(y - 1, pol) * (1 - mort[pol.sex][pol.issue_age+y]) / (1 + pol.COLA)
    end
end

function res(y::Int, pol::Policy)
    s = 0.0
    if y >= 1 && y <= 120 - pol.issue_age
        for t in (pol.issue_age+y):120
            prem = 0.0
            if y <= pol.pp
                prem = pol.prem
            end
            s += sur(t - pol.issue_age - y, pol) * mort[pol.sex][t] * pol.benefit_base - prem
        end
    end
    s
end
```

28. Sensitivity Analysis

```
    end
end
s
end

res (generic function with 1 method)
```

28.4. Common Sensitivity Analysis Methodologies

28.4.1. Finite Differences

Define a customized finite difference function with respect to the COLA, rippled by a small difference.

```
function res_wrt_r_fd(y::Int, pol::Policy, r::Float64, h=1e-3)
    p_+, p_- = deepcopy(pol), deepcopy(pol)
    p_+.COLA, p_-.COLA = r + h, r - h
    (res(y, p_-) - res(y, p_+)) / (2res(y, pol))
end

res_wrt_r_fd(2, policies[1], 0.03) # changes in reserve at year 2 when the interest rate at 3%
```

0.021366520936389077

28.4.2. Regression Analyses

```
using GlobalSensitivity

function r1_wrt_r(r)
    p = deepcopy(policies[1])
    p.COLA = r[2]
    p.prem = r[3]
    res(Int(floor(r[1])), p)
end

# reserve @ year 1/2, interest rate @ 0.03 ± 0.01, prem @ 1000.0 ± 0.1
reg_anal = gsa(r1_wrt_r, RegressionGSA(), [[1, 2], [0.029, 0.031], [999.9, 1000.1]], samples=1000)
@show reg_anal.pearson

reg_anal.pearson = [0.002266826033036182 -0.9999603237826435 -0.003255946696849575]
```

28.4. Common Sensitivity Analysis Methodologies

```
1×3 Matrix{Float64}:
0.00226683 -0.99996 -0.00325595
```

The Pearson Spearman coefficients show the correlation coefficient matrix between inputs and outputs.

28.4.3. Sobol Indices

Sobol is a variance-based method, and it decomposes the variance of the output of the model or system into fractions which can be attributed to inputs or sets of inputs. This helps to get not just the individual parameter's sensitivities, but also gives a way to quantify the affect and sensitivity from the interaction between the parameters.

$$Y = f_0 + \sum_{i=1}^d f_i(X_i) + \sum_{i < j}^d f_{ij}(X_i, X_j) + \dots + f_{1,2,\dots,d}(X_1, X_2, \dots, X_d)$$

$$Var(Y) = \sum_{i=1}^d V_i + \sum_{i < j}^d V_{ij} + \dots + V_{1,2,\dots,d}$$

The Sobol Indices are “ordered”, the first order indices given by $S_i = \frac{V_i}{Var(Y)}$, the contribution to the output variance of the main effect of X_i . Therefore, it measures the effect of varying X_i alone, but averaged over variations in other input parameters. It is standardized by the total variance to provide a fractional contribution. Higher-order interaction indices S_{ij}, S_{ijk} and so on can be formed by dividing other terms in the variance decomposition by $Var(Y)$.

```
using QuasiMonteCarlo, GlobalSensitivity
```

```
# reserve @ year 1/2, interest rate @ 0.03 ± 0.01, prem @ 1000.0 ± 0.1
L, U = QuasiMonteCarlo.generate_design_matrices(1000, [1, 0.029, 999.9], [2, 0.031, 1000.1], S
s = gsa(r1_wrt_r, Sobol()), L, U)
@show s.S1
@show s.ST
```

```
Warning: The `generate_design_matrices(n, d, sampler, R = NoRand(), num_mats)` method does not p
└── Prefer using randomization methods such as `R = Shift()`, `R = MatousekScrambling()`, etc., se
└── @ QuasiMonteCarlo ~/.julia/packages/QuasiMonteCarlo/KvLfb/src/RandomizedQuasiMonteCarlo/itera
s.S1 = [0.0, 1.2308143537488936, -0.00010730838653271882]
s.ST = [0.0, 1.0014202549551285, 5.83151055643934e-6]
```

28. Sensitivity Analysis

```
3-element Vector{Float64}:
0.0
1.0014202549551285
5.83151055643934e-6
```

The output shows the first order and total order of variations in different input parameters.

28.4.4. Morris Method

The Morris method also known as Morris's OAT method where OAT stands for One At a Time can be described in the following steps:

$$EE_i = \frac{f(x_1, x_2, \dots, x_i + \Delta, \dots, x_k) - y}{\Delta}$$

We calculate local sensitivity measures known as “elementary effects”, which are calculated by measuring the perturbation in the output of the model on changing one parameter.

These are evaluated at various points in the input chosen such that a wide “spread” of the parameter space is explored and considered in the analysis, to provide an approximate global importance measure. The mean and variance of these elementary effects is computed. A high value of the mean implies that a parameter is important, a high variance implies that its effects are non-linear or the result of interactions with other inputs. This method does not evaluate separately the contribution from the interaction and the contribution of the parameters individually and gives the effects for each parameter which takes into consideration all the interactions and its individual contribution.

```
using GlobalSensitivity

# reserve @ year 1/2, interest rate @ 0.03 ± 0.01, prem @ 1000.0 ± 0.1
m = gsa(r1_wrt_r, Morris(), [[1, 2], [0.029, 0.031], [999.9, 1000.1]])
@show m.means
@show m.variances

m.means = [4977.452086301237 -725208.084911562 -17.279540384169234]
m.variances = [6.191280398282986e8 2.2266513365354595e8 0.016318702371709272]

1x3 Matrix{Float64}:
 6.19128e8  2.22665e8  0.0163187
```

From the means it can be observed which variables are more important, and the variances imply higher degree of nonlinearity or interactions with other variables.

28.4.5. Fourier Amplitude Sensitivity Tests

FAST offers a robust, especially at low sample size, and computationally efficient procedure to get the first and total order indices as discussed in Sobol. It utilizes monodimensional Fourier decomposition along a curve, exploring the parameter space. The curve is defined by a set of parametric equations,

$$EE_i = \frac{f(x_1, x_2, \dots, x_i + \Delta, \dots, x_k) - y}{\Delta}$$

where s is a scalar variable varying over the range $-\infty < s < +\infty$, G_i are transformation functions and $w_i, \forall i = 1, 2, \dots, N$ is a set of different (angular) frequencies, to be properly selected, associated with each factor for all N (samples) number of parameter sets.

```
using GlobalSensitivity
```

```
# reserve @ year 1/2, interest rate @ 0.03 ± 0.01, prem @ 1000.0 ± 0.1
fast = gsa(r1_wrt_r, eFAST(), [[1, 2], [0.029, 0.031], [999.9, 1000.1]], samples=1000)
@show fast.S1
@show fast.ST

fast.S1 = [5.131843109942656e-12 0.997696367809449 5.796050388482397e-6]
fast.ST = [7.349891162355959e-7 0.999993803041509 0.0023004274889842913]

1x3 Matrix{Float64}:
 7.34989e-7  0.999994  0.00230043
```

The output shows the first order and total order of variations in different input parameters.

28.4.6. Automatic Differentiation

By applying the chain rule repeatedly on elementary operations of computer calculations, automatic differentiation can be applied to measure impacts of small differences. More details in Chapter 16 on automatic differentiation.

28. Sensitivity Analysis

28.4.7. Scenario Analyses

Scenarios can be generated following scenario generation methodologies to evaluate impacts. More details in [?@sec-scegen](#) on scenario generation.

When scenarios are generated to evaluate sensitivities, one may need to take the following into consideration.

- Reverse stress testing. Reverse stress testing in scenario analysis involves identifying extreme scenarios that could potentially lead to catastrophic outcomes for a financial institution or a system. Unlike traditional sensitivity testing to simulate the impact of adverse events on the system, reverse stress testing starts with a catastrophic outcome and works backwards to determine the combination of events or circumstances that could lead to such an outcome.

One typically follows these steps to do reverse stress testing. – Define a critical failure point (e.g., bankruptcy, system outage, regulatory breach). – Analyze the combinations of events or variables that could cause the failure. – Model the path from normal conditions to the adverse outcome.

Potential benefits that reverse stress testing could bring include: – Focusing on Vulnerabilities: Highlights specific scenarios to avoid at all costs. – Enhancing Resilience: Strengthens systems against extreme risks. – Regulatory Compliance: Often required in highly regulated industries like banking and energy.

- Stylistic scenarios. Developing stylistic scenarios in scenario analysis involves creating narratives or storylines that describe plausible future states or situations. These scenarios are crafted to capture key uncertainties, trends, and factors that could significantly impact the organization, industry, or environment under study.
- Backtesting against historical data. Backtesting in scenario analysis involves an iterative process of using past data to validate the effectiveness and accuracy of scenarios developed for forecasting future outcomes. Scenarios are first defined and applied on selective historical data, and refined after any discrepancies of scenario outcomes versus historical results are identified.

One typically follows these steps to do backtesting against historical data. – Define Scenarios: Establish hypothetical scenarios (e.g., market crashes, changes in interest rates, or operational disruptions), and ensure scenarios cover a range of possibilities, such as best-case, base-case, and worst-case scenarios. – Collect Historical Data: Gather relevant historical data for key variables (e.g., stock prices, interest rates, production metrics), and ensure data spans periods where similar events occurred in the past. – Model Scenario Impacts: Use historical data to simulate the impacts of the scenarios on key metrics or performance indicators. – Compare Results: Compare the modeled results of the scenarios with the actual historical outcomes, and assess how well the scenarios

28.4. Common Sensitivity Analysis Methodologies

predict or explain the observed data. – Adjust and Refine: If the scenarios do not align with historical outcomes, refine the assumptions or parameters in the scenario models, and incorporate lessons learned from historical patterns to improve future scenario analyses.

Some considerations in incorporating historical data. – Data Quality: Ensure historical data is accurate, complete, and relevant to the scenarios being tested. – Model Limitations: Scenario models are based on assumptions that might not fully capture real-world complexities. – Overfitting: Avoid fine-tuning scenarios to perfectly match historical outcomes, as this reduces their applicability to future events. – Changing Dynamics: Historical events may not fully represent future possibilities due to changes in market conditions, regulations, or technology.

29. Portfolio Optimization

Yun-Tien Lee

“Portfolio optimization is not about finding the perfect mix, but about managing risk and return in a way that aligns with your goals and circumstances.”
— Unknown

29.1. In This Chapter

Optimization in a portfolio context with examples of asset selection under different constraints and objectives.

29.2. The Data

