Software Engineering

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Contents

1	Software models	4
	1.1 The "Amazon Prime Video" article	4
	1.2 Formal theory	5
	1.2.1 Markov chain	5
	1.2.2 Markov decision process	5
	1.2.3 Example	6
	1.3 Tips and tricks	7
	1.3.1 Average	7
	1.3.2 Welford's online algorithm (standard deviation)	7
	1.3.3 Euler's method for differential equations	
2	· C++	9
_	2.1 The random library	-
	2.1.1 Random engines	
	2.1.2 Operator overloading (quick note)	
	2.1.3 Distributions	
	2.2 Dynamic structures	
	2.2.1 Manual memory allocation (and how to avoid it)	
	2.2.2 std::vector <t>()</t>	
	2.2.3 std::deque <t>()</t>	
	2.2.4 Sets	
	2.2.5 Maps	
	2.3 I/O	
	2.3.1 Standard I/O	
	2.3.2 Files	
		14
3	Debugging with gdb	15
4	Examples	16
	4.1 First examples	16
	4.1.1 A simple Markov decision process [1100]	16
	4.1.2 Markov decision processes network pt.1 [1200]	16
	4.1.3 Markov decision processes network pt.2 [1300]	17
	4.1.4 Markov decision processes network pt.3 [1400]	18
	4.2 Traffic light [2000]	18
	4.3 Control center	18
	4.3.1 No network [3100]	18
	4.3.2 Network monitor	18
	4.3.3 Faults & repair [3400]	18
	4.3.4 Faults & repair + correct protocol [3500]	18
	4.4 Statistics	
	4.4.1 Expected value [4100]	19
	4 4 2 Probability [4200]	19

4.5	Development process simulation	20
	4.5.1 Random transition matrix [5100]	20
	4.5.2 [5200] Software development & error detection	21
	4.5.3 Optimizing costs for the development team [5300]	24
	4.5.4 Key performance index [5400]	24
4.6	Complex systems	24
	4.6.1 Insulin pump [6100]	24
	4.6.2 Buffer [6200]	24
	4.6.3 Server [6300]	24
5 Ex	vom	25
	Development team (time & cost)	
	Backend load balancing	
5.2	5.2.1 Env	
	5.2.2 Dispatcher, Server and Database	
	5.2.3 Response time	
5.9	3 Heater simulation	
	Task management	
5.4	task management	20
	OCC library	26
6.1	Observer Pattern	26
6.1	· ·	26
6.1 6.2	Observer Pattern	26 26
6.1 6.2 7 E 2	Observer Pattern	26 26 27
6.1 6.2 7 E 2	Observer Pattern	26 26 27 27
6.1 6.2 7 E 2	Observer Pattern	26 26 27 27 27
6.1 6.2 7 E 2 7.1	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs	26 26 27 27 27 27
6.1 6.2 7 E 2 7.1	Observer Pattern C++ generics & virtual methods Ctras VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs C Advanced testing techinques (in Rust & C)	26 26 27 27 27 27 27
6.1 6.2 7 E 2 7.1	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall)	26 26 27 27 27 27 27 27
6.1 6.2 7 E 2 7.1	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs 2 Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall) 7.2.2 Fuzzying (cargo-fuzz)	26 26 27 27 27 27 27 27 27 27
6.1 6.2 7 E 2 7.1	Observer Pattern C++ generics & virtual methods Atras VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall) 7.2.2 Fuzzying (cargo-fuzz) 7.2.3 Property-based testing	26 26 27 27 27 27 27 27 27 27
6.1 6.2 7 E 2 7.1	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs 2 Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall) 7.2.2 Fuzzying (cargo-fuzz)	26 26 27 27 27 27 27 27 27 27 27 27
6.1 6.2 7 E 2 7.1 7.2	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall) 7.2.2 Fuzzying (cargo-fuzz) 7.2.3 Property-based testing 7.2.4 Test augmentation (Miri, Loom, Valgrind) 7.2.5 Performance testing	26 26 27 27 27 27 27 27 27 27 27 27 27
6.1 6.2 7 E 2 7.1 7.2	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall) 7.2.2 Fuzzying (cargo-fuzz) 7.2.3 Property-based testing 7.2.4 Test augmentation (Miri, Loom, Valgrind) 7.2.5 Performance testing 8 UI testing?	26 26 27 27 27 27 27 27 27 27 27 27 27
6.1 6.2 7 E 2 7.1 7.2	Observer Pattern C++ generics & virtual methods VDM (Vienna Development Method) 7.1.1 It's cool, I promise 7.1.2 VDM++ to design valid UMLs Advanced testing techinques (in Rust & C) 7.2.1 Mocking (mockall) 7.2.2 Fuzzying (cargo-fuzz) 7.2.3 Property-based testing 7.2.4 Test augmentation (Miri, Loom, Valgrind) 7.2.5 Performance testing	26 26 27 27 27 27 27 27 27 27 27 27 27

