University of Copenhagen

MSC IN ACTUARIAL MATHEMATICS

Complete Theory

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 $\begin{array}{c} \textit{Date:} \\ \textit{February 13, 2023} \end{array}$



Abstract

This document contain a comprehensive outline of theory on probability theory and mathematical statistics applied in finance, life insurance and non-life insurance.

 $\label{lem:condition} \begin{tabular}{ll} Keywords: probability theory, insurance mathematics, life insurance, non-life insurance, stochastic differential equations. \end{tabular}$

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Introduction

1.1 Abbreviations

Below is given the abbreviations used when referencing to books:

Chapter	Abbreviation	Source
Basic Life Insurance Mathematics Stochastic Processes in Life Insurance Mathematics		
Life Insurance Mathematics	Asmussen	Risk and Insurance: A Graduate Text by Soren Asmussen and Mogens Steffensen (2020).
	Bladt	Notes from lectures in Liv2.
Topics in Life Insurance Mathematics	Asmussen	Risk and Insurance: A Graduate Text by Soren Asmussen and Mogens Steffensen (2020).
Continuous Time Finance	Bjork	Arbitrage Theory in Continuous Time (Fourth edition) by Thomas Bjork, Oxford University Press (2019).
Basic Non-Life Insurance Mathematics Stochastic Processes in Life Insurance Mathematics		
Topics in Non-Life Insurance Mathematics		
Probabilistic Machine Learning Quantative Risk Management	None	Slides from lectures.
Measure Theory	Bjork	Arbitrage Theory in Continuous Time (Fourth edition) by Thomas Bjork, Oxford University Press (2019).
	Protter	Probability Essentials (2. edition) by Jean Jacod and Philip Protter (2004).
Random Variables	Bjork	Arbitrage Theory in Continuous Time (Fourth edition) by Thomas Bjork, Oxford University Press (2019).
	Hansen	Stochastic Processes (2. edition) by Ernst Hansen (2021).
Discrete Time Stochastic Processes	Hansen	Stochastic Processes (2. edition) by Ernst Hansen (2021).

Chapter	Abbreviation	Source
Continuous Time Stochastic Processes	Bjork	Arbitrage Theory in Continuous Time (Fourth edition) by Thomas Bjork, Oxford University Press (2019).
Stochastic Calculus	Bjork	Arbitrage Theory in Continuous Time (Fourth edition) by Thomas Bjork, Oxford University Press (2019).
Linear Algebra	Bladt Wiki	Notes from lectures in Liv2. Wikipedia

1.2 To-do reading

Chapter	Note	Progress
Liv2	Properties of product integral	Done
	Relation to kolmogorov and markov chains	Bevis 2.3 og 2.12
ML	Risk measure and loss (First week in general)	
QRM	Finish risk measures	
StokLiv	Week 1	
StokLiv	Week 2	
SkadeStok	Week 1	
SkadeStok	Week 2	
Liv1	Week 1	
Liv1	Week 2	
Nonlife1	Week 1	
Nonlife1	Week 2	
MI	Insert notes	
Stok	Insert notes	

Basic Life Insurance Mathematics

Noget indhold

Stochastic Processes in Life Insurance Mathematics

Noget indhold

Topics in Life Insurance Mathematics

4.1 Markov Jump Processes

Life insurance mathematics revolves arrow Markov processes in continuous time on some discrete (or at least countable) state space. This may in the simplest model be the life-death model or it can be a generalized life-death model with interest levels. In any case, we use Markov processes to determine the possible payments that may occur in the future or have occurred in the past. Let us start by define the Markov jump process as follows.

Definition 2.1. (Bladt) A continuous-time stochastic process $\{X_t\}_{t\geq 0}$ taking values in a countable space E is called a Markov jump process with state space E if for all E is the value of E in the value of E in the value of E is the value of E in the value of E in the value of E in the value of E is the value of E in the value of E

$$P(X_{t_n}=i_n \mid X_{t_{n-1}}=i_{n-1},...,X_{t_1}=i_1,X_0=i_0) = P(X_{n-1}=i_n \mid X_{n-1}=i_{n-1}).$$

It is seen that process have some property of "forgetability" in the sense that the jump to a fixed state at some future time only depends on the current state, not the sample path of X. We may then define the transition probabilities as

$$p_{ij}(s,t) = P(X_t = j \mid X_s = i),$$

being the probability of the process being in state j at time t given that the process is in state i at time s. This gives the natural transition matrix

$$\mathbf{P}(t,s) = \big\{p_{ij}(s,t)\big\}_{i,j \in E}, \quad s \leq t.$$

In general, we study the time in-homogeneous models where the above is possibly different for any pair $s \le t$. There does however exist the special case, where the probabilities only depends on the time h = t - s. We call these time homogeneous Markov jump processes and they have the form

$$p_{ij}(t,s) = p_{ij}(t-s). \label{eq:pij}$$

I.e. the probability of jumping from i to j is the same so long the interval is the same length. Let us now define the transition intensities as below.

Definition 2.2. (Bladt) Assume that the limit

$$\mathbf{M}(s) = \left\{ \mu_{ij}(s) \right\}_{i,j \in E} = \lim_{h \to 0_+} \frac{\mathbf{P}(s,s+h) + \mathbf{I}}{j}$$

exist for all $s \ge 0$. The matrices $\mathbf{M}(s)$ are called intensity matrices. The the special case of time-homogenous processes we have $\mathbf{M}(s) = \mathbf{M}$ is a constant matrix.

We may interpret the intensity matrices in the infinitesimal interpretation as

$$\mathbf{P}(s, s+h) = \mathbf{I} + h\mathbf{M}(s) + o(h),$$

i.e.

$$p_{ij}(s, s + h) = \delta_{ij} + \mu_{ij}(s)h + o(h),$$

where $\delta_{ij} = 1_{i=j}$ is the Kronecker delta. We furthermore see, that the matrix $\mathbf{M}(s)$ has non-positive diagonal and non-negative upper and lower triangle. Furthermore, the matrix has zero row sums and so it follows that

$$\mu_{ii}(s) = -\sum_{j \neq i} \mu_{ij}(s) := -\mu_i(t),$$

for all $i \in E$. Returning to the infinitesimal interpretation we may write the dynamics of μ_{ij} in the following manner

$$d\mu_{ij}(t) = \mu_{ij}(t) \ dt = P(X_{t+dt} = j \mid X_t = i)$$

i.e. the probability of a jump from state i to state j during [t, t + dt). On the other hand we have

$$1 - \mu_{ii}(t) dt = P(X_{t+dt} = i \mid X_t = i),$$

i.e. the probability that no jumps occur during [t, t + dt). We can then define the conditional probability that a jump, if one occurs, is to state j from i as

$$\begin{split} q_{ij}(t) &= P(X_{t+dt} = j \mid X_t = i, X_{t+dt} \neq i) = \frac{P(X_{t+dt} = j \mid X_t = i)}{P(X_{t+dt} \neq i \mid X_t = i)} \\ &= \frac{\mu_{ij}(t)}{\mu_i(t)} = -\frac{\mu_{ij}(t)}{\mu_{ii}(t)}. \end{split}$$

We can now look at the time until the next jump as the random variable

$$T(s) = \inf\{u > 0 : X_{s+u} \neq X_s\},\$$

i.e. T(s) is a positive random variable. If we condition on $X_s = i$ then $T(s) \mid X_s = i$ is the time until X jumps out of state i into some other state $j \neq i$. If we set $S_i(t) = P(T(s) > t \mid X_s = i)$ and $f_i(t)$ is the density of $T(s) \mid X_s = i$, then we have

$$\mu_i(u) = \frac{f_i(u)}{S_i(u)} = -\frac{S_i'(u)}{S_i(u)} = -\frac{d}{du}\log(S_i(u))$$

which is solved for

$$S_i(t) = P(T(s) > t \mid X_s = i) = \exp\left(-\int_s^t \mu_i(u) \ du\right)$$

and then

$$f_i(t) = \exp\left(-\int_s^t \mu_i(u) \ du\right) \mu_i(t).$$

Theorem 2.3. (Bladt) (Kolmogorov' Differential Equations) Relation between $\mathbf{M}(s)$ and $\mathbf{P}(s,t)$ are given by Kolmogorov' forward and backward differential equations:

$$\frac{\partial}{\partial s} \mathbf{P}(s,t) = -\mathbf{M}(s) \mathbf{P}(s,t)$$

and

$$\frac{\partial}{\partial t} \mathbf{P}(s,t) = \mathbf{P}(s,t) \mathbf{M}(t).$$

the soluction of which is given by

$$\mathbf{P}(s,t) = \prod_{s}^{t} (\mathbf{I} + \mathbf{M}(x) \ dx).$$

In the time-homogeneous case we have that

$$\frac{d}{dt}\mathbf{P}(t) = \mathbf{MP}(t) = \mathbf{P}(t)\mathbf{M}$$

with solution

$$\mathbf{P}(t) = \exp(\mathbf{M}t).$$

Proof.

We may subdivide the interval [s,t] into [s,u] and [u,t] and use theorem 1.5 from the chapter on Product Integrals to obtain the Chapman-Kolmogorov's equation

$$\mathbf{P}(s,t) = \mathbf{P}(s,u)\mathbf{P}(u,t)$$

and in the time homogeneous case we have

$$\mathbf{P}(s+t) = \mathbf{P}(s)\mathbf{P}(t).$$

Finally, let us define a stopping time and introduce the strong markov property.

Definition 2.4. (Bladt) (Stopping time) Let X be a Markov jump process and let $\mathcal{F}_t = \sigma(X_s: s \leq t)$. A nonnegative random variable τ is called a stopping time for X of $\{\tau \leq t\} \in \mathcal{F}_t$ for all $t \setminus T$ he σ -algebra of the process up to a stopping time τ , \mathcal{F}_{τ} , is defined by the measurable sets A for which

$$\forall t \geq 0 : A \cap \{\tau \leq t\} \in \mathcal{F}_t.$$

We see that a stopping time in an intuitive sense is a random variable which at any time t we know either that the stopping time is in the future i.e. $\tau > t$ or that the stopping time already occurred and so τ becomes known at time t. Furthermore, we see that $\{\tau \leq t\}$ is \mathcal{F}_t measurable and so we only need information from the Markov jump process in order to determine the above.

Theorem 2.5. (Bladt) (Strong Markov property) Every Markov jump process satisfies the strong Markov property, i.e., for all $0 \le h_1 \le h_2 \le \dots \le h_n$, we have that

$$P(X_{\tau+h_1} = i_1, ..., X_{\tau+h_n} = i_n \mid \mathcal{F}_t) = P(X_{\tau+h_1} = i_1, ..., X_{\tau+h_n} = i_n \mid X_\tau)$$

on $\{\tau < \infty\}$. For time-homogeneous processes this further reduces to

$$P(X_{\tau+h_1} = i_1, ..., X_{\tau+h_n} = i_n \mid \mathcal{F}_t) = P_{X_\tau}(X_{\tau+h_1} = i_1, ..., X_{\tau+h_n} = i_n)$$

also on $\{\tau < \infty\}$.

4.2 Phase-type distributions

In life insurance models we often, that is always, have a model with one absorbing state, namely, death. This means that the state space E will have the form of $p \ge 1$ transient states (states that may interact in both ways) and one absorbing state (a state that is never moved away from). In this case we would often like to study the distribution of

$$\inf\{t \ge 0 : X(t) = p + 1\},\$$

where p+1 is the absorbing state and 1, ..., p are the transient states. We will assume that P(X(0) = p+1) = 0 i.e. the above is at least zero and never $-\infty$. Let us now formalise this setup.

Consider a time-inhomogeneous Markov jump process $\{X_t\}_{t\geq 0}$ on the finite state-space $E=\{1,...,p,p+1\}$ where the states 1,...,p are transient states and p+1 is the only absorbing state. This implies that the intesity matrix of $\{X_t\}_{t\geq 0}$ take the form

$$\Lambda(t) = \begin{bmatrix} \mathbf{T}(t) & \mathbf{t}(t) \\ \mathbf{0} & 0 \end{bmatrix}.$$

In the above $\mathbf{T}(t)$ is a $p \times p$ matrix function and $\mathbf{t}(t)$ is a $p \times 1$ matrix function. We now define the initial distribution of X_0 as

$$\{P(X_0=i)\}_{i\in E}\,_{\{p+1\}}= \ = \begin{pmatrix} \pi_1 \\ \vdots \\ \pi_p \end{pmatrix}^\top.$$

By assumption we have $P(X_0 = p + 1) = 0$ and so $\sum_{i=1}^{p} \pi_i = 1$ and $\$ is a proper distribution.

Definition 2.10. (Bladt) (Phase-Type distribution) Let

$$\tau = \inf\{t > 0 : X(t) = p + 1\}$$

denote the time until absorption of X. The distribution of τ is then said to be an inhomogeneous phase-type distribution with representation $(\ ,\mathbf{T}(t))$ and we write $\tau \sim IPH(\ ,\mathbf{T}(t))$.

Lemma 2.11. (Bladt) We have the following decomposition:

$$\mathbf{P}(s,t) = \prod_{s}^{t} (\mathbf{I} + \ (u) \ du) = \begin{bmatrix} \prod_{s}^{t} (\mathbf{I} + \mathbf{T}(u) \ du) & \mathbf{e} - \prod_{s}^{t} (\mathbf{I} + \mathbf{T}(u) \ du) \mathbf{e} \\ \mathbf{0} & 1 \end{bmatrix},$$

where $\mathbf{e} = (1, 1, ..., 1)^{\top}$.

Theorem 2.12. (Bladt) Assume that $\tau \sim IPH(\ , \mathbf{T}(t))$. Then the density f and the distribution function F of τ are given by

$$f(x) = \prod_{0}^{x} (\mathbf{I} + \mathbf{T}(u) \ du) \mathbf{t}(x),$$

$$F(x) = 1 - \prod_{0}^{x} (\mathbf{I} + \mathbf{T}(u) \ du) \mathbf{e}.$$

Proof.

Theorem 2.13. (Bladt) If $\tau \sim IPH(\cdot, \mathbf{T}(t))$ then

$$P(\tau > s + t \mid \tau > s) = \frac{\prod_{0}^{s} (\mathbf{I} + \mathbf{T}(u) \ du)}{\prod_{0}^{s} (\mathbf{I} + \mathbf{T}(u) \ du) \mathbf{e}} \prod_{s}^{t} (\mathbf{I} + \mathbf{T}(u) \ du) \mathbf{e}$$

so that

$$\tau - s \ | \ \{\tau > s\} \sim IPH(\ , \mathbf{S}(\cdot)),$$

where $\mathbf{S}(u) = \mathbf{T}(s+u)$ and

$$= \frac{\prod_{0}^{s} (\mathbf{I} + \mathbf{T}(u) \ du)}{\prod_{0}^{s} (\mathbf{I} + \mathbf{T}(u) \ du) \mathbf{e}}.$$

Corollary 2.14. (Bladt) If $\tau \sim IPH(\ , \mathbf{T}(t))$ and $\mathbf{T}(t_1)$ and $\mathbf{T}(t_2)$ commute for all $t_1, t_2 \geq 0$, then the density f and the distribution function F of τ are given by

$$f(x) = \exp\left(\int_0^x \mathbf{T}(u) \ du\right) \mathbf{t}(x),$$

$$F(x) = 1 - \exp\left(\int_0^x \mathbf{T}(u) \ du\right) \mathbf{e}.$$

Example (Approximation in time-inhomogeneous case).

Consider the Markov jump process on the state space $E = \{1, 2, 3\}$ where 1 is the state "alive" (working), 2 is "disabled" and 3 is the state "death". We assume that $\Lambda(t)$ has the structure:

$$\begin{split} &\Lambda(t) = \begin{bmatrix} \mu_{11}(t) & \mu_{12}(t) & \mu_{13}(t) \\ \mu_{21}(t) & \mu_{22}(t) & \mu_{23}(t) \\ \mu_{31}(t) & \mu_{32}(t) & \mu_{33}(t) \end{bmatrix} \\ &= \begin{bmatrix} -\mu_{12}(t) - \mu_{13}(t) & 0.000015 + 10^{(4.6-10+0.015\cdot t)} & 0.00005 + 10^{(4.6-10+0.05\cdot t)} \\ 0.000005 + 10^{(4.6-10+0.015\cdot t)} & -\mu_{21}(t) - \mu_{23}(t) & 0.0001 + 10^{(4.6-10+0.05\cdot t)} \\ 0 & 0 & 0 & 0 \end{bmatrix} \end{split}$$

We can now implement this intensity matrix as a function in R:

```
mu12 <- function(t) {</pre>
  0.000015 + 10^{(4.6-10+0.015*t)}
mu13 <- function(t) {</pre>
  0.00005 + 10^{(4.6-10+0.05*t)}
mu11 <- function(t) {</pre>
  -mu12(t)-mu13(t)
mu21 <- function(t) {</pre>
  0.000005 + 10^{(4.6-10+0.015*t)}
}
mu23 <- function(t) {</pre>
  0.0001 + 10^{(4.6-10+0.05*t)}
mu22 <- function(t) {</pre>
  -mu21(t)-mu23(t)
}
mu31 <- function(t) {0}</pre>
mu32 <- function(t) {0}</pre>
mu33 <- function(t) {0}</pre>
M <- function(t) {</pre>
  matrix(
    c(mu11(t),mu21(t),mu31(t),
       mu12(t), mu22(t), mu32(t),
       mu13(t),mu23(t),mu33(t)),
    ncol = 3
  )
}
M(0)
```

```
## [,1] [,2] [,3]
## [1,] -7.296214e-05 1.898107e-05 5.398107e-05
## [2,] 8.981072e-06 -1.129621e-04 1.039811e-04
## [3,] 0.000000e+00 0.000000e+00 0.000000e+00
```

We see that the intensities is choosen such that the following holds for any time interval [t, t+dt)

- 1. The transition $1 \to 3$ is less likely than $2 \to 3$,
- 2. The transition $2 \to 1$ is less likely than $1 \to 2$,
- 3. The transition $1 \to 2$ is less likely than $1 \to 3$.

We now wish to compute the density and distribution of

$$\tau = \inf\{t \ge 0 : X(t) = 3\}.$$

We assume that $\pi = (1,0)$ i.e. the person is alive. Calculating the distribution of τ is then the distribution of the length of a newborn child lifespan. From theorem 2.12 this ammount to calculating the product integral of

$$\mathbf{T}(t) = \{\mu_{ij}(t)\}_{i,j \in \{1,2\}}$$

i.e.

$$\prod_{0}^{x} (\mathbf{I} + \mathbf{T}(u) \ du).$$

To du this we simply approximate with stepsize h = 1/n (for some $n \ge 1$) by

$$\prod_{0}^{x+h}(\mathbf{I}+\mathbf{T}(u)\ du)=h\prod_{0}^{x}(\mathbf{I}+\mathbf{T}(u)\ du)\mathbf{T}(x)+\prod_{0}^{x}(\mathbf{I}+\mathbf{T}(u)\ du),$$

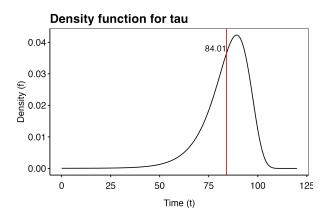
and

$$\prod_{0}^{0} (\mathbf{I} + \mathbf{T}(u) \ du) = \mathbf{I}.$$

This is done in the code below

```
n \leftarrow 12 \#monthly
h \leftarrow 1/n
N <- 120 #max number years
T_matrix <- function(t) {</pre>
  matrix(
    c(mu11(t),mu21(t),
      mu12(t),mu22(t)),
    ncol = 2
  )
}
T_matrix(0)
##
                   [,1]
                                    [,2]
## [1,] -7.296214e-05 1.898107e-05
## [2,] 8.981072e-06 -1.129621e-04
t_matrix <- function(t) {</pre>
  matrix(
    c(mu13(t),mu23(t)),
    ncol=1
  )
}
t_matrix(0)
##
                  [,1]
## [1,] 5.398107e-05
## [2,] 1.039811e-04
library(rlist)
\#Initial condition t=0
T_product_integral <- list(diag(c(1,1)))</pre>
pi \leftarrow matrix(c(1,0),ncol = 1)
f <- t(pi) %*% diag(c(1,1)) %*% t_matrix(0)
F \leftarrow 1-t(pi) %*% diag(c(1,1)) %*% matrix(c(1,1),ncol = 1)
for (i in 1:(N*n)) {
  x_1 <- i/n
  x_0 \leftarrow (i-1)/n
  T_0 \leftarrow T_{matrix}(x_0)
  M_0 <- T_product_integral[[i]]</pre>
```

```
M_1 <- h*M_0 %*% T_0 + M_0
T_product_integral <- list.append(T_product_integral,M_1)
f <- c(f,t(pi) %*% M_1 %*% t_matrix(x_1))
F <- c(F,1-t(pi) %*% M_1 %*% matrix(c(1,1),ncol = 1))
}</pre>
```



We see that by calculating

$$E[\tau\mid X_0=1] = \int_0^\infty \tau f(\tau)\ d\tau \approx \int_0^{120} \tau f(\tau)\ d\tau \approx \frac{1}{n} \sum_{i=0}^{120 \cdot n} i/n \cdot f(i/n).$$

that the life expectation is 84.01 years. We have that we may approximate f and F with an arbitrary precision by choosing n appropriately.

Continuous Time Finance

This topic revolves around the theory of the Brownian motion and martingale processes. Other main topics are the binomial model and an introduction to financial derivatives. Financial derivatives is contingent on the outcome of a stochastic process at some future time t=T and often is a function Φ of some assets price S_t . As such the derivative will give a stochastic payout, at time t=T of the size $X_T=\Phi(S_T)$. Naturally we want to say something about the fair price of the derivative in the form of

$$\Pi_t(X_T) = \mathbb{E}\left[\Phi(S_T) \mid \mathcal{F}_t\right],$$

where $\mathcal{F}_t \subset \mathcal{F}$ is the available information at time t. We will by defualt interest the times t=0 as today and t=T as tomorrow. This indeed require some fundamental understanding of the behaviour of the asset price S_t . This lead us over to discussing the process in center of the Black-Scholes model: the Brownian motion.

5.1 Discrete time models

5.1.1 One-period time models

The study of this course is the **European call** option (and put option). This financial derivative is an agreement between two parties where the holder of the option has the right to "exercise" the derivative, at a future time t = T. Exercising means buying an asset at a certain agreed opon price-strike K. In the case of the put-option: the holder has the right (but not obligation) to sell the asset at the strike price K. As such the derivative has the payoff

Call option:
$$\Phi(S_T) = (S_T - K)^+$$
, Put option: $\Phi(S_T) = (K - S_T)^+$.

Our objective is to understand when an arbitrage exist and to find the fair price of these derivative. The strategy in pricing is finding a replicating portfolio with the same payoff as the option (with probability one) and then price the derivative accordingly.

5.1.1.1 Model description

In the one-period model we consider the simplest possible market. We have two distinct times t=0 (today) and t=1 (tomorrow) and we may buy any portfolio as a mixture of bonds and one stock. We denote the bonds price by B_t and the stocks price by S_t and we assume the following:

$$B_0 = 1, \ B_1 = 1 + R, \qquad S_0 = s, \ S_1 = \begin{cases} s \cdot u, & with \ probability \ p_u. \\ s \cdot d, & with \ probability \ p_d. \end{cases}$$

We may introduce Z as the random variable

$$Z = u \cdot (I) + d \cdot (1 - I),$$

for an bernoulli variable I with succes probability p_u . Naturally, we assume $d \leq (1+R) \leq u$ (this is imperative to ensure no arbitrage as we will see).

5.1.1.2 Portfolios and arbirtage

We study any portfolio on the (B,S) market as a vector h=(x,y) where x is the amount of bonds and y is the amount of stock held in the portfolio. Notice that we allow for shorting, that is x < 0 or y < 0. As such, we have that $h \in \mathbb{R}^2$. In this we have made some unrealistic, but attractable assumptions included in the assumptions:

- We allow short positions and fractional holding, i.e. $h \in \mathbb{R}^2$,
- We assume no spread between ask and bids,
- No transaction costs and
- A completely liquid market i.e. we may borrow and buy as much stock and bonds as wanted.

Given that we have chosen a portfolio h we may introduce the value process.

Definition 2.1. (Bjork) The value process of the portfolio $h \in \mathbb{R}^2$ is the stochastic process

$$V_t^h = xB_t + yS_t, \ t = 0, 1.$$

Given this notation we may define what an arbitrage is.

Definition 2.2. (Bjork) An arbitrage is a portfolio h with the properties: 1) $V_0^h = 0$, 2) $P(V_1^h \ge 0) = 1$ and 3) $P(V_1^h > 0) > 0$.

That is h is an deterministic money-machine where we at least never loose any money. Granted the bonds give a deterministic non-negative return, but an arbitrage does not require any money out of pocket. With the notion of an arbitrage we will show the first proposition regarding the choice of R, u, d as defined above.

Proposition 2.3. (Bjork) The one-period binomial model is arbitrage free if and only if the following inequality hold:

$$d \le (1+R) \le u. \tag{2.1}$$

Proof.

The statement is proofed by contradiction. Assume that d > 1 + R holds. Then by definition u > d > 1 + R. Notice that any portfolio satisfying $V_0^h = 0$ must satisfy

$$0 = xB_0 + yS_0 = x + ys \iff x = -ys$$

That is for some choice y the only arbitrage candidate is the portfolio h = (-ys, y). Calculating the value at time t = 1 we have

$$V_1^h = -ys \cdot (1+R) + y \cdot s \cdot Z = ys(Z-1-R)$$

However since $Z \ge d$ we have $Z - (1 + R) \ge 0$ and therefore an arbitrage (for y > 0). The other inequality 1 + R > u follows analog steps. Simply choose some y < 0 and the result follows.

From inequality (2.1) we see that since 1+R is between u and d we may find a pair $q_d, q_u \ge 0$ with $q_d+q_u=1$ such that

$$1+R=q_u\cdot u+q_d\cdot d.$$

This yields the important risk neutral valuation formula as summed op in the following definition

Definition 2.4. (Bjork) A probability measure Q is called a martingale measure if the following condition holds:

$$S_0 = \frac{1}{1+R} E^Q[S_1].$$

The above measure Q is the measure $Q(Z=d)=q_d$ and $Q(Z=u)=q_u$ for the binomial model. This does in fact yield the risk neautral valuation formula:

$$\begin{split} S_0 &= \frac{1}{1+R} E^Q[S_1] = \frac{1}{1+R} (Q(Z=d) \cdot d \cdot s + Q(Z=u) \cdot u \cdot s) \\ &= s \frac{1}{1+R} (q_d \cdot d + q_u \cdot u) = s, \end{split}$$

where we simply use $1+R=q_d\cdot d+q_u\cdot u$. We call this the risk neautral valuation formula because it in some sense gives an expected discounted value of the future stock price. We end this endavour with reformulating the arbitrage proposition and determining the values of the Q-measure.

Proposition 2.5. (Bjork) The one-period binomial model is arbitrage free if and only if there exists a martingale measure Q.

Proposition 2.6. (Bjork) The one-period binomial model has martingale probabilities given by:

$$\begin{cases} q_u = \frac{(1+R)-d}{u-d}, \\ q_u = \frac{u-(1+R)}{u-d}. \end{cases}$$

5.1.1.3 Contingent Claims

This chapter revolves around the financial derivative and we start by stating the definition of the financial derivative.

Definition 2.7. (Bjork) A contingent claim (financial derivative) is any stochastic variable X of the form $\Phi(Z)$, where Z is the stochastic varible driving the stock price process.

We may also call the function Φ the **contract function** as it states how the contract is resolved once the stochastic variable Z has been realised. Our objective is now to study, what a buyer of said contract would have to pay at any given time t. We call the fair price of X at time t: $\Pi_t[X]$. As such it is easy to see that the fair price at the time of maturity T is simply the payout X i.e. $\Pi_T[X] = X$. Our strategy is to find a replicating portfolio h and determine the price of said portfolio.

Definition 2.8. (Bjork) A contingent claim X can be **replicated**, or said to be **reachable** if there exist a portfolio h such that

$$V_1^h = X$$
,

with probability one. In that case, we say that the portfolio h is a **hedging** portfolio or a **replicating** portfolio. If all claims can be replicated we say that the market is **complete**.

Our pricing strategy is then to determine the value process of the replicating portfolio and then by the first pricing principle below we say that the price is imply the value of the replicating portfolio.

Pricing principle 1. If a clain X is reachable with replicating portfolio h, then the only reasonable price process for X is given by

$$\Pi_t[X] = V_t^h$$
.

Notice, that this assumes that a replicating portfolio exist and even so we have a uniqueness statement to solve. We end this section by writing two important results.

Proposition 2.9. (Bjork) Suppose that a claim X is reachable with replicating portfolio h. Then any price at time $t \geq 0$ of the claim X other than the value process of h will lead to an arbitrage on the extended market (B, S, X).

Proposition 2.10. (Bjork) If the one-period binomial model is free of arbitrage, then it is also complete.

The hedging portfolio in the one-period binomial model is given by the portfolio (x, y) below

$$x = \frac{1}{1+R} \cdot \frac{u\Phi(d) - d\Phi(u)}{u - d},$$
(2.2)

$$y = \frac{1}{s} \cdot \frac{\Phi(u) - \Phi(d)}{u - d}.\tag{2.3}$$

5.1.1.4 Risk Neutral Valuation

We see that since the one-period model is complete we can price any contingent claim and we see that

$$\begin{split} \Pi_0[X] &= \frac{1}{1+R} \cdot \frac{u\Phi(d) - d\Phi(u)}{u-d} + s\frac{1}{s} \cdot \frac{\Phi(u) - \Phi(d)}{u-d} \\ &= \frac{1}{1+R} \left\{ \frac{u\Phi(d) - d\Phi(u)}{u-d} + (1+R)\frac{\Phi(u) - \Phi(d)}{u-d} \right\} \\ &= \frac{1}{1+R} \left\{ \frac{(1+R) - d}{u-d} \Phi(u) + \frac{u - (1+R)}{u-d} \Phi(d) \right\} \\ &= \frac{1}{1+R} E^Q[X]. \end{split}$$

i.e. the price at time t = 0 should simply be the expected discounted payout according to the martingale measure. This leads to the important pricing proposition:

Proposition 2.11. (Bjork) If the one-period binomial model is free of arbitrage, then the arbitrage free price of a contingent claim X is given by

$$\Pi_0[X] = \frac{1}{1+R} E^Q[X]. \tag{2.4}$$

Here the martingale measure Q is uniquely determined by the relation

$$S_0 = \frac{1}{1+R} E^Q[S_1],\tag{2.5}$$

and the explicit expressions for q_u and q_d are given in proposition 2.6. Furthermore the claim X can be replicated using the portfolio

$$x = \frac{1}{1+R} \cdot \frac{u\Phi(d) - d\Phi(u)}{u - d},$$
(2.6)

$$y = \frac{1}{s} \cdot \frac{\Phi(u) - \Phi(d)}{u - d}.\tag{2.7}$$

5.1.2 Multi-period model

The one-period binomial model can easily be extended to a multi-period model, by assuming that the bond and stock pricess evolve by the processes:

$$t \ge 1: B_t = (1+R)B_{t-1}$$
 and $B_0 = 1$,

$$t \ge 1: S_t = Z_{t-1}S_{t-1}$$
 and $S_0 = s$,

where we obviously have that $B_t = (1+R)^t$ for $t \ge 0$. In the above Z_t is u with probability p_u and d with probability p_d . In this context, we need to define a portfolio in terms of a strategy.

Definition 2.13. (Bjork) A portfolio strategy is a stochastic process on $\{1, ..., T\}$

$$h = \{h_t = (x_t, y_t); \ t = 1, ..., T\}$$

such that h_t is a function of $S_0, S_1, ..., S_{t-1}$. For a given portfolio strategy h we set $h_0 = h_1$ by convention. The associated value process corresponding to the portfolio h is defined by

$$V_t^h = x_t(1+R) + y_t S_t. \label{eq:Vh}$$

Given this notation we may define what an arbitrage is, but first we introduce the notion of a self-financing portfolio. A self-financing portfolio in an intuative sense is a portfolio that is not withdrawn from or deposited into.

Definition 2.14. (Bjork) A portfolio strategy h is said to be **self-financing** if the following condition holds for all t = 0, ..., T - 1:

$$x_t(1+R) + y_t S_t = x_{t+1} + y_{t+1} S_t$$

The above equation says that the portfolio purchased at time t and helt until t+1 (x_{t+1},y_{t+1}) can only be financed by the market value of the portfolio held from [t-1,t) i.e. (x_t,y_t) . We now define an arbitrage.

Definition 2.15. (Bjork) An arbitrage is a self-financing portfolio h with the properties: 1) $V_0^h = 0$, 2) $P(V_T^h \ge 0) = 1$ and 3) $P(V_T^h \ge 0) > 0$.

The multiperiod binomial model has an just like the oneperiod model a result regarding when an arbitrage exists.

Lemma 2.16. (Bjork) If $d \le (1+R) \le u$ (eq. 2.8) then the multiperiod model is arbitrage-free.

As one can see, the multiperiod model is rather similar to the one period model. We will in the following summarise equivalent statements for the multiperiod model as the ones in the oneperiod model.

Definition 2.17. (Bjork) The martingale probabilities q_u and q_d are defined as the probabilities for which the relation below holds.

$$s = \frac{1}{1+R} E^{Q}[S_{t+1} \mid S_t].$$

Proposition 2.18. (Bjork) The martingale probabilities q_u and q_d are given by

$$\begin{cases} q_u = \frac{(1+R)-d}{u-d}, \\ q_u = \frac{u-(1+R)}{u-d}. \end{cases}$$

Definition 2.19. (Bjork) A contingent claim is a stochastic variable X of the form

$$X = \Phi(S_T),$$

where the contract function is some given real valued function.

Definition 2.20. (Bjork) A given contingent claim X is said to be **reachable** if there exists a self-financing portfolio h such that

$$V_T^h = X$$
,

with probability one. In that case we say that the portfolio h is a **hedging** portfolio or a **replicating** portfolio. If all claims can be replicated we say that the market is (dynamically) **complete**.

Pricing principle 2. (Bjork) If a claim X is reachable with replicating portfolio h, then the only reasonable price process for X os given by

$$\Pi_t[X] = V_t^h, \ t = 0, 1, ..., T.$$

Proposition 2.21. (Bjork) Assume X is reachable by h, then any price other than V_t^h for some $t \geq 0$ leads to an arbitrage opportunity.