```
μ = [0.1, 0.15, 0.12] # returns
ρ = [0.1 0.05 0.03;
      0.05 0.12 0.04;
      0.03 0.04 0.08] # covariances
n_a = length(μ) # number of assets
```

3

29.3. Theory

Harry Markowitz introduced the modern portfolio theory in 1952. The main idea is that investors are pursuing to maximize their expected return of a portfolio given a certain amount of risk. By definition any portfolio yielding a higher return must have higher amount of risk, so there is a trade-off between desired expected returns and allowable risks. The risk versus maximized expected return relationship can be plotted out as a curve, a.k.a. the efficient frontier.

29. Portfolio Optimization

29.4. Mathematical tools

29.4.1. Mean-variance optimization model

Mean-variance optimization is a mathematical framework that seeks to maximize expected returns while minimizing portfolio variance (or standard deviation). It involves calculating the expected return and risk of individual assets and finding the optimal combination of assets to achieve the desired risk-return tradeoff.

$$\begin{aligned} \text{minimize} \quad & w^T \Sigma w \\ \text{subject to} \quad & R^T \geq \mu_{\text{target}} \\ & 1^T w = 1 \\ & w \geq 0 \end{aligned}$$

```
using JuMP, Ipopt
```

```
# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Objective: minimize portfolio variance
@objective(model, Min, sum(w[i] * rho[i, j] * w[j] for i in 1:n_a, j in 1:n_a))
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or positive
@constraint(model, sum(w) == 1)
# May also add additional constraints
# target_return = 0.1
# @constraint(model, dot(mu, w) >= target_return)
# Solve the optimization problem
optimize!(model)
# Print results
@show "Optimal Portfolio Weights:"
for i = 1:n_a
    @show ("Asset ", i, ": ", value.(w)[i])
end
```

```
*****
This program contains Ipopt, a library for large-scale nonlinear optimization.
Ipopt is released as open source code under the Eclipse Public License (EPL).
For more information visit https://github.com/coin-or/Ipopt
*****
```

```
"Optimal Portfolio Weights:" = "Optimal Portfolio Weights:"
("Asset ", i, ":", value.(w)[i]) = ("Asset ", 1, ":", 0.3333333012309821)
("Asset ", i, ":", value.(w)[i]) = ("Asset ", 2, ":", 0.16666675086886984)
("Asset ", i, ":", value.(w)[i]) = ("Asset ", 3, ":", 0.4999999479001481)
```

In mean-variance portfolio optimization, incorporating a cost of risk-based capital on assets is a practical consideration that reflects the additional capital required to support riskier assets in a portfolio. This approach ensures that the optimization process not only maximizes returns relative to risk but also considers the regulatory or internal cost implications associated with holding riskier assets.

$$\begin{aligned} & \text{maximize} && w^T R_{adj} \\ & \text{subject to} && w^T \Sigma w \leq \sigma_{max}^2 \\ & && 1^T w = 1 \\ & && w \geq 0 \end{aligned}$$

where $R_{adj} = [(\mu_1 - \lambda_1), (\mu_2 - \lambda_2), \dots, (\mu_N - \lambda_N)]$ is the adjusted expected returns.

using JuMP, Ipopt

```
# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
r = μ .- [0.01, 0.02, 0.05] # risk adjusted returns
σ²_max = 0.1 # maximum portfolio variance
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Objective: minimize portfolio variance
@objective(model, Max, sum(w[i] * r[i] for i in 1:n_a))
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or positive
@constraint(model, sum(w) == 1)
# Constraints: Sum of allowable portfolio variance is limited
@constraint(model, sum(w[i] * ρ[i, j] * w[j] for i in 1:n_a, j in 1:n_a) <= σ²_max)
# May also add additional constraints
# target_return = 0.1
# @constraint(model, dot(μ, w) >= target_return)
# Solve the optimization problem
optimize!(model)
# Print results
@show "Optimal Portfolio Weights:"
for i = 1:n_a
```

29. Portfolio Optimization

```
@show ("Asset ", i, ": ", value.(w)[i])
end

"Optimal Portfolio Weights:" = "Optimal Portfolio Weights:"
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 1, ": ", 0.16666454953827514)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 2, ": ", 0.8333339906581219)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 3, ": ", 1.4598036030472386e-6)
```

29.4.2. Efficient frontier analysis

The efficient frontier represents the set of portfolios that offer the highest expected return for a given level of risk or the lowest risk for a given level of return. Efficient frontier analysis involves plotting risk-return combinations for different portfolios and identifying the optimal portfolio on the frontier.

```
using JuMP, Ipopt, CairoMakie, LinearAlgebra

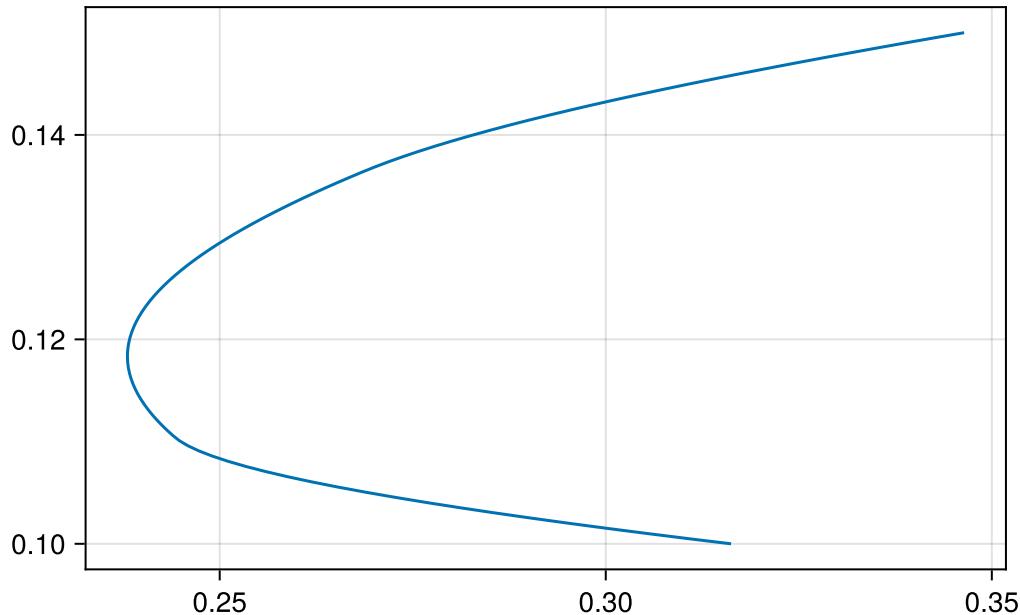
# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Define objective function: minimize portfolio variance
portfolio_variance = w'ρ * w
@objective(model, Min, portfolio_variance)
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or p
@constraint(model, sum(w) == 1)
# Generate a range of target returns
points = 100
target_returns = range(minimum(μ), maximum(μ), length=points)

efficient_frontier = []
for target_return in target_returns
    # Add additional constraint for target return
    @constraint(model, c, dot(μ, w) == target_return)
    # Solve the problem
    optimize!(model)
    # Show solution
    if termination_status(model) == MOI.LOCALLY_SOLVED
        push!(efficient_frontier, (sqrt(objective_value(model)), target_return))
    end
    unregister(model, :c)
    delete(model, c)
```

```

end
# Plot Efficient Frontier
fig = Figure()
Axis(fig[1, 1])
lines!(map(x → x[1], efficient_frontier), map(x → x[2], efficient_frontier))
fig

```



29.4.3. Black-Litterman

The Black-Litterman model combines the views of investors with market equilibrium assumptions to generate optimal portfolios. It starts with a market equilibrium portfolio and adjusts it based on investor views and confidence levels. The model incorporates subjective opinions while maintaining diversification and risk management principles.

$$\begin{aligned}
 & \text{maximize} \quad \mu^T w - \lambda \cdot \frac{1}{2} w^T \Sigma w \\
 & \text{subject to} \quad \sum_{i=1}^N w_i = 1 \\
 & \quad \quad \quad w_i \geq 0, \quad \forall i
 \end{aligned}$$

29. Portfolio Optimization

```

using JuMP, Ipopt

λ = 2.5 # risk aversion
rfr = 0.02 # risk free rate
# Market equilibrium parameters (prior)
μ_market = [0.08, 0.08, 0.08] # Market equilibrium return
Σ_market = ρ # Market equilibrium covariance matrix
# Investor views
Q = μ # Expected returns on assets according to investor views
P = [1 0 0; 0 1 0; 0 0 1] # Pick matrix specifying which assets views are on
Ω = [0.001^2 0.0 0.0; 0.0 0.002^2 0.0; 0.0 0.0 0.003^2] # Views uncertainty (covariance matrix)

# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Black-Litterman expected return adjustment
Σ_prior_inv = inv(Σ_market)
τ = 0.05 # Scaling factor
# Calculate the posterior expected returns
μ_posterior = Σ_prior_inv * (τ * Σ_market * (Σ_prior_inv + P' * inv(Ω) * P)) \
    (τ * Σ_market * (Σ_prior_inv * μ_market + P' * inv(Ω) * Q) + Σ_prior_inv * μ_mark
# Objective: maximize sharpe ratio
sr = (w' * μ_posterior - rfr) / (λ / 2 * w' * Σ_market * w)
@objective(model, Max, sr)
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or p
@constraint(model, sum(w) == 1)
# Solve the optimization problem
optimize!(model)
# Print results
v = sqrt(value.(w)' * Σ_market * value.(w))
@show "Optimal Portfolio Weights, Expected Portfolio Return, Portfolio Volatility:", v
for i = 1:n_a
    @show ("Asset ", i, ": ", value.(w)[i], value.(w)[i] * μ_posterior[i])
end

```