1 Software models

Software projects require **design choices** that often can't be driven by experience or reasoning alone. That's why a **model** of the project is needed to compare different solutions. In this course **Markov decision processes** are used to describe software systems.

1.1 The "Amazon Prime Video" article

If you were tasked with designing the software architecture for **Amazon Prime Video** (a live streaming service for Amazon), how would you go about it? What if you had the to keep the costs minimal? Would you use a distributed architecture or a monolith application?

More often than not, monolith applications are considered **more costly** and **less scalable** than the counterpart due to an inefficient usage of resources. But, in a recent article, a Senior SDE at Prime Video describes how they "reduced the cost of the audio/video monitoring infrastructure by 90%" [1] by using a monolith architecture.

There isn't a definitive way to answer these type of questions, but one way to go about it is building a model of the system to compare the solutions. In the case of Prime Video, "the audio/video monitoring service consists of three major components:" [1]

- the **media converter** converts input audio/video streams
- the **defect detectors** analyze frames and audio buffers in real-time
- the **orchestrator** controls the flow in the service

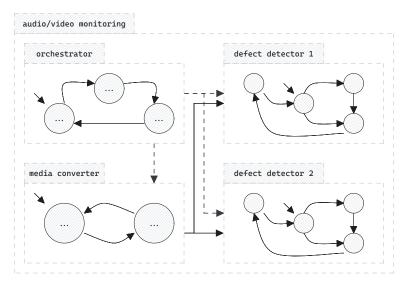


Figure 1: audio/video monitoring system

To derive conclusions the system can be **simulated** by modeling its components as **Markov decision processes**.

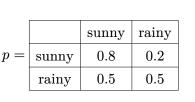
1.2 Formal theory

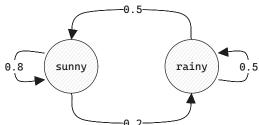
1.2.1 Markov chain

A Markov chain is defined by a set of states S and the transition **probability** $p: S \times S \rightarrow [0,1]$ such that p(s'|s) is the probability to transition to state s' if the current state is s, with the constraint that

$$\forall s \in S \ \sum_{s' \in S} p(s'|s) = 1 \tag{1}$$

For example, the weather can be modeled with $S = \{\text{sunny}, \text{rainy}\}\$ and p such that





If a Markov chain transitions at discrete time steps, i.e. the time steps $t_0, t_1, t_2, ...$ are a **countable**, then it's called a DTMC (discrete-time Markov chain), otherwise it's called a CTMC (continuous-time Markov chain).

This kind model is **limited**. To describe complex systems (e.g. servers, rockets, medical devices) the concepts of **input** and **output** are needed.

1.2.2 Markov decision process

A Markov decision process (MDP) is **different** from a Markov chain. A MDP M is a tuple (U, X, Y, p, q) s.t.

- U is the set of input values
- X is the set of states
- Y is the set of **output values**
- $p: X \times X \times U \to [0,1]$ is such that p(x'|x,u) is the probability to **transition** from state x to state x' when the **input value** is u
- $g: X \to Y$ is the **output function**
- and let $x_0 \in X$ be the initial state

The same constrain in Equation 1 holds for the MDP, with an important difference: **for each input value**, the sum of the transition probabilities for **that input value** must be 1.

$$\forall x \in X \ \forall u \in U \ \sum_{x' \in X} p(x'|x, u) = 1$$
 (2)

To be more precise, let \boldsymbol{x}_t be the state at the time step t, then

$$p(x'|x,u) = p(x_{t+1} = x' \mid x_t = x, u_t = u)$$
(3)

if the MDP is discrete-time.

1.2.3 Example

The development process of a company can be modeled as a MDP M=(U,X,Y,p,g) s.t.

