# Notes on Quantum Field Theory 

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## Chapter 1

## Introduction

Quantum Field Theory (QFT) is the fundamental tool that is currently used for the description of physics at very short distances and high energies. Since high energy implies relativistic motion, QFT has to nicely combine special relativity with quantum mechanics. The description in terms of "fields", indeed, arise to have a manifestly relativistic invariant description, where time and space are treated (almost) on equal footing. One might think that including special relativity in quantum mechanics should be possible without drastic consequences. This is not true, and its reason is intuitively clear. At very short time scales, the energy-time uncertainty principle tells us that particles with energy $E \geq m c^{2}$ could be created from the vacuum for a time $t \sim \hbar / E$ (virtual particles), before disappearing again in the vacuum. This effect is totally negligible in studying physical systems at low energies and long time scales, but it becomes relevant for physical processes whose time scale is $t \sim \hbar / E$ or less. A quantum description in terms of single-particle wave function is then inadequate and a more powerful description is needed. This formulation is in fact Quantum Field Theory. It leads to striking consequences, such as the prediction of anti-particles and an understanding of the spin-statistic relation between particles. The experimental successes of QFT are impressive, in particular when applied to the description of electrodynamics, giving rise to the Quantum Electro Dynamics (QED).

Most of the considerations in these lectures are devoted to the study of fields which are weakly interacting, namely in which the interactions can be studied in a perturbative fashion, starting from the description of free fields.

These notes do assume that the reader has a basic knowledge of QFT. Quantizations of spin 0 , spin $1 / 2$ and abelian spin 1 fields are assumed, as well as basic notions of the path integral formulation of QFT (including Berezin integration for fermions), Feynman rules, basic knowledge of renormalization and the notion of functional generators of disconnected,
connected and one-particle irreducible (1PI) Green functions.
In writing these notes we have often consulted refs. [1] and especially [2]. Some parts of these notes follow in part either ref. [1] or ref. [2]. When this is the case, we warn the reader with a footnote. Notice that in these notes the metric convention is mostly minus, like in ref. [1], in contrast to ref.[2], where it is mostly plus. This implies a multitude of sign changes with respect to ref. [2]. Moreover, we warn the reader that we do not always follow the same notation as refs. [1, 2], in order to have a common and light notation throughout the notes.

These notes cover most of the QFT course given at SISSA by M.S., in collaboration with Roberto Iengo until the academic year 2010-2011, in collaboration with Andrea Gambassi until 2014-2015, and in collaboration with Joan Elias-Miró in the years 2016-2017 and 2017-2018. The tutorials and exercises, essential parts of the QFT course, are not reported in these notes. This implies that some technical topics closely associated with the exercises, such as details of the renormalization of QED, Yukawa and other theories, are not currently included in these lecture notes. The notes are far from being comprehensive. Due to lack of time, several important topics are not covered at all, or just mentioned. Infra-red divergences, the deep inelastic scattering and the operator product expansion, connection to critical phenomena and statistical field theory are examples of important topics currently missing. Phenomenological applications are limited to a minimum, since they are more systematically considered in the Standard Model course.

The topics marked with a $*$ in the text are optional for students in the Astroparticle curriculum. The topics marked with $* *^{*}$ are optional for all the students.

These notes are preliminary and surely contain many typos, imprecisions, etc.
We hope that the students will help us in improving the notes and in spotting the many mistakes in there.

## Acknowledgments

I thank Roberto Iengo and Andrea Gambassi for their contribution in the elaboration of part of these lecture notes. I thank Roberto Iengo, Andrea Gambassi and Joan EliasMiró for their precious collaboration over the years in giving this course. I thank Matteo Bertolini, Marcello D. Caio, Lorenzo Di Pietro, Matthijs Hogervorst, Marco Letizia, Himanshu Raj, Andrea Romanino, Giacomo Sberverglieri for useful comments and help in debugging these notes. I am particularly grateful to Marco Gorghetto for having found so many typos!

## Chapter 2

## General Properties of QFT

Quantum field theory aims at describing the processes which occur as a result of the interaction between elementary particles, mediated by a certain number of force carriers. The very same requirement of Lorentz covariance (i.e., compatibility with the theory of special relativity) is a fundamental ingredient of the theory and actually motivated the introduction of concepts such as antiparticles which were later on detected experimentally. Most of our knowledge of the interaction between elementary particles comes from the prototypical scattering experiment in which incoming particles are well-separated in space before getting close to each other in such a way to interact and scatter. After scattering has occurred, the particles resulting from the process typically separate again in space in such a way that their interaction is negligible and their properties as isolated particles, e.g., mass, can be determined via suitable detectors. In this chapter we discuss QFT keeping in mind the framework of scattering experiments in order to define the concept of asymptotic states and assign a specific meaning to the mass of particles from their dynamical properties (e.g., the propagator). Then we discuss some important properties of QFT which derive from the existence of a unitary mapping (the $S$-matrix) between the incoming and outgoing particle states, such as the optical theorem, and we discuss how scattering amplitude and decay rates can be inferred from the correlation function of the fields naturally described by QFT.

### 2.1 The Källén-Lehmann Spectral Representation

### 2.1.1 Asymptotic Theory: a Brief Reminder

In a nutshell, quantum field theory can be thought of as based on the following elements: (a) the possible (quantum) states of the theory are generated from a (unique) vacuum
state $|0\rangle$ by the action of free fields $\phi_{\text {in }}(x)$, which generates the Fock space of states (and which we assume to be a real scalar); (b) physical observables - such as the interacting field $\phi(x)$ can be all expressed in terms of $\phi_{i n}(x)$. The basic idea behind this setting is that $\phi(x)$ as well as all the other observables can be actually obtained from the free fields $\phi_{\text {in }}(x)$ by switching adiabatically on and off the interaction as $\left|x_{0}\right| \rightarrow \infty$. This construction is clearly relevant to scattering processes which originate from particles (wavepackets) wellseparated in space. Starting from $\phi\left(x_{0}, \vec{x}\right)$ one should recover $\phi_{\text {in }}\left(x_{0}, \vec{x}\right)$ as $x_{0} \rightarrow-\infty$, but this generally occurs up to the wave-function renormalization constant $Z_{i n}^{1 / 2}$ :

$$
\begin{equation*}
\phi\left(x_{0} \rightarrow-\infty, \vec{x}\right) \rightarrow Z_{i n}^{1 / 2} \phi_{i n}\left(x_{0}, \vec{x}\right) . \tag{2.1.1}
\end{equation*}
$$

Analogously, $\phi\left(x_{0} \rightarrow+\infty, \vec{x}\right) \rightarrow Z_{\text {out }}^{1 / 2} \phi_{\text {out }}\left(x_{0}, \vec{x}\right)$, where $\phi_{\text {out }}$ is a free field. The fields $\phi_{\text {in/out }}$ satisfy the free Klein-Gordon equation

$$
\begin{equation*}
\left(\square+m^{2}\right) \phi_{\text {in } / \text { out }}(x)=0, \tag{2.1.2}
\end{equation*}
$$

while the interacting field does not, $\left(\square+m^{2}\right) \phi(x) \neq 0$. In and out free fields can be expressed in terms of the corresponding creation and annihilation operators as

$$
\begin{equation*}
\phi_{\text {in } / \text { out }}(x)=\int \frac{d^{3} \vec{p}}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}}}}\left(a(\vec{p})_{\text {in/out }} e^{-i p x}+a^{\dagger}(\vec{p})_{\text {in } / \text { out }} e^{i p x}\right), \tag{2.1.3}
\end{equation*}
$$

where $p x=\omega_{\vec{p}} x_{0}-\vec{p} \cdot \vec{x}, p_{0}=\omega_{\vec{p}}=\sqrt{\vec{p}^{2}+m^{2}}$. The canonical quantization condition

$$
\begin{equation*}
\left[\dot{\phi}_{\text {in/out }}(\vec{x}, t), \phi_{\text {in/out }}(\vec{y}, t)\right]=i \delta^{(3)}(\vec{x}-\vec{y}), \tag{2.1.4}
\end{equation*}
$$

where the dot stands for a time derivative, implies the commutation relations

$$
\begin{equation*}
\left[a(\vec{p})_{\text {in/out }}, a^{\dagger}(\vec{q})_{\text {in/out }}\right]=\delta^{3}(\vec{p}-\vec{q}) \tag{2.1.5}
\end{equation*}
$$

Inverting eq.(2.1.3), we can express the annihilation and creation operators in terms of $\phi$ and $\dot{\phi}$ :

$$
\begin{equation*}
a(\vec{p})_{\text {in/out }}=i \int d^{3} \vec{x} \underbrace{\frac{e^{i p x}}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}}}}}_{h_{\vec{p}}(x)} \stackrel{\leftrightarrow}{\partial_{0}} \phi_{\text {in }(\text { out })}(x) \tag{2.1.6}
\end{equation*}
$$

where $\overleftrightarrow{\partial_{0}} \equiv \overrightarrow{\partial_{0}}-\overleftarrow{\partial_{0}}$. The hermitian conjugate of the right-hand side of eq.(2.1.6) clearly gives $a^{\dagger}(\vec{p})_{\text {in/out }}$. Notice that the time-dependence of this expression is only apparent. In general, for any function $h(x)$ satisfying the free Klein-Gordon equation $\left(\square+m^{2}\right) h(x)=0$,

$$
\begin{equation*}
\int d^{3} \vec{x} h(x) \stackrel{\leftrightarrow}{\partial_{0}} \phi_{i n(o u t)}(x) \tag{2.1.7}
\end{equation*}
$$

is time-independent. This is easily shown by taking a time derivative of this expression, the use of the Klein-Gordon equation and an integration by parts. The same result applies if we substitute $\phi_{\text {in }(o u t)}(x)$ by some other solution of the free Klein-Gordon equation.

The incoming multiparticle states are generated from the corresponding vacuum state $|0\rangle_{\text {in }}$ by the action of the creation operators $a^{\dagger}(\vec{p})_{\text {in }}$ of the free theory, e.g., $|\vec{p}\rangle_{\text {in }}=$ $a^{\dagger}(\vec{p})_{i n}|0\rangle_{\text {in }}$ for the single-particle state with three-momentum $\vec{p} .{ }^{1}$ Analogous construction can be done for the outgoing states, starting from the corresponding vacuum $|0\rangle_{\text {out }}$ : $|\vec{p}\rangle_{\text {out }}=a^{\dagger}(\vec{p})_{\text {out }}|0\rangle_{\text {out }}$. Given that both the incoming and outgoing particle states are nothing but different representations of the free theories, there should be an unitary isomorphism between them, the so-called $S$-matrix: $S|i\rangle_{\text {out }}=|i\rangle_{\text {in }}$. We will discuss the consequences of the general properties of the $S$-matrix in secs. 2.3 and 2.4. Accordingly, the transition amplitude from an initial state $i\left(|i\rangle_{\text {in }}\right)$ to a final state $f\left(|f\rangle_{\text {out }}\right)$ can be expressed only in terms of incoming states as

$$
\begin{equation*}
{ }_{\text {out }}\langle f \mid i\rangle_{\text {in }}={ }_{\text {in }}\langle f| S|i\rangle_{\text {in }} . \tag{2.1.8}
\end{equation*}
$$

The stability of the vacuum requires that $|0\rangle_{\text {in }}=|0\rangle_{\text {out }} \equiv|0\rangle$ while the kinematic stability of the massive one-particle state implies, in addition, $|\vec{p}\rangle_{\text {in }}=|\vec{p}\rangle_{\text {out }} \equiv|\vec{p}\rangle$. The isomorphism induced by $S$ implies that, at the level of field operators, $\phi_{\text {in }}(x)=S \phi_{\text {out }}(x) S^{-1}$. The $S$ matrix can be used in order to express all the states in terms of the incoming ones, which we always refer to in what follows, unless specified differently.

Despite the interacting field does not satisfy the free Klein-Gordon equation, the amplitude $\langle 0| \phi(x)|\vec{p}\rangle$ does. Indeed, invariance under translations implies

$$
\begin{equation*}
\phi(x)=e^{i \hat{p} x} \phi(0) e^{-i \hat{p} x} \tag{2.1.9}
\end{equation*}
$$

where $\hat{p}$ is the four-momentum operator, and thus

$$
\begin{equation*}
\langle 0| \phi(x)|\vec{p}\rangle=\langle 0| \phi(0)|\vec{p}\rangle e^{-i p x} \tag{2.1.10}
\end{equation*}
$$

It then follows that

$$
\begin{equation*}
\left(\square+m^{2}\right)\langle 0| \phi(x)|\vec{p}\rangle=0, \tag{2.1.11}
\end{equation*}
$$

even if $\left(\square+m^{2}\right) \phi(x) \neq 0$. By using eq.(2.1.7) we notice that

$$
\begin{equation*}
\int d^{3} \vec{x} h_{\vec{p}}(x) \overleftrightarrow{\partial_{0}}\langle 0| \phi(x)|\vec{p}\rangle \tag{2.1.12}
\end{equation*}
$$

[^0]is time-independent. Equation (2.1.3) fixes the analogue amplitude for in and out fields to be
\[

$$
\begin{equation*}
\langle 0| \phi_{i n}(x)|\vec{p}\rangle=\frac{e^{-i p x}}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}}}} . \tag{2.1.13}
\end{equation*}
$$

\]

Taking the limit $x^{0} \rightarrow \pm \infty$ of eq.(2.1.12) and using eq.(2.1.13) we get

$$
\begin{equation*}
Z_{\text {in }}=Z_{\text {out }} \equiv Z \tag{2.1.14}
\end{equation*}
$$

and hence

$$
\begin{equation*}
\langle 0| \phi(x)|\vec{p}\rangle=Z^{1 / 2} \frac{e^{-i p x}}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}}}} \tag{2.1.15}
\end{equation*}
$$

Note that $\phi_{i n}(x)|0\rangle$ generates only the one-particle state, whereas $\phi(x)|0\rangle$ generates, in addition, multiparticle states because of the interaction. As a result, one expects $\left.|\langle\vec{p}| \phi(x)| 0\rangle\left.\right|^{2} \leq\left|\langle\vec{p}| \phi_{\text {in }}(x)\right| 0\right\rangle\left.\right|^{2}$ and therefore $|Z| \leq 1$ and, in particular, $|Z|$ can be interpreted as the probability of generating the one-particle state when applying the interacting field $\phi(x)$ to the vacuum state $|0\rangle$. We will determine $Z$ in sec. 2.1.2.

The limit in eq. (2.1.1) has to be understood in a weak sense because at best it can hold at the level of matrix elements between fields well separated in space (in the form of suitable wavepackets), while it cannot hold for operators. Indeed, both the interacting and free fields $\phi$ have to satisfy the equal-time canonical commutation relations $\left[\dot{\phi}\left(x_{0}, \vec{x}\right), \phi\left(x_{0}, \vec{x}^{\prime}\right)\right]=-i \delta^{3}\left(\vec{x}-\vec{x}^{\prime}\right)$ which would imply $Z=1$ by taking the limit $x_{0} \rightarrow-\infty$.

### 2.1.2 Spectral Representation

Important properties of an interacting quantum field theory emerge from the so-called spectral representation of vacuum expectation values of suitable quantities. Here we consider both the commutator $\langle 0|\left[\phi(x), \phi\left(x^{\prime}\right)\right]|0\rangle$ and the time-ordered product $\langle 0| T \phi(x) \phi\left(x^{\prime}\right)|0\rangle$ of the fields, i.e., the propagator of the theory, where the time-ordering is defined by $T \phi(x) \phi\left(x^{\prime}\right)=\theta\left(x_{0}-x_{0}^{\prime}\right) \phi(x) \phi\left(x^{\prime}\right)+\theta\left(x_{0}^{\prime}-x_{0}\right) \phi\left(x^{\prime}\right) \phi(x)$, being $\theta\left(x_{0}>0\right)=1$ and 0 otherwise (note that in case of Fermionic fields, the exchange of their order introduces an additional - sign). Both these quantities can be expressed in terms of

$$
\begin{align*}
\langle 0| \phi(x) \phi\left(x^{\prime}\right)|0\rangle & \left.=\sum_{n}\langle 0| \phi(x)|n\rangle\langle n| \phi\left(x^{\prime}\right)|0\rangle=\sum_{n} e^{-i p_{n}\left(x-x^{\prime}\right)}|\langle 0| \phi(0)| n\right\rangle\left.\right|^{2} \\
& \left.=\int d^{4} q e^{-i q\left(x-x^{\prime}\right)} \sum_{n}|\langle 0| \phi(0)| n\right\rangle\left.\right|^{2} \delta^{4}\left(q-p_{n}\right)  \tag{2.1.16}\\
& =\int \frac{d^{4} q}{(2 \pi)^{3}} e^{-i q\left(x-x^{\prime}\right)} \rho(q)
\end{align*}
$$

where in the first line we introduced the completeness relation $\mathbb{\square}=\sum_{n}|n\rangle\langle n|$ in terms of the multi-particle states $|n\rangle$ with definite (total) momentum $p_{n}$ (such that $\hat{p}|n\rangle=p_{n}|n\rangle$ )
and we used eq. (2.1.9); $\sum_{n}$ stands both for the sum over the particles number and by the integral on the corresponding phase space. In the second line we introduced the identity $1=\int d^{4} q \delta^{4}\left(q-p_{n}\right)$ and, in the last, the so-called spectral density

$$
\begin{equation*}
\left.\rho(q) \equiv(2 \pi)^{3} \sum_{n}|\langle 0| \phi(0)| n\right\rangle\left.\right|^{2} \delta^{4}\left(q-p_{n}\right) \tag{2.1.17}
\end{equation*}
$$

which is a Lorentz-invariant, positive scalar and therefore is actually a function of $q^{2}$. In addition, $p_{n}=\sum_{i=1}^{n} p_{n, i}$ where $p_{n, i}$ is the momentum of the $i$-th particle within the $n$-particle state $|n\rangle$ and therefore $\left(p_{n, i}\right)_{0}>0$ with $p_{n, i}^{2}>0$; this implies that $\left(p_{n}\right)_{0}>0$ and $p_{n}^{2}>0$ and, in turn, that $\rho(q)$ vanishes in the backward light cone: $\rho(q) \mapsto \rho\left(q^{2}\right) \theta\left(q_{0}\right)$ with $\rho\left(q^{2}<\right.$ $0)=0$. In eq. (2.1.17), it is convenient to isolate the contribution of the one-particle state $n=1$, which can be readily calculated by taking into account the normalization (2.1.15): $\left.\left.\sum_{n=1}|\langle 0| \phi(0)| n\right\rangle\left.\right|^{2} \delta^{4}\left(q-p_{n}\right)=\int d^{3} \vec{p}|\langle 0| \phi(0)| \vec{p}\right\rangle\left.\right|^{2} \delta^{4}(q-p)=Z \theta\left(q_{0}\right) \delta\left(q^{2}-m^{2}\right) /(2 \pi)^{3}$. The resulting expression for $\rho$ is

$$
\begin{equation*}
\left.\rho\left(q^{2}\right) \theta\left(q_{0}\right)=Z \delta\left(q^{2}-m^{2}\right) \theta\left(q_{0}\right)+(2 \pi)^{3} \sum_{n>1}|\langle 0| \phi(0)| n\right\rangle\left.\right|^{2} \delta^{4}\left(q-p_{n}\right) \tag{2.1.18}
\end{equation*}
$$

which reveals its generic structure: it consists of an isolated $\delta$ corresponding to the mass of the asymptotic field (including the self-interaction) and of a continuum starting from the threshold $m_{t h}^{2}$ for the production of multi-particle states. In addition, there might be some bound states in the gap, as depicted in fig. 2.1(a).

In terms of the spectral density, $\langle 0|[\phi(x), \phi(0)]|0\rangle$ can be written as

$$
\begin{align*}
\langle 0|[\phi(x), \phi(0)]|0\rangle & =\int \frac{d^{4} q}{(2 \pi)^{3}} \rho\left(q^{2}\right) \theta\left(q_{0}\right)\left[e^{-i q x}-e^{i q x}\right] \\
& =\int_{0}^{\infty} d \sigma \rho(\sigma) \underbrace{\int \frac{d^{4} q}{(2 \pi)^{3}} \delta\left(q^{2}-\sigma\right) \theta\left(q_{0}\right)\left[e^{-i q x}-e^{i q x}\right]}_{i \Delta_{c}^{0}(x ; \sigma)} \tag{2.1.19}
\end{align*}
$$

where we introduced the identity in the form $1=\int_{0}^{\infty} d \sigma \delta\left(q^{2}-\sigma\right)$. In the previous equation one recognizes the expression of $\langle 0|\left[\phi_{0}(x), \phi_{0}(0)\right]|0\rangle \equiv i \Delta_{c}^{0}\left(x ; m^{2}\right)$ for a free scalar field $\phi_{0}(x)$ of mass $m$. The analogous expression for $\langle 0| T \phi(x) \phi(0)|0\rangle$ is

$$
\begin{align*}
\langle 0| T \phi(x) \phi(0)|0\rangle & =\int \frac{d^{4} q}{(2 \pi)^{3}} \rho\left(q^{2}\right) \theta\left(q_{0}\right)\left[\theta\left(x_{0}\right) e^{-i q x}+\theta\left(-x_{0}\right) e^{i q x}\right] \\
& =\int_{0}^{\infty} d \sigma \rho(\sigma) \underbrace{\int \frac{d^{4} q}{(2 \pi)^{3}} \delta\left(q^{2}-\sigma\right) \theta\left(q_{0}\right)\left[\theta\left(x_{0}\right) e^{-i q x}+\theta\left(-x_{0}\right) e^{i q x}\right]}_{i \Delta_{F}^{0}(x ; \sigma)} \tag{2.1.20}
\end{align*}
$$



Figure 2.1: Sketches (a) of the spectral density $\rho\left(q^{2}\right)$ as a function of $q^{2}$, which highlights the presence of a $\delta$ corresponding to the physical mass, of possible additional peaks related to bound states (dashed) and of a continuum associated to multi-particle states. (b) Analytic structure of the propagator $G^{(2)}(q)$ of a scalar field, as a function of $q^{2} \in \mathbb{C}$ which highlights the presence of an isolated pole corresponding to the physical mass, possible additional poles due to bound states and a branch cut related to multi-particle states.
and involves, instead, the Feynman propagator $i \Delta_{F}^{0}\left(x ; m^{2}\right) \equiv\langle 0| T \phi_{0}(x) \phi_{0}(0)|0\rangle$ of the same field. Taking into account eq. (2.1.18) one finds

$$
\begin{align*}
\langle 0|[\phi(x), \phi(0)]|0\rangle & =Z i \Delta_{c}^{0}\left(x ; m^{2}\right)+\int_{m_{t h}^{2}}^{\infty} d \sigma \rho(\sigma) i \Delta_{c}^{0}(x ; \sigma),  \tag{2.1.21}\\
\langle 0| T \phi(x) \phi(0)|0\rangle & =Z i \Delta_{F}^{0}\left(x ; m^{2}\right)+\int_{m_{t h}^{2}}^{\infty} d \sigma \rho(\sigma) i \Delta_{F}^{0}(x ; \sigma), \tag{2.1.22}
\end{align*}
$$

which provide the Källén-Lehmann spectral representation of the commutator and the propagator, respectively. Occasionally, it might be useful to introduce renormalized fields $\phi_{R}(x) \equiv Z^{-1 / 2} \phi(x)$ and the renormalized spectral density $\rho_{R}$ via $\rho=Z \rho_{R}$. The consequences of these relations can be easily worked out: for example, by taking the derivative $\partial_{0}=\partial / \partial x_{0}$ of the l.h.s. of eq. (2.1.21) and by eventually setting $x_{0}=0$ one recovers the canonical commutation relation for the field $\phi$; the same relation for the free field $\phi_{0}$ implies that $\left.\partial_{0} i \Delta_{c}^{0}\left(x ; m^{2}\right)\right|_{x_{0}=0}=-i \delta^{3}(\vec{x})$, which appears on the r.h.s. of eq. (2.1.21) and implies

$$
\begin{equation*}
1=Z\left[1+\int_{m_{t h}^{2}}^{\infty} d \sigma \rho_{R}(\sigma)\right] . \tag{2.1.23}
\end{equation*}
$$

This expression shows that $Z \leq 1\left(\rho_{R} \geq 0\right)$ and allows one to calculate $Z$ from the spectral density. Consider now the Fourier transform of the renormalized propagator
$G_{R}^{(2)}(x)=\langle 0| T \phi_{R}(x) \phi_{R}(0)|0\rangle=G^{(2)}(x) / Z$. Recalling that for the free theory

$$
\begin{equation*}
\Delta_{F}^{0}\left(x, m^{2}\right)=\int \frac{d^{4} q}{(2 \pi)^{4}} \frac{e^{-i q x}}{q^{2}-m^{2}+i \epsilon} \tag{2.1.24}
\end{equation*}
$$

one finds

$$
\begin{equation*}
G_{R}^{(2)}\left(q^{2}\right)=\frac{i}{q^{2}-m^{2}+i \epsilon}+\int_{m_{t h}^{2}}^{\infty} d \sigma \rho_{R}(\sigma) \frac{i}{q^{2}-\sigma+i \epsilon} \tag{2.1.25}
\end{equation*}
$$

As a function of $q^{2} \in \mathbb{C}, G_{R}^{(2)}\left(q^{2}\right)$ is characterized by an isolated pole for $q^{2}=m^{2}$, with residue $i$ and by a branch cut on the positive real axis starting from $q^{2}=m_{t h}^{2}$ and controlled by the spectral density, with possible additional poles due to bound states, as depicted in fig. 2.1(b). Note in particular that $G_{R}^{(2)}(z)$ is analytic in the complex plane away from the real axis. The $i \epsilon$ prescription in the propagator tells us that the physical sheet lies above the branch-cut. In sec. 2.4 we provide a physical interpretation of the discontinuity which arises in $G_{R}^{(2)}(q)$ upon crossing the branch cut. Another relevant relation between the propagator of the theory and the spectral density can be found from eq. (2.1.25), taking into account that

$$
\begin{equation*}
\frac{1}{x+i \epsilon}=P \frac{1}{x}-i \pi \delta(x) \quad \text { for } \quad \epsilon \rightarrow 0^{+} \tag{2.1.26}
\end{equation*}
$$

(in the sense of distributions, being $P$ the principal part) and therefore

$$
\begin{equation*}
\operatorname{Im} i G_{R}^{(2)}\left(q^{2}\right)=\pi \rho_{R}\left(q^{2}\right) \tag{2.1.27}
\end{equation*}
$$

### 2.1.3 Spectral Representation for Fermions**

Here we briefly outline how to construct the spectral representation for Fermions. The asymptotic theory discussed in sec. 2.1.1 carries over to this case, with the following normalizations [to be compared with eq. (2.1.15)]:

$$
\begin{equation*}
\langle 0| \psi(x)|\vec{p}, s\rangle=Z_{2}^{1 / 2} u_{s}(\vec{p}) \frac{e^{-i p x}}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}}}}, \quad\langle 0| \bar{\psi}(x)|\vec{p}, s, c\rangle=Z_{2}^{1 / 2} \bar{v}_{s}(\vec{p}) \frac{e^{-i p x}}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}}}}, \tag{2.1.28}
\end{equation*}
$$

with $p=\left(\omega_{\vec{p}}, \vec{p}\right), \omega_{\vec{p}}=\sqrt{\vec{p}^{2}+m^{2}}$, and where the single-particle state carries also an additional spin index $s$ and $c$ indicates the charge conjugate. In view of the canonical (anti)comutation relations, in this case one focuses on $\langle 0|\left\{\psi_{\alpha}(x), \bar{\psi}_{\beta}(0)\right\}|0\rangle$ (where $\alpha$ and $\beta$ are spinor indices) and on $\langle 0| T \psi_{\alpha}(x) \bar{\psi}_{b}(0)|0\rangle$ (remember that $T$ carries a - sign). As in the case of the scalar field one derives for $\langle 0| \psi_{\alpha}(x) \bar{\psi}_{\beta}\left(x^{\prime}\right)|0\rangle$ the analogous expression of eq. (2.1.16) with a spectral density

$$
\begin{equation*}
\rho_{\alpha \beta}(q) \equiv(2 \pi)^{3} \sum_{n}\langle 0| \psi_{\alpha}(0)|n\rangle\langle n| \bar{\psi}_{\beta}(0)|0\rangle \delta^{4}\left(q-p_{n}\right), \tag{2.1.29}
\end{equation*}
$$

which is a 4 matrix in the spinor space. Accordingly, it can be expanded in the natural basis provided by (a) the identity ! (scalar component), (b) the 4 gamma matrices $\gamma_{\mu}$ (vector), (c) the 6 matrices $\sigma_{\mu \nu}=i\left[\gamma_{\mu}, \gamma_{\nu}\right] / 2$ (tensor), (d) the matrix $\gamma_{5}$ associated to parity (pseudoscalar) and (e) the 4 matrices $\gamma_{5} \gamma_{\mu}$ (pseudovector):

$$
\begin{equation*}
\rho(q)=\rho_{S}(q) \square+\rho_{V}^{\mu}(q) \gamma_{\mu}+\rho_{T}^{\mu \nu}(q) \sigma_{\mu \nu}+\rho_{P S}(q) \gamma_{5}+\rho_{P V}^{\mu}(q) \gamma_{5} \gamma_{\mu} \tag{2.1.30}
\end{equation*}
$$

Note that, while the spectral density of the scalar field was a Lorentz scalar, in the present case one can verify that $\rho$ transforms under a boost $\Lambda$ as $\rho(\Lambda q)=S(\Lambda) \rho(q) S^{-1}(\Lambda)$, where $S(\Lambda)=e^{-i \sigma_{\mu \nu} \omega^{\mu \nu} / 4}$, being $\omega^{\mu \nu}$ the parameters of the boost. The last requirement of covariance implies that the coefficients of the expansion in eq. (2.1.30) are, respectively, scalar, vectors, tensors, pseudoscalar, and pseudovectors under Lorentz transformations and being only functions of a single vector $q^{\mu}$ can be written as

$$
\begin{equation*}
\rho_{S, P S}(q)=f_{S, P S}\left(q^{2}\right), \quad \rho_{V, P V}^{\mu}(q)=q^{\mu} f_{V, P V}\left(q^{2}\right), \quad \rho_{T}^{\mu \nu}(q)=q^{\mu} q^{\nu} f_{T}\left(q^{2}\right) \tag{2.1.31}
\end{equation*}
$$

and therefore

$$
\begin{equation*}
\rho(q)=f_{S}\left(q^{2}\right) \mathbb{\square}+f_{V}\left(q^{2}\right) d q+f_{P S}\left(q^{2}\right) \gamma_{5}+f_{P V}\left(q^{2}\right) \gamma_{5} \not q . \tag{2.1.32}
\end{equation*}
$$

As in the case of the scalar, one readily conclude on the basis of kinematics that the support of the various coefficients is in the forward light-cone and therefore, generically $f_{i}\left(q^{2}<0\right)=0$ and $f_{i}\left(q^{2}\right) \rightarrow f_{i}\left(q^{2}\right) \theta\left(q_{0}\right)$. Depending on possible additional symmetries of the theory, one can further constrain the coefficients of the expansion in eq. (2.1.32). For example, if parity is a symmetry of the theory (as we assume below), then one can formally prove $f_{P S}=f_{P V} \equiv 0$ and therefore

$$
\begin{equation*}
\langle 0| \psi_{\alpha}(x) \bar{\psi}_{\beta}(0)|0\rangle=\int \frac{d^{4} q}{(2 \pi)^{3}}\left[f_{V}\left(q^{2}\right) i \not \partial+f_{S}\left(q^{2}\right)\right]_{\alpha \beta} e^{-i q\left(x-x^{\prime}\right)} . \tag{2.1.33}
\end{equation*}
$$