Proposition 2.22. (Bjork) The multiperiod model is complete, i.e. every claim can be replicated by a self-financing portfolio.

Proposition 2.24. (Bjork) (Binomial algorithm) Consider a T-claim $X = \Phi(S_T)$. Then this claim can be replicated using af self-financing portfolio. If $V_t(k)$ denotes the value of the portfolio at the node (t,k) (k referring to k amount of up-moves for the stock), then $V_t(k)$ can be computed recursively by the scheme

$$\begin{cases} V_t(k) = \frac{1}{1+R} \left\{ q_u V_{t+1}(k+1) + q_d V_{t+1}(k) \right\}, \\ V_T(k) = \Phi(su^k d^{T-k}). \end{cases}$$

where the martingale probabilities q_u and q_d are given by

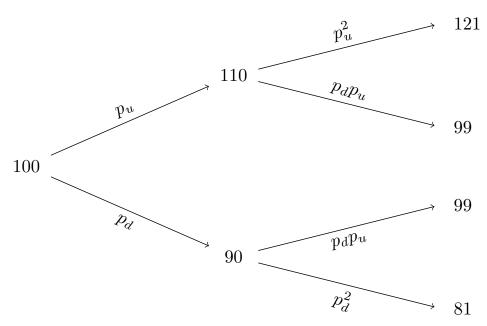
$$\begin{cases} q_u = \frac{(1+R)-d}{u-d}, \\ q_u = \frac{u-(1+R)}{u-d}. \end{cases}$$

With the notation as above, the hedging portfolio is given by

$$\begin{cases} x_t(k) = \frac{1}{1+R} \cdot \frac{uV_t(k) - dV_t(k+1)}{u - d}, \\ y_t(k) = \frac{1}{S_{t-1}} \cdot \frac{V_t(k+1) - V_t(k)}{u - d}. \end{cases}$$

In particular, the arbitrage free price of the claim at t=0 is given by $V_0(0)$.

Example.



Consider R = 0.04, s = 100, u = 1.1, d = 0.9, $p_u = 0.6$ and $p_d = 0.4$. We consider a model of length T = 2 and we want to evaluate the price of the european call option with srike K = 90 that is the contingent claim

$$X = (S_T - K)^+, \qquad \Phi(s) = (s - K)^+.$$

For each time t we know the replicating portfolio, if we know the payoff the following period. Therefore we start from the leaves of the tree and work towards the root. Since the strike price is K = 90 the end result will be the following payoffs:

$$u^2$$
: $(121 - 90)^+ = 31$
 ud : $(99 - 90)^+ = 9$
 du : $(99 - 90)^+ = 9$
 d^2 : $(81 - 90)^+ = 0$

Therefore by the risk neautral valuation formula with $q_u = \frac{(1+R)-d}{u-d} = 0.7$ and $q_d = \frac{u-(1+R)}{u-d} = 0.3$ we have that the cost of the replicating portfolio at time t=1 is respectively

$$\begin{split} u: & \frac{1}{1+R} \left\{ 31 \cdot q_u + 9 \cdot q_d \right\} \approx 23.46 \\ d: & \frac{1}{1+R} \left\{ 9 \cdot q_u + 0 \cdot q_d \right\} \approx 6.06 \end{split}$$

To replicate this payoff at time t = 1 we can use the risk neutral valuation formula once more to find the base cost of the replicating portfolio i.e. the price of X at time t = 0

$$\frac{1}{1+R} \left\{ 23.46 \cdot q_u + 6.06 \cdot q_d \right\} \approx 17.54.$$

Working from the root to the leaves we can now calculate the hedging portfolio at time t = 0, 1 for each path. For time t = 0 we calculate

$$x = \frac{1}{1+R} \cdot \frac{u \cdot 6.06 - d \cdot 23.46}{u - d} \approx -69.46,$$

$$y = \frac{1}{s} \cdot \frac{23.46 - 6.06}{u - d} \approx 0.87$$

We see by calculations that this does indeed replicate the payoff at time t=1:

$$u: V_1^h = (1+R) \cdot x + 110 \cdot y \approx 23.46,$$

 $d: V_1^h = (1+R) \cdot x + 90 \cdot y \approx 6.06.$

We also see by calculation that the initial portfolio does cost the expected 17.54 as

$$x \cdot 1 + y \cdot 100 = 87 - 69.46 = 17.54.$$

Following these steps at time t=1 the portfolios (-86.54,1) (for the up-scenario) and (-38.94,0.5) (for the down-scenario) would arise. Notice when calculating y one has to use the current price $S_1=S_0\cdot Z$ not S_0 . One should also check by similar calculations as above, that these portfolios does indeed replicate the payoff of the contingent claim X. \square

Proposition 2.25. (Bjork) The arbitrage free price at t = 0 of a T-claim X is given by

$$\Pi_0[X] = \frac{1}{(1+R)^T}E^Q[X]$$

where Q denotes the martingale measure, or more explicitly

$$\Pi_0[X] = \frac{1}{(1+R)^T} \sum_{k=0}^T \binom{T}{k} q_u^k q_d^{T-k} \Phi(su^k d^{T-k}).$$

Example.

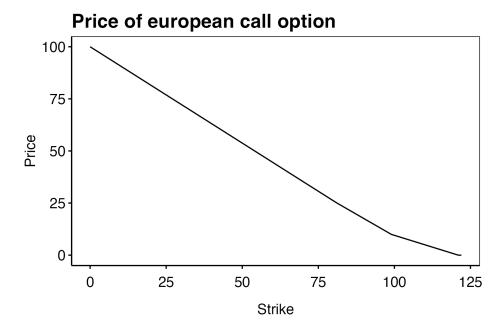


Figure 5.1: The pricing function of the European call option.

We follow an analog example as the one after proposition 2.24. Let K = 90 and we see that

$$\begin{split} &\Pi_0[X] \\ &= \frac{1}{(1+0.04)^2} \sum_{k=0}^2 \binom{2}{k} \cdot 0.7^k \cdot 0.3^{2-k} \cdot \Phi(100 \cdot 1.1^k \cdot 0.9^{2-k}) \\ &= 0.9245562 \cdot \left(\underbrace{1 \cdot 1 \cdot 0.09 \cdot 0}_{k=0} + \underbrace{2 \cdot 0.7 \cdot 0.3 \cdot 9}_{k=1} + \underbrace{1 \cdot 0.49 \cdot 1 \cdot 31}_{k=2}\right) \\ &= 0.9245562 \cdot (0+3.78+15.19) \\ &= 17.53883 \end{split}$$

Since we know that K must meaningfully range in [0,121] we could try to calculate the price of the contingent claim at time t=0 for all integers in this interval. We see that the price range between S_0 and 0 as expected. One can also see that the price changes slope at the prices 99 and 121 as the function is linear in Φ and som realisations loose any effect on the price when the strike is higher than the outcome. \Box

Proposition 2.26. (Bjork) The condition d < (1+R) < u is necessary and sufficient condition for absence of arbitrage.

5.1.3 Generelised one-period model

In the previous we had the simpel model where we only had one stochastic asset S and only one stochastic variable Z determining the future stock price. Now we will generalise this model by introducing N assets and introducing som stochastic behaviour to the system.

5.1.3.1 Model specification

We consider the market consisting of a collection of stochastic prices assets i = 1, ..., N with N-dimensional price process.

$$S_t = \begin{bmatrix} S_t^1 \\ \vdots \\ S_t^N \end{bmatrix}$$

We now assume that S_t is defined on a background space with finite sample space $\Omega = \{\omega_1, ..., \omega_M\}$ with associated probabilities $p_j = P(\omega_j), \ j = 1, ..., M$. We can then for eact time t = 1, ..., T define the $N \times M$ matrix D_t as such

$$D_t = \begin{bmatrix} S^1_t(\omega_1) & \cdots & S^1_t(\omega_M) \\ \vdots & \ddots & \vdots \\ S^N_t(\omega_1) & \cdots & S^M_t(\omega_M) \end{bmatrix}.$$

We will assume that $S_0^1 > 0$ and $S_1^1(\omega_j) > 0$, j = 1, ..., M.

5.1.3.2 Absence of Arbitrage

We now define a **portfolio** as an N-dimensional row vector

$$h = [h^1, \dots, h^N]$$

representing the amount of assets held at time t = 0 and held until t = 1. The value process is then

$$V_t^h = h \cdot S_t = \sum_{i=1}^N h^i S_t^i, \ t = 0, 1. \tag{3.1}$$

For a given $\omega_j \in \Omega$ we have the realisation

$$V_t^h = hS_t(\omega_j) = hd_j = (hD)_j.$$

Definition 3.1. (Bjork) The portfolio h is an arbitrage portfolio fil it satisfies the conditions: $V_0^h = 0$, $P(V_1^h \ge 0) = 1$ and $P(V_1^h > 0) > 0$.

Lemma 3.2. (Bjork) (Farkas' Lemma) Suppose that $d_0, d_1, ..., d_M$ are column vectors in \mathbb{R}^N . Then exactly one of the following problems possesses a solution.

- Problem 1: There exist $\lambda_1,...,\lambda_M \geq 0$ such that $d_0 = \sum_{j=1}^M \lambda_j d_j$.
- Problem 2: There exist $h \in \mathbb{R}^N$ such that $h^{\top}d_0 < 0$ and $h^{\top}d_j \geq 0$ for j = 1, ..., M.

We now investegate this system for any possible arbitrage portfolios. However first we acknowledge that there exist a nominal price system S_t and a normalised price system Z_t . The latter we define as the nominal pricess under the numeraire S_t^1 that is

$$Z_t = \begin{bmatrix} S_t^1/S_t^1 \\ S_t^2/S_t^1 \\ \vdots \\ S_t^N/S_t^1 \end{bmatrix} = \begin{bmatrix} 1 \\ S_t^2/S_t^1 \\ \vdots \\ S_t^N/S_t^1 \end{bmatrix}.$$

The reason for introducing the normalized price system is that we can without much effort translate results in this system to the nominal system and the normalised system is easier to analyze. For this, however, we need af few results.

Lemma 3.3. (Bjork) With notation as above, the following hold.

1. The Z_t value process i related to the S_t value process by

$$V_t^{h,Z} = hZ_t = \frac{1}{S_t^1} V_t^h.$$

- 2. A portfolio is an arbitrage in the S_t system if and only if there is an arbitrage in the Z_t system. 3. In the Z_t price system, the numeraie asset Z^1 has unit constant prices i.e. $Z_t^1 = 1$ for all $t \geq 0$.

One of the reason that the normalised system is attractable is that the numeraire asset is constant i.e. risk free in the normalised system. Let us formulate our first main result.

Proposition 3.4. (Bjork) The market is arbitrage free if and only if there exists strictly positive real numbers $q_1, ..., q_M \ge 0$ with $q_1 + \cdots + q_M = 1$ (eq. 3.2) (probability vector) such that the following vector equality holds

$$\begin{bmatrix} Z_0^1 \\ \vdots \\ Z_N^1 \end{bmatrix} = \begin{bmatrix} Z_1^1(\omega_1) \\ \vdots \\ Z_1^N(\omega_1) \end{bmatrix} q_1 + \dots + \begin{bmatrix} Z_1^1(\omega_M) \\ \vdots \\ Z_1^N(\omega_M) \end{bmatrix} q_M. \tag{3.3}$$

5.1.3.3 Martingale Measures

Definition 3.5. (Bjork) Given the objective probability measure P on (Ω, \mathcal{F}, P) , we say that another probability measure Q defined on Ω is **equivalent** to P if

$$\forall A \in \mathcal{F} : P(A) = 0 \iff Q(A) = 0,$$

or equivalently

$$\forall A \in \mathcal{F} : P(A) = 1 \iff Q(A) = 1.$$

Definition 3.7. (Bjork) Consider the market model above and set S^1 as the numeraire asset. We say that a probability measure Q defined on Ω is a martingale measure if it satisfies the following conditions:

- 1. Q is equivalent to P, i.e. $Q \sim P$.
- 2. For every i = 1, ..., N, the normalized asset price process

$$Z_t^i = \frac{S_t^i}{S_t^1},$$

is martingale under the measure Q.

Theorem 3.8. (Bjork) (First Fundamental Theorem) Given a fixed numeraire, ther market is free of arbitrage possibilities if and only if there exists a martingale measure Q.

By assuming that the numeraire asset is risk free (i.e. does not depend on ω) then by scaling we can derive the short interest rate as

$$1 + R = \frac{S_1^1}{S_0^1}.$$

With this in mind we can formulate theorem 3.8 in its more widely used form.

Theorem 3.9. (Bjork) (First Fundamental Theorem) Assume that there exist a risk free asset, and denote the corresponding risk free interest rate by R. Then the market is arbitrage free if and only if there exist a measure $Q \sim P$ such that

$$S_0^i = \frac{1}{1+R} E^Q[S_1^i], \quad \text{for all } i = 1, ..., N.$$
 (3.9)

5.1.3.4 Martingale Pricing

Moving forward we will assume that there exist a risk free asset and we will denote it by B_t ($B_t = S_t^1/S_0^1$).

Definition 3.10. (Bjork) A contingent claim is any random variable X, defined on the sample space Ω .

To ensure no arbitrage in the extended market containing the N assets and the contingent claim we can apply the first fundamental pricing theorem on the extended market.

Proposition 3.11. (Bjork) Consider a given claim X. In order to avoid arbitrage, X must then be priced according to the formula

$$\Pi_0[X] = \frac{1}{1+R} E^Q[X],\tag{3.10}$$

where Q is a martingale measure for the underlying market $(\Pi, S^1, ..., S^N)$.

5.1.3.5 Completeness

Given that a market is arbitrage-free we may run into a uniqueness issue when determining the price of a contingent claim. If a martingale measure exist we will very much like it to be unique as this will ensure that the price from the risk neutral valuation formula is unique. To this we need the market to be complete.

Definition 3.12. (Bjork) Consider a contingent claim X. If there exists a portfolio h, based on the underlying assets, such that

$$V_1^h = X$$
, with probability 1 (3.11)

i.e.

$$V_1^h(\omega_j) = X(\omega_j), \ j = 1, ..., M,$$
 (3.12)

then we say that X is **replicated**, or **hedged** by h. Such a portfolio h is called a replicating, or hedging portfolio. If every contingent claim can be replicated, we say that the market is **complete**.

We can now formulate a proposition on when the market is complete in terms of the matrix D.

Proposition 3.13. (Bjork) The market is complete if and only if the rows of the matrix D span \mathbb{R}^M , i.e. if and only if D has rank M.

Now we formulate the second fundamental pricing theorem in terms of the martingale measure Q.

Proposition 3.14. (Bjork) (Second Fundamental Theorem) Assume that the model is arbitrage free i.e. Q exist. Then the market is unique if and only if the martingale measure is unique.

5.1.3.6 Stochastic Discount Factors

Definition 3.16. (Bjork) The random variable L on Ω is defined by

$$L(\omega_i) = \frac{q_i}{p_i}, \qquad i=1,...,M.$$

Definition 3.17. (Bjork) Assume the absence of arbitrage, and fix a martingale measure Q. With notation as above, the **stochastic discount factor** (or "state price deflator") is the random variable Λ on Ω by

$$\mathbf{M}(\omega) = \frac{1}{1+R} \cdot L(\omega). \tag{3.19}$$

Proposition 3.18. (Bjork) The arbitrage free price of any claim X is given by the formula

$$\Pi_0[X] = E^P[\mathbf{M} \cdot X] \tag{3.20}$$

where M is a stochastic discount factor.

5.2 Self-financing portfolios

We move forward in this chapter by first defining a self-financing portfolio in discrete time and then by letting the step length tend to zero obtain the continuous time analogue.

5.2.1 Discrete time SF portfolio

We consider N different adapted price processes $S^1, ..., S^N$. We use the following definition.

Definition 6.1. (Bjork) We use the following definitions.

- S_n^i is th price of asset i at time n,
- h_n^i is the number of units of asset i held during [n, n+1), that is bought at time n,
- d_n^i is the dividends from asset i in the time-interval [n-1,n), that is received at time n,
- h_n is the portfolio $(h_n^1,...,h_n^N)$ held during [n,n+1),
- ullet c_n is the consumption i.e. withdrawel at time n (negative being deposits/saving),
- V_n is the value of the portfolio just before time n i.e. of the portfolio h_{n-1} at time n.

We are now ready to define the self-financing portfolio

Definition 6.2. (Bjork) A self-financing portfolio supporting the consumption stream c is a portfolio adhering to the budget constraint given as

$$h_{n+1}S_{n+1} + c_{n+1} = h_nS_{n+1} + h_nd_{n+1}$$

The interpretation being, that we may only use funds obtained from selling the old portfolio h_n and recieved in dividends to buy the new portfolio h_{n+1} and consume the amount c_{n+1} .

Before studying the self-financing portfolio we define the operator Δ (in definition 6.3) as the increment $\Delta x_n = x_{n+1} - x_n$ of a countable sequence $(x_n)_{n \in \mathbb{N}_0}$. Notice that we define the increment forward so the increment n is the increment over the time period [n, n+1) with the first increment being [0, 1). Using this notation we can derive the lemma below.

Lemma 6.4. (Bjork) For any pair of sequences of real numbers $(x_n)_{n\in\mathbb{N}_0}$ and $(y_n)_{n\in\mathbb{N}_0}$ we have the relations

$$\Delta(xy)_n = x_n \Delta y_n + y_{n+1} \Delta x_n, \tag{6.5}$$

$$\Delta(xy)_n = y_n \Delta x_n + x_{n+1} \Delta y_n, \tag{6.6}$$

$$\Delta(xy)_n = x_n \Delta y_n + y_n \Delta x_n + \Delta x_n \Delta y_n. \tag{6.7}$$

This is also valid if the sequences are N-dimensional, where we interpret the products above as scalar products (xy^{\top}) .

Using these definitions and the lemma above we see that the dynamics of the self-financing portfolio is given below.

Proposition 6.6. (Bjork) The dynamics of any self-financing portfolio supporting the consumption stream c are given by

$$\Delta V_n = h_n \Delta S_n + h_n d_{n+1} - c_{n+1}, \tag{6.11}$$

or, in more detail

$$\Delta V_n = \sum_{i=1}^{N} h_n^i (\Delta S_n^i + d_{n+1}^i) - c_{n+1}. \tag{6.12}$$

We may rewrite the dividends as accumulating dividends $D_n^i = \sum_{k=1}^n d_k^i$ and see that $d_{n+1}^i = \Delta D_n^i$ and so the above condition is equivalent with.

Proposition 6.8. (Bjork) The dynamics of any self-financing portfolio supporting the consumption stream c are given by

$$\Delta V_n = h_n \Delta S_n + h_n \Delta D_n - c_{n+1}, \tag{6.15}$$

or, in more detail

$$\Delta V_n = \sum_{i=1}^{N} h_n^i (\Delta S_n^i + \Delta D_n^i) - c_{n+1}. \tag{6.16}$$

5.2.2 Continuous time SF portfolio

Formulating the dynamics of the self-financing portfolio in continuous time is easy work given the discrete setup above. However since we now are in continuous time we will change the n with a t and cosider the behavour $V_{t+dt} - V_t$ as we let $dt \to 0$. First we formulate some basic notation.

Definition 6.9. (Bjork) We use the following definitions.

- S_t^i is th price of asset i at time t,
- h_t^i is the number of units of asset i held at time t,
- D_t^i is the cumulative dividend process for asset i,
- h_t is the portfolio $(h_t^1, ..., h_t^N)$ held at time t,
- ullet c_t is the consumption rate at time n (negative being deposits/saving),
- V_t is the value of the portfolio at time t i.e. of the portfolio h_t at time t.

Given these definitions we may define a portfolio strategy that is self-financing.

Definition 6.10. (Bjork) Let S be and adapted N-dimensional price process. We define the following

- 1. A portfolio strategy is any adapted N-dimensional process h.
- 2. The value process V^h corresponding to the portfolio h is given by

$$V_t^h = \sum_{i=1}^N h_t^i S_t^i. (6.17)$$

- 3. A consumption process is any adapted one-dimensional process c.
- 4. A portfolio-consumption pair (h,c) is called **self-financing** if the value process V^h satisfies the condition

$$dV_t^h = \sum_{i=1}^N h_t^i (dS_t^i + dD_t^i) - c_t dt,$$
(6.18)

i.e. if

$$dV_t^h = h_t dS_t + h_t dD_t - c_t dt.$$

5. The gain process G is defined by

$$G_t = S_t + D_t \tag{6.19}$$

so we can write the self-financing condition as

$$dV_t = h_t \ dG_t - c_t \ dt. \tag{6.20}$$

6. The portfolio h is said to be Markovian if it is of the form

$$h_t = h(t, S_t),$$

for some function $h: \mathbb{R}_+ \times \mathbb{R}^N \to \mathbb{R}^N$.

5.2.3 Portfolio weights

Definition 6.11. (Bjork) For a given portfolio h the corresponding relative portfolio or portfolio weights w are defined by

$$w_t^i = \frac{h_t^i S_t^i}{V_t^h}, \ i = 1, ..., N, \tag{6.21}$$

so, in particular, we have $\sum_{i=1}^{N} w_i = 1$.

Lemma 6.12. (Bjork) A portfolio-consumption pair (h, c) is self-financing if and only if

$$dV_t^h = V_t^h \sum_{i=1}^N w_t^i \frac{dS_t^i + dD_t^i}{S_t^i} - c_t dt$$
 (6.22)

or equivalently with the absolute weights

$$dV^h_t = \sum_{i=1}^N h^i_t (dS^i_t + dD^i_t) - c_t \ dt.$$

Lemma 6.13. (Bjork) Consider the case with no dividends. Let c be a consumption process, and assume that there exist a scalar process Z and a vector process $q = (q^1, ..., q^N)$ such that

$$dZ_t = Z_t \sum_{i=1}^{N} q_t^i \frac{dS_t^i}{S_t^i} - c_t dt, (6.23)$$

and $\sum_{i=1}^{N} qq^{i} = 1$ (eq. 6.24). Now define a portfolio h by

$$h_t^i = \frac{q_t^i Z_t}{S_t^i}. (6.25)$$

Then the value process V^h is given by $V^h = Z$, the pair (h, c) is self-financing, and the corresponding relative portfolio w is given by w = q.

5.3 Black-Scholes PDE

The Black-Scholes model revolves arround SDE's as seen above. In this model we have two assets a risk free asset B and a stochastic priced asset S. We therefore start by defining what we mean by a quote-on-quote risk free asset.

Definition 7.1. (Bjork) The price process B is the price of a risk free asset if it has the dynamics

$$dB_t = r_t B_t \ dt, \tag{7.1}$$

where r is any \mathcal{F}_t adapted process.

We see from this definition that the meaning of "risk free" is the property, that B is priced locally deterministic in the sence that r is adapted and therefore known at time t and we therefore know the yield on a short term basis. This is also why we may call r the **short interest rate**. Given the dynamics above, we know that B in fact is represented by the process

$$B_t = B_0 e^{\int_0^t r_s \ ds}$$

for some B_0 initial value. We will moving forward assume that $B_0 = 1$. The stochastic asset S has dynamics.

$$dS_t = \mu(t, S_t) dt + \sigma(t, S_t) dW_t, \tag{7.2}$$

where as usual μ and σ are deterministic functions and W_t is a standard Brownian motion. Note that the risk free asset has a similarly process with $\sigma = 0$. We may now include this in the definition of the Black-Scholes model.

Definition 7.2. (Bjork) The Black-Scholes model consists of two assets with dynamics given by

$$dB_t = rB_t dt, (7.3)$$

$$dS_t = \mu S_t \ dt + \sigma S_t \ dW_t, \tag{7.4}$$

where r, μ, σ are deterministic constants.

Definition 7.3. (Bjork) A zero coupon bond with maturity T (henceforth "T-bond") is an asset which pays the holder the face value 1 dollar at time T. The price at time n of a T-bond is denoted by p(n,T).

Definition 7.4. The (possible stochastic) discrete **short rate** r_n , for the period [n, n+1], is defined as

$$p(n, n+1) = \frac{1}{1+r_n}. (7.6)$$

From this short rate we may derive the dynamics of the bank account recieving zero-coupon rates for each distinct time interval.

Definition 7.5. (Bjork) The dynamics of the bank account are given by

$$\Delta B_n = r_n B_n. \tag{7.7}$$

5.3.1 Contingent Claims and Arbitrage

Definition 7.6. (Bjork) A European call option with exercise price (or strike price) K and time of maturity (exercise date) T on the underlying asset S is a contract defined by the following clauses:

• The holder of the option has, at time T, the right to buy one share of the underlying stock at the price K dollars from the underwriter of the option.

- The holder of the option is in no way obliged to buy the underlying stock.
- The right to buy the underlying stock at the price K can only be exercised at the precise time T.

Obviously, we also have the **european put** option which gives the owner the right to sell an asset at price K at time T. Let os formally define a contingent claim.

Definition 7.7. Consider a financial market with vector price process S. A contingent claim with date of maturity T, also called a T-claim, is any random variable $X \in \mathcal{F}_T^S$. A contingent claim X is called a simple claim if it is of the form $X = \Phi(S_t)$. The function Φ is called the contract function.

Definition 7.8. (Bjork) An arbitrage possibility on a financial market is a self-financed portfolio h such that

$$V^h(0) = 0, (7.13)$$

$$P(V_T^h \ge 0) = 1, (7-14)$$

$$P(V_T^h > 0) > 0. (7.15)$$

We say that the market is arbitrage free if there are no arbitrage possibilities.

Definition 7.9. (Bjork) Suppose that there exists a self-financing portfolio h, such that the value process V^h has the dynamics

$$dV_t^h = k_t V_t^h dt, (7.16)$$

where k is an adapted process. Then it must hold that $k_t = r_t$ for all t, ore there exists an arbitrage possibility.

Theorem 7.10. (Bjork) (Black-Scholes equation) Assume that the market is specified by the equations

$$dB_t = rB_t dt, (7.18)$$

$$dS_t = \mu(t, S_t)S_t dt + \sigma(t, S_t)S_t dW_t, \tag{7.19}$$

and that we want to price a contingent claim of the form $\mathcal{X} = \Phi(S_t)$ (eq. 7.20). Then the only pricing function of the form $\Pi_t[\Phi(S_t)] = F(t, S_t)$ (eq. 7.21) which is consistent with the absence of arbitrage in the market $[B_t, S_t, \Pi_t]$ is when F is the solution of the following boundary value problem in the domain $[0, T] \times \mathbb{R}_+$:

$$F_t(t,s) + rsF_s(t,s) + \frac{1}{2}s^2\sigma^2(t,s)F_{ss}(t,s) - rF(t,s) = 0,$$

$$F(T,s) = \Phi(s).$$

5.3.2 Risk Neutral Valuation

Theorem 7.11. (Bjork) (Risk Neutral Valuation) The arbitrage free price of the claim $\Phi(S_t)$ is given by $\Pi_t[\Phi] = F(t, S_t)$, where F is given by the formula

$$F(t,s) = e^{-r(T-t)} E_{t,s}^{Q}[\Phi(S_T)], \tag{7.43}$$

where the Q-dynamics of S are those of

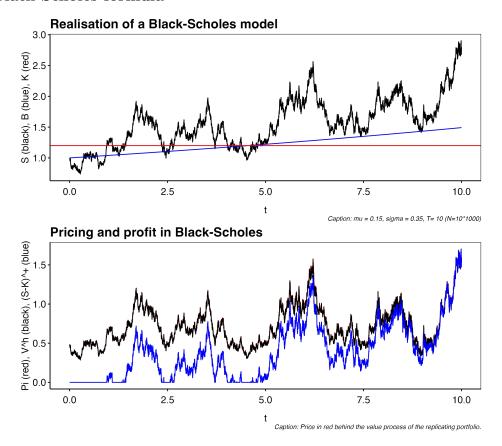
$$dS_t = rS_t dt + S_t \sigma(t, S_t) dW_t^Q. (7.42)$$

Property 7.12. (Bjork) (The Martingale Property) In the Black-Scholes model, the price process Π_t for every traded asset, be it the underlying or derivate asset, has the property the the normalized price process

$$Z_t = \frac{\Pi_t}{B_t},$$

(including S_t/B_t) is a martingale under the measure Q.

5.3.3 Black-Scholes formula



This chapter will center on deriving the famous Black-Scholes formula. We start by laying out the assumptions of the model. We have a market consiting of two assets: a stochastic prices asset S and a risk free asset B. The prices processes have dynamics:

$$dS_t = \mu S_t \ dt + \sigma S_t \ dW_t, \tag{7.45}$$

$$dB_t = rB_t dt, (7.44)$$

where $S_0 = s$ and $B_0 = 1$ (by assumption). Now from Feynmann-Kac and the definition of arbitrage we know that a simple claim $\Phi(S_t)$ has the arbitrage free price given by the risk neutral valueation formula.

$$F(t,s) = e^{-r(T-t)} E_{t,s}^{Q}[\Phi(S_T)], \tag{7.46}$$

where Q is a probability measure, namely a Martingale measure, such that the dynamics of S under this measure is

$$dS_t = rS_t dt + \sigma S_t dW_t^Q, (7.47)$$

with W_t^Q being a Brownian motion wrt. to the probability measure Q (not P). The above still has the initial condition $S_0 = s$. Given these assumptions we may formulate the Black-Scholes formula.

Theorem 7.13. (Bjork) (Black-Scholes formula) The price of the european call option with strikeprice K and maturity T (contract function $\Phi(S_t) = (S_t - K)^+$) takes the form $\Pi_t = F(t, s)$, where

$$F(t,s) = sN(d_1(t,s)) - e^{-r(T-t)}KN(d_2(t,s)), \tag{7.52} \label{eq:fitting}$$

where N is the distribution-function for an $\mathcal{N}(0,1)$ -distributed random variable and

$$d_1(t,s) = \frac{1}{\sigma\sqrt{T-t}} \left(\log\left(\frac{s}{K}\right) + \left(r + \frac{1}{2}\sigma^2\right) (T-t) \right),\tag{7.53}$$

$$d_2(t,s) = d_1(t,s) - \sigma\sqrt{T-t}. (7.54)$$

Proof.

We let the market be given in terms of the price processes S and B with dynamics.

$$dS_t = \mu S_t dt + \sigma S_t dW_t,$$

$$dB_t = rB_t dt,$$

with $B_t = 1$ and $S_t = s$. We assume that μ, σ, r are deterministic real numbers. Consider the contingent claim

$$\Phi(S_t) = \left(S_t - K\right)^+,\,$$

that is the European call option. Let Q be a martingale measure such that the dynamics of S may be written as

$$dS_t = rS_t \ dt + \sigma S_t \ dW_t^Q,$$

then S_t is clearly a GBM wrt. the measure Q. Therefore we know the solution given in terms of the increment of the Brownian motion W^Q as follows

$$S_{u} = s \cdot \exp \left\{ \left(r - \frac{1}{2}\sigma^{2}\right)\left(u - t\right) + \sigma\left(W_{u}^{Q} - W_{t}^{Q}\right)\right\},$$

for some initial condition $S_t = s$. From theorem 7.10 we know that the only pricing function which takes the form

$$\Pi_t[\Phi(S_T)] = F(t, S_t),$$

can only be consistent with the absence of arbitrage if F is the solution the the boundary value problem

$$F_t(t,s) + rsF_s(t,s) + \frac{1}{2}s^2\sigma^2F_{ss}(t,s) - rF(t,s) = 0,$$

$$F(T,s) = \Phi(s).$$

From Feynmann-Kac we then know that the stochastic representation of such a solution take the form

$$F(t,s) = e^{-r(T-t)} E_{t,s}^Q [\Phi(S_T)].$$

Here the superscript refers to taking mean value with respect to the measure Q. This gives the solution to the pricing function

$$F(t,s) = e^{-r(T-t)} \, \int \Phi(S_T) \; dQ.$$

Under the measure Q we have that for $u \geq t$:

$$Z_u = \log(S_u/s) \sim \mathcal{N}\left(\left(r - \frac{1}{2}\sigma^2\right)(u - t), \sigma\sqrt{u - t}\right)$$

Hence we may set u = T and observe that

$$\begin{split} F(t,s) &= e^{-r(T-t)} \int_{-\infty}^{\infty} \Phi(se^z) f(z) \ dz \\ &= e^{-r(T-t)} \int_{-\infty}^{\infty} (se^z - K)^+ f(z) \ dz \\ &= e^{-r(T-t)} \int_{\log\left(\frac{K}{s}\right)}^{\infty} (se^z - K) f(z) \ dz \\ &= e^{-r(T-t)} \left(s \int_{\log\left(\frac{K}{s}\right)}^{\infty} e^z f(z) \ dz - K \int_{\log\left(\frac{K}{s}\right)}^{\infty} f(z) \ dz \right), \end{split}$$

where we used that f is the distribution function of a normal distributed random variable with mean $(r - \sigma^2/2)(T - t)$ and variance $\sigma\sqrt{T - t}$ and that

$$(se^z - K)^+ \ge 0 \iff se^z \ge K \iff z \ge \log\left(\frac{K}{s}\right)$$

Using that the MGF of a $X \sim \mathcal{N}(\alpha, \beta^2)$ variable is

$$E[e^{tX}] = e^{\alpha t + \frac{1}{2}\beta^2 t^2}.$$

and the shorthand N(t) for the distribution function of the standard normal distribution, we have

$$\begin{split} F(t,s) &= e^{-r(T-t)} \left(sE\left[e^{Z_T} \mathbf{1}_{Z_T \geq \log\left(\frac{K}{s}\right)} \right] - KP\left(Z_T \geq \log\left(\frac{K}{s}\right) \right) \right) \\ &= e^{-r(T-t)} s \exp\left\{ \left(r - \frac{1}{2}\sigma^2 \right) (T-t) + \frac{1}{2}\sigma^2 (T-t) \right\} E\left[\mathbf{1}_{Z_T \geq \log\left(\frac{K}{s}\right)} \right] \\ &- e^{-r(T-t)} KP\left(X \geq \frac{1}{\sigma\sqrt{T-t}} \left(\log\left(\frac{K}{s}\right) - (r-\sigma^2/2)(T-t) \right) \right) \\ &= sE\left[\mathbf{1}_{Z_T \geq \log\left(\frac{K}{s}\right)} \right] - e^{-r(T-t)} KP\left(X \leq \frac{1}{\sigma\sqrt{T-t}} \left(\log\left(\frac{s}{K}\right) + (r-\sigma^2/2)(T-t) \right) \right) \\ &= sN(d_1(s,t)) - e^{-r(T-t)} KN\left(d_2(s,t) \right), \end{split}$$

as desired. \blacksquare

5.4 Completeness and Hedging

We derived the pricing function of the european call option above and introduced the theory around boundary value problems and Feynmann-Kac solution to the partial differential stochastic equation. Now we want to see if a portfolio exists such that it gives the payout $\Phi(S_T)$ with probability one.