```

("Optimal Portfolio Weights, Expected Portfolio Return, Portfolio Volatility:", v) = ("Optimal Po
("Asset ", i, ": ", value.(w)[i], value.(w)[i] * μ_posterior[i]) = ("Asset ", 1, ": ", 2.0824146208
("Asset ", i, ": ", value.(w)[i], value.(w)[i] * μ_posterior[i]) = ("Asset ", 2, ": ", 0.2311617288
("Asset ", i, ": ", value.(w)[i], value.(w)[i] * μ_posterior[i]) = ("Asset ", 3, ": ", 0.7688382502

```

29.4.4. Risk Parity

Risk parity is an asset allocation strategy that allocates capital based on risk rather than traditional measures such as market capitalization or asset prices. It aims to balance risk contributions across different assets or asset classes to achieve a more stable portfolio. Risk parity portfolios often include assets with different risk profiles, such as stocks, bonds, and commodities.

$$\begin{aligned} \text{minimize} \quad & \sum_{i=1}^N (w_i \cdot \sqrt{\sigma_i})^2 \\ \text{subject to} \quad & \sum_{i=1}^N w_i = 1 \\ & w_i \geq 0, \quad \forall i \end{aligned}$$

using JuMP, Ipopt

```
# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Objective: minimize portfolio variance
portfolio_variance = w'ρ * w
margin = (ρ * w ./ sqrt(portfolio_variance)) .* w
risk_contributions = margin ./ sum(margin)
target = repeat([1.0 / n_a], n_a)
@objective(model, Max, sum((risk_contributions .- target) .^ 2))
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or positive
@constraint(model, sum(w) == 1)
# Solve the optimization problem
optimize!(model)
# Print results
@show "Optimal Portfolio Weights:"
for i = 1:n_a
    @show ("Asset ", i, ": ", value.(w)[i])
end

"Optimal Portfolio Weights:" = "Optimal Portfolio Weights:"
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 1, ": ", -6.957484531612737e-9)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 2, ": ", 1.0000000131375544)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 3, ": ", -6.180069741122123e-9)
```

29. Portfolio Optimization

29.4.5. Sharpe Ratio Maximization

The Sharpe ratio measures the risk-adjusted return of a portfolio and is calculated as the ratio of excess return to volatility. Maximizing the Sharpe ratio involves finding the portfolio allocation that offers the highest risk-adjusted return. This approach focuses on achieving the best tradeoff between risk and return.

$$\begin{aligned} & \text{maximize} && \frac{E[R_p] - R_f}{\sigma_p} \\ & \text{subject to} && \sum_{i=1}^N w_i = 1 \\ & && w_i \geq 0, \quad \forall i \end{aligned}$$

```
using JuMP, Ipopt

# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Objective: minimize portfolio variance
rfr = 0.05 # risk free rate
@objective(model, Max, (dot(mu, w) - rfr) / sqrt(sum(w[i] * rho[i, j] * w[j] for i in 1:n_a, j in 1:n_a)))
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or positive
@constraint(model, sum(w) == 1)
# Solve the optimization problem
optimize!(model)
# Print results
@show "Optimal Portfolio Weights:"
for i = 1:n_a
    @show ("Asset ", i, ": ", value.(w)[i])
end

"Optimal Portfolio Weights:" = "Optimal Portfolio Weights:"
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 1, ": ", 0.010841995514843134)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 2, ": ", 0.5352292318109132)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 3, ": ", 0.45392877267424375)
```

29.4.6. Robust Optimization

Robust optimization techniques aim to create portfolios that are resilient to uncertainties and fluctuations in market conditions. These techniques consider a range of possible sce-

narios and optimize portfolios to perform well across different market environments. A robust parameter in robust portfolio optimization is typically chosen to ensure the portfolio's performance remains stable and satisfactory under different market conditions or variations in input data. Robust optimization may involve incorporating stress tests, scenario analysis, or robust risk measures into the portfolio construction process.

$$\begin{aligned} & \text{minimize} && w^T \Sigma w + \gamma \|w - w_0\|_2^2 \\ & \text{subject to} && \sum_{i=1}^N w_i = 1 \\ & && w_i \geq 0, \quad \forall i \\ & && \|(\Sigma^{1/2}(w - w_0))\|_2 \leq \epsilon \end{aligned}$$

using JuMP, Ipopt

```
# Create an optimization model
model = Model(optimizer_with_attributes(Ipopt.Optimizer, "print_level" => 0))
# Set up weights as variables to optimize
@variable(model, w[1:n_a] >= zero(0.0))
# Objective: minimize portfolio variance
ε = 0.05 # Uncertainty level
γ = 0.1 # Robustness parameter
w₀ = [0.3, 0.4, 0.3] # expected weights
@objective(model, Min, dot(w, ρ * w) + γ * sum((w[i] - w₀[i])^2 for i in 1:n_a))
# Constraints: Sum of portfolio weights should equal to 1, and all weights should be zero or p
@constraint(model, sum(w) == 1)
@constraint(model, sum((ρ[i, j] * (w[i] - w₀[i]) * (w[j] - w₀[j]))) for i in 1:n_a, j in 1:n_a) <
# Solve the optimization problem
optimize!(model)
# Print results
@show "Optimal Portfolio Weights:"
for i = 1:n_a
    @show ("Asset ", i, ": ", value.(w)[i])
end

"Optimal Portfolio Weights:" = "Optimal Portfolio Weights:"
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 1, ": ", 0.31250000098314346)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 2, ": ", 0.31250000376951037)
("Asset ", i, ": ", value.(w)[i]) = ("Asset ", 3, ": ", 0.37499999524734623)
```

29.4.7. Asset weights from different methodologies

29. Portfolio Optimization

Table 29.1.: Optimized asset weights from different methodologies

| Methodology | Asset weights |
|--|--|
| Standard mean variance
(with RBC costs) | [0.33, 0.17, 0.50]
[0.17, 0.83, 0.00] |
| Black-Litterman | [0.00, 0.23, 0.77] |
| Risk parity | [0.00, 1.00, 0.00] |
| Sharpe ratio | [0.01, 0.54, 0.45] |
| Robust | [0.31, 0.31, 0.38] |

Seeing from the asset weights from a standard mean variance approach, due to RBC costs asset weights shifted to ones with higher yields. Asset weights from Sharpe ratio approach aligns with the Sharpe ratio for each asset. Those from Robust approach seek results not far away from expected weights under different conditions.

29.5. Practical considerations

29.5.1. Fractional purchases of assets

In traditional portfolio optimization, fractional purchases of assets refer to the ability to allocate fractions or percentages of capital to individual assets. However, in certain contexts or practical implementations, fractional purchases may not be allowed or considered.

- Practical constraints. Some investment vehicles or platforms may restrict investors from purchasing fractions of shares or assets. For instance, certain mutual funds, exchange-traded funds (ETFs), or other investment products may require whole units of shares to be purchased.
- Simplicity and cost-effectiveness. Handling fractional shares can add complexity and operational costs to portfolio management, especially in terms of transaction fees, administrative overhead, and reconciliation processes.
- Market liquidity. Some assets may have limited liquidity or trading volumes, making it impractical or difficult to execute fractional purchases without significantly impacting market prices or transaction costs.
- Regulatory considerations. Regulations in certain jurisdictions may impose restrictions on fractional share trading or ownership, potentially limiting the ability to include fractional purchases in portfolio optimization strategies.

29.5.2. Large number of assets

In portfolio optimization, a penalty factor for a large volume of assets typically refers to a mechanism or adjustment applied to the optimization process to mitigate the potential biases or challenges that arise when dealing with a large number of assets. This concept is particularly relevant in the context of mean-variance optimization and other optimization frameworks where computational efficiency and practical portfolio management considerations come into play. Too many assets may have the following issues.

- Dimensionality. As the number of assets (or dimensions) increases in a portfolio, traditional optimization methods may become computationally intensive or prone to overfitting. This is because the complexity of the optimization problem grows exponentially with the number of assets.
- Sparsity and concentration. In practice, not all assets may contribute equally to portfolio performance. Some assets may have negligible impact on the overall portfolio characteristics (such as risk or return) due to low weights or correlations with other assets.
- Penalizing excessive complexity. A penalty factor can be introduced to penalize portfolios that overly diversify or allocate small weights to a large number of assets. This encourages the optimization process to focus on more significant assets or reduce the complexity of the portfolio structure.

There are various ways to implement a penalty factor for a large volume of assets:

- Regularization techniques. Techniques like Lasso (L1 regularization) or Ridge (L2 regularization) regression can penalize small weights or excessive diversification by adding a penalty term to the objective function.
- Subset selection. Methods that explicitly select a subset of assets based on their contribution to portfolio performance, rather than including all assets indiscriminately.
- Heuristic adjustments. Introducing heuristic rules or adjustments based on practical portfolio management principles or empirical observations.

30. Bayesian Mortality Modeling

"After a year of intense mental struggle, however, [Arthur Bailey] realized to his consternation that actuarial sledgehammering worked. He even preferred [the Bayesian underpinnings of credibility theory] to the elegance of frequentism. He positively liked formulae that described 'actual data. . . . I realized that the hard-shelled underwriters were recognizing certain facts of life neglected by the statistical theorists.' He wanted to give more weight to a large volume of data than to the frequentists' small sample; doing so felt surprisingly 'logical and reasonable.' He concluded that only a 'suicidal' actuary would use Fisher's method of maximum likelihood, which assigned a zero probability to nonevents." - Excerpt From The Theory That Would Not Die Sharon Bertsch McGrayne

30.1. In This Chapter

An example of using a Bayesian MCMC approach to fitting a mortality curve to sample data, with multi-level models and censored data.

30.2. Generating fake data

The problem of interest is to look at mortality rates, which are given in terms of exposures (whether or not a life experienced a death in a given year).

We'll grab some example rates from an insurance table, which has a "selection" component: When someone enters observation, say at age 50, their mortality is path dependent (so for someone who started being observed at 50 will have a different risk/mortality rate at age 55 than someone who started being observed at 45).

Additionally, there may be additional groups of interest, such as:

- high/medium/low risk classification
- sex
- group (e.g. company, data source, etc.)
- type of insurance product offered

30. Bayesian Mortality Modeling

The example data will start with only the risk classification above.