In order to proceed to the spectral representation of the propagator and anticommutator, it is necessary to represent also $\langle 0| \bar{\psi}_{\beta}\left(x^{\prime}\right) \psi_{\alpha}(x)|0\rangle$ because, due to the Lorentz structure, this is not trivially related to $\langle 0| \psi_{\alpha}(x) \bar{\psi}_{\beta}\left(x^{\prime}\right)|0\rangle$ by a coordinate exchange, as it was the case for the scalar field. However, these two quantities can be related by using the fundamental symmetry CPT: involving a charge conjugation it changes $\psi$ into $\psi^{\dagger}$, which is what one needs in order to connect the two expressions. This symmetry is implemented in terms of an antiunitary operator $\theta$, such that: $\theta \psi_{\alpha}(x) \theta^{-1}=i\left(\gamma_{5}\right)_{\alpha \beta} \psi_{\beta}^{\dagger}(-x)$. Using the invariance of the vacuum under $\theta$, one can conclude (after a bit of algebra to be done carefully) that $\langle 0| \bar{\psi}_{\beta}\left(x^{\prime}\right) \psi_{\alpha}(x)|0\rangle=-\left(\gamma_{5}\langle 0| \psi(-x) \bar{\psi}\left(-x^{\prime}\right)|0\rangle \gamma_{5}\right)_{\alpha \mathrm{e} t a}$ and therefore it has the same spectral
representation as the one in eq. (2.1.33) but with an overall minus sign. As a result

$$
\begin{align*}
& \langle 0|\left\{\psi_{\alpha}(x), \bar{\psi}_{\beta}(0)\right\}|0\rangle=\int \frac{d^{4} q}{(2 \pi)^{3}}\left[f_{V}\left(q^{2}\right) i \not \subset+f_{S}\left(q^{2}\right)\right]_{\alpha \beta}\left(e^{-i q x}-e^{i q x}\right),  \tag{2.1.34}\\
& \langle 0| T \psi_{\alpha}(x) \bar{\psi}_{\beta}(0)|0\rangle=\int \frac{d^{4} q}{(2 \pi)^{3}}\left[f_{V}\left(q^{2}\right) i \notin+f_{S}\left(q^{2}\right)\right]_{\alpha \beta}\left[\theta\left(x_{0}\right) e^{-i q x}-\theta\left(-x_{0}\right) e^{i q x}\right], \tag{2.1.35}
\end{align*}
$$

to be compared with eqs. (2.1.19) and (2.1.20) and the rest of the discussion proceeds as in the case of the scalar field, taking into account that, for free fields,

$$
\begin{equation*}
\langle 0|\left\{\psi_{0, \alpha}(x), \bar{\psi}_{0, \beta}(0)\right\}|0\rangle=(i \not \partial+m)_{\alpha \beta} i \Delta_{c}^{0}\left(x ; m^{2}\right) \equiv i S_{\alpha \beta}^{0}(x, m), \tag{2.1.36}
\end{equation*}
$$

whereas the propagator is given by

$$
\begin{equation*}
\langle 0| \psi_{0, \alpha}(x) \bar{\psi}_{0, \beta}(0)|0\rangle=(i \not \partial+m)_{\alpha \beta} i \Delta_{F}^{0}\left(x ; m^{2}\right) \tag{2.1.37}
\end{equation*}
$$

where $i \Delta_{c}^{0}$ and $i \Delta_{F}^{0}$ are the same as for the scalar. For example, by introducing the additional constraint as in eq. (2.1.19) (with, for convenience $\sigma \mapsto \mu^{2}$ ), one finds

$$
\begin{align*}
\langle 0|\left\{\psi_{\alpha}(x), \bar{\psi}_{\beta}(0)\right\}|0\rangle & =\int_{0}^{\infty} d \mu^{2}\left[f_{V}\left(\mu^{2}\right) i \not \partial+f_{S}\left(\mu^{2}\right)\right]_{\alpha \beta} i \Delta_{c}^{0}\left(x ; \mu^{2}\right)  \tag{2.1.38}\\
& =\int_{0}^{\infty} d \mu^{2}\left\{f_{V}\left(\mu^{2}\right) i S_{\alpha \beta}^{0}(x ; \mu)+\left[f_{S}\left(\mu^{2}\right)-\mu f_{V}\left(\mu^{2}\right)\right] \delta_{\alpha \beta} i \Delta_{c}^{0}\left(x ; \mu^{2}\right)\right\}
\end{align*}
$$

the canonical anticommutation relation $\left.\left\{\psi_{\alpha}(x), \bar{\psi}_{\beta}\left(x^{\prime}\right)\right\}\right|_{x_{0}=x_{0}^{\prime}}=\gamma_{\alpha \beta}^{0} \delta^{3}\left(\vec{x}-\vec{x}^{\prime}\right)$ (which applies both to the free and interacting field) implies

$$
\begin{equation*}
\int_{0}^{\infty} d \mu^{2} f_{V}\left(\mu^{2}\right)=1 \tag{2.1.39}
\end{equation*}
$$

As in the case of the scalar field, one can isolate from the spectral density $\rho$ the contribution of the one-particle state and introduce the renormalized density (and therefore the renormalized $f_{V, S}$, with the result that

$$
\begin{equation*}
1=Z_{2}\left[1+\int_{m_{t h}^{2}}^{\infty} d \mu^{2} f_{V, R}\left(\mu^{2}\right)\right] \tag{2.1.40}
\end{equation*}
$$

In order to conclude that $0<Z_{2}<1$ - as we expect from an interpretation of $Z_{2}$ analogous to the one given for scalar fields - one need to prove that: (a) $f_{V, S}$ are real, (b) $f_{V} \geq 0$, and (c) $\mu f_{V}\left(\mu^{2}\right)-f_{S}\left(\mu^{2}\right) \geq 0$. These conclusions can actually be drawn with a bit of algebra, by using the fact that $\rho_{\alpha \beta}^{*}=\left[\gamma^{0} \rho \gamma^{0}\right]_{\beta \alpha}$ implies (a), $\operatorname{Tr}\left[\gamma^{0} \rho\right] \geq 0$ implies (b) and that, after having formed the modulus square of $(i \not \partial-\mu) \psi$, one infers that $\operatorname{Tr}\left[\gamma^{0}(d-\mu) \rho(q)(d-\mu)\right] \geq 0$ implies (c).

By repeating the analysis for the propagator of the field, one can conclude that

$$
\begin{equation*}
G^{(2)}(q)=Z_{2} i \frac{\not q+m}{q^{2}-m^{2}+i \epsilon}+\int_{m_{t h}^{2}}^{\infty} d \mu^{2} i Z_{2} \frac{f_{V, R}\left(\mu^{2}\right) d+f_{S, R}\left(\mu^{2}\right)}{q^{2}-\mu^{2}+i \epsilon} \tag{2.1.41}
\end{equation*}
$$

to be compared with eq. (2.1.25).

### 2.2 The Cluster Decomposition Principle and the Connected $S$-Matrix

In the previous section we briefly mentioned that the the transition amplitudes from an initial state $|i\rangle$ to a final state $|f\rangle$ can be calculated as a matrix element of the $S$-matrix, as in eq. (2.1.8). ${ }^{2}$ Let us denote by $S_{m \rightarrow n}$ a generic $S$-matrix element for the transition from an $m$-particle incoming state to an $n$-particle outgoing state. A fundamental principle in physics states that two spatially separated events should have no correlation among each other. ${ }^{3}$ In the context of $S$-matrix, this principle is called cluster decomposition. If the initial and final $m$ and $n$ particles are grouped in sets $\left\{m_{1}, \ldots m_{N}\right\},\left\{n_{1}, \ldots, n_{N}\right\}$ and the process $m \rightarrow n$ consists of a set of $N$ different sub-processes $m_{i} \rightarrow n_{i}(i=1, \ldots, N)$, occurring all far away from each other, then we should demand that

$$
\begin{equation*}
S_{m \rightarrow n}=\prod_{i=1}^{N} S_{m_{i} \rightarrow n_{i}} \tag{2.2.1}
\end{equation*}
$$

We can rephrase eq.(2.2.1) in another way by defining the connected part of the $S$-matrix, $S^{C}$. In perturbation theory, $S^{C}$ is defined in terms of connected Feynman diagrams, namely $S_{m \rightarrow n}^{C}$ is given, order by order in perturbation theory, by the diagrams where all $m$ and $n$ particles are connected with each other. A non-perturbative definition for $S_{m \rightarrow m}^{C}$ can be given iteratively starting from the $1 \rightarrow 1$ process for which

$$
\begin{equation*}
S_{p \rightarrow p^{\prime}}=S_{p \rightarrow p^{\prime}}^{C}=\delta\left(p-p^{\prime}\right) \tag{2.2.2}
\end{equation*}
$$

where $\delta\left(p-p^{\prime}\right) \equiv \delta\left(\vec{p}-\vec{p}^{\prime}\right) \delta_{s, s^{\prime}} \delta_{n, n^{\prime}}$ includes Kronecker or Dirac delta's in momentum, spin and particle species. The $2 \rightarrow 2$ connected matrix element $S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}}^{C}$ is given as

$$
\begin{align*}
S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}} & =S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}}^{C}+S_{p_{1} \rightarrow p_{1}^{\prime}}^{C} S_{p_{2} \rightarrow p_{2}^{\prime}}^{C} \pm S_{p_{1} \rightarrow p_{2}^{\prime}}^{C} S_{p_{2} \rightarrow p_{1}^{\prime}}^{C} \\
& =S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}}^{C}+\delta\left(p_{1}-p_{1}^{\prime}\right) \delta\left(p_{3}-p_{3}^{\prime}\right) \pm \delta\left(p_{1}-p_{2}^{\prime}\right) \delta\left(p_{2}-p_{1}^{\prime}\right) \tag{2.2.3}
\end{align*}
$$

[^1]where $\pm$ refers to an even or odd exchange of fermion particles. Similarly, the $3 \rightarrow 3$ connected matrix element $S_{p_{1}, p_{2}, p_{3} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}, p_{3}^{\prime}}^{C}$ is given as
\[

$$
\begin{align*}
S_{p_{1}, p_{2}, p_{3} \rightarrow p_{1}, p_{2}, p_{3}}= & S_{p_{1}, p_{2}, p_{3} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}, p_{3}^{\prime}}+\left(S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}}^{C} \delta\left(p_{3}-p_{3}^{\prime}\right) \pm 8 \text { perms. }\right)  \tag{2.2.4}\\
& +\left(\delta\left(p_{1}-p_{1}^{\prime}\right) \delta\left(p_{2}-p_{2}^{\prime}\right) \delta\left(p_{3}-p_{3}^{\prime}\right) \pm 5 \text { perms. }\right),
\end{align*}
$$
\]

where perms. refer to all other possible ways to combine the in-coming and the out-coming particles. By further iterating, we can define all higher connected matrix elements $S_{m \rightarrow m}^{C}$ with $m>3$. Similarly, we could define the "off-diagonal" connected elements $S_{m \rightarrow n}^{C}$ with $m \neq n$. In terms of $S^{C}$, cluster decomposition implies that

$$
\begin{equation*}
S_{m \rightarrow n}^{C}=0 \tag{2.2.5}
\end{equation*}
$$

if any among the $m+n$ particles is far away from the remaining ones. For instance, suppose that a $3 \rightarrow 3$ scattering consists of an actual $2 \rightarrow 2$ scattering involving particles 1 and 2 , with particle 3 far away from the interaction place. In this case, we would expect from eq.(2.2.2)

$$
\begin{equation*}
S_{3 \rightarrow 3}=S_{2 \rightarrow 2} S_{1 \rightarrow 1}=S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}} \delta\left(p_{3}-p_{3}^{\prime}\right), \tag{2.2.6}
\end{equation*}
$$

with $S_{p_{1}, p_{2} \rightarrow p_{1}^{\prime}, p_{2}^{\prime}}$ given by eq.(2.2.3). When we impose eq.(2.2.5), namely we demand that all connected $S$ matrix elements where particle 3 enters together with particles 1 and 2 vanish, leaves only three terms in eq.(2.2.4), that indeed recombine in the right-hand-side of eq.(2.2.6).

Since cluster decomposition is formulated in configuration space, we can define a configuration space connected $S$-matrix $S_{x, x^{\prime}}^{C}$ by taking the Fourier transform of $S_{m \rightarrow n}^{C}$ :

$$
\begin{equation*}
S_{m \rightarrow n}^{C}\left(x_{1}, \ldots x_{m}, x_{1}^{\prime}, \ldots x_{n}^{\prime}\right) \equiv \int \prod_{i=1}^{m} d^{3} p_{i} \prod_{j=1}^{n} d^{3} p_{j}^{\prime} e^{i \overrightarrow{p_{i}} \cdot \overrightarrow{x_{i}}-i \overrightarrow{p_{j}^{\prime}} \cdot \overrightarrow{x_{j}^{\prime}}} S_{m \rightarrow n}^{C}\left(p_{1}, \ldots p_{m}, p_{1}^{\prime}, \ldots p_{n}^{\prime}\right) \tag{2.2.7}
\end{equation*}
$$

Invariance under space-time translations imply that $S_{m \rightarrow n}^{C}\left(x, x^{\prime}\right)$ cannot vary if we shift by a constant four-vector all coordinates. Correspondingly the connected $S$-matrix should be proportional to a delta function imposing energy-momentum conservation:

$$
\begin{equation*}
S_{m \rightarrow n}^{C}\left(p_{1}, \ldots p_{m}, p_{1}^{\prime}, \ldots p_{n}^{\prime}\right)=(2 \pi)^{4} \delta^{(4)}\left(\sum_{i} p_{i}-\sum_{j} p_{j}^{\prime}\right) A^{C}\left(p_{1}, \ldots p_{m}, p_{1}^{\prime}, \ldots p_{n}^{\prime}\right) \tag{2.2.8}
\end{equation*}
$$

with $A^{C}$ encoding the remaining part of the scattering process. Cluster decomposition implies that

$$
\begin{equation*}
S_{m \rightarrow n}^{C}\left(x_{1}, \ldots x_{m}, x_{1}^{\prime}, \ldots x_{n}^{\prime}\right) \rightarrow 0 \tag{2.2.9}
\end{equation*}
$$

if we move apart the in-coming particles $\left|\overrightarrow{x_{i}}-\overrightarrow{x_{j}}\right| \rightarrow \infty, i=1,2, \ldots, m_{1}, j=1,2, \ldots m-$ $m_{1}$, and their associated out-going particles $\left|\overrightarrow{x_{k}}{ }^{\prime}-\overrightarrow{x_{l}}\right| \rightarrow \infty, k=1,2, \ldots, n_{1}\left(m_{1}\right), l=$ $1,2, \ldots n-n_{1}\left(m_{1}\right)$, for any choice of $m_{1}$.

No other delta functions imposing further constraints on the initial and outgoing momenta can enter in $A^{C}$. If such a constraint would exist, we could move all the particles subject to the constraint away from the others without affecting eq.(2.2.7), violating the cluster decomposition property (2.2.9). In contrast, the disconnected element of the full $S$-matrix can and do have multiple delta functions. This shows that the connected and disconnected elements of the $S$-matrix cannot interfere among each other. We will mostly focus in what follows to the connected part of the $S$-matrix, but it should be emphasized that disconnected $S$-matrix elements are not negligible, since in actual experiments the scattering process is often of this type.

### 2.3 The Reduction Formula for Connected S-Matrix Elements

Scattering experiments are determined by the $S$-matrix. However, quantum field theory naturally provides access to time-ordered products of fields and operators and therefore it is important to understand how to extract from the latter the $S$-matrix elements. This connection is provided by the so-called Lehmann-Symanzik-Zimmermann (LSZ) reduction formulas $[4,5]$ which we discuss here for the case of a neutral scalar field.

Consider a generic $S$-matrix element for the transition from the $m$-particle incoming state $\left|\vec{p}_{1}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}$ to the $n$-particle outgoing state $\left|\vec{q}_{1}, \ldots \vec{q}_{n}\right\rangle_{\text {out }}$ :

$$
\begin{equation*}
S_{m \rightarrow n}={ }_{\text {out }}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n} \mid \vec{p}_{1}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}={ }_{\text {in }}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right| S\left|\vec{p}_{1}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}, \tag{2.3.1}
\end{equation*}
$$

where $\left|\vec{p}_{1}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}=a^{\dagger}\left(\vec{p}_{1}\right)_{\text {in }}\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}$ (see sec. 2.1.1). In turn, $a(\vec{p})_{\text {in (out) }}$ (and $\left.a^{\dagger}(\vec{p})_{\text {in(out) }}\right)$ can be expressed in terms of the incoming (outgoing) free field $\phi_{\text {in(out) }}$ as in eq.(2.1.6). We now write $a_{\vec{p}, \text { in }}^{\dagger}=\left(a_{\vec{p}, \text { in }}^{\dagger}-a_{\vec{p}, \text { out }}^{\dagger}\right)+a_{\vec{p}, \text { out }}^{\dagger}$ and therefore
$S_{m \rightarrow n}={ }_{\text {out }}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right|\left(a^{\dagger}\left(\vec{p}_{1}\right)_{\text {in }}-a^{\dagger}\left(\vec{p}_{1}\right)_{\text {out }}\right)\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}+{ }_{\text {out }}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right| a^{\dagger}\left(\vec{p}_{1}\right)_{\text {out }}\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}$,
where the second term on the r.h.s. does not vanish only if $\vec{p}_{1} \in\left\{\vec{q}_{1}, \ldots \vec{q}_{n}\right\}$, i.e., if it coincides with one of the outgoing momenta. This contribution represents a process in which one of the incoming particles does not participate in the scattering and can be neglected if we are only interested in connected $S$-matrix elements. In the first contribution, instead, one can use eq. (2.1.6) in order to express $a^{\dagger}\left(\vec{p}_{1}\right)_{\text {in }}$ and $a^{\dagger}\left(\vec{p}_{1}\right)_{\text {out }}$ in terms of $\phi_{\text {in }}(x)$ and $\phi_{\text {out }}(x)$, respectively, and the latter, in turn, as $\lim _{x_{0} \rightarrow-\infty} Z^{-1 / 2} \phi(x)$ and
$\lim _{x_{0} \rightarrow+\infty} Z^{-1 / 2} \phi(x)$ [see eq. (2.1.1)]:

$$
\begin{equation*}
S_{m \rightarrow n}^{C}=-i Z^{-1 / 2} \int d^{3} \vec{x}\left(\lim _{x_{0} \rightarrow-\infty}-\lim _{x_{0} \rightarrow+\infty}\right){ }_{\text {out }}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right| h_{\vec{p}_{1}}^{*}(x) \stackrel{\leftrightarrow}{\partial_{0}} \phi(x)\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }} \tag{2.3.3}
\end{equation*}
$$

The difference between the limits can be obtained as $-\int_{-\infty}^{+\infty} d x_{0} \partial_{0}$ which renders an integration over space-time $\int d^{4} x$, while $\partial_{0}\left[h_{\vec{p}_{1}}^{*}(x) \overleftrightarrow{\partial_{0}} \phi(x)\right]=h_{\vec{p}_{1}}^{*}(x)\left(\square_{x}+m^{2}\right) \phi(x)$ and therefore

$$
\begin{equation*}
S_{m \rightarrow n}^{C}=i Z^{-1 / 2} \int d^{4} x h_{\vec{p}_{1}}^{*}(x)\left(\square_{x}+m^{2}\right)_{o u t}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right| \phi(x)\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }} \tag{2.3.4}
\end{equation*}
$$

which has been now reduced to the matrix element of the field $\phi(x)$ with one less particle in the incoming state. An analogous reduction can be done for the outgoing state by expressing out $\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right|$ as ${ }_{\text {out }}\left\langle\vec{q}_{2}, \ldots \vec{q}_{n}\right| a\left(\vec{q}_{1}\right)_{\text {out }}$ and by using the fact that $a\left(\vec{q}_{1}\right)_{\text {out }} \phi(x)=$ $\left[a\left(\vec{q}_{1}\right)_{\text {out }} \phi(x)-\phi(x) a\left(\overrightarrow{q_{1}}\right)_{\text {in }}\right]+\phi(x) a\left(\overrightarrow{q_{1}}\right)_{\text {in }}$; as before, when inserted in eq. (2.3.4), the second term on the r.h.s. of this identity corresponds to a disconnected amplitude and can be neglected. The first, instead, can be simplified by expressing $a\left(\vec{q}_{1}\right)_{o u t / i n}$ via eq. (2.1.6) and then by introducing the suitable (weak) limits in order to replace $\phi$ by $\phi_{\text {in/out }}$

$$
\begin{align*}
& a\left(\vec{q}_{1}\right)_{\text {out }} \phi(x)-\phi(x) a\left(\overrightarrow{q_{1}}\right)_{\text {in }} \\
& \quad=i Z^{-1 / 2} \int d^{3} \vec{y}\left[\lim _{y_{0} \rightarrow+\infty} h_{\vec{q}_{1}}(y) \overleftrightarrow{\partial_{y_{0}}} \phi(y) \phi(x)-\lim _{y_{0} \rightarrow-\infty} h_{\vec{q}_{1}}(y) \overleftrightarrow{\partial_{y_{0}}} \phi(x) \phi(y)\right] ; \tag{2.3.5}
\end{align*}
$$

due to the presence of the limits, both $\phi(y) \phi(x)$ and $\phi(x) \phi(y)$ can be replaced by their timeordered product $T \phi(x) \phi(y)$ and therefore the difference between the limits, as before, can be expressed as $\int_{-\infty}^{+\infty} d y_{0} \partial_{y_{0}}$, while $\partial_{y_{0}}\left[h_{\vec{q}_{1}}(y) \overleftrightarrow{\partial_{y_{0}}} T \phi(x) \phi(y)\right]=h_{\vec{q}_{1}}(y)\left(\square_{y}+m^{2}\right) T \phi(x) \phi(y)$, with the result

$$
\begin{align*}
& { }_{\text {out }}\left\langle\vec{q}_{1}, \ldots \vec{q}_{n}\right| \phi(x)\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}= \\
& \quad=i Z^{-1 / 2} \int d^{4} y h_{\vec{q}_{1}}(y)\left(\square_{y}+m^{2}\right)_{\text {out }}\left\langle\vec{q}_{2}, \ldots \vec{q}_{n}\right| T \phi(x) \phi(y)\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }} \tag{2.3.6}
\end{align*}
$$

in which the number of particles in the outgoing state has been reduced by one. Together with eq. (2.3.4) one finds

$$
\begin{array}{r}
S_{m \rightarrow n}^{C}=\left(i Z^{-1 / 2}\right)^{2} \int d^{4} x d^{4} y h_{\vec{p}_{1}}^{*}(x) h_{\vec{q}_{1}}(y)\left(\square_{x}+m^{2}\right)\left(\square_{y}+m^{2}\right)  \tag{2.3.7}\\
\times{ }_{\text {out }}\left\langle\overrightarrow{q_{2}}, \ldots \vec{q}_{n}\right| T \phi(x) \phi(y)\left|\vec{p}_{2}, \ldots \vec{p}_{m}\right\rangle_{\text {in }}
\end{array}
$$

in which the procedure outlined above can be iteratively repeated till all the particles in the incoming and outgoing states are reduced to zero and one is left with the vacuum state

$$
\begin{align*}
& |0\rangle_{\text {in }}=|0\rangle_{\text {out }}=|0\rangle: \\
& S_{m \rightarrow n}^{C}=\left(i Z^{-1 / 2}\right)^{m+n} \int \prod_{i=1}^{m} d^{4} x_{i} h_{\vec{p}_{i}}^{*}\left(x_{i}\right)\left(\square_{x_{i}}+m^{2}\right) \prod_{j=1}^{n} d^{4} y_{j} h_{\vec{q}_{j}}\left(y_{j}\right)\left(\square_{y_{j}}+m^{2}\right)  \tag{2.3.8}\\
& \\
& \times\langle 0| T \phi\left(x_{1}\right) \cdots \phi\left(x_{m}\right) \phi\left(y_{1}\right) \cdots \phi\left(y_{n}\right)|0\rangle,
\end{align*}
$$

which is the LSZ reduction formula. ${ }^{4}$
In order to connect the r.h.s. of this equation to the $N$-point correlation function $\langle 0| T \phi\left(x_{1}\right) \cdots \phi\left(x_{N}\right)|0\rangle$ of the fields which is usually calculated in quantum field theory, we focus on its Fourier transform

$$
\begin{equation*}
G^{(N)}\left(p_{1}, \ldots, p_{N}\right)=\int \prod_{k=1}^{N} d^{4} x_{k} e^{-i p_{k} x_{k}}\langle 0| T \phi\left(x_{1}\right) \cdots \phi\left(x_{N}\right)|0\rangle \tag{2.3.9}
\end{equation*}
$$

Due to the space-time translational invariance of the theory, the correlation function on the r.h.s. depends only on the difference of the coordinates and therefore one can always factor out $(2 \pi)^{4} \delta^{4}\left(\sum_{k=1}^{N} p_{k}\right)$ and actually indicate by $G^{(N)}$ the remaining factor. Taking into account the definition of $h_{\vec{p}}(x)$ in eq. (2.1.6) one easily recognizes in eq. (2.3.8) $(2 \pi)^{4} \delta^{4}\left(\sum_{i=1}^{m} p_{i}-\sum_{j=1}^{n} q_{j}\right) G^{(n+m)}\left(p_{1}, \ldots, p_{m},-q_{1}, \ldots,-q_{n}\right)$, while $\square_{x_{i}}$ and $\square_{y_{j}}$ turn into $-p_{i}^{2}$ and $-q_{j}^{2}$ when acting to the left on $h_{\vec{p}_{i}}^{*}\left(x_{i}\right)$ and $h_{\vec{q}_{j}}\left(y_{j}\right)$, respectively. Note, however, that according to eq. (2.1.6) $p_{i}^{2}=q_{j}^{2}=m^{2}$, being the incoming and outgoing particles on-shell, and therefore the factors $\square_{x_{i} / y_{j}}+m^{2}$ vanish; it is then convenient to think of them as limits:

$$
\begin{align*}
S_{m \rightarrow n}^{C} & =(2 \pi)^{4} \delta^{4}\left(\sum_{i=1}^{m} p_{i}-\sum_{j=1}^{n} q_{j}\right) \prod_{i=1}^{m} \frac{1}{\sqrt{(2 \pi)^{3} 2 \omega_{p_{i}}}} \prod_{j=1}^{n} \frac{1}{\sqrt{(2 \pi)^{3} 2 \omega_{q_{j}}}}  \tag{2.3.10}\\
& \times \prod_{i=1}^{m} \lim _{p_{i}^{2} \rightarrow m^{2}} \frac{p_{i}^{2}-m^{2}}{i Z^{1 / 2}} \prod_{j=1}^{n} \lim _{q_{j}^{2} \rightarrow m^{2}} \frac{q_{j}^{2}-m^{2}}{i Z^{1 / 2}} G^{(n+m)}\left(p_{1}, \ldots, p_{m},-q_{1}, \ldots,-q_{n}\right)
\end{align*}
$$

This equation clearly shows that connected $S$-matrix elements are basically given by the residues of the poles of the Fourier transform of the $T$-product of fields (i.e., of their Green functions) as the incoming and outgoing particles go on-shell. As we discussed in sec. 2.1.2, $G^{(2)}\left(p^{2} \rightarrow m^{2}\right)=i Z /\left(p^{2}-m^{2}\right)$, which allows one to identify the factors $\lim _{p^{2} \rightarrow m^{2}}\left(p^{2}-m^{2}\right) /\left(i Z^{1 / 2}\right)$ in eq. (2.3.10) as $\sqrt{Z}\left[G^{(2)}\left(p^{2} \rightarrow m^{2}\right)\right]^{-1}$. In view of this fact it is convenient to introduce the notion of amputated Green function $G_{a m p}^{(N)}$, which in terms of Feynman diagrams, is obtained by stripping the diagram corresponding to $G^{(N)}$ of the

[^2]external propagators (the latter being the cause of the presence of the poles mentioned above):
\[

$$
\begin{equation*}
G^{(N)}\left(p_{1}, \ldots, p_{N}\right)=\left[\prod_{i=1}^{N} G^{(2)}\left(p_{i}\right)\right] G_{a m p}^{(N)}\left(p_{1}, \ldots, p_{N}\right) \tag{2.3.11}
\end{equation*}
$$

\]

Accordingly, eq. (2.3.10) becomes

$$
\begin{align*}
& S_{m \rightarrow n}^{C}=(2 \pi)^{4} \delta^{4}\left(\sum_{i=1}^{m} p_{i}-\sum_{j=1}^{n} q_{j}\right) \prod_{i=1}^{m} \frac{1}{\sqrt{(2 \pi)^{3} 2 \omega_{\vec{p}_{i}}}} \prod_{j=1}^{n} \frac{1}{\sqrt{(2 \pi)^{3} 2 \omega_{q_{j}}}}  \tag{2.3.12}\\
& \times\left. Z^{(n+m) / 2} G_{a m p}^{(n+m)}\left(p_{1}, \ldots, p_{m},-q_{1}, \ldots,-q_{n}\right)\right|_{p_{i}^{2}=q_{j}^{2}=m^{2}} .
\end{align*}
$$

Quantum field theory provides access (e.g., perturbatively) to $G_{a m p}^{(N)}$ on the r.h.s. and therefore, via this relation, to $S_{m \rightarrow n}^{C}$ which has to be supplemented by appropriate phase space (i.e., kinematic) factors in order to calculate cross sections.

An important consequence of eq. (2.3.12) is the crossing symmetry: in fact, incoming particles are distinguished from the outgoing ones only by the sign of the four-momentum as the argument of $G_{a m p}^{(n+m)}$. When the theory involves also antiparticles, the construction can be straightforwardly generalized in order to conclude that an incoming antiparticle with momentum $\vec{p}$ [and four-momentum $p=\left(-\omega_{\vec{p}}, \vec{p}\right)$ ] is equivalent to an outgoing particle with momentum $-\vec{p}$ [i.e., with four momentum $\left.\left(\omega_{\vec{p}},-\vec{p}\right)=-p\right]$.

### 2.4 The Optical Theorem

The $S$-matrix is a unitary mapping between the space of the incoming asymptotic states and the one of the outgoing states. As such, it satisfies $S^{\dagger} S=S S^{\dagger}=1$. This unitarity has important consequences, which we investigate here.