In order to do this, we return to the concept of hedge and replication.

Definition 8.1. (Bjork) We say that a T-claim \mathcal{X} can be **replicated**, alternatively the it is **reachable** or **hedgeable**, if there exists a self-financing portfolio h such that

$$V_T^h = \mathcal{X}, \ P - \text{a.s.}$$
 (8.1)

In this case we say that h is a **hedge** against \mathcal{X} . Alternatively, h is called a **replicating** or **hedging** portfolio. If every contingent claim is reachable we say that the market is **complete**.

If we can find a portfolio h that reaches \mathcal{X} in value over the time period [t,T] it must mean, that holding the portfolio is equivalent with holding the contract itself. We therefore have the natural assumption that the price process must satisfie $\Pi_t[\mathcal{X}] = V_t^h$ for all $t \geq 0$. How this relates to the absence of arbitrage is given below.

Proposition 8.2. (Bjork) Suppose \mathcal{X} is hedged using the portfolio h. Then the only price process $\Pi_t[\mathcal{X}]$ which is consistent with no arbitrage is given by $\Pi_t[\mathcal{X}] = V_t^h$. Furthermore, if \mathcal{X} can be hedged by both h and g then $V_t^g = V_t^h$ for all t with probability one.

5.4.1 Completeness in Black-Scholes

The Black-Scholes model will be investigated in the following. We start by stating the important theorem.

Theorem 8.3. (Bjork) Consider the Black-Scholes model given by

$$dS_t = \mu(t, S_t)S_t dt + \sigma(t, S_t)S_t dW_t, \tag{8.2}$$

$$dB_t = rB_t dt, (8.3)$$

The model above is complete.

The following lemma gives us replicability of a **simple** claim (which we will restrict ud to).

Lemma 8.4. (Bjork) Suppose that there exist an adapted process V and an adapted process $w = [w^B, w^S]$ with $w_t^B + w_t^S = 1$ (eq. 8.4) for all $t \ge 0$, such that

$$dV_{t} = V_{t}(w_{t}^{B}r + w_{t}^{S}\mu(t, S_{t})) dt + V_{t}w_{t}^{S}\sigma(t, S_{t}) dW_{t},$$
(8.5)

$$V_t = \Phi(S_t). \tag{8.5}$$

Then the claim $\mathcal{X} = \Phi(S_t)$ can be replicated using w as the relative portfolio. The corresponding value process is given by the process V and the absolute portfolio h is given by

$$h_t^B = \frac{w_t^B V_t}{B_t},\tag{8.6}$$

$$h_t^S = \frac{w_t^S V_t}{S_t}. (8.7)$$

Doing some heuristics we come up with some clever weights, which turns on to adhere to the boundary value problem formulated in the Black-Scholes equation. Given that the weights gives rise to the desired value process, we have successfully found the portfolio weight (see lemma above).

Theorem 8.5. (Bjork) Consider the Black-Scholes model given in (8.3)-(8.4), and a simple contingent claim $\mathcal{X} = \Phi(S_t)$. Define F as the solution to the boundary value problem

$$F_t(t,s) + rsF_s(t,s) + \frac{1}{2}s^2\sigma^2F_{ss}(t,s) - rF(t,s) = 0, \tag{8.17}$$

$$F(T,s) = \Phi(s). \tag{8.17}$$

Then X can be replicated by the relative portfolio

$$w_t^B = \frac{F(t, S_t) - S_t F_s(t, S_t)}{F(t, S_t)},$$
(8.18)

$$w_t^S = \frac{S_t F_s(t, S_t)}{F(t, S_t)}. (8.19)$$

The corresponding absolute portfolio is given by

$$h_t^B = \frac{F(t, S_t) - S_t F_s(t, S_t)}{B_t},$$
(8.20)

$$h_t^S = F_s(t, S_t), \tag{8.21}$$

and the value process V^h is given by

$$V_t^h = F(t, S_t). (8.22)$$

Proposition 8.6. (Bjork) Consider the Black-Scholes model given in (8.3)-(8.4), and a contingent claim on the form $\mathcal{X} = \Phi(S_T, Z_T)$ (eq. 8.29). We define the process Z_t as

$$Z_{t} = \int_{0}^{t} g(u, S_{u}) \ du, \tag{8.30}$$

for some choice of the deterministic function g. Then \mathcal{X} can be replicated using a relative portfolio given by

$$w_{t}^{B} = \frac{F(t, S_{t}, Z_{t}) - S_{t}F_{s}(t, S_{t}, Z_{t})}{F(t, S_{t}, Z_{t})}, \tag{8.31} \label{eq:8.31}$$

$$w_t^S = \frac{S_t F_s(t, S_t, Z_t)}{F(t, S_t, Z_t)}. (8.32)$$

where F is the solution to the boundary value problem

$$F_t(t,s,z) + rsF_s(t,s,z) + \frac{1}{2}s^2\sigma^2F_{ss}(t,s,z) - rF(t,s,z) = 0, \tag{8.33}$$

$$F(T, s, z) = \Phi(s, z).$$
 (8.33)

The corresponding value process is given by $V_t = F(t, S_t, Z_t)$, and F has the stochastic representation

$$F(t,s,z) = e^{-r(T-t)} E_{t,s,z}^{Q} [\Phi(S_T, Z_T)], \tag{8.34}$$

where the Q-dynamics are given by

$$dS_u = rS_u \ du + S_u \sigma(u, S_u) \ dW_u^Q, \tag{8.35}$$

$$S_t = s, (8.36)$$

$$dZ_u = g(u, S_u) \ du, \tag{8.37}$$

$$Z_t = z. (8.38)$$

5.4.2 Absence of Arbitrage

In general we have conflicting forces when evaluating when a certain market is arbitrage free and/or complete. We have in simple terms the non-rigorous theorem below.

Meta-theorem 8.3.1. (Bjork) Let N denote the number of underlying traded assets in the model excluding the risk free asset, and let R denote the number of random sources driving the price system. Generically we then have the following statements.

- The model is arbitrage free if and only if $N \leq R$.
- The model is complete if and only if $N \geq R$.
- The model is arbitrage free and complete if and only if N = R.

5.4.3 Incomplete Markets

We assume a market with a risk free asset and one risky assets with dynamics

$$dX_t = \mu(t, X_t) \ dt + \sigma(t, X_t) \ dW_t. \tag{9.1} \label{eq:9.1}$$

We want to find a unique price of a derivative on a functional form of the risky asset. We assume that we cannot invest in the asset representing the process X_t and so we can solely write contracts based on the observation X_T . The problem here is that we can only short or long the risk free asset and so no derivative is replicable.

The way we solve this problem is by having the market set the price of risk and universally price derivatives based on this given price process. We then have the assumptions

Assumption 9.2.1 We have the market given with the only investable asset B with dynamics

$$dB_t = rB_t dt. (9.2)$$

We furthermore, have an empirically observable stochastic process X which is **not** the price process of any traded asset. The P-dynamics of X is given by

$$dX_t = \mu(t, X_t) dt + \sigma(t, X_t) dW_t$$

Assumption 9.2.2 There is a liquid market for every contingent claim.

Assumption 9.2.3 We assume that

- There is a liquid, frictionless market for each of the contingent claims \mathcal{Y} and \mathcal{Z} .
- The market prices of the claims are of the form

$$\Pi_t[\mathcal{Y}] = F(t, X_t),$$

$$\Pi_t[\mathcal{Z}] = G(t, X_t),$$

where F and G are smooth real valued function.

From Ito's formula we have the dynamics

$$dF = \mu_F F \ dt + \sigma_F F \ dW, \tag{9.4}$$

$$dG = \mu_G G \ dt + \sigma_G G \ dW. \tag{9.5}$$

Where the processes μ_F and σ_F are given by

$$\mu_F = \frac{F_t + \mu F_x + \frac{1}{2}\sigma^2 F_{xx}}{F},$$

$$\sigma_F = \frac{\sigma F_x}{F}.$$

By forming a portfolio of the two contracts we lead to the relation.

$$\frac{\mu_F-r}{\sigma_F} = \frac{\mu_G-r}{\sigma_G}.$$

This gives the important insight.

Proposition 9.1. (Bjork) Assume that the market for derivatives is free of arbitrage. Then there exists a universal process $\lambda(t, X_t)$ such that, with probability one, and for all t, we have

$$\frac{\mu_F(t,X_t)-r}{\sigma_F(t,X_t)}=\mu(t,X_t), \tag{9.7}$$

regardless of the specific choice of the derivative F.

Proposition 9.2. (Bjork) Assume absence of arbitrage, the pricing function F(t,x) of the T-claim $\Phi(X_T)$ solves the following boundary value problem.

$$F_t(t,x) + \mathcal{A}F(t,x) - rF(t,x) = 0, \qquad (t,x) \in (0,T) \times \mathbb{R}, \qquad (9.8)$$

$$F(T,x) = \Phi(x), \qquad x \in \mathbb{R}, \qquad (9.9)$$

$$F(T,x) = \Phi(x), \qquad (0,x) \in (0,T) \times \mathbb{R}, \qquad (0,0)$$

where

$$\mathcal{A}F(t,x) = \left\{\mu(t,x) - \lambda(t,x)\sigma(t,x)\right\}F_x(t,x) + \frac{1}{2}\sigma^2(t,x)F_{xx}(t,x).$$

Proposition 9.3. (Bjork) (Risk neutral valuation) Assuming absence of arbitrage, the pricing function F(t,x) of the T-claim $\Phi(X_T)$ is given by the formula

$$F(t,x) = e^{-r(T-t)} E_{t,x}^{Q}[\Phi(X_T)]. \tag{9.11}$$

The dynamics of X under the martingale measure Q are given by

$$dX_t = \left\{ \mu(t,x) - \lambda(t,x) \sigma(t,x) \right\} F_x(t,x) + \sigma(t,x) \ dW_t^Q,$$

where W^Q is a Q-Brownien motion.

5.5 Parity relations

5.5.1 Put-call Parity

The notion of continuous rebalancing the replicating portfolio require leads to problems in the real world. Trading does cost some money (typical in fractions) and so continuous balancing would make the portfolio go to 0 rather quickly. Why? The Brownian motion has unbounded variation and so we would have to sell and buy the portfolio uncountable many time in any interval and the shift in weight is not neglible. Because of this we would like to see which claims we can replicate by buying and holding a combination of assets and derivatives.

Proposition 10.1. (Bjork) Let Φ and Ψ be contract functions for the T-claims $\mathcal{X} = \Phi(S_T)$ and $\mathcal{Y} = \Psi(S_T)$. Then for any real numbers α and β we have the following price relation:

$$\Pi_t[\alpha \Phi + \beta \Psi] = \alpha \Pi_t[\Phi] + \beta \Pi_t[\Psi]. \tag{10.1}$$

If we consider the basic contract functions

$$\Phi_S(x) = x,\tag{10.2}$$

$$\Phi_B(x) = 1,\tag{10.3}$$

$$\Phi_{C,K}(x) = (x - K)^+, \tag{10.4}$$

$$\Phi_{PK}(x) = (K - x)^{+}.$$

That is a contract paying (respectively): the price of one stock, 1 dollar, one european call and one european put both with strike K. It is clear that the following prices are

$$\Pi_t[\Phi_S] = S_t, \tag{10.5}$$

$$\Pi_t[\Phi_B] = e^{-r(T-t)},$$
(10.6)

$$\Pi_t[\Phi_{C,K}] = c(t, S_t; K, T),$$
(10.7)

$$\Pi_t[\Phi_{P,K}] = p(t, S_t; K, T).$$

Where $c(t, s, K, T, r, \sigma)$ and $p(t, s, K, T, r, \sigma)$ are the pricing function of the european call and put option. We see that we can replicate these payouts by: buying the stock today and selling at time T, buying a zero coupon T-bond with face value 1, buying the call and put option.

Then we can by choosing $\alpha, \beta, \gamma_1, ..., \gamma_n$ form a portfolio consisting of α stocks, β T-bonds and γ_i call options with maturity T and strike K_i . The price is then a linear combination given the choice (se proposition 10.1).

The put option is not included in the above portfolio as we have the put-call parity below

Proposition 10.2. (Bjork) (Put-call parity) Consider a European call and a European put, both with strike K and time of maturity T. Then we have the relation.

$$p(t,s) = Ke^{-r(T-t)} + c(t,s) - s. (10.11)$$

In particular the put option can be replicated by a constant portfolio consisting of K zero coupon T-bonds, a European call option and a single short position in the underlying stock.

We now have the pleasing proposition given the class of claims we can reach with the buy-and-hold portfolio with T-bonds, stock and call options

Proposition 10.3. (Bjork) Fix an arbitrary continuous contract function Φ with compact support. Then the corresponding contract can be replicated with arbitrary precision (in sup-norm) using a constant portfolio consisting only of bonds, call options and the underlying stock.

5.5.2 The Greeks

When holding a portfolio we may denote the pricing function by P(t, s). Here we only have one **underlying** asset with price process S_t . We now have two types of risk:

- Price changes in the underlying asset.
- Misspecifications in the model parameters.

These two risk give rise to "the greeks" as defined below.

Definition 10.4. (Bjork) The greeks of a portfolio is given by

$$\Delta = \frac{\partial P}{\partial s}, \ \Gamma = \frac{\partial^2 P}{\partial s^2}, \ \rho = \frac{\partial P}{\partial r}, \ \Theta = \frac{\partial P}{\partial t}, \ \mathcal{V} = \frac{\partial P}{\partial s}.$$

For the call option in particular we have the following derivatives.

Proposition 10.5. (Bjork) The greeks of a portfolio consisting of a single European call option with maturity T and strike price K have the following greeks (φ denoting the density function of a $\mathcal{N}(0,1)$ -variable):

$$\Delta = N(d_1),\tag{10.17}$$

$$\Gamma = \frac{\varphi(d_1)}{s\sigma\sqrt{T-t}},\tag{10.18}$$

$$\rho = K(T - t)e^{-r(T - t)}N(d_2), \tag{10.19}$$

$$\Theta = -\frac{s\varphi(d_1)\sigma}{2\sqrt{T-t}} - rKe^{-r(T-t)}N(d_2), \tag{10.20} \label{eq:equation:equation}$$

$$\mathcal{V} = s\varphi(d_1)\sqrt{T - t}.\tag{10.21}$$

5.6 Fundamental pricing theorem I and II

We start by stating the following theorem.

Theorem 11.1. (Bjork) If at least on of the assets $S^1, ..., S^N$ has diffusion term which is non-zero at all times, and if naive portfolio strategies are admitted, then the model admits arbitrage.

We will go as follows. Derive the fundamental pricing theorem 1 and 2 in a setting with zero interest rate. Then we will extend the result in general by choosing a simple numeraire. We start by defining some basic notation.

Definition 11.2. (Bjork) Define the process h as

$$h = [h^0, h^S] := [h^0, h^1, ..., h^N]$$

We define the following.

• For a process h, its value process V_t^h is defined by

$$V_t^h = h_t^0 \cdot 1 + \sum_{i=1}^N h_t^i S_t^i, \tag{11.3}$$

or in compact form

$$V_t^h = h_t^0 \cdot 1 + h_t^S S_t \tag{11.4}$$

• An adapted process h^S is called **admissible** if there exists a non-negative real number α (which may depend on the choice of h^S) such that

$$\int_0^t h_u^S dS_u \ge -\alpha,\tag{11.5}$$

for all $t \in [0,T]$. A process h, is called an **admissible portfolio** process if h^S is admissible.

An admissible portfolio is said to be **self-financing**, if

$$V_t^h = V_0^h + \int_0^t h_u^S dS_u, (11.6)$$

i.e. if

$$dV_t^h = h_t^S dS_t. (11.7)$$

Lemma 11.3. (Bjork) For any adapted process h^S satisfying the admissibility condition above, and for any real number x, there exists a unique adapted process h^0 , such that:

- The process h defined by $h = [h^0, h^S]$ is self-financing.
- The value process is given by

$$V_t^h = x + \int_0^t h_u^S dS_u. (11.8)$$

In particular, the space \mathcal{K}_0 of portfolio values, reachable at time T by means of a self-financing portfolio with zero initial cost is given by

$$\mathcal{K}_0 = \left\{ \int_0^T h_t^S dS_u : h^S \text{ is admissible} \right\}. \tag{11.9}$$

Definition 11.4. (Bjork) A probability measure Q on \mathcal{F}_T is called equivalent martingale measure for the market model, the numeraire S^0 , and the time interval [0,T], if it has the following properties:

- Q ~ P on \$\mathcal{F}_T\$, so P and Q are equivalent.
 All price processes \$S^0, S^1, ..., S^N\$ are martingales under Q on the time interval [0, T].

An equivalent martingale measure will often be referred to as just "a martingale measure" or as "an EMM". If $Q \sim P$ has the property that $S^0, S^1, ..., S^N$ are local martingales, then Q is called a **local martingale** measure.

Theorem 11.5. (Bjork) (The First Fundamental Theorem) The model is arbitrage free "essentially" if and only if there exists a (local) martingale measure Q.

Definition 11.6. (Bjork) With the notation above, we say that the model admits

• No Arbitrage (NA) if

$$\mathcal{C} \cap L_+^{\infty} = \{0\},\tag{11.21}$$

• No Free Lunch with Vanishing Risk (NFLVR) if

$$\tilde{\mathcal{C}} \cap L^{\infty}_{+} = \{0\},\tag{11.22}$$

where $\tilde{\mathcal{C}}$ denotes the closure of \mathcal{C} in L^{∞} .

Theorem 11.7. (Bjork) (Kreps-Yan Separation Theorem) If C is weak* closed, and if

$$\mathcal{C} \cap L^{\infty}_{+} = \{0\},\$$

then there exists a random variable $L \in L^1$ such that L is P almost surely strictly positive, and

$$E^P[L \cdot X] \le 0,$$

for all $X \in \mathcal{C}$.

Proposition 11.8. (Bjork) If the asset price processes are uniformly bounded, then the condition NFLVR implies that C is weak* closed.

Theorem 11.9. (Bjork) (First Fundamental Theorem) Assume that the asset price process S is bounded. Then there exists an equivalent martingale measure if and only if the model satisfies NFLVR.

Theorem 11.10. (Bjork) (First Fundamental Theorem) Assume that the asset price process S is locally bounded. Then there exists an equivalent martingale measure if and only if the model satisfies NFLVR.

Assumption 11.4.1. (Bjork) We assume that $S_t^0 > 0$ P-a.s. for all $t \ge 0$.

Definition 11.11. (Bjork) The normalized economy (also referred to as the "Z-economy") is defined by the price vector process Z, where

$$Z_t = \frac{S_t}{S_t^0}.$$

Definition 11.12. (Bjork)

• A portfolio stragegy is any adapted (N+1)-dimensional process

$$h_t = [h_t^0, h_t^1, ..., h_t^N].$$

- The S-value process V_t^S corresponding to the portfolio h is $h_t S_t$. The Z-value process V_t^Z corresponding to the portfolio h is $h_t Z_t$.
- A portfolio is said to be admissible if it is admissible as an Z portfolio.
- An admissible portfolio is **S-self-balancing** if

$$dV_t^S = \sum_{i=0}^{N} h_t^i \ dS_t^i \tag{11.26}$$

• An admissible portfolio is **Z-self-balancing** if

$$dV_t^Z = \sum_{i=0}^{N} h_t^i \ dZ_t^i. \tag{11.28}$$

Lemma 11.13. (Bjork) (Invariance Lemma) With assumptions as above, the following hold.

- i. A portfolio h is S-self-financing if and only if it is Z-self-financing.
- ii. The value processes V^S and V^Z are connected by

$$V_t^Z = \frac{1}{S_t^0} \cdot V_t^S.$$

iii. A claim Y is S-replical if and only if the claim

$$\frac{\mathcal{Y}}{S_T^0}$$

is Z-replicable.

iv. The model is S arbitrage free if and only if it is Z arbitrage free.

Theorem 11.14. (Bjork) (The First Fundamental Theorem) Consider the market model $S^0, S^1, ..., S^N$ where we assume that $S_t^0 > 0$, P-a.s. for all $t \ge 0$. Assume furthermore that $S^0, S^1, ..., S^N$ are locally bounded. Then the followin conditions are equivalent:

- The model satisfies NFLVR.
- There exists a measure $Q \sim P$ such that the processes

$$Z^0, Z^1, ..., Z^N,$$

are local martingales under Q.

5.6.1 Completeness

Lemma 11.15. (Bjork) Consider a given T-claim X. Fix a martingale measure Q and assume that the normalized claim X/S_T^0 is integrable. If the Q-martingale M, defined by

$$M_t = E^Q \left[\left. \frac{X}{S_T^0} \right| \mathcal{F}_t \right], \tag{11.34}$$

admits an integral representation of the form

$$M_t = x + \sum_{i=1}^{N} \int_0^t h_s^i \ dZ_s^i, \tag{11.35}$$

then X can be hedged in the S-economy. Furthermore, the replicating portfolio $(h^0, h^1, ..., h^N)$ is given by the above for h^i , i=1,...,N and $h^0_t=M_t-\sum_{i=1}^N h^i_t Z^i_t$.

Theorem 11.16. (Bjork) (Jacod) Let \mathcal{M} denote the convex set of equivalent martingale measures. Then, for any fixed $Q \in \mathcal{M}$, the following statements are equivalent:

• Every Q local martingale M has dynamics of the form

$$dM_t = \sum_{i=1}^N h_s^i \ dZ_s^i.$$

• Q is an extremal point of \mathcal{M} .

Theorem 11.17. (Bjork) (The Second Fundamental Theorem) Assume that the market is arbitrage free and consider a fixed numeraire asset S^0 . Then the market is complete if and only if the martingale measure Q, corresponding to the numeraire S^0 , is unique.

5.6.2 Risk Neutral Valuation Formula

We have the setting of a market consisting of the assets $S^0, ..., S^N$ of N+1 assets. We consider the numeraire S^0 being a risk free asset. We introduce a price of contingent claim X, such that the extended market consisting of the price process of X and the N+1 assets is arbitrage free. Alternatively, we can, equivalently, find a replicating portfolio h such that $V_T^h = X$ with probability one.

Theorem 11.18. (Bjork) (General Pricing Equation) The arbitrage free price process for the T-claim X is given by

$$\Pi_t[X] = S_t^0 E^Q \left[\left. \frac{X}{S_T^0} \right| \mathcal{F}_t \right], \tag{11.41}$$

where Q is the (not necessarily unique) martingale measure for the a priori given market S^0 , S^1 , ..., S^N , with S^0 as the numeraire.

If if we assume that the bank account takes the form

$$S_t^0 = S_0^0 e^{-\int_0^t r(s) \ ds},$$

where r is the short rate, then we have the familier risk neutral valuation formula.

Theorem 11.19. (Bjork) (Risk Neutral Valuation Formula) Assuming the existence of a short rate, the pricing formula takes the form

$$\Pi_t[X] = E^Q \left[e^{-\int_0^T r(s) \ ds} X \middle| \mathcal{F}_t \right], \tag{11.42}$$

where Q is the (not necessarily unique) martingale measure with the bank account as the numeraire.

Definition 11.20. (Bjork) A zero coupon bond with maturity date T, also called a T-bond, is a contract which guarantees the holder one dollar to be paid on the date T. The price at time t of a bond with maturity date T is denoted by p(t,T).

Proposition 11.21. (Bjork) The price of a zero coupon T-bond is given by

$$p(t,T) = E^Q \left[\left. e^{-\int_t^T r(s) \ ds} \right| \mathcal{F}_t \right], \tag{11.43}$$

and in particular we have p(T,T) = 1 for all $T \ge 0$ (eq. 11.44).

5.6.3 Stochastic Discount Factors

Definition 11.22. (Bjork) Assume the existence of a short rate r. For any fixed martingale measure Q, let the likelihood process L be defined by

$$L_t = \frac{dQ}{dP}, \text{ on } \mathcal{F}_t. \tag{11.48}$$

The stochastic discount factor (SDF) process M, corresponding to Q, is defined as

$$\mathbf{M}_t = e^{-\int_0^t r(s) \ ds} L_t \quad \left(= \frac{1}{B_t} \cdot L_t \right). \tag{11.49/50}$$

Proposition 11.23. (Bjork) Assume absence of arbitrage. With notation as above, the following hold:

• For any sufficiently integrable T-claim X, the arbitrage free price is given by

$$\Pi_t[X] = E^P \left[\frac{\mathbf{M}_T}{\mathbf{M}_t} X \mid \mathcal{F}_t \right]. \tag{11.51}$$

- For any arbitrage free asset price process S (derivative or underlying) the process $\mathbf{M}_t S_t$ is a (local) P-martingale.
- The P-dynamics of M are given by

$$d\mathbf{M}_t = -r_t \mathbf{M}_t \ dt + \frac{1}{B_t} \ dL_t. \tag{11.53}$$

5.6.4 Summary

Theorem 11.24. (Bjork) (First Fundamental Theorem) The market model is free of arbitrage if and only if there exists a martingale measure, i.e. a measure $Q \sim P$ such that the processes

$$\frac{S_t^0}{S_t^0}, \frac{S_t^1}{S_t^0}, ..., \frac{S_t^N}{S_t^0}$$

are (local) martingales under Q.

Proposition 11.25. (Bjork) If the numeraire S^0 is the money account, i.e.

$$S_t^0 = e^{\int_0^t r(s) \ ds},$$

where r is the (possibly stochastic) short rate, and if we assume that all processes are Brownian driven, then a measure $Q \sim P$ is a martingale measure if and only if all assets $S^0, S^1, ..., S^N$ have the short rate as their local rates of return, i.e. if the Q-dynamics are of the form

$$dS_t^i = S_t^i r_t dt + S_t^i \sigma_t^i dW_t^Q, (11.54)$$

where W^Q is a (multidimensional) Q-Brownian motion.

Theorem 11.26. (Bjork) (Second Fundamental Theorem) Assuming absence of arbitrage, the market model is complete if and only if the martingale measure Q is unique.

Proposition 11.27. (Bjork)

1. In order to avoid arbitrage, X must be priced according to the formula

$$\Pi_t[X] = S_t^0 E^Q \left[\frac{X}{S_T^0} \middle| \mathcal{F}_t \right], \tag{11.55}$$

where Q is a martingale measure for $[S^0, S^1, ..., S^N]$, with S^0 as the numeraire.

2. In particular, we can choose the bank account B_t , as the numeraire. Then B has the dynamics

$$dB_t = r_t B_t \ dt, \tag{11.56}$$

where r is the (possibly stochastic) short rate process. In this case the pricing formula above reduces to

$$\Pi_t[X] = E^Q \left[e^{-\int_t^T r(s) \ ds} X \mid \mathcal{F}_t \right]. \tag{11.57}$$

3. As a special case, the price of a zero coupon T-bond is given by

$$p(t,T) = E^{Q} \left[e^{-\int_{t}^{T} r(s) \ ds} \ \middle| \ \mathcal{F}_{t} \right]. \tag{11.58}$$

4. Defining the stochastic discount factor \mathbf{M} by $\mathbf{M}_t = B_t^{-1} L_t$ we also have the pricing formula.

$$\Pi_t[X] = E^Q \left[\frac{\mathbf{M}_T}{\mathbf{M}_t} X \mid \mathcal{F}_t \right]. \tag{11.59}$$

CHAPTER 5. CONTINUOUS TIME FINANCE 5.6. FUNDAMENTAL PRICING THEOREM I AND II

5. Different choices of Q will generically give rise to different price processes for a fixed claim X. However, if X is attainable then all choices of Q will produce the same price process, which then is given by

$$\Pi_t[X] = V_t^h, \tag{11.60}$$

where h is the hedging portfolio. Different choices of hedging portfolios (if such exist) will produce the same price process.

 $6. \ In \ particular, for \ every \ replicable \ claim \ X \ it \ holds \ that$

$$V_t^Q = E^Q \left[e^{-\int_t^T r(s) \ ds} X \mid \mathcal{F}_t \right]. \tag{11.61}$$

Mathematics of the martingale approach 5.7

Martingale representation theorem

Theorem 12.1. (Bjork) (Representation of Brownian Functionals) Let W be a d dimensional Brownian motions, and let X be a random variable such that

- $X \in \mathcal{F}_T^W$, $E[|X|] < \infty$.

Then there exist uniquely determined \mathcal{F}_t^W -adapted processes $h^1,...,h^d$, such that X has the representation

$$X = E[X] + \sum_{i=1}^{d} \int_{0}^{T} h_{s}^{i} dW_{s}^{i}.$$
 (12.2)

Under the additional assumption

$$E[X^2] < \infty$$
,

then $h^1, ..., h^d$ are in \mathcal{L}^2 .

Theorem 12.2. (Bjork) (The Martingale Representation Theorem) Let W be a d dimensional Brownian motions, and assume that the filtration \mathbf{F} is defined as

$$\mathcal{F}_t = \mathcal{F}_t^W, \qquad t \in [0, T].$$

Let M be any \mathcal{F}_t -adapted martingale. Then there exist uniquely determined \mathcal{F}_t -adapted processes $h^1, ..., h^d$ such that M has the representation

$$M_t = M_0 + \sum_{i=1}^d \int_0^t h_s^i \ dW_s^i, \qquad t \in [0, T].$$
 (12.9)

If the martingale M is square integrable, then $h^1, ..., h^d$ are in \mathcal{L}^2 .

5.7.2Girsanov theorem

Theorem 12.3. (Bjork) (The Girsanov Theorem) Let W be a d dimensional P-Brownian motion on $(\Omega, \mathcal{F}, P, \mathbf{F})$ and let φ be any d-dimensional adapted column vector process. Choose a fixed T and define the process L on [0,T] by

$$dL_t = \varphi_t^\top L_t \ dW_t, \tag{12.16}$$

$$L_0 = 1, (12.17)$$

i.e.

$$L_t = \exp\left\{\int_0^t \varphi_s^\top \ dW_s - \frac{1}{2} \int_0^t \|\varphi_s\|^2 \ ds\right\}.$$

Assume that

$$E^{P}[L_{T}] = 1, (12.18)$$

and define the new probability measure Q on \mathcal{F}_T by

$$L_T = \frac{dQ}{dP}, \quad on \ \mathcal{F}_T. \tag{12.19}$$

Then

$$dW_t = \varphi \ dt + dW_t^Q, \tag{12.20}$$

where W^Q is a d dimensional Q-Brownian motion or equivalently

$$W_t^Q = W_t - \int_0^t \varphi_s \, ds \tag{12.21}$$

is a standard Q-Brownian motion.

We will often refere to φ as the **Girsanov kernel** of the measure transformation. Furthermore, we have written on component form above and the L dynamics will have the form

$$dL_t = L_t \sum_{i=1}^d \varphi_t^i \ dW_t^i,$$

and L will have the explicit form

$$L_t = \exp\left\{ \sum_{i=1}^d \int_0^t \varphi_s^i \ dW_s^i - \frac{1}{2} \int_0^d \sum_{i=1}^d (\varphi_s^i)^2 \ ds \right\}.$$

The conclusion of the Girsanov Theorem is thwn that we can write

$$dW_t^i = \varphi_t^i \ dt + dW_t^{Q,i},$$

for i=1,...,d where $W_t^{Q,1},...,W_t^{Q,d}$ are independent standard Brownian motions under Q.

Definition 12.4. (Bjork) For any Brownian motion W and any kernel process φ , the **Doleans exponential** process \mathcal{E} is defined by

$$\mathcal{E}(\varphi \bullet W)_t = \exp\left\{ \int_0^t \varphi_s^\top \ dW_s - \frac{1}{2} \int_0^t \|\varphi\|_s^2 \ ds \right\}. \tag{12.24}$$

Lemma 12.5. (Bjork) (The Novikov Condition) Assume that the Girsanov kernel φ is such that

$$E^{P}\left[e^{\frac{1}{2}\int_{0}^{T}\|\varphi_{t}\|^{2}}dt\right]<\infty. \tag{12.27}$$

Then L is a martingale and in particular $E^P[L_T] = 1$.

Theorem 12.6. (Bjork) (The Converse of the Girsanov Theorem) Let W be a d-dimensional standard P-Brownian motion on $(\Omega, \mathcal{F}, P, \mathbf{F})$ and assume that

$$\mathcal{F}_t = \mathcal{F}_t^W, \ \forall t.$$

Assume that there exists a probability measure Q such that $Q \ll P$ on \mathcal{F}_T . Then there exists an adapted process φ such that the likelihood process L has the dynamics

$$dL_t = L_t \varphi_t^\top \ dW_t,$$

$$L_0 = 1.$$

This gives us a recipe to transform dynamics of Ito processes under the measure Q as we may rewrite the dynamics of the Brownian motion. We therefore have for an Ito process X with dynamics

$$dX_t = \mu(t, X_t)X_t dt + \sigma(t, X_t)X_t dW_t,$$

may be transformed under Q as

$$\begin{split} dX_t &= \mu(t, X_t) X_t \ dt + \sigma(t, X_t) X_t \ dW_t \\ &= \mu(t, X_t) X_t \ dt + \sigma(t, X_t) X_t \ (\varphi_t \ dt + dW_t^Q) \\ &= (\mu(t, X_t) + \varphi_t) \ X_t \ dt + \sigma(t, X_t) X_t \ dW_t^Q. \end{split}$$

This may lead us into deducing that

$$\mu(t,X_t) + \sigma(t,X_t)\varphi_t = r_t \iff \varphi_t = \frac{r_t - \mu(t,X_t)}{\sigma(t,X_t)}.$$

We furthermore have the Levy characterisation of a Brownian motion.

Theorem. (Remark FinKont) (Levy Characterisation of Brownian motion) Let X_t be an Ito process with $X_0 = 0$. Then X_t is a Brownian motion if and only if the two processes X_t and $X_t^2 - t$ are continuous martingales.