```
using MortalityTables
using Turing
using DataFramesMeta
using MCMCChains
using LinearAlgebra
using CairoMakie
using StatsBase

n = 10_000
inforce = map(1:n) do i
    (
        issue_age=rand(30:70),
        risk_level=rand(1:3),
    )
end

10000-element Vector{@NamedTuple{issue_age::Int64, risk_level::Int64}:]:
  (issue_age = 50, risk_level = 2)
  (issue_age = 69, risk_level = 3)
  (issue_age = 43, risk_level = 2)
  (issue_age = 67, risk_level = 2)
  (issue_age = 41, risk_level = 1)
  (issue_age = 45, risk_level = 2)
  (issue_age = 50, risk_level = 2)
  (issue_age = 68, risk_level = 1)
  (issue_age = 46, risk_level = 1)
  (issue_age = 30, risk_level = 3)
  :
  (issue_age = 42, risk_level = 3)
  (issue_age = 45, risk_level = 1)
  (issue_age = 70, risk_level = 2)
  (issue_age = 51, risk_level = 1)
  (issue_age = 65, risk_level = 1)
  (issue_age = 61, risk_level = 1)
  (issue_age = 48, risk_level = 3)
  (issue_age = 39, risk_level = 2)
  (issue_age = 53, risk_level = 2)

base_table = MortalityTables.table("2001 VBT Residual Standard Select and Ultimate - Male Nons
```

function tabular_mortality(params, issue_age, att_age, risk_level)

```

q = params.ultimate[att_age]
if risk_level == 1
    q *= 0.7
elseif risk_level == 2
    q = q
else
    q *= 1.5
end
end

tabular_mortality (generic function with 1 method)

function model_outcomes(inforce, assumption, assumption_params; n_years=5)

outcomes = map(inforce) do pol
    alive = 1
    sim = map(1:n_years) do t
        att_age = pol.issue_age + t - 1
        q = assumption(
            assumption_params,
            pol.issue_age,
            att_age,
            pol.risk_level
        )
        if rand() < q
            out = (att_age=att_age, exposures=alive, death=1)
            alive = 0
            out
        else
            (att_age=att_age, exposures=alive, death=0)
        end
    end
    filter!(x → x.exposures == 1, sim)
end

df = DataFrame(inforce)

df.outcomes = outcomes
df = flatten(df, :outcomes)

df.att_age = [x.att_age for x in df.outcomes]

```

30. Bayesian Mortality Modeling

```
df.death = [x.death for x in df.outcomes]
df.exposures = [x.exposures for x in df.outcomes]
select!(df, Not(:outcomes))

end

exposures = model_outcomes(inforce, tabular_mortality, base_table)
data = combine(groupby(exposures, [:issue_age, :att_age])) do subdf
    (exposures=nrow(subdf),
     deaths=sum(subdf.death),
     fraction=sum(subdf.death) / nrow(subdf))
end

data2 = combine(groupby(exposures, [:issue_age, :att_age, :risk_level])) do subdf
    (exposures=nrow(subdf),
     deaths=sum(subdf.death),
     fraction=sum(subdf.death) / nrow(subdf))
end
```

30.3. 1: A single binomial parameter model

| | issue_age | att_age | risk_level | exposures | deaths | fraction |
|-----|-----------|---------|------------|-----------|--------|-----------|
| | Int64 | Int64 | Int64 | Int64 | Int64 | Float64 |
| 1 | 30 | 30 | 1 | 74 | 0 | 0.0 |
| 2 | 30 | 30 | 2 | 94 | 0 | 0.0 |
| 3 | 30 | 30 | 3 | 66 | 1 | 0.0151515 |
| 4 | 30 | 31 | 1 | 74 | 0 | 0.0 |
| 5 | 30 | 31 | 2 | 94 | 0 | 0.0 |
| 6 | 30 | 31 | 3 | 65 | 0 | 0.0 |
| 7 | 30 | 32 | 1 | 74 | 0 | 0.0 |
| 8 | 30 | 32 | 2 | 94 | 0 | 0.0 |
| 9 | 30 | 32 | 3 | 65 | 0 | 0.0 |
| 10 | 30 | 33 | 1 | 74 | 0 | 0.0 |
| 11 | 30 | 33 | 2 | 94 | 0 | 0.0 |
| 12 | 30 | 33 | 3 | 65 | 0 | 0.0 |
| 13 | 30 | 34 | 1 | 74 | 0 | 0.0 |
| 14 | 30 | 34 | 2 | 94 | 1 | 0.0106383 |
| 15 | 30 | 34 | 3 | 65 | 0 | 0.0 |
| 16 | 31 | 31 | 1 | 65 | 0 | 0.0 |
| 17 | 31 | 31 | 2 | 86 | 0 | 0.0 |
| 18 | 31 | 31 | 3 | 78 | 0 | 0.0 |
| 19 | 31 | 32 | 1 | 65 | 0 | 0.0 |
| 20 | 31 | 32 | 2 | 86 | 0 | 0.0 |
| 21 | 31 | 32 | 3 | 78 | 0 | 0.0 |
| 22 | 31 | 33 | 1 | 65 | 0 | 0.0 |
| 23 | 31 | 33 | 2 | 86 | 0 | 0.0 |
| 24 | 31 | 33 | 3 | 78 | 0 | 0.0 |
| ... | ... | ... | ... | ... | ... | ... |

30.3. 1: A single binomial parameter model

Estiamte q , the average mortality rate, not accounting for any variation within the population/sample. Our model is defines as:

$$q \sim Beta(1, 1) \\ p(\text{death}) \sim Binomial(q)$$

```
@model function mortality(data, deaths)
    q ~ Beta(1, 1)
    for i = 1:nrow(data)
        deaths[i] ~ Binomial(data.exposures[i], q)
    end
end
```

30. Bayesian Mortality Modeling

```
m1 = mortality(data, data.deaths)
```

| Row | issue_age | att_age | exposures | deaths | fraction |
|-----|-----------|---------|-----------|--------|------------|
| | Int64 | Int64 | Int64 | Int64 | Float64 |
| 1 | 30 | 30 | 234 | 1 | 0.0042735 |
| 2 | 30 | 31 | 233 | 0 | 0.0 |
| 3 | 30 | 32 | 233 | 0 | 0.0 |
| 4 | 30 | 33 | 233 | 0 | 0.0 |
| 5 | 30 | 34 | 233 | 1 | 0.00429185 |
| 6 | 31 | 31 | 229 | 0 | 0.0 |
| 7 | 31 | 32 | 229 | 0 | 0.0 |
| 8 | 31 | 33 | 229 | 0 | 0.0 |
| : | : | : | : | : | : |
| 199 | 69 | 72 | 226 | 6 | 0.0265487 |
| 200 | 69 | 73 | 220 | 6 | 0.0272727 |
| 201 | 70 | 70 | 268 | 6 | 0.0223881 |
| 202 | 70 | 71 | 262 | 12 | 0.0458015 |
| 203 | 70 | 72 | 250 | 11 | 0.044 |
| 204 | 70 | 73 | 239 | 21 | 0.0878661 |
| 205 | 70 | 74 | 218 | 14 | 0.0642202 |

30.3.1. Sampling from the posterior

We use a No-U-Turn-Sampler (NUTS) technique to sample multiple chains at once:

```
num_chains = 4  
chain = sample(m1, NUTS(), MCMCThreads(), 400, num_chains)
```

Chains MCMC chain (400×13×4 Array{Float64, 3}):

```
Iterations      = 201:1:600
Number of chains = 4
Samples per chain = 400
Wall duration    = 2.9 seconds
```

30.3. 1: A single binomial parameter model

```

Compute duration = 2.34 seconds
parameters      = q
internals       = lp, n_steps, is_accept, acceptance_rate, log_density, hamiltonian_energy, hamil

Summary Statistics
parameters      mean      std      mcse    ess_bulk   ess_tail     rhat    e ...
Symbol          Float64  Float64  Float64  Float64   Float64  Float64   ...
q               0.0087  0.0004  0.0000  713.1171  854.4636  1.0011  ...
                                                               ...
                                                               1 column omitted

Quantiles
parameters      2.5%     25.0%    50.0%    75.0%    97.5%
Symbol          Float64  Float64  Float64  Float64  Float64
q               0.0079  0.0084  0.0087  0.0089  0.0095

```

Here, we have asked for the outcomes to be modeled via a single parameter for the population. We see that the posterior distribution of q is very close to the overall population mortality rate:

```
sum(data.deaths) / sum(data.exposures)
```

```
0.008676392896330312
```

However, We can see that the sampling of possible posterior parameters doesn't really fit the data very well since our model was so simplified. The lines represent the posterior binomial probability.

This is saying that for the observed data, if there really is just a single probability p that governs the true process that came up with the data, there's a pretty narrow range of values it could possibly be:

```

let
  data_weight = sqrt.(data.exposures) / 2
  f = Figure(title="Parametric Bayesian Mortality"
  )
  ax = Axis(f[1, 1],
    xlabel="age",
    ylabel="mortality rate",
    limits=(nothing, nothing, -0.01, 0.10),
  )
  scatter!(ax,

```

30. Bayesian Mortality Modeling

```
data.att_age,
data.fraction,
markersize=data_weight,
color=(:blue, 0.5),
label="Experience data point (size indicates relative exposure quantity),)