In order to factor out from the $S$-matrix the contribution corresponding to the absence of interaction between the incoming and the outgoing particles, one usually introduces the $T$-matrix:

$$
\begin{equation*}
S=1+i T \tag{2.4.1}
\end{equation*}
$$

in terms of which, the condition of unitarity becomes

$$
\begin{equation*}
T-T^{\dagger}=i T^{\dagger} T \tag{2.4.2}
\end{equation*}
$$

Consider now a scattering process from an initial state $|i\rangle$ to a final state $|f\rangle$ : the previous relation implies that

$$
\begin{equation*}
T_{f, i}-T_{f, i}^{\dagger}=i \sum_{n} T_{f, n}^{\dagger} T_{n, i} \tag{2.4.3}
\end{equation*}
$$

for the $T$-matrix element $T_{f, i} \equiv\langle f| T|i\rangle$ between these states, where the sum runs over all possible intermediate states with definite momentum. Note that the conservation of energy and momentum requires generically

$$
\begin{equation*}
T_{a, b}=\langle a| T|b\rangle=(2 \pi)^{4} \delta^{4}\left(p_{a}-p_{b}\right) A_{a, b} \tag{2.4.4}
\end{equation*}
$$

where $p_{a}$ and $p_{b}$ are the momenta of states $|a\rangle$ and $|b\rangle$, respectively, and $A_{a, b}$ is the remaining factor usually calculated via Feynman diagrams. In terms of $A_{a, b}$, eq. (2.4.3) becomes

$$
\begin{equation*}
A_{f, i}-A_{i, f}^{*}=i \sum_{n}(2 \pi)^{4} \delta^{4}\left(p_{n}-p_{f(i)}\right) A_{n, f}^{*} A_{n, i} \tag{2.4.5}
\end{equation*}
$$

This relation has a very simple interpretation for $f=i$, i.e., when focussing on the forward scattering amplitude because in this case the l.h.s. is $2 i \operatorname{Im} A_{i, i}$, while the r.h.s. involves $\left|A_{n, i}\right|^{2}$ :

$$
\begin{equation*}
2 \operatorname{Im} A_{i, i}=\sum_{n}(2 \pi)^{4} \delta^{4}\left(p_{n}-p_{i}\right)\left|A_{n, i}\right|^{2} \tag{2.4.6}
\end{equation*}
$$

Consider the case in which the initial state $|i\rangle=\left|\vec{k}_{1}, \vec{k}_{2}\right\rangle$ is a two-particle state: the r.h.s of eq. (2.4.6) is then the sum over all possible final states of the probability to obtain these states from the scattering of the two initial particles. This is related to the total cross section $\sigma_{T}$. In our conventions, we have ${ }^{5}$

$$
\begin{equation*}
\sigma_{T}=\frac{(2 \pi)^{6} 2 E_{1} 2 E_{2}}{\sqrt{\left(p_{1} \cdot p_{2}\right)^{2}-m_{1}^{2} m_{2}^{2}}} \sum_{n} \prod_{i=1}^{n} \int d^{3} p_{i}^{\prime}(2 \pi)^{4} \delta\left(p_{1}+p_{2}-\sum_{i} p_{i}^{\prime}\right)\left|A_{i, n}\right|^{2} \tag{2.4.7}
\end{equation*}
$$

In the center of mass frame of the two particles we get

$$
\begin{equation*}
\frac{4\left|\vec{k}_{c m}\right|}{(2 \pi)^{6} E_{c m}} \sigma_{T}=\sum_{n}(2 \pi)^{4} \delta^{4}\left(p_{1}+p_{2}-p_{n}^{\prime}\right)\left|A_{n, i}\right|^{2} \tag{2.4.8}
\end{equation*}
$$

where we have rewritten more compactly the phase space integral as $\sum_{n}$. Accordingly, the optical theorem implies that

$$
\begin{equation*}
2 \operatorname{Im} A_{i, i}=\frac{4\left|\vec{k}_{c m}\right|}{(2 \pi)^{6} E_{c m}} \sigma_{T} \tag{2.4.9}
\end{equation*}
$$

Further below we shall check explicitly this relation. On this specific case of a two-particle initial state we can discuss a general property of the amplitude $A_{i, i}$ as a function of the Mandelstam variable $s=E_{c m}^{2}$ which we let be a complex number. First of all we note that the diagrams which contribute to $A_{i, i}$ in a perturbative expansion are generically real as

[^3]long as one can forget about the Feynman prescription of adding $+i \epsilon$ at the denominators of the propagators of the particles involved (think, for example, of scalars). This is the case if none of the intermediate virtual particles involved in the diagram goes on-shell: in fact, being the denominator of the corresponding propagator always non-vanishing, one can safely take $\epsilon \rightarrow 0$, which makes the imaginary part of the corresponding diagram vanish. Accordingly, the amplitude $A_{i, i}(s)$ is a real function of $s$ for real $s<m_{t h}^{2}$, where $m_{t h}^{2}$ is the energy threshold for the production of on-shell multiparticle states:
\[

$$
\begin{equation*}
A_{i, i}^{*}(s)=A_{i, i}\left(s^{*}\right) \tag{2.4.10}
\end{equation*}
$$

\]

Since $A_{i, i}$ is real on the real axis for $s<m_{t h}^{2}$, its continuation satisfies eq. (2.4.10), i.e.,

$$
\left\{\begin{array}{l}
\operatorname{Re} A_{i, i}(s)=\operatorname{Re} A_{i, i}\left(s^{*}\right)  \tag{2.4.11}\\
\operatorname{Im} A_{i, i}(s)=-\operatorname{Im} A_{i, i}\left(s^{*}\right)
\end{array}\right.
$$

As a consequence of the continuation outside the real axis, $A_{i, i}(s)$ acquires an imaginary part: upon approaching the real axis $s \in \mathbb{R}$ from above with $s+i \epsilon$ and $\epsilon=0^{+}$or from below with $s-i \epsilon$, eq. (2.4.11) implies that $A_{i, i}(s)$ has opposite imaginary parts $\operatorname{Im} A_{i, i}(s+i \epsilon)=-\operatorname{Im} A_{i, i}(s-i \epsilon)$ and therefore, upon crossing the real axis, there is a discontinuity

$$
\begin{equation*}
\operatorname{Disc} A_{i, i}(s)=A_{i, i}(s+i \epsilon)-A_{i, i}(s-i \epsilon)=2 i \operatorname{Im} A_{i, i}(s+i \epsilon) \tag{2.4.12}
\end{equation*}
$$

which vanishes along the real axis with $s<m_{t h}^{2}$. Accordingly, the imaginary part of the amplitude $A_{i, i}$ is alternatively given by its discontinuity upon crossing the real axis, which might be easier to determine in actual calculations. Note that the presence of this discontinuity signals the presence of a branch cut running along the real axis for $s>m_{t h}^{2}$, which is analogous to the one highlighted in fig. 2.1.

### 2.4.1 Perturbative Unitarity

The results obtained so far are exact, in the sense that they do not rely on a perturbative expansion. Exact results in QFT are however quite rare (at least in $d=4$ space-time dimensions). In practice we often demand perturbative unitarity, namely we expand in some small coupling constant both the left and right hand side of eq.(2.4.9) (or more generally of eq.(2.4.6)) and demand the equality of both terms at each order in perturbation theory. In order to calculate directly the discontinuity of an amplitude across the cut, one can take advantage of the so-called Cutkosky (or cutting) rules:


Figure 2.2: (a) Lowest-order correction to the two-particle scattering amplitude $A_{i, i}$ in a $\lambda \phi^{4}$ theory. (b) Schematic representation of the optical theorem (2.4.6) applied to the diagram in panel (a): the imaginary part of the amplitude in panel (a) can be obtained via the cutting rules by cutting the diagram as indicated in (b) by the dashed line.

1. consider all possible ways of cutting the internal propagators of a diagram in such a way that the diagram is cut into two disconnected parts and that the particles whose propagators have been cut can go simultaneously on-shell;
2. in each cut propagator with momentum $p$ substitute the factor $1 /\left(p^{2}-m^{2}+i \epsilon\right)$ with $-2 \pi i \delta\left(p^{2}-m^{2}\right)$ and calculate the resulting diagram;
3. sum over all possible cuts the contributions obtained above.

As an illustrative example, consider the scattering amplitude $i A_{2,2}\left(k^{2}\right)$ for $2-2$ particles with a $\lambda \phi^{4}$ interaction. At the lowest (tree) non-trivial order in perturbation theory, the amplitude is proportional to the coupling $\lambda$ and is manifestly always real. This is in agreement with eq.(2.4.9), since an imaginary part of order $\lambda$ would not match with the right hand side which is of order $\lambda^{2}$. An imaginary part in the amplitude should then arise starting from one-loop. The relevant diagram contribution $i \mathcal{D}\left(k^{2}\right)$ is depicted in fig. 2.2(a). According to the LSZ formula (2.3.12), $i A_{2,2}\left(k^{2}\right)=i \mathcal{D}\left(k^{2}\right) /\left[(2 \pi)^{6} k^{2}\right]$ (with $k=k_{1}+k_{2}$ ), where

$$
\begin{equation*}
i \mathcal{D}\left(k^{2}\right)=\frac{(i \lambda)^{2}}{2} \int \frac{d^{4} q}{(2 \pi)^{4}} \frac{i}{\left(\frac{k}{2}-q\right)^{2}-m^{2}+i \epsilon} \frac{i}{\left(\frac{k}{2}+q\right)^{2}-m^{2}+i \epsilon} \tag{2.4.13}
\end{equation*}
$$

One can easily realize that among the three possible scattering channels $s, t$, and $u$ only in the channel $s$ it is possible to exceed the threshold for particle production such that the corresponding amplitude acquires an imaginary part. We shall focus on this case. By

Lorentz invariance $A_{2,2}$ can be calculated in the center-of-mass frame with $k=\left(k_{0}, \overrightarrow{0}\right)$ and it is convenient for the present purposes to perform first the integration over $q_{0}$ using Cauchy's theorem, closing the contour at infinity. The integrand is characterized by 4 simple poles at

$$
\begin{equation*}
\frac{k_{0}}{2}+\omega_{\vec{q}}-i \epsilon, \quad-\frac{k_{0}}{2}+\omega_{\vec{q}}-i \epsilon, \quad-\frac{k_{0}}{2}-\omega_{\vec{q}}+i \epsilon, \quad \frac{k_{0}}{2}-\omega_{\vec{q}}+i \epsilon \tag{2.4.14}
\end{equation*}
$$

After the integration one finds

$$
\begin{equation*}
\mathcal{D}\left(k^{2}\right)=\frac{\lambda^{2}}{2} \int \frac{d^{3} \vec{q}}{(2 \pi)^{3}} \frac{1}{4 k_{0} \omega_{\vec{q}}}\left(\frac{1}{\omega_{\vec{q}}-\frac{k_{0}}{2}-i \epsilon}-\frac{1}{\omega_{\vec{q}}+\frac{k_{0}}{2}-i \epsilon}\right) \tag{2.4.15}
\end{equation*}
$$

Using eq.(2.1.26), an imaginary part can arise only when one of the denominators vanishes. It is evident that only the first of the two terms on the r.h.s. can possibly vanish (we take $\left.k^{0}>0\right)$ when $\omega_{\vec{q}}=k_{0} / 2$ for some values of $|\vec{q}|$. This is possible only if $k_{0} \geq 2 m \equiv m_{t h}$, as expected on kinematic reasons. We then get

$$
\begin{align*}
\operatorname{Im} \mathcal{D}\left(k^{2}\right) & =\frac{\lambda^{2}}{16 \pi^{2} k_{0}} \int_{m}^{\infty} d \omega_{\vec{q}} \sqrt{\omega_{\vec{q}}^{2}-m^{2}} \times \pi \delta\left(\omega_{\vec{q}}-\frac{k_{0}}{2}\right)  \tag{2.4.16}\\
& =\frac{\lambda^{2}}{16 \pi k_{0}} \sqrt{\left(k_{0} / 2\right)^{2}-m^{2}} \theta\left(k_{0}-2 m\right)
\end{align*}
$$

where we used the fact that $d^{3} \vec{q}=4 \pi q^{2} d q=4 \pi \sqrt{\omega_{\vec{q}}^{2}-m^{2}} \omega_{\vec{q}} d \omega_{\vec{q}}$. Note also that for $k^{0}<2 m$ the location of the poles (2.4.14) allows us to perform a Wick's rotation $q_{0}=i q_{E}$. The two propagators are now complex conjugates of each other and the amplitude $A_{2,2}$ is manifestly real, in agreement with eq.(2.4.10).

The same result can be obtained by using Cutvosky rules. In this case we have

$$
\begin{equation*}
\operatorname{Disc} \mathcal{D}\left(k^{2}\right)=\frac{\lambda^{2}}{2 i} \int \frac{d^{4} q}{(2 \pi)^{4}}(-2 \pi i)^{2} \delta\left((k / 2-q)^{2}-m^{2}\right) \delta\left((k / 2+q)^{2}-m^{2}\right) \tag{2.4.17}
\end{equation*}
$$

where, as before, it is convenient to integrate first on $q_{0}$ in the center-of-mass frame. Using the fact that we can express the second $\delta$ above as

$$
\begin{equation*}
\delta\left(\left(k_{0} / 2+q_{0}\right)^{2}-\omega_{\vec{q}}^{2}\right)=\frac{\delta\left(q_{0}+k_{0} / 2-\omega_{\vec{q}}\right)}{2 \omega_{\vec{q}}}+\frac{\delta\left(q_{0}+k_{0} / 2+\omega_{\vec{q}}\right)}{2 \omega_{\vec{q}}} \tag{2.4.18}
\end{equation*}
$$

one immediately sees that the product of the first $\delta$ on the r.h.s. of eq. (2.4.17) with the second term above can never vanish, while the product with the first term, after integration on $q_{0}$, is equivalent to $\delta\left(k_{0}\left(k_{0}-2 \omega_{\vec{q}}\right)\right)=\delta\left(\omega_{\vec{q}}-k_{0} / 2\right) /\left(2 k_{0}\right)$, such that eq. (2.4.17) becomes

$$
\begin{align*}
\operatorname{Disc} \mathcal{D}\left(k^{2}\right) & =i \frac{\lambda^{2}}{8 \pi k_{0}} \int_{m}^{\infty} d \omega_{\vec{q}} \sqrt{\omega_{\vec{q}}^{2}-m^{2}} \delta\left(\omega_{\vec{q}}-k_{0} / 2\right)  \tag{2.4.19}\\
& =2 i \operatorname{Im} \mathcal{D}\left(k^{2}\right)
\end{align*}
$$

where on the second line we have taken into account the first line of eq. (2.4.16). This example shows the effectiveness of the rules mentioned above for the calculation of the imaginary part of an amplitude or, alternatively, of its discontinuity across the cut.

As a final check, let us consider the total cross section $\sigma_{T}$ for the lowest-order scattering process of these two particles, which is characterized by the amplitude $i \mathcal{D}_{0}=-i \lambda$. According to the optical theorem (2.4.9) and to (2.4.12), the imaginary part of the diagram in fig. 2.2(a) - which can be calculated (up to an $i$ ) by cutting the diagram as indicated by the dashed line fig. 2.2(b) - is equal to the modulus square of the tree-level scattering amplitude (up to the corresponding phase-space), as schematically indicated by fig. 2.2(b). By using the standard expression for the differential cross-section in a two-body scattering (see, e.g., eq. (4.99) in ref. [1]), one has

$$
\begin{equation*}
\left(\frac{d \sigma}{d \Omega}\right)_{c m}=\frac{|\lambda|^{2}}{64 \pi^{2} E_{c m}^{2}} \times \frac{1}{2} \tag{2.4.20}
\end{equation*}
$$

where the factor $1 / 2$ accounts for the fact that the particles are assumed indistinguishable and which gives $\sigma_{T}=\lambda^{2} /\left(32 \pi E_{c m}^{2}\right)$ for the total cross section. By forming the combination (with $E_{c m}=k_{0}$ )

$$
\begin{equation*}
\frac{4\left|\vec{k}_{c m}\right|}{(2 \pi)^{6} E_{c m}} \sigma_{T}=\frac{1}{(2 \pi)^{6} k_{0}^{2}} \times \frac{\lambda^{2}}{8 \pi k_{0}} \sqrt{\left(k_{0} / 2\right)^{2}-m^{2}} \tag{2.4.21}
\end{equation*}
$$

and by comparing with eq. (2.4.16), one recognizes $2 \operatorname{Im} \mathcal{D}\left(k_{0}^{2}\right) /\left[(2 \pi)^{6} k_{0}^{2}\right]=2 \operatorname{Im} A_{2,2}$ on the r.h.s., in agreement with the optical theorem (2.4.9). Alternatively, one can verify the validity of the optical theorem written in the form of eq. (2.4.6) by comparing the expression of its l.h.s. in eq. (2.4.16) with the one obtained by specializing its r.h.s. to the case in which $G_{R}^{(4)}=i \lambda$ (and therefore $A_{2, i}=\lambda /\left[(2 \pi)^{3} k_{0} \sqrt{(2 \pi)^{3} \omega_{\vec{p}_{1}}} \sqrt{(2 \pi)^{3} \omega_{\overrightarrow{p_{1}}}}\right]$ from eq. (2.3.12)) and by taking into account that the two-particle invariant phase space is given by $\int \frac{d^{3} \overrightarrow{p_{1}}}{(2 \pi)^{3} 2 \omega_{\overrightarrow{p_{1}}}} \frac{d^{3} \overrightarrow{p_{2}}}{(2 \pi)^{3} 2 \omega_{\overrightarrow{p_{2}}}}(2 \pi)^{4} \delta^{4}\left(p_{n}-k\right)=\left|\vec{k}_{c m}\right| /\left(4 \pi k_{0}\right)$ (see, e.g., eq. (A.58) of ref. [1]).

### 2.5 Unstable Particles

The formalism above is also useful in understanding how unstable particles should be considered in QFT. These particles cannot appear as asymptotic states in the Hilbert space and hence their treatment is a bit trickier. ${ }^{6}$ We can distinguish between two kinds of unstable particles. Those that turn into stable asymptotic states when interactions are switched off: this is the typical instance of weakly coupled unstable states, that turn

[^4]into free stable states in the limit, and correspond to all unstable elementary particles we know. It also applies to bound states, when their instability is not governed by the same interactions that are at the basis of their formation. For instance the pion mesons in QCD are stable in the limit in which we switch off the electroweak interactions. A more complicated situation occurs with bound states that decay due to the same interactions responsible for their formation. In this case there is not a simple limit where we can consider them as stable. This is for instance the case of the positronium, a bound state composed of an electron and a positron in QED. Let us consider the former simpler case of unstable states admitting a limit in which they are stable. In this case we expect on general grounds that the particle develops a propagator with a Breit-Wigner form. For $p^{2} \sim m^{2}$ we have
\[

$$
\begin{equation*}
G_{R}^{(2)}\left(p^{2}\right) \sim \frac{i}{p^{2}-m^{2}+\Sigma\left(p^{2}\right)} \tag{2.5.1}
\end{equation*}
$$

\]

where $m$ is the mass of the stable particle when interactions are switched off. In presence of interactions $\Sigma$ develops an imaginary part. In the particle rest frame eq.(2.5.1) implies an "effective mass"

$$
\begin{equation*}
M \simeq m-i \operatorname{Im} \Sigma\left(m^{2}\right) /(2 m), \tag{2.5.2}
\end{equation*}
$$

where we have neglected $\operatorname{Re} \Sigma$ and assumed $\operatorname{Im} \Sigma\left(m^{2}\right) \ll m$. The wave function $\Psi$ of the particle evolves as $\Psi(t)=e^{-i M t} \Psi(0)$, such that the probability $|\Psi(t)|^{2} \propto e^{-\Gamma t}$ decreases exponentially in time with a decay rate

$$
\begin{equation*}
\Gamma=\frac{\operatorname{Im} \Sigma\left(p^{2}=m^{2}\right)}{m} . \tag{2.5.3}
\end{equation*}
$$

In the rest frame of the decaying particle the total decay width equals

$$
\begin{equation*}
\Gamma=(2 \pi)^{3} \sum_{n}(2 \pi)^{4} \delta^{4}\left(p_{n}-p\right)\left|A_{n, 1}\right|^{2}, \tag{2.5.4}
\end{equation*}
$$

where $A_{1, n}$ are the amplitudes for the decay in $n$ particles: $1 \rightarrow n$. Despite it does not make sense to talk of $S$-matrix elements for a single-particle state, the link between $\operatorname{Im} \Sigma$ and $\Gamma$ can be formally seen as a consequence of eq.(2.4.6), with $|i\rangle$ is a single particle state.

When the interactions responsible for the decay are switched off, $G_{R}^{(2)}$ has a simple isolated pole, as expected from the representation (2.1.25), and the operator $\phi(x)$ associated to this particle has vanishing matrix elements with the decaying product particles. When interactions are switched on, the operator $\phi(x)$ has non-vanishing matrix elements with the decaying particles. As a consequence the threshold mass in eq.(2.1.25), which is given by the lowest energy to produce multi-particle states, decreases and the previously isolated pole falls within the branch-cut singularity associated to the production of the
decaying particles. Moreover, the simple pole moves away from the real axis and for small $\Sigma$ lies close to the branch-cut singularity. We see from eqs.(2.5.1) and (2.5.3) that the pole sits approximately at $p^{2} \sim m^{2}-i m \Gamma$ and moves in the lower complex half-plane. At first sight, this result is in contradiction with the analyticity properties derived from the Källén-Lehmann spectral representation (2.1.25), according to which $G_{R}^{(2)}$ can have no singularity away from the real axis. Indeed, this complex pole does not appear as a pole of $G_{R}^{(2)}$ but shows up in its analytic continuation below the branch-cut singularity. In other words, the pole is not on the physical Riemann sheet, but in the second (or higher) Riemann sheet. This is understood by recalling that the physical momenta sit above the cut. Since the pole lies instead below the cut, in moving around the cut $\operatorname{Im} \Sigma$ picks up a phase. With such phase the propagator has no longer a pole. In order to not get any phase we have to cross through the cut, which unavoidably leads us to another Riemann sheet. A concrete example of this phenomenon is described in section 2.7.

### 2.6 Causality and Analyticity

In subsection 2.1.2 we have briefly discussed the behaviour of the exact propagator $G_{R}^{(2)}\left(q^{2}\right)$ for complex values of $q^{2}$ and have seen that it can be analytically continued for complex values of $q^{2}$. A similar analytic continuation was assumed in section 2.4 where we briefly discussed the elastic 2-particle scattering. More in general, one could ask whether and in which sense S-matrix elements or the Green functions $G^{(N)}\left(p_{i}\right)$ can be analytically continued for unphysical values of the momenta. We will see that there is a fundamental connection between causality and analyticity, which points towards a positive answer to this question.

The key points are best understood from classical considerations. Let $A_{\text {inc }}(z, t)$ be a wave packet traveling along the $z$ direction:

$$
\begin{equation*}
A_{\mathrm{inc}}(z, t)=\int_{-\infty}^{+\infty} d \omega a_{\mathrm{inc}}(\omega) e^{i \omega(z-t)} \tag{2.6.1}
\end{equation*}
$$

and scattered by a particle at rest at $z=0$. The asymptotic forward scattered wave packet will be of the form

$$
\begin{equation*}
A_{\mathrm{scatt}}(z, t)_{|z| \rightarrow \infty} \approx \frac{1}{z} \int_{-\infty}^{+\infty} d \omega a_{\mathrm{scatt}}(\omega) e^{i \omega(z-t)} \tag{2.6.2}
\end{equation*}
$$

where $a_{\text {scatt }}(\omega)=a_{\text {inc }}(\omega) f(\omega)$, with $f(\omega)$ the forward scattering amplitude. Suppose that this packet represents a signal which vanishes for $t<z$, so in particular $A_{\text {inc }}(0, t)=0$ for $t<0$. By taking the inverse Fourier transform of eq.(2.6.1), we get

$$
\begin{equation*}
a_{\mathrm{inc}}(\omega)=\frac{1}{2 \pi} \int_{-\infty}^{+\infty} d t A_{\mathrm{inc}}(0, t) e^{i \omega t} \tag{2.6.3}
\end{equation*}
$$



Figure 2.3: Thanks to the analyticity of $f\left(\omega^{\prime}\right)$, the contour of integration $\mathcal{C}_{z}$ can be blown up in a large semi-circle approaching the real line from above and going at infinity $\left(\mathcal{C}_{\infty}\right)$ in the upper half-plane.

The function $a_{\text {inc }}(\omega)$ can be analytically continued for complex $\omega$ in the upper half-plane. Indeed, for $t>0$, the phase factor is exponentially suppressed, while for $t<0$, where we would have an exponentially increasing factor, the wave packet vanishes. We can then consider eq.(2.6.1) in the complex $\omega$-plane. If we assume that $a_{\text {inc }}(\omega)$ does not grow too much at infinity, so as to compensate the exponential suppression given by $e^{i \omega(z-t)}$, we can close the contour at infinity. This is automatically satisfied if we demand that $a_{\text {inc }}(\omega)$ grows at most as a polynomial in $\omega$. Such condition is often denoted as polynomial boundedness and is typically assumed for QFT amplitudes. For $t<0$, we get

$$
\begin{equation*}
\oint_{\mathcal{C}} d \omega a_{\mathrm{inc}}(\omega) e^{i \omega(z-t)}=0 \tag{2.6.4}
\end{equation*}
$$

implying that $a_{\text {inc }}(\omega)$ should be an analytic function in the upper $\omega$ half-plane.
The causality principle requires that the scattered wave should also vanish for $t<z$, since no wave can be ahead of the incident wave packet:

$$
\begin{equation*}
A_{\text {scatt }}(z, t)=0, \quad \text { for } t<z \tag{2.6.5}
\end{equation*}
$$

We can now take the inverse Fourier transform of eq.(2.6.2) and repeat the argument above to conclude that the causality condition (2.6.5) implies the analyticity of $a_{\text {scatt }}(\omega)$

- and hence of $f(\omega)$ - in the upper half-plane. ${ }^{7}$

Let us see some implications coming from such analyticity properties. Given a point $z$ in the upper half-plane we can use Cauchy theorem to write

$$
\begin{equation*}
f(z)=\frac{1}{2 i \pi} \int_{\mathcal{C}_{z}} \frac{d \omega^{\prime} f\left(\omega^{\prime}\right)}{\omega^{\prime}-z} \tag{2.6.6}
\end{equation*}
$$

where $\mathcal{C}_{z}$ is a small circle around the point $z$. Since $f$ is analytic over the whole upper $\omega^{\prime}$ half-plane, we can blow the contour $\mathcal{C}_{z}$ until it approaches the real axis and closes at infinity, like in fig.2.3. When $z$ approaches the real line from above, $z=\omega+i \epsilon$, we have

$$
\begin{align*}
f(\omega) & \equiv \lim _{\epsilon \rightarrow 0} f(\omega+i \epsilon)=\lim _{\epsilon \rightarrow 0} \frac{1}{2 i \pi} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} f\left(\omega^{\prime}\right)}{\omega^{\prime}-\omega-i \epsilon}+\frac{1}{2} f_{\infty}  \tag{2.6.7}\\
& =\frac{1}{2 i \pi} \mathrm{P} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} f\left(\omega^{\prime}\right)}{\omega^{\prime}-\omega}+\frac{1}{2} f(\omega)+\frac{1}{2} f_{\infty}
\end{align*}
$$

In the last relation we have used eq.(2.1.26), P stands for the principal part of the integral and $f_{\infty} / 2$ denotes the contribution to the integral coming from $\mathcal{C}_{\infty}$. If we assume for the moment that $f(\omega) \rightarrow 0$ at infinity, so that $f_{\infty}=0$, eq.(2.6.7) gives

$$
\begin{equation*}
f(\omega)=\frac{1}{i \pi} \mathrm{P} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} f\left(\omega^{\prime}\right)}{\omega^{\prime}-\omega} \tag{2.6.8}
\end{equation*}
$$

The real and imaginary parts of eq.(2.6.8) give

$$
\begin{equation*}
\operatorname{Re} f(\omega)=\frac{1}{\pi} \mathrm{P} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} \operatorname{Im} f\left(\omega^{\prime}\right)}{\omega^{\prime}-\omega}, \quad \operatorname{Im} f(\omega)=-\frac{1}{\pi} \mathrm{P} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} \operatorname{Re} f\left(\omega^{\prime}\right)}{\omega^{\prime}-\omega} \tag{2.6.9}
\end{equation*}
$$

Another commonly used rewriting of these equations is

$$
\begin{equation*}
f(\omega)=\lim _{\epsilon \rightarrow 0} \frac{1}{\pi} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} \operatorname{Im} f\left(\omega^{\prime}\right)}{\omega^{\prime}-\omega-i \epsilon} \tag{2.6.10}
\end{equation*}
$$

whose real part coincides with the first relation in eq.(2.6.9). We see that, by knowing the imaginary part of the forward scattering amplitude, we can reconstruct the whole amplitude. The original derivation of eq.(2.6.10) dates back to Kramers and Kronig that applied it in classical optics with $f(\omega)$ being the index of refraction. A frequency dependence of the index of refraction leads to the dispersion of light (like in a rainbow) and hence eq.(2.6.10) (and their quantum generalizations) is called a dispersion relation. An imaginary component of the refractive index leads to dissipation, so the dispersion relations allow us to compute the refractive index of a material (i.e. Re from its dissipation

[^5]properties $(\operatorname{Im} f)$. The latter is also related, through a classical optical theorem, to the total absorption cross section.

If $f(\omega)$ does not vanish at infinity, as we assumed, but rather

$$
\begin{equation*}
f(\omega)_{|\omega| \rightarrow \infty} \propto|\omega|^{N-1} \tag{2.6.11}
\end{equation*}
$$

for some $N \geq 0$, we can divide $f(\omega)$ by a given polynomial $P_{n}(\omega)$ of degree $n=[N]$ and apply Cauchy theorem to $f / P_{n}$. By construction, the contribution at infinity vanishes, but we have to be careful because the zeros of $P_{n}$ give rise to poles whose contribution should be considered. Choosing for definiteness $P_{n}=\omega^{n}$ and repeating the same steps as before, we get

$$
\begin{equation*}
f(\omega)=\omega^{n} \lim _{\epsilon \rightarrow 0} \frac{1}{\pi} \int_{-\infty}^{+\infty} \frac{d \omega^{\prime} \operatorname{Im} f\left(\omega^{\prime}\right)}{\left(\omega^{\prime}-\omega-i \epsilon\right)\left(\omega^{\prime}\right)^{n}}+Q_{n}(\omega) \tag{2.6.12}
\end{equation*}
$$

In eq.(2.6.12), the residues due to the poles of $P_{n}$ are encoded in the factor $Q_{n}$, which is a polynomial of degree $n$ in $\omega$ with coefficients determined in terms of the values of $f(\omega)$ and its first $(n-1)$-derivatives at $\omega=0$ (more generally, at the specific values of $\omega$ where $P_{n}=0$ ). The function $f$ is thus no longer determined by $\operatorname{Im} f$ only, but requires also the knowledge of such "boundary" data. This is the price we have to pay in order to have a sufficiently well-behaved function at infinity. Eq.(2.6.12) is called a dispersion relation with $n$ subtractions, or $n$-subtracted dispersion relation.