5.8 Black-Scholes model - martingale approach

We consider the standard Black-Scholes model with a single risk free asset and risky asset with dynamics

$$dS_t = \mu S_t \ dt + \sigma S_t \ dW_t, \tag{13.1}$$

$$dB_t = rB_t dt. (13.2)$$

We want check whether the model is arbitrage free on any time interval [0, T], and find a (perhaps unique) martingale measure such that we may apply the fundamental pricing theorem 1 and 2. From Girsanov this endavour is equivalent with searching for a (perhaps unique) Girsanov kernel φ . We therefore define as usual the likelihood process

$$dL_t = \varphi_t L_t \ dW_t,$$

and setting $dQ = L_T dP$ on \mathcal{F}_T , we know from Girsanov theorem that

$$dW_t = \varphi_t \ dt + dW_t^Q.$$

Inserting into the Black-Scholes model we have

$$dS_t = S_t(\mu + \sigma \varphi_t) dt + \sigma S_t dW_t^Q.$$

We know that for Q to be a martingale measure we know that the local rate of return under Q of S must be the short rate r i.e. we have

$$\mu + \sigma \varphi_t = r \iff \varphi_t = \frac{r - \mu}{\sigma} = -\frac{\mu - r}{\sigma},$$
 (13.3)

and so we see the Girsanov kernel is **constant and deterministic**. The process has the economic interpretation that the Girsanov kernel is the risk premium per unit volatility.

Lemma 13.1. (Bjork) The Girsanov kernel φ is given by

$$\varphi = -\lambda$$

where the market price of risk λ is defined by

$$\lambda = \frac{r - \mu}{\sigma}.$$

We therefore have determined a martingale and so we have the result.

Theorem 13.2. (Bjork) The Black-Scholes model above is arbitrage free.

We could in general have that μ, σ, r are adapted processes. If this is the case we would have to show the Nivokov condition.

Pricing then of any T-claim X then is given by the risk neatral pricing formula

$$\Pi_t[X] = e^{-r(T-t)} E^Q[X \mid \mathcal{F}_t], \tag{13.7}$$

where the Q dynamics of S has local drift r and volatility from the Q-brownian motion W^Q .

Theorem 13.3. (Bjork) The Black-Scholes model above is complete. This also holds for the more general model where r, μ, σ are adapted processes.

Hedging is the possible by considering a T claim with

$$E^Q\left[\frac{X}{B_T}\right] < \infty.$$

Notice the numeraire B_t as in the normalized Z-economy. Consider now the Q-martingale

$$M_t = E^Q \left[\frac{X}{B_T} \middle| \mathcal{F}_t \right], \tag{13.9}$$

and it now follows from lemma 11.15 the the model is complete if we can find a process h_t^1 such that

$$dM_t = h_t^1 \ dZ_t^i. {13.10}$$

In order to prove existance of such a process h^1 we use the Martingale Representation Theorem 12.2, which says that there *exists* a process g_t such that

$$dM_t = g_t \ dW_t^Q. \tag{13.11}$$

We can now combine these two equation by the following Q dynamics

$$dZ_t^1 = Z_t^1 \sigma \ dW_t^Q. \tag{13.12}$$

Hence we have

$$dM_t = h_t^1 Z_t^1 \sigma \ dW_t^Q = g_t \ dW_t^Q \ \Rightarrow \ h_t^1 = \frac{g_t}{Z_t^1 \sigma}.$$

Theorem 13.4. (Bjork) In the Black-Scholes model every T-claim X satisfying

$$E^Q\left[\frac{X}{B_T}\right]<\infty$$

can be replicated. The replicating portfolio is given by

$$h_t^1 = \frac{g_t}{Z_t^1 \sigma},\tag{13.13}$$

$$h_t^0 = M_t - h_t^1 Z_t^1, \tag{13.14}$$

where M is defined by the above and g is defined by above.

If the T-claim is simple that is $X = \Phi(S_T)$ we may solve a boundary value problem with Feynmann-Kac to arrive at the familiar result.

Proposition 13.5. (Bjork) In the Black-Scholes model every T-claim on the form $X = \Phi(S_T)$. Then X can be replicated by the portfolio

$$h_t^0 = \frac{F(t, S_t) - S_t \frac{\partial F}{\partial s}(t, S_t)}{B_t},\tag{13.15}$$

$$h_t^1 = \frac{\partial F}{\partial s}(t, S_t), \tag{13.15}$$

where F solves the Black-Scholes equation

$$\frac{\partial F}{\partial t}(t,s) + rs\frac{\partial F}{\partial s}(t,s) + \frac{1}{2}\sigma^2 s^2 \frac{\partial^2 F}{\partial s^2}(t,s) - rF(t,s) = 0, \tag{13.16}$$

$$F(T,s) = \Phi(s). \tag{13.16}$$

Furthermore the value process for the replicating portfolio is given by

$$V_t = F(t, S_t).$$

5.9 Multidimensional models

We specify the general model by the assumptions below.

Assumption 14.0.1 We assume the following:

- There are n risky assets $S^1, ..., S^n$.
- Under the objective probability measure P, the S-dynamics are given by

$$dS_t^i = \mu_t^i S_t^i \ dt + S_t^i \sum_{j=1}^N \sigma_t^{ij} \ dW_t^j, \tag{14.1}$$

for i = 1, ..., n.

- The coefficients processes μ^i and σ^{ij} above are assumed to be adapted.
- We have a standard risk free asset with price process B with dynamics

$$dB_t = r_t B_t \ dt, \tag{14.2}$$

where the short rate process r is assumed to be an adapted stochastic process.

We can use the representation of μ^i , σ^{ij} and S^i as vectors and matrices on the form.

$$\mu = \begin{bmatrix} \mu^1 \\ \vdots \\ \mu^n \end{bmatrix}, \ \sigma = \begin{bmatrix} \sigma^{1,1} & \cdots & \sigma^{i,N} \\ \vdots & \ddots & \vdots \\ \sigma^{n,1} & \cdots & \sigma^{n,N} \end{bmatrix}, \ D(S) = \begin{bmatrix} S^1 & \cdots & 0 \\ \vdots & \ddots & \vdots \\ 0 & \cdots & S^n \end{bmatrix}.$$

And so we have the model on compact form.

$$dS_t = D(S_t)\mu_t dt + D(S_t)\sigma_t dW_t, (14.3)$$

$$dB_t = r_t B_t \ dt. ag{14.4}$$

Now using Girsanov Theorem we can define the prospective likelihood process L by

$$dL_t = L_t \varphi_t^\top dW_t, \tag{14.5}$$

$$L_0 = 1,$$
 (14.6)

where φ is an adapted N-dimensional process. Then our candidate martingale measure Q is given by $dQ = L_t \ dP$ on \mathcal{F}_t and the Girsanov theorem give the dynamics

$$dW_t = \varphi_t \ dt + dW_t^Q, \tag{14.7}$$

where W^Q is a standard Q-Brownian motion. Inserting into the P-dynamics we obtain.

$$dS_t = D(S_t)[\mu_t + \sigma_t \varphi_t] dt + D(S_t)\sigma_t dW_t^Q.$$
(14.8)

The from (11.54) we know that Q is a martingale measure if and only if the local rate of return (the dt-term) is the short interest rate i.e. if and only if

$$\mu_t + \sigma_t \varphi = \mathbf{r}_t, \tag{14.9}$$

where $\mathbf{r}_t = [r, ..., r]^{\top} \in \mathbb{R}^n$. We then have that

$$\sigma_t \varphi_t = \mathbf{r}_t - \mu_t, \tag{14.11}$$

where we want to solve for φ . Thus we may write a condition for the absence of arbitrage in linear algebra terms.

Proposition 14.1. (Bjork) A necessary condition for absence of arbitrage is that

$$\mathbf{r}_t - \mu_t \in Im[\sigma_t]$$

with probability one for each t. A sufficient condition for absence of arbitrage is that there exists a process φ which solves (14.11) and such that L is a martingale.

Note that it is not enought for φ to solce (14.11). We also need that L is a martingale.

Definition 14.2. (Bjork) A Girsanov kernel φ is said to be admissible if it generates a martingale measure, i.e. it solves (14.11) and L is a true martingale.

Definition 14.3. (Bjork) The model above is said to be generically arbitrage free if it is arbitrage free for every choice of μ .

Proposition 14.4. (Bjork) Disregarding integrability problems the model is generically arbitrage free if and only if, for each $t \leq T$ and P-a.s., the mapping

$$\sigma_t: \mathbb{R}^B \to \mathbb{R}^n$$

is surjective, i.e. if and only if the volatility matrix σ_t has rank n.

Proposition 14.5. (Bjork) Assume that the model i generically arbitrage free and that the filtration F is defined by

$$\mathcal{F}_t = \mathcal{F}_t^W. \tag{14.14}$$

Then, disregarding integrability problems, the model is complete if and only if n = N and the volatility matrix σ_t is invertible P-a.s. for each $t \leq T$.

Assumption 14.3.1. (Bjork) We assume that the model is generically free of arbitrage, i.e. that

$$Im[\sigma_t] = \mathbb{R}^n, \tag{14.16}$$

for all t and with probability one. We also assume that the model is purely Brownian driven, i.e. that $\mathcal{F}_t = \mathcal{F}_t^W$.

Proposition 14.6. (Bjork) Under assumption 14.3.1 the model is complete if and only if

$$Im[\sigma_t^\top] = \mathbb{R}^N. \tag{14.23}$$

If the model is complete then, using the notation of chapter 11, the replicating portfolio $[h^0, h^S]$ is given by

$$\begin{split} h_t^S &= g_t \sigma_t^{-1} D^{-1}(Z_t), \\ h_t^2 &= M_t - h_t Z_t. \end{split} \tag{14.24}$$

$$h_t^2 = M_t - h_t Z_t. (14.25)$$

With M_t defined by

$$M_t = E^Q \left[\frac{\mathcal{X}}{B_T} \middle| \mathcal{F}_t \right]. \tag{14.17}$$

Theorem 14.7. (Bjork) (The Second Fundamental Theorem) Under assumptions 14.3.1 the model is complete if and only if the martingale measure is unique. This is equivalent with the statements: $Ker[\sigma_t] = \{0\}$, $Im[\sigma_t^{\top}] = \mathbb{R}^N$ and σ_t^{-1} exists (i.e. σ_t is invertible).

Pricing of any T-claim $\mathcal X$ is now given by the risk neutral valuation formula

$$\Pi_t[\mathcal{X}] = E^Q \left[e^{-\int_t^T r_u \ du} \mathcal{X} \mid \mathcal{F}_t \right], \tag{14.27}$$

where Q is some choice of martingale measure. Alternatively we can write the price as

$$\Pi_t[\mathcal{X}] = E^Q \left[\frac{\mathbf{M}_T}{\mathbf{M}_t} \mathcal{X} \mid \mathcal{F}_t \right], \tag{14.29}$$

where \mathbf{M} is the stochastic discount factor, defined by

$$\mathbf{M}_t = \frac{1}{B_t} L_t.$$

If we have a simple claim i.e. of the form $\mathcal{X} = \Phi(S_t)$ and if S is Markovian we have

$$e^{-r(T-t)}E^{Q}[\Phi(S_{t}) \mid \mathcal{F}_{t}] = e^{-r(T-t)}E^{Q}[\Phi(S_{t}) \mid S_{t}],$$

and then the pricing process must be of the form $\Pi_t[\Phi] = F(t, S_t)$. We then have F to be the solutions to the PDE

$$F_t(t,s) + \sum_{i=1}^n r s_i F_i(t,s) + \frac{1}{2} \text{tr} \left\{ \sigma^\top D(S) F_{ss} D(S) \sigma \right\} - r F(t,s) = 0, \tag{14.31}$$

$$F(T,s) = \Phi(s), \tag{14.31}$$

where $F_i = \frac{\partial F}{\partial s_i}$ and F_{ss} denotes the Hessian matrix. Furthermore, $\mathrm{tr}(A)$ denotes the trace of A i.e. the sum of the diagonal. We have that the hedging portfolio has value process $V_t^h = F(t,S_t)$ with dynamics

$$dV_t^h = \sum_{i=1}^n F_i(t, S_t) \ dS_t^i.$$

Then we must gave the solution

$$h_t^i = \frac{\partial F}{\partial s_i}(t, S_t), \quad i = 1, ..., n, \tag{14.32}$$

$$h_t^0 = \frac{1}{B_t} \left\{ F(t, S_t) - \sum_{i=1}^n \frac{\partial F}{\partial s_i}(t, S_t) \ S_t^i \right\}.$$
 (14.33)

Proposition 14.8. (Bjork) With L-dynamics as in $dL_t = L_t \varphi_t^{\top} dW_t$, the M-dynamics are

$$d\mathbf{M}_t = -r_t \mathbf{M}_t \ dt + \mathbf{M}_t \varphi_t^{\top} \ dW_t, \tag{14.39}$$

or alternatively in terms of the market price of risk $\lambda_t = -\varphi_t$

$$d\mathbf{M}_t = -r_t \mathbf{M}_t \ dt - \mathbf{M}_t \lambda_t^\top \ dW_t. \tag{14.40}$$

Proposition 14.9. (Bjork) (The Hansen-Jagannathan Bounds) Assume generic absence of arbitrage. Then the following holds for all assets, underlying or derivative, and for all admissible Girsanov kernels φ , and market prices of risk λ .

$$\frac{|\mu_t^p - r_t|}{\|\sigma_t^p\|} \le \|\varphi_t\|, \qquad \frac{|\mu_t^p - r_t|}{\|\sigma_t^p\|} \le \|\lambda_t\|. \tag{14.42}$$

Basic Non-Life Insurance Mathematics

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Probabilistic Machine Learning

9.1 Supervised Learning

In this chapter we restrict our selves to the area of Supervised Learning as we use numerical methods and machine learning algorithms to estimate models in a restricted framework. Take for instance the random forest, this algorithm's estimation method is perfectly capable of being written recursively and so no "leaning" is done in the sense, that the calculations are predetermined from the algorithm. We therefore call the area of study supervised learning instead of the wider area of study machine learning. Let us define what we mean by supervised learning.

Definition. (Supervised Learning) Supervised learning is a field in machine learning that works with labeled data, i.e. data consisting of a set of features X, and a response Y. The goal is to learn a function m^* that maps a given input x to an output y.

We will in this chapter only use data in the form of a spread sheet e.g.

$$\mathcal{D}_n = \left(X_i, Y_i\right)_{i=1,\dots,n} = \left[\begin{array}{cccc} X_{11} & X_{12} & \cdots & X_{1p} & Y_1 \\ X_{21} & X_{22} & \cdots & X_{2p} & Y_2 \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ X_{n1} & X_{n2} & \cdots & X_{nn} & Y_n \end{array} \right],$$

we could however consider any data that may be interpreted by computer software.

The setting is then; assume that we have n independent copies of the random variable $D=(X,Y)\in\mathcal{X}\times\mathcal{Y}$, where X is p-dimensional and Y is one-dimensional. We make no assumption on whether $X_j, j=1,...,p$ and Y are discrete or continuous, however in concrete cases this will be specified. We combine the sample of the n observations in the matrix $\mathcal{D}_n=(X_i,Y_i)_{i=1,...,n}$ being a $n\times p+1$ matrix as in the above. We call \mathcal{D}_n the training data.

This specification does indeed imply that $D_i = (X_i, Y_i)$ are iid. This is actually a bit controversial as we would expect that the distribution of D will shift over time. For instance, the distribution of ages in a population changes over time and have been more right skewed as humanity advances. This may be accounted for by transforming the data such that the distribution becomes the same. This may be done in a variety of ways some example include: 1) transforming to uniform variable with the time dependent distribution F_t , 2) normalizing using a price index and so forth. One should therefore start any analysis by ensuring that the data a given algorithm is trained on is iid.

The job becomes finding a good estimator such that we may predict Y given X i.e. $Y \mid X$. Let us define what an estimator is.

Definition. (Estimator) Consider a training dataset \mathcal{D}_n . A estimator m is a function-valued mapping that takes \mathcal{D}_n as input and associates a function $m_n : \mathcal{X} \to \mathcal{Y}$ i.e. m takes the form

$$m(\mathcal{D}_n) = m_n : \mathcal{X} \to \mathcal{Y}$$

the class of estimators is called \mathcal{G} .

9.1.1 What is a good estimator?

It is easy to construct an estimator \hat{m} for instance by maximum likelihood esitmation, bayes optimization or simply by taking conditional expectations. We are however interested in two problems:

- 1. What is the best class of estimators $\mathcal{G}_0 \subset \mathcal{G}$,
- 2. In the subset of estimators $m \in \mathcal{G}_0$, what is then the best estimator.

We will now discuss the meaning of being the "best" estimator. Obviously, the first type of consideration is regarding the inherent restrictions of some algorithm, where the second is the problem of error comming from the restriction that we do not have infinite observations. We therefore have two problems namely the **inductive bias** from the class \mathcal{G}_0 and the **estimation error** from the available data. A typical problem is that the larger a class \mathcal{G}_0 gives low inductive bias we then may not be able to estimate anything and so the estimation error will be large. The converse also applies.

Definition. Let D = (X, Y) be a random variable on the background space (Ω, \mathcal{F}, P) . Then we define the following:

- A decision rule is a deterministic function $m: \mathcal{X} \to \mathcal{Y}$,
- A loss function is a deterministic function $L: \mathcal{Y} \times \mathcal{Y} \to \mathbb{R}_+$,
- The **risk** of a decision rule m is given a loss function L is r(m) = E[L(Y, m(X))].

Notice that in the definition of risk we see that m is included inside the expectation. This means in particular that the training data \mathcal{D}_n is also accounted for i.e. by the tower rule we have

$$r(m) = \mathbb{E}[L(Y, m(X))] = \mathbb{E}\left[\mathbb{E}\Big[L(Y, m_n(X)) \ \Big| \ \mathcal{D}_n\right]\right] := \mathbb{E}\left[R(m)\right].$$

Some widely used loss function include

- • Quadratic loss function: $L(y_1,y_2)=(y_1-y_2)^2,$
- Poisson Deviance: $L(y_1,y_2)=2\left(y_1\log\frac{y_1}{y_2}-y_1+y_2\right),$
- Binary loss function: $L(y_1, y_2) = 1_{y_1 \neq y_2}$.

Given a loss function L we may find a (possibly non-unique) solution m^* that minimizes R(m). We call this the **Bayes estimator**. The quantity $R(m^*)$ is called the **Bayes risk**. On some special case loss functions we may determine the unique solution.

Lemma. Assume Y is L_2 (square integrable), then the decision function that minimized the risk for the quadratic loss function is

$$m^* = \mathop{\rm argmin}_{m} \; \mathbb{E}[(Y - m(X))^2] = E[Y \mid X = x]$$

i.e. the conditional expectation.

Lemma. Assume $\mathcal{Y} = \{1, ..., K\}$, the decision function that minimizes the risk for the binary loss function satisfies

$$m^* = \operatorname*{argmin}_{m} \ \mathbb{E}[1_{Y \neq m(X)}] = \operatorname*{argmin}_{m} \ \mathbb{P}(Y \neq m(X)) = \operatorname*{argmax}_{k=1,\ldots,K} \mathbb{P}(Y = k \mid X = x).$$

We can now define the prediction risk and the generalization error which relates to the balance of a sufficiently large class \mathcal{G}_0 and how effective the optimal estimator is conditional on the class \mathcal{G}_0 .

Definition. (Conditional risk) Let \mathcal{D}_n be some training data. Given an estimator \hat{m}_n we call

$$R(\hat{m}_n) = \mathbb{E}[L(Y, \hat{m}_n(X)) \ | \ \mathcal{D}_n]$$

the prediction risk or conditional generalized error.

Definition. (Risk) We call

$$r(\hat{m}_n) = \mathbb{E}[R(\hat{m}_n)],$$

the prediction risk or generalized error.

9.1.2 Excess risk

Definition. (Excess Risk) Consider the set \mathcal{G} be the set of all measurable estimators. Fix a subset $\mathcal{G}_0 \subset \mathcal{G}$. Given some training data \mathcal{D}_n consider the Bayes estimator restricted to \mathcal{G}_0 denoted by \hat{m}_n and the unconditional Bayes estimator restricted to \mathcal{G} we define the quantity

$$R(\hat{m}_n) - r(m^*)$$

or

$$\mathbb{E}[R(\hat{m}_n) \mid X_1, ..., X_n] - r(m^*),$$

as the excess risk. This is the difference between the generalization error and the risk obtained by an optimal decision function.

In the context of the above definition we can decompose the risk associated with the optimal estimator $\hat{m}_n \in \mathcal{G}_0$ into the estimation error and the inductive bias.

$$R(\hat{m}_n) - r(m^*) = \underbrace{\left[R(\hat{m}_n) - \inf_{m \in \mathcal{G}_0} R(m)\right]}_{\text{estimation error}} + \underbrace{\left[\inf_{m \in \mathcal{G}_0} R(m) - R(m^*)\right]}_{\text{inductions bias/approximation error}}.$$

where we have to balance the trade-off with a larger \mathcal{G}_0 infer a lower induction bias but larger estimation error and a smaller class \mathcal{G}_0 infer a lower estimation error but larger induction bias.

Definition. (Empirical risk and empirical risk minimizer) Given training data \mathcal{D}_n and a loss function L, we call

$$\hat{R}_n(m) := \sum_{i=1}^n L(Y_i, m(X_i))$$

the empirical risk. Given an additional function class \mathcal{G}_0 ,

$$\underset{m \in \mathcal{G}_0}{\operatorname{argmin}} \ \hat{R}_n(m) = \underset{m \in \mathcal{G}_0}{\operatorname{argmin}} \ \sum_{i=1}^n L(Y_i, m(X_i))$$

is called empirical risk minimizer or (standard leaner).

For larger function classes \mathcal{G}_0 the empirical risk minimizer might not be unique and possibly too noisy. In this case one sometimes adds a penalty term $J_{\lambda}:\mathcal{G}\to\mathbb{R}_+$ that penaltizes the complexity of m and minimizes the penalized empirical risk:

$$\underset{m \in \mathcal{G}_0}{\operatorname{argmin}} \ \hat{R}_{n,\lambda} := \underset{m \in \mathcal{G}_0}{\operatorname{argmin}} \ \sum_{i=1}^n L(Y_i, m(X_i)) + J_{\lambda}(m).$$

If J_{λ} and R_n is convex one can show that

$$\underset{m \in \mathcal{G}_0}{\operatorname{argmin}} \ \hat{R}_{n,\lambda} = \underset{m \in \mathcal{G}_n}{\operatorname{argmin}} \ \hat{R}_n$$

for the class $\mathcal{G}_{\eta} = \{m \in \mathcal{G}_0 \mid J_{\lambda}(m) \leq \eta\}$. Some penalty terms could be

- $\begin{array}{ll} \bullet & J_{\lambda}(m) = \lambda \int m''(x) \; dx, \\ \bullet & J_{\lambda}(m) = \lambda \int |m'(x)| \; dx, \\ \bullet & J_{\lambda}(m) = \lambda \int (m(x))^2 \; dx. \end{array}$

Proposition. (Probability bounds) Let $\tilde{m} = \operatorname{argmin} r(m)$. We have

$$r(\hat{m}_n) - r(\tilde{m}) \leq 2 \sup_{m \in \mathcal{G}_0} \Big| \hat{R}_n(m) - r(m) \Big|,$$

and for all $\lambda \in \Lambda$,

$$r(\hat{m}_{n,\lambda}) - r(\tilde{m}) \leq 2 \sup_{m \in \mathcal{G}_0} \left| \hat{R}_{n,\lambda}(m) - r(m) \right| + J_{\lambda}(\tilde{m}) - J_{\lambda}(\hat{m}_n).$$

Definition. We say \hat{m}_n is ε -accurate with probability $1 - \delta$, if

$$P\Big(R(\hat{m}_n) - \inf_{m \in \mathcal{G}} r(m) > \varepsilon\Big) < \delta.$$

9.2 Training, Validating and Testing

When deciding which method too choose for a given task, one may like to pick the method with the smallest generalization error. However, most machine learning methods depend on hyper parameters and one may first need to decide which hyper parameters are best for the given task. In short: We would like to compare the generalization error of optimally tuned machine learning methods given our data.

Definition. (Training and test set) One often randomly divides the given data into training data and test data:

- $\mathcal{D}_n = (X_i, Y_i)_{i=1,...,n}$ (training data)
- $\mathcal{T}_m = (\tilde{X}_i, \tilde{Y}_i)_{i=1,...,m}$ (test data)

with $n \in [0.8m, 0.95m]$.

Definition. (Training and test error) The empirical risk on the training data is called training error and on the test data test error.

Definition. (Validation set) To tune the hyper parameters for an algorithm, one often randomly divides the given training data into training data (yes: also called training data.) and validation data:

- $\mathcal{D}_{n_1} = (X_i, Y_i)_{i=\tau(1),...,\tau(n_1)}$ (traning data)
- $\mathcal{V}_{n_2} = (\tilde{X}_i, \tilde{Y}_i)_{i=\tau(n_1+1),\dots,\tau(n)}$ (validation data)

where τ is a randomly picked permutation of $\{1,...,n\}$ and $n_1 + n_2 = n$ is respectively the size of the training set and the validation set.

One can now compare different methods via the following simple algorithm:

- 1. Split your data into train, validation and test set.
- 2. For a given method and a rich set of hyper parameter configurations train the method on the training set and compare performance via empirical risk on the validation set.
- 3. For every method, pick the hyper parameter with the smallest empirical risk.
- 4. Compare different methods with the chosen hyper parameters on the test set (trained on training+validation set) and pick the method with smallest empirical risk.

Notice that the procedure above is stochastic and has bias and variance both for selecting the optimal hyper parameters and for selecting the optimal method.

- Bias: Bias occurs because the sample sizes used for learning the hyper parameters n_1 are smaller than the actual training size n and also the full data size n + m.
- Variance: The results are stochastic because the validation and test set are not of infinite size.

Variance can be reduced by repeating steps 1-4 several times. The most popular method for doing so is (nested) cross validation.

9.2.1 Estimating risk

We want to estimate the generalization error of a method $\hat{m}_{n,\lambda}$ that depends on a fixed hyper parameter λ .

Definition. (M-fold Cross validation) Given the indices of a data set $S = \{1, ..., n\}$, M-fold cross validation follows the following steps:

- 1. Divide the data into M disjoint sets $S_1, ..., S_M$ of same size. Define $S_{-l} = \bigcup_{k \neq l} S_k$ being the complement to S_l . $(\#S_l = n/M \text{ and } \#S_{-l} = (M-1)n/M)$
- 2. For each subdivision l=1,...,M train the algorithm on S_{-l} and denote the estimator $\hat{m}_{\lambda}(S_{-l})$.
- 3. Calculate the cross validated empirical risk

$$CV(\hat{m}_{n,\lambda}) = \frac{1}{M} \sum_{l=1}^{M} \frac{1}{|S_{-l}|} \sum_{i \in S_{-l}} L(Y_i, \hat{m}_{\lambda}(S_{-l})(X_i))$$

It is a non-trivial discussion what $CV(\hat{m}_{n,\lambda})$ is estimating. Consider the heuristic

$$\begin{split} CV(\hat{m}_{n,\lambda}) &= \frac{1}{M} \sum_{l=1}^{M} \frac{1}{|S_{-l}|} \sum_{i \in S_{-l}} L(Y_i, \hat{m}_{\lambda}(S_{-l})(X_i)) \\ &\approx \frac{1}{M} \sum_{l=1}^{M} R(Y, \hat{m}_{\lambda}(S_{-l})(X)) \\ &\approx \mathbb{E}\left[R(Y, \hat{m}_{\lambda}(S_{-l})(X))\right]. \end{split}$$

Where we used the law of large numbers in the first approximation. The last approximation is not so clear because the summands are dependent for M > 2.

The bias is minimal for large M since estimation is based on training the algorithm on (M-1)n/M data points. In the case that M=n, M-fold cross validation is also known as leave-one-our cross validation. It is however often not practical because of the computational cost. In practice, setting M=5 or M=10 is common choices.

When deciding on an optimal hyper parameter we would like to pick

$$\underset{\lambda \in \Lambda}{\operatorname{argmin}} \ CV(\hat{m}_{n,\lambda}).$$

But the hyper parameter space Λ is often multi-dimensional and partly continuous. Hence it is infeasible to try out all parameters. Common practice are

- Grid search
- Random search (e.g. pick 200 parameters uniformly random from the parameter space)
- Advanced optimization techniques (i.e. techniques that aim to find the minimzer of a function (here: cross validated empirical risk) without the requirement of knowing the analytical form of the function to optimize.

Above we have discussed how we can use cross validation to pick an optimal parameter for a given method and data set. But how to choose between different methods? One popular way is nested cross validation comprising an inner loop for hyper parameter selection (tuning) and an outer loop for method comparison. Assume that we want to compare J methods $\hat{m}_{n,j}$, j=1,...,J. then we may use nested cross validation.

Definition. (Nested $M_1 - M_2$ Cross-validation) Given the indices of a data set $S = \{1, ..., n\}$, nested $M_1 - M_2$ cross validation follows the following steps:

- 1. Divide the data into M_1 disjoint sets $S_1,...,S_{M_1}$ of same size. Define $S_{-l}=\cup_{k\neq l}S_k$ being the complement to S_l . $(\#S_l=n/M_1$ and $\#S_{-l}=(M_1-1)n/M_1)$
- 2. For each $l=1,...,M_1$, run M_2 -fold cross validation on S_{-l} for all J methods, returning optimal hyper parameters $\hat{\lambda}(j,l),\ j=1,...,J$ and $l=1,...,M_1$. (the one with lowest CV is choosen for each (j,l))
- 3. Calculate the cross validated empirical risk

$$CV(\hat{m}_{n,j}) = \frac{1}{M_1} \sum_{l=1}^{M_1} \frac{1}{|S_{-l}|} \sum_{i \in S_l} L\big(Y_i, \hat{m}_{\lambda(j,l)}(S_{-l})(X_i)\big)$$

4. Pick the method with the smallest risk (and possibly tune again for fitting)

9.3 Linear Models

We may take the linear model as a case study of the methods introduced in the above chapter. As such we consider the squaed loss $L(y_1, y_2) = (y_1 - y_2)^2$ for which we already know the Bayes rule:

$$m^*(x) = \underset{m}{\operatorname{argmin}} \ R(m) = \mathbb{E}[Y \mid X = x].$$

The linear model has the following assumptions. There exists paramaters $\beta_0^*, \beta_1^*, ..., \beta_p^*$, with

$$m^*(x) = \beta_0^* + \sum_{j=1}^p \beta_j^* x_j.$$

In other words, we assume that m^* is a linear function, i.e.,

$$m^* \in \mathcal{G} = \{f: \mathbb{R}^p \to \mathbb{R} \mid f(x) = \beta^\top x\}.$$

Given iid training data $(X_i,Y_i)_{i=1,...,n}$ we have an additive noise model

$$Y_i = \beta_0^* + \sum_{i=1}^p \beta_j^* X_{ij} + \varepsilon_i,$$

with $\varepsilon_i = Y_i - m^*(X_i)$ and hence iid with $\mathbb{E}[\varepsilon_i \mid X_i] = 0$.

Notice, that since we are assuming $m^* \in \mathcal{G}$ we have by assumption no inductive bias and we therefore only consider estimation error i.e.

$$R(\hat{m}_n) - r(m^*).$$

Given the training data we may approximate the coefficients using the following.

Lemma. (Coefficients in the linear model) Under the Linear model assumption we have for j = 1, ..., p

$$\beta_j^* = \frac{\operatorname{Cov} \left(X_{1j}, Y_1 - \sum_{k \in \{1, \dots, p\}} {\{j\}} \, \beta_k^* X_{1k} \right)}{\operatorname{Var}(X_{1j})}$$

In particular, if the components of X are uncorrelated, we have

$$\beta_j^* = \frac{\operatorname{Cov}(X_{1j}, Y_1)}{\operatorname{Var}(X_{1j})}.$$

Lemma. (Bayes risk in the linear model) Under the Linear model assumption we have

- 1. $r(m^*) = \mathrm{Var}(\varepsilon_i)$
- 2. For $m(x) = \beta_0 + \sum_{j=1}^{p} \beta_j x_j$,

$$r(m) - r(m^*) = \|\Sigma^{1/2}(\beta - \beta^*)\|_2^2.$$

Moving forward we will assume $\beta_0^* = 0$ since we can always translate the data and make the centered around 0. We will furthermore use the notation:

$$\mathbf{X} = \begin{bmatrix} X_{11} & \cdots & X_{1p} \\ \vdots & \ddots & \vdots \\ X_{n1} & \cdots & X_{np} \end{bmatrix}, \quad \mathbf{Y} = \begin{pmatrix} Y_1 \\ \vdots \\ Y_n \end{pmatrix}, \quad \boldsymbol{\varepsilon} = \begin{pmatrix} \boldsymbol{\varepsilon}_1 \\ \vdots \\ \boldsymbol{\varepsilon}_n \end{pmatrix}.$$

In this cases the empirical risk takes the form

$$\hat{R}_n(m) = \frac{1}{n} \|\mathbf{Y} - \mathbf{X}\boldsymbol{\beta}\|_2^2.$$

Lemma. (Least squares estimator)

- It holds that $(\mathbf{X}^{\top}\mathbf{X})\hat{\beta} = \mathbf{X}^{\top}\mathbf{Y}$,
- If X has full rank, then

$$\hat{\beta}_n^{LS} = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{Y}.$$

Theorem. (Excess risk Least squares estimator) If $\mathbf{X}^{\top}\mathbf{X}$ is invertible, then

$$\mathbb{E}[R(\hat{m}_n^{LS}) \mid \mathbf{X}] - r(m^*) = \frac{\sigma^2}{n} \cdot \operatorname{tr}\left(\Sigma \hat{\Sigma}^{-1}\right)$$

with $\Sigma = \mathbb{E}[\mathbf{X}^{\top}\mathbf{X}]$ and $\hat{\Sigma} = \frac{1}{n}\mathbf{X}^{\top}\mathbf{X}$.

Chapter 10

Quantative Risk Management

10.1 The Loss Variable

We describe the financial risk with random variables defined on a filtered probability space $(\Omega, \mathcal{F}, P, \{\mathcal{F}_t\}_{t \in \mathbb{R}_+})$. We assume that the value process V_t , denoting the market value of the portfolio at time t, is adapted to the filtration i.e. may be determined at time t or from information available at time t. We will be considering discrete time jumps of size Δt for simplicity and assume that

- the portfolio remains fixed over the time horizon $[t, t + \Delta t)$,
- there are no income or payments made during the time period.

This means in particular that the value

$$\Delta V_{t+\Delta t} = V_{t+\Delta t} - V_t,$$

only depends on the valuation of the components in the portfolio. We will henceforth be using the notation t and t+1 and so forth denoting the intervals $[t+n\Delta t, t+(n+1)\Delta t)$ i.e. t may be any time and the integer representing how many time jumps of size Δt has been made since t.