# show n samples from the posterior plotted on the graph
n = 300
ages = sort!(unique(data.att_age))

q_posterior = sample(chain, n)[:q]

for i in 1:n

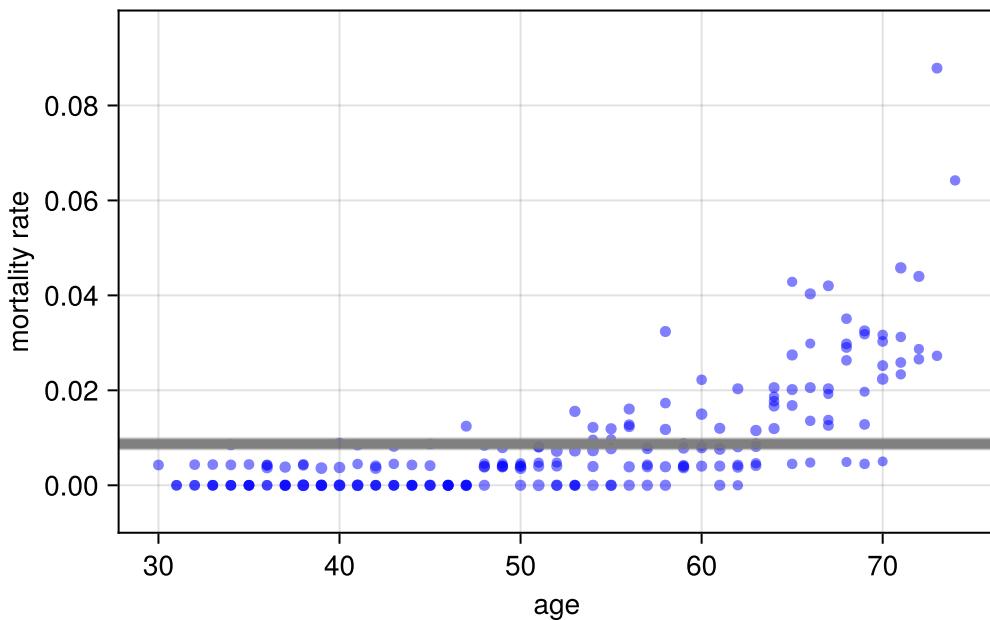
    hlines!(ax, [q_posterior[i]], color=(:grey, 0.1))
end

# Need to simulate at individual level and then aggregate?

sim05 = Float64[]
sim95 = Float64[]
for r in eachrow(data)
    outcomes = map(1:n) do i
        rand(Binomial(r.exposures, q_posterior[i]), 500)
    end
    push!(sim05, quantile(Iterators.flatten(outcomes), 0.05) / r.exposures)
    push!(sim95, quantile(Iterators.flatten(outcomes), 0.95) / r.exposures)
end

f
end
```

30.3. 1: A single binomial parameter model



```
let
    n = 300
    q_posterior = sample(chain, n)[:q]

end

2-dimensional AxisArray{Float64,2,...} with axes:
    :iter, 1:300
    :chain, 1:1
And data, a 300×1 Matrix{Float64}:
0.008874147042778021
0.009143536361329889
0.008574010582901878
0.00877679333926885
0.008900258745216263
0.008477314489455998
0.008924412320845891
0.00869410413509644
0.008057701578419414
0.008174086987897523
⋮
0.008310737056777122
0.008946449734788926
```

30. Bayesian Mortality Modeling

```
0.00893346519817041  
0.008314889642732496  
0.008288823097788747  
0.008542127285064486  
0.008756511359966412  
0.008864318706374172  
0.008736117817141393
```

30.4. 2. Parametric model

In this example, we utilize a MakehamBeard parameterization because it's already very similar in form to a logistic function. This is important because our desired output is a probability (ie the probability of a death at a given age), so the value must be constrained to be in the interval between zero and one.

The **prior** values for a,b,c, and k are chosen to constrain the hazard (mortality) rate to be between zero and one.

This isn't an ideal parameterization (e.g. we aren't including information about the select underwriting period), but is an example of utilizing Bayesian techniques on life experience data. "

```
@model function mortality2(data, deaths)  
    a ~ Exponential(0.1)  
    b ~ Exponential(0.1)  
    c = 0.0  
    k ~ truncated(Exponential(1), 1, Inf)  
  
    # use the variables to create a parametric mortality model  
    m = MortalityTables.MakehamBeard(; a, b, c, k)  
  
    # loop through the rows of the dataframe to let Turing observe the data  
    # and how consistent the parameters are with the data  
    for i = 1:nrow(data)  
        age = data.att_age[i]  
        q = MortalityTables.hazard(m, age)  
        deaths[i] ~ Binomial(data.exposures[i], q)  
    end  
end  
  
mortality2 (generic function with 2 methods)
```

30.4. 2. Parametric model

We combine the model with the data and sample from the posterior using a similar call as before:

```
m2 = mortality2(data, data.deaths)

chain2 = sample(m2, NUTS(), MCMCThreads(), 400, num_chains)
```

Chains MCMC chain (400×15×4 Array{Float64, 3}):

```
Iterations          = 201:1:600
Number of chains   = 4
Samples per chain  = 400
Wall duration      = 78.35 seconds
Compute duration   = 77.94 seconds
parameters         = a, b, k
internals          = lp, n_steps, is_accept, acceptance_rate, log_density, hamiltonian_energy, hamil

Summary Statistics
parameters      mean     std    mcse  ess_bulk  ess_tail    rhat  e ...
Symbol    Float64  Float64  Float64  Float64  Float64  Float64  ...
a        0.1529  0.2648  0.1311  6.9459  15.4685  1.6026  ...
b        1.7521  2.8654  1.4183  6.9013  11.2360  1.5987  ...
k        1.8425  0.6345  0.0950  27.2549  644.2088  1.1030  ...
                                         1 column omitted

Quantiles
parameters      2.5%    25.0%    50.0%    75.0%    97.5%
Symbol    Float64  Float64  Float64  Float64  Float64
a        0.0000  0.0000  0.0000  0.1529  0.6113
b        0.0884  0.0956  0.1011  1.7649  6.7136
k        1.0307  1.3025  1.8237  2.2080  3.3722
```

30.4.1. Plotting samples from the posterior

We can see that the sampling of possible posterior parameters fits the data well:

```
let
    data_weight = sqrt.(data.exposures) / 2

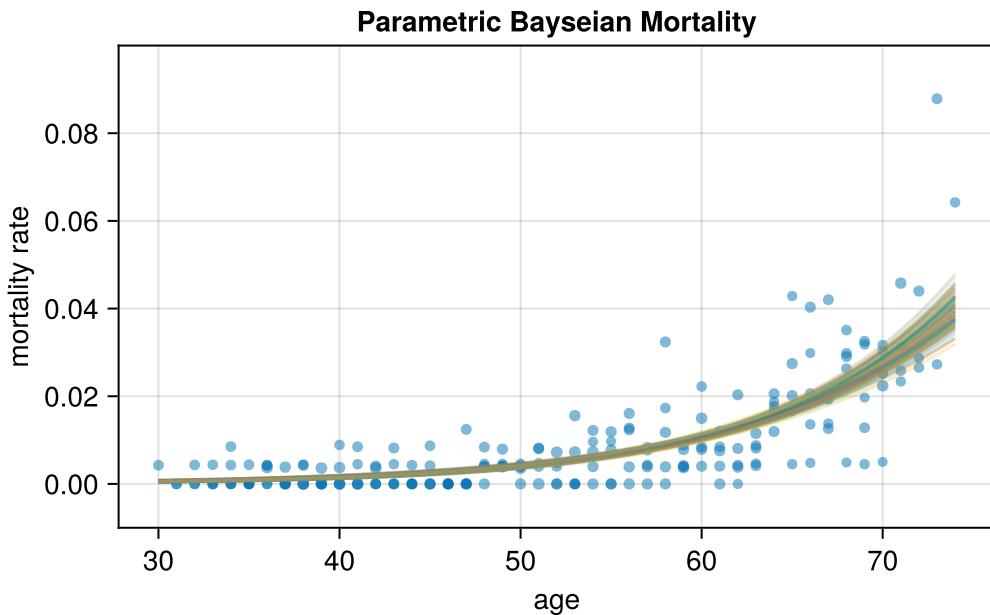
    p = scatter(
        data.att_age,
```

30. Bayesian Mortality Modeling

```
data.fraction,
markerSize=data_weight,
alpha=0.5,
label="Experience data point (size indicates relative exposure quantity)",
axis=(
    xlabel="age",
    limits=(nothing, nothing, -0.01, 0.10),
    ylabel="mortality rate",
    title="Parametric Bayesian Mortality"
)
)

# show n samples from the posterior plotted on the graph
n = 300
ages = sort!(unique(data.att_age))

for i in 1:n
    s = sample(chain2, 1)
    a = only(s[:a])
    b = only(s[:b])
    k = only(s[:k])
    c = 0
    m = MortalityTables.MakehamBeard(; a, b, c, k)
    lines!(ages, age → MortalityTables.hazard(m, age), alpha=0.1, label="")
end
p
end
```



```

let
    data_weight = sqrt.(data.exposures) / 2
    f = Figure(title="Parametric Bayesian Mortality")
)
ax = Axis(f[1, 1],
    xlabel="age",
    ylabel="mortality rate",
    limits=(nothing, nothing, -0.01, 0.10),
)
scatter!(ax,
    data.att_age,
    data.fraction,
    markersize=data_weight,
    color=(:blue, 0.5),
    label="Experience data point (size indicates relative exposure quantity),")