- $U = \{\varepsilon\}$ because U can't be empty
- $X = \{0, 1, 2, 3, 4\}$
- $Y = \text{Cost} \times \text{Duration}$
- $x_0 = 0$

$$g(x) = \begin{cases} (0,0) & x = 0 \lor x = 4\\ (20000,2) & x = 1 \lor x = 3\\ (40000,4) & x = 2 \end{cases} \tag{4}$$

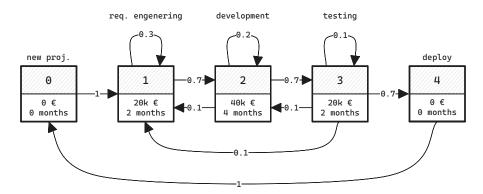


Figure 2: the model of a team's development process

	ε	0	1	2	3	4
	0	0	1	0	0	0
n —	1	0	.3	.7	0	0
p =	2	0	.1	.2	.7	0
	3	0	.1	.1	.1	.7
	4	1	0	0	0	0

Note that only **1 transition matrix** is defined, because |U| = 1 (there's 1 input value). If $U = \{\text{apple, banana, orange}\}$ 3 transition matrices must be defined, one **for each input value**.

1.3 Tips and tricks

1.3.1 Average

Given a set of values $X=\{x_1,...,x_n\}\subset\mathbb{R}$ the average $\overline{x}_n=\frac{\sum_{i=0}^n x_i}{n}$ can be computed with a simple procedure

```
float average(std::vector<float> X) {
   float sum = 0;
   for (auto x_i : X)
      sum += x_i;

return sum / X.size();
}
```

The problem with this procedure is that, by adding up all the values before the division, the **numerator** could **overflow**, even if the value of \overline{x}_n fits within the IEEE-754 limits. Nonetheless, \overline{x}_n can be calculated incrementally.

$$\overline{x}_{n+1} = \frac{\sum_{i=0}^{n+1} x_i}{n+1} = \frac{\left(\sum_{i=0}^{n} x_i\right) + x_{n+1}}{n+1} = \frac{\sum_{i=0}^{n} x_i}{n+1} + \frac{x_{n+1}}{n+1} = \frac{\left(\sum_{i=0}^{n} x_i\right)n}{(n+1)n} + \frac{x_{n+1}}{n+1} = \frac{\sum_{i=0}^{n} x_i}{n} \cdot \frac{n}{n+1} + \frac{x_{n+1}}{n+1} = \frac{\overline{x}_{n+1}}{n} = \frac{\overline{x}_{n+1}}{n+1}$$

$$(5)$$

With this formula the numbers added up are smaller: \overline{x}_n is multiplied by $\frac{n}{n+1} \sim 1$, and, if x_{n+1} fits in IEEE-754, then $\frac{x_{n+1}}{n+1}$ can also be encoded.

```
float average_r(std::vector<float> X) {
   float average = 0;
   for (size_t n = 0; n < X.size(); n++)
        average =
            average * ((float)n / (n + 1)) + X[n] / (n + 1);
   return average;
}</pre>
```

In examples/average.cpp the procedure average_r() is able to calculate the average and average() results in Inf.

1.3.2 Welford's online algorithm (standard deviation)

In a similar fashion, it could be faster and require less memory to calculate the **standard deviation** incrementally. Welford's online algorithm can be used for this purpose.

"It is often useful to be able to compute the variance in a single pass, inspecting each value x_i only once; for example, when the data is being

collected without enough storage to keep all the values, or when costs of memory access dominate those of computation." (Wikpedia)

$$\begin{split} M_{2,n} &= \sum_{i=1}^{n} (x_i - \overline{x}_n)^2 \\ M_{2,n} &= M_{2,n-1} + (x_n - \overline{x}_{n-1})(x_N - \overline{x}_n) \\ \sigma_n^2 &= \frac{M_{2,n}}{n} \\ s_n^2 &= \frac{M_{2,n}}{n-1} \end{split} \tag{6}$$

Given M_2 , the standard deviation can be calculated as $\sqrt{\frac{M_{2,n}}{n}}$ if n > 0.

```
real_t Stat::stddev_welford() const {
    return sqrt(n > 0 ? m_2__ / n : 0);
}
```

Listing 1: mocc/stat.hpp

1.3.3 Euler's method for differential equations

Got it from here [2]. Useful if a differential equation can't be solved analitically.

$$y_{n+1} = y_n + h \cdot f(t_n, y_n) \tag{7}$$

- TODO: approximation graphs, maybe with gnuplot
- gnuplot -e "set terminal png size 400,300; set output 'xyz.png'; plot [-4:4] exp(-x**2 / 2)"

2 C++

This section will cover the basics for the exam.