The classical causality condition (2.6.5) in QFT is replaced by the microcausality condition [6]

$$
\begin{equation*}
\left[\mathcal{O}_{1}(x), \mathcal{O}_{2}(y)\right]=0, \quad \text { for }(x-y)^{2}<0 \tag{2.6.13}
\end{equation*}
$$

where $\mathcal{O}_{1}$ and $\mathcal{O}_{2}$ are arbitrary bosonic local operators in the QFT. Using eq.(2.6.13) and the LSZ reduction formulae, the analyticity properties of certain amplitudes can be proved axiomatically, i.e. at the non-perturbative level and using first principles only, such as unitarity and Lorentz invariance (in addition to microcausality itself, of course). Unfortunately, an axiomatic proof for general Green functions in QFT is still not available. In these cases one has to rely on perturbation theory, where analytic properties can be directly checked looking at the non-analyticities of Feynman diagrams. All these studies seem compatible with the conjecture that Green functions are the real-boundary values of analytic functions. For concreteness, let us consider the elastic scattering of two identical particles of mass $m$, that we assume to be the lightest particles present in the theory. The scalar amplitude $\mathcal{M}$ is a function of the four particle momenta $p_{i}(i=1,2,3,4)$. Particles 1 and 2 are in-going, particles 3 and 4 are out-going. Momentum conservation and the on-shell conditions fix $p_{1}^{\mu}+p_{2}^{\mu}=p_{3}^{\mu}+p_{4}^{\mu}, p_{i}^{2}=m^{2}$. Lorentz invariance requires that the
scalar quantity $\mathcal{M}$ can only depend on Lorentz invariant combinations of the momenta. A useful parametrization is given by the Mandelstam variables

$$
\begin{equation*}
s=\left(p_{1}+p_{2}\right)^{2}, \quad t=\left(p_{1}-p_{3}\right)^{2}, \quad u=\left(p_{1}-p_{4}\right)^{2} . \tag{2.6.14}
\end{equation*}
$$

It is immediate to verify that $s+t+u=4 m^{2}$, so that the amplitude is effectively a function of two independent variables, that we can take to be $s$ and $t: \mathcal{M}=\mathcal{M}(s, t)$. Going to the center of mass frame of the two particles, it is straightforward to see that the kinematically allowed range of the Mandelstam variables is

$$
\begin{equation*}
s \geq 4 m^{2}, \quad t \leq 0 \quad u \leq 0, \quad(1+2 \rightarrow 3+4) \tag{2.6.15}
\end{equation*}
$$

Studying the analyticity properties of a function of two variables is a difficult task, so let us take $t$ fixed and consider $\mathcal{M}$ as a function of $s$. Some analytic properties of $\mathcal{M}$ can be deduced from perturbation theory. We can have simple poles at $s=\mu^{2}$, where $\mu$ is the mass of the possible particle exchanged in the 1-2 interaction, and then a branch-cut singularity starting from the 2 -particle threshold production, i.e. $s \geq 4 m^{2}$. Using crossing symmetry, the same amplitude $\mathcal{M}$ also describes the scattering process $1+\overline{3} \rightarrow \overline{2}+4$ and $1+\overline{4} \rightarrow \overline{2}+3$. In terms of the variables (2.6.14), the allowed range of $s, t$ and $u$ in the different channels is

$$
\begin{array}{lll}
t \geq 4 m^{2}, & s \leq 0 & u \leq 0, \\
u \geq 4 m^{2}, & s \leq 0 & t \leq 0, \tag{2.6.16}
\end{array} \quad(1+\overline{3} \rightarrow \overline{2}+4),
$$

This implies that, at fixed $t$, if there is a pole in $\mathcal{M}(s)$ at $s=\mu^{2}$ in the $1+2 \rightarrow 3+4$ scattering process ( $s$-channel), a pole should also occur at $s=4 m^{2}-\mu^{2}-t$, which corresponds to the same pole $\mu^{2}$ occurring at $4 m^{2}-s-t=u=\mu^{2}$ in the $1+\overline{4} \rightarrow \overline{2}+3$ process ( $u$-channel). ${ }^{8}$ Similarly, the branch-cut at $u \geq 4 m^{2}$ in the $1+\overline{4} \rightarrow \overline{2}+3$ process implies a branch-cut at $s=-t$ in $\mathcal{M}(s)$. The analyticity assumption based on causality as described above assume that all the singularities of $\mathcal{M}(s)$ are associated to physical processes and, as such, they all lie on the real $s$-axis. In particular, we might have other poles at $s=\mu_{1}^{2}, \mu_{2}^{2}, \ldots$ associated to the exchange of other particles (bound states in general), or branch-cut singularities due to the opening of other multi-particle processes. For example, another branch-cut is expected to occur for $s \geq 9 m^{2}$ in association to the production of three-particles (if allowed by the symmetries of the system, of course). These other branch-cuts will occur on top of the first one that starts at $s \geq 4 m^{2}$ (and $s \leq-t$ ). The expected analyticity domain of $\mathcal{M}(s)$, at fixed $t$, is depicted in fig.2.4. The physical

[^6]

Figure 2.4: The expected singularity structure of the $2 \rightarrow 2$ scattering amplitude $\mathcal{M}(s)$ at fixed $t$ in the complex $s$-plane. The black points indicate the simple poles associated to the particles (bound states) exchanged in the $s$-channel $\left(\mu^{2}\right)$ and in the $u$-channel $\left(4 m^{2}-\mu^{2}-t\right)$. The crosses indicate the branch-cut singularities, starting at $s=4 m^{2}$ in the $s$-channel and at $s=-t$ in the $u$-channel. Poles associated to the possible presence of other bound states have not been indicated.
region in the $s$ channel for $s \geq 4 m^{2}$ lies above the cut. This can be easily seen from the $i \epsilon$ prescription that amounts to a shift $m^{2} \rightarrow m^{2}-i \epsilon$. Since $s+t+u=4 m^{2}$, the physical region in the $u$-channel sits below the cut $s<-t$ in fig.2.4.

As we said, the above picture can be axiomatically proved only in certain cases, but being associated to a fundamental principle such as causality, it is often taken for granted. Once we know the analyticity domain of $\mathcal{M}(s, t)$, dispersion relations like eq.(2.6.10) can be written down. Among other things, these allow us to relate $\mathcal{M}$ to the total cross section, that through the optical theorem can be expressed in terms of $\operatorname{Im} \mathcal{M}$.

The analytical properties of Green functions received great attention starting from the late 50 's until about the early 70 's, in an attempt to describe the strong interactions from first principles, without relying on a weakly coupled Lagrangian description (S-matrix bootstrap). This activity quickly faded away after the discovery of Quantum CromoDynamics (QCD) as the theory of strong interactions, but left us with several important results. In the impossibility to review here all such results, let us just mention a general bound on the cross-section (Froissart-Martin bound), an understanding of certain high energy scattering limits in strongly coupled gauge theories in terms of so called Regge trajectories, and the birth of what we now call string theory. More recently, among other developments, arguments based on the analyticity of amplitudes have allowed us to put constraint on effective field theories [7] and played an essential role in proving an important property satisfied by the renormalization group flow of QFTs in four space-time
dimensions (the so called a-theorem) [8].
A proper understanding of the analytic structure of amplitudes in QFT would also be important to rigorously establish important properties of amplitudes, such as crossing symmetry, that are at the moment assumed. ${ }^{9}$ It is not excluded that a fresh look at some of the old problems left open in the S-matrix bootstrap (together with the vastly increased numerical power we have today with respect to half-century ago) would lead to a revival and further progress in this direction.

### 2.7 Bound States and Resonances**

We have seen in the previous sections that simple poles of the two-point function or of the scattering amplitude ( $S$-matrix) are a signal of one-particle states. These can arise also in correspondence of bound states. We have also seen at the end of section 2.4 that unstable states, resonances, correspond to simple poles with a non-vanishing imaginary component, i.e. complex poles. In this section we will show that

- bound states correspond to simple poles of the scattering amplitude
- resonances give rise to poles hidden below the branch-cut.

In order to keep the analysis as simple as possible, we will just work out a simple example in quantum mechanics (relativistic bound states are a notoriously difficult subject).

The system we consider is a one-dimensional quantum system in a potential well, see fig.2.5. The Schrödinger equation is

$$
\begin{equation*}
\left.\left(\frac{-1}{2 m} \frac{d^{2}}{d x^{2}}+V(x)\right) \psi_{n}(x)\right)=E_{n} \psi_{n}(x) \tag{2.7.1}
\end{equation*}
$$

where $\psi_{n}$ and $E_{n}$ are the eigenfunctions and eigenvalues of the system, respectively. The system is invariant under the $\mathbf{Z}_{2}$ parity transformation $x \rightarrow-x$, so we can separately consider even and odd eigenstates. Let us first discuss the bound states, i.e. the states with energy $-V_{0}<E_{n}<0$, where $V_{0}>0$ is the depth of the well. For the different regions in fig. 2.5 we have, for the parity even states,

$$
\begin{align*}
\psi_{\mathrm{I}}(x) & =C e^{k x} \\
\psi_{\mathrm{II}}(x) & =A \cos (p x)  \tag{2.7.2}\\
\psi_{\mathrm{III}}(x) & =C e^{-k x}
\end{align*}
$$

[^7]

Figure 2.5: One-dimensional potential well.
where

$$
\begin{equation*}
k=\sqrt{-2 m E}, \quad p=\sqrt{2 m\left(E+V_{0}\right)} \tag{2.7.3}
\end{equation*}
$$

and $A$ and $C$ are undetermined constants so far. Demanding the continuity of the wave function and its first derivative at $x=a / 2$ gives the equation that the even eigenfunctions should satisfy:

$$
\begin{equation*}
k=p \tan \frac{p a}{2}, \quad(\text { even }) \tag{2.7.4}
\end{equation*}
$$

Analogously, the equation for the odd eigenfunctions is

$$
\begin{equation*}
k=-p \cot \frac{p a}{2}, \quad(\text { odd }) . \tag{2.7.5}
\end{equation*}
$$

For illustration, we plot in fig.2.6 the solutions of eqs.(2.7.4) and (2.7.5) for some values of $m, a$, and $V_{0}$. At fixed $m$ and $a$, the number of bound states depends on $V_{0}$ and increases with $V_{0}$, as expected. The deepest bound state (i.e. with the largest negative energy) is always parity even. For small enough $V_{0}$ there are no odd bound states and only one even. Its energy can be computed analytically by expanding eq.(2.7.4). For small $V_{0}$, the argument of the tangent is small, $\tan x \approx x$. Taking the square of that equation, we see that $E \sim \mathcal{O}\left(V_{0}^{2}\right)$ and hence at leading order we get the bound state energy

$$
\begin{equation*}
E \approx-\frac{m a^{2} V_{0}^{2}}{2} \tag{2.7.6}
\end{equation*}
$$

Let us now consider the states with $E>0$. The spectrum is a continuum of waves and we look for solutions with an incident wave coming from $x \rightarrow-\infty$. In the different


Figure 2.6: Solutions of the bound states equations (2.7.4) and (2.7.5) for even and odd parity, respectively. We plot in red the function $k$ and in blue $p \tan (p a / 2)$ (even) and $-p \cot (p a / 2)$ (odd). We have taken $m=a=1$ and $V_{0}=100$.
regions we have

$$
\begin{align*}
\psi_{\mathrm{I}}(x) & =e^{i q x}+R e^{-i q x} \\
\psi_{\mathrm{II}}(x) & =B_{1} e^{i p x}+B_{2} e^{-i p x}  \tag{2.7.7}\\
\psi_{\mathrm{III}}(x) & =T e^{i q(x-a)}
\end{align*}
$$

where $q=\sqrt{2 m E}$. The coefficients $B_{1,2}, R$ and $T$ are to be determined, with the latter two being identified with the reflection and transmission coefficients of the wave, respectively. Demanding the continuity of the wave function and its first derivative at $x= \pm a / 2$ allows us, after some algebra, to determine these coefficients. In particular, we are interested here to the transmission coefficient that reads

$$
\begin{equation*}
T(E)=\left[\cos (p a)-\frac{i}{2}\left(\frac{q}{p}+\frac{p}{q}\right) \sin (p a)\right]^{-1} \tag{2.7.8}
\end{equation*}
$$

which is the one-dimensional analogue of a scattering amplitude. Having the luxury of an explicit expression, the analytic properties of $T(E)$ in the complex $E$-plane are easily determined. The transmission coefficient is analytic over the whole complex plane except a branch-cut at $E=0$ given by the variable $q^{10}$ and possible poles whenever

$$
\begin{equation*}
\cos p a=\frac{i}{2}\left(\frac{q}{p}+\frac{p}{q}\right) \sin p a \tag{2.7.9}
\end{equation*}
$$

When properly normalized, poles in $T(E)$ corresponds to a transmitted wave without

[^8]

Figure 2.7: Analytic structure of the transmission coefficient in the complex $E$-plane. The black dots for $E<0$ represents the stable bound states of the system, while the red points are the resonances, that appear as poles in the second Riemann sheet when crossing the branch-cut.
the corresponding incident wave, that is bound states. Dividing by $\cos (p a)$ and using the trigonometric relation $\tan 2 x=2(\cot x-\tan x)^{-1}$, we can recast eq.(2.7.9) in the form

$$
\begin{equation*}
\cot \frac{p a}{2}-\tan \frac{p a}{2}=i\left(\frac{q}{p}+\frac{p}{q}\right) \tag{2.7.10}
\end{equation*}
$$

which has two possible set of solutions:

$$
\begin{equation*}
\cot \frac{p a}{2}=i \frac{p}{q}, \quad \cot \frac{p a}{2}=i \frac{q}{p} . \tag{2.7.11}
\end{equation*}
$$

Let us now write

$$
\begin{equation*}
E=|E| e^{i \phi} \tag{2.7.12}
\end{equation*}
$$

and take the branch of the square root where

$$
\begin{equation*}
\sqrt{E}=\sqrt{|E|} e^{i \phi / 2} \tag{2.7.13}
\end{equation*}
$$

In this branch, negative real values of $E$ correspond to $\phi=\pi$, and thus

$$
\begin{equation*}
q=\sqrt{2 m E} \rightarrow i \sqrt{2 m|E|}=i k \tag{2.7.14}
\end{equation*}
$$

with $k$ as in eq.(2.7.3). It is now immediate to see that eqs.(2.7.11) turn into the bound state equations (2.7.4) and (2.7.5), confirming the expectations that simple poles of the scattering amplitude are associated to stable bound states in the system.

Let us finally discuss unstable resonances. The analogue of the $S$-matrix is given by

$$
\begin{equation*}
S(E) \equiv|T|^{2}=\left(1+\frac{\sin ^{2} p a}{4\left(\frac{E}{V_{0}}\right)\left(1+\frac{E}{V_{0}}\right)}\right)^{-1} \tag{2.7.15}
\end{equation*}
$$

Resonances occur when $S$ is maximal, in this case at

$$
\begin{equation*}
\sin p a=0, \quad \text { with } E>0 \tag{2.7.16}
\end{equation*}
$$

corresponding to

$$
\begin{equation*}
a \sqrt{2 m\left(E+V_{0}\right)}=n \pi \rightarrow E_{n}=\frac{\pi^{2} n^{2}}{2 a^{2} m}-V_{0}>0 \tag{2.7.17}
\end{equation*}
$$

for all positive integer values of $n$ for which the last relation in eq(2.7.17) holds. For such values of $E_{n}$ the reflection coefficient vanishes and the the wave is totally transmitted. The transmission coefficient can be written as

$$
\begin{equation*}
T(E)=\frac{1}{\cos p a}\left(1-\frac{i}{2} f(E)\right)^{-1} \tag{2.7.18}
\end{equation*}
$$

where $f(E)=(q / p+p / q) \tan p a$. Expanding around $E_{n}$, we have

$$
\begin{equation*}
f(E)=0+\left.\frac{d f}{d E}\right|_{E_{n}}\left(E-E_{n}\right)+\mathcal{O}\left(E-E_{n}\right)^{2}, \quad \cos p a=1+\mathcal{O}\left(E-E_{n}\right)^{2} \tag{2.7.19}
\end{equation*}
$$

where

$$
\begin{equation*}
\left.\frac{d f}{d E}\right|_{E_{n}}=\left.\left(\frac{q}{p}+\frac{p}{q}\right) \frac{d p}{d E}\right|_{E_{n}} \equiv \frac{4}{\Gamma_{n}} \tag{2.7.20}
\end{equation*}
$$

is real and positive. Plugging in $T(E)$, we finally have, for $E \approx E_{n}$,

$$
\begin{equation*}
T(E) \approx \frac{1}{1-\frac{i}{2} \frac{4}{\Gamma_{n}}\left(E-E_{n}\right)}=\frac{i \Gamma / 2}{\left(E-E_{n}\right)+i \Gamma_{n} / 2} \tag{2.7.21}
\end{equation*}
$$

which reproduces the expected Breit-Wigner form of a resonance, with $\Gamma_{n}$ its decay width. For $\Gamma_{n} \ll E_{n}$, eq.(2.7.21) would indicate that $T(E)$ has simple poles at

$$
\begin{equation*}
E=E_{n}-\frac{i}{2} \Gamma_{n} \tag{2.7.22}
\end{equation*}
$$

i.e. in the complex plane just below the branch-cut, see fig.2.7. However, a closer inspection shows that these poles are not associated to the original function $T(E)$, but to its analytic continuation below the cut, namely they sit at the second Riemann sheet of the function. Indeed, the pole equation for $T$ is given by eq.(2.7.9). Expanding around $E_{n}$, where $\sin p_{n} a=0$, we get

$$
\begin{equation*}
1+\mathcal{O}\left(E-E_{n}\right)^{2}=\left.\frac{i}{2}\left(\frac{q}{p}+\frac{p}{q}\right) \frac{d p}{d E}\right|_{E_{n}}\left(E-E_{n}\right)+\mathcal{O}\left(E-E_{n}\right)^{2} \tag{2.7.23}
\end{equation*}
$$

With the choice (2.7.13), a pole in the first Riemann sheet just below the positive real axis is reached by taking $\phi=2 \pi$, which would give in eq.(2.7.23) $q \rightarrow-q$. In this way, using eqs.(2.7.20) and (2.7.22), the r.h.s. of eq.(2.7.23) would give -1 instead of +1 . The equation is instead satisfied if we make a clock-wise rotation with $\phi$ small and negative, crossing in this way the branch-cut at $E>0$. In conclusion, the complex poles associated to resonances occur in the second (more in general higher) Riemann sheet of scattering amplitude functions.

## Chapter 3

## Renormalization Theory

Loop amplitudes in quantum field theories are generally divergent. There are at least three kinds of divergences in QFT: infra-red, collinear and ultra-violet divergences. Here we only deal with the last kind, the ultra-violet (UV) divergences. They occur when the momenta of the virtual particles running in the loops go to infinity (hence the name UV). In order to make sense of these otherwise ill-defined amplitudes, we have to "renormalize" the theory. Schematically, this process requires two steps. First, we change the theory so that the amplitudes become finite (regularization) and then we redefine the parameters of the Lagrangian in order to "hide" the divergences in unphysical quantities. All this process is called renormalization of the theory. Before discussing the details of this procedure, it is useful to first classify which amplitudes are potentially divergent. ${ }^{1}$ In this chapter we will consider a generic QFT, with an arbitrary number of scalar, fermion and gauge fields. The latter are taken in the unbroken phase, when they are massless. ${ }^{2}$

### 3.1 Superficial Degree of Divergence

Consider a generic graph $\mathcal{G}$ contributing to a given connected one-particle irreducible (1PI) amplitude. Our aim will be to compute the superficial degree of divergence of $\delta(\mathcal{G}) .{ }^{3} \delta$ is defined as the degree of divergence of the graph when all virtual momenta go to infinity at the same time. Denoting by $q_{i}$ all the virtual momenta running and by $\mathcal{G}\left(q_{i}\right)$ the integrand

[^9]of the $L$-loop graph, we can define $\delta$ as
\[

$$
\begin{equation*}
\lim _{\lambda \rightarrow \infty} \lambda^{4 L} \mathcal{G}\left(\lambda q_{i}\right) \propto \lambda^{\delta(\mathcal{G})} \tag{3.1.1}
\end{equation*}
$$

\]

Notice that the rescaling $q_{i} \rightarrow \lambda q_{i}$ has to be performed also to the loop integrals, and this explains the $\lambda^{4 L}$ factor in eq. (3.1.1). For instance, a one-loop graph of the form

$$
\begin{equation*}
\int \frac{d^{4} q}{(2 \pi)^{4}} \frac{i}{(p+q)^{2}-m^{2}} \frac{i}{q^{2}-m^{2}} \tag{3.1.2}
\end{equation*}
$$

gives

$$
\begin{equation*}
\lim _{\lambda \rightarrow \infty} \lambda^{4} \frac{i}{(p+\lambda q)^{2}-m^{2}} \frac{i}{\lambda^{2} q^{2}-m^{2}} \propto \lambda^{0} \tag{3.1.3}
\end{equation*}
$$

and hence $\delta=0$. A graph is superficially divergent if $\delta \geq 0$ and superficially convergent if $\delta<0$. In particular, $\delta=0$ corresponds to a logarithmic divergence, while $\delta>0$ gives rise to a power-like divergence. The superficial degree of divergence of a graph depends clearly on the structure of the graph, number of internal propagators, vertices, etc. In order to be as general as possible, let us introduce some notation. We label with $i=1, \ldots, n_{V}$ the $n_{V}$ different interactions present in the Lagrangian of the theory, and by $d_{i}$ the number of derivatives present in each interaction vertex $i$. We also denote by $f=e, \gamma, \ldots$ the different fields present in the theory and by $n_{i f}$ the number of fields of type $f$ present in the interaction vertex $i$. Any graph $\mathcal{G}$ is then characterized by the following quantities:

$$
\begin{align*}
I_{f} & \equiv \text { number of propagators of type } f \\
E_{f} & \equiv \text { number of external lines of type } f  \tag{3.1.4}\\
N_{i} & \equiv \text { number of interaction vertices } i
\end{align*}
$$

The momentum space propagator behaves as

$$
\begin{equation*}
\Delta_{f}(p) \propto p^{2 s_{f}-2} \tag{3.1.5}
\end{equation*}
$$

where $s_{f}$ is the spin of the particle type $f$. This formula is correct for scalars $\left(s_{f}=0\right)$ and spin $1 / 2$ fermions $\left(s_{f}=1 / 2\right)$. For gauge fields with $s_{f}=1$, eq. (3.1.5) would give $\Delta_{f}(p) \sim$ constant, which is in fact the correct result for a massive vector, but it is not the correct one, due to gauge invariance, for massless gauge fields, where one has $s_{f}=0$. In the following we will use eq. (3.1.5), with the understanding that $s_{f}=0$ for massless gauge fields. The powers of virtual momenta entering $\mathcal{G}$ is easily determined. We have

$$
\begin{equation*}
\delta(\mathcal{G})=\sum_{f} I_{f}\left(2 s_{f}-2\right)+\sum_{i} d_{i} N_{i}+4\left(\sum_{f} I_{f}-\sum_{i} N_{i}+1\right) . \tag{3.1.6}
\end{equation*}
$$

The first, second and third terms in the right-hand side of eq. (3.1.6) represent the propagator, the vertex and the loop contributions, respectively. The loop contribution has been
written by considering each internal propagator with an independent virtual momentum and imposing momentum conservation at each vertex. The +1 takes into account the overall momentum conservation of the external lines. We can rewrite eq. (3.1.6) as

$$
\begin{equation*}
\delta(\mathcal{G})=\sum_{f} I_{f}\left(2 s_{f}+2\right)+\sum_{i} N_{i}\left(d_{i}-4\right)+4 \tag{3.1.7}
\end{equation*}
$$

Since each internal (propagator) and external line has to end to a vertex, the internal to both ends, the external to one end only, the following identities must hold, for any field type $f$ :

$$
\begin{equation*}
2 I_{f}+E_{f}=\sum_{i} N_{i} n_{i f} \tag{3.1.8}
\end{equation*}
$$

We can use eq. (3.1.8) to solve for $I_{f}$ and plug the result in eq. (3.1.7). In so doing we get

$$
\begin{equation*}
\delta(\mathcal{G})=4-\sum_{f} E_{f}\left(s_{f}+1\right)-\sum_{i} N_{i}\left[4-d_{i}-\sum_{f} n_{i f}\left(s_{f}+1\right)\right] . \tag{3.1.9}
\end{equation*}
$$

We now show that the expression in square brackets in the second term of the right-hand side of eq. (3.1.9) coincides with the dimensionality of the coupling constant, call it $g_{i}$, multiplying the vertex $i$ in the Lagrangian. The latter, indeed, has, by definition, the schematic form

$$
\begin{equation*}
\mathcal{L} \supset g_{i} \partial^{d_{i}} \prod_{f} \phi_{f}^{n_{i f}}, \tag{3.1.10}
\end{equation*}
$$

omitting unnecessary details, such as the Lorentz indices and how the derivatives act on the various fields. Since

$$
\begin{equation*}
\left[\phi_{f}\right]=s_{f}+1 \tag{3.1.11}
\end{equation*}
$$

we immediately get

$$
\begin{equation*}
\left[g_{i}\right] \equiv \Delta_{i}=4-d_{i}-\sum_{f} n_{i f}\left(s_{f}+1\right) \tag{3.1.12}
\end{equation*}
$$

Using eq. (3.1.12) we finally arrive to the desired expression for the superficial degree of divergence of a graph:

$$
\begin{equation*}
\delta(\mathcal{G})=4-\sum_{f} E_{f}\left(s_{f}+1\right)-\sum_{i} N_{i} \Delta_{i} . \tag{3.1.13}
\end{equation*}
$$

Remarkably, the key factors determining $\delta$ are the dimensionality of the coupling constants $\Delta_{i}$. Notice that we are here computing the degree of divergence of single graphs. A given connected 1PI Green function $\Gamma_{E_{f}}$ is uniquely determined by $E_{f}$. In perturbation theory, $\Gamma_{E_{f}}=\sum \mathcal{G}$, where the sum runs over all graphs with $E_{f}$ external lines. A Green function $\Gamma\left(E_{f}\right)$ is superficially finite only if all the graphs $\mathcal{G}$ entering this sum have $\delta(\mathcal{G})<0$. We immediately learn from eq. (3.1.13) that increasing the number of external lines tends to decrease $\delta$. We classify all QFT's as follows:

- If $\Delta_{i} \geq 0$ for all $i=1, \ldots, n_{V}$, only a finite number of Green functions $\Gamma_{E_{f}}$, those for which at most $4-\sum_{f} E_{f}\left(s_{f}+1\right) \geq 0$, is divergent. We call such theories renormalizable.
- If at least one $\Delta_{i}<0$, an infinite number of Green functions $\Gamma\left(E_{f}\right)$ is divergent. We call such theories non-renormalizable.

Among the renormalizable theories, a further distinction is possible: if $\Delta_{i}>0$ for all $i=1, \ldots, n_{V}$, only a finite number of individual graphs diverges. We call such theories super-renormalizable. When this distinction occurs, a renormalizable theory is defined as one in which at least one coupling is dimensionless. The terminology "renormalizable" and "non-renormalizable" applies also to individual operators or couplings. We denote a coupling constant as renormalizable if its dimension is positive and non-renormalizable if it is strictly negative. Similarly, an operator is denoted renormalizable if its dimension is less or equal to four and non-renormalizable if it is higher than four.

Let us list the possible renormalizable couplings. In a QFT with scalars only, we can have $\phi^{3}$ and $\phi^{4}$ interactions only. No renormalizable derivative interactions are allowed, compatibly with Lorentz invariance. In a QFT with scalars and fermions, the only possible renormalizable interactions are the Yukawa-like of the form $\phi \bar{\psi} \psi$ or $\phi \bar{\psi} \gamma_{5} \psi$. Again, no Lorentz invariant renormalizable derivative interactions are allowed. When gauge fields are included, we have $\bar{\psi} \gamma^{\mu} \psi A_{\mu}, \bar{\psi} \gamma^{\mu} \gamma_{5} \psi A_{\mu},\left(\phi^{\dagger} \partial_{\mu} \phi-\phi \partial_{\mu} \phi^{\dagger}\right) A^{\mu}$ and $\phi^{2} A_{\mu} A^{\mu}$. Other possibly renormalizable interactions like $\left(A_{\mu} A^{\mu}\right)^{2}$ or $\left(\phi^{\dagger} \partial_{\mu} \phi+\phi \partial_{\mu} \phi^{\dagger}\right) A^{\mu}$ are forbidden by gauge invariance.

Multi-loop graphs can be divergent, despite being superficially convergent. This happens when the divergence arises from a sub-set of all virtual momenta becoming large, keeping fixed the others. As an example, consider the two two-loop contributions to the Compton scattering in QED in figure 3.1. In QED we have $f=e, \gamma$ and no index $i$ is needed, since we have one interaction only, with $\Delta=0$ (the theory is renormalizable). Applying eq. (3.1.13) to these graphs, we get

$$
\begin{equation*}
\delta\left(\mathcal{G}_{a}\right)=\delta\left(\mathcal{G}_{b}\right)=4-2 \times 1-2 \times \frac{3}{2}=-1 \tag{3.1.14}
\end{equation*}
$$

Both graphs are superficially convergent, but while graph (a) is actually convergent, graph (b) is divergent. This is seen by keeping fixed the virtual momentum running over the big external loop while letting the virtual momenta in the small loop go to infinity. In this situation, the small loop behaves as a sub-graph with effectively $E_{e}=2$ and $E_{\gamma}=0$, that has $\delta=1$. Let us see this in some further detail. Neglecting unnecessary factors, the


Figure 3.1: Two superficially convergent graphs contributing at two-loop level to the Compton scattering. Graph (a) is absolutely convergent, graph (b) is divergent, due to the subintegration associated to the sub-graph enclosed in the dashed red rectangle.
amplitude associated to the graphs $(a)$ and (b) in fig. 3.1 is schematically given by

$$
\begin{align*}
& \mathcal{M}_{a} \sim \int d^{4} q_{1} d^{4} q_{2} \frac{1}{p_{1}+q_{1}} \frac{1}{q_{1}^{2}} \frac{1}{p_{1}+\phi_{1}+q_{2}} \frac{1}{q_{2}^{2}} \frac{1}{p_{1}+\not p_{2}+\phi_{1}+q_{2}} \frac{1}{p_{4}+\phi_{1}+q_{2}} \frac{1}{p_{4}+\phi_{1}},  \tag{3.1.15}\\
& \mathcal{M}_{b} \sim \int d^{4} q_{1} d^{4} q_{2} \frac{1}{\not p_{1}+\not q_{1}} \frac{1}{q_{1}^{2}} \frac{1}{\left(\not p_{1}+\not p_{2}+\not q_{1}\right)^{2}} \frac{1}{q_{2}^{2}} \frac{1}{\not p_{1}+\not p_{2}+\not q_{1}+\not q_{2}} \frac{1}{\not p_{4}+\not q_{1}} .
\end{align*}
$$

We have denoted by $q_{1}$ and $q_{2}$ the virtual momenta of the big and small photon loops, respectively, by $p_{1}$ and $p_{4}$ the electron momenta and by $p_{2}$ and $p_{3}$ the photon momenta, and used momentum conservation to write eq. (3.1.15) in terms of $p_{1}, p_{2}$ and $p_{4}$. By rescaling $q_{1,2} \rightarrow \lambda q_{1,2}$ in eq. (3.1.15) one immediately recovers $\delta=-1$ for both graphs. However, if one keeps $q_{1}$ fixed, $\mathcal{M}_{b}$ is still finite when integrating in $q_{2}$, while $\mathcal{M}_{a}$ diverges, with an effective degree of divergence of one, in agreement with the analysis above.

In general, a graph is convergent if $\delta<0$ not only when all the virtual momenta are taken large, but also when any combination of virtual momenta is taken large, with the remaining ones kept fixed. When this is the case, we say that the graph is absolutely convergent.

It is straightforward to generalize to an arbitrary number of space-time dimensions the computation of the superficial degree of divergence of a graph. Equation (3.1.6) is still valid provided we replace the factor of 4 with $D$. In so doing we get

$$
\begin{equation*}
\delta(\mathcal{G})=D-\sum_{f} E_{f}\left(s_{f}-1+\frac{D}{2}\right)-\sum_{i} N_{i}\left(D-d_{i}-\sum_{f} n_{i f}\left(s_{f}-1+\frac{D}{2}\right)\right) \tag{3.1.16}
\end{equation*}
$$

In $D$ space-time dimensions

$$
\begin{equation*}
\left[\phi_{f}\right]=s_{f}-1+\frac{D}{2} \tag{3.1.17}
\end{equation*}
$$

and consequently the interactions (3.1.10) have dimensions

$$
\begin{equation*}
\left[g_{i}\right] \equiv \Delta_{i}=D-d_{i}-\sum_{f} n_{i f}\left(s_{f}-1+\frac{D}{2}\right) \tag{3.1.18}
\end{equation*}
$$

and eq. (3.1.13) becomes

$$
\begin{equation*}
\delta(\mathcal{G})=D-\sum_{f} E_{f}\left(s_{f}-1+\frac{D}{2}\right)-\sum_{i} N_{i} \Delta_{i} \tag{3.1.19}
\end{equation*}
$$

The renormalizability of a theory in $D$ dimensions is still characterized by the dimensions of their couplings $\Delta_{i}$. If all $\Delta_{i} \geq 0$ the theory is renormalizable, otherwise it is nonrenormalizable. We notice that when $D>4$ most four-dimensional renormalizable theories turn into non-renormalizable ones, because the dimension of the couplings $\Delta_{i}$ decreases when $D$ increases. For instance, in $D=5$ dimensions, the gauge couplings constants have dimension $\Delta=-1 / 2$. All gauge theories are non-renormalizable for $D>4$. The only renormalizable interaction up to $D=6$ is the trilinear $\phi^{3}$ vertex. For $D>6$ there are no renormalizable vertices. On the contrary, when $D<4$, non-renormalizable interactions in $D=4$ turn into renormalizable ones. For instance, $\phi^{6}$ in $D=3$. Gauge theories become super-renormalizable for $D<4$.