Under the assumptions above it is reasonable to assume that V_t is Markovian in the sense that there exist $d \ge 1$ random sources $\mathbf{Z}_t = (Z_{t,1},...,Z_{t,d})^{\top}$ such that

$$V_t = f(t, \mathbf{Z}_t) \tag{2.2}$$

for some measurable function $f: \mathbb{R}_+ \times \mathbb{R}^d \to \mathbb{R}$. We will call \mathbf{Z}_t the *risk factors* associated with the portfolio. The could be for instance the stockprice of a stock held in the portfolio. We may now define the *loss variable* as $L_{t+1} := -(V_{t+1} - V_t)$ and *risk-factor changes* as $\mathbf{X}_{t+1} := \mathbf{Z}_{t+1} - \mathbf{Z}_t$ and see that

$$L_{t+1} = -\left(f(t+1, \mathbf{Z}_t + \mathbf{X}_{t+1}) - f(t, \mathbf{Z}_t)\right),\tag{2.3}$$

if one assume differentiability of f we may approximate L_{t+1} as

$$L_{t+1}^{\Delta} := -\left(\frac{\partial f}{\partial t}(t, \mathbf{Z}_t) + \sum_{i=1}^d \frac{\partial f}{\partial z_i}(t, \mathbf{Z}_t) \mathbf{X}_{t+1, i}\right). \tag{2.4}$$

This is obviously a nice linear but the approximation error may be large if Δt is large. We are interested in the the distribution of L_{t+1} such that we may determine sufficiently capital holding in realtion to the risk associated with the assets and liabilities held on the firm's balance sheet.

10.1.1 Risk measures

In the general sense, a risk measure is simply a measurable function $\rho : \mathbb{L} \to \mathbb{R}$ taking a loss variable $L \in \mathbb{L}$ as input and associating a number $\rho(L)$ as output. In this setting we let \mathbb{L} be the set of all real-valued random variables. We may interpret this as the riskyness of the portfolio with associated loss variable L. That is say we have two loss variable from the value processes V and V' i.e. L and L' we say that V is riskier than V' if and only if $\rho(L) \geq \rho(L')$.

We may now consider the fundamental definition of a coherent risk measures as in *Coherent Measures of Risk* by Artzner, Delbean, Eber and Heath (1999), by first stating the definition of a risk measure.

Definition. (Risk Measure) Let \mathbb{L} be the set of all real valued random variable i.e.

$$\mathbb{L} = \{X \mid X : (\Omega, P, \mathcal{F}) \to (\mathbb{R}, \mathbb{B})\}\$$

Any mapping $\rho: \mathbb{L} \to \mathbb{R}$ is called a measure of risk.

We now want some properties to be satisfied by the mapping ρ such that it is a sensable measure of risk. To this we define four axioms as.

Definition. (Coherent Risk Measure) Let ρ be a measure of risk. We say that ρ is a coherent risk measure if and only if ρ satisfies the axioms below.

- 1. Translation invariance: Let $\alpha \in \mathbb{R}$ and r be a predictable process we have $\rho(X + \alpha \cdot r) = \rho(X) + \alpha$.
- $2. \ \, \textbf{Subadditivity:} \ \, Let \, X_1, X_2 \in \mathbb{L}, \, then \, \, \rho(X_1 + X_2) \leq \rho(X_1) + \rho(X_2).$
- 3. Positive homogeneity: Let c > 0, then $\rho(cX) = c\rho(X)$.
- 4. Monotonicity: Let $X, Y \in \mathbb{L}$ such that $X \leq Y$ P-a.s., then $\rho(X) \leq \rho(Y)$.

Remark: We see that axiom (1) gives the equation $\rho(X+\rho(X)\cdot r)=0$ i.e. we may hedge the risk by holding $\rho(X)$ in a asset with rate of return r. The axiom (2) ensures that we take into account the correlation of multiple assets i.e. we will in general not experience the maximal loss of two sources at the same time (although this is possible). If X_1 and X_2 satisfies $\operatorname{Corr}(X_1,X_2)=1$ then $\rho(X_1+X_2)=\rho(X_1)+\rho(X_2)$.

There exist alot of different tangible measures of risk, some being coherent others non-coherent. The most well studied include Value at Risk VaR and Expected Shortfall ES. These two measures are in the realm of the loss distribution approach, however before studying these we introduce a few other worthy mentions: Factor sensitivity measures and scenario based risk measures:

Factor sensitivity measures are measures on the form $\rho = g(\nabla L)$ where g is some d-dimensional measurable function. In this ∇L is the gradient i.e.

$$\nabla L = \nabla \Big(f(t, \mathbf{Z}_t) - f(t+1, \mathbf{Z}_{t+1}) \Big) = \left(\frac{\partial L}{\partial z_1}, ..., \frac{\partial L}{\partial z_d} \right).$$

Scenario based risk measures are measures where one assume that a collection $\mathbf{x}=(x_1,...,x_N)$ of events $x_i\in\Omega$ such that $\sum_{i=1}^N P(x_i)=1-\varepsilon$ for some small $\varepsilon>0$. We may then associate the measures ψ as

$$\psi(L) = \max \{ w_1 L(x_1), ..., w_N L(x_N) \},\,$$

with $\mathbf{w} \geq 0$ and $\sum_{i=1}^N w_i = 1$ being weights. That is ψ gives the largest weighted loss. A natural way of choosing w_i is such that $w_i \approx P(x_i)$, but one may also weight certain events with a larger weight for a more prudent measure. Notice also that the criteria $\sum_{i=1}^N P(x_i) = 1 - \varepsilon$ does not necessarily have to be satisfied when considering the worst possible outcomes.

10.1.1.1 Value at Risk

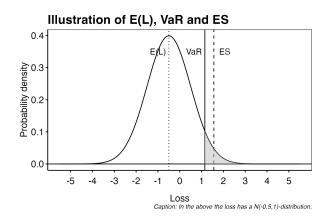
The Value at Risk, henceforth VaR, is a quantile measure of the loss distribution F_L associated with L. We define VaR as such:

Definition 2.8. (McNeil) (Value-at-Risk) Let $\alpha \in (0,1)$ (α being large) we define the VaR of a portfolio with loss variable L at the confidence level α is a given by

$$\begin{split} \operatorname{VaR}_{\alpha}(L) &= \inf \left\{ l \in \mathbb{R} \ : \ P(L > l) \leq 1 - \alpha \right\} \\ &= \inf \left\{ l \in \mathbb{R} \ : \ F_L(l) \geq \alpha \right\} \\ &= F_L^{\leftarrow}(\alpha), \end{split}$$

where F_L^{\leftarrow} is the generalized inverse of F_L .

There exist alot of different tangible measures of risk, some being coherent others non-coherent. The most well studied include Value at Risk VaR and Expected Shortfall ES. These two measures are in the realm of the loss distribution approach, however before studying these we introduce a few other worthy mentions: Factor sensitivity measures and scenario based risk measures:



Chapter 11

Measure theory

Axioms of Probability 11.1

Som udgangspunkt betragtes rummet (Ω, \mathcal{A}) udstyret med en brolægning $\mathcal{A} \subseteq 2^{\Omega}$, hvor $2^{\Omega} = \mathcal{P}(\Omega)$ er mængden af alle delmængder af Ω . Typisk anvendes brolægningerne 1) en algebra eller 2) en σ -algebra.

Definition 2.1. (Protter) (Algebra) En brolægning $A \subseteq 2^{\Omega}$ kaldes en algebra, hvis

- $2. \ \forall A \in \mathcal{A} \Rightarrow A^c = \Omega \ \ A \in \mathcal{A} \ (\textit{lukket under komplementer}),$
- 3. $A_1, A_2, ..., A_n \in \mathcal{A} \Rightarrow \bigcup_{i=1}^n A_i \in \mathcal{A}$ (lukket under endelige foreninger).

Definition 2.2. (Protter) (-algebra) En brolægning $A \subseteq 2^{\Omega}$ kaldes en σ -algebra, hvis

- 2. $\forall A \in \mathcal{A} \Rightarrow A^c \in \mathcal{A}$ (lukket under komplementer),
- 3. $(A_n)_{n\in\mathbb{N}}\subset\mathcal{A}\Rightarrow \cup_{i\in\mathbb{N}}A_i\in\mathcal{A}$ (lukket under tællelige foreninger).

 $\textit{Egenskaber for en σ-algebra: 1)} \ \emptyset \in \mathcal{A}, \ 2) \ A, B \in \mathcal{A} \Rightarrow A \cup B \in \mathcal{A}, \ \text{og 3)} \ (A_i)_{i \in \mathbb{N}} \subseteq \mathcal{A} \Rightarrow \cap_{i \in \mathbb{A}} A_i \in \mathcal{A}.$

Eksempel: Borel-sigma-algebraen $\mathcal{B}(\mathbb{R}^n) = \sigma(\mathcal{C})$, hvor \mathcal{C} kan være en af følgende frembrigersystem. ($\sigma(\mathcal{C})$ benævner den mindste σ -algebra på Ω , hvor $\mathcal{C} \in \mathcal{A}$).

- i. De åbne mængder dvs. $\mathcal{C} = \mathcal{O}^n$
- ii. De lukkede mængder dvs. $\mathcal{C} = \{A^c : A \in \mathcal{O}^n\}$
- iii. Halvåbne mængder/bokse
- iv. Uendelige intervaller som $\mathcal{C} = \{(-\infty, a] : a \in \mathbb{R}\}$

Definition 2.2. (Protter) (Sandsynlighedsmål) Lad $\mathcal{A} \subseteq 2^{\Omega}$ være en σ -algebra. $P: \mathcal{A} \to [0,1]$ kaldes et sandsynlighedsmål hvis

- 1. $P(\Omega) = 1$,
- 2. For $(A_n)_{n\in\mathbb{N}}$ af parvist disjunkte delmængder af Ω gælder $P\left(\bigcup_{n\in\mathbb{N}}A_n\right)=\sum_{n\in\mathbb{N}}P(A_n)$.

Bemærkning: Husk et mål $\mu: \mathcal{A} \to [0, \infty)$ på \mathcal{A} opfylder at 1) $\mu(\emptyset) = 0$ og 2) for en parvis disjunkt familie $(A_n)_{n\in\mathbb{N}}$ gælder $\mu\left(\bigcup_{n\in\mathbb{N}}A_n\right)=\sum_{n\in\mathbb{N}}\mu(A_n).$

Konsekvenser: Umildbare konsekvenser ved definition 2.3 er følgende

- i. $\sum_{i=1}^n P(A_i)=P\left(\bigcup_{i=1}^n A_i\right)$, hvis alle A_i er pavist disjunkte ii. $0\leq P(A)\leq 1,$ for alle $A\in\mathcal{A}$
- iii. $P(A^c) = 1 P(A)$, for alle $A \in \mathcal{A}$
- iv. $P(A) \leq P(B)$ hvis $A \subseteq B$

Theorem 2.3. (Protter) (Opad- og nedadkontinuitet) Lad $P : \mathcal{A} \to [0,1]$ være et ssh. mål på (Ω, \mathcal{A}) . Da gælder

iii.
$$P(A_n) \uparrow P(A)$$
 hvis $A_n \uparrow A$, iv. $P(A_n) \downarrow P(A)$ hvis $A_n \downarrow A$

11.2 Conditional Probability and Independence

Definition 3.1. (Protter) (Uafhængighed) Lad (Ω, \mathcal{A}, P) være et ssh. rum.

- a. Lad $A, B \in \mathcal{A}$ være to hændelser. A og B kaldes uafhængige hvis $P(A \cap B) = P(A)P(B)$.
- b. Lad $A_i \in \mathcal{A}$ for en indeksmængde $i \in I$ (ikke et krav om endelighed eller tællelig). Hændelserne A_i kaldes uafhængige hvis $P(\cap_{i \in J} A_i) = \prod P(A_i)$ for et $J \subseteq I$ med $\#J < +\infty$.

Bemærkning. Der gælder at 1) \emptyset og Ω er uafhængige af alle $A \in \mathcal{A}$ samt $A \in \mathcal{A}$ er uafhængig med sig selv hvis og kun hvis $P(A) = \{0, 1\}$.

Theorem 3.1. (Protter) Lad (Ω, \mathcal{A}, P) være et ssh. rum. Antag $A, B \in \mathcal{A}$ være uafhængige, så er følgende par uafhængige: $(A^c, B), (A, B^c)$ og (A^c, B^c) .

Definition 3.2. (Protter) (Betinget sandsynlighed) Lad $A, B \in \mathcal{A}$ være hændelser og P(B) > 0. Den betingede sandsynlighed A givet B er $P(A|B) = P(A \cap B)/P(B)$.

Theorem 3.2. (Protter) Lad $A, B \in \mathcal{A}$ være hændelser og P(B) > 0.

- a. A og B er uafhængige hvis og kun hvis P(A|B) = P(A).
- b. Funktionen $P(\cdot|B): \mathcal{A} \to [0,1]$ definerer et nyt ssh. mål på \mathcal{A} , kaldet **den betingede sandsynlighed** givet B.

Andet kan tilføjes f.eks. Bayes'.

11.3 Equivalent Probability Measures

11.3.1 The Radon-Nikodym Theorem

Definition A.50. (Bjork) Consider a measurable space (X, \mathcal{F}) on which there are defined two separate measures μ and ν :

• If, for all $A \in \mathcal{F}$, it holds that

$$\mu(A) = 0 \Rightarrow \nu(A) = 0, \tag{A.7}$$

t hen ν is said to be absolutely continuous with respect to μ on \mathcal{F} and we write this as $\nu \ll \mu$.

- If we have both $\mu \ll \nu$ and $\nu \ll \mu$, then μ and ν said to be **equivalent** and we write $\mu \sim \nu$.
- If there exists two events, A and B such that:
 - $-A \cup B = X,$
 - $-A \cap B = \emptyset$.
 - $-\mu(B) = 0$, and $\nu(A) = 0$,

then ν and μ are said to be mutually **singular**, and we write $\mu \perp \nu$.

Theorem A.52. (Bjork) (The Radon-Nikodym Theorem) Consider the measure space (X, \mathcal{F}, μ) , where we assume that μ is finite, i.e. that $\mu(X) < \infty$. Let ν be a measure on (X, \mathcal{F}) such that $\nu << \mu$ on

 \mathcal{F} . Then there exists a non-negative function $f: X \to \mathbb{R}$ such that:

$$f$$
 is \mathcal{F} -measurable (A.9)

$$\int_{X} f(x) \ d\mu(x) < \infty, \tag{A.10}$$

$$\nu(A) = \int_A f(x) \ d\mu(x), \text{ for all } A \in \mathcal{F}. \tag{A.11}$$

The function f is called the **Radon-Nikodym derivative** of ν w.r.t. μ . It is uniquely determined μ -a.e. and we write

$$f(x) = \frac{d\nu(x)}{d\mu(x)},\tag{A.12}$$

or alternatively

$$d\nu(x) = f(x) \ d\mu(x). \tag{A.13}$$

11.3.2 Equivalent Probability Measures

Lemma B.38. (Bjork) For two probability measures P and Q, the relation $P \sim Q$ on \mathcal{F} holds if and only if P(A) = 1 if and only if Q(A) = 1 for all $A \in \mathcal{F}$.

Proposition B.39. (Bjork) Assume that $Q \ll P$ on \mathcal{F} and that $\mathcal{G} \subseteq \mathcal{F}$. Then the Radon-Nikodym derivatives $L^{\mathcal{F}}$ and $L^{\mathcal{G}}$ are related by

$$L^{\mathcal{G}} = E^{P}[L^{\mathcal{F}} \mid \mathcal{G}]. \tag{B.17}$$

Proposition B.41. (Bjork) (Bayes' Theorem) Assume that X is a random variable on (Ω, \mathcal{F}, P) , and let Q be another probability measure on (Ω, \mathcal{F}) the Radon-Nikodym derivative

$$L = \frac{dQ}{dP}$$

on \mathcal{F} . Assume that $X \in L^1(\Omega, \mathcal{F}, Q)$ and \mathcal{G} is a sigma-algebra with $\mathcal{G} \subseteq \mathcal{F}$. Then

$$E^{Q}[X \mid \mathcal{G}] = \frac{E^{P}[L \cdot X \mid \mathcal{G}]}{E^{P}[L \mid \mathcal{G}]}, \ Q - a.s.$$
 (B.18)

11.3.3 Likelihood processes

Proposition C.12. (Bjork) Consider a filtered probability space $(\Omega, \mathcal{F}, P, \mathcal{F}_t)$ on a compact interval [0, T]. Suppose L_T is some non-negative integrable random variable in \mathcal{F}_T . We can then define a new measure Q on \mathcal{F}_T by setting

$$dQ = L_T dP$$

on \mathcal{F}_T and if $E^P[L_T] = 1$ the measure Q will also be a probability measure. The likelihood process L, defined by

$$L_t = \frac{dQ}{dP}, \text{ on } \mathcal{F}_t, \tag{C.8}$$

is a (P, \mathcal{F}_t) -martingale.

Proposition C.13. (Bjork) A process M is a Q-martingale if and only if the process $L \cdot M$ is a P-martingale.

Chapter 12

Random Variables

12.1 Introduction

Definition 1.1. (Hansen) A real-valued random variable X on a probability space (Ω, \mathbb{F}, P) is a measurable map $X : (\Omega, \mathbb{F}) \to (\mathbb{R}, \mathbb{B})$.

We never specify the background space (Ω, \mathbb{F}, P) however we always assume X is $\mathbb{F} - \mathbb{B}$ measurable. This assumption implies $(X \in A) \in \mathbb{F}$ for every $A \in \mathbb{B}$. We may want to show measurability for constructed variables and so it surfises to show measurability for generaters for \mathbb{B} such as checking $(X \le a) \in \mathbb{F}$ for every $a \in \mathbb{R}$.

Definition 1.2. (Hansen) The distribution of a real-valued random variable X, defined on a probability space (Ω, \mathbb{F}, P) , is the collection of probability values

$$P(X \in A)$$
 for $A \in \mathbb{B}$. (1.3)

In other words: the distribution of X is the image measure X(P) on (\mathbb{R}, \mathbb{B}) .

Lemma 1.3. (Hansen) Let X and X' be two real-valued random variables on a probability space (Ω, \mathbb{F}, P) . If

$$P(X = X') = 1$$

then X and X' has the same distribution.

An often used way of summarizing the distribution is through the **distribution function** $F(x) = P(X \le x)$ for some $x \in \mathbb{R}$.

Definition 1.4. (Hansen) A real-valued random variable X has a **discrete** distribution if there is a countable set $S \subset \mathbb{R}$ such that $P(X \in S) = 1$.

Usually S is one of $\mathbb{N}, \mathbb{Z}, \mathbb{Q}$ or a subset of these. We may in the discrete case define the distribution by the point probabilities P(X = x) = p(x) for $x \in S$.

Definition 1.5. (Hansen) A real-valued random variable X has a distribution with **density** $f : \mathbb{R} \to [0, \infty)$ if

$$P(X \in A) = \int_{A} f(x)dx \quad \text{for } A \in \mathbb{B}.$$
 (1.5)

If this is the case we will write $X(P) = f \cdot m$ or $X \sim f \cdot m$.

Definition 1.6. (Hansen) A real-valued random variable X defined on a probability space (Ω, \mathbb{F}, P) is said to have p'th moment for som p > 0 if

$$E|X|^p < \infty \tag{1.12}$$

The collection of all variables that satisfies (1.12) is denoted by $\mathcal{L}^p(\Omega, \mathbb{F}, P)$.

Recall the definition of the **expectation** of X by

$$EX = \int XdP \in R \cup \{-\infty, +\infty\}. \tag{1.11}$$

We recall that for any measurable function $f: \mathbb{R} \to \mathbb{R}$ that are continuous on a set $A \in \mathbb{B}$ such that $P(X \in A) = 1$ we may change variable simply by computing

$$E f(X) = \int f \circ X dP = \int f(x) dX(P)(x).$$

Lemma 1.7. (Hansen) (Markov's inequality) Let X be a non-negative random variable. For any c > 0 it holds that

$$P(X \ge c) \le \frac{EX}{c} \left(\le \frac{EX^n}{c^n} \text{ or } \le \frac{E(\varphi(X))}{\varphi(c)} \right).$$
 (1.14)

for some φ non-negative monotome increasing function.

Some other versions of Markov's inequality can be found in the form of **Chebyshev's inequality**, **Chebyshev-Cantelli's inequality** or **Jensen's inequality** repectively: Let X be a real-valued random variable in $\mathcal{L}^2(\Omega, \mathbb{F}, P)$ it holds for any $\varepsilon > 0$.

$$P(|X - EX| \ge \varepsilon) \le \frac{VX}{\varepsilon^2}$$
(1.15)

$$P(X - EX \ge \varepsilon) \le \frac{VX}{VX + \varepsilon^2}$$
 (prob. 1.13(c))
 $\varphi(EX) \le E(\varphi(X))$

for some convex function φ .

Lemma 1.8. (Hansen) Let X be a non-negative random variable. It holds that

$$EX = \int_0^\infty P(X > t)dt. \tag{1.16}$$

where the integral on the right hand side is with respect to Lebesgue measure.

Definition 1.9. (Hansen) The joint distribution of real-valued random variables $X_1, ..., X_k$, defined on a probability space (Ω, \mathbb{F}, P) , is the collection of probability values

$$P(\mathbf{X} \in A) \quad \text{for } A \in \mathbb{B}_{k}.$$
 (1.21)

In other words: the joint distribution of $X_1, ..., X_k$ (or simply: the distribution of \mathbf{X}) is the image measure $\mathbf{X}(P)$ on $(\mathbb{R}^k, \mathbb{B}_k)$.

Definition 1.11. (Hansen) Real-valued random variables $X_1, ..., X_k$, defined on a probability space (Ω, \mathbb{F}, P) , are jointly independent if

$$P(X_1 \in A_1, ..., X_k \in A_k) = \prod_{i=1}^k P(X_i \in A_i) \quad \text{for } A_1, ..., A_k \in \mathbb{B}.$$
 (1.23)

In other words: the variables are independent if the joint distribution $\mathbf{X}(P)$ equals the product measure $X_1(P) \otimes ... \otimes X_k(P)$.

Theorem 1.12. (Hansen) Let $X_1,...,X_k$ be real-valued random variables defined on a probability space (Ω,\mathbb{F},P) . If the variables are independent and if $E|X_i|<\infty$ for i=1,...,k, then the product $X_1\cdot...\cdot X_k$ has first moment and

$$E(X_1 \cdot ... \cdot X_k) = \prod_{i=1}^k EX_i$$
 (1.24)

The equality only holds for two independent variables. However the **Cauchy-Schwarz inequality** which closely resembles (1.24) holds wether or not X or Y are independent:

$$(E|XY|)^2 \le EX^2EY^2 \text{ or } E|XY| \le \sqrt{EX^2}\sqrt{EY^2}.$$
 (1.25)

Furthermore the theorem give rise to a measure for dependence i.e. the **covariance** between two variables X and Y

$$Cov(X,Y) = E((X - EX)(Y - EY)) = E(XY) - (EX)(EY)$$
(1.26)

with $Cov(X,Y) \neq 0$ if and only if X and Y are dependent. With Cov(X,Y) = 0 the test is inconlusive. However independence implies Cov(X,Y) = 0.

12.2 Conditional expectation

The theory of conditional expectation is well-known from courses on the bachelor. Because of this we will only summarise the most important results.

We consider a background space (Ω, \mathcal{F}, P) and a sub-sigma algebra $\mathcal{G} \subseteq \mathcal{F}$. We assume that some stochastic variable is \mathcal{F} -measurable, that is the mapping $X:(\Omega,\mathcal{F},P)\to(\mathbb{R},\mathbb{B},m)$ is $\mathcal{F}-\mathbb{B}$ -measurable i.e. $\forall B\in\mathbb{B}:\{X\in B\}\in\mathcal{F}$. For some random variable Z defined on the subspace (Ω,\mathcal{G},P) , we say that Z is the conditional expectation of X given \mathcal{G} if

$$\forall G \in \mathcal{G}: \int_G Z(\omega) \ dP(\omega) = \int_G X(\omega) \ dP(\omega).$$

This fact is summed up in the definition below.

Definition B.27. (Bjork) (Conditional expectation) Let (Ω, \mathcal{F}, P) be a probability space and X a random variable in $L^1(\Omega, \mathcal{F}, P)$ (|X| is integrable). Let furthermore \mathcal{G} be a sigma-algebra such that $\mathcal{G} \subseteq \mathcal{F}$. If Z is a random variable with the properties that:

- i. Z is \mathcal{G} -measurable.
- ii. For every $G \in \mathcal{G}$ it holds that

$$\int_{G} Z(\omega) \ dP(\omega) = \int_{G} X(\omega) \ dP(\omega). \tag{B.5}$$

Then we say that Z is the conditional expectation of X given the sigma-algebra \mathcal{G} . In that case we denote Z by the symbol $E[X \mid \mathcal{G}]$.

We see that from the above it always holds that X satisfies (ii). It does not, however, always hold that X is \mathcal{G} -measurable. Given this fact it is not trivial that a random variable $E[X \mid \mathcal{G}]$ even exists. This nontriviality is fortunatly resolved by the Radon-Nikodym theorem.

Theorem B.28. (Bjork) (Existence and uniqueness of Conditional expectation) Let (Ω, \mathcal{F}, P) , X and \mathcal{G} be given as in the definition above. Then the following holds:

- There will always exist a random variable Z satisfying conditions (i)-(ii) above.
- The variable Z is unique, i.e. if both Y and Z satisfy (i)-(ii) then Y = Z P-a.s.

This result ensures that we may condition on any sigma-algebra for instance $\mathcal{G} = \sigma(Y)$ in that case we (pure notation) write

$$E[X \mid \sigma(Y)] = E[X \mid Y], \qquad \sigma(Y) = \sigma\left(\{Y \in A, A \in \mathbb{B}\}\right).$$

In the above $\sigma(Y)$ is simply the smallest sigma-algebra containing all the pre-images of Y, that is the smallest sigma-algebra making Y measurable! Giving this foundation there are a few properties conditional expectation have which is rather useful (for instance the tower property).

Below we assume: Let (Ω, \mathcal{F}, P) be a probability space and X, Y be random variables in $L^1(\Omega, \mathcal{F}, P)$.

Proposition B.29. (Monotinicity/Linearity of Conditional expectation) The following holds:

$$X \le Y \Rightarrow E[X \mid \mathcal{G}] \le E[Y \mid \mathcal{G}], \qquad P - \text{a.s.}$$
 (B.6)

$$E[\alpha X + \beta Y \mid \mathcal{G}] = \alpha E[X \mid \mathcal{G}] + \beta E[Y \mid \mathcal{G}], \qquad \forall \alpha, \beta \in \mathbb{R}. \tag{B.7}$$

Proposition B.30. (Bjork) (Tower property) Assume that it holds that $\mathcal{H} \subseteq \mathcal{G} \subseteq \mathcal{F}$. Then the following hold:

$$E[E[X|\mathcal{G}]|\mathcal{H}] = E[X|\mathcal{H}],\tag{B.8}$$

$$E[X] = E[E[X|\mathcal{G}]]. \tag{B.9}$$

Proposition B.31. (Bjork) Assume X is \mathcal{G} and that both X, Y and XY are in L^1 (only assuming Y is \mathcal{F} -measurable), then

$$E[X|\mathcal{G}] = X, \qquad P - \text{a.s.}$$
 (B.11)

$$E[XY|\mathcal{G}] = XE[Y|\mathcal{G}], \qquad P - \text{a.s.}$$
 (B.12)

Proposition B.32. (Bjork) (Jensen inequality) Let $f : \mathbb{R} \to \mathbb{R}$ be a convex (measurable) function and assume f(X) is in L^1 . Then

$$f(E[X|\mathcal{G}]) \leq E[f(X)|\mathcal{G}], \qquad P-\text{a.s.}$$

Proposition B.37. (Bjork) Let (Ω, \mathcal{F}, P) be a given probability space, let \mathcal{G} be a sub-sigma-algebra of \mathcal{F} , and let X be a square integrable random variable. Consider the problem of minimizing

$$E\left[(X-Z)^2\right]$$

where Z is allowed to vary over the class of all square integrable $\mathcal G$ measurable random variables. The optimal solution $\hat Z$ is then given by.

$$\hat{Z} = E[X|\mathcal{G}].$$

Proof.

Let $X \in L^2(\Omega, \mathcal{F}, P)$ be a random variable. Now consider an arbitrary $Z \in L^2(\Omega, \mathcal{G}, P)$. Recall that $\mathcal{G} \subset \mathcal{F}$ and so X is also in $Z \in L^2(\Omega, \mathcal{G}, P)$, as it is bothe square integrable and \mathcal{G} -measurable. Then

$$E[Z \cdot (X - E[X|\mathcal{G}])] = E[Z \cdot X] - E[Z \cdot E[X|\mathcal{G}]].$$

Then by using the law of total expectation and secondly that Z is \mathcal{G} -measurable we have that

$$E[Z \cdot X] = E[E[Z \cdot X|\mathcal{G}]] = E[Z \cdot E[X|\mathcal{G}]].$$

Combining the two equations gives the desired result. Obviously, we have that

$$X - Z = X - Z + E[X|\mathcal{G}] - E[X|\mathcal{G}] = (X - E[X|\mathcal{G}]) + (E[X|\mathcal{G}] - Z).$$

Then squaring the terms gives

$$(X-Z)^2 = (X - E[X|\mathcal{G}])^2 + (E[X|\mathcal{G}] - Z)^2 + 2(X - E[X|\mathcal{G}])(E[X|\mathcal{G}] - Z)$$

Taking expectation on each side and using linearity of the expectation we have that

$$E[(X-Z)^2] = E\left[(X-E[X|\mathcal{G}])^2\right] + E\left[(E[X|\mathcal{G}]-Z)^2\right] + 2E\left[(X-E[X|\mathcal{G}])(E[X|\mathcal{G}]-Z)\right].$$

We can now use that $E[X|\mathcal{G}] - Z$ is \mathcal{G} -measurable with the above result on the last term.

$$E[(X-Z)^2] = E\left[(X-E[X|\mathcal{G}])^2\right] + E\left[(E[X|\mathcal{G}]-Z)^2\right].$$

Now since X is given the term $E\left[(X-E[X|\mathcal{G}])^2\right]$ is simply a constant not depending on the choice of Z. The optimal choice of Z is then $E[X|\mathcal{G}]$ since this minimizes the second term. The statement is then proved.

Independence 12.3

Definition 3.1. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space. Two events $A, B \in \mathbb{F}$ are independent if

$$P(A \cap B) = P(A)P(B) \tag{3.1}$$

Definition 3.4. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space and let $\mathbb{G}, \mathbb{H} \subset \mathbb{F}$ be two classes of measurable sets. We sat that \mathbb{G} and \mathbb{H} are independent, written $\mathbb{G} \perp \!\!\!\perp \mathbb{H}$, if

$$P(A \cap B) = P(A)P(B)$$
 for all $A \in \mathbb{G}, B \in \mathbb{H}$. (3.2)

Lemma 3.5. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space and let $\mathbb{G}, \mathbb{H} \subset \mathbb{F}$ be two classes of measurable sets. Let $\mathbb{G}_1 \subset \mathbb{G}$ and $\mathbb{H}_1 \subset \mathbb{H}$ be two subclasses. If $\mathbb{G} \perp \!\!\! \perp \mathbb{H}$ then it holds that $\mathbb{G}_1 \perp \!\!\! \perp \mathbb{H}_1$.

Definition 3.6. (Hansen) A class \mathbb{H} of subsets of Ω is a **Dynkin class** if

Lemma 3.7. (Hansen) (Dynkin) Let $\mathbb{D} \subset \mathbb{H}_0 \subset \mathbb{H}$ be three nested classes of subsets of Ω . if

- 1. $\sigma(\mathbb{D}) = \mathbb{H}$,
- $2. A, B \in \mathbb{D} \implies A \cap B \in \mathbb{D}$
- 3. \mathbb{H}_0 is a Dynkin class.

then it holds that $\mathbb{H}_0 = \mathbb{H}$.

Lemma 3.8. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $A \in \mathbb{F}$ be a fixed event. The class

$$\mathbb{H} = \{B \in \mathbb{F} \mid A \perp \!\!\!\perp B\}$$

is a Dynkin class.

Theorem 3.9. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, \mathbb{G}_2 \subset \mathbb{F}$ be two sigma-algebras. Let \mathbb{D}_1 and \mathbb{D}_2 be two classes such that $\sigma(\mathbb{D}_i) = \mathbb{G}_i$ for i = 1, 2. If both \mathbb{D}_1 and \mathbb{D}_2 are \cap -stable then it holds

$$\mathbb{D}_1 \perp\!\!\!\perp \mathbb{D}_2 \quad \Rightarrow \quad \mathbb{G}_1 \perp\!\!\!\perp \mathbb{G}_2.$$

Definition 3.10. (Hansen) Two real-valued random variable X and Y on a background space (Ω, \mathbb{F}, P) are **independent**, written $X \perp \!\!\! \perp Y$, if the corresponding sigma-algebras $\sigma(X)$ and $\sigma(Y)$ are independent.

Definition 3.15. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, ..., \mathbb{G}_n \subset \mathbb{F}$ be finitely many classes of measurable sets. We say that $\mathbb{G}_1, ..., \mathbb{G}_n$ are **jointly independent**, written $\mathbb{G}_1 \perp \!\!\! \perp ... \perp \!\!\! \perp \mathbb{G}_n$, if

$$P(A_1 \cap \dots \cap A_n = \prod_{i=1}^n P(A_i) \quad \text{ for } A_1 \in \mathbb{G}_1, \dots, A_n \in \mathbb{G}_n.$$
 (3.8)

Lemma 3.16. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, ..., \mathbb{G}_n \subset \mathbb{F}$ be finitely many classes of measurable sets. It holds that

$$\mathbb{G}_1 \perp \!\!\! \perp \ldots \perp \!\!\! \perp \mathbb{G}_n \quad \Rightarrow \quad \mathbb{G}_1 \perp \!\!\! \perp \ldots \perp \!\!\! \perp \mathbb{G}_{n-1}$$

provided that $\Omega \in \mathbb{G}_n$.

Theorem 3.17. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, ..., \mathbb{G}_n \subset \mathbb{F}$ be sigma-algebras. Let $\mathbb{D}_1,...,\mathbb{D}_n$ be classes such that $\sigma(\mathbb{D}_i)=\mathbb{G}_1$ for i=1,...,n. Suppose that for all lengths k=2,...,n an all choices of indices $\leq j_1 < \dots < j_k \leq n$ it holds that

$$\mathbb{D}_{j_1} \perp \dots \perp \!\!\!\perp \mathbb{D}_{j_k} \tag{3.9}$$

If all the generators \mathbb{D}_i are \cap -stable, then it holds that $\mathbb{G}_1 \perp \!\!\! \perp \ldots \perp \!\!\! \perp \mathbb{G}_n$.