2.1 The random library

The C++ standard library offers powerful tools to easily implement MDPs.

2.1.1 Random engines

In C++ there are many ways to **generate random numbers** [3]. Generally it's **not recommended** to use **random()** ①. It's recommended to use a **random generator** ④, because it's fast, deterministic (given a **seed**, the sequence of generated numbers is the same) and can be used with **distributions**. To generate the seed a **random_device** is used: it's non deterministic because it uses a **hardware entropy source** (if available) to generate the random numbers.

```
#include <iostream>
#include <random>
int main() {
    std::cout << random() ① << std::endl;

    std::random_device random_device; ②
    std::cout << random_device() ③ << std::endl;

    std::default_random_engine r_engine(random_device() ④ );
    std::cout << r_engine() ⑤ << std::endl;
}</pre>
```

Listing 2: examples/random.cpp

The typical course of action is to instantiate a random_device ②, and use it to generate a seed ④ for a random_engine. Given that random engines can be used with distributions, they're really useful to implement MDPs.

From this point on, std::default_random_engine will be reffered to as urng_t (uniform random number generator type).

```
#include <random>
// works like typedef in C
using urng_t = std::default_random_engine;
int main() {
    std::random_device random_device;
    urng_t urng(random_device());
}
```

2.1.2 Operator overloading (quick note)

In Listing 2, to generate a random number, random_device() ③ and r_engine() ⑤ are used like functions with the () operator, but they aren't functions, they're instances of a class. That's because in C++ you can define how a certain operator (like +, +=, <<, >>, [], () etc..) should behave when used on a instance of the class. It's called **operator overloading** and other languages have it too:

- in Python operation overloading is done by implementing methods with special names, like __add__() [4]
- in Rust it's done by implementing the Trait associated with the operation, like std::ops::Add [5].
- Java and C don't have operator overloading

For example, std::cout is an instance of the std::basic_ostream class, which overloads the method "operator<<()" [6].

2.1.3 Distributions

Just the capability to generate random numbers isn't enough, these numbers need to be manipulated to fit certain needs. Luckly, C++ covers **basically all of them**. For example, the MDP in Figure 2 can be easily simulated with the following code code:

```
#include <iostream>
#include <random>
using urng t = std::default random engine;
int main() {
    std::random_device random_device;
    urng_t urng(random_device());
    std::discrete distribution<> transition matrix[] = {
        \{0, 1\},
        \{0, .3, .7\},\
        \{0, .2, .2, .6\},\
        \{0, .1, .2, .1, .6\},\
        {1},
    };
    size t state = 0;
    for (size_t step = 0; step < 15; step++) {</pre>
        state = transition_matrix[state](urng);
        std::cout << state << std::endl;</pre>
    }
    return 0;
}
```

Listing 3: examples/transition_matrix.cpp

2.1.3.1 Uniform discrete [7]

Let's consider a simple exercise

To test a system S it's required to build a generator that sends value v_t to S every T_t seconds. For each send, the value of T_t is an **integer** chosen uniformly in the range [20, 30].

The C code to compute T_t would be T = 20 + rand() % 11;, which is very **error prone**, hard to remember and has no semantic value.

In C++ the same can be done in a **simpler** and **cleaner** way:

```
std::uniform_int_distribution<> random_T(20, 30); (1)
size_t T = (2) random_T(urng);
```

The interval T_t can be easily generated ② without needing to remember any formula or trick. The behaviour of T_t is defined only once ①, so it can be easily changed without introducing bugs or inconsistencies. It's also worth to take a look at the implementation of the exercise above (with the addition that $v_t = T_t$), as it comes up very often in software models.

```
#include <iostream>
#include <random>
using urng_t = std::default_random_engine;
int main() {
    std::random_device random_device;
    urng_t urng(random_device());
    std::uniform_int_distribution<> random_T(20, 30);
    size_t T = random_T(urng), next_request_time = T;
    for (size_t time = 0; time < 1000; time++) {</pre>
        if (time < next_request_time)</pre>
            continue;
        std::cout << T << std::endl;</pre>
        T = random T(urng);
        next_request_time = time + T;
    }
    return 0;
}
```