### 3.2 Cancellation of divergences and Local Counterterms

Any graph $\mathcal{G}$ with $E_{f}$ external lines is a function of the momenta of the external particles. If $\mathcal{G}$ is divergent, one might naively argue that this function is ill-defined and does not make sense. On the contrary, there is a way to disentangle the divergent part from the rest, which remains finite and calculable. In order to understand how divergences can be disentangled, it is enough to consider one-dimensional divergent integrals, toy versions of the more involved integrals appearing in real amplitudes. Take for instance

$$
\begin{equation*}
I(q)=\int_{0}^{\infty} \frac{d k}{q+k} \tag{3.2.1}
\end{equation*}
$$

The function $I(q)$ is divergent, with $\delta=0$. Differentiating with respect to the "external momentum" $q$ gives a finite integral, with $\delta=-1$ : $^{4}$

$$
\begin{equation*}
I^{\prime}(q)=-\int_{0}^{\infty} \frac{d k}{(q+k)^{2}}=-\frac{1}{q} \tag{3.2.2}
\end{equation*}
$$

Integrating back, we get

$$
\begin{equation*}
I(q)=-\log q+c \tag{3.2.3}
\end{equation*}
$$

[^10]Evidently, from eq. (3.2.3), the divergence of $I(q)$ is all encoded in the constant $c$, while the $\log q$ term is finite and calculable. Consider another example, with $\delta=1$ :

$$
\begin{equation*}
\hat{I}(q)=\int_{0}^{\infty} \frac{k d k}{q+k} \tag{3.2.4}
\end{equation*}
$$

We now need to differentiate twice with respect to $q$ to get a finite result. Integrating back twice gives

$$
\begin{equation*}
\hat{I}(q)=q \log q+a q+b \tag{3.2.5}
\end{equation*}
$$

As before all the divergences are encoded in the two coefficients $a$ and $b$, while the nonpolynomial $q \log q$ term is perfectly finite. The generalization of these results is obvious. A graph with $\delta>0$ can be made convergent by taking $\delta+1$ derivatives with respect to the external momenta and all the divergences are encoded in the coefficients of a polynomial of degree $\delta$ in the external momenta. Notice that in general the graph will be only superficially, and not absolutely, convergent. We will ignore for the moment this problem, postponing its solution for later. Terms polynomial in momenta can be written, in configuration space, as derivatives of a local Lagrangian density. In other words, divergences only affect terms that can be written as local interactions ${ }^{5}$ in a Lagrangian. The divergences of a graph with degree of divergence $\delta$ can be reabsorbed by adding to the Lagrangian $\delta+1$ vertices with $n_{i f}=E_{f}$, of the form

$$
\begin{equation*}
\mathcal{L} \supset a_{0} \prod_{f} \phi_{f}^{E_{f}}+a_{1} \partial \prod_{f} \phi_{f}^{E_{f}}+\ldots+a_{\delta} \partial^{\delta} \prod_{f} \phi_{f}^{E_{f}} \tag{3.2.6}
\end{equation*}
$$

If the theory already contains such interactions, the divergences simply amount to a redefinitions of the corresponding couplings, or to a redefinition of the field themselves for 2-point functions. If they were not present, they are induced by quantum effects. In a non-renormalizable theory, where Green functions eventually diverge for any $E_{f}$, cancellation of the divergences imposes to us to add to the Lagrangian all possible terms of the form (3.2.6). Of course, not every possible term is generated, but only those compatible with the possible global symmetries of the Lagrangian. In a renormalizable theory, the number of terms in eq. (3.2.6) is finite. Moreover, the dimension of the couplings $a_{i}$ in eq. (3.2.6) induced by the divergences is

$$
\begin{equation*}
\left[a_{n}\right]=4-\sum_{f} E_{f}\left(s_{f}+1\right)-n, \quad n=0, \ldots, \delta \tag{3.2.7}
\end{equation*}
$$

For a renormalizable theory, we have $\delta \leq 4-\sum_{f} E_{f}\left(s_{f}+1\right)$ and hence

$$
\begin{equation*}
\left[a_{n}\right] \geq \delta-n \geq 0, \quad n=0, \ldots, \delta \tag{3.2.8}
\end{equation*}
$$

[^11]The couplings induced by quantum effects in a renormalizable theory are themselves renormalizable. These theories are self-consistent at the quantum level, provided we write down all the couplings with $\Delta_{i} \geq 0$ allowed by symmetries. In renormalizable theories the finite number of local counter-terms $a_{n}$ is enough to remove all divergences and make the theory finite. Since these terms were already present in the original Lagrangian, we see that all the divergences can be buried into a redefinition of the interactions, masses and fields.

### 3.3 Regularization and Renormalization: QED Case

Dealing with divergent Green functions can lead to subtleties and misleading results. For this reason it is highly desirable to modify the original theory so that divergences no longer appear. This process is called regularization. There are various ways to do that and we will not list all of them. The regularization process should be done in a way such that the modified Lagrangian has all the symmetries present in the original one. ${ }^{6}$

We only mention two relevant types of regularization, that will be used throughout the course: cut-off regularization and dimensional regularization. The first is certainly the simplest and most physical way of regularizing a theory. It amounts of truncating the integration over the virtual momenta, in principle running up to infinity, to a limiting value $\Lambda$. In this way all UV divergences are clearly removed and Green functions become $\Lambda$-dependent. divergences will now appear as power-like or logarithmic terms in $\Lambda$ in an expansion for $\Lambda \rightarrow \infty$. The physical and very reasonable assumption underlying cutoff regularization is that at short distances $L \sim 1 / \Lambda$ our theory is no longer valid, and we parametrize our ignorance by cutting the momenta at that scale. The regularized theory differs from the original one only in the UV, while their IR behaviours are similar, a welcome feature. Cut-off regularization has however a serious drawback: it does not respect gauge invariance.

Dimensional regularization (DR) is a bit more exotic: it amounts in changing the space-time dimensions from 4 to $4-\epsilon$, with $\epsilon \ll 1$. We make sense of the notion of a noninteger number of dimensions by analytic continuation. In the dimensionally regulated Lagrangian UV divergences no longer appear. In an expansion for small $\epsilon$ they are given by poles in $1 / \epsilon^{n}$, with $n \geq 1$. Dimensional regularization respects gauge invariance and is practically one of the simplest regulators to use. Most of the computations in the literature, mainly when dealing with gauge theories, are performed using DR . In addition,

[^12]DR has the great virtue of respecting dimensional analysis and is hence the most suitable regularization in the context of effective field theories. More on this point in section 7.7. Dimensional regularization has a (minor) drawback: it modifies the theory to all scales, not only in the UV, and for this reason it has to be used with some care, when mass scales are present. More on this point in section 5.6. Further technical details on how to use DR are discussed in subsection 3.4. Of course, introducing a regulator is not the whole story to deal with divergences, since divergences strike back when we recover the original theory! However, in the regularized theory we can meaningfully manipulate the Green functions, now finite. As we have seen in section 3.2, divergences only appear in local terms in Green functions and can be removed by redefining the fields, couplings and masses of the original Lagrangian. This process is called renormalization. It is clear that such a redefinition implies that the masses and coupling constants appearing in the initial Largangian (sometimes called bare parameters) cannot be the physical ones. We have seen that the Källen-Lehmann representation for the exact propagators in an interacting theory provides us with a definition of physical mass and properly normalized quantum field. There is not a generic and unique procedure for defining the coupling constants, that depends on the specific theory we are dealing with. As we will see in great detail in the rest of these lectures, coupling constants are commonly defined by demanding that the 1PI Green functions associated to them have a specific value (determined from the experiment) in a given point in momentum space.

Let us consider as an example the renormalization of QED, namely one charged massive fermion (electron) interacting with a photon. First of all, let us use eq. (3.1.13) to classify the possible divergent Green functions. We have only one dimensionless coupling constant and $f=e, \gamma$, so

$$
\begin{equation*}
\delta=4-\frac{3}{2} E_{e}-E_{\gamma} . \tag{3.3.1}
\end{equation*}
$$

The divergent Green functions contain graphs with $\delta \geq 0$. Given eq. (3.3.1), they are $\left(E_{e}=0, E_{\gamma}=0,1,2,3,4\right),\left(E_{e}=1, E_{\gamma}=0,1,2\right),\left(E_{e}=2, E_{\gamma}=0,1\right)$. Lorentz invariance forbids Green functions with an odd number of electron lines, while charge conjugation implies that ( $E_{e}=0, E_{\gamma}=1,3$ ) vanishes. We neglect the quartically divergent vacuum amplitude $\left(E_{0}=E_{\gamma}=0\right)$ since it is irrelevant in a QFT where gravity is not dynamical. In total, we are left with four cases: $\left(E_{e}=0, E_{\gamma}=2,4\right),\left(E_{e}=2, E_{\gamma}=0,1\right)$.
$\mathbf{E}_{\mathbf{e}}=\mathbf{2}, \mathbf{E}_{\gamma}=\mathbf{1}$ We denote this Green function by $\Gamma_{\mu}^{a b}\left(p, p^{\prime}\right)$, where $a, b$ are spinor indices and $p$ and $p^{\prime}$ are the momenta of the electron lines. It corresponds to the basic QED interaction given by minimal coupling. Since $\delta\left(\Gamma_{\mu}\right)=0$, the divergence can only appear in a constant term, independent of $p$ and $p^{\prime}$. Given the Lorentz structure of $\Gamma_{\mu}$, we necessarily
have (omitting spinor indices from now on)

$$
\begin{equation*}
i \Gamma_{\mu}\left(p, p^{\prime}\right)=i e L \gamma^{\mu}+i \Gamma_{\mu}^{\text {finite }}\left(p, p^{\prime}\right) \tag{3.3.2}
\end{equation*}
$$

where $L$ is a divergent constant, possibly reabsorbed in the definition of the electric charge.
$\mathbf{E}_{\mathbf{e}}=\mathbf{2}, \mathbf{E}_{\gamma}=\mathbf{0}$ We denote this Green function by $\Sigma(p)$. It corresponds to the electron propagator. We have $\delta(\Sigma)=1$, so divergences proportional to a constant and to a term linear in momentum can appear. The Lorentz structure implies

$$
\begin{equation*}
i \Sigma(p)=i A+i B \not p+i \Sigma^{\text {finite }}(p) \tag{3.3.3}
\end{equation*}
$$

where $A$ and $B$ are divergent constants. Gauge invariance imposes the identity $B=L$ to all orders in perturbation theory.
$\mathbf{E}_{\mathbf{e}}=\mathbf{0}, \mathbf{E}_{\gamma}=\mathbf{2}$ We denote this Green function by $\Pi_{\mu \nu}(q)$. It corresponds to the photon propagator. We have $\delta(\Pi)=2$, so divergences up to terms quadratic in $q$ might appear. Lorentz invariance forbids terms linear in $q$, so we have

$$
\begin{equation*}
i \Pi_{\mu \nu}(q)=i C_{1} \eta_{\mu \nu}+i C_{2} q^{2} \eta_{\mu \nu}+i C_{3} q_{\mu} q_{\nu}+i \Pi_{\mu \nu}^{f i n i t e}(q) \tag{3.3.4}
\end{equation*}
$$

Dimensional arguments only would imply the appearance of three independent divergent constants. However, gauge invariance implies the following identity: ${ }^{7}$

$$
\begin{equation*}
q^{\mu} \Pi_{\mu \nu}(q)=q^{\nu} \Pi_{\mu \nu}(q)=0 \tag{3.3.5}
\end{equation*}
$$

Using eq. (3.3.5) we have

$$
\begin{align*}
C_{1} & =0, \quad C_{2}=-C_{3} \equiv C \\
\Pi_{\mu \nu}^{\text {finite }}(q) & =\left(\eta_{\mu \nu} q^{2}-q_{\mu} q_{\nu}\right) \Pi^{\text {finite }}\left(q^{2}\right) \tag{3.3.6}
\end{align*}
$$

and hence

$$
\begin{equation*}
i \Pi_{\mu \nu}(q)=\left(\eta_{\mu \nu} q^{2}-q_{\mu} q_{\nu}\right)\left(i C+i \Pi^{\text {finite }}\left(q^{2}\right)\right) \tag{3.3.7}
\end{equation*}
$$

We notice how symmetries (gauge invariance) are crucial to establish relations among divergences and to possibly lower the degree of divergence of a Green function from its power counting value (in our case, from $\delta=2$ to $\delta=0$ ).

[^13]$\mathbf{E}_{\mathbf{e}}=\mathbf{0}, \mathbf{E}_{\gamma}=\mathbf{4}$ We denote this Green function by $M_{\mu \nu \rho \sigma}\left(p_{i}\right)$, with $p_{i}(i=1,2,3,4)$ the momenta of the external photons. It corresponds to the $2 \rightarrow 2$ photon scattering. We have $\delta(M)=0$, implying a possible divergence in the constant factor. Lorentz invariance and Bose symmetry of the external legs fix the Green function to be
\[

$$
\begin{equation*}
M_{\mu \nu \rho \sigma}\left(p_{i}\right)=K\left(\eta_{\mu \nu} \eta_{\rho \sigma}+\eta_{\mu \rho} \eta_{\nu \sigma}+\eta_{\mu \sigma} \eta_{\nu \rho}\right)+M_{\mu \nu \rho \sigma}^{f i n i t e}\left(p_{i}\right) \tag{3.3.8}
\end{equation*}
$$

\]

with $K$ the divergent constant. Once again, gauge invariance imposes other constraints and requires that ${ }^{8}$

$$
\begin{equation*}
p_{1}^{\mu} M_{\mu \nu \rho \sigma}\left(p_{i}\right)=p_{2}^{\nu} M_{\mu \nu \rho \sigma}\left(p_{i}\right)=p_{3}^{\rho} M_{\mu \nu \rho \sigma}\left(p_{i}\right)=p_{4}^{\sigma} M_{\mu \nu \rho \sigma}\left(p_{i}\right)=0 \tag{3.3.9}
\end{equation*}
$$

The above relation implies that $K=0$ and that a relation similar to eq. (3.3.9) applies with $M \rightarrow M^{\text {finite } .} 9$

Summarizing, taking gauge invariance into account, three Green functions diverge in QED, the photon and electron propagators and the interaction vertex, with a total of three independent divergences, encoded in the constants $L, A$ and $C$ in eqs. (3.3.2), (3.3.3), and (3.3.7).

Let us now see how these divergences can be cancelled, to all orders in perturbation theory, by suitable local counter-terms. The QED Lagrangian reads

$$
\begin{equation*}
\mathcal{L}_{Q E D}=-\frac{1}{4} F_{B, \mu \nu} F_{B}^{\mu \nu}+\bar{\psi}_{B}\left(i \not D_{B}\left(A_{B}\right)-m_{B}\right) \psi_{B} \tag{3.3.10}
\end{equation*}
$$

Here $D_{B, \mu}\left(A_{B}\right)=\partial_{\mu}-i e_{B} A_{B, \mu}$ is the covariant derivative and the subscript $B$ reminds us that the fields, mass and coupling constant are the bare ones. The physical fields are defined as

$$
\begin{equation*}
\psi_{B}=\sqrt{Z_{2}} \psi, \quad A_{B, \mu}=\sqrt{Z_{3}} A_{\mu} \tag{3.3.11}
\end{equation*}
$$

We then have

$$
\begin{equation*}
\mathcal{L}_{Q E D}=-\frac{1}{4} Z_{3} F_{\mu \nu} F^{\mu \nu}+Z_{2} \bar{\psi} i \not \partial \psi-m_{B} Z_{2} \bar{\psi} \psi+e_{B} Z_{2} \sqrt{Z_{3}} \bar{\psi} A_{\mu} \gamma^{\mu} \psi \tag{3.3.12}
\end{equation*}
$$

and define renormalized mass and electric charge as

$$
\begin{equation*}
e_{B} Z_{2} \sqrt{Z_{3}}=Z_{1} e, \quad m_{B} Z_{2}=m+\delta m \tag{3.3.13}
\end{equation*}
$$

[^14]

Figure 3.2: Feynman rules for the QED counter-terms defined in eq. (3.3.14).

In terms of the physical fields and parameters, the QED Lagrangian (3.3.10) can be rewritten as

$$
\begin{align*}
\mathcal{L}_{Q E D}= & -\frac{1}{4} F_{\mu \nu} F^{\mu \nu}+\bar{\psi}(i \not D(A)-m) \psi  \tag{3.3.14}\\
& -\frac{1}{4}\left(Z_{3}-1\right) F_{\mu \nu} F^{\mu \nu}+\left(Z_{2}-1\right) \bar{\psi} i \not \partial \psi-\delta m \bar{\psi} \psi+e\left(Z_{1}-1\right) \bar{\psi} A_{\mu} \gamma^{\mu} \psi
\end{align*}
$$

The first line in eq. (3.3.14) is the QED Lagrangian written now in terms of the physical field and parameters, and the second line contains the counter-terms. The latter are unphysical local terms which are there just to cancel the divergences appearing in the loop diagrams. They also include terms quadratic in the fields. However, contrary to the canonical kinetic terms in the first line, they should be considered as new effective vertices. This is clear by noticing that for $e=0$ (free theory), no divergences occur and all counter-terms vanish. For $e \ll 1$, all the counter-terms can be Taylor expanded and are given by positive powers of $e$. As such, they should be considered as interaction terms. In evaluating Feynman diagrams, we should take into account the new contributions given by the counter-terms. The Feynman rules associated to these interactions is straightforwardly derived and reported in fig. 3.2. The four counter-terms $Z_{1}, Z_{2}, Z_{3}$ and $\delta m$ in the Lagrangian (3.3.14) are not all independent. In particular, gauge invariance requires

$$
\begin{equation*}
Z_{1}=Z_{2} \tag{3.3.15}
\end{equation*}
$$

which is equivalent to the condition $B=L$ mentioned before. This will be proved in section 4.4.3. Order by order in perturbation theory, the values of the counter-terms $Z_{1}=Z_{2}$, $Z_{3}$ and $\delta m$ is suitably chosen to cancel the divergences appearing in $L=B, C$ and $A$, respectively. For instance, at one-loop level, we have $L=1+e^{2} L_{(1)}+\ldots, C=e^{2} C_{(1)}+\ldots$,
$A=e^{2} A_{(1)}+\ldots$ Adding the counter-term contributions, the tree-level + one-loop Green functions read

$$
\begin{align*}
\Gamma_{\mu}\left(p, p^{\prime}\right) & =e \gamma^{\mu}\left(1+e^{2} L_{(1)}+Z_{1}-1\right)+\Gamma_{\mu}^{\text {finite }}\left(p, p^{\prime}\right), \\
\Sigma(p) & =\left(1+e^{2} L_{1}+Z_{2}-1\right) p-m+e^{2} A_{(1)}-\delta m+\Sigma^{\text {finite }}(p),  \tag{3.3.16}\\
\Pi_{\mu \nu}(q) & =\left(\eta_{\mu \nu} q^{2}-q_{\mu} q_{\nu}\right)\left(e^{2} C_{(1)}-Z_{3}+1+\Pi^{\text {finite }}\left(q^{2}\right)\right) .
\end{align*}
$$

from which we get the values of the counter-terms up to order $e^{2}$ :

$$
\begin{equation*}
Z_{1}-1=-e^{2} L_{(1)}, \quad Z_{2}=Z_{1}, \quad Z_{3}=1+e^{2} C_{(1)}, \quad \delta m=e^{2} A_{(1)} \tag{3.3.17}
\end{equation*}
$$

It is now time to come back to a problem we encountered in section 3.2, related to the fact that in general taking derivatives with respect to external momentum does not necessarily make the graph absolutely convergent. An example of a graph of this sort is given by the two-loop graph (b) in fig. 3.1. If we take derivatives with respect to the external momenta of $\mathcal{M}_{b}$ in eq. (3.1.15), say $p_{1}$ or $p_{4}$, we see that when the derivative acts on the $q_{2}$-independent propagators we do not improve the convergence of the $q_{2}$-loop integration, that in fact remains divergent. This problem is easily solved by noticing that, together with the graph (b) of fig. 3.1, we should also consider the graph in which the one-loop enclosed in the red rectangle is replaced by the fermion propagator counter-term. The divergence in the small loop is cancelled when summing the two diagrams. This example makes clear that counter-terms have to be considered in evaluating Feynman diagrams like ordinary interaction vertices and should be inserted in graphs consistently to the order in perturbation theory one is considering. In this way divergences are removed even in the more tricky situation of divergent sub-graphs that share a common line, the so called overlapping divergences. An example of this sort in QED is given by the two-loop correction to the photon propagator shown in fig. 3.3. In order to remove all divergences from this graph, one has to consider the one-loop graphs where one vertex is replaced by a counter-term, up to order $e^{2}$. Then a remaining local divergence is finally cancelled by considering the photon propagator counter-term at order $e^{4}$.

The considerations made here apply of course to any local QFT and are not restricted to the QED case, that we have taken as an example. The proof that all divergences are cancelled by the counter-terms is not trivial at all and is known as the BPHZ theorem, from the authors Bogoliubov, Parasiuk, Hepp and Zimmermann that at different stages have proved this statement.

Renormalizability puts stringent constraints on the possible interactions appearing in the Lagrangian. However, upon including all possible interactions, all theories become renormalizable, in the sense that we can get rid of all possible infinities. The key difference


Figure 3.3: Two-loop graph contributing to the photon propagator and its additional graphs including counter-terms canceling all the divergences. The $e^{2}$ and $e^{4}$ factors denote the order in which the counter-terms $Z_{2}$ and $Z_{3}$ enter the various graphs.
between renormalizable and non-renormalizable theories is in their predictivity. There is nothing wrong with non-renormalizable theories. If we have a coupling constant $g$ with scaling dimension $\Delta<0$, we can always write it as

$$
\begin{equation*}
g=\frac{1}{M^{|\Delta|}} . \tag{3.3.18}
\end{equation*}
$$

In scattering amplitudes, by dimensional analysis, the coupling $g$ will always appear as $(E / M)^{|\Delta|}$, where $E$ is the energy scale of the process under consideration. For $E \ll M$, the coupling $g$ is irrelevant (in fact, non-renormalizable couplings are also called irrelevant couplings, see section 5.1). Despite the fact that the theory is formally non-renormalizable and requires an infinite number of counter-terms, in practice at sufficiently low energies most of these interactions are negligible and make the non-renormalizable theory useful and predictive enough. A relevant example of non-renormalizable theory is gravity. Seen as a QFT, gravity is among the most non-renormalizable theories. The mass scale entering the non-renormalizable couplings like in eq. (3.3.18) is $M_{P l} \sim 10^{18} \mathrm{GeV}$. We do not know which is the quantum theory of gravity (although we have some candidates, most notably string theory). Unfortunately (or fortunately, depending on the point of view), in all practical cases, $M_{P l}$ is so large that gravity effects are always negligible in high energy physics. Non-renormalizable theories might also appear from renormalizable ones when some massive fields are integrated out. A relevant case is the four-fermion electroweak interaction, approximation of the Standard Model when one integrates out the $W$ and $Z$ bosons. At sufficiently low energies $E \ll M_{W}, M_{Z}$, the Fermi theory provides an excellent description of the physics.

In the modern approach to QFT we consider all theories except, maybe, the possible ultimate theory of nature (string theory?) as effective field theories, namely theories valid up to some energy scales. More on effective field theories in chapter 7.

### 3.4 Dimensional Regularization

Dimensional regularization can be formally introduced via a set of axioms, as explained, e.g., in ref. [9]. Here we briefly discuss its practical aspects, starting from the simple example of the (Euclidean) integral (tadpole of $\phi^{4}$ theory)

$$
\begin{equation*}
T_{d} \equiv \int \frac{d^{d} p}{(2 \pi)^{d}} \frac{1}{p^{2}+m^{2}}=m^{d-2} \frac{\Gamma(1-d / 2)}{(4 \pi)^{d / 2}}, \tag{3.4.1}
\end{equation*}
$$

(where $m$ is the mass of the scalar) which converges in the IR as long as $m^{2} \neq 0$ and in the UV only for $d<2$. In this case, the r.h.s. is easily calculated by recalling that $1 / A=\int_{0}^{\infty} d t e^{-t A}$, which turns $T_{d}$ into a Gaussian integral in $d$-dimensions that, once calculated, allows to express $T_{d}$ in terms of $\Gamma(x)=\int_{0}^{\infty} d t t^{x-1} e^{-t}$, which converges for $x>0$. Note that, while in the intermediate steps of the calculation $d$ has been treated as an integer number, the r.h.s. of eq. (3.4.1) is well defined (via the analytic continuation of the $\Gamma$ function ${ }^{10}$ ) for any complex $d$ which is not an even integer. This simple observation constitutes, in practice, a regularization of the integral $T_{d}$ because it now assumes a finite value (given by the r.h.s. of eq. (3.4.1)) for (almost) all $d$, analogously to what happens when the regularization is done by other techniques, such as the introduction of a large-momentum cutoff $\Lambda$. However, a reminder of the divergent nature of the original integral is in the fact that the r.h.s. itself diverges if $d$ approaches even integer values $\geq 2$. An important and general property of this regularization method is that the regularized integral (3.4.1) in the massless case $m \rightarrow 0$ vanishes for $d>2$, leading to the relation

$$
\begin{equation*}
\int \frac{d^{d} p}{(2 \pi)^{d}} \frac{1}{p^{2}}=0 \tag{3.4.2}
\end{equation*}
$$

This is due to the fact that, because of the absence of a mass scale, the IR and UV behaviors of the original integral are the same and therefore there is no value of $d$ for which the integral converges in the absence of suitable cutoffs. In general, DR sets consistently to 0 all the integrals of this kind, i.e., those which are homogeneous functions of the momenta and therefore lack a mass scale.

As stated above, dimensionally regularized integrals are characterized by dimensional poles. In order to understand their meaning consider $T_{d}$ in $d=4$ with a cutoff regularization: power counting (or a direct analysis) suggests that $T_{4}^{(\Lambda)}(\Lambda \gg m)=a_{2} \Lambda^{2}+$ $a_{0} m^{2} \ln (\Lambda / m)+$ finite, where $\Lambda$ is the cutoff. In DR, instead, $T_{d}$ shows (among other terms) a dimensional pole $\sim 1 / \epsilon$ for $d=4-\epsilon \rightarrow 4$. Interestingly enough, on the basis of

[^15]the approach discussed further below one can show that the residue at this pole is $a_{0} m^{2}$, i.e., the coefficient of $\ln (\Lambda / m)$ in $T_{4}^{(\Lambda)}$. In this sense, dimensional poles appear in an integral in DR if and only if the same integral regularized with a cutoff shows a logarithmic term $\sim \ln (\Lambda / m)$, either as a leading or subleading dependence on $\Lambda \gg m$. This is the reason why $T_{d}$ has also a dimensional pole in $d=2$ : in fact, one can easily see that $T_{2}^{(\Lambda)}$ has a leading dependence on $\ln (\Lambda / m)$, the coefficient of which is given by the residue of $T_{d}$ for $d \rightarrow 2$. This is also the reason why an integral such as the one in eq. (3.4.2) vanishes: if calculated with a cutoff one has $\int d^{d} p 1 / p^{2} \propto \Lambda^{d-2}$ which, for generic dimensionality $d$, shows no logarithmic dependence on $\Lambda$ and therefore the corresponding dimensionally regularized integral has no poles. The absence of poles (plus additional requirements of boundedness) is heuristically the reason why the corresponding dimensionally regularized integral can be taken vanishing.

Dimensional regularization without $\Gamma$ functions-Here we discuss a practical way of determining the leading dimensional pole of a generic loop integral without expressing it in terms of the $\Gamma$ functions which are typically associated with DR (see, e.g., ref. [1]). In order to be concrete, consider again $T_{d}$ in eq. (3.4.1) which can be expressed in terms of the $\Gamma$ function, as briefly discussed above. Being $\Gamma(x)$ an analytic function with isolated simple poles for $x=0,-1,-2, \ldots$, one concludes that $T_{d}$ develops a dimensional pole $\simeq 2 m^{2} /(4 \pi)^{2} \times 1 / \epsilon$ for $\epsilon=4-d \rightarrow 0$. This conclusion can actually be drawn without invoking the properties of $\Gamma(x)$, as we explain below. Additional details can be found in ref. [9].

The main point of DR is to assign a mathematical meaning to integrals performed in a non-integer (or even complex) number of dimensions $d$, i.e., to quantities like

$$
\begin{equation*}
I_{d} \equiv \int d^{d} p f(|\vec{p}|) \tag{3.4.3}
\end{equation*}
$$

where $f$ is a sufficiently smooth function. For integer $d=1,2, \ldots$, the integral can be done in spherical coordinates and therefore

$$
\begin{equation*}
I_{d}=\Omega_{d} \int_{0}^{\infty} d p p^{d-1} f(p) \tag{3.4.4}
\end{equation*}
$$

where

$$
\begin{equation*}
\Omega_{d}=\frac{2 \pi^{d / 2}}{\Gamma(d / 2)} \tag{3.4.5}
\end{equation*}
$$

is the solid angle in $d$-dimensions. ${ }^{11}$ The r.h.s. of eq. (3.4.4) is actually well-defined also for complex values of $d \in \mathbb{C}$, as long as the remaining integral converges. In fact, $\Omega_{d}$ in

[^16]eq. (3.4.5) is also well-defined for $d \in \mathbb{C}$, due to the properties of the $\Gamma$ function. Upon increasing the value of $\operatorname{Re} d$, however, the "radial" integral
\[

$$
\begin{equation*}
R_{d} \equiv \int_{0}^{\infty} d p p^{d-1} f(p) \tag{3.4.6}
\end{equation*}
$$

\]

might not converge for $p \rightarrow \infty$. In order to be concrete assume, for example, that

$$
\begin{equation*}
f(p \rightarrow \infty)=a_{-1} p^{-1}+a_{-2} p^{-2}+\ldots \quad \text { with } \quad \text { finite } f(0) \tag{3.4.7}
\end{equation*}
$$

Accordingly, $R_{d}$ converges within the strip $S_{1}$ in the complex plane with $0<\operatorname{Re} d<1$. For larger values of Re $d$, one can simply define $R_{d}$ via its analytic continuation, which can be determined by observing that the following identity

$$
\begin{equation*}
R_{d}=\int_{0}^{A} d p p^{d-1} f(p)+\int_{A}^{\infty} d p p^{d-1}\left[f(p)-a_{-1} p^{-1}\right]-\frac{a_{-1}}{d-1} A^{d-1} \tag{3.4.8}
\end{equation*}
$$

holds for arbitrary $A>0$ and $d \in S_{1}$, the latter being a necessary condition for the last integral to converge. In fact, the last term above is nothing but $\int_{A}^{\infty} d p p^{d-1} \times a_{-1} p^{-1}$, subtracted from the preceding term. The original integration $\int_{0}^{\infty}$ has been split as $\int_{0}^{A}+\int_{A}^{\infty}$ because, otherwise, the subtraction from $f(p)$ of the leading term $a_{-1} p^{-1}$ of its expansion for large $p$ would have spoiled the convergence of the integral for $p \rightarrow 0$. Now we note that the first integral on the r.h.s. of eq. (3.4.8) is finite as long as $\operatorname{Re} d>0$, while the integrand of the second integral behaves as $\sim a_{-2} p^{-2}$ for large $p$ and therefore it converges as long as $d$ is within the strip $S_{2}$ with $0<\operatorname{Re} d<2$, which extends up to $\operatorname{Re} d=2$. The third term, instead, is characterized by an isolated pole (on the real axis) for $d=1$, with a residue $-a_{-1}$ independent of the value of $A$. Overall, $R_{d}$ written as in eq. (3.4.8) converges within the strip $S_{2}$, which is larger than the original one $S_{1}$. As a consequence of this extension a dimensional pole at $d=1$ with residue $-a_{-1}$ emerges within $S_{2}$, at the boundary of $S_{1}$. The procedure outlined above can be repeated: in fact, the domain of convergence of the integral of $f(p)-a_{-1} p^{-1}$ in eq. (3.4.8) can be further extended by subtracting (and adding) the leading behavior of the integrand for $p \rightarrow \infty$, i.e., $a_{-2} p^{-2}$. As a result, within the strip $S_{2}, R_{d}$ in eq. (3.4.8) can also be written as
$R_{d}=\int_{0}^{A} d p p^{d-1} f(p)+\int_{A}^{\infty} d p p^{d-1}\left[f(p)-a_{-1} p^{-1}-a_{-2} p^{-2}\right]-\frac{a_{-1}}{d-1} A^{d-1}-\frac{a_{-2}}{d-2} A^{d-2}$,
which actually converges within the larger strip $S_{3}$ extending up to Red=3. As in the previous case, this further extension leaves behind a dimensional pole at $d=2$, with a residue $-a_{-2}$ that is independent of $A$. Proceeding further one can analytically continue $R_{d}$ in eq. (3.4.6) to arbitrarily large values of $\operatorname{Re} d$, leaving behind a set of poles for $d$ equal
to the opposite powers of $p$ which appear in the large- $p$ expansion of $f(p)$ (see eq. (3.4.7)); the residue at each of these poles is the opposite of the corresponding coefficient of this expansion. Given that $I_{d}=\Omega_{d} R_{d}$ (compare eqs. (3.4.4) and (3.4.6)), the same applies to $I_{d}$, the residues of which carry the additional factor $\Omega_{d}$. These are actually the dimensional poles that are encoded in the $\Gamma$ functions in terms of which, very often, one is able to cast the integral of the form (3.4.3) which are encountered in field-theoretical calculations.