Lemma 3.18. (Hansen) (Grouping) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, ..., \mathbb{G}_n \subset \mathbb{F}$ be sigma-algebras. It holds that

$$\mathbb{G}_1 \perp \!\!\! \perp \ldots \perp \!\!\! \perp \mathbb{G}_n \quad \Rightarrow \quad \mathbb{G}_1 \perp \!\!\! \perp \ldots \perp \!\!\! \perp \mathbb{G}_{n-2} \perp \!\!\! \perp \sigma(\mathbb{G}_{n-1},\mathbb{G}_n).$$

Definition 3.19. (Hansen) The real-valued random variables $X_1, ..., X_n$ on a background space (Ω, \mathbb{F}, P) are **jointly independent**, written $X_1 \perp \!\!\! \perp ... \perp \!\!\! \perp X_n$, if the corresponding sigma-algebras $\sigma(X_1), ..., \sigma(X_n)$ are jointly independent.

Definition 3.20. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $(\mathbb{G}_i)_{i \in I}$ be a family of classes of measurable sets. We say that the family $(\mathbb{G}_i)_{i \in I}$ is **jointly independent** if any finite subfamily is jointly independent.

Theorem 3.21. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, \mathbb{G}_2, ... \subset \mathbb{F}$ be sigma-algebras. Let $\mathbb{D}_1, \mathbb{D}_2, ...$ be classes such that $\sigma(\mathbb{D}_n) = \mathbb{G}_n$ for all $n \in \mathbb{N}$. Suppose that for all lengths $k \in \mathbb{N}$ and all choices of indices $1 \leq j_1 < ... < j_k$ it holds that

$$\mathbb{D}_{j_1} \perp \!\!\!\perp \ldots \perp \!\!\!\perp \mathbb{D}_{j_k}. \tag{3.14}$$

If all the generators \mathbb{D}_n are \cap -stable, then it holds that $\mathbb{G}_1 \perp \!\!\! \perp \mathbb{G}_2 \perp \!\!\! \perp \dots$

Lemma 3.22. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space, and let $\mathbb{G}_1, \mathbb{G}_2, ... \subset \mathbb{F}$ be sigma-algebras. It holds that

$$\mathbb{G}_1 \perp\!\!\!\perp \mathbb{G}_2 \perp\!\!\!\perp \dots \quad \Rightarrow \quad \mathbb{G}_1 \perp\!\!\!\perp \dots \perp\!\!\!\perp \mathbb{G}_n \perp\!\!\!\perp \sigma(\mathbb{G}_{n+1},\mathbb{G}_{n+2},\dots).$$

Definition 3.23. (Hansen) The real-valued random variables $(X_i)_{i\in I}$ on a background space (Ω, \mathbb{F}, P) are **jointly independent** if the corresponding sigma-algebras $(\sigma(X_i))_{i\in I}$ are jointly independent.

Definition 3.28. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space. A sigma-algebra $\mathbb{G} \subset \mathbb{F}$ satisfies a **zero-one** law if

$$P(A) \in \{0, 1\}$$
 for all $A \in \mathbb{G}$.

Theorem 3.29. (Hansen) Let (Ω, \mathbb{F}, P) be a probability space and let $\mathbb{G} \subset \mathbb{F}$ be a sigma-algebra. The following three conditions are equivalent:

- 1. For any sigma-algebra $\mathbb{H} \subset \mathbb{F}$ it holds that $\mathbb{G} \perp \!\!\! \perp \mathbb{H}$,
- 2. It holds that $\mathbb{G} \perp \!\!\!\perp \mathbb{G}$,
- 3. \mathbb{G} satisfies a 0-1 law.

Definition 3.30. (Hansen) Let $X_1, X_2, ...$ be real-valued random variables on a background space (Ω, \mathbb{F}, P) . The tail sigma-algebra of the process is defined as

$$\mathbb{J}(X_1,X_2,\ldots)=\bigcap_{n=1}^\infty\sigma(X_n,X_{n+1},\ldots).$$

Theorem 3.32. (Hansen) (Kolmogorov's zero-one law) Let $X_1, X_2, ...$ be real-valued random variables on a background space (Ω, \mathbb{F}, P) . If $X_1 \perp \!\!\! \perp X_2 \perp \!\!\! \perp ...$ then the tail-algebra $\mathbb{J}(X_1, X_2, ...)$ satisfies a 0-1 law.

Lemma 3.35. (Hansen) (2nd half of Borel-Cantelli) Let (Ω, \mathbb{F}, P) be a probability space, and let $A_1, A_2, ...$ be a sequence of \mathbb{F} -measurable sets. If $A_1 \perp \!\!\!\perp A_2 \perp \!\!\!\perp ...$ then it holds that

$$\sum_{n=1}^{\infty} P(A_n) < \infty \iff P(A_n \text{ i.o.}) = 0.$$

12.4 Moment generating function

Let X be a random variable with distribution function $F(x) = P(X \le x)$ and Y be a random variable with distribution function $G(y) = P(Y \le y)$.

Definition. (Ex. FinKont) The moment generating function or Laplace transform of X is

$$\psi_X(\lambda) = E\left[e^{\lambda X}\right] = \int_{-\infty}^{\infty} e^{\lambda x} dF(x)$$

provided the expectation is finite for $|\lambda| < h$ for some h > 0.

The MGF uniquely determine the distribution of a random variable, due to the following result.

Theorem. (Ex. FinKont) (Uniqueness) If $\psi_X(\lambda) = \psi_Y(\lambda)$ when $|\lambda| < h$ for some h > 0, then X and Y has the same distribution, that is, F = G.

There is also the following result of independence for Moment generating functions.

Theorem. (Ex. FinKont) (Independence) If

$$E\left[e^{\lambda_1X+\lambda_2Y}\right]=\psi_X(\lambda_1)\psi_Y(\lambda_2)$$

for $|\lambda_i| < h$ for i = 1, 2 for some h > 0, then X and Y are independent random variables.

12.5 Standard distributions

12.5.1 Normal disribution

The following gives a comprehensive table of the standard some properties.

Description	Symbol	Normal distribution
Definition	~	$X \sim \mathcal{N}(\mu, \sigma^2)$
Parameters	$ heta \in \Theta$	$ heta = (\mu, \sigma^2) \in \mathbb{R} \times \mathbb{R}_+$
Support	$\operatorname{Im}(X)$	$x\in\mathbb{R}$
Density	f	$rac{1}{\sqrt{2\pi\sigma^2}}e^{-\left(rac{x-\mu}{\sqrt{2\sigma^2}} ight)^2} \ rac{1}{2}\left(1+N\left(rac{x-\mu}{\sqrt{2\sigma^2}} ight) ight)$
Distribution	F	$\frac{1}{2}\left(1+N\left(\frac{x-\mu}{\sqrt{2-2}}\right)\right)$
Mean value	E[X]	
Variance	$\operatorname{Var}(X)$	$rac{\mu}{\sigma^2}$
MGF*	$\psi_X = E[e^{\lambda X}]$	$e^{\mu\lambda+rac{1}{2}\lambda^2\sigma^2} \ e^{it\mu-rac{1}{2}\sigma^2t^2}$
Characteristic function	$\varphi_X(t) = E[e^{itX}]$	$e^{it\mu-rac{1}{2}\sigma^2t^2}$

In the table above we used the abbreviations: *MGF = Moment Generating function.

We also used the shorthand: N being the distribution of a standard normal distributed variable $\mathcal{N}(0,1)$.

Chapter 13

Discrete Time Stochastic Processes

13.1 Convergence concepts

We start this chapter by refering to a sequence X_1, X_2, \ldots of real-valued random variables as a **proces**. Consider the event $(X_n \to X) = \{\omega \in \Omega \mid X_m(\omega) \to X(\omega) \text{ for } n \to \infty\}$. We want to study such convergence in detail. However first we check measurability. Consider a family $(A_i)_{i \in I} \subset \Omega$ and observe that

$$\left\{\omega \in \Omega \mid \forall i \in I : \omega \in A_i\right\} = \bigcap_{i \in I} A_i, \tag{2.1}$$

$$\left\{\omega\in\Omega\mid\exists i\in I:\omega\in A_i\right\}=\bigcup_{i\in I}A_i. \tag{2.2}$$

From the standard N, ε definition of a convergent sequence $(x_n)_{n \in \mathbb{N}}$ we may formulate this convergens in the stochastic setting:

$$\begin{split} (X_n \to X) &= \Big\{ \omega \in \Omega \ | \ \forall \varepsilon > 0 \\ \exists N \in \mathbb{N} \\ \forall n \geq N : |X_n(\omega) - X(\omega)| < \varepsilon \Big\} \\ &= \Big(\forall \varepsilon > 0 \\ \exists N \in \mathbb{N} \\ \forall n \geq N : |X_n - X| < \varepsilon \Big) \\ &= \bigcap_{\epsilon \in \mathbb{R}^+} \Big(\exists N \in \mathbb{N} \\ \forall n \geq N : |X_n - X| < \varepsilon \Big) \\ &= \bigcap_{\epsilon \in \mathbb{R}^+} \bigcup_{N=1}^{\infty} \bigcap_{n=N}^{\infty} \Big(|X_n - X| < \varepsilon \Big) \in \mathbb{F} \quad \text{ for all } \varepsilon > 0. \end{split}$$

Hence $(X_n \to X)$ lies in $\mathbb F$ since $(|X_n - X| < \varepsilon)$ lies in $\mathbb F$ since $X_n - X$ is measurable.

Lemma 2.1. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . It holds that

$$(X_n \to X) \in \mathbb{F}.$$

Definition 2.2. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . We say that X_n converges to X almost surely, written $X_n \overset{a.s.}{\to} X$, if

$$P(X_n \to X) = 1. \tag{2.6}$$

Lemma 2.3. (Hansen) Let $X, X', X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . If $X_n \overset{a.s.}{\to} X$ and $X_n \overset{a.s.}{\to} X'$ then X = X' almost surely.

Lemma 2.7. (Hansen) Let $X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . Then

$$(X_n)$$
 is Cauchy $\in \mathbb{F}$.

Lemma 2.8. (Hansen) Let $X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . If $P(X_n)$ is Cauchy = 1 then there exists and \mathbb{F} -measurable real-valued random variable X such that $X_n \stackrel{a.s.}{\to} X$.

Theorem 2.10. (Hansen) Let $X_1, X_2, ...$ and $Y_1, Y_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . Assume that the X-process and the Y-process have the same distribution in the sense that $(X_1, ..., X_n)$ has the same distribution ad $(Y_1, ..., Y_n)$ for all $n \in \mathbb{N}$. If $X_n \stackrel{a.s.}{\to} X$ for som limit variable X, there is a limit variable Y such that $Y_n \stackrel{a.s.}{\to} Y$.

Definition 2.11. (Hansen) Let $\mathbf{X}, \mathbf{X}_1, \mathbf{X}_2, \dots$ be \mathbb{R}^k valued random variables on (Ω, \mathbb{F}, P) . We say that \mathbf{X}_n converges to \mathbf{X} almost surely, written $\mathbf{X}_n \overset{a.s.}{\to} \mathbf{X}$, if

$$|\mathbf{X}_n - \mathbf{X}| \stackrel{\text{a.s.}}{\to} 0.$$
 (2.15)

Lemma 2.12. (Hansen) Let $\mathbf{X}, \mathbf{X}_1, \mathbf{X}_2, ...$ be \mathbb{R}^k valued random variables on (Ω, \mathbb{F}, P) such that $\mathbf{X}_n \overset{a.s.}{\to} \mathbf{X}$. Let $f: \mathbb{R}^k \to \mathbb{R}^m$ be a measurable map. Assume that there is a set $A \in \mathbb{B}_k$ such that f is continuous on A and such that $P(\mathbf{X} \in A) = 1$. Then it holds that $f(\mathbf{X}_n) \overset{a.s.}{\to} f(\mathbf{X})$.

Definition 2.13. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . We say that X_n converges to X in probability, written $X_n \stackrel{P}{\to} X$, if

$$\forall \varepsilon > 0: \quad P(|X_n - X| \ge \varepsilon) \to 0 \quad \text{for } n \to \infty.$$
 (2.17)

Lemma 2.14. (Hansen) Let $X, X', X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . If $X_n \stackrel{P}{\to} X$ and $X_n \stackrel{P}{\to} X'$ then X = X' almost surely.

Lemma 2.14. (Hansen) Let $X, X', X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . If $X_n \stackrel{a.s.}{\to} X$, then $X_n \stackrel{P}{\to} X$.

Definition 2.17. (Hansen) Let $\mathbf{X}, \mathbf{X}_1, \mathbf{X}_2, \dots$ be \mathbb{R}^k valued random variables on (Ω, \mathbb{F}, P) . We say that \mathbf{X}_n converges to \mathbf{X} in probability, written $\mathbf{X}_n \overset{P}{\to} \mathbb{X}$, if

$$|\mathbf{X}_n - \mathbf{X}| \stackrel{\mathrm{P}}{\to} 0.$$
 (2.23)

Lemma 2.18. (Hansen) Let $X, Y, X_1, Y_1, X_2, Y_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . It holds that

$$\begin{pmatrix} X_n \\ Y_n \end{pmatrix} \stackrel{P}{\to} \begin{pmatrix} X \\ Y \end{pmatrix} \quad \iff \quad X_n \stackrel{P}{\to} X \text{ and } Y_n \stackrel{P}{\to} Y. \tag{2.24}$$

Definition 2.19. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables in $\mathcal{L}^p(\Omega, \mathbb{F}, P)$ for some $p \geq 1$. We say that X_n converges to X in \mathcal{L}^p , written $X_n \overset{\mathcal{L}^p}{\to} X$, if

$$||X_n - X||_p \to 0.$$
 (2.27)

Where the p'th norm is defined as the mapping $\|\cdot\|_p:\Omega\to[0,\infty)$ given by $X\mapsto\left(\int_\Omega|X|^p\;dP\right)^{1/p}$.

One might also define convergence in \mathcal{L}^p by simply saying if $X_n \stackrel{\mathcal{L}^p}{\to} X$ then $E \|X_n - X\|_p \to 0$.

Lemma 2.20. (Hansen) (Extended Cauchy-Schwarz inequality) Let $X, Y \in \mathcal{L}^p(\Omega, \mathbb{F}, P)$ for some $p \geq 1$. For any $a \in [0, p]$ it holds that

$$E|X|^a|Y|^{p-a} \le \left(E|X|^p\right)^{\frac{a}{p}} \left(E|Y|^p\right)^{\frac{p-a}{p}}.$$
 (2.29)

Theorem 2.21. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables in $\mathcal{L}^p(\Omega, \mathbb{F}, P)$ for some $p \in \mathbb{N}$. If $X_n \stackrel{\mathcal{L}^p}{\to} X$, then it holds that $E X_n^p \to E X^p$.

Lemma 2.22. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables in $\mathcal{L}^p(\Omega, \mathbb{F}, P)$ for some $p \geq 1$. If $X_n \stackrel{\mathcal{L}^p}{\to} X$, then $X_n \stackrel{P}{\to} X$.

Lemma 2.25. (Hansen) (Borel-Cantelli) Let (Ω, \mathbb{F}, P) be a probability space, and let A_1, A_2, \ldots be a sequence of \mathbb{F} -measurable sets. It holds that

$$\sum_{n=1}^{\infty} P(A_n) < \infty \quad \Rightarrow \quad P(A_n \text{ i.o.}) = 0.$$

Let A_1, A_2, \dots be a sequence of subsets of Ω . We define

$$(A_n \text{ i.o.}) = \bigcap_{n=1}^{\infty} \bigcup_{m=n}^{\infty} A_m, \quad (A_n \text{ evt.}) = \bigcup_{n=1}^{\infty} \bigcap_{m=n}^{\infty} A_m.$$

One might also define $Y=\sum_{n=1}^{\infty}1_{A_n}$ and realise that $(A_n \text{ i.o.})=(Y=\infty)$ and $(A_n \text{ evt.})=(Y<\infty)$. Also by de Morgan's law it follows that $(A_n \text{ evt.})^c=(A_n^c \text{ i.o.})$.

Theorem 2.26. (Hansen) Let $X, X_1, X_2, ...$ be real-valued random variables on (Ω, \mathbb{F}, P) . If

$$\forall \varepsilon > 0: \quad \sum_{n=1}^{\infty} P(|X_n - X| \ge \varepsilon) < \infty, \tag{2.32}$$

then it holds that $X_n \stackrel{a.s.}{\to} X$.

Theorem 2.27. (Hansen) Let X, X_1, X_2, \ldots be real-valued random variables on (Ω, \mathbb{F}, P) . If $X_n \overset{P}{\to} X$, then there is a subsequence X_{n_1}, X_{n_2}, \ldots such that $X_{n_k} \overset{a.s.}{\to} X$ for $k \to \infty$.

Lemma 2.28. (Hansen) Let $\mathbf{X}, \mathbf{X}_1, \mathbf{X}_2, ...$ be \mathbb{R}^k -valued random variables on (Ω, \mathbb{F}, P) such that $\mathbf{X}_n \overset{P}{\to} \mathbf{X}$. Let $f: \mathbb{R}^k \to \mathbb{R}^m$ be a measurable map. Assume that there is a set $A \in \mathbb{B}_k$ such that f is continuous on A and such that $P(\mathbf{X} \in A) = 1$. Then it holds that $f(\mathbf{X}_n) \overset{P}{\to} f(\mathbf{X})$.

Lemma. (Fatou's Lemma) Let (Ω, \mathbb{F}, P) be a measure space (here probability space). Let $f_n : \mathcal{X} \to [0, \infty]$, with $\mathcal{X} \in \mathbb{F}$, be a sequence of non-negative measurable functions. Assume f_n converge pointwise to $f : \mathcal{X} \to [0, \infty)$. Then

$$\int_{\mathcal{X}} \liminf_{n \to \infty} f_n \ dP \le \liminf_{n \to \infty} \int_{\mathcal{X}} f_n \ dP.$$

Lemma. (Holder's Inequality) Let (Ω, \mathbb{F}, P) be a measure space (here probability space). Let f and g be real-valued (or complex-valued) functions defined on Ω . Assume f and g are measurable. For any $p, q \geq 1$ such that $\frac{1}{p} + \frac{1}{q} = 1$ it holds that

$$\left(\int_{\Omega}|fg|^1\ dP\right)^1\leq \left(\int_{\Omega}|f|^p\ dP\right)^{1/p}\left(\int_{\Omega}|g|^q\ dP\right)^{1/q}$$

13.1.1 Sums and average processes

Lemma 4.1. (Hansen) Let $X_1, ..., X_n$ be independent real-valued random variables with $E[X_i^4] < \infty$ for all i. If $E[X_i] = 0$ for all i then it holds that

$$E\left(\sum_{i=1}^{n}X_{i}\right)^{4} = \sum_{i=1}^{n}E\,X_{i}^{4} + 6\sum_{i=1}^{n-1}\sum_{j=i+1}^{n}E\,X_{i}^{2}\,E\,X_{j}^{2}.$$

Theorem 4.2. (Hansen) (SLLN, weak form) Let $X_1, X_2, ...$ be a sequence of independent and identically distributed real-valued random variables. If $E[X_1^4] < \infty$ it holds that

$$\frac{1}{n} \sum_{i=1}^{n} X_i \stackrel{\text{a.s.}}{\to} EX_1. \tag{4.3}$$

Theorem 4.10. (Hansen) (Etemadi's maximal inequality) Let $X_1, ..., X_n$ be independent real-valued random variables. Consider the cumulative sums

$$S_k = \sum_{i=1}^k X_i$$
 for $k = 1, ..., n$.

For any $\alpha > 0$ it holds that

$$P\left(\max_{j=1,\dots,n} |S_j| \ge 3\alpha\right) \le 3 \max_{j=1,\dots,n} P(|S_j| \ge \alpha). \tag{4.11}$$

Theorem 4.11. (Hansen) (Levy's maximal inequality) Let $X_1, ..., X_n$ be independent real-valued random variables, each with a symmetric distribution. Consider the cumulative sums

$$S_k = \sum_{i=1}^k X_i$$
 for $k = 1, ..., n$.

For any $\alpha > 0$ it holds that

$$P\left(\max_{j=1,...,n} S_j \ge \alpha\right) \le 2P(S_j \ge \alpha). \tag{4.13}$$

Corollary 4.12. (Hansen) Let $X_1, ..., X_n$ be independent real-valued random variables, each with a symmetric distribution. Consider the cumulative sums

$$S_k = \sum_{i=1}^k X_i$$
 for $k = 1, ..., n$.

For any $\alpha > 0$ it holds that

$$P\left(\max_{j=1,\dots,n} |S_j| \ge \alpha\right) \le 2P(|S_j| \ge \alpha). \tag{4.14}$$

Theorem 4.13. (Hansen) (Skorokhod) Let $X_1, X_2, ...$ be a sequence of independent real-valued random variables, and consider the cumulative sums $S_k = \sum_{i=1}^k X_i$. Let S be a potential limit variable. It holds that

$$S_n \stackrel{\mathrm{P}}{\to} S \quad \Rightarrow \quad S_n \stackrel{\mathrm{a.s.}}{\to} S.$$

Corollary 4.14. (Hansen) (Khintchine-Kolmogorov) Let $X_1, X_2, ...$ be a sequence of independent real-valued random variables. Assume $E X_n^2 < \infty$ and that $E X_n = 0$ for every $n \in \mathbb{N}$. Consider the cumulative sums $S_k = \sum_{i=1}^k X_i$. If

$$\sum_{n=1}^{\infty} E X_n^2 < \infty \tag{4.18}$$

then there exist a limit variable S such that $S_n \to S$ almost surely and in \mathcal{L}^2 . The limit variable satisfies

$$ES = 0$$
 and $VS = \sum_{n=1}^{\infty} VX_n$.

Theorem 4.17. (Hansen) Let X_1, X_2, \ldots be a sequence of independent real-valued random variables, and consider the cumulative sums $S_k = \sum_{i=1}^k X_i$. Let S be a potential limit variable. Assume that there is a constant c > 0 such that $P(|X_n| \le c) = 1$ for all n, and assume that $E[X_n] = 0$ for all n. The the three statements 1. $S_n \stackrel{P}{\to} S$, 2. $S_n \stackrel{\text{a.s.}}{\to} S$ 3. $S_n \stackrel{\mathcal{L}^2}{\to} S$ are equivalent.

Lemma 4.18. (Hansen) Let X_1, X_2, \ldots be a sequence of independent real-valued random variables. Assume that there is a constant c>0 such that $P(|X_n|\leq c)=1$ for all n. If the associated random walk $S_n=\sum_{i=1}^n X_i$ satisfies that $S_n\to S$ almost surely for some limit variable then it holds that

1)
$$\sum_{n=1}^{N} E X_n$$
 converges in \mathbb{R} for $N \to \infty$, (13.1)

$$2) \quad \sum_{n=1}^{\infty} V(X_n) < \infty. \tag{13.2}$$

Theorem 4.19. (Hansen) (Kolmogorov's 3-series theorem) Let $X_1, X_2, ...$ be a sequence of independent real-valued random variables. Consider the assoiciated random walk $S_n = \sum_{i=1}^n X_i$. If there is a cut-off value c > 0 such that the capped variables $\tilde{X}_n = 1_{|X_n| \le c} X_n$ satisfies that

$$1)\quad \sum_{n=1}^{\infty}P(X_{n}\neq \tilde{X}_{n})<\infty,$$

2)
$$\sum_{n=1}^{N} E \tilde{X}_n$$
 converges in \mathbb{R} for $N \to \infty$,

$$3)\quad \sum_{n=1}^{\infty}V(\tilde{X}_{n})<\infty,$$

then there is a real-valued limit variable S such that $S_n \to S$ almost surely. Conversely, if $(S_n)_{n \in \mathbb{N}}$ is almost surely convergent, then the three series above converge for any cut-off value c > 0.

Lemma 4.20. (Hansen) Let $(x_n)_{n\in\mathbb{N}}$ be a real-valued sequence, and let c be a real number. It holds that

$$x_n \to c \text{ for } n \to \infty \quad \Rightarrow \quad \frac{1}{n} \sum_{i=1}^n x_i \to x \text{ for } n \to \infty.$$

Lemma 4.21. (Hansen) (Kronecker) Let $(x_n)_{n\in\mathbb{N}}$ be real-valued sequence, and let c be a real number. It holds that

$$\sum_{i=1}^n \frac{x_i}{i} \to c \quad \Rightarrow \quad \frac{1}{n} \sum_{i=1}^n x_i \to 0$$

for $n \to \infty$.

Lemma 4.23. (Hansen) Let $X_1, X_2, ...$ be a sequence of identically distributed real-valued random variables, and let $\tilde{X}_n = 1_{\{|X_n| \le n} X_n$. If $E[X_1] < \infty$ it holds that

$$\sum_{n=1}^{\infty} \frac{E \ \tilde{X}_n^2}{n^2} < \infty.$$

Theorem 4.24. (Hansen) (SLLN, strong version) Let $X_1, X_2, ...$ be sequence of independent and identically distributed real-valued random variables. If $E|X_1| < \infty$ it holds that

$$\frac{1}{n} \sum_{n=1}^{\infty} X_i \stackrel{\text{a.s.}}{\to} E X_1. \tag{4.24}$$

Theorem 4.25. (Hansen) (SLLN, \mathcal{L}^p -version) Let $X_1, X_2, ...$ be sequence of independent and identically distributed real-valued random variables. If $E|X_1|^p < \infty$ for some $p \ge 1$, then it holds that

$$\frac{1}{n} \sum_{n=1}^{\infty} \stackrel{\mathcal{L}^p}{\to} E X_1. \tag{4.26}$$

Lemma 4.26. (Hansen) Let $X_1, X_2, ...$ be a sequence of pairwise independent, identically distributed real-valued random variables with $E|X_1| < \infty$. Let $n_1 < n_2 < ...$ be a sequence of natural numbers. If there are constants $C_1, C_2 > 0$ and $\alpha > 1$ such that

$$C_1 \alpha^k \le n_k \le C_2 \alpha^k \quad \text{for } k \to \infty$$
 (4.27)

then it holds that

$$\frac{1}{n_k} \sum_{i=1}^{n_k} X_i \overset{\text{a.s.}}{\to} E X_1 \quad \text{ for } k \to \infty.$$

Theorem 4.27. (Hansen) (Etemahdi's version) Let $X_1, X_2, ...$ be a sequence of pairwise independent, identically distributed real-valued random variables. If $E|X_1| < \infty$ it holds that

$$\frac{1}{n} \sum_{i=1}^{n} X_i \stackrel{\text{a.s.}}{\to} E X_1. \tag{4.30}$$

13.1.2 Ergodic Theory

Definition 5.3. (Hansen) Let $(\mathcal{X}, \mathbb{E})$ be a measurable space and let $T : \mathcal{X} \to \mathcal{X}$ be measurable. A probability measure μ on $(\mathcal{X}, \mathbb{E})$ is **invariant** under T if

$$\mu(T^{-1}(A)) = \mu(A) \quad \text{for all } A \in \mathbb{E}$$
 (5.5)

In this case we call the quadruple $(\mathcal{X}, \mathbb{E}, \mu, T)$ a measure-preserving dynamical system.

We say that a set $A \in \mathbb{E}$ is an **invariant set** if $T^{-1}(A) = A$ i.e. the orbit of all $x \in A$ stays in A.

Definition 5.5. (Hansen) A measure-preserving dynamical system $(\mathcal{X}, \mathbb{E}, \mu, T)$ is ergodic if

$$T^{-1}(A) = A, \ A \in \mathbb{E} \quad \Rightarrow \quad \mu(A) \in \{0, 1\}. \tag{5.7}$$

Definition 5.6. (Hansen) A measure-preserving dynamical system $(\mathcal{X}, \mathbb{E}, \mu, T)$ is **mixing** if

$$\mu(A \cap T^{-n}(B)) \to \mu(A)\mu(B)$$
 for all $A, B \in \mathbb{E}$. (5.8)

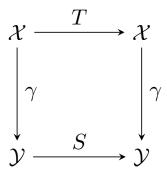
Lemma 5.7. (Hansen) If a measure-preserving dynamical system $(\mathcal{X}, \mathbb{E}, \mu, T)$ is mixing then it is also ergodic.

Lemma 5.8. (Hansen) Let $(\mathcal{X}, \mathbb{E}, \mu, T)$ be a measure-preserving dynamical system. Let \mathbb{D} be an \cap -stable generator for \mathbb{E} . If

$$\mu(A \cap T^{-n}(B)) \to \mu(A)\mu(B) \quad \text{for all } A, B \in \mathbb{D}.$$
 (5.10)

then the system is mixing. (and ergodic)

Lemma 5.9. (Hansen) Let $(\mathcal{X}, \mathbb{E}, \mu)$ be a probability space, and let $T : \mathcal{X} \to \mathcal{X}$ be a measure-preserving map. Let $(\mathcal{Y}, \mathbb{G})$ be another measurable space, and let $S : \mathcal{Y} \to \mathcal{Y}$ be a measurable map. Suppose there is a measurable map $\gamma : \mathcal{X} \to \mathcal{Y}$ such that the following diagram commutes:



Then $(\mathcal{Y}, \mathbb{G}, \gamma(\mu), S)$ is a measure-preserving dynamical system.

Lemma 5.10. (Hansen) Let $(\mathcal{X}, \mathbb{E}, \mu, T)$ and $(\mathcal{Y}, \mathbb{G}, \nu, S)$ be two measure-preserving dynamical systems. Suppose there is a measurable map $\gamma : \mathcal{X} \to \mathcal{Y}$ such that $\nu = \gamma(\mu)$ and such that the diagram in lemma 5.9 commutes. If $(\mathcal{X}, \mathbb{E}, \mu, T)$ is ergodic then $(\mathcal{Y}, \mathbb{G}, \nu, S)$ is also ergodic.

Lemma 5.11. (Hansen) (Maximal Ergodic Lemma) Let $(\mathcal{X}, \mathbb{E}, \mu, T)$ be a measure-preserving dynamical system, and let $f: \mathcal{X} \to \mathbb{R}$ be Borel measurable. If $f \in \mathcal{L}^1(\mu)$ then it holds that

$$\int_{(M_n>0)} f \ d\mu \ge 0 \tag{5.14}$$

where $M_n = \max\{0, S_1, S_2, ..., S_n\}$ from the sequence

$$\Big(f(x),f\circ T(x),f\circ T^2(x),f\circ T^3(x),\ldots\Big)\quad\text{with}\quad S_n=\sum_{i=0}^{n-1}f\circ T^i.$$

Theorem 5.12. (Hansen) (Birkhoff's ergodic theorem) Let $(\mathcal{X}, \mathbb{E}, \mu, T)$ be an ergodic system. For $f \in \mathcal{L}^1(\mu)$ it holds that

$$\frac{1}{n} \sum_{i=0}^{n-1} f \circ T^i \stackrel{\text{a.s.}}{\to} \int f \ d\mu. \tag{5.16}$$

Theorem 5.13. (Hansen) (Ergodic theorem, \mathcal{L}^p -version) Let $(\mathcal{X}, \mathbb{E}, \mu, T)$ be an ergodic system. If $f \in \mathcal{L}^p(\mu)$ for some $p \geq 1$ then it holds that

$$\frac{1}{n} \sum_{i=0}^{n-1} f \circ T^i \stackrel{\mathcal{L}^p}{\to} \int f \ d\mu. \tag{5.21}$$

Lemma 5.14. (Hansen) Let $(\mathcal{X}, \mathbb{E})$ be a measurable space. The measurable finite dimensional product sets in $\mathcal{X}^{\mathbb{N}}$ form an \cap -stable generator for $\mathbb{E}^{\otimes \mathbb{N}}$.

An element of the space $\mathcal{X}^{\mathbb{N}}$ is a countable set of coordinates x_n for $n \in \mathcal{N}$ with $x_n \in \mathcal{X}$. A finite dimensional product set in $\mathcal{X}^{\mathbb{N}}$ is set on the form

$$A_1 \times ... \times A_k \times \mathcal{X} \times \mathcal{X} \times ...$$

where $A_1,...,A_k\subset\mathcal{X}$. We also define the **projection sigma-algebra** $\mathbb{E}^{\otimes\mathbb{N}}$ as $\sigma\left(\left(\hat{X}_n(\mathcal{X}^n)\right)_{n\in\mathbb{N}}\right)$ where $\hat{X}_n(x_1,...,x_n)=x_n$.

Definition 5.15. (Hansen) Let $X_1, X_2, ...$ be a sequence of $(\mathcal{X}, \mathbb{E})$ -valued random variable, defined on a background space (Ω, \mathbb{F}, P) , and let $\mathbb{X} = (X_1, X_2, ...)$ be their bundling. The **distribution** of the process is the image measure $\mathbb{X}(P)$ on $(\mathcal{X}^{\mathbb{N}}, \mathbb{E}^{\otimes \mathbb{N}})$.

Lemma 5.16. (Hansen) Let $\mathbb{X}=(X_1,X_2,...)$ and $\mathbb{Y}=(Y_1,Y_2,...)$ be two (\mathcal{X},\mathbb{E}) -valued stochastic process, defined on a common background space. The two processes \mathbb{X} and \mathbb{Y} have the same distribution if and only if the have the same fidis. This can be checked by showing that

$$P(X_1 \in A_i, ..., X_k \in A_k) = P(Y_1 \in A_1, ..., Y_k \in A_k)$$
(5.25)

for any $k \in \mathbb{N}$ and any choice of $A_1, ..., A_k \in \mathbb{E}$.

Definition 5.18. (Hansen) Let $X_1, X_2, ...$ be a sequence of $(\mathcal{X}, \mathbb{E})$ -valued random variable, defined on a background space (Ω, \mathbb{F}, P) , and let $\mathbb{X} = (X_1, X_2, ...)$ be their bundling. Then we define: 1. The proces \mathbb{X} is **stationary** if the distribution $\mathbb{X}(P)$ is an S-invariant probability on $(\mathbb{R}^{\mathbb{N}}, \mathbb{B}^{\otimes \mathbb{N}})$, 2. The proces \mathbb{X} is **ergodic** if it is stationary and if the dynamical system $(\mathbb{R}^{\mathbb{N}}, \mathbb{B}^{\otimes \mathbb{N}}, \mathbb{X}(P), S)$ is ergodic. 3. The proces \mathbb{X} is **mixing** if it is stationary and if the dynamical system $(\mathbb{R}^{\mathbb{N}}, \mathbb{B}^{\otimes \mathbb{N}}, \mathbb{X}(P), S)$ is mixing. with S being the **shift** map defined as $S(x_1, x_2, ...) = (x_2, x_3, ...)$.