Listing 4: examples/interval_generator.cpp

The uniform_int_distribution has many other uses, for example, it could uniformly generate a random state in a MDP. Let STATES_SIZE be the number of states

```
uniform int distribution >> random state(0, STATES SIZE - 1 (1));
```

random_state generates a random state when used. Be careful! Remember to use STATES_SIZE - 1 ①, because uniform_int_distribution is inclusive... forgettig it can lead to very sneaky bugs: random segfaults at different points of the code. It's very hard to debug unless using gdb. The uniform_int_distribution can also generate negative integers, for example $z \in \{x \mid x \in \mathbb{Z} \land x \in [-10, 15]\}$.

2.1.3.2 Uniform continuous [8]

It's the same as above, with the difference that it generates **real** numbers \mathbb{R} , in the range [a, b).

2.1.3.3 Bernoulli [9]

Let's consider the following exercise

```
To model a network protocol P it's required to model a request. The request can randomly fail with probability p=0.001.
```

Traditionally, a random real numer $r \in [0,1]$ is generated, and it's checked if r is above or below p.

```
std::uniform_real_distribution<> random_r(0, 1);
r = random_r(urng);
if (r > 0.001)
    fail();
```

The ${\tt std::bernoulli_distribution}$ is a better tool to implement this specification

```
std::bernoulli_distribution random_fail(0.001);
if (random_fail(urng))
    fail();
```

2.1.3.4 Normal

2.1.3.5 Exponential

The Exponential distribution is very useful when simulating user requests (generally, the interval between requests to a servers is described by a Exponential distribution, you just have to specify λ)

2.1.3.6 Poisson

2.1.3.7 Geometric

2.1.3.8 Discrete distribution

Let's consider the following exercise.

To choose the architecture for an e-commerce it's required to implement a model C of the customers that simulates the requests. After interviewing 678 people it's determined that 232 of them would buy a hat, 158 would buy a hoodie and the other 288 would buy a mug.

A discrete distribution can be used for this case. Let's say that "hat" = 0, "hoodie" = 1 and mug = 2.

```
#include <random>
#include <iostream>
int main() {
  std::random_device random_device;
  urng t urng(random device());
  std::discrete_distribution<> random_item = {232, 158, 288};
  for (int request = 0; request < 1000; request++) {</pre>
    size_t item = random_item(urng);
    switch (item) {
      case 0:
        std::cout << "hat" << std::endl;</pre>
        break:
      case 1:
        std::cout << "hoodie" << std::endl;</pre>
        break;
      case 2:
        std::cout << "mug" << std::endl;</pre>
        break:
  }
  return 0;
```

With the discrete distribution, the generated items are proportional to the data.

2.2 Dynamic structures

2.2.1 Manual memory allocation (and how to avoid it)

If you allocate with new, you must deallocate with delete, you can't mixup them with malloc() and free()

To avoid manual memory allocation, most of the time it's enough to use the structures in the standard library, like std::vector<T>.

2.2.2 std::vector<T>()

You don't have to allocate memory, basically never! You just use the structures that are implemented in the standard library, and most of the time they are enough for our use cases. They are really easy to use.

2.2.3 std::deque<T>()

2.2.4 Sets

Not needed as much

2.2.5 Maps

Could be useful

2.3 I/O

2.3.1 Standard I/O

2.3.2 Files

Working with files is way to easy in C++

```
#include <ofstream>
#include <ifstream>

int main(){
    ofstream output("output.txt");
    output << "some text" << std::endl;
    output.close();

    ifstream inputs("inputs.txt");
    int number;
    while (inputs >> number) {
        // do stuff with number...
    }
    inputs.close();

    return 0;
}
```

3 Debugging with gdb

It's super useful! Trust me, if you learn this everything is way easier

First of all, use the -ggdb flags to compile the code. Remember to not use any optimization like -03... using optimizations makes the program harder to debug.

```
DEBUG_FLAGS := -lm -std=c++11 -ggdb3 -Wall -Wextra -pedantic
```

Then it's as easy as running gdb ./main

- TODO: could be useful to write a script if too many args
- TODO: just bash code to compile and run
- TODO (just the most useful stuff, the chnically not enough):
 - r
 - c
 - n
 - ► c 10
 - enter (last instruction)
 - **▶** b
 - on lines
 - on symbols
 - on specific files
 - ightharpoonup clear
 - display
 - set print pretty on

4 Examples

Each example has 4 digits xxxx that are the same as the ones in the software folder in the course material.