As a simple application of the procedure discussed above, consider the integral in eq. (3.4.1): by expanding the integrand for $p^{2} \gg m^{2}$, one has

$$
\begin{equation*}
\frac{1}{p^{2}+m^{2}}=\frac{1}{p^{2}}-\frac{m^{2}}{p^{4}}+\frac{m^{4}}{p^{6}}+O\left(p^{-8}\right) \tag{3.4.10}
\end{equation*}
$$

and therefore the integral has dimensional poles for $d=2,4,6, \ldots$, with residues $-\Omega_{2} /(2 \pi)^{2}, m^{2} \Omega_{4} /(2 \pi)^{4},-m^{4} \Omega_{6} /(2 \pi)^{6}, \ldots$ Taking into account the expression of $\Omega_{d}$ in eq. (3.4.5), one readily verifies that these are the same poles as those of the $\Gamma$ function on the r.h.s. of eq. (3.4.1).

As a less trivial example, consider the integral that one encounters in the calculation of the one-loop vacuum polarization in QED. After having introduced the Feynman parameterization and done the Wick's rotation, it takes the form

$$
\begin{equation*}
\int \frac{d^{d} p}{(2 \pi)^{d}} \frac{A p^{2}+B}{\left(p^{2}+\Delta\right)^{2}} \quad \text { with } \quad \frac{A p^{2}+B}{\left(p^{2}+\Delta\right)^{2}}=\frac{A}{p^{2}}+\frac{B-2 A \Delta}{p^{4}}+O\left(p^{-6}\right) \tag{3.4.11}
\end{equation*}
$$

According to the discussion above, the dimensional pole for $d=4-\epsilon$ is simply given by

$$
\begin{equation*}
\frac{1}{-\epsilon} \times \frac{\Omega_{4}}{(2 \pi)^{4}} \times(-1)(B-2 A \Delta) \tag{3.4.12}
\end{equation*}
$$

without the need of invoking $\Gamma$ functions. Similarly, when calculating the renormalization of the QED vertex at one loop, one encounters the integral

$$
\begin{equation*}
\int \frac{d^{4} k}{(2 \pi)^{4}} \frac{-i}{k^{2}} \gamma_{\nu} i \frac{\not q+\not k+m}{(q+k)^{2}-m^{2}} \gamma_{\mu} i \frac{\not p+\not k+m}{(p+k)^{2}-m^{2}} \gamma^{\nu} \tag{3.4.13}
\end{equation*}
$$

In order to determine the dimensional pole for $d \rightarrow 4$, we need to know the coefficient of the term $\sim 1 / k^{4}$ in the large- $k$ expansion of the integrand, which, in this case, coincides with its leading term:

$$
\begin{equation*}
\frac{-i}{k^{2}} \gamma_{\nu} i \frac{\nmid k}{k^{2}} \gamma_{\mu} i \frac{\nmid k}{k^{2}} \gamma^{\nu} f i \gamma_{\mu} \frac{k^{2}}{\left(k^{2}\right)^{3}}, \tag{3.4.14}
\end{equation*}
$$

where we used the algebra of $\gamma$ matrices and the fact that $k_{\mu} k_{\nu} f\left(k^{2} / 4\right) \eta_{\mu \nu}$. After a Wick's rotation, the dimensional pole for $\epsilon=4-d \rightarrow 0$ of the integral (3.4.13) is therefore given by

$$
\begin{equation*}
i \times i \gamma_{\mu} \times \frac{1}{-\epsilon} \times \frac{\Omega_{4}}{(2 \pi)^{4}} \times(-1)=-\gamma_{\mu} \frac{1}{\epsilon} \frac{1}{2 \pi^{2}} \tag{3.4.15}
\end{equation*}
$$

## Chapter 4

## External Fields and Generating Functionals

It is often useful to study the properties of QFT's coupled with external sources. Consider a set of classical sources $J_{r}(x)$ coupled to some fields $\phi_{r}(x)$. The fields $\phi_{r}$ can be scalars, fermions or vectors, and will correspondingly couple to scalar, fermion or vector currents $J_{r}$. In general $\phi_{r}$ can also be composite fields, although we take them elementary for simplicity. We can define the $J$-dependent vacuum functional

$$
\begin{equation*}
\langle 0 \mid 0\rangle_{J} \equiv Z[J]=\int \mathcal{D} \phi_{r} e^{i S(\phi)+i \int d^{4} x J_{r}(x) \phi_{r}(x)} \tag{4.0.1}
\end{equation*}
$$

All correlation functions of the theory will be given by functional derivatives of $Z[J]$ with respect to $J_{r}$, evaluated at $J_{r}=0$ :

$$
\begin{equation*}
\left.\frac{1}{Z} \frac{\delta^{n} Z[J]}{\delta J_{r_{1}}\left(x_{1}\right) \ldots \delta J_{r_{n}}\left(x_{n}\right)}\right|_{J_{r}=0}=i^{n}\langle 0| T\left[\phi_{r_{1}}\left(x_{1}\right) \ldots \phi_{r_{n}}\left(x_{n}\right)\right]|0\rangle \tag{4.0.2}
\end{equation*}
$$

The functional $Z[J]$ is the generator of all amplitudes, connected and disconnected. We can define a similar functional generating connected amplitudes only, $W[J]$. This is simply given by

$$
\begin{equation*}
W[J]=-i \log Z[J] . \tag{4.0.3}
\end{equation*}
$$

Connected Green functions are defined similarly to eq.(4.0.2):

$$
\begin{equation*}
\left.\frac{\delta^{n} W[J]}{\delta J_{r_{1}}\left(x_{1}\right) \ldots \delta J_{r_{n}}\left(x_{n}\right)}\right|_{J_{r}=0}=i^{n-1}\langle 0| T\left[\phi_{r_{1}}\left(x_{1}\right) \ldots \phi_{r_{n}}\left(x_{n}\right)\right]|0\rangle_{c} \tag{4.0.4}
\end{equation*}
$$

For instance, for a single field $\phi$, we have

$$
\begin{equation*}
\left.\frac{\delta W[J]}{\delta J(x)}\right|_{J=0}=\left.(-i) \frac{1}{Z} \frac{\delta Z}{\delta J(x)}\right|_{J=0}=\langle 0| \phi(x)|0\rangle \tag{4.0.5}
\end{equation*}
$$

and hence

$$
\begin{equation*}
\langle 0| \phi(x)|0\rangle_{c}=\langle 0| \phi(x)|0\rangle . \tag{4.0.6}
\end{equation*}
$$

Taking one further functional derivative gives

$$
\begin{align*}
\left.\frac{\delta^{2} W[J]}{\delta J(x) \delta J(y)}\right|_{J=0} & =\left.(-i)\left(\frac{1}{Z} \frac{\delta^{2} Z}{\delta J(x) \delta J(y)}-\frac{1}{Z^{2}} \frac{\delta Z}{\delta J(x)} \frac{\delta Z}{\delta J(y)}\right)\right|_{J=0}  \tag{4.0.7}\\
& =i(\langle 0| T[\phi(x) \phi(y)]|0\rangle-\langle 0| \phi(x)|0\rangle\langle 0| \phi(y)|0\rangle)
\end{align*}
$$

from which we find the connected time-ordered two-point function

$$
\begin{equation*}
\langle 0| T[\phi(x) \phi(y)]|0\rangle_{c}=\langle 0| T[\phi(x) \phi(y)]|0\rangle-\langle 0| \phi(x)|0\rangle\langle 0| \phi(y)|0\rangle . \tag{4.0.8}
\end{equation*}
$$

Continuing in this way iteratively allows us to find all connected higher point functions.

### 4.1 The 1PI Effective Action

We can proceed further and define a functional $\Gamma[\Phi]$ that generates 1-particle irreducible amplitudes only. ${ }^{1}$ The source $\Phi(x)$ is defined as (omitting for simplicity the index $r$ from now on)

$$
\begin{equation*}
\Phi(x) \equiv \frac{\delta W[J]}{\delta J(x)} \tag{4.1.1}
\end{equation*}
$$

and is then a (in general very complicated) functional of the external sources $J .{ }^{2}$ The 1PI generator is defined as

$$
\begin{equation*}
\Gamma[\Phi]=W[J]-\int d^{4} x \Phi(x) J(x) \tag{4.1.2}
\end{equation*}
$$

It is the Legendre transform of $W[J]$, very much like the Lagrangian is the Legendre transform of the Hamiltonian. In eq. (4.1.2) it is understood that we have inverted eq. (4.1.1) so that $J=J[\Phi]$. We have

$$
\begin{equation*}
\frac{\delta \Gamma}{\delta \Phi(x)}=\int d^{4} y\left(\frac{\delta W}{\delta J(y)}-\Phi(y)\right) \frac{\delta J(y)}{\delta \Phi(x)}-J(x)=-J(x) \tag{4.1.3}
\end{equation*}
$$

When $J=0$, eq. (4.1.3) can be seen as an equation of motion of $\Phi$ with action $\Gamma$. $\Gamma$ is in fact the quantum generalization of the action and is also denoted the quantum effective action. It is an effective action in the sense that all loop amplitudes can be calculated as sum of connected tree diagrams obtained using $\Gamma[\Phi]$ instead of $S[\phi] .^{3}$ Thus, all loop effects

[^17]are encoded in $\Gamma[\Phi]$. In order to see this, let us define the functional generator $W_{\Gamma}[J]$ as the generator of connected amplitudes obtained starting from an action given by $\Gamma$ :
\[

$$
\begin{equation*}
e^{i W_{\Gamma}[J, g]}=\int \mathcal{D} \phi \exp \left(\frac{i}{g}\left(\Gamma[\phi]+\int d^{4} x J(x) \phi(x)\right)\right) . \tag{4.1.4}
\end{equation*}
$$

\]

In (4.1.4) we have introduced a parameter $g$, that plays the role of $\hbar$ and is a loop-counting parameter. For connected graphs the number of loops $L$ is given by

$$
\begin{equation*}
L=I-V+1 \tag{4.1.5}
\end{equation*}
$$

where $I$ is the total number of propagators and $V$ is the total number of vertices in the graph. Indeed, this coincides with the number of independent virtual momenta entering the graph, the +1 coming from the conservation of the overall external momenta. We have already used this relation in eq. (3.1.6). We see from eq. (4.1.4) that the propagator of each field is proportional to $g$ (inverse of the action), while any vertex goes like $1 / g$. Thus the $L$-loop contribution to $W_{\Gamma}, W_{\Gamma}^{(L)}$, goes like $g^{I-V}=g^{L-1}$. In a loop expansion we can write

$$
\begin{equation*}
W_{\Gamma}[J, g]=\sum_{L=0}^{\infty} g^{L-1} W_{\Gamma}^{(L)}[J] . \tag{4.1.6}
\end{equation*}
$$

Consider now in eq. (4.1.4) the tree-level contribution only, which can be selected by taking the limit $g \rightarrow 0$. By using a saddle point approximation we have

$$
\begin{equation*}
\lim _{g \rightarrow 0} e^{i W_{\Gamma}[J, g]}=e^{\frac{i}{g} W_{\Gamma}^{(0)}[J]}=\exp \left(\frac{i}{g}\left(\Gamma[\Phi]+\int d^{4} x J(x) \Phi(x)\right)\right) \tag{4.1.7}
\end{equation*}
$$

where $\Phi(x)$ extremizes the exponential, namely

$$
\begin{equation*}
\left.\frac{\delta \Gamma}{\delta \phi(x)}\right|_{\phi=\Phi}+J(x)=0 \tag{4.1.8}
\end{equation*}
$$

Comparing eq. (4.1.2) with eq. (4.1.7) and eq. (4.1.3) with (4.1.8), we conclude that

$$
\begin{equation*}
W_{\Gamma}^{(0)}[J]=W[J] \tag{4.1.9}
\end{equation*}
$$

Stated differently, we can write

$$
\begin{equation*}
i W[J]=\int_{\substack{\text { Connected } \\ \text { tree graphs }}} \mathcal{D} \phi e^{i \Gamma(\phi)+i \int J \phi} . \tag{4.1.10}
\end{equation*}
$$

Since any connected graph is a tree with 1PI vertices, we conclude that $\Gamma(\phi)$ generates all 1PI amplitudes:

$$
\begin{equation*}
\Gamma(\phi)=\sum_{n=0}^{\infty} \frac{1}{n!} \int d^{4} x_{1} \ldots \int d^{4} x_{n} \Gamma^{(n)}\left(x_{1}, \ldots, x_{n}\right) \phi\left(x_{1}\right) \ldots \phi\left(x_{n}\right) \tag{4.1.11}
\end{equation*}
$$

The quantum effective action $\Gamma$ can also be computed directly starting from the action $S$, without passing through $W[J]$. One has

$$
\begin{equation*}
e^{i \Gamma\left(\phi_{0}\right)}=\int_{1 P I} \mathcal{D} \phi e^{i S\left(\phi_{0}+\phi\right)}, \tag{4.1.12}
\end{equation*}
$$

where 1PI means that we should only consider 1PI graphs in evaluating the path integral and $\phi_{0}$ is an arbitrary field configuration. This formula is a direct consequence of the fact that $\Gamma$ is the quantum version of $S$. Given a classical configuration $\phi_{0}$, the quantum action of $\phi_{0}$ is given by integrating out quantum fluctuations $\phi$ around the classical configuration $\phi_{0}$. We invite the reader to check the validity (and the usefulness) of eq. (4.1.12) by formally computing $\Gamma(\phi)$ at one-loop level using eq. (4.1.12) and directly from the definition (4.1.2).

### 4.2 The Coleman-Weinberg Effective Potential

The effective potential $V_{\text {eff }}$ is defined as the quantum action $\Gamma$, evaluated for constant field configurations $\phi_{0}:^{4}$

$$
\begin{equation*}
\Gamma\left(\phi=\phi_{0}\right)=-V_{4} V_{e f f}\left(\phi_{0}\right), \tag{4.2.1}
\end{equation*}
$$

where $V_{4}$ is the volume of space-time, coming from $\int d^{4} x$. It is the quantum analogue of the classical potential $V_{0}$ defined as $S\left(\phi=\phi_{0}\right)=-V_{4} V_{0}\left(\phi_{0}\right)$. Let us explicitly compute $V_{e f f}$ at one-loop level for the simple scalar theory

$$
\begin{equation*}
\mathcal{L}(\phi)=\frac{1}{2}\left(\partial_{\mu} \phi_{B}\right)\left(\partial^{\mu} \phi_{B}\right)-\frac{1}{2} m_{B}^{2} \phi_{B}^{2}-\frac{\lambda_{B}}{4!} \phi_{B}^{4} . \tag{4.2.2}
\end{equation*}
$$

First of all, we express the Lagrangian in terms of the physical parameters and fields:

$$
\begin{equation*}
\phi_{B}=\sqrt{Z} \phi, \quad m_{B}^{2} Z=m^{2}+\delta m^{2}, \quad \lambda_{B} Z^{2}=Z_{\lambda} \lambda \tag{4.2.3}
\end{equation*}
$$

so that

$$
\begin{equation*}
\mathcal{L}(\phi)=\frac{1}{2}\left(\partial_{\mu} \phi\right)\left(\partial^{\mu} \phi\right)-\frac{1}{2} m^{2} \phi^{2}-\frac{\lambda}{4!} \phi^{4}+\mathcal{L}_{\text {c.t. }}(\phi), \tag{4.2.4}
\end{equation*}
$$

where the counter-term Lagrangian reads

$$
\begin{equation*}
\mathcal{L}_{\text {c.t. }}(\phi)=\frac{1}{2}(Z-1)\left(\partial_{\mu} \phi\right)\left(\partial^{\mu} \phi\right)-\frac{1}{2} \delta m^{2} \phi^{2}-\frac{\lambda}{4!}\left(Z_{\lambda}-1\right) \phi^{4} . \tag{4.2.5}
\end{equation*}
$$

We compute the effective potential starting from eq. (4.1.12) with $\phi_{0}$ constant. We get

$$
\begin{align*}
\mathcal{L}\left(\phi_{0}+\phi\right)= & -\frac{1}{2} m^{2} \phi_{0}^{2}-\frac{\lambda}{4!} \phi_{0}^{4}+\phi\left(-m^{2} \phi_{0}-\frac{\lambda}{6} \phi_{0}^{3}\right) \\
& +\frac{1}{2}\left(\partial_{\mu} \phi\right)\left(\partial^{\mu} \phi\right)-\frac{1}{2} \phi^{2}\left(m^{2}+\frac{\lambda}{2} \phi_{0}^{2}\right)-\frac{\lambda}{6} \phi_{0} \phi^{3}-\frac{\lambda}{4!} \phi^{4}+\mathcal{L}_{\text {c.t. }}\left(\phi_{0}+\phi\right) . \tag{4.2.6}
\end{align*}
$$

[^18]We have ordered the terms according to the powers of quantum fluctuations $\phi$. The first three terms in the first row of eq. (4.2.6) reproduce the initial tree-level potential. Terms cubic or quartic in $\phi$ cannot contribute at one-loop level and start contributing at twoloops. The term linear in $\phi$ can be neglected altogether: it can only enter one-particle reducible graphs and never 1PI amplitudes. The counter-terms are non-vanishing only at loop order, so at one-loop level we should only keep $\mathcal{L}_{\text {c.t. }}\left(\phi_{0}\right)$. The relevant terms of the Lagrangian (4.2.6) are then

$$
\begin{equation*}
\mathcal{L}\left(\phi_{0}+\phi\right) \supset-\frac{1}{2} m^{2} \phi_{0}^{2}-\frac{\lambda}{4!} \phi_{0}^{4}+\frac{1}{2}\left(\partial_{\mu} \phi\right)\left(\partial^{\mu} \phi\right)-\frac{1}{2} \mu^{2}\left(\phi_{0}\right) \phi^{2}+\mathcal{L}_{\text {c.t. }}\left(\phi_{0}\right), \tag{4.2.7}
\end{equation*}
$$

where we have defined the $\phi_{0}$-dependent mass

$$
\begin{equation*}
\mu^{2}\left(\phi_{0}\right) \equiv m^{2}+\frac{\lambda}{2} \phi_{0}^{2} . \tag{4.2.8}
\end{equation*}
$$

Up to one-loop level we have

$$
\begin{equation*}
V_{e f f}\left(\phi_{0}\right)=V_{0}\left(\phi_{0}\right)+V_{1}\left(\phi_{0}\right)+V_{\text {c.t. }}\left(\phi_{0}\right), \tag{4.2.9}
\end{equation*}
$$

with

$$
\begin{equation*}
V_{\text {c.t. }}\left(\phi_{0}\right)=\frac{1}{2} \delta m^{2} \phi_{0}^{2}+\frac{\lambda}{4!}\left(Z_{\lambda}-1\right) \phi_{0}^{4} \tag{4.2.10}
\end{equation*}
$$

and

$$
\begin{align*}
e^{-i V_{4} V_{1}\left(\phi_{0}\right)} & =\int \mathcal{D} \phi e^{i \int d^{4} x\left(\frac{1}{2}\left(\partial_{\mu} \phi\right)\left(\partial^{\mu} \phi\right)-\frac{1}{2} \mu^{2}\left(\phi_{0}\right) \phi^{2}\right)}  \tag{4.2.11}\\
& =\mathcal{N} \operatorname{det}^{-\frac{1}{2}}\left(\square+\mu^{2}\left(\phi_{0}\right)\right)=\mathcal{N} e^{-\frac{1}{2} \operatorname{Tr} \log \left(\square+\mu^{2}\left(\phi_{0}\right)\right)}
\end{align*}
$$

The determinant or the trace of the differential operator $\square+\mu^{2}\left(\phi_{0}\right)$ is easily computed in momentum space, where the operator is diagonal. We have

$$
\begin{equation*}
\log \left(\square+\mu^{2}\left(\phi_{0}\right)\right)(p, q)=\delta(p-q) \log \left(-p^{2}+\mu^{2}\left(\phi_{0}\right)\right) \tag{4.2.12}
\end{equation*}
$$

and hence

$$
\begin{equation*}
-i V_{4} V_{1}\left(\phi_{0}\right)=-\frac{1}{2} \delta^{(4)}(0) \int d^{4} p \log \left(-p^{2}+\mu^{2}\left(\phi_{0}\right)\right) . \tag{4.2.13}
\end{equation*}
$$

The $\delta^{(4)}(0)$ accounts for the space-time volume, $\delta^{(4)}(0) \rightarrow V_{4} /(2 \pi)^{4}$. Wick rotating $p_{0}=$ $i p_{4}$, we finally get

$$
\begin{equation*}
V_{1}\left(\phi_{0}\right)=\frac{1}{2} \int \frac{d^{4} p_{E}}{(2 \pi)^{4}} \log \left(p_{E}^{2}+\mu^{2}\left(\phi_{0}\right)\right), \tag{4.2.14}
\end{equation*}
$$

where $p_{E}^{2}=p_{1}^{2}+p_{2}^{2}+p_{3}^{2}+p_{4}^{2}$. The integral in eq. (4.2.14) is highly divergent. In terms of the degree of divergence defined in section 3.1 it has $\delta=4$. We can lower this divergence
by taking derivative with respect to $\mu^{2}\left(\phi_{0}\right)$. Since each derivative lowers $\delta$ by two, we have to derive three times to get a finite result:

$$
\begin{equation*}
\frac{d^{3} V_{1}\left(\phi_{0}\right)}{\left(d \mu^{2}\right)^{3}}=\int \frac{d^{4} p_{E}}{(2 \pi)^{4}} \frac{1}{\left(p_{E}^{2}+\mu^{2}\left(\phi_{0}\right)\right)^{3}}=\int_{0}^{\infty} \frac{d p_{E}}{8 \pi^{2}} \frac{p_{E}^{3}}{\left(p_{E}^{2}+\mu^{2}\left(\phi_{0}\right)\right)^{3}}=\frac{1}{32 \pi^{2}} \frac{1}{\mu^{2}\left(\phi_{0}\right)} \tag{4.2.15}
\end{equation*}
$$

Integrating back three times gives, modulo an irrelevant constant factor,

$$
\begin{equation*}
V_{1}\left(\phi_{0}\right)=\frac{1}{64 \pi^{2}} \mu^{4}\left(\phi_{0}\right) \log \mu^{2}\left(\phi_{0}\right)+A \mu^{2}\left(\phi_{0}\right)+B \mu^{4}\left(\phi_{0}\right) \tag{4.2.16}
\end{equation*}
$$

In analogy to the case of loop amplitudes considered in section 3.2, the effective potential is the sum of a non-analytic (in $\phi_{0}$ ) and finite piece, the first term in eq.(4.2.16), plus local divergent terms. The divergences hidden in the constants $A$ and $B$ are cancelled by a proper choice of the counter-terms $\delta m^{2}$ and $Z_{\lambda}$ in eq.(4.2.10). Putting all together, we finally get the total tree+one-loop level potential:

$$
\begin{equation*}
V_{e f f}\left(\phi_{0}\right)=\frac{1}{2} m^{2} \phi_{0}^{2}+\frac{\lambda}{4!} \phi_{0}^{4}+\frac{1}{64 \pi^{2}} \mu^{4}\left(\phi_{0}\right) \log \mu^{2}\left(\phi_{0}\right) . \tag{4.2.17}
\end{equation*}
$$

When the mass vanishes, the potential (4.2.17) becomes $^{5}$

$$
\begin{equation*}
V_{e f f}\left(\phi_{0}\right)=\frac{\lambda}{4!} \phi_{0}^{4}+\frac{\lambda^{2}}{256 \pi^{2}} \phi_{0}^{4} \log \phi_{0}^{2} \tag{4.2.18}
\end{equation*}
$$

The effective potential, often also denoted by Coleman-Weinberg (CW) potential, from the authors that first discussed it [10], is a useful tool in QFT. It is in particular crucial to determine the correct vacuum of a theory when the tree level potential does not fix it. We will come back to the CW potential in section 5.10 and in the final chapter 11 .

### 4.3 A Subtlety about Effective Potentials**

In this section we will discuss a subtlety regarding effective potentials. In eq.(4.1.2) we have defined the 1PI action $\Gamma[\Phi]$ as the Legendre transform of $W[J]$. It can be shown that the effective potential, defined as in eq.(4.2.1) with $\Gamma[\Phi]$ above, must necessarily be a convex function [11, 12]. This is however puzzling, since the classical potential might clearly be non-convex to start with. If the system is parametrically weakly coupled, it is hard to imagine that quantum corrections could significantly modify the classical potential. The resolution of the puzzle is a bit tricky and is related to the fact that in general the 1PI

[^19]action defined in eq.(4.1.2) does not correspond to that defined in eq.(4.1.12). In particular, the effective potential $V_{1 \text { PI }}$ coming from eq.(4.1.12) is the natural extension of the classical potential $V_{0}$ and can in general be non-convex, while the effective potential $V_{e f f}$ defined as the constant field configurations of eq.(4.1.2) is the one that must necessarily be convex. Starting from eq.(4.1.12), we can define $W[J]$ as the Legendre transform of $\Gamma[\Phi]_{1 \mathrm{PI}}$, which is the $\Gamma$ appearing on the left-hand-side of eq.(4.1.12). Taking now the Legendre transform of the so derived $W[J]$ we get $\Gamma[\Phi]_{e f f}$. We then see that $V_{\text {eff }}$ is the double Legendre transform of $V_{1 \mathrm{PI}}$, also called its convex envelope. Only for convex functions we are ensured that $V_{\text {eff }}=V_{1 \mathrm{PI}}$. Physically, $V_{\text {eff }}\left(\phi_{0}\right)$ represents the expectation value of the energy density in the state $|\Psi\rangle$ that minimizes $\langle\Psi| H|\Psi\rangle$ and such that $\langle\Psi| \phi(x)|\Psi\rangle=\phi_{0}$. This quantity is always real and, as we said, must be convex everywhere. The key point on how $V_{\text {eff }}\left(\phi_{0}\right)$ is convex starting form a non-convex classical potential is understood from a purely classical analysis, where $\langle\Psi| \phi(x)|\Psi\rangle$ represents the spatial average of the field $\phi(\vec{x})$. Consider a classical potential $V_{c}$ like that depicted in the left panel of fig.4.1, with two minima at $\phi(\vec{x})= \pm \sigma$. For $\left|\phi_{0}\right| \geq \sigma$, the field configuration minimizing the energy is given by $\phi(\vec{x})=\phi_{0}$ over the whole space, and $V_{\text {eff }}\left(\phi_{0}\right)=V_{c}\left(\phi_{0}\right)$. For $\left|\phi_{0}\right|<\sigma$ the mininum energy is given by an inhomogeneous state where $\phi(\vec{x})=\sigma$ in a fraction $\left(\sigma+\phi_{0}\right) / 2 \sigma$ of space, and $\phi(\vec{x})=-\sigma$ in the remaining $\left(\sigma-\phi_{0}\right) / 2 \sigma$ fraction of space, so that the spatial average of the field is
\[

$$
\begin{equation*}
\sigma \times \frac{\sigma+\phi_{0}}{2 \sigma}+(-\sigma) \times \frac{\sigma-\phi_{0}}{2 \sigma}=\phi_{0}, \tag{4.3.1}
\end{equation*}
$$

\]

as it should be. The energy density in such inhomogeneous state is all concentrated in the regions between the two phases. In the infinite volume limit the boundary energy contribution becomes negligible, and hence $V_{e f f}\left(\phi_{0}\right)=0$ for $\left|\phi_{0}\right|<\sigma$. The potential $V_{e f f}$, depicted in the right panel of fig.4.1 is then everywhere convex.

In the quantum case, the inhomogeneous state is replaced by a linear combination of the two vacua at $\pm \sigma$, call them $| \pm\rangle$. For $\left|\phi_{0}\right|<\sigma$ the mininum energy is given by a state of the form $\alpha|+\rangle+\beta|-\rangle$, with $|\alpha|^{2}+|\beta|^{2}=1$, and $|\alpha|^{2}-|\beta|^{2}=\phi_{0}$. The ending result is the same, giving rise to a flat effective potential $V_{\text {eff }}$ for $\left|\phi_{0}\right|<\sigma$. As we will see in more detail in chapter 8 in the context of spontaneous symmetry breaking, inhomogeneous states of the form $\alpha|+\rangle+\beta|-\rangle$ violate cluster decomposition and should not be considered viable vacua for a QFT. If the vacua are degenerate, like in fig.4.1, the system will be unstable under any external small perturbation and will fall in either the vacuum $|+\rangle$ or
 by $V_{1 \mathrm{PI}}$, in agreement with our naive intuition.


Figure 4.1: Classical (left) and effective (right) potentials $V_{c}$ and $V_{e f f}$.

### 4.4 Functional Relations

The formalism based on the generating functions introduced and discussed above is particularly useful for practical applications because it can make the derivation of important relationships rather straightforward. This is particularly true if one proceeds heuristically, as we will do, without taking into account the effects of renormalization. In other words, we will effectively be dealing in this section with bare quantities, without introducing counterterms. This is justified by the fact that eventually the same results apply for the renormalized expressions, but in a much simpler way! See e.g. ref.[9] for a proper derivation where renormalization is taken into account.