Theorem 5.20. (Hansen) (Khintchine's ergodic theorem) Let $X_1, X_2, ...$ be a stationary and ergodic sequence of real-valued random variables. If $E|X_1| < \infty$ it holds that

$$\frac{1}{n} \sum_{i=1}^{n} X_i \stackrel{\text{a.s.}}{\to} E X_1. \tag{5.27}$$

If $E|X_1|^p < \infty$ for some $p \ge 1$, the convergence is also in \mathcal{L}^p .

Theorem 5.21. (Hansen) (Ergodic transformation theorem) Let $X_1, X_2, ...$ be a sequence of real-valued random variables. For a measurable function $\phi: (\mathbb{R}^{\mathbb{N}}, \mathbb{B}^{\otimes \mathbb{N}}) \to (\mathbb{R}, \mathbb{B})$ we define new real-valued random variables

$$Y_n = \phi(X_n, X_{n+1}, \dots) = \phi \circ S^{n-1}(\mathbb{X})$$
 for $n \in \mathbb{N}$.

If $X_1, X_2, ...$ is stationary and ergodic then $Y_1, Y_2, ...$ is also stationary and ergodic.

Definition 5.23. (Hansen) Let $(X_n)_{n\in\mathbb{Z}}$ be a two-sided sequence of real-valued random variables, defined on a background space (Ω, \mathbb{F}, P) , and let \mathbb{X} be their bundling. 1. The proces \mathbb{X} is **stationary** if the distribution $\mathbb{X}(P)$ is an S-invariant probability on $(\mathbb{R}^{\mathbb{Z}}, \mathbb{B}^{\otimes \mathbb{Z}})$, 2. The proces \mathbb{X} is **ergodic** if it is stationary and if the dynamical system $(\mathbb{R}^{\mathbb{Z}}, \mathbb{B}^{\otimes \mathbb{Z}}, \mathbb{X}(P), S)$ is ergodic. 3. The proces \mathbb{X} is **mixing** if it is stationary and if the dynamical system $(\mathbb{R}^{\mathbb{Z}}, \mathbb{B}^{\otimes \mathbb{Z}}, \mathbb{X}(P), S)$ is mixing.

Theorem 5.25. (Hansen) (Khintchine's ergodic theorem, two-sided version) Let $(X_n)_{n\in\mathbb{Z}}$ be a two-sided sequence of real-valued random variables. If the sequence is stationary and ergodic and if $E|X_1| < \infty$ then it holds that

$$\frac{1}{n} \sum_{i=1}^{n} X_i \stackrel{\text{a.s.}}{\to} E X_1. \tag{5.30}$$

If $E|X_1|^p < \infty$ for some $p \ge 1$, the convergence is also in \mathcal{L}^p .

Theorem 5.26. (Hansen) (Ergodic transformation theorem) Let $(X_n)_{n\in\mathbb{Z}}$ be a two-sided sequence of real-valued random variables. For a measurable function $\phi:(\mathbb{R}^{\mathbb{Z}},\mathbb{B}^{\otimes\mathbb{Z}})\to(\mathbb{R},\mathbb{B})$ we define new real-valued random variables

$$Y_n = \phi \circ S^n(\mathbb{X}) \quad \text{for } n \in \mathbb{Z}.$$

If $(X_n)_{n\in\mathbb{Z}}$ is stationary and ergodic then $(Y_n)_{n\in\mathbb{Z}}$ is also stationary and ergodic.

13.1.3 Weak Convergence

Definition 6.1. (Hansen) A sequence of probability measures $\nu_1, \nu_2, ...$ on (\mathbb{R}, \mathbb{B}) is said to converge weakly to a limit probability measure ν if

$$\int f \ d\nu_n \to \int f \ d\nu \quad \text{ for every } f \in C_b(\mathbb{R})$$
 (6.2)

We write $\nu_n \stackrel{wk}{\rightarrow} \nu$ to denote weak convergence.

Theorem 6.4. (Hansen) (Scheffe's) Let $\nu_1, \nu_2, ...$ and ν be probability measures on (\mathbb{R}, \mathbb{B}) . Assume that for some choice of basic measure μ it holds that $\nu_n = f_n \cdot \mu$ for every n and $\nu = f \cdot \mu$ for suitable density functions f_n and f. If

$$f_n(x) \to f(x)$$
 μ -almost surely

then it holds that $\nu_n \stackrel{wk}{\rightarrow} \nu$.

Lemma 6.8. (Hansen) Let μ and ν be two probability measures on (\mathbb{R}, \mathbb{B}) . If

$$\int f \ d\mu = \int f \ d\nu \quad \text{ for all } f \in C_b(\mathbb{R})$$
 (6.7)

then it holds that $\mu = \nu$.

Theorem 6.9. (Hansen) Let $\nu_1, \nu_2, ...$ be a sequence of probability measures on (\mathbb{R}, \mathbb{B}) and let μ and ν be two extra probability measures. If

$$\nu_n \stackrel{\mathrm{wk}}{\to} \mu$$
 and $\nu_n \stackrel{\mathrm{wk}}{\to} \nu$

then $\mu = \nu$.

Definition 6.10. (Hansen) A sequence of real-valued variables $X_1, X_2, ...,$ defined on a common background space (Ω, \mathbb{F}, P) is said to **converge in distribution** to a limit variable X if

$$\int f(X_n) \ dP \to \int f(X) \ dP \quad \text{ for every } f \in C_b(\mathbb{R}). \tag{6.9}$$

We will write $X_n \stackrel{\mathcal{D}}{\to} X$ to denote convergence in distribution.

Lemma 6.11. (Hansen) Let $X_1, X_2, ...$ and X be real-valued random variables. It holds that

$$X_n \stackrel{\mathrm{P}}{\to} X \quad \Rightarrow \quad X_n \stackrel{\mathcal{D}}{\to} X.$$

Lemma 6.12. (Hansen) Let $X_1, X_2, ...$ be real-valued random variable and let $x_0 \in \mathbb{R}$. It holds that

$$X_n \xrightarrow{\mathcal{D}} x_0 \quad \Rightarrow \quad X_n \xrightarrow{\mathrm{P}} x_0.$$

Theorem 6.13. (Hansen) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) . Let \mathcal{H} be a class of bounded, non-negative and measurable functions with the following approximation property: For $f \in C_b(\mathbb{R})$ with $f \geq 0$ there is a sequence $h_1, h_2, ...$ of \mathcal{H} -functions such that $h_n \nearrow f$. If

$$\int h \ d\nu_n \to \int h \ d\nu \quad \text{for all } h \in \mathcal{H}.$$
 (6.11)

then it holds that $\nu_n \stackrel{wk}{\rightarrow} \nu$.

Theorem 6.14. (Hansen) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) . If

$$\int f \ d\nu_n \to \int f \ d\nu \quad \text{for all } f \in C_c(\mathbb{R})$$
 (6.13)

then it holds that $\nu_n \stackrel{wk}{\to} \nu$.

The class $C_c(\mathbb{R})$ is denoted as the set of all continuous real-valued functions with compact support i.e. there exist a M > 0 such that f(x) = 0 for all |x| > M.

Lemma 6.15. (Hansen) Let ν, ν_1, ν_2, \dots be probability measures on (\mathbb{R}, \mathbb{B}) , and let F, F_1, F_2, \dots be the corresponding distribution functions. If $\nu_n \stackrel{wk}{\to} \nu$ then it holds that

$$F_n(x_0) \to F(x_0),$$

whenever x_0 is a point of continuity for F.

Theorem 6.17. (Hansen) (Helly-Bray) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) , and let $F, F_1, F_2, ...$ be the corresponding distribution functions. It holds that $\nu_n \stackrel{wk}{\to} \nu$ if and only if there is a dense subset $A \subset \mathbb{R}$ such that

$$F_n(x) \to F(x)$$
 for every $x \in A$. (6.16)

Theorem 6.18. (Hansen) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) . Let $F, F_1, F_2, ...$ be the corresponding distribution functions, and let $q, q_1, q_2, ...$ be the corresponding quantile functions. If $\nu_n \stackrel{wk}{\to} \nu$ then it holds that

$$q_n(p) \to q(p)$$

for any $p \in (0,1)$ such that the equation F(x) = p hast at most one solution.

Theorem 6.19. (Hansen) (Skorokhod's representation theorem) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) . If $\nu_n \stackrel{wk}{\to} \nu$ then it is possible to find random variables $X, X_1, X_2, ...$ on a background space (Ω, \mathbb{F}, P) such that

$$X(P) = \nu, \ X_1(P) = \nu_1, \ X_2(P) = \nu_2, \dots$$

and such that $X_n \stackrel{a.s.}{\to} X$.

Corollary 6.20. (Hansen) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) such that $\nu_n \stackrel{wk}{\to} \nu$. Let $h : \mathbb{R} \to \mathbb{R}$ be a bounded and measurable function. If there is a Boral-measurable set $C \subset \mathbb{R}$ such that h is continuous in every point of C and such that $\nu(C) = 1$, then it holds that

$$\int h \ d\nu_n \to \int h \ d\nu. \tag{6.20}$$

Corollary 6.21. (Hansen) (Portmanteau's lemma) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) such that $\nu_n \stackrel{wk}{\to} \nu$. For any open set $G \subset \mathbb{R}$ it holds that

$$\liminf \nu_n(G) \ge \nu(G) \tag{6.21}$$

Definition 6.22. (Hansen) The characteristic function for a probability measure ν on (\mathbb{R}, \mathbb{B}) is the function $\phi : \mathbb{R} \to \mathbb{C}$ given by

$$\phi(\theta) = \int e^{ix\theta} d\nu(x). \quad \text{for } \theta \in \mathbb{R}.$$
 (6.23)

Some useful observations include $|e^{ix\theta}|=1$ hence $\phi(\theta)\leq 1$ for all $\theta\in\mathbb{R}$. We may also split the ϕ into an imaginary part and a real part with Euler's cartesian form

$$\phi(\theta) = \int \cos(x\theta) \ d\nu(x) + i \int \sin(x\theta) \ d\nu(x)$$
 (6.24)

And lastly we have the implication

$$\nu_n \overset{\mathrm{wk}}{\to} \nu \quad \Rightarrow \quad \phi_n(\theta) \to \phi(\theta) \quad \text{for all } \theta \in \mathbb{R}.$$

Furthermore, if $Y = \xi + \sigma X$ and $X \sim \mathcal{N}(0,1)$ we have

$$\phi_V(\theta) = e^{i\xi\theta} e^{-\sigma^2\theta^2/2}.$$

Theorem 6.28. (Hansen) The characteristic function for any probability measure ν on (\mathbb{R}, \mathbb{B}) is uniformly continuous.

Theorem 6.29. (Hansen) Let ν be a probability measure on (\mathbb{R}, \mathbb{B}) . If

$$\int |x|^k \ d\nu(x) < \infty$$

for some $k \in \mathbb{N}$, then the characteristic function ϕ is C^k and it holds that

$$\phi^{(k)}(\theta) = i^k \int x^k e^{i\theta x} \ d\nu(x) \quad \text{for } \theta \in \mathbb{R}.$$
 (6.31)

Definition 6.30. (Hansen) The **convolution** of two probability measures ν and ξ on (\mathbb{R}, \mathbb{B}) is the image measure

$$\nu * \xi = \kappa(\nu \otimes \xi) \tag{6.33}$$

where $\kappa: \mathbb{R}^2 \to \mathbb{R}$ is the addition map $\kappa(x,y) = x + y$.

Theorem 6.31. (Hansen) Let ν and ξ be two probability measures on \mathbb{R} . If $\xi = f \cdot m$, then the convolution $\nu * \xi$ will have a density with respect to m. The density is given as

$$g(x) = \int f(x - y) \ d\nu(y) \quad \text{ for } x \in \mathbb{R}. \tag{6.35}$$

Lemma 6.31. (Hansen) Let ν_1 and ν_2 be two probability measures on (\mathbb{R}, \mathbb{B}) with characteristic functions ϕ_1 and ϕ_2 . The convolution $\nu_1 * \nu_2$ has characteristic function γ given by

$$\gamma(\theta) = \phi_1(\theta)\phi_2(\theta) \quad \text{for } \theta \in \mathbb{R}.$$
 (6.37)

Theorem 6.34. (Hansen) Let $\xi, \nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) . It holds that

$$\nu_n \overset{\text{wk}}{\to} \nu \quad \Rightarrow \quad \nu_n * \xi \overset{\text{wk}}{\to} \nu * \xi.$$

Definition 6.35. (Hansen) A probability measure $\nu = f \cdot m$ on (\mathbb{R}, \mathbb{B}) with density f with respect to Lebesgue measure is of **Polya class** if $f \in C_b(\mathbb{R})$ and if the Fourier transform \hat{f} is m-integrable.

Lemma 6.39. (Hansen) Let ν and ξ be two probability measures on \mathbb{R} . If ξ is of Polya class then the convolution $\nu * \xi$ is also of Polya class.

Theorem 6.40. (Hansen) (Inversion theorem) Let $\nu = f \cdot m$ be a probability measure on (\mathbb{R}, \mathbb{B}) of Polya class. It holds that

$$f(x) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{-i\theta x} \hat{f}(\theta) \ d\theta, \ x \in \mathbb{R}.$$
 (6.39)

Theorem 6.41. (Hansen) Let ν_1 and ν_2 be two probability measures on (\mathbb{R}, \mathbb{B}) with characteristic functions ϕ_1 and ϕ_2 . If

$$\phi_1(\theta) = \phi_2(\theta), \ \forall \theta \in \mathbb{R}$$

then $\nu_1 = \nu_2$.

Lemma 6.42. (Hansen) Let $f : \mathbb{R} \to \mathbb{R}$ be bounded and uniformly continuous. For every $\varepsilon > 0$ there is a probability measure ξ of Polya class with the property that

$$\left| f(x) - \int f(x+y) \ d\xi(y) \right| < \varepsilon, \ \forall x \in \mathbb{R}.$$
 (6.43)

Theorem 6.43. (Hansen) (Continuity theorem) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on (\mathbb{R}, \mathbb{B}) with characteristic functions $\phi, \phi_1, \phi_2, ...$ If

$$\phi_n(\theta) \to \phi(\theta), \ \theta \in \mathbb{R},$$
 (6.45)

then it holds that $\nu_n \stackrel{wk}{\rightarrow} \nu$.

Definition 6.44. (Hansen) A sequence of probability measures $\nu_1, \nu_2, ...$ on $(\mathbb{R}^k, \mathbb{B}_k)$ is said to converge weakly to a limit probability measure ν if

$$\int f(x) \ d\nu_n(x) \to \int f(x) \ d\nu(x), \forall f \in C_b(\mathbb{R}^k). \tag{6.46}$$

We will write $\nu_n \stackrel{wk}{\to} \nu$ to denote weak convergence.

Theorem 6.45. (Hansen) (Continuity theorem) Let $\nu, \nu_1, \nu_2, ...$ be probability measures on $(\mathbb{R}^k, \mathbb{B}_k)$ with characteristic functions $\phi, \phi_1, \phi_2, ...$ If

$$\phi_n(\theta) \to \phi(\theta), \ \theta \in \mathbb{R}^k,$$
 (6.47)

then it holds that $\nu_n \stackrel{wk}{\to} \nu$.

Lemma 6.46. (Hansen) Let \mathbf{X} be an \mathbb{R}^k -valued random variable with characteristic function $\phi_{\mathbf{X}}$, and let \mathbf{Y} be an \mathbb{R}^m -valued random variable with characteristic function $\phi_{\mathbf{Y}}$. If $\mathbf{X} \perp \!\!\! \perp \mathbf{Y}$ then the bundle (\mathbf{X}, \mathbf{Y}) is an \mathbb{R}^{k+m} -valued random variable with

$$\phi_{(\mathbf{X},\mathbf{Y})}(\theta_1,\theta_2) = \phi_{\mathbf{X}}(\theta_1)\phi_{\mathbf{Y}}(\theta_2), \ \theta_1 \in \mathbb{R}^k, \theta_2 \in \mathbb{R}^m. \tag{6.49}$$

Theorem 6.47. (Hansen) (Continuous mapping theorem) Let $\mathbf{X}_1, \mathbf{X}_2, ...$ and \mathbf{X} be random variables with values in \mathbb{R}^k , and let $h: \mathbb{R}^k \to \mathbb{R}^m$ be continuous. If $\mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{X}$, then it holds that $h(\mathbf{X}_n) \overset{\mathcal{D}}{\to} h(\mathbf{X})$.

Theorem 6.48. (Hansen) (Cramer-Wold's device) Let $X_1, X_2, ...$ and X be random variables with values in \mathbb{R}^k . If

$$\mathbf{v}^{\top} \mathbf{X}_n \stackrel{\mathcal{D}}{\to} \mathbf{v}^{\top} \mathbf{X}, \tag{6.51}$$

for any fixed vector $\mathbf{v} \in \mathbb{R}^k$, then it holds that $\mathbf{X}_n \stackrel{\mathcal{D}}{\to} \mathbf{X}$.

Lemma 6.49. (Hansen) Let $X, X_1, X_2, ...$ be random variables with values in \mathbb{R}^k , let $Y_1, Y_2, ...$ be random variables in \mathbb{R}^m , and let y be a vector in \mathbb{R}^m . If it holds that

$$\mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{X}, \quad \ \mathbf{Y}_n \overset{\mathrm{P}}{\to} \mathbf{y}$$

then the bundle $(\mathbf{X}_n, \mathbf{Y}_n)$ in \mathbb{R}^{k+m} will satisfy that

$$(\mathbf{X}_n, \mathbf{Y}_n) \stackrel{\mathcal{D}}{\to} (\mathbf{X}, \mathbf{y}).$$

Corollary 6.50. (Hansen) (Slutsky's lemma) Let $X, X_1, X_2, ...$ and $Y_1, Y_2, ...$ be random variables with values in \mathbb{R}^k . It holds that

$$\mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{X}, \quad \mathbf{Y}_n \overset{\mathrm{P}}{\to} \mathbf{0} \quad \Rightarrow \quad \mathbf{X}_n + \mathbf{Y}_n \overset{\mathcal{D}}{\to} \mathbf{X}.$$

Corollary 6.51. (Hansen) Let $X_1, X_2, ...$ and X be random variables with values in \mathbb{R}^k and let $Y_1, Y_2, ...$ be real-valued random variables. It holds that

$$\mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{X}, \quad Y_n \overset{\mathrm{P}}{\to} \mathbf{1} \quad \Rightarrow \quad Y_n \mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{X}.$$

Corollary 6.52. (Hansen) Let $X_1, X_2, ...$ and X be random variables with values in \mathbb{R}^k and let $Y_1, Y_2, ...$ be real-valued random variables. If

$$\mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{X}, \quad Y_n \overset{\mathrm{P}}{\to} 0 \quad \Rightarrow \quad Y_n \mathbf{X}_n \overset{\mathrm{P}}{\to} \mathbf{0}.$$

Definition 6.53. (Hansen) The \mathbb{R}^k -valued random variable $\mathbf{Z} = (Z_1, ..., Z_k)$ has a multivariate Gaussian distribution if and only if the real-valued random variable $\sum_{j=1}^k c_j Z_j$ has a one-dimensional Gaussian distribution for every choice of $c_1, ..., c_k \in \mathbb{R}$.

Theorem 6.54. (Hansen) Let $\mathbf{Z} = (Z_1, ..., Z_k)$ have a multivariate Gaussian distribution with $E \mathbf{Z} = \xi$ and $V \mathbf{Z} = \Sigma$. Then the characteristic function is

$$\phi_{\mathbf{Z}}(\theta) = e^{i\theta^\top \xi} \exp\left(-\frac{1}{2}\theta^\top \Sigma \theta\right), \ \theta \in \mathbb{R}^k. \tag{6.53}$$

Conversly, if **Z** has characteristic function given by (6.53) for some $\xi \in \mathbb{R}^k$ and some symmetric, positive semi-definite $k \times k$ matrix Σ then **Z** has a multivariate Gaussian distribution with $E \mathbf{Z} = \xi$ and $V \mathbf{Z} = \Sigma$.

Corollary 6.55. (Hansen) Let **Z** be an \mathbb{R}^k -valued random variable, let $\mathbf{a} \in \mathbb{R}^m$ and let B be an $m \times k$ matrix. It holds that

$$\mathbf{Z} \sim \mathcal{N}(\xi, \Sigma) \quad \Rightarrow \quad \mathbf{a} + B\mathbf{Z} \sim \mathcal{N}\left(\mathbf{a} + B\xi, B\Sigma B^{\top}\right).$$

Lemma 6.56. (Hansen) Let $\mathbf{X} = (X_1, ..., X_k)$ and $\mathbf{Y} = (Y_1, ..., Y_m)$ be random variables with values in \mathbb{R}^k respectively \mathbb{R}^m . If both \mathbf{X} and \mathbf{Y} have multivariate Gaussian distributions and if \mathbf{X} and \mathbf{Y} are independent, then the \mathbb{R}^{k+m} -valued bundle (\mathbf{X}, \mathbf{Y}) has a multivariate Gaussian distribution.

Lemma 6.58. (Hansen) Let $\mathbf{X} = (X_1, ..., X_k)$ and $\mathbf{Y} = (Y_1, ..., Y_m)$ be random variables with values in \mathbb{R}^k respectively \mathbb{R}^m . If the \mathbb{R}^{k+m} -valued bundle (\mathbf{X}, \mathbf{Y}) has a multivariate Gaussian distribution, and if

$$Cov(X_i, Y_l) = 0, \ \forall j \text{ and } l$$

then $X \perp \!\!\!\perp Y$.

Definition 6.59. (Hansen) Let $\mathbf{X}_1, \mathbf{X}_2, ...$ be \mathbb{R}^k -valued random variables. Let $\xi \in \mathbb{R}^k$ be a vector and let Σ be a symmetric, positive semi-definite $k \times k$ matrix. We sat that \mathbf{X}_n has an asymptotic normal distribution with parameters $(\xi, \frac{1}{n}\Sigma)$, written

$$\mathbf{X}_n \overset{\text{a.s.}}{\sim} \mathcal{N}\left(\xi, \frac{1}{n}\Sigma\right),$$

if it holds that

$$\sqrt{n}(\mathbf{X}_n - \xi) \stackrel{\mathcal{D}}{\to} \mathcal{N}(0, \Sigma).$$

Lemma 6.60. (Hansen) Let $\mathbf{X}_1, \mathbf{X}_2, ...$ and \mathbf{Y} be random variables with values in \mathbb{R}^k . If it holds that $\mathbf{X}_n \stackrel{a.s.}{\sim} \mathcal{N}(\xi, \frac{1}{n}\Sigma)$ then it follows that $\mathbf{X}_n \stackrel{P}{\to} \xi$.

Lemma 6.61. (Hansen) Let $\mathbf{X}_1, \mathbf{X}_2, ...$ and \mathbf{Y} be \mathbb{R}^k -valued random variables, and assume that $\sqrt{n}\mathbf{X}_n \overset{\mathcal{D}}{\to} \mathbf{Y}$. Let $g: \mathbb{R}^k \to \mathbb{R}^m$ be a measurable map. Assume that $g(\mathbf{0}) = \mathbf{0}$ and that g is differentiable in $\mathbf{0}$ with deriviate $Dg(\mathbf{0}) = A$. Then it holds that

$$\sqrt{n}g(\mathbf{X}_n) \stackrel{\mathcal{D}}{\to} A \mathbf{Y}.$$

Lemma 6.62. (Hansen) (Delta method) Let $\mathbf{X}_1, \mathbf{X}_2, ...$ be \mathbb{R}^k -valued random variables, and let $f : \mathbb{R}^k \to \mathbb{R}^m$ be measurable. If f is differentiable in ξ , then it holds that

$$\mathbf{X}_n \overset{\text{a.s.}}{\sim} \mathcal{N}\left(\xi, \frac{1}{n}\Sigma\right) \quad \Rightarrow \quad f(\mathbf{X}_n) \overset{\text{a.s.}}{\sim} \mathcal{N}\left(f(\xi), \frac{1}{n}Df(\xi)\Sigma Df(\xi)^\top\right).$$

13.1.4 Central Limit Theorems

Lemma 7.1. (Hansen) Let $z_1,...,z_n$ and $w_1,...,w_n$ be complex numbers. If $|z_i| \le 1$ and $|w_i| \le 1$ for all i = 1,...,n then it holds that

$$\left| \prod_{i=1}^{n} z_i - \prod_{i=1}^{n} w_i \right| \le \sum_{i=1}^{n} |z_i - w_i|. \tag{7.1}$$

Theorem 7.2. (Hansen) (Basic CLT) Let $X_1, X_2, ...$ be independent and identically distributed real-valued random variables. Assume that $E(X_1) = 0$ and $E(X_1) = 1$. Then it holds that

$$\frac{1}{\sqrt{n}} \sum_{i=1}^{n} X_i \xrightarrow{\mathcal{D}} \mathcal{N}(0,1) \tag{7.3}$$

Theorem 7.3. (Hansen) (Laplace's CLT) Let $X_1, X_2, ...$ be independent and identically distributed real-valued random variables with $E X_1^2 < \infty$. It holds that

$$\frac{1}{n} \sum_{i=1}^{n} X_i \stackrel{\mathcal{D}}{\to} \mathcal{N}\left(E \ X_1, \frac{V \ X_1}{n}\right) \tag{7.4}$$

Theorem 7.7. (Hansen) (Laplace's CLT, multivariate version) Let $\mathbf{X}_1, \mathbf{X}_2, ...$ be independent and identically distributed random variables with values in \mathbb{R}^k . Assume that $E|\mathbf{X}_1|^2 < \infty$. It holds that

$$\frac{1}{n} \sum_{i=1}^{n} \mathbf{X}_{i} \stackrel{\mathcal{D}}{\to} \mathcal{N}\left(E \ \mathbf{X}_{1}, \frac{1}{n} V \ \mathbf{X}_{1}\right) \tag{7.7}$$

Definition 7.10. (Hansen) Let (X_{nm}) be a centralized array of real-valued random variables. a. The array satisfies the vanishing variance condition if

$$\max_{m=1,\dots,n} E X_{nm}^2 \to 0. \tag{7.8}$$

b. The array satisfies Lindeberg's condition if

$$\forall c > 0: \qquad \sum_{m=1}^{n} \int_{(|X_{nm}| > c)} X_{nm}^2 \ dP \to 0. \tag{7.9}$$

c. The array satisfies **Lyapounov's condition** of order $\alpha > 2$ if

$$\sum_{m=1}^{n} E |X_{nm}|^{\alpha} \to 0. \tag{7.10}$$

Lemma 7.11. (Hansen) Lyapounov's condition of order $\alpha > 2$ implies Lindeberg's condition.

Lemma 7.12. (Hansen) Lindeberg's condition implies the vanishing variance condition.

Theorem 7.14. (Hansen) (Lindeberg's CLT) Let (X_{nm}) be a centralized array of real-valued random variables with $E X_{nm}^2 < \infty$. Assume that the array satisfies that

$$E X_{nm} = 0, \ \forall n, m,$$

and that

$$\sum_{m=1}^{n} E X_{nm}^{2} = 1. (7.13)$$

Assume that the array has independence within rows i.e. $X_{i1} \perp \!\!\! \perp ... \perp \!\!\! \perp X_{ii}$ for all i=1,...,n and satisfies Lindeberg's condition. Then it holds that

$$\sum_{m=1}^n X_{nm} \overset{\mathcal{D}}{\to} \mathcal{N}(0,1).$$

Theorem 7.19. (Hansen) (Lindeberg's CLT, multivariate version) Let (\mathbf{X}_{nm}) be a triangular array of random variables with values in \mathbb{R}^k with $E |\mathbf{X}_{nm}|^2 < \infty$ for all n, m. Assume that

$$E \mathbf{X}_{nm} = \mathbf{0}, \ \forall n, m$$

and

$$\sum_{m=1}^{n} V \mathbf{X}_{nm} \to \Sigma$$

for a fixed $k \times k$ matrix Σ . Assume that the array has independence within rows, and assume that the associated real-valued array $(|\mathbf{X}_{nm}|)$ satisfies Lindeberg's condition. Then it holds that

$$\sum_{m=1}^n \mathbf{X}_{nm} \xrightarrow{\mathcal{D}} \mathcal{N}(\mathbf{0}, \Sigma)$$

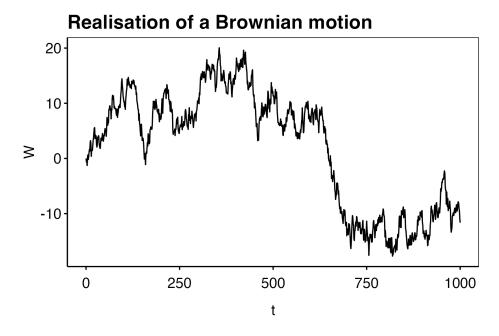
Chapter 14

Continuous Time Stochastic Processes

14.1 Brownian Motion

Definition 4.1. (Bjork) A stochastic process W is called a Brownian motion or Wiener process if the following conditions hold

- 1. $W_0 = 0$.
- 2. The process W has independent increments, i.e. if $r < s \le t < u$ then $W_u W_t$ and $W_s W_r$ are independent random variables.
- 3. For s < t the random variable $W_t W_s$ has the Gaussian distribution $\mathcal{N}(0, t s)$.
- 4. W has continuous trajectories i.e. $s \mapsto W(s; \omega)$ i continuous for all $\omega \in \Omega$.



As one can see from the simulated sample path on the right, the Brownian motion is rather irratic. In fact, the process varies infinitely on any interval with length greater than 0. This gives some of the characteristics of the process including that: W is continuous and W is non-differential everywhere. This irratic behaviour is summed up in the theorem.

Theorem 4.2. (Bjork) A Brownian motions trajectory $t \mapsto W_t$ is with probability one nowhere differential,

 $and\ it\ has\ locally\ infinite\ total\ variation.$

14.2 Filtration

Filtrations is widely used in stochastic processes, as they allow for the concept of knowledge/information. This is useful when considering mean-values of future states but in an increasing information setting. For this we introduce the term adapted processes.

Definition B.17. (Bjork) (Adapted process) Let $(\mathcal{F}_t)_{t\geq 0}$ be a filtration on the probability space $(\mathcal{F}_t)_{t\geq 0}$. Furthermore, let $(X_t)_{t\geq 0}$ be a stochastic process on the same space. We say that X_t is adapted to the filtration \mathbf{F} if

$$X_t$$
 is \mathcal{F}_t – measurable, $\forall t \geq 0$.

Obviously, we may introduce the **natural filtration** \mathcal{F}_t^X given by the tragetory of the process X_t :

$$\mathcal{F}_t^X = \sigma(\{X_s, \ s \le t\}).$$

Indeed, X_t is adapted to this filtration.

14.3 Martingale

Definition. Let M_t be a stochastic process defined on a background space (Ω, \mathcal{F}, P) . Let $(\mathcal{F}_t)_{t\geq 0}$ be a filtration. If M_t is adapted to the filtration \mathcal{F}_t , $E|M_t|<\infty$ and

$$E[M_t|\mathcal{F}_s] = M_s, \qquad P-\text{a.s.}$$

holds for any t > s we say that M_t is a martingale (F-martingale). If the above has $\leq or \geq we$ say that M_t is either a submartingale or supermartingale respectively.

Naturally, this defintions may easily be extended to discrete models and we have the trivial equality:

$$E[M_t - M_s \mid \mathcal{F}_s] = 0.$$

Martingales is useful, when proofing probalistic statements as the posses tractable properties. A useful technique often include the construction of the martingale

$$M_t = E[X \mid \mathcal{F}_t].$$

Chapter 15

Stochastic calculus

15.1 Stochastic Integrals

We want to formulate financial markets in continuous time and the most elegant theory is obtained from processes that can be defined in terms of **stochastic differential equations** or in other words by their dynamics. We may call them **diffusion processes**, as they may be approximated by a stochastic difference equation:

$$X_{t+\Delta t} - X_t = \mu(t, X_t) \Delta t + \sigma(t, X_t) Z_t. \tag{4.1}$$

Above Z_t is a normally distributed random variable (a disturbance). In this formulation we say that S_t is driven by two forces: on one hand a locally deterministic velocity or drift $\mu(t, X_t)$ and on the other hand a Gaussian term amplified by the deterministic factor $\sigma(t, X_t)$.

15.1.1 Information

We consider a primary process X_t and we introduce the notion of information generated by X_t in terms of the natural filtration. The idea can be summed up in the following definition.

Definition 4.3. (Bjork) The symbol $\mathcal{F}_t^X \subseteq \mathcal{F}$ denotes "the information generated by X_t on the interval [0,t]", or alternatively "what has happened to X_t over ther interval [0,t]".

1. If, based upon observations of the trajectory $\{X_s; 0 \le s \le t\}$, it is possible to decide whether a given event A has occurred or not, then we write this as

$$A \in \mathcal{F}_t^X$$

or say that "A is \mathcal{F}_t^X -measurable".

2. If the value of a given random variable Z can be completely determined given observations of the tragectory $\{X_s; 0 \le s \le t\}$, then we also write

$$Z \in \mathcal{F}_t^X$$
. (Z is \mathcal{F}_t^X -measurable)

3. If Y is a stochastic process such that we have

$$Y_t \in \mathcal{F}_t^X$$

for all $t \geq 0$ then we say that Y_t is **adapted** to the **filtration** $\{\mathcal{F}^X_t\}_{t\geq 0}$. For brevity of notation, we will sometimes write the filtration as $\{\mathcal{F}^X_t\}_{t\geq 0} = \mathbf{F}$.

15.1.2 Stochastic Integrals

We will now formulate the theory of stochastic integrals, that is, processes written in terms of stochastic processes with stochastic integrator and/or stochastic integrant. We will consider some given standard Brownian motion W_t and another stochastic process X_t . We need som integrability condition on X_t in order to do the calculations. We therefore determine a selection of suitable stochastic processes X must be contained in.

Definition 4.4. (Bjork) Let X_t be a stochastic process, then

i. We say that X_t belongs to the class $\mathcal{L}^2[a,b]$ if X_t is adapted to the filtration \mathcal{F}_t^X and the following holds

$$\int_{a}^{b} E[X_{s}^{2}] \ ds < \infty$$

ii. We say that X_t belongs to the class \mathcal{L}^2 if $X_t \in \mathcal{L}^2[0,t]$ for all t > 0.