4.1 First examples

This section puts together the **formal definitions** and the C++ knowledge to implement some simple MDPs.

4.1.1 A simple Markov decision process [1100]

```
The first MDP M = (U, X, Y, p, g) is such that
• U = \{\varepsilon\} (see Section 1.2.3)
• X = [0,1] \times [0,1], each state is a pair 3 of real numbers 1
• Y = [0,1] \times [0,1]
• p: X \times X \times U \to X = \mathcal{U}(0,1) \times \mathcal{U}(0,1), the transition probability is
   a uniform continuous distribution ②
• g: X \to Y: (r_0, r_1) \mapsto (r_0, r_1) outputs the current state \textcircled{4}
• x_0 = (0,0) is the initial state 3
  #include <random>
  using real_t 1 = double;
   const size_t HORIZON = 10;
   int main() {
       std::random_device random_device;
       urng_t urng(random_device());
       std::uniform_real_distribution<real_t> uniform(0, 1); (2)
       std::vector<real_t> state(2, 0); 3
       std::ofstream log("log");
       for (size_t time = 0; time <= HORIZON; time++) {</pre>
            for (auto &r : state)
                 r = uniform(urng); (2)
            log << time << ' ';
            for (auto r : state) log << r << ' '; t 4</pre>
            log << std::endl;</pre>
       }
       log.close();
       return 0;
  }
```

 ${
m Listing} \ 5: \ {
m software/ll00/main.cpp}$

4.1.2 Markov decision processes network pt.1 [1200]

This example has 2 MDPs M_0, M_1 like the one in the first example Section 4.1.1, with the difference that, and $U_i = [0, 1] \times [0, 1]$:

- $U_0(t+d) = Y_1(t)$
- $U_1(t+d) = Y_0(t)$

TODO: formula to get input from other stuff and calculate the state..., maybe define

$$p: X \times X \times U \to [0,1] \\ (x_0, x_1), (x_0', x_1'), (u_0, u_1) \mapsto \dots$$
 (8)

Lorem ipsum dolor sit amet, consectetur adipiscing elit, sed do eiusmod tempor incididunt ut labore et dolore magnam aliquam quaerat.

```
const size_t HORIZON = 100;
struct DTMC { real_t state[2]; };

int main() {
    std::vector<DTMC> mdps(2, {0, 0});

    for (size_t time = 0; time <= HORIZON; time++) {
        for (size_t r = 0; r < 2; r++) {
            mdps[0].state[r] = mdps[1].state[r] * uniform(urng);
            mdps[1].state[r] = mdps[0].state[r] + uniform(urng);
        }
    }
}</pre>
```

Listing 6: software/1200/main.cpp

4.1.3 Markov decision processes network pt.2 [1300]

The same as above, but with a different connection

 $\operatorname{Listing} 7: \mathsf{software/1300/main.cpp}$

4.1.4 Markov decision processes network pt.3 [1400]

The same as above, but with a twist (in the original uses variables to indicate each input... which is sketcy... I can do it with MOCC)

4.2 Traffic light [2000]

In this example we want to model a **traffic light**. The three versions of the system on the drive (2100, 2200 and 2300) do the same thing with a different code structure.

```
const size t HORIZON = 1000;
enum Light { GREEN = 0, YELLOW = 1, RED = 2 };
int main() {
    auto random_timer_duration =
        std::uniform_int_distribution<>(60, 120);
    Light traffic light = Light::RED;
    size_t timer = random_timer_duration(random_engine);
    for (size t time = 0; time <= HORIZON; time++) {</pre>
        if (timer > 0) {
            timer--;
            continue;
        }
        traffic_light =
            (traffic_light == RED
                 : (traffic_light == GREEN ? YELLOW : RED));
        timer = random_timer_duration(random_engine);
    }
}
```

Listing 8: software/2000/main.cpp

4.3 Control center

- 4.3.1 No network [3100]
- 4.3.2 Network monitor
- 4.3.2.1 No faults [3200]
- 4.3.2.2 Faults & no repair [3300]
- 4.3.3 Faults & repair [3400]
- 4.3.4 Faults & repair + correct protocol [3500]

4.4 Statistics

4.4.1 Expected value [4100]

In this one we just simulate a development process (phase 0, phase 1, and phase 2), and we calculate the average ...