### 4.4.1 Schwinger-Dyson Equation

The Schwinger-Dyson equations are the quantum mechanical analogue of the classical equations of motion. They are derived from the observation (based on the extension of the concept of ordinary differentiation and integration to the case of functional differentiation and integration) that the functional integral of a functional derivative vanishes if one can neglect boundary terms, which is always assumed here. Accordingly, taking the exponential in eq. (4.0.1) as the function to be differentiated, one concludes that

$$
\begin{equation*}
0=\int \mathcal{D} \phi \frac{\delta}{\delta \phi(x)} e^{i S(\phi)+i \int d^{4} x^{\prime} J\left(x^{\prime}\right) \phi\left(x^{\prime}\right)}=\int \mathcal{D} \phi\left[i \frac{\delta S(\phi)}{\delta \phi(x)}+i J(x)\right] e^{i S(\phi)+i \int d^{4} x^{\prime} J\left(x^{\prime}\right) \phi\left(x^{\prime}\right)} \tag{4.4.1}
\end{equation*}
$$

Setting $J=0$ in eq.(4.4.1) gives

$$
\begin{equation*}
\left\langle\frac{\delta S}{\delta \phi(x)}\right\rangle=0 \tag{4.4.2}
\end{equation*}
$$

which is the strict quantum-mechanical analogue of the classical equations of motion. Applying $n$ functional derivatives with respect to the source $J$ to eq.(4.4.1), and then
setting $J=0$, gives the Schwinger-Dyson equations

$$
\begin{equation*}
i\left\langle\frac{\delta S}{\delta \phi(x)} \phi\left(x_{1}\right) \ldots \phi\left(x_{n}\right)\right\rangle=-\sum_{i=1}^{n} \delta^{(4)}\left(x-x_{i}\right)\left\langle\phi\left(x_{1}\right) \ldots \hat{\phi}\left(x_{i}\right) \ldots \phi\left(x_{n}\right)\right\rangle, \tag{4.4.3}
\end{equation*}
$$

where the hat means that that operator should be omitted from the correlator. We see that at the quantum level the equations of motion turn into an infinite set of relations among different Green functions. We will make use of eqs.(4.4.3) later on in these lectures.

### 4.4.2 Symmetries and Ward-Takahashi Identities

Symmetries are important in QFT because they constrain the form of amplitudes, correlation functions, and establish relationships among them. This can be rather easily seen within the functional formalism, as we discuss below in generality. It is important to realize, however, that the formal manipulations presented here assume that it is possible to regularize the theory - therefore providing a meaning to the functional integral - in such a way that the original symmetry of the theory is preserved.

Let us consider a local field theory characterized by a Lagrangian $\mathcal{L}$. An infinitesimal transformation

$$
\begin{equation*}
\phi(x) \mapsto \phi^{\prime}(x)=\phi(x)+\epsilon \Delta \phi(x) \tag{4.4.4}
\end{equation*}
$$

parameterized by a parameter $\epsilon$ is a symmetry of the theory if, correspondingly, the action $S=\int d^{4} x \mathcal{L}\left(\phi, \partial_{\mu} \phi\right)$ associated with $\mathcal{L}$ does not change. In order to be so, the variation of the Lagrangian induced by eq. (4.4.4) has to take the form of a total derivative. This requirement can be translated in requiring that there exists some "current" $\mathcal{J}^{\mu}$ such that

$$
\begin{equation*}
\mathcal{L}\left(\phi^{\prime}, \partial_{\mu} \phi^{\prime}\right)=\mathcal{L}\left(\phi, \partial_{\mu} \phi\right)+\epsilon \partial_{\mu} \mathcal{J}^{\mu}+\mathcal{O}\left(\epsilon^{2}\right) \tag{4.4.5}
\end{equation*}
$$

(Note that $\mathcal{J}^{\mu}$ can also be a Lorentz tensor, as in the specific case discussed in sec. 9.7.) A direct calculation of the variation of the Lagrangian induced by the transformation shows that

$$
\begin{equation*}
\partial_{\mu} \mathcal{J}^{\mu}=\frac{\delta \mathcal{L}}{\delta \phi} \Delta \phi+\frac{\delta \mathcal{L}}{\delta \partial_{\mu} \phi} \partial_{\mu} \Delta \phi \tag{4.4.6}
\end{equation*}
$$

By rewriting the second term on the r.h.s. as $\partial_{\mu}\left[\Delta \phi \delta \mathcal{L} / \delta \partial_{\mu} \phi\right]-\Delta \phi \partial_{\mu} \delta \mathcal{L} / \delta \partial_{\mu} \phi$ one finds that

$$
\begin{equation*}
\partial_{\mu} j^{\mu}=-\Delta \phi\left[\frac{\delta \mathcal{L}}{\delta \phi}-\partial_{\mu} \frac{\delta \mathcal{L}}{\delta \partial_{\mu} \phi}\right], \tag{4.4.7}
\end{equation*}
$$

where we introduced the current

$$
\begin{equation*}
j^{\mu}(x) \equiv \frac{\delta \mathcal{L}}{\delta \partial_{\mu} \phi} \Delta \phi-\mathcal{J}^{\mu} \tag{4.4.8}
\end{equation*}
$$

The r.h.s. of eq. (4.4.7) vanishes on the field configuration which minimizes the action (i.e., without quantum fluctuations) and which satisfies Euler-Lagrange equation: accordingly $j_{\mu}$ is nothing but the classical Noether's current associated with the symmetry (4.4.4) and $j_{\mu}$ is the corresponding conserved current from which one can derive the conserved charge. In the case of QFT, where quantum fluctuations are inevitably present, relations such as eq. (4.4.7) have to be properly interpreted and acquire a definite meaning only when inserted in correlation functions. For the purpose of understanding the consequence of the symmetry (4.4.4) on the generating functions, it is convenient to consider how the Lagrangian density changes when the parameter $\epsilon$ in eq. (4.4.4) is assumed to be space-dependent (and, in general, the local transformation is no longer a symmetry of the theory): by direct calculation one can verify that

$$
\begin{align*}
\mathcal{L}\left(\phi^{\prime}, \partial_{\mu} \phi^{\prime}\right) & =\mathcal{L}\left(\phi, \partial_{\mu} \phi\right)+\epsilon(x) \partial_{\mu} \mathcal{J}^{\mu}+\frac{\delta \mathcal{L}}{\delta \partial_{\mu} \phi} \Delta \phi \partial_{\mu} \epsilon(x)+\mathcal{O}\left(\epsilon^{2}\right)  \tag{4.4.9}\\
& f \mathcal{L}\left(\phi, \partial_{\mu} \phi\right)+j^{\mu}(x) \partial_{\mu} \epsilon(x)+\mathcal{O}\left(\epsilon^{2}\right)
\end{align*}
$$

where $f$ indicates that the equality is valid after integration over the space, i.e., at the level of the action (under the assumption of vanishing boundary terms). Though the discussion so far is completely general, we shall focus below on the case in which the symmetry is linearly realized, meaning that $\Delta \phi$ is an arbitrary linear function of $\phi$. This is the case, for example of global $U(1)$ or other internal symmetries for multiplets such as

$$
\begin{equation*}
\phi_{l}(x) \mapsto \phi_{l}^{\prime}(x)=\phi_{l}(x)+\epsilon^{\alpha}\left(t_{\alpha}\right)_{l}^{m} \phi_{m}(x), \tag{4.4.10}
\end{equation*}
$$

where $\left\{t_{\alpha}\right\}_{\alpha}$ are the generators of the symmetry in a suitable representation. Alternatively, for infinitesimal space-time translations, $\epsilon \Delta \phi$ in eq. (4.4.4) is replaced by $\epsilon^{\mu} \partial_{\mu} \phi$. This case is discussed in detail in sec. 9.7. In order to work out the consequences of the symmetry, consider again the generating function $Z[J]$ in eq. (4.0.1) and denote by $\phi^{\prime}$ the "integration variable" in the functional integral, with the aim of eventually performing a change of variable towards the field $\phi$ which $\phi^{\prime}$ is connected to via eq. (4.4.4) with a space-dependent parameter $\epsilon=\epsilon(x)$ :

$$
\begin{align*}
Z[J] & =\int \mathcal{D} \phi^{\prime} e^{i S\left(\phi^{\prime}\right)+i \int d^{4} x J(x) \phi^{\prime}(x)}=\int \mathcal{D} \phi^{\prime} e^{i S(\phi)+i \int d^{4} x j^{\mu} \partial_{\mu} \epsilon(x)+i \int d^{4} x J(x)[\phi(x)+\epsilon \Delta \phi(x)]} \\
& =\int \mathcal{D} \phi\left(1+i \int d^{4} x \epsilon(x)\left[-\partial_{\mu} j^{\mu}(x)+J(x) \Delta \phi(x)\right]+\mathcal{O}\left(\epsilon^{2}\right)\right) e^{i S(\phi)+i \int d^{4} x J(x) \phi(x)} \\
& =Z[J]+i \int d^{4} x \epsilon(x)\left(-\left\langle\partial_{\mu} j^{\mu}(x)\right\rangle_{J}+J(x)\langle\Delta \phi(x)\rangle_{J}\right)+\mathcal{O}\left(\epsilon^{2}\right), \tag{4.4.11}
\end{align*}
$$

where on the second line we expanded up to first order in $\epsilon$ and used the fact that, being the symmetry realized linearly, the Jacobian of the change of variable $\phi^{\prime} \mapsto \phi$ in the functional integration measure is independent of the fields and can be absorbed in the overall normalization of the measure. The previous equation has to be valid for an arbitrary choice of $\epsilon(x)$ and therefore it implies that

$$
\begin{equation*}
\left\langle\partial_{\mu} j^{\mu}(x)\right\rangle_{J}=J(x)\langle\Delta \phi(x)\rangle_{J} . \tag{4.4.12}
\end{equation*}
$$

This relation is an example of a Ward-Takahashi (WT) identity associated with the symmetry (4.4.4) and constitutes the equivalent of the Noether theorem in QFT. By taking functional derivatives with respect to $J$, this equation implies an infinite set of relationships between correlation functions of the fields $\phi$ on the r.h.s. and those of the fields with an insertion of the current operator $j_{\mu}(x)$ on the l.h.s. As an explicit example, consider the internal symmetry (4.4.10), where the associated current is given by $j_{\alpha, \mu}=\left(\partial_{\mu} \phi\right)_{l}\left(t_{\alpha}\right)_{l}^{m} \phi_{m}$. By differentiating eq.(4.4.12) with respect to $J_{n_{1}}\left(x_{1}\right)$ and $J_{n_{2}}\left(x_{2}\right)$ and by then setting $J=0$ we get

$$
\begin{align*}
& i \partial_{\mu}\left\langle j_{\alpha}^{\mu}(x) \phi_{n_{1}}\left(x_{1}\right) \phi_{n_{2}}\left(x_{2}\right)\right\rangle= \\
& \quad=\delta\left(x-x_{1}\right)\left(t_{\alpha}\right)_{n_{1}}^{m}\left\langle\phi_{m}\left(x_{1}\right) \phi_{n_{2}}\left(x_{2}\right)\right\rangle+\delta\left(x-x_{2}\right)\left(t_{\alpha}\right)_{n_{2}}^{m}\left\langle\phi_{m}\left(x_{2}\right) \phi_{n_{1}}\left(x_{1}\right)\right\rangle \tag{4.4.13}
\end{align*}
$$

We point out that, from the point of view of QFT, the current $j_{\mu}(x)$ defined in eq. (4.4.8) is a so-called composite operator because it involves the product of various quantum fields taken at the same point in space-time. After the renormalization of the theory, correlation functions with the insertion of composite fields still show ultraviolet divergences and need to be additionally renormalized by introducing suitable renormalization constants for these operators. As we shall discuss in more detail in sec. 5.9, the presence of these non-trivial additional renormalization constants implies that the composite operators acquire an anomalous dimension. However, conserved currents $j_{\mu}(x)$ associated to symmetries of the theory do not need to be renormalized and therefore they do not acquire an anomalous dimension. This is clear from the example in eq.(4.4.13). Correlation functions with one insertion of $j_{\mu}$ can be expressed in terms of correlation functions without the current. When the latter are properly renormalized by the introduction of suitable counter-terms and renormalization constants, the former will also be finite. In other words, a conserved current into a correlation function does not cause additional divergences to appear.

It is important to point out here that the effective action $\Gamma$ is not always invariant under the same field transformations for which the classical action $S$ is. The transformations of $S$ and $\Gamma$ coincide only when they act at most linearly on fields. Let us discuss in more
detail this important point. Consider a transformation of the form

$$
\begin{equation*}
\phi(x) \mapsto \phi^{\prime}(x)=\phi(x)+\epsilon F[\phi(x)] \tag{4.4.14}
\end{equation*}
$$

where $F[\phi]$ is a generic functional of the field $\phi$, not necessarily linear. We shall encounter non-linear transformations in section 6.3 when we will study the BRST symmetries in the context of non-abelian gauge theories. Assume now that both the action $S(\phi)$ and the integration measure are invariant under the transformation $\phi \rightarrow \phi^{\prime}$, such that $S(\phi)=S\left(\phi^{\prime}\right)$ and $\mathcal{D} \phi^{\prime}=\mathcal{D} \phi$. In terms of the generating function $Z[J]$ one finds

$$
\begin{align*}
Z[J] & =\int \mathcal{D} \phi^{\prime} e^{i S\left(\phi^{\prime}\right)+i \int d^{4} x J(x) \phi^{\prime}(x)}=\int \mathcal{D} \phi e^{i S(\phi)+i \int d^{4} x J(x)\{\phi(x)+\epsilon F[\phi(x)]\}} \\
& =\int \mathcal{D} \phi\left\{1+i \epsilon \int d^{4} x J(x) F[\phi(x)]+\mathcal{O}\left(\epsilon^{2}\right)\right\} e^{i S(\phi)+i \int d^{4} x J(x) \phi(x)} \tag{4.4.15}
\end{align*}
$$

which implies

$$
\begin{equation*}
\int d^{4} x\langle F[\phi(x)]\rangle_{J} J(x)=0 . \tag{4.4.16}
\end{equation*}
$$

Using eq.(4.1.3), we can rewrite eq.(4.4.16) as

$$
\begin{equation*}
\int d^{4} x\langle F[\phi(x)]\rangle_{J(\Phi)} \frac{\delta \Gamma}{\delta \Phi(x)}=0 \tag{4.4.17}
\end{equation*}
$$

which implies that the 1PI action is invariant for $\Phi \rightarrow \Phi+\epsilon\langle F[\phi(x)]\rangle_{J(\Phi)}$ :

$$
\begin{equation*}
\Gamma\left(\Phi+\epsilon\langle F[\phi(x)]\rangle_{J(\Phi)}\right)=\Gamma(\Phi)+\epsilon \int d^{4} x\langle F[\phi(x)]\rangle_{J(\Phi)} \frac{\delta \Gamma}{\delta \Phi(x)}=\Gamma(\Phi) . \tag{4.4.18}
\end{equation*}
$$

Notice that, in general,

$$
\begin{equation*}
\langle F[\phi(x)]\rangle_{J(\Phi)} \neq F(\Phi) . \tag{4.4.19}
\end{equation*}
$$

They coincide only for transformations that are at most linear in the fields, namely for ${ }^{6}$

$$
\begin{equation*}
F(\phi)=s(x)+\int d^{4} y t(x, y) \phi(y) \tag{4.4.20}
\end{equation*}
$$

In this case, and only in this case, we have

$$
\begin{equation*}
\langle F[\phi(x)]\rangle_{J(\Phi)}=s(x)+\int d^{4} y t(x, y)\langle\phi(x)\rangle_{J(\Phi)}=s(x)+\int d^{4} y t(x, y) \Phi(y)=F(\Phi), \tag{4.4.21}
\end{equation*}
$$

where the second identity is a consequence of the definition of $\Phi$.
The general approach described above can be used in order to derive the consequences of the global $U(1)$ symmetry of QED on the correlation functions of the field. This will be the subject of next section.

[^20]
### 4.4.3 WT Identities in QED

Gauge invariance in QED requires the addition of a gauge-fixing term

$$
\begin{equation*}
\mathcal{L}_{\text {g.f. }}=-\frac{1}{2 \xi}\left(\partial_{\mu} A^{\mu}\right)^{2}, \tag{4.4.22}
\end{equation*}
$$

where $\xi$ is an arbitrary real parameter. ${ }^{7}$ The total Lagrangian density is then

$$
\begin{equation*}
\mathcal{L}\left(A_{\mu}, \psi, \bar{\psi}\right)=\mathcal{L}_{Q E D}\left(A_{\mu}, \psi, \bar{\psi}\right)+\mathcal{L}_{g . f .}\left(A_{\mu}\right) \tag{4.4.23}
\end{equation*}
$$

where $\mathcal{L}_{Q E D}$ is given in eq.(3.3.14) and the functional integration is done both on the vector field $A_{\mu}$ and the spinors $\bar{\psi}$ and $\psi$. The source term at the exponential has a density of the form

$$
\begin{equation*}
J^{\mu}(x) A_{\mu}(x)+\bar{J}(x) \psi(x)+\bar{\psi}(x) J(x), \tag{4.4.24}
\end{equation*}
$$

where $J$ and $\bar{J}$ are Grassmann-valued (and hence anticommuting) functions. If we make the infinitesimal change of variable associated to a gauge transformation,

$$
\begin{equation*}
\psi(x) \rightarrow \psi(x)+i e \epsilon(x) \psi(x), \quad \bar{\psi}(x) \rightarrow \bar{\psi}(x)-i e \epsilon(x) \bar{\psi}(x), \quad A_{\mu}(x) \rightarrow A_{\mu}(x)+\partial_{\mu} \epsilon(x) \tag{4.4.25}
\end{equation*}
$$

the measure and $\mathcal{L}_{Q E D}$ remain invariant, while $\mathcal{L}_{\text {g.f. }}$ and the source term will change:

$$
\begin{align*}
& \mathcal{L}\left(A_{\mu}, \psi, \bar{\psi}\right) \rightarrow \mathcal{L}\left(A_{\mu}, \psi, \bar{\psi}\right)-\frac{\partial^{\mu} A_{\mu}}{\xi} \square \epsilon \\
& J^{\mu} A_{\mu}+\bar{J} \psi+\bar{\psi} J \rightarrow J^{\mu} A_{\mu}+\bar{J} \psi+\bar{\psi} J+J^{\mu} \partial_{\mu} \epsilon+i e \bar{J} \psi \epsilon-i e \bar{\psi} J \epsilon . \tag{4.4.26}
\end{align*}
$$

By bringing down the $\mathcal{O}(\epsilon)$ terms from the exponential and taking a functional derivative with respect to $\epsilon(x)$, we get the identity

$$
\begin{equation*}
i \partial_{\mu} J^{\mu}(x) Z=e\langle\bar{\psi}(x)\rangle_{J} J(x)-e \bar{J}(x)\langle\psi(x)\rangle_{J}-\frac{i}{\xi} \square \partial_{\mu}\left\langle A^{\mu}(x)\right\rangle_{J} . \tag{4.4.27}
\end{equation*}
$$

We also have ${ }^{8}$

$$
\begin{equation*}
\langle\bar{\psi}(x)\rangle_{J}=i \frac{\delta Z}{\delta J(x)}, \quad\langle\psi(x)\rangle_{J}=-i \frac{\delta Z}{\delta \bar{J}(x)},\left\langle A^{\mu}(x)\right\rangle_{J}=-i \frac{\delta Z}{\delta J_{\mu}(x)} \tag{4.4.28}
\end{equation*}
$$

and thus we can rewrite eq.(4.4.27) in terms of $W=-i \log Z$ as

$$
\begin{equation*}
i \partial_{\mu} J^{\mu}(x)=-e \frac{\delta W}{\delta J(x)} J(x)-e \bar{J}(x) \frac{\delta W}{\delta \bar{J}(x)}-\frac{i}{\xi} \square \partial_{\mu} \frac{\delta W}{\delta J_{\mu}(x)} . \tag{4.4.29}
\end{equation*}
$$

[^21]We can go further and write the WT Identity (4.4.29) in terms of the 1PI generator $\Gamma$ defined as

$$
\begin{equation*}
\Gamma=W-\int d^{4} x\left(\mathcal{A}_{\mu}(x) J^{\mu}(x)+\bar{J}(x) \Psi(x)+\bar{\Psi}(x) J(x)\right) \tag{4.4.30}
\end{equation*}
$$

where

$$
\begin{equation*}
\mathcal{A}_{\mu}(x)=\frac{\delta W}{\delta J^{\mu}(x)}, \quad \Psi(x)=\frac{\delta W}{\delta \bar{J}(x)}, \quad \bar{\Psi}(x)=-\frac{\delta W}{\delta J(x)}, \tag{4.4.31}
\end{equation*}
$$

and correspondingly

$$
\begin{equation*}
J^{\mu}(x)=-\frac{\delta \Gamma}{\delta \mathcal{A}_{\mu}(x)}, \quad J(x)=-\frac{\delta \Gamma}{\delta \bar{\Psi}(x)}, \quad \bar{J}(x)=\frac{\delta \Gamma}{\delta \Psi(x)} . \tag{4.4.32}
\end{equation*}
$$

In terms of $\Gamma$, eq.(4.4.28) reads

$$
\begin{equation*}
i \partial_{\mu} \frac{\delta \Gamma}{\delta \mathcal{A}_{\mu}(x)}=e \bar{\Psi}(x) \frac{\delta \Gamma}{\delta \bar{\Psi}(x)}+e \frac{\delta \Gamma}{\delta \Psi(x)} \Psi(x)+\frac{i}{\xi} \square \partial_{\mu} \mathcal{A}^{\mu}(x) . \tag{4.4.33}
\end{equation*}
$$

By taking arbitrary functional derivatives of eqs.(4.4.27), (4.4.29) and (4.4.33) with respect to the sources we get an infinite set of WT identities between connected and 1PI amplitudes, respectively.

Let us see more closely the WT identities that will allow us to prove three assertions in section 3.3 when studying the renormalizability of this theory:

1. A relation linking the electron two-point function with the vertex, eventually proving eq.(3.3.15).
2. The transversality of the photon propagator, eq.(3.3.5).
3. Decoupling of unphysical photon polarization states, eq.(3.3.9) (and its generalization to any other scattering amplitude involving at least one photon).

Relation 1 is obtained by taking one functional derivative with respect to $\Psi$ and $\bar{\Psi}$ of eq.(4.4.33). In this way one has

$$
\begin{equation*}
i \partial_{\mu} \frac{\delta^{3} \Gamma}{\delta \Psi\left(x_{1}\right) \delta \bar{\Psi}\left(x_{2}\right) \delta \mathcal{A}_{\mu}(x)}=e \delta\left(x-x_{2}\right) \frac{\delta^{2} \Gamma}{\delta \Psi\left(x_{1}\right) \delta \bar{\Psi}(x)}-e \delta\left(x-x_{1}\right) \frac{\delta^{2} \Gamma}{\delta \Psi(x) \delta \bar{\Psi}\left(x_{2}\right)} . \tag{4.4.34}
\end{equation*}
$$

In momentum space eq.(4.4.34) becomes,

$$
\begin{equation*}
q^{\mu} \Gamma_{\mu}(p, q)=e \Gamma^{(2)}(q+p)-e \Gamma^{(2)}(p) \tag{4.4.35}
\end{equation*}
$$

where $\Gamma_{\mu}$ is precisely the function vertex defined in eq.(3.3.2) and $\Gamma^{(2)}=\not p-m+\Sigma(p)$ is the inverse of the electron propagator, with $\Sigma$ defined in eq.(3.3.3). Using eq.(4.4.35) and


Figure 4.2: Diagrammatic representation of the WT identity for QED, which generalizes eq. (4.4.35) for connected ampltudes involving additional $2 m$ fermionic and an arbitrary number $n$ of photon external lines. The extra photon line in the diagram in the l.h.s. is amputated.
the above definitions, it is straightforward to derive the identity (3.3.15). We see that a similar identity halso holds for the finite parts:

$$
\begin{equation*}
q^{\mu} \Gamma_{\mu}^{\text {finite }}(p, q)=\Sigma^{\text {finite }}(p+q)-\Sigma^{\text {finite }}(p) \tag{4.4.36}
\end{equation*}
$$

Relation 2 is obtained by setting $\Psi=\bar{\Psi}=0$ and taking one derivative with respect to $\mathcal{A}_{\mu}$ of eq.(4.4.33). This gives, in momentum space

$$
\begin{equation*}
q^{\nu} \Gamma_{\mu \nu}(q)=\frac{1}{\xi} q^{2} q_{\mu} \tag{4.4.37}
\end{equation*}
$$

where $\Gamma_{\mu \nu}$ is the inverse photon propagator. We decompose $\Gamma_{\mu \nu}(q)=\Gamma_{\mu \nu}^{(0)}(q)+\Pi_{\mu \nu}(q)$ in terms of its tree-level and quantum corrections, where $\Pi_{\mu \nu}(q)$ is the same entering eq.(3.3.5). From the classical action we readily get

$$
\begin{equation*}
\Gamma_{\mu \nu}^{(0)}(q)=\eta_{\mu \nu} q^{2}-q_{\mu} q_{\nu}+\frac{1}{\xi} q_{\mu} q_{\nu} \tag{4.4.38}
\end{equation*}
$$

We notice that $\Gamma_{\mu \nu}^{(0)}$ alone saturates the identity (4.4.37), automatically implying eq.(3.3.5) to all orders in perturbation theory. The fundamental importance of this conclusion relies on the fact that it ensures that the photon does not acquire a mass as a consequence of interactions, being protected by gauge invariance. An analogous conclusion is drawn in chapter 6 for non-abelian gauge theories, where a relation very similar to eq.(4.4.37), eq.(6.3.33), will be proved starting from the so called BRST symmetries.

Let us now consider relation 3. In order to relate the resulting WT identities directly with $S$-matrix amplitudes, that are associated to amputated connected Green functions, it is better to start from eq.(4.4.29). In absence of fermion fields, we can set $J=\bar{J}=0$ and take an arbitrary number $n$ of functional derivatives of eq.(4.4.29) with respect to $J^{\mu}$. In this way the l.h.s. and the first two terms in the r.h.s. of eq.(4.4.29) vanish and we trivially get

$$
\begin{equation*}
q^{2} q^{\mu} G_{\mu \alpha_{1} \ldots \alpha_{n}}\left(q, p_{1}, \ldots, p_{n}\right)=0 \tag{4.4.39}
\end{equation*}
$$

for the $n+1$ connected amplitudes involving photons only. The factor of $q^{2}$ in eq.(4.4.39) is removed when we amputate the amplitude, see below. In this way we get the generalization of eq.(3.3.9) for any number of external photon fields.

When external fermions are present, the analysis is slightly more complicated. If $2 m$ external fermion lines are present, we have to take $m$ functional derivatives of eq.(4.4.29) with respect to $J$ and $\bar{J}$ (in addition to the ones necessary for the photon ones as above). In this way we derive WT identities for connected correlation functions involving an arbitrary number of gauge and fermion fields. The term in the l.h.s. still vanishes, but the first two terms in the r.h.s. are non-zero now. In momentum space we have

$$
\begin{equation*}
\frac{i}{\xi} q^{2} q^{\mu} G_{\mu}^{(2 m, n+1)}\left(p_{i}, k_{i}, q_{j}\right)=e \sum_{i=1}^{m}\left(G^{(2 m, n)}\left(p_{i}+q, k_{i}, q_{j}\right)-G^{(2 m, n)}\left(p_{i}, k_{i}-q, q_{j}\right)\right) \tag{4.4.40}
\end{equation*}
$$

where $G^{(2 m, n)}$ schematically represent the connected Green functions for $2 m$ external fermions and $n$ external photons and we have made explicit in $G^{(2 m, n+1)}$ only the Lorentz index contracted with $q_{\mu}$. As we mentioned, physical $S$-matrix amplitudes are related to amputated connected Green functions. Let us start by amputating the Green function $G^{(2 m, n+1)}$ in the l.h.s. of eq.(4.4.40) of its additional photon line. One has

$$
\begin{equation*}
G_{\mu}^{(2 m, n+1)}=G_{\mu \nu}^{\gamma}(q) G_{a m p}^{\nu(2 m, n+1)} \tag{4.4.41}
\end{equation*}
$$

where $G_{\mu \nu}^{\gamma}$ is the (all-orders) photon propagator $G_{\mu \nu}^{\gamma} \Gamma^{\nu \rho}=-i \delta_{\mu}^{\rho}$ and $G_{a m p}^{\nu(2 m, n+1)}$ refers to the amputation of this photon line only. Given eq.(4.4.37), one immediately gets

$$
\begin{equation*}
q^{2} q^{\mu} G_{\mu \nu}^{\gamma}(q)=-i \xi q_{\nu} \tag{4.4.42}
\end{equation*}
$$

Plugging eq.(4.4.42) in eq.(4.4.40) we get

$$
\begin{equation*}
q^{\mu} G_{\mu, a m p}^{(2 m, n+1)}\left(p_{i}, k_{i}, q_{j}\right)=e \sum_{i=1}^{m}\left(G^{(2 m, n)}\left(p_{i}+q, k_{i}, q_{j}\right)-G^{(2 m, n)}\left(p_{i}, k_{i}-q, q_{j}\right)\right) \tag{4.4.43}
\end{equation*}
$$

This relation is diagrammatically represented in fig. 4.2, where $n$ physical photons and $2 m$ fermions are indicated by wiggly and straight lines, respectively, and $q^{\mu}$ denotes the wiggly
line of the amputated photon with polarization $\epsilon_{\mu}(q)=q_{\mu}$. The grey circle stands for all possible diagrams which can be constructed with the specified external legs. By using the LSZ reduction formula, the correlation functions (4.4.43) are directly related to the physical $S$-matrix amplitudes. The latter is in particular proportional to the total residue given by the $2 m+n$ poles in the correlation function (see eq. (2.3.10)), since we still have to amputate the additional $2 m+n$ external states. However, we now notice that on the r.h.s. of eq.(4.4.43) one fermion momentum is always shifted from its on-shell value. If $p_{i}^{2}=k_{i}^{2}=m^{2}$, generically $\left(p_{i}+q\right)^{2},\left(k_{i}-q\right)^{2} \neq m^{2}$. Hence the correlation functions on the r.h.s. can only have $2 m+n-1$ poles and hence eq. (2.3.10) necessarily vanishes. We conclude that any physical amplitude involving an arbitrary number of photon and fermion fields must vanish when any of the external photon polarizations is taken proportional to the photon momentum:

$$
\begin{equation*}
\epsilon_{\mu_{1}}\left(p_{1}\right) \ldots p_{\mu_{j}} \ldots \epsilon_{\mu_{n+1}}\left(p_{n+1}\right) \mathcal{M}^{\mu_{1} \ldots \mu_{n+1}}=0 \tag{4.4.44}
\end{equation*}
$$

## Chapter 5

## The Renormalization Group

The renormalization group is a key concept in quantum field theory. It essentially tells us that instead of describing a physical system by some constant parameters in a Lagrangian, it is more convenient to let the parameters vary and keep only some of them, depending on the energy scale at which we are looking at the system. Intuitively this is quite obvious and is at the basis of the usual reductionism used in physics. We do not need the SM Lagrangian to study the energy levels of the hydrogen atom! The latter are well described by a much simpler Schrödinger equation, which captures the effective dynamics entering at the eV scale, namely the Coulomb potential between the electron and the proton. In general, however, the microscopic short distance behaviour of a system is not completely negligible. When this is the case, if we are interested in processes occurring at some energy scale $E$, we can "integrate out", rather than simply neglect, all states with higher frequencies and retain only the effective degrees of freedom of interest. Historically, the renormalization group was developed by Gell-Mann and Low in ref.[13] as a way to improve the perturbative expansion in QED. Although, as we will see, in the context of particle physics the original approach of ref.[13] is essentially still used today, the relation to the idea of integrating out degrees of freedom and the use of effective theories was pioneered several years after by K.G. Wilson [14, 15].