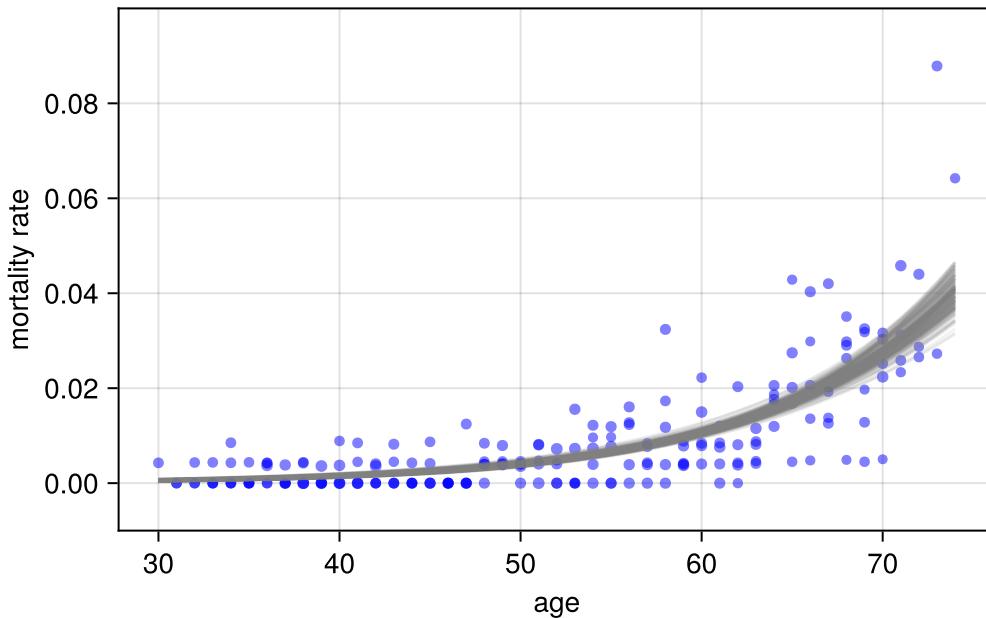
# show n samples from the posterior plotted on the graph
n = 300
ages = sort!(unique(data.att_age))

for i in 1:n
    s = sample(chain2, 1)
    a = only(s[:a])
    b = only(s[:b])

```

30. Bayesian Mortality Modeling

```
k = only(s[:k])
c = 0
m = MortalityTables.MakehamBeard(; a, b, c, k)
qs = MortalityTables.hazard.(m, ages)
lines!(ax, ages, qs, color=(:grey, 0.1))
end
f
end
```



Recall that the lines are not plotting the possible outcomes of the claims rates, but the *mean* claims rate for the given age.

30.5. 3. Multi-level model

This model extends the prior to create a multi-level model. Each risk class (`risk_level`) gets its own a parameter in the `MakhamBeard` model. The prior for a_i is determined by the hyper-parameter \bar{a} .

```
@model function mortality3(data, deaths)
    risk_levels = length(levels(data.risk_level))
    b ~ Exponential(0.1)
    ā ~ Exponential(0.1)
    a ~ filldist(Exponential(ā), risk_levels)
```

30.5. 3. Multi-level model

```

c = 0
k ~ truncated(Exponential(1), 1, Inf)

# use the variables to create a parametric mortality model

# loop through the rows of the dataframe to let Turing observe the data
# and how consistent the parameters are with the data
for i = 1:nrow(data)
    risk = data.risk_level[i]

    m = MortalityTables.MakehamBeard(; a=a[risk], b, c, k)
    age = data.att_age[i]
    q = MortalityTables.hazard(m, age)
    deaths[i] ~ Binomial(data.exposures[i], q)
end
end

m3 = mortality3(data2, data2.deaths)

chain3 = sample(m3, NUTS(), 1000)

summarize(chain3)

```

| parameters | mean | std | mcse | ess_bulk | ess_tail | rhat | e ... |
|------------|---------|---------|---------|----------|----------|---------|-------|
| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 | Float64 | ... |
| b | 0.0990 | 0.0053 | 0.0005 | 122.7749 | 260.7200 | 1.0125 | ... |
| ā | 0.0001 | 0.0001 | 0.0000 | 238.5899 | 272.1710 | 1.0245 | ... |
| a[1] | 0.0000 | 0.0000 | 0.0000 | 122.4680 | 298.1728 | 1.0144 | ... |
| a[2] | 0.0000 | 0.0000 | 0.0000 | 129.2373 | 290.2938 | 1.0125 | ... |
| a[3] | 0.0000 | 0.0000 | 0.0000 | 119.5696 | 286.8880 | 1.0154 | ... |
| k | 1.7292 | 0.7470 | 0.0269 | 505.5544 | 397.1778 | 1.0028 | ... |

1 column omitted

```

let data = data2

data_weight = sqrt.(data.exposures)
color_i = data.risk_level
cm = CairoMakie.Makie.wong_colors()

p, ax, _ = scatter(
    data.att_age,

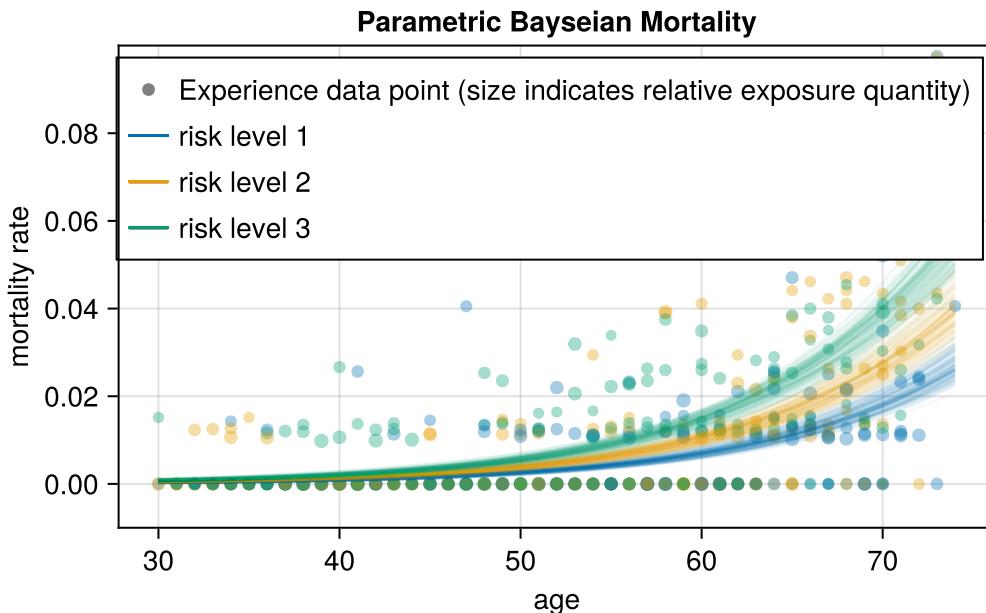
```

30. Bayesian Mortality Modeling

```
data.fraction,
markersize=data_weight,
alpha=0.5,
color=[(CairoMakie.Makie.wong_colors())[c], 0.7) for c in color_i],
colormap=CairoMakie.Makie.wong_colors(),
label="Experience data point (size indicates relative exposure quantity)",
axis=(
    xlabel="age",
    limits=(nothing, nothing, -0.01, 0.10),
    ylabel="mortality rate",
    title="Parametric Bayseian Mortality"
)
)

# show n samples from the posterior plotted on the graph
n = 100

ages = sort!(unique(data.att_age))
for r in 1:3
    for i in 1:n
        s = sample(chain3, 1)
        a = only(s[Symbol("a[$r]")])
        b = only(s[:b])
        k = only(s[:k])
        c = 0
        m = MortalityTables.MakehamBeard(; a, b, c, k)
        lines!(ages, age → MortalityTables.hazard(m, age), label="risk level $r", alpha=0
            end
        end
        axislegend(ax, merge=true)
    p
end
```



Again, the lines are not plotting the possible outcomes of the claims rates, but the *mean* claims rate for the given age and risk class.

30.6. Handling non-unit exposures

The key is to use the Poisson distribution, which is a continuous approximation to the Binomial distribution:

```
@model function mortality4(data, deaths)
    risk_levels = length(levels(data.risk_level))
    b ~ Exponential(0.1)
    ā ~ Exponential(0.1)
    a ~ filldist(Exponential(ā), risk_levels)
    c ~ Beta(4, 18)
    k ~ truncated(Exponential(1), 1, Inf)

    # use the variables to create a parametric mortality model

    # loop through the rows of the dataframe to let Turing observe the data
    # and how consistent the parameters are with the data
    for i = 1:nrow(data)
        risk = data.risk_level[i]
```

30. Bayesian Mortality Modeling

```

m = MortalityTables.MakehamBeard(; a=a[risk], b, c, k)
age = data.att_age[i]
q = MortalityTables.hazard(m, age)
deaths[i] ~ Poisson(data.exposures[i] * q)
end
end

m4 = mortality4(data2, data2.deaths)

chain4 = sample(m4, NUTS(), 1000)

Chains MCMC chain (1000×19×1 Array{Float64, 3}):

Iterations          = 501:1:1500
Number of chains   = 1
Samples per chain  = 1000
Wall duration      = 29.4 seconds
Compute duration   = 29.4 seconds
parameters         = b, ā, a[1], a[2], a[3], c, k
internals          = lp, n_steps, is_accept, acceptance_rate, log_density, hamiltonian_energy, hamil

Summary Statistics
parameters      mean     std    mcse    ess_bulk    ess_tail    rhat    e ...
Symbol          Float64  Float64  Float64  Float64  Float64  Float64  ...
b              0.1280  0.0113  0.0008  201.4085  273.8571  1.0004  ...
ā              0.0000  0.0000  0.0000  272.6510  336.1233  0.9997  ...
a[1]            0.0000  0.0000  0.0000  197.3968  263.0450  1.0004  ...
a[2]            0.0000  0.0000  0.0000  200.2965  293.7501  0.9997  ...
a[3]            0.0000  0.0000  0.0000  203.1390  281.7629  0.9996  ...
c              0.0014  0.0004  0.0000  285.2823  397.9634  1.0020  ...
k              1.9532  0.9831  0.0405  375.1733  224.7325  1.0042  ...
                                         1 column omitted