4.4.2 Probability [4200]

In this one we simulate a more complex software developmen process, and we calculate the average cost (Wait, what? Do we simulate it multiple times?)

4.5 Development process simulation

An MDP can be implemented by using a **transition matrix** (like in Section 1.2.3). The simplest implementation can be done by using a std::discrete_distribution by using the trick in Listing 3.

4.5.1 Random transition matrix [5100]

This example builds a random transition matrix.

```
const size t HORIZON = 20, STATES SIZE = 10;
int main() {
    std::random_device random_device;
    urng_t urng(random_device());
    auto random_state = 1
        std::uniform int distribution<>(0, STATES SIZE - 1);
    std::uniform real distribution<> random real 0 1(0, 1);
    std::vector<std::discrete_distribution<>>
        transition matrix(STATES SIZE); (2)
    std::ofstream log("log.csv");
    for (size_t state = 0; state < STATES_SIZE; state++) {</pre>
        std::vector<real_t> weights(STATES_SIZE); 3
        for (auto &weight : weights)
            weight = random_real_0_1(urng);
        transition_matrix[state] = 4
            std::discrete_distribution<>(weights.begin(),
                                          weights.end());
    }
    size_t state = random_state(urng);
    for (size_t time = 0; time <= HORIZON; time++) {</pre>
        log << time << " " << state << std::endl;</pre>
        state = transition_matrix[state 5] (urng); 6
    }
    log.close();
    return 0;
}
```

Listing 9: software/5100/main.cpp

A transition matrix is a vector<discrete_distribution<>> ② just like in Listing 3. Why can we do this? First of all, the states are numbered from 0 to STATES_SIZE - 1, that's why we can generate a random state ① just by generating a number from 0 to STATES_SIZE - 1.

The problem with using a simple uniform_int_distribution is that we don't want to choose the next state uniformly, we want to do something like in Figure 3.

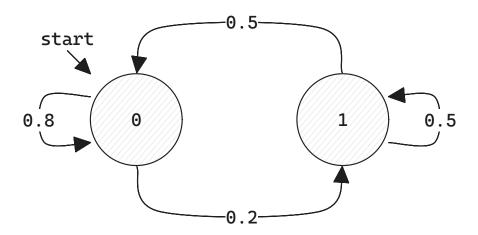


Figure 3: A simple Markov Chain

Luckly for us std::discrete_distribution<> does exactly what we want. It takes a list of weights $w_0, w_1, w_2, ..., w_n$ and assigns each index i the probability $p(i) = \frac{\sum_{i=0}^n w_i}{w_i}$ (the probability is proportional to the weight, so we have that $\sum_{i=0}^n p(i) = 1$ like we would expect in a Markov Chain).

To instantiate the discrete_distribution 4, unlike in Listing 3, we need to first calculate the weights 3, as we don't know them in advance.

To randomly generate the next state ③ we just have to use the discrete distribution assigned to the current state ⑤.

4.5.2 [5200] Software development & error detection

Our next goal is to model the software development process of a team. Each phase takes the team 4 days to complete, and, at the end of each phase the testing team tests the software, and there can be 3 outcomes:

- **no error** is introduced during the phase (we can't actually know it, let's suppose there is an all-knowing "oracle" that can tell us there aren't any errors)
- **no error detected** means that the "oracle" detected an error, but the testing team wasn't able to find it
- error detected means that the "oracle" detected an error, and the testing team was able to find it

If we have **no error**, we proceed to the next phase... the same happens if **no error was detected** (because the testing team sucks and didn't find any errors). If we **detect an error** we either reiterate the current phase (with a certain probability, let's suppose 0.8), or we go back to

one of the previous phases with equal probability (we do this because, if we find an error, there's a high chance it was introduced in the current phase, and we want to keep the model simple).