We now want to define what we mean by

$$\int_{a}^{b} X_{t} \ dW_{s}$$

for some process $X_t \in \mathcal{L}^2$. We see that a way to go about this problem is to start by defining the concept for a simple stochastic process X_t . By *simple* we mean a process X_t that is constant on between some deterministic points in time $a = t_0 < t_1 < \dots < t_n = b$. In that case we may define the integral as

$$\int_{a}^{b} X_{s} \ dW_{s} = \sum_{k=0}^{n-1} X_{t_{k}} [W_{t_{k+1}} - W_{t_{k}}]. \tag{4.8}$$

In the more general setting we may follow the following approach:

1. Approximate X with a sequence $\{X^n\}_{n\in\mathbb{N}}$ of simple processes such that the following convergence criteria hold

$$\int_{a}^{b} E[(X_{s}^{n}-X_{s})^{2}] ds \rightarrow 0, \ n \rightarrow \infty$$

- 2. For each n the integral $\int_a^b X_s^n \ dW_s := Z^n$ is well defined and it is possible to prove, using DCT, that a variable Z exists such that $Z^n \to Z$ that is in L^2 .
- 3. We now define the stochastic integral by the limit

$$\int_{a}^{b} X_s \ dW_s = \lim_{n \to \infty} \int_{a}^{b} X_s^n \ dW_s. \tag{4.9}$$

Obviously the hardest step is finding the processes X^n . This stochastic har some properties we will use.

Proposition 4.5. (Bjork) Let $X_t \in \mathcal{L}^2$, then

$$E\left[\int_{a}^{b} X_{s} \ dW_{s}\right] = 0. \tag{4.12}$$

$$E\left[\left(\int_a^b X_s \ dW_s\right)^2\right] = \int_a^b E[X_s^2] \ dW_s. \tag{4.13}$$

$$\int_{a}^{b} X_{s} \ dW_{s} \text{ is } \mathcal{F}_{b}^{W}\text{-measurable.}$$

$$\tag{4.14}$$

15.1.3 Martingales

Definition 4.7. (Bjork) Let M_t be a stochastic process defined on a background space (Ω, \mathcal{F}, P) . Let $(\mathcal{F}_t)_{t\geq 0}$ be a filtration. If M_t is adapted to the filtration \mathcal{F}_t , $E|M_t| < \infty$ and

$$E[M_t|\mathcal{F}_s] = M_s, \qquad P - \text{a.s.}$$

holds for any t > s we say that M_t is a martingale (F-martingale). If the above has $\geq or \leq we$ say that M_t is either a **submartingale** or **supermartingale** respectively.

Proposition 4.8. (Bjork) For any process $X_t \in \mathcal{L}^2[s,t]$ the following hold:

$$E\left[\left.\int_{s}^{t}X_{s}\ dW_{s}\right|\mathcal{F}_{s}^{W}\right]=0$$

Corollary 4.9. (Bjork) For any process $X_t \in \mathcal{L}^2$ the process

$$M_t = \int_s^t X_s \ dW_s,$$

is an (\mathcal{F}_t^W) -martingale. In other words, modulo an integrability condition, **every stochastic integral is a** martingale.

Lemma 4.10. (Bjork) Let M_t be a stochastic process with stochastic differential, then M_t is a martingale if and only if the stochastic differential has the form $dM_t = X_t dW_t$ i.e. M_t as no dt-term.

15.1.4 Stochastic Calculus and the Ito Formula

Given this breif introduction to stochastic integrals we may formulate som simple calculus revolving around Ito's formula. We consider the stochastic process X_t and we suppose that there exist a real number X_0 and adapted processes μ and σ wrt. \mathcal{F}^W_t such that for all $t \geq 0$ we have

$$X_{t} = X_{0} + \int_{0}^{t} \mu_{s} \, ds + \int_{0}^{t} \sigma_{s} \, dW_{s}, \tag{4.16}$$

where W_t is a standard Brownian motion. We know from earlier courses that the above may be written in terms of the dynamics (pure notation):

$$\begin{cases} dX_t = \mu_t \ dt + \sigma_t \ dW_t, \\ X_0 = X_0. \end{cases}$$
 (4.17/18)

Here we interest the above as X_t has boundary condition X_0 and evolves with a drift μ_t dt amplified and distorted by the drift σ_t dW_t . We say that X_t has **stochastic differential** dX_t and initial condition X_0 .

We want to understand how transformation of such an integral behaves and therefore we introduce some calculus which will tell how for instance $f(t, X_t)$ (for some $C^{1,2}$ -function) behaves. This insight is given by the important Ito's formula.

Thoerem 4.11. (Bjork) (Ito's formula, one-dimensional) Assume that the process X has a stochastic differential form given by

$$dX_t = \mu_t \ dt + \sigma_t \ dW_t, \tag{4.28}$$

where μ and σ are adapted processes, and let $f: \mathbb{R}_+ \times \mathbb{R} \to \mathbb{R}$ be a $C^{1,2}$ -function. Define the process Z by $Z_t = f(t, X_t)$. Then Z has stochastic differential given by

$$df(t, X_t) = \left(\frac{\partial f}{\partial t}(t, X_t) + \mu_t \frac{\partial f}{\partial x}(t, X_t) + \frac{1}{2}\sigma_t^2 \frac{\partial^2 f}{\partial x^2}(t, X_t)\right) dt + \sigma_t \frac{\partial f}{\partial x}(t, X_t) dW_t. \tag{4.29}$$

Proposition 4.12. (Bjork) (Ito's formula, one-dimensional) With assumptions as in theorem 4.11, df is given by

$$df = \frac{\partial f}{\partial t} dt + \frac{\partial f}{\partial x} dX_t + \frac{1}{2} \frac{\partial^2 f}{\partial x^2} (dX_t)^2, \tag{4.31}$$

where we use the following table

$$\begin{cases} (dt)^2 = 0, \\ dt \cdot dW_t = 0, \\ (dW_t)^2 = dt. \end{cases}$$

Lemma 4.18. (Bjork) Let $\sigma(t)$ be deterministic function of time and define the process X by

$$X_t = \int_0^t \sigma(s) \ dW_s. \tag{4.37}$$

Then

$$X_t \sim \mathcal{N}\left(0, \int_0^t \sigma^2(s) \ ds\right).$$

15.1.5 The multidimensional Ito Formula

Consider a vector process $X = (X^1, ..., X^n)^{\top}$ where each component X^i has stochastic differential

$$dX_t^i = \mu_t^i \ dt + \sum_{j=1}^d \sigma_t^{ij} \ dW_t^j$$

where $W^1,...,W^d$ is independent Brownian motions. Then we have respectively the drift vector process μ_t in n dimensions, the vector Brownian motion in d dimensions and a $n \times d$ -dimensional **diffusion matrix** σ_t given as below

$$\mu_t = \begin{bmatrix} \mu_t^1 \\ \vdots \\ \mu_t^n \end{bmatrix}, \quad W_t = \begin{bmatrix} W_t^1 \\ \vdots \\ W_t^d \end{bmatrix}, \quad \sigma_t = \begin{bmatrix} \sigma_t^{11} & \cdots & \sigma_t^{1d} \\ \vdots & \ddots & \vdots \\ \sigma_t^{n1} & \cdots & \sigma_t^{nd} \end{bmatrix}.$$

Given this we may write the dynamics of X as

$$dX_t = \mu_t \ dt + \sigma_t \ dW_t \in \mathbb{R}^n.$$

Consider now a function $f: \mathbb{R}_+ \times \mathbb{R}^n \to \mathbb{R}$ which is a $C^{1,2}$ -mapping. We want to study the dynamics of the process

$$Z_t = f(t, X_t).$$

The dynamics is given in the multidimensional version of Ito's formula.

Thoerem 4.19. (Bjork) (Ito's formula, multi-dimensional) Let X be given as above. Then the following holds:

• The process $f(t, X_t)$ has the stochastisc differential given by

$$df(t,X_t) = \left(\frac{\partial f}{\partial t}(t,X_t) + \sum_{i=1}^n \mu_t^i \frac{\partial f}{\partial x^i}(t,X_t) + \frac{1}{2} \sum_{i,j=1}^n C_t^{ij} \frac{\partial^2 f}{\partial x^i \partial x^j}(t,X_t)\right) \ dt + \sum_{i=1}^n \sigma_t^i \frac{\partial f}{\partial x^i}(t,X_t) \ dW_t.$$

Here the row vector σ_t^i is the i'th row of the matrix σ_t and the matrix C is defined by $C = \sigma \sigma^{\top}$.

• Alternatively, the differential is given by the formula

$$df(t,X_t) = \frac{\partial f}{\partial t}(t,X_t) \ dt + \sum_{i=1}^n \frac{\partial f}{\partial x^i}(t,X_t) \ dX_t^i + \frac{1}{2} \sum_{i,j=1}^n \frac{\partial^2 f}{\partial x^i \partial x^j}(t,X_t) \ dX_t^i dX_t^j,$$

with the formal multiplication table

$$\begin{cases} (dt)^2 = 0, \\ dt \cdot dW_t^i = 0, & i = 1, ..., d, \\ (dW_t^i)^2 = dt, & i = 1, ..., d, \\ dW_t^i \cdot dW_t^i = 0, & i \neq j. \end{cases}$$

Obviously, one can write the differential in Ito's formula in many other ways including a matrix-wise version using the Hessian matrix $H_{ij} = \frac{\partial^2 f}{\partial x^i \partial x^j}$.

15.1.6 Correlated Brownian motions

In the previous section the d-dimensional Brownian was assumed to have independent Brownian motions. However we may instead consider a variation where we have some dependence between the Brownian motions.

This section has not been finished.

15.2 Discrete Stochastic Integrals

This section has not been finished.

15.3 Stochastic Differential Equations

We start the chapter by formalising some used objects. We consider the following objects.

- M(n,d) denotes the class of $n \times d$ -matrices.
- W is a d-dimensional Brownian motion
- μ is a \mathbb{R}^n -valued function with arguments (t, X_t) with X_t being a n-dimensional stochastic process.
- σ a M(n,d)-valued function with arguments as in μ .
- x_0 a \mathbb{R}^n -valued vector.

We want then to understand when the following has a solution

$$dX_t = \mu(t, X_t) \ dt + \sigma(t, X_t) \ dW_t, \quad X_0 = x_0. \tag{5.1/2}$$

We call such an equation the **stochastic differential equation** or simply SDE. We know that the above is loosely notation for the integral form as

$$X_{t} = x_{0} + \int_{0}^{t} \mu(s, X_{s}) ds + \int_{0}^{t} \sigma(s, X_{s}) dW_{s}, \tag{5.3}$$

for all $t \ge 0$. The following proposition tells us when an solution exist to the problem above. In the below $\|\cdot\|$ is usual euclidian norm

$$||x|| = \sqrt{\sum_{i=1}^{n} x_i^2}.$$

Proposition 5.1. (Bjork) Suppose that there exists a constant K such that the following conditions are satisfied for all x, y and t.

$$\|\mu(t,x) - \mu(t,y)\| \le K\|x - y\|,\tag{5.6}$$

$$\|\sigma(t,x) - \sigma(t,y)\| \le K\|x - y\|,$$
 (5.7)

$$\|\mu(t,x)\| + \|\sigma(t,x)\| \le K(1+\|x\|). \tag{5.8}$$

Then there exists a unique solution to the SDE above. Furthermore, the solution has the properties

- 1. X is \mathcal{F}_t^W -adapted.
- 2. X has continuous trajectories.
- 3. X is a Markov process.
- 4. There exists a constant C such that

$$E[\|X_t\|^2] \le Ce^{Ct}(1 + \|x_0\|^2). \tag{5.9}$$

In general the solution to an SDE is so complicated, that it in practical terms is unsolvable and may only be approximated on a finely subdividet grid as jumps. There does however exist som nontrivial cases where we may infer a analytical solution. One is the rather important **Geometric Brownian motion**.

Proposition 5.2. (Bjork) Consider the SDE

$$dX_t = \alpha X_t \ dt + \sigma X_t \ dW_t, \tag{5.13}$$

with $X_0 = x_0$. Then the solution is given as

$$X_t = x_0 \cdot \exp\left\{\left(\alpha - \frac{\sigma^2}{2}\right)t + \sigma W_t\right\}. \tag{5.15}$$

The expected value of X is given as $E[X_t] = x_0 e^{\alpha t}$ (eq. 5.16).

One other generalisation that is analytically solvable is the Linear SDE.

Proposition 5.3. (Bjork) Consider the SDE

$$dX_t = (AX_t + b_t) dt + \sigma_t dW_t, (5.19)$$

with $X_0=x_0$ and $A\in M(n,n)$ and b_t being a real-valued function. Then the solution is given as

$$X_t = e^{At}x_0 + \int_0^t e^{A(t-s)}b_s \ ds + \int_0^t e^{A(t-s)}\sigma_s \ dW_s. \tag{5.20}$$

Where we define the exponential of a matrix as below

$$e^{At} = \sum_{k=0}^{\infty} A^k \frac{1}{k!} t^k.$$

In general with the SDE we have a partial differential operator \mathcal{A} called the **infinitesimal operator** of X which has some interesting analytical properties regarding X.

Definition 5.4. (Bjork) Consider the SDE

$$dX_t = \mu(t, X_t) dt + \sigma(t, X_t) dW_t. \tag{5.21}$$

The partial differential operator \mathcal{A} is defined, for any function $h \in C^2(\mathbb{R}^n)$, by

$$\mathcal{A}h(t,x) = \sum_{i=1}^n \mu_i(t,x) \frac{\partial h}{\partial x_i}(x) + \frac{1}{2} \sum_{i,j=1}^n (\sigma(t,x)\sigma(t,x)^\top)_{ij} \frac{\partial^2 h}{\partial x_i \partial x_j}(x).$$

We see that in terms of Ito's formula the operator is included as such

$$df(t,X_t) = \left\{ \frac{\partial f}{\partial t}(t,X_t) + \mathcal{A}f(t,x) \right\} \ dt + [\nabla_x f](t,X_t) \sigma(t,X_t) \ dW_t,$$

where ∇_x is the gradient for function $h \in C^1(\mathbb{R}^n)$ as

$$\nabla_x h(x) = \left[\frac{\partial h}{\partial x_1}(x),...,\frac{\partial h}{\partial x_n}(x)\right].$$

15.4 Partial differential equations

Proposition 5.5. (Bjork) (Feynmann-Kac) Assume that F is a solution to the boundary value problem

$$\frac{\partial F}{\partial t}(t,x) + \mu(t,x) \frac{\partial F}{\partial x}(x,t) + \frac{1}{2}\sigma^2(t,x) \frac{\partial^2 F}{\partial x^2}(t,x) = 0,$$

with boundary condition $F(T,x) = \Phi(x)$. Assume furthermore that the process

$$\sigma(s,X_s)\frac{\partial F}{\partial x}(s,X_s)\in\mathcal{L}^2$$

as per definition 4.4, where X is defined below. Then F has the representation

$$F(t,x) = E_{t,x}[\Phi(X_T)] = E[\Phi(X_T) \mid X_t = x], \tag{5.29}$$

where X satisfies the SDE

$$dX_s = \mu(s, X_s) ds + \sigma(s, X_s) dW_s, \tag{5.30}$$

with boundary condition $X_t = x$.

Proposition 5.6. (Bjork) (Feynmann-Kac) Assume that F is a solution to the boundary value problem

$$\frac{\partial F}{\partial t}(t,x) + \mu(t,x)\frac{\partial F}{\partial x}(x,t) + \frac{1}{2}\sigma^2(t,x)\frac{\partial^2 F}{\partial x^2}(t,x) - rF(t,x) = 0, \tag{5.34}$$

with boundary condition $F(T,x) = \Phi(x)$. Assume furthermore that the process

$$e^{-rs}\sigma(s,X_s)\frac{\partial F}{\partial x}(s,X_s)\in\mathcal{L}^2$$

as per definition 4.4, where X is defined below. Then F has the representation

$$F(t,x) = e^{-r(T-t)}E_{t,x}[\Phi(X_T)] = e^{-r(T-t)}E[\Phi(X_T) \mid X_t = x], \tag{5.36}$$

where X satisfies the SDE

$$dX_s = \mu(s, X_s) ds + \sigma(s, X_s) dW_s, \tag{5.37}$$

with boundary condition $X_t = x$.

Proposition 5.8. (Bjork) (Feynmann-Kac) Assume that F is a solution to the boundary value problem

$$\frac{\partial F}{\partial t}(t,x) + \sum_{i=1}^n \mu_i(t,x) \frac{\partial F}{\partial x}(x,t) + \frac{1}{2} \sum_{i,j=1}^n C_{ij}(t,x) \frac{\partial^2 F}{\partial x^2}(t,x) - rF(t,x) = 0,$$

with boundary condition $F(T,x) = \Phi(x)$ and $C_{ij} = \sigma \sigma^{\top}$. Assume furthermore that the process

$$e^{-rs}\sum_{i=1}^n\sigma_i(s,X_s)\frac{\partial F}{\partial x}(s,X_s)\in\mathcal{L}^2$$

as per definition 4.4, where X is defined below. Then F has the representation

$$F(t,x) = e^{-r(T-t)} E_{t,x}[\Phi(X_T)], \tag{5.39}$$

where X satisfies the SDE

$$dX_s = \mu(s, X_s) ds + \sigma(s, X_s) dW_s, \tag{5.40}$$

with boundary condition $X_t = x$.

Proposition 5.9. (Bjork) Consider as given a vector process X with generator A, and a function F(t,x). Then, modulo some integrability condition, the following hold:

• The process $F(t, X_t)$ is a martingale relative to the filtration \mathcal{F}^X if and only if F satisfies the PDE

$$\frac{\partial F}{\partial t} + \mathcal{A}F = 0.$$

• The process $F(t, X_t)$ is a martingale relative to the filtration \mathcal{F}^X if and only if, for every (t, x) and $T \geq t$, we have

$$F(t,x) = E_{t,x}[F(T,X_T)].$$

15.5 The Product Integral

Consider the (stochastic) function $Y : \mathbb{R} \to \mathbb{R}^{n \times n}$, where $\mathbb{R}^{n \times n}$ is the set of all $n \times n$ real-valued matrices, that is Y has the matrix-representation

$$Y(t) = \begin{bmatrix} Y_{11}(t) & \cdots & Y_{1n}(t) \\ Y_{21}(t) & \cdots & Y_{2n}(t) \\ \vdots & \ddots & \vdots \\ Y_{n1}(t) & \cdots & Y_{nn}(t) \end{bmatrix}.$$

Assume furthermore that for each coordinate function $Y_{ij}(t)$ the equation

$$\frac{dY_{ij}}{dt}(t) = Y_{i1}(t)A_{1j}(t) + \dots + Y_{in}(t)A_{nj}(t), \qquad Y_{ij}(s) = C_{ij},$$

is satisfied. That is on matrix form the linear system of differential equation

$$\frac{dY}{dt}(t) = Y(t)A(t), \qquad Y(s) = C,$$

for some function $A: \mathbb{R} \to \mathbb{R}^{n \times n}$ and initial condition $C \in \mathbb{R}^{n \times n}$. If A is continuous we know that Y is well-defined and absolutely continuous. We may then state the following theorem regarding uniqueness and existence of such a function above.

Theorem 1.1. (Bladt) (Uniqueness) Consider the homogeneous system of linear differential equations

$$\mathbf{Y}'(t) = \mathbf{Y}(t)\mathbf{A}(t), \qquad \mathbf{Y}(s) = \mathbf{C}, \tag{1}$$

where $\mathbf{Y}(t)$, $\mathbf{A}(t)$ and \mathbf{C} are $n \times n$ -matrices and $\mathbf{A}(t)$ is continuous on [s,t]. Then (1) has at most one solution.

Theorem 1.2. (Bladt) (Existence) The matrix function

$$\mathbf{Y}(t) = \sum_{k=0}^{\infty} \mathbf{Y}_k(t), \tag{3}$$

converges uniformly and absolutely on finite intervals, and solves the differential equation

$$\mathbf{Y}'(t) = \mathbf{Y}(t)\mathbf{A}(t), \quad \mathbf{Y}(s) = \mathbf{C}.$$

Now, we know that for any system as in (1) with a continuous function $\mathbf{A}(t)$ we can always construct the solution as some converging series as per theorem 1.2 then theorem 1.1 gives that the solution is unique. We can then with a piece of mind define a symbol for such a solution, without care for the *exact* solution.

Definition 1.3. (Bladt) (The Product Integral) For any continuous matrix function $\mathbf{A}(t)$ we define the product integral as

$$\prod_{s}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx)$$

as the unique solution $\mathbf{Y}(t)$ to

$$\mathbf{Y}'(t) = \mathbf{Y}(t)\mathbf{A}(t), \quad \mathbf{Y}(s) = \mathbf{I}.$$

From a simple integral argument we may construct a converging series not containing \mathbf{Y} itself. We see that

$$\begin{split} \prod_{s}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx) &= \mathbf{I} + \int_{s}^{t} \mathbf{Y}'(x) \ dx = \mathbf{I} + \int_{s}^{t} \mathbf{Y}(x) \mathbf{A}(x) \ dx \\ &= \mathbf{I} + \int_{s}^{t} \left[\mathbf{I} + \int_{s}^{x_{1}} \mathbf{Y}'(x_{2}) \ dx_{2} \right] \mathbf{A}(x_{1}) \ dx_{1} \\ &= \mathbf{I} + \int_{s}^{t} \left[\mathbf{I} + \int_{s}^{x_{1}} \mathbf{Y}(x_{2}) \mathbf{A}(x_{2}) \ dx_{2} \right] \mathbf{A}(x_{1}) \ dx_{1} \\ &= \mathbf{I} + \int_{s}^{t} \mathbf{A}(x_{1}) \ dx_{1} + \int_{s}^{t} \int_{s}^{x_{1}} \mathbf{Y}(x_{2}) \mathbf{A}(x_{2}) \mathbf{A}(x_{1}) \ dx_{2} \ dx_{1}. \end{split}$$

One can continue indefinitely and see that we have the following representation.

Corollary 1.4. (Bladt) (Peano Representation) The prduct integral has series representation given by

$$\begin{split} \prod_s^t (\mathbf{I} + \mathbf{A}(x) \ dx) &= \mathbf{I} + \sum_{i=1}^\infty \int_s^t \int_s^{x_1} \cdots \int_s^{x_n} A(x_n) \cdots \mathbf{A}(x_2) \mathbf{A}(x_1) \ dx_n \cdots dx_2 \ dx_1 \\ &= \mathbf{I} + \int_s^t \mathbf{A}(x_1) \ dx_1 + \sum_{i=2}^\infty \int_s^t \int_s^{x_1} \cdots \int_s^{x_n} A(x_n) \cdots \mathbf{A}(x_2) \mathbf{A}(x_1) \ dx_n \cdots dx_2 \ dx_1. \end{split}$$

15.5.1 Properties of the Product Integral

The product integral is the fundamental solution in the sense that if

$$\mathbf{Y}'(t) = \mathbf{Y}(t)\mathbf{A}(t), \quad \mathbf{Y}(s) = \mathbf{C},$$

then

$$\mathbf{Y}(t) = \mathbf{C} \prod_{s}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx).$$

We furthermore have for any s, t, u have

$$\mathbf{Y}(t) = \mathbf{Y}(s) \prod_{x=0}^{u} (\mathbf{I} + \mathbf{A}(x) \ dx) \prod_{x=0}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx).$$

From uniqueness we then get the following theorem.

Theorem 1.5. (Bladt) For any s, t, u we have that

$$\prod_{x=0}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx) = \prod_{x=0}^{u} (\mathbf{I} + \mathbf{A}(x) \ dx) \prod_{x=0}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx).$$

By choosing u = t in the above we get the following corollary.

Corollary 1.6. (Bladt) The inverse of the product integral is

$$\left[\prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \right]^{-1} = \prod_{t}^{s} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right).$$

Theorem 1.7. (Bladt) If A(x) commutes withall B(x = i.e. A(x)B(x) = B(x)A(x), then

$$\prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \prod_{s}^{t} \left(\mathbf{I} + \mathbf{B}(x) \ dx \right) = \prod_{s}^{t} \left(\mathbf{I} + \left(\mathbf{A}(x) + \mathbf{B}(x) \right) \ dx \right).$$

Corollary 1.8. (Bladt) Let r be a real-valued function, then

$$\exp\left(-\int_s^t r(x)\ dx\right) \prod_s^t \left(\mathbf{I} + \mathbf{A}(x)\ dx\right) = \prod_s^t \left(\mathbf{I} + (\mathbf{A}(x) - r(x)\mathbf{I})\ dx\right).$$

Theorem 1.9. (Bladt) If $\mathbf{A}(x)$ commutes for all x i.e. $\mathbf{A}(y)\mathbf{A}(x) = \mathbf{A}(x)\mathbf{A}(y)$, then

$$\prod_{s}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx) = e^{\int_{s}^{t} \mathbf{A}(x) \ dx}.$$

In particular, if $\mathbf{A}(x) = \mathbf{A}\lambda(x)$ for some real-valued function λ and a $n \times n$ matrix \mathbf{A} then

$$\prod_{s}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx) = e^{\mathbf{A} \int_{s}^{t} \lambda(x) \ dx}.$$

Taking the function $\lambda = 1$ we have that $\int_s^t \lambda(x) \ dx = (t - s)$ and so we get the result.

Theorem 1.10. (Bladt) If A(x) = A is constant then

$$\prod_{s}^{t} (\mathbf{I} + \mathbf{A}(x) \ dx) = e^{\mathbf{A}(t-s)}.$$

Theorem 1.11. (Bladt) The product integral satisfies

$$\frac{\partial}{\partial s} \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) = -\mathbf{A}(s) \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right).$$

On the other hand, the solution to the system of differential equations

$$\frac{\partial}{\partial s} \mathbf{X}(s) = -\mathbf{A}(s) \mathbf{X}(s)$$

with initial condition $\mathbf{X}(t) = \mathbf{I}$ is the function

$$s \mapsto \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right)$$

We have a very useful result for insurance mathematics and markov chains given by the Van Loan's theorem giving the product integral of a block matrix with zero lower left block.

Theorem 1.12. (Bladt) Let A(x), B(x) and C(x) be continuous matrix functions such that

$$\mathbf{D}(x) = \begin{bmatrix} \mathbf{A}(x) & \mathbf{B}(x) \\ \mathbf{0} & \mathbf{C}(x) \end{bmatrix},$$

is a square matrix. Then the product integral of D is

$$\prod_{s}^{t} \left(\mathbf{I} + \mathbf{D}(x) \ dx \right) = \begin{bmatrix} \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) & \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_{u}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \\ \mathbf{0} & \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \end{bmatrix}.$$

Proof.

Start by defining $\mathbf{X}(t)$ as below

$$\begin{split} \mathbf{X}(t) &= \begin{bmatrix} \prod_s^t \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) & \int_s^t \prod_s^u \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_u^t \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \\ \mathbf{0} & \prod_s^t \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \end{bmatrix} \\ &:= \begin{bmatrix} \mathbf{X}_{11}(t) & \mathbf{X}_{12}(t) \\ \mathbf{X}_{21}(t) & \mathbf{X}_{22}(t) \end{bmatrix}. \end{split}$$

The derivative is then

$$\frac{d}{dt}\mathbf{X}(t) = \begin{bmatrix} \frac{d}{dt}\mathbf{X}_{11}(t) & \frac{d}{dt}\mathbf{X}_{12}(t) \\ \frac{d}{dt}\mathbf{X}_{21}(t) & \frac{d}{dt}\mathbf{X}_{22}(t) \end{bmatrix}.$$

with

$$\begin{split} &\frac{d}{dt}\mathbf{X}_{11}(t) = \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx\right) \mathbf{A}(t), \\ &\frac{d}{dt}\mathbf{X}_{12}(t) = \frac{d}{dt} \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx\right) \mathbf{B}(u) \prod_{u}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx\right) \ du, \\ &\frac{d}{dt}\mathbf{X}_{21}(t) = \mathbf{0}, \\ &\frac{d}{dt}\mathbf{X}_{22}(t) = \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx\right) \mathbf{C}(t). \end{split}$$

Consider that

$$\prod_{u}^{t} (\mathbf{I} + \mathbf{C}(x) \ dx) = \prod_{u}^{t} (\mathbf{I} + \mathbf{C}(x) \ dx) \underbrace{\prod_{t}^{s} (\mathbf{I} + \mathbf{C}(x) \ dx) \prod_{s}^{t} (\mathbf{I} + \mathbf{C}(x) \ dx)}_{=\mathbf{I}}$$

$$= \prod_{u}^{s} (\mathbf{I} + \mathbf{C}(x) \ dx) \underbrace{\prod_{t}^{t} (\mathbf{I} + \mathbf{C}(x) \ dx)}_{s},$$

hence using af Riemann approximation of the integral $\mathbf{X}_{12}(t)$ we may write

$$\begin{split} \mathbf{X}_{12}(t) &= \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_{u}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \\ &= \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_{u}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \\ &= \lim_{n \to \infty} \frac{t - s}{n} \sum_{i=0}^{n} \prod_{s}^{s + (t - s)i/n} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(s + (t - s)i/n) \prod_{s + (t - s)i/n}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \\ &= \left\{ \lim_{n \to \infty} \frac{t - s}{n} \sum_{i=0}^{n} \prod_{s}^{s + (t - s)i/n} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(s + (t - s)i/n) \prod_{s + (t - s)i/n}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \right\} \\ &\prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \\ &= \left\{ \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_{u}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \right\} \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right). \end{split}$$

We then get that

$$\begin{split} \frac{d}{dt}\mathbf{X}_{12}(t) &= \frac{d}{dt} \left\{ \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_{u}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \right\} \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \\ &= \left\{ \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(t) \prod_{t}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \right\} \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \\ &+ \left\{ \int_{s}^{t} \prod_{s}^{u} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(u) \prod_{u}^{s} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \ du \right\} \prod_{s}^{t} \left(\mathbf{I} + \mathbf{C}(x) \ dx \right) \mathbf{C}(t) \\ &= \prod_{s}^{t} \left(\mathbf{I} + \mathbf{A}(x) \ dx \right) \mathbf{B}(t) + \mathbf{X}_{12}(t) \mathbf{C}(t). \end{split}$$

We hope to show that $\frac{d}{dt}\mathbf{X}(t) = \mathbf{X}(t)\mathbf{D}(t)$. Calculating the right side we have

$$\begin{split} \mathbf{X}(t)\mathbf{D}(t) &= \begin{bmatrix} \mathbf{X}_{11}(t) & \mathbf{X}_{12}(t) \\ \mathbf{0} & \mathbf{X}_{22}(t) \end{bmatrix} \begin{bmatrix} \mathbf{A}(x) & \mathbf{B}(x) \\ \mathbf{0} & \mathbf{C}(x) \end{bmatrix} \\ &= \begin{bmatrix} \mathbf{X}_{11}(t)\mathbf{A}(x) & \mathbf{X}_{11}(t)\mathbf{B}(x) + \mathbf{X}_{12}(t)\mathbf{C}(x) \\ \mathbf{0} & \mathbf{X}_{22}(t)\mathbf{C}(x) \end{bmatrix}, \end{split}$$

given that **X** satisfies the desired differential equation and so it follows that **X** is the product integral $\prod_{s=0}^{t} (\mathbf{I} + \mathbf{D}(x) \ dx)$ as desired.

Chapter 16

Linear Algebra

16.1 Invertible matrices

This sections study some fundamental properties of the *invertible matrix*. We start by defining what an invertible matrix is.

Definition. Let A be an $n \times m$ matrix and B be an $m \times n$ matrix. We say that A is invertible if

$$AB = I_{\min(m,n)}.$$

In general, we only consider square matrices. If the above holds we say that B is A's inverse and we write $B = A^{-1}$.

We can consider some equivalence statements regarding invertible matrices.

Theorem. (The Invertible Matrix Theorem) Let A be a $n \times n$ matrix over a field K (\mathbb{R}^n), then the following statements are equivalent

- 1. There exists an $n \times n$ matrix B such that $AB = I_n = BA$.
- 2. There exist either a left inverse B or a right inverse C i.e. $BA = I_n = AC$. In this case, B = C.
- 3. A has an inverse and is nonsingular and is nondegenerate.
- 4. A is row-equivalent to I_n .
- 5. A is column-equivalent to I_n .
- 6. A has n pivot positions.
- 7. A has full rank i.e. rank(A) = n (spans K).
- 8. The equation Ax = 0 $(x \in K)$ has only the trivial solution x = 0.
- 9. The equation Ax = b has only one solution x.
- 10. The kernal of A is trivial i.e. $ker(A) = \{0\}.$
- 11. The columns of A are linearly independent.
- 12. The columns of A span K.
- 13. span(A) = K.
- 14. The columns of A form a basis of K.
- 15. The linear transformation Ax is a bijection from K to K.
- 16. A has non-zero determinant i.e. $det(A) \neq 0$.
- 17. A has not 0 as an eigenvalue.
- 18. The transpose of A is invertible.
- 19. A can be expressed as a finite product of elemeary matrices.

We futhermore have some properties.

Proposition. (Properties) Let A be an $n \times n$ invertible matrix. Then

- $\begin{array}{l} 1. \ \, (A^{-1})^{-1} = A \\ 2. \ \, (kA)^{-1} = k^{-1}A^{-1} \ \, with \, \, k \neq 0. \end{array}$
- 3. $(Ax)^+ = x^+A^{-1}$ if A has orthonormal columns. $(\cdot)^+$ denotes the Moore-Penrose inverse and x is a
- 4. If B is an $n \times n$ invertible matrix then $(AB)^{-1} = B^{-1}A^{-1}$.
- 5. $\det(A^{-1}) = (\det(A))^{-1}$

The property 2 is especially useful in some settings. Consider for instance

$$A = \begin{bmatrix} \sigma & 0 \\ \sigma & \sigma \end{bmatrix} = \sigma \begin{bmatrix} 1 & 0 \\ 1 & 1 \end{bmatrix} = \sigma \tilde{A}.$$

Then we simply find the inverse of \tilde{A} and multiply by σ^{-1} . That is,

$$A^{-1} = \frac{1}{\sigma} \tilde{A}^{-1} = \frac{1}{\sigma} \begin{bmatrix} 1 & 0 \\ -1 & 1 \end{bmatrix} = \begin{bmatrix} 1/\sigma & 0 \\ -1/\sigma & 1/\sigma \end{bmatrix}.$$

We have an easy propositions regarding diagonal matrices.

Proposition. If A is an diagonal matrix, then A is invertible. In particular,

$$A^{-1} = \operatorname{diag}(A_{11}^{-1}, ..., A_{nn}^{-1}).$$

Chapter 17

Coding

17.1 R-Packages

17.1.1 mlr3

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