Quantiles
parameters      2.5%     25.0%    50.0%    75.0%    97.5%
Symbol          Float64  Float64  Float64  Float64  Float64
b              0.1073  0.1203  0.1269  0.1355  0.1506
ā              0.0000  0.0000  0.0000  0.0000  0.0001
a[1]            0.0000  0.0000  0.0000  0.0000  0.0000
a[2]            0.0000  0.0000  0.0000  0.0000  0.0000
a[3]            0.0000  0.0000  0.0000  0.0000  0.0000

```

30.6. Handling non-unit exposures

```

c      0.0006    0.0011    0.0014    0.0016    0.0021
k      1.0193    1.2613    1.6520    2.2800    4.7314

risk_factors4 = [mean(chain4[Symbol("a[$f]")]) for f in 1:3]

risk_factors4 ./ risk_factors4[2]

let data = data2

data_weight = sqrt.(data.exposures) / 2
color_i = data.risk_level

p, ax, _ = scatter(
    data.att_age,
    data.fraction,
    markersize=data_weight,
    alpha=0.5,
    color=color_i,
    label="Experience data point (size indicates relative exposure quantity)",
    axis=(xlabel="age",
          limits=(nothing, nothing, -0.01, 0.10),
          ylabel="mortality rate",
          title="Parametric Bayesian Mortality"
    )
)

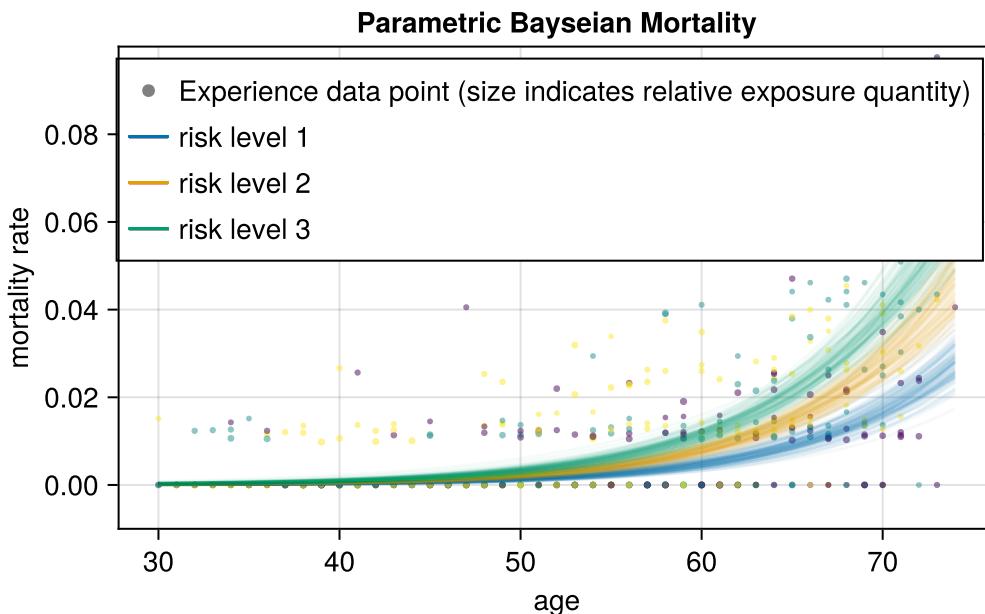
# show n samples from the posterior plotted on the graph
n = 100

ages = sort!(unique(data.att_age))
for r in 1:3
    for i in 1:n
        s = sample(chain4, 1)
        a = only(s[Symbol("a[$r]")])
        b = only(s[:b])
        k = only(s[:k])
        c = 0
        m = MortalityTables.MakehamBeard(; a, b, c, k)
        lines!(ages, age → MortalityTables.hazard(m, age), label="risk level $r", alpha=0.5)
    end
end
axislegend(ax, merge=true)
p

```

30. Bayesian Mortality Modeling

end



30.7. Model Predictions

We can generate predictive estimates by passing a vector of `missing` in place of the outcome variables and then calling `predict`.

We get a table of values where each row is the prediction implied by the corresponding chain sample, and the columns are the predicted value for each of the outcomes in our original dataset.

```
preds = predict(mortality4(data2, fill(missing, length(data2.deaths))), chain4)
```

Chains MCMC chain (1000×615×1 Array{Float64, 3}):

```
Iterations      = 1:1:1000
Number of chains = 1
Samples per chain = 1000
parameters     = deaths[1], deaths[2], deaths[3], deaths[4], deaths[5], deaths[6],
internals       =
```

```
Summary Statistics
parameters     mean       std      mcse    ess_bulk   ess_tail    rhat    ...

```

30.7. Model Predictions

| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 | Float64 | ... |
|------------|---------|---------|---------|-----------|-----------|---------|-----|
| deaths[1] | 0.1100 | 0.3377 | 0.0108 | 935.2083 | 902.0928 | 0.9992 | ... |
| deaths[2] | 0.1530 | 0.3972 | 0.0121 | 1060.9060 | 884.9780 | 0.9991 | ... |
| deaths[3] | 0.1130 | 0.3322 | 0.0107 | 951.8825 | 965.5019 | 0.9994 | ... |
| deaths[4] | 0.1100 | 0.3225 | 0.0110 | 860.9006 | 857.5899 | 1.0046 | ... |
| deaths[5] | 0.1610 | 0.3990 | 0.0123 | 1050.1385 | 1029.1752 | 0.9990 | ... |
| deaths[6] | 0.1280 | 0.3518 | 0.0111 | 1000.3078 | 892.5468 | 1.0019 | ... |
| deaths[7] | 0.1250 | 0.3654 | 0.0119 | 946.2305 | 960.6542 | 0.9996 | ... |
| deaths[8] | 0.1550 | 0.3939 | 0.0125 | 985.8479 | 793.6999 | 0.9990 | ... |
| deaths[9] | 0.1190 | 0.3478 | 0.0111 | 979.2443 | 980.3293 | 0.9990 | ... |
| deaths[10] | 0.1270 | 0.3592 | 0.0119 | 900.8917 | 897.1038 | 0.9993 | ... |
| deaths[11] | 0.1500 | 0.3894 | 0.0123 | 1000.5973 | 998.3315 | 1.0012 | ... |
| deaths[12] | 0.1080 | 0.3201 | 0.0104 | 942.7675 | 944.2663 | 0.9990 | ... |
| deaths[13] | 0.0940 | 0.3054 | 0.0097 | 978.1857 | 962.5454 | 0.9991 | ... |
| deaths[14] | 0.1770 | 0.4193 | 0.0137 | 942.1112 | 907.4458 | 0.9990 | ... |
| deaths[15] | 0.1260 | 0.3439 | 0.0106 | 1053.0661 | 1004.0241 | 1.0010 | ... |
| deaths[16] | 0.1100 | 0.3286 | 0.0116 | 795.0206 | 795.0903 | 0.9999 | ... |
| deaths[17] | 0.1330 | 0.3513 | 0.0113 | 964.0478 | 947.5567 | 0.9992 | ... |
| : | : | : | : | : | : | : | .. |

1 column and 598 rows omitted

Quantiles

| parameters | 2.5% | 25.0% | 50.0% | 75.0% | 97.5% |
|------------|---------|---------|---------|---------|---------|
| Symbol | Float64 | Float64 | Float64 | Float64 | Float64 |
| deaths[1] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[2] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[3] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[4] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[5] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[6] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[7] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[8] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[9] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[10] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[11] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[12] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[13] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[14] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[15] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[16] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |
| deaths[17] | 0.0000 | 0.0000 | 0.0000 | 0.0000 | 1.0000 |

30. Bayesian Mortality Modeling

```
⋮ ⋮ ⋮ ⋮ ⋮ ⋮  
598 rows omitted
```

31. Other Useful Techniques

31.1. In this chapter

Other useful techniques are surveyed, such as: memoization to avoid repeated computations, psuedo-monte carlo, creating a model office, and tips on modeling a complete balance sheet. Also covered are elements of practical review such as static and dynamic validations, and implied rate analysis.

31.2. Conceptual Techniques

31.2.1. Taking things to the Extreme

Consider what happens if something is taken to an extreme. For example, what happens in the model if we input negative rates? Where should negative rates be allowed and can the model handle them?

31.2.2. Range Bounding

Sometimes you just need to know that an outcome is within a certain range - if you can develop a "high" and "low" estimate by making assumptions that you know are outside of feasible ranges, then you can determine whether something is reasonable or within tolerances.

To take an example from the pages of interview questions: say you need to determine if a mortgaged property's value is greater than the amount of the outstanding loan (say \$100,000). You don't have an appraisal, but know that it's in reasonable condition and that (1) a comparable house with many more issues sold for \$100 per square foot. You also don't know the square footage of the house, but know from the number of rooms and layout that it must be at least 1000 square feet. Therefore you know that the value should at least be greater than:

$$\frac{\$100}{\text{sq. ft}} \times 1000 \text{sq. ft} = \$100,000$$

31. Other Useful Techniques

We'd then conclude that the value of the house very likely exceeds the outstanding balance of the loan and resolves our query without complex modeling or expensive appraisals.

31.3. Modeling Techniques

31.3.1. Serialization

31.4. Model Validation

31.4.1. Static and dynamic validation

Static validation typically involves splitting the dataset into training and testing sets, where the testing set is held out and not used during model training. The model is trained on the training set and then evaluated on the held-out testing set to assess its performance. This approach helps to measure how well the model generalizes to unseen data.

Dynamic validation, on the other hand, involves using a rolling or expanding window to train and test the model iteratively over time. In each iteration, the model is trained on past data and tested on future data, simulating how the model would perform in a real-world scenario where new data becomes available over time. This approach helps to assess the model's ability to adapt to changing patterns and trends in the data.

The following example shows how to do a static validation in Julia.