In this exercise we take the parameters for each phase (the probability to introduce an error and the probability to not detect an error) from a file.

```
#include <...>
using real_t = double;
const size_t HORIZON = 800, PHASES_SIZE = 3;
enum Outcome (1) {
    NO ERROR = 0,
    NO ERROR DETECTED = 1,
    ERROR DETECTED = 2
};
int main() {
    std::random_device random_device;
    std::default_random_engine urng(random_device());
    std::uniform_real_distribution<> uniform_0_1(0, 1);
    std::vector<std::discrete distribution<>>
        phases_error_distribution;
    {
        std::ifstream probabilities("probabilities.csv");
        real_t probability_error_introduced,
            probability_error_not_detected;
        while (probabilities >> probability error introduced >>
               probability error not detected)
            phases_error_distribution.push_back(
                (2) std::discrete_distribution<>({
                    1 - probability_error_introduced,
                    probability_error_introduced *
                        probability_error_not_detected,
                    probability error introduced *
                        (1 - probability_error_not_detected),
                }));
        probabilities.close();
        assert(phases_error_distribution.size() ==
               PHASES_SIZE);
    }
    real_t probability_repeat_phase = 0.8;
    size t phase = 0;
```

```
std::vector<size_t> progress(PHASES_SIZE, 0);
   std::vector<Outcome> outcomes(PHASES_SIZE, NO_ERROR);
    for (size_t time = 0; time < HORIZON; time++) {</pre>
        progress[phase]++;
        if (progress[phase] == 4) {
            outcomes[phase] = static_cast<Outcome>(
                phases_error_distribution[phase](urng));
            switch (outcomes[phase]) {
            case NO ERROR:
            case NO_ERROR_DETECTED:
                phase++;
                break;
            case ERROR_DETECTED:
                if (phase > 0 && uniform_0_1(urng) >
                                      probability_repeat_phase)
                    phase = std::uniform_int_distribution<>(
                        0, phase - 1)(urng);
                break;
            }
            if (phase == PHASES_SIZE)
                break;
            progress[phase] = 0;
        }
   }
    return 0;
}
```

Listing 10: software/5300/main.cpp

TODO: class enum vs enum. We can model the outcomes as an enum ①... we can use the discrete_distribution trick to choose randomly one of the outcomes ②. The other thing we notice is that we take the probabilities to generate an error and to detect it from a file.

4.5.3 Optimizing costs for the development team [5300]

If we want we can manipulate the "parameters" in real life: a better experienced team has a lower probability to introduce an error, but a higher cost. What we can do is:

- 1. randomly generate the parameters (probability to introduce an error and to not detect it)
- 2. simulate the development process with the random parameters

By repeating this a bunch of times, we can find out which parameters have the best results, a.k.a generate the lowest development times (there are better techinques like simulated annealing, but this one is simple enough for us).

4.5.4 Key performance index [5400]

We can repeat the process in exercise [5300], but this time we can assign a parameter a certain cost, and see which parameters optimize cost and time (or something like that? Idk, I should look up the code again).

4.6 Complex systems

- 4.6.1 Insulin pump [6100]
- 4.6.2 Buffer [6200]
- 4.6.3 Server [6300]

- 5 Exam
- 5.1 Development team (time & cost)
- 5.2 Backend load balancing
- $5.2.1~\mathrm{Env}$
- 5.2.2 Dispatcher, Server and Database
- 5.2.3 Response time
- 5.3 Heater simulation
- 5.4 Task management

6 MOCC library

Model CheCking

6.1 Observer Pattern

Basically: the "Observer Pattern" [10] can be used because a MDP is like an entity that "is notified" when something happens (receives an input, in fact, in the case of MDPs, another name for input is "action"), and notifies other entities (output, or reward)

6.2 C++ generics & virtual methods

Generics allow to connect MDPs more safely, as the inputs and outputs are typed! (It's still not fault-proof)

7 Extras

7.1 VDM (Vienna Development Method)

7.1.1 It's cool, I promise

• Alloy? Maybe it's a good alternative, haven't tried it enough

7.1.2 VDM++ to design valid UMLs

7.2 Advanced testing techinques (in Rust & C)

- TODO: cite "Rust for Rustaceans"
- TODO: unit tests aren't the only type of test

7.2.1 Mocking (mockall)

- 7.2.2 Fuzzying (cargo-fuzz)
- 7.2.3 Property-based testing
- 7.2.4 Test augmentation (Miri, Loom, Valgrind)

7.2.5 Performance testing

- Rust is very focused on performance
- TODO: non-functional requirements

7.3 UI testing?

7.3.1 Playwright

7.4 Model checking with Bevy (Rust)

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