```
using Statistics

# Generate synthetic time series data
num_samples = 100
data = rand(num_samples)
X = [ones(num_samples) data]
y = 2data .+ 1 .+ 0.1 * randn(num_samples, 1) # dependent variable with noise
# Train the model on the training set
θ = X \ y
# Predictions
y_pred = θ[2] .* data .+ θ[1]
# Compute evaluation metrics
mse = mean((y_pred .- y) .^ 2)
mae = mean(abs.(y_pred .- y))

println("Static validation results:")
```

31.4. Model Validation

```
println("Mean Squared Error (MSE): ", mse)
println("Mean Absolute Error (MAE): ", mae)

Static validation results:
Mean Squared Error (MSE): 0.007494982616634391
Mean Absolute Error (MAE): 0.06583730697076114
```

The following example shows how to do a dynamic validation in Julia.

```
using Statistics

# Dynamic validation to update model over time and evaluate
num_updates = 5
mse_dyn = Float64[]
mae_dyn = Float64[]
for i in 1:num_updates
    data = rand(num_samples)
    X = [ones(num_samples) data]
    y = 2data .+ 1 .+ 0.1 * randn(num_samples, 1) # dependent variable with noise
    # Train the model on the training set
    θ = X \ y
    # Predictions
    y_pred = θ[2] .* data .+ θ[1]
    # Compute evaluation metrics
    mse = mean((y_pred .- y) .^ 2)
    mae = mean(abs.(y_pred .- y))
    push!(mse_dyn, mse)
    push!(mae_dyn, mae)
end

println("Dynamic validation results:")
println("Mean Squared Error (MSE): ", mean(mse_dyn))
println("Mean Absolute Error (MAE): ", mean(mae_dyn))

Dynamic validation results:
Mean Squared Error (MSE): 0.010216023499729613
Mean Absolute Error (MAE): 0.08193562774367437
```

31.4.2. Implied rate analysis

Implied rates are rates that are derived from the prices of financial instruments, such as bonds or options. For example, in the context of bonds, the implied rate is the interest

31. Other Useful Techniques

rate that equates the present value of future cash flows from the bond (coupons and principal) to its current market price.

```
using Zygote

# Define the bond cash flows and prices
cash_flows = [100, 100, 100, 100, 1000] # Coupons and principal
prices = [950, 960, 1010, 1020, 1050] # Market prices

# Define a function to calculate the present value of cash flows given a rate
function present_value(rate, cash_flows)
    pv = 0
    for (i, cf) in enumerate(cash_flows)
        pv += cf / (1 + rate)^i
    end
    return pv
end

# Define a function to calculate the implied rate using bisection method
function implied_rate(cash_flows, price)
    f(rate) = present_value(rate, cash_flows) - price
    return rootassign(f, 0.0, 1.0)
end

function rootassign(f, l, u)
    # Define an initial value
    x = 0.05
    # tolerance of difference in value
    tol = 1e-6
    # maximum number of iteration of the algorithm
    max_iter = 100
    iter = 0
    while abs(f(x)) > tol && iter < max_iter
        x -= f(x) / gradient(f, x)[1]
        iter += 1
    end
    if iter < max_iter && l < x < u
        return x
    else
        return -1.0
    end
end

# Calculate implied rates for each bond
implied_rates = [implied_rate(cash_flows, price) for price in prices]
```

31.4. Model Validation

```
# Print the results
for (i, rate) in enumerate(implied_rates)
    println("Implied rate for bond $i: $rate")
end
```

```
Implied rate for bond 1: 0.09658339166435045
Implied rate for bond 2: 0.09380219311021369
Implied rate for bond 3: 0.08046244727376842
Implied rate for bond 4: 0.0779014164014789
Implied rate for bond 5: 0.07041724037694008
```


Part VIII.

Appendices

32. The Julia Ecosystem Today

A tour of relevant available packages as of 2023.

The Julia ecosystem favors composability and interoperability, enabled by multiple dispatch. In other words, because it's easy to automatically specialize functionality based on the type of data being used, there's much less need to bundle a lot of features within a single package.

As you'll see, Julia packages tend to be less vertically integrated because it's easier to pass data around. Counterexamples of this in Python and R:

- Numpy-compatible packages that are designed to work with a subset of numerically fast libraries in Python
- special functions in Pandas to read CSV, JSON, database connections, etc.
- The Tidyverse in R has a tightly coupled set of packages that works well together but has limitations with some other R packages

Julia is not perfect in this regard, but it's neat to see how frequently things *just work*. It's not magic, but because of Julia features outside the scope of this article it's easy for package developers (and you!) to do this.

Julia also has language-level support for documentation, so packages can follow a consistent style of help-text and have the docs be auto-generated into web pages available locally or online.

The following highlighted packages were chosen for their relevance to typical actuarial work, with a bias towards those used regularly by the authors. This is a small sampling of the over 6000 registered Julia Packages¹

32.0.1. Data

Julia offers a rich data ecosystem with a multitude of available packages. Perhaps at the center of the data ecosystem are `CSV.jl` and `DataFrames.jl`. `CSV.jl` is for reading and writing files text files (namely CSVs) and offers top-class read and write performance. `DataFrames.jl` is a mature package for working with dataframes, comparable to Pandas or dplyr.

¹(`time?`) is a simple, built-in function. For true benchmarking purposes, see `?@sec-benchmarking`.

32. The Julia Ecosystem Today

Other notable packages include `ODBC.jl`, which lets you connect to any database (given you have the right drivers installed), and `Arrow.jl` which implements the Apache Arrow standard in Julia.

Worth mentioning also is `Dates`, a built-in package making date manipulation straightforward and robust.

Check out [JuliaData.org](https://juliadata.org) for more packages and information.

32.0.2. Plotting

`Plots.jl` is a meta-package providing an interface to consistently work with several plotting backends, depending if you are trying to emphasize interactivity on the web or print-quality output. You can very easily add animations or change almost any feature of a plot.

`StatsPlots.jl` extends `Plots.jl` with a focus on data visualization and compatibility with dataframes.

`Makie.jl` supports GPU-accelerated plotting and can create very rich, beautiful visualizations, but its main downside is that it has not yet been optimized to minimize the time-to-first-plot.

32.0.3. Statistics

Julia has first-class support for missing values, which follows the rules of three-valued logic so other packages don't need to do anything special to incorporate missing values.

`StatsBase.jl` and `Distributions.jl` are essentials for a range of statistics functions and probability distributions respectively.

Others include:

- `Turing.jl`, a probabilistic programming (Bayesian statistics) library, which is outstanding in its combination of clear model syntax with performance.
- `GLM.jl` for any type of linear modeling (mimicking R's `glm` functionality).
- `LsqFit.jl` for fitting data to non-linear models.
- `MultivariateStats.jl` for multivariate statistics, such as PCA.

You can find more packages and learn about them [here](#).

32.0.4. Machine Learning

Flux, Gen, Knet, and MLJ are all very popular machine learning libraries. There are also packages for PyTorch, Tensorflow, and SciKitML available. One advantage for users is that the Julia packages are written in Julia, so it can be easier to adapt or see what's going on in the entire stack. In contrast to this design, PyTorch and Tensorflow are built primarily with C++.

Another advantage is that the Julia libraries can use automatic differentiation to optimize on a wider range of data and functions than those built into libraries in other languages.

32.0.5. Differentiable Programming

Sensitivity testing is very common in actuarial workflows: essentially, it's understanding the change in one variable in relation to another. In other words, the derivative!

Julia has unique capabilities where almost across the entire language and ecosystem, you can take the derivative of entire functions or scripts. For example, the following is real Julia code to automatically calculate the sensitivity of the ending account value with respect to the inputs:

```
julia> using Zygote

julia> function policy_av(pol)
    COIs = [0.00319, 0.00345, 0.0038, 0.00419, 0.0047, 0.00532]
    av = 0.0
    for (i,coi) in enumerate(COIs)
        av += av * pol.credit_rate
        av += pol.annual_premium
        av -= pol.face * coi
    end
    return av           # return the final account value
end

julia> pol = (annual_premium = 1000, face = 100_000, credit_rate = 0.05);

julia> policy_av(pol)      # the ending account value
4048.08

julia> policy_av'(pol)     # the derivative of the account value with respect to the inputs
(annual_premium = 6.802, face = -0.0275, credit_rate = 10972.52)
```

32. The Julia Ecosystem Today

When executing the code above, Julia isn't just adding a small amount and calculating the finite difference. Differentiation is applied to entire programs through extensive use of basic derivatives and the chain rule. **Automatic differentiation**, has uses in optimization, machine learning, sensitivity testing, and risk analysis. You can read more about Julia's autodiff ecosystem [here](#).

32.0.6. Utilities

There are also a lot of quality-of-life packages, like `Revise.jl` which lets you edit code on the fly without needing to re-run entire scripts.

`BenchmarkTools.jl` makes it incredibly easy to benchmark your code - simply add `@benchmark` in front of what you want to test, and you will be presented with detailed statistics. For example:

```
julia> using ActuaryUtilities, BenchmarkTools

julia> @benchmark present_value(0.05,[10,10,10])

BenchmarkTools.Trial: 10000 samples with 994 evaluations.
Range (min ... max): 33.492 ns ... 829.015 ns      GC (min ... max): 0.00% ... 95.40%
Time   (median):    34.708 ns                      GC (median):    0.00%
Time   (mean ± σ):  36.599 ns ± 33.686 ns       GC (mean ± σ):  4.40% ±  4.55%
                                                 ┌─────────────────────────────────────────────────────────────────────────┐
                                                 └────────────────────────────────────────────────────────────────────────┘
  33.5 ns           Histogram: log(frequency) by time        45.6 ns <
```

Memory estimate: 112 bytes, allocs estimate: 1.

`Test` is a built-in package for performing testsets, while `Documenter.jl` will build high-quality documentation based on your inline documentation.

`ClipData.jl` lets you copy and paste from spreadsheets to Julia sessions.

32.0.7. Other packages

Julia is a general-purpose language, so you will find packages for web development, graphics, game development, audio production, and much more. You can explore packages (and their dependencies) at <https://juliahub.com/>.

32.0.8. Actuarial packages

Saving the best for last, the next article in the series will dive deeper into actuarial packages, such as those published by JuliaActuary for easy mortality table manipulation, common actuarial functions, financial math, and experience analysis.

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