### 5.1 Relevant, Marginal and Irrelevant Couplings

In this section we consider the renormalization group in the spirit of Wilson's original idea, focusing on a particular model, the $\phi^{4}$ theory in four space-time dimensions, Wick rotated in euclidean space. ${ }^{1}$ Wilson's approach has the main advantage of being conceptually very clear. It assumes the presence of a physical cut-off $\Lambda$ in the theory, above which no mode

[^22]can be excited. Imagine we are interested in processes occurring at scales of order $b \Lambda$, with $b<1$. We label all modes $\phi(k)$ as "heavy" and "light", depending on their momentum $k$. We write
\[

$$
\begin{equation*}
\phi(k)=\phi_{L}(k) \theta(b \Lambda-|k|)+\phi_{H}(k) \theta(|k|-b \Lambda), \quad|k|<\Lambda \tag{5.1.1}
\end{equation*}
$$

\]

and we correspondingly decompose the action of the $\phi^{4}$ theory as follows: ${ }^{2}$

$$
\begin{equation*}
S=S_{L}+S_{H}^{0}+S_{i n t} \tag{5.1.2}
\end{equation*}
$$

where

$$
\begin{align*}
S_{L} & =\int d^{4} x\left(\frac{1}{2}\left(\partial \phi_{L}\right)^{2}+\frac{1}{2} m^{2} \phi_{L}^{2}+\frac{\lambda}{4!} \phi_{L}^{4}\right), \\
S_{H}^{0} & =\int d^{4} x\left(\frac{1}{2}\left(\partial \phi_{H}\right)^{2}+\frac{1}{2} m^{2} \phi_{H}^{2}\right), \\
S_{\text {int }} & =\frac{\lambda}{4!} \int d^{4} x\left(\phi_{H}^{4}+4 \phi_{H} \phi_{L}^{3}+4 \phi_{H}^{3} \phi_{L}+6 \phi_{H}^{2} \phi_{L}^{2}\right) . \tag{5.1.3}
\end{align*}
$$

It is clear from eq. (5.1.1) that the quadratic terms of the form $\phi_{H} \phi_{L}$ vanish. The modes $\phi_{H}$ cannot be excited for processes at scales below or of order $b \Lambda$, so at first approximation they can be ignored, in which case we just recover the $\phi^{4}$ theory $S_{L}$ for the light mode only. Quantum mechanically, however, the modes $\phi_{H}$ contribute as virtual particles. Instead of neglecting them, we should more properly integrate them out, getting in this way an effective action for the light modes $\phi_{L}$ :

$$
\begin{equation*}
e^{-S_{e f f}\left(\phi_{L}\right)}=e^{-S_{L}\left(\phi_{L}\right)} \int \mathcal{D} \phi_{H}(k) e^{-S_{H}^{0}\left(\phi_{H}\right)-S_{i n t}\left(\phi_{L}, \phi_{H}\right)} \tag{5.1.4}
\end{equation*}
$$

The effective action $S_{\text {eff }}\left(\phi_{L}\right)$ takes into account of the effects of the heavy fields at the full quantum level. In this sense, it resembles the 1PI action $\Gamma(\phi)$ defied in section 4.1, but only for the heavy fields. This makes a crucial difference: while the functional $\Gamma(\phi)$ is in general non-local and is expanded as in eq.(4.1.11), the action $S_{\text {eff }}\left(\phi_{L}\right)$, admits an expansion in terms of a local action, though in general containing an infinite number of terms. In eq. (5.1.4) the modes $\phi_{L}$ act like external fields. It is not difficult to see that the tree-level exchange of the $\phi_{H}$ modes generates effective $\phi_{L}$ couplings of the form $\left(\lambda^{2} / m^{2}\right) \phi_{L}^{6}$ (from two $\phi_{H} \phi_{L}^{3}$ vertices) plus an infinite set of higher derivative couplings involving six $\phi_{L}$ fields. Similar considerations can be made for all other couplings. At some order in perturbation theory all possible couplings compatible with the symmetries

[^23]will be generated. The effective action $S_{\text {eff }}\left(\phi_{L}\right)$ reads as
$S_{e f f}\left(\phi_{L}\right)=\int d^{4} x\left(\frac{1}{2}(1+\delta Z)\left(\partial \phi_{L}\right)^{2}+\frac{1}{2}\left(m^{2}+\delta m^{2}\right) \phi_{L}^{2}+\frac{\lambda+\delta \lambda}{4!} \phi_{L}^{4}+\delta Z_{1} \phi_{L}^{6}+\delta Z_{2}\left(\partial \phi_{L}\right)^{4}+\ldots\right)$
where ... stands for higher dimensional operators. Recall that $\phi_{L}$ contains only momenta $k \leq b \Lambda$. The rescaled momentum $k^{\prime}=k / b$ satisfies the same constraint as the original system, $k^{\prime}<\Lambda$. Correspondingly we redefine coordinates as $x^{\prime}=x b$. We also have to redefine the field $\phi$ so that it has a canonically normalized kinetic term:
\[

$$
\begin{equation*}
\phi_{L}^{c a n}=b^{-1} \sqrt{1+\delta Z} \phi_{L} . \tag{5.1.6}
\end{equation*}
$$

\]

The action $S_{\text {eff }}$ reads now (redefining $x^{\prime} \rightarrow x$ and $\phi_{L}^{\text {can }} \rightarrow \phi_{L}$ )

$$
\begin{equation*}
S_{e f f}\left(\phi_{L}\right)=\int d^{4} x\left(\frac{1}{2}\left(\partial \phi_{L}\right)^{2}+\frac{1}{2} m^{2}(b) \phi_{L}^{2}+\frac{\lambda(b)}{4!} \phi_{L}^{4}+\delta Z_{1}(b) \phi_{L}^{6}+\delta Z_{2}(b)\left(\partial \phi_{L}\right)^{4}+\ldots\right) \tag{5.1.7}
\end{equation*}
$$

where ${ }^{3}$

$$
\begin{align*}
m^{2}(b)=\frac{1}{b^{2}} \frac{m^{2}+\delta m^{2}}{1+\delta Z}, & \lambda(b)=\frac{1}{b^{0}} \frac{\lambda+\delta \lambda}{(1+\delta Z)^{2}} \\
\delta Z_{1}(b)=b^{2} \frac{\delta Z_{1}}{(1+\delta Z)^{3}}, & \delta Z_{2}(b)=b^{4} \frac{\delta Z_{2}}{(1+\delta Z)^{2}} \tag{5.1.8}
\end{align*}
$$

The process of integrating out heavy degrees of freedom and rescale the momentum is called "renormalization group" (RG).

The action (5.1.7) is the proper action for describing processes at scales $E \leq b \Lambda$. It is straightforward to see that a coupling $c_{\mathcal{O}}$ of a generic operator $\mathcal{O}$ of dimension $\Delta$ (in mass) scales as

$$
\begin{equation*}
c_{\mathcal{O}}(b)=b^{\Delta-4} c_{\mathcal{O}} . \tag{5.1.9}
\end{equation*}
$$

At lower and lower energies (smaller and smaller $b$ ) among all infinite operators appearing in eq. (5.1.7), only the finite subset of those with $\Delta \leq 4$ do actually matter, all the others being "irrelevant". This is a very remarkable result, which dramatically simplifies the physical description of a system. We define as

- Irrelevant the operators with $\Delta>4$
- Relevant the operators with $\Delta<4$
- Marginal the operators with $\Delta=4$

[^24]Correspondingly, we define as

- Irrelevant the couplings with mass dimensions $<0$
- Relevant the couplings with mass dimensions $>0$
- Marginal the couplings with mass dimensions $=0$

Relevant couplings grow in the IR and dominate the physics. In our $\phi^{4}$ example, the only relevant coupling is the mass term, which is indeed the most important parameter governing the dynamics of a particle at low energies (in perturbation theory). When all relevant couplings vanish, the IR theory is controlled by marginal couplings only, in which case we say that the theory is (classically, see below) scale invariant, namely the form of the action does not change under the RG flow.

The fate of marginal operators under the renormalization group cannot be deduced from classical scaling and requires a quantum computation. In the $\phi^{4}$ example, the only marginal operator is $\lambda$ and its $b$-dependence is determined by that of $\delta Z$ and $\delta \lambda$. It is very useful to determine $\lambda(b)$ at the lowest non-trivial order in perturbation theory, which is one-loop. At one-loop level the exchange of $\phi_{H}$ fields modifies the $\phi_{L}^{4}$ coupling by means of two $\phi_{L}^{2} \phi_{H}^{2}$ interactions. No correction arises in $\delta Z$ at one-loop order, so that we can neglect it. By denoting the heavy fields with a wavy line, the relevant Feynman graph to consider is

$$
\begin{equation*}
\operatorname{lnc}_{p+k}^{k}+2 \text { perms. } \equiv-\delta_{\lambda}=\frac{3 \lambda^{2}}{2} \int_{b \Lambda \leq k \leq \Lambda} \frac{d^{4} k}{(2 \pi)^{4}} \frac{1}{k^{2}+m^{2}} \frac{1}{(p+k)^{2}+m^{2}} \text {, } \tag{5.1.10}
\end{equation*}
$$

where $p$ is the incoming external momentum and "perms." refer to two other diagrams obtained by permuting the external lines. When the external momentum and the mass $m$ are much smaller than $b \Lambda, \delta_{\lambda}$ is easily computed:

$$
\begin{equation*}
\delta_{\lambda}=-\frac{3 \lambda^{2}}{32 \pi^{4}} \int_{b \Lambda \leq k \leq \Lambda} \frac{d^{4} k}{k^{4}}=-\frac{3 \lambda^{2}}{16 \pi^{2}} \log \frac{1}{b} \tag{5.1.11}
\end{equation*}
$$

The effective quartic interaction between the light modes is then

$$
\begin{equation*}
\lambda(b)=\lambda+\delta_{\lambda}=\lambda+\frac{3 \lambda^{2}}{16 \pi^{2}} \log b \tag{5.1.12}
\end{equation*}
$$

As we can seen from eq. (5.1.12), $\lambda(b)$ decreases at larger distances (smaller $b$ ), so we say that the coupling $\lambda$ is marginally irrelevant, meaning in this way that it decreases in the IR, although not as quickly as a classically irrelevant operator. Similarly, we call
marginally relevant a classically marginal coupling that increases at larger distance due to quantum effects. Classically marginal couplings that remain marginal at the quantum level are called exactly marginal.

At the quantum level, as we have just seen, all the terms $\delta Z, \delta m^{2}$ etc., appearing in eqs. (5.1.7) and (5.1.8), can depend on $b$. This implies that the scaling dimension of an operator in general receives corrections with respect to its classical value determined from the powers of $b$ appearing in eq. (5.1.7). The difference between the quantum contribution to the scaling dimension of an operator from its classical value is called the anomalous dimension of the operator. We will come back to it later on in this chapter.

We started our analysis from the UV action (5.1.2) but it should be now clear that if we had started already at the UV with the most general (non-renormalizable) action of the form (5.1.5), under the renormalization group flow we would have always ended up with the usual $\phi^{4}$ theory, plus an infinite number of irrelevant operators. From the Wilsonian RG point of view, then, renormalizable theories can be seen as critical surfaces, in parameter space, where a much larger class of theories flow to. It is important to stress here that our classification of the operators in relevant, irrelevant and marginal is based on the classical dimension of fields, valid at parametrically weak coupling in perturbation theory around free field theories. At strong coupling, it might happen that, say, an operator that is classically irrelevant becomes marginal due to non-perturbative effects.

The Wilsonian picture of the RG flow is very intuitive and physical but, pragmatically speaking, it is not the best way to proceed in high energy physics. Distinguishing light and heavy modes in a single field can give rise to complicated expressions and in Minkowski space is not a Lorentz invariant notion. ${ }^{4}$ Cut-off regularization is often unavailable, like in gauge theories. Moreover, the Wilsonian RG flow requires, as a starting point, some UVregulated Lagrangian, while in high energy physics we prefer to hide our ignorance about the UV physics in the renormalization of the parameters entering into the Lagrangian. For all these reasons, in the remaining of this chapter we will change perspective and consider the RG flow from an "high energy physics" point of view. As we will see, in a perturbative context, it essentially gives us a way of improving the perturbative expansion, in the spirit of ref.[13].

[^25]
### 5.2 The Sliding Scale and the Summation of Leading Logs

The RG flow technique is very useful in standard four dimensional (weakly interacting) theories in high energy physics. In particular, it improves the perturbative expansion, allowing us to sum whole series of higher order effects. An example will clarify the problem and its resolution.

Consider the 1PI 4-point function at one-loop level in our usual $\phi^{4}$ theory in $d=4$ space-time dimensions. Using a cut-off $\Lambda$ in the momenta, we get

where $s, t$ and $u$ are the Mandelstam variabled defined in eq.(2.6.14). As we already discussed at length, we need a renormalization condition that fixes the counter-terms needed to remove the logarithmic divergence in eq. (5.2.1). Let us define the coupling constant $\lambda$ as the value of $\Gamma^{(4)}\left(p_{i}\right)$ at the (unphysical) symmetric point $s=t=u=4 m^{2} / 3$ :

$$
\begin{equation*}
i \Gamma^{(4)}\left(s=t=u=4 m^{2} / 3\right) \equiv-i \lambda \tag{5.2.2}
\end{equation*}
$$

In this way, the finite, renormalized 1PI four-point function reads

$$
\begin{equation*}
\Gamma^{(4)}(s, t, u)=-\lambda+\frac{\lambda^{2}}{32 \pi^{2}} \int_{0}^{1} d x \log \frac{m^{2}(1-4 x(1-x) / 3)}{m^{2}-s x(1-x)}+(s \rightarrow t)+(s \rightarrow u) \tag{5.2.3}
\end{equation*}
$$

At high energies, when the absolute values of the Mandelstam variables are much greater than $m$, large $\operatorname{logs} \sim \log m^{2} / E^{2}$ appear. ${ }^{5}$ If $E$ is sufficiently high, it can happen that the $\log m^{2} / E^{2}$ term is so large to compensate for the one-loop suppression given by $\sim \lambda /\left(16 \pi^{2}\right)$, breaking down the perturbative expansion. Similarly, at two-loop level, terms of the form $\lambda^{3} \log ^{2} m^{2} / E^{2}$ and $\lambda^{3} \log m^{2} / E^{2}$ appear. In general, at $l$-loop level, all terms of the form $\lambda^{l+1} \log ^{l} m^{2} / E^{2}, \lambda^{l+1} \log ^{l-1} m^{2} / E^{2}, \ldots, \lambda^{l+1} \log m^{2} / E^{2}$ can appear. It is clear that if $\lambda \ll 1$, but $\lambda \log E^{2} / m^{2} \sim 1$, the terms of the form $\lambda\left(\lambda^{l} \log ^{l} m^{2} / E^{2}\right)$ are all of the same order. Logs of this form are called leading logs (LL) for obvious reasons. The terms of the form $\lambda^{2}\left(\lambda^{l} \log ^{l} m^{2} / E^{2}\right)$ are denoted next-to-leading logs (NLL) and so on. There is

[^26]actually a simple way to avoid the explicit appearance of large logs in $\Gamma^{(4)}$. The idea is based on the fact that we can choose the renormalization condition for the coupling as we wish. If we define the coupling constant at an energy scale $\mu \sim E$, replacing eq. (5.2.2) by
\[

$$
\begin{equation*}
i \Gamma^{(4)}\left(s=t=u=-\mu^{2}\right) \equiv-i \lambda(\mu), \tag{5.2.4}
\end{equation*}
$$

\]

the finite, 1PI four-point function (5.2.3) would now read

$$
\begin{equation*}
\Gamma^{(4)}(s, t, u)=-\lambda(\mu)+\frac{\lambda^{2}(\mu)}{32 \pi^{2}} \int_{0}^{1} d x \log \frac{m^{2}+\mu^{2} x(1-x)}{m^{2}-s x(1-x)}+(s \rightarrow t)+(s \rightarrow u) \tag{5.2.5}
\end{equation*}
$$

and no large log term appears anymore. The arbitrary scale $\mu$ is denoted the sliding or renormalization scale. The coupling $\lambda(\mu)$ is determined by noting that the physics cannot depend on our arbitrary choice of scale $\mu$. We must require that

$$
\begin{equation*}
\mu \frac{d \Gamma^{(4)}}{d \mu}=\left[\mu \frac{\partial}{\partial \mu}+\beta\left(\lambda, \frac{m}{\mu}\right) \frac{\partial}{\partial \lambda}\right] \Gamma^{(4)}=0 \tag{5.2.6}
\end{equation*}
$$

where we have defined the function $\beta$ as

$$
\begin{equation*}
\beta\left(\lambda, \frac{m}{\mu}\right) \equiv \mu \frac{d \lambda}{d \mu} . \tag{5.2.7}
\end{equation*}
$$

The behaviour of $\lambda$ as a function of $\mu$, as given by the first-order differential equation (5.2.7), is called the renormalizaion group (RG) flow of $\lambda$. It is straightforward to compute $\beta$ given the explicit form (5.2.5) of $\Gamma^{(4)}$. We get

$$
\begin{equation*}
\beta\left(\lambda, \frac{m}{\mu}\right)=\frac{3 \lambda^{2}}{16 \pi^{2}} \int_{0}^{1} d x \frac{\mu^{2} x(1-x)}{m^{2}+\mu^{2} x(1-x)}+\mathcal{O}\left(\lambda^{3}\right) . \tag{5.2.8}
\end{equation*}
$$

The coupling $\lambda(\mu)$ is determined by the first-order differential equation (5.2.7). In the extreme high (UV) and low (IR) energy regimes $\mu \gg m, \mu \ll m$, eq. (5.2.8) simplifies considerably, giving ${ }^{6}$

$$
\begin{align*}
\beta_{U V} & \simeq \frac{3 \lambda^{2}}{16 \pi^{2}} \\
\beta_{I R} & \simeq \frac{\lambda^{2}}{32 \pi^{2}}\left(\frac{\mu}{m}\right)^{2} \simeq 0 \tag{5.2.9}
\end{align*}
$$

whose solutions are simply

$$
\begin{array}{llr}
\lambda_{U V}(\mu) \simeq \frac{\lambda\left(\mu_{0}\right)}{1-\frac{3 \lambda\left(\mu_{0}\right)}{16 \pi^{2}} \log \frac{\mu}{\mu_{0}}}, & \mu_{0}, \mu \gg m \\
\lambda_{I R}(\mu) & \simeq \mathrm{constant}, & \mu \ll m \tag{5.2.11}
\end{array}
$$

[^27]Strictly speaking, eqs. (5.2.10) and (5.2.11) are valid only in the asymptotic UV and IR regimes, but as a first crude approximation we can take $\mu_{0}=m$ in eq. (5.2.10) and $\lambda(\mu) \simeq \lambda(m)$ for any $\mu \leq m$ in eq. (5.2.11). We will perform a refined approximation in section 5.3, but for the moment this suffices to understand the main point of the RG analysis. Let us compare the two equivalent 1PI 4-point functions (5.2.3) and (5.2.5) in the euclidean UV point $s=t=u \equiv-E^{2} \gg m^{2}$ :

$$
\begin{align*}
\Gamma^{(4)}(E) & \simeq-\lambda-\frac{3 \lambda^{2}}{16 \pi^{2}} \log \frac{E}{m}+\mathcal{O}\left(\lambda^{2}\right), \quad(\text { no RG })  \tag{5.2.12}\\
\Gamma^{(4)}(E) & \simeq-\lambda(E)=-\frac{\lambda}{1-\frac{3 \lambda}{16 \pi^{2}} \log \frac{E}{m}}+\mathcal{O}\left(\lambda^{2}\right), \quad(\mathrm{RG}) \tag{5.2.13}
\end{align*}
$$

where $\lambda \simeq \lambda(m)$. Expanding in $\lambda$, we see that eq. (5.2.13) reproduces the one-loop result in eq. (5.2.12) but, in addition, automatically gives us all the LL logs of the form $\lambda \lambda^{l} \log ^{l} E / m$ (this will be proved in section 5.7). This is the key point of the RG evolution in perturbation theory: a powerful way to improve the perturbative expansion.

Various approximations have been made in comparing eqs. (5.2.3) and (5.2.5), such as $\lambda(m) \simeq \lambda(2 m / \sqrt{3})$ and $\mu_{0}=m$. All these approximations change $\Gamma(E)$ at $\mathcal{O}\left(\lambda^{2} \log ^{0} E / M\right)$ and are hidden in the $\mathcal{O}\left(\lambda^{2}\right)$ terms in eqs. (5.2.12) and (5.2.13). When $E \gg m$, these terms can consistently be neglected, being sub-leading with respect to the $\lambda^{2} \log E / M$ terms. They are important if we want to go beyond the LL approximation, in which case they are the first terms in the $\lambda^{2}\left(\lambda^{l-1} \log ^{l-1} E / M\right)$ (NLL) series. Resumming these logs require the knowledge of the $\lambda^{3} \log E / M$ term, i.e. a two-loop perturbative computation.

The sliding scale $\mu$ and the RG evolution of the coupling are also useful in the IR. For instance, when $m=0$, we immediately see that there is an IR singularity in the 1PI 4 -point function (5.2.3), singularity which is avoided in eq. (5.2.5), which is well defined for $m=0$.

Going back to eqs. (5.2.10) and (5.2.11), the RG evolution implies that the effective coupling constant in a QFT (i.e the one which does not give rise to large logs in amplitudes) depends on the energy scale, namely it is a "running" coupling constant. In particular eq. (5.2.10) implies that $\lambda\left(\mu_{1}\right)>\lambda\left(\mu_{0}\right)$ for $\mu_{1}>\mu_{0}$ (in agreement with the results derived using the Wilsonian RG flow) and predicts a pole at the scale

$$
\begin{equation*}
\mu_{L}=\mu_{0} e^{\frac{16 \pi^{2}}{3 \lambda\left(\mu_{0}\right)}} \tag{5.2.14}
\end{equation*}
$$

where the coupling diverges (Landau pole). ${ }^{7}$ The scale $\mu_{L}$ should not be taken too seriously, since at energies below $\mu_{L}$, when the denominator in eq. (5.2.10) starts to significantly differ from one, perturbation theory breaks down and higher loop corrections are

[^28]no longer negligible. However, eq. (5.2.14) indicates that, no matter how small $\lambda\left(\mu_{0}\right)$ is in the IR, at energies of order $\mu_{L}$ perturbation theory breaks down. On the other hand, perturbativity improves at low energies. When $m=0$, and eq. (5.2.10) is valid for any range of $\mu_{0}$ and $\mu$, we see that $\lambda(\mu) \rightarrow 0$ as $\mu \rightarrow 0$. In other words, provided the theory is in a perturbative regime at some scale $\mu_{0}$, in the far IR it asymptotes a free field theory. Theories of this sort are denoted IR free. When $m \neq 0$, the running instead essentially stops for $\mu \lesssim m$. This is in agreement with the fact that at low energies no large logs to be resummed appear, see eq. (5.2.3). This is a manifestation of a much more general principle, actually a theorem (Appelquist-Carazzone), according to which the effects of massive particles at low energies should be negligible and go to zero when $m \rightarrow \infty .^{8}$ This is a key principle in QFT (and physics in general), allowing to reliably describe physical processes at some scale without necessarily knowing the "true" (if any) physical theory underlying all processes. In this sense any QFT should always be seen as an "effective" theory. We will come back to this point in chapter 7 , where we will systematically study effective field theories.

### 5.3 Asymptotic Behaviours of $\beta$-Functions

The $\beta$-function is the crucial object to determine the evolution of a coupling constant. In general, in a theory with $n$ couplings $g_{i}$, we have to solve a set of coupled differential equations of the kind

$$
\begin{equation*}
\mu \frac{d g_{i}}{d \mu}=\beta_{i}, i=1, \ldots, n \tag{5.3.1}
\end{equation*}
$$

where the $\beta_{i}$ in eq. (5.3.1) depend on all the other couplings and masses of the theory. We can get rid of the masses by focusing only on the universal UV relevant coefficients, so that $\beta_{i}=\beta_{i}\left(g_{j}\right)$. Yet, it is not possible to describe the main properties of the solutions $g_{i}=g_{i}(\mu)$, because the system (5.3.1) is in general too complicated. We can further simplify the situation by considering a single coupling $g$. In perturbation theory, $\beta(g)$ admits an expansion as follows:

$$
\begin{equation*}
\beta(g)=\beta_{0} g^{2}+\beta_{1} g^{3}+\mathcal{O}\left(g^{4}\right) \tag{5.3.2}
\end{equation*}
$$

Note that it is always possible, by a proper coupling redefinition, to write the $\beta$-function expansion as in eq. (5.3.2). For instance, in QED $\beta(e)$ starts at cubic, rather than quadratic, order in the coupling, but it is enough to consider $\beta=\beta\left(e^{2}\right)$ rather than $\beta(e)$ to put the

[^29]

Figure 5.1: (Left panel) Schematic picture of the $\beta$-function of an IR stable fixed point (Right panel) Schematic picture of the $\beta$-function of an UV stable fixed point. In both cases we have taken $g^{*}=3 / 2$. The units in both axes are arbitrary and irrelevant.
$\beta$ in the form (5.3.2). At leading order, i.e. neglecting $\beta_{1}$, eq. (5.3.2) is solved by

$$
\begin{equation*}
g(\mu)=\frac{g_{0}}{1-g_{0} \beta_{0} \log \left(\mu / \mu_{0}\right)}, \tag{5.3.3}
\end{equation*}
$$

where $g_{0}=g\left(\mu_{0}\right)$. The fate of $g(\mu)$ is entirely governed by the sign of $\beta_{0}$. When $\beta_{0}>0$, like in the $\phi^{4}$ or QED cases, $g(\mu)$ is marginally irrelevant, it increases in the UV and at some high energy scale the theory is no longer perturbative.

When $\beta_{0}<0$, on the other hand, the opposite happens. The coupling is marginally relevant and it decreases in the UV. In the limit of infinite energy, the coupling vanishes. Theories with (all) couplings of this kind are called asymptotically free. The low energy regime of these theories (including the spectrum of particles) is not perturbatively accessible. The coupling formally diverges at the scale

$$
\begin{equation*}
\Lambda=\mu e^{\frac{1}{g(\mu) \beta_{0}}} \tag{5.3.4}
\end{equation*}
$$

This scale, which is RG invariant, is said to be a dynamically generated scale, since there is no trace of it in the classical theory. It is a purely quantum effect. The most famous theory belonging to this class of theories is QCD.

It is useful to consider two other more exotic forms of $\beta$-function, that correspond to a perturbative expansion around a non-trivial value of the coupling (see figs. 5.1). Let us assume that there exists a critical value of the coupling, $g^{*}$, such that $\beta\left(g^{*}\right)=0$. If $g$ is sufficiently close to $g^{*}$, we can expand $\beta$ as follows:

$$
\begin{equation*}
\beta(g)=\beta\left(g^{*}\right)+\beta^{*}\left(g-g^{*}\right)+\mathcal{O}\left(g-g^{*}\right)^{2}=\beta^{*}\left(g-g^{*}\right)+\mathcal{O}\left(g-g^{*}\right)^{2} . \tag{5.3.5}
\end{equation*}
$$

The solution of eq. (5.3.1), with $\beta$ as in eq. (5.3.5), is

$$
\begin{equation*}
g(\mu)=g^{*}+\left(g\left(\mu_{0}\right)-g^{*}\right)\left(\frac{\mu}{\mu_{0}}\right)^{\beta^{*}}+\mathcal{O}\left(g-g^{*}\right)^{2} \tag{5.3.6}
\end{equation*}
$$

Independently of $g\left(\mu_{0}\right)$, when $\beta^{*}<0$ the coupling approaches $g^{*}$ in the UV, while for $\beta^{*}>0$ it approaches $g^{*}$ in the IR. In the former case, $g^{*}$ is called an ultraviolet stable fixed point of the RG flow, in the latter an infrared stable fixed point of the RG flow. When $g=g^{*}$ and $\beta$ vanishes, no running occurs. In this case the theory is invariant under scaling transformations. More in general, it can be shown that it is invariant also under conformal transformations. A theory of this sort is called a Conformal Field Theory (CFT). Trivial CFT's are free theories of massless particles, for which $g=0$ and $\beta$ vanishes. Non-trivial CFT's are generally strongly coupled, because a non-trivial zero of a $\beta$-function requires a cancellation among different orders in the perturbative expansion. But this signals a break down of perturbation theory, where by definition a term of order $n+1$ should be parametrically smaller than the one of order $n$. There is however a way to possibly get weakly coupled CFTs. Suppose that in a theory the coefficient $\beta_{0}$ of the one-loop $\beta$ function is accidentally small. If so, setting to zero eq.(5.3.2) we get, aside from $g=0$, the non-trivial solution

$$
\begin{equation*}
g^{\star}=-\frac{\beta_{0}}{\beta_{1}} \tag{5.3.7}
\end{equation*}
$$

This solution (5.3.6) can be trusted only when $\beta_{0} / \beta_{1} \ll 1$, our working hypothesis, otherwise higher order terms would destabilize it and bring $g^{\star}$ at a generically strongly coupled value. The fixed point (5.3.6) is called of the Banks-Zaks kind [16]. In non-abelian gauge theories with a large gauge group and a large number of matter fields one can tune $\beta_{0}$ to be parametrically small. With a proper number of matter fields, a Banks-Zaks fixed point can be obtained also in QCD. The clearest example is the one of QCD with 16 fermions, in which case $\beta_{0}<0$ and in modulus is the smallest possible. Lattice simulations show that, contrary to "real world" QCD, at large distances this theory does not confine and approaches a fixed-point, with a $\beta$-function qualitatively as the one depicted in the left panel of fig.5.1.

We now show that the coefficients $\beta_{0}$ and $\beta_{1}$ in eq. (5.3.2) do not depend on the renormalization scheme chosen, while higher order terms are scheme dependent. The coupling constants in two different schemes, call them schemes $g$ and $\tilde{g}$, are equal at lowest order, but they start to differ at higher orders in the coupling (see eq. (5.6.12) for a concrete relation between coupling constants defined in different renormalization schemes). In general we have

$$
\begin{equation*}
\tilde{g}(g)=g+a g^{2}+\mathcal{O}\left(g^{3}\right), \tag{5.3.8}
\end{equation*}
$$

where $a$ is a constant. Their associated $\beta$-functions are related as

$$
\begin{equation*}
\tilde{\beta}(\tilde{g})=\mu \frac{d \tilde{g}}{d \mu}=\mu \frac{d g}{d \mu} \frac{d \tilde{g}}{d g}=\beta(g) \frac{d \tilde{g}}{d g} . \tag{5.3.9}
\end{equation*}
$$

Taking $\beta(g)$ as in eq. (5.3.2) we have

$$
\begin{align*}
\tilde{\beta}(\tilde{g}) & =\left(\beta_{0} g^{2}+\beta_{1} g^{3}+\mathcal{O}\left(g^{4}\right)\right)\left(1+2 a g+\mathcal{O}\left(g^{2}\right)\right) \\
& =\left(\beta_{0}\left(\tilde{g}-a \tilde{g}^{2}\right)^{2}+\beta_{1} \tilde{g}^{3}+\mathcal{O}\left(\tilde{g}^{4}\right)\right)\left(1+2 a \tilde{g}+\mathcal{O}\left(\tilde{g}^{2}\right)\right) \\
& =\beta_{0} \tilde{g}^{2}+\beta_{1} \tilde{g}^{3}+\mathcal{O}\left(\tilde{g}^{4}\right) \tag{5.3.10}
\end{align*}
$$

where the $\mathcal{O}\left(\tilde{g}^{4}\right)$ terms are different from the $\mathcal{O}\left(g^{4}\right)$ terms in $\beta(g)$. Hence we have proved the scheme independence of the coefficients $\beta_{0}$ and $\beta_{1}$.

We will see in the following that even at one- (and two-) loop level the detailed form of a $\beta$-function in a theory depends on the renormalization scheme chosen, although its UV behaviour is universal and governed by the coefficients $\beta_{0}$ and $\beta_{1}$ above.

The generalization of eqs. (5.3.2) and (5.3.8) for multiple couplings are

$$
\begin{align*}
& \beta_{i}\left(g_{j}\right)=\beta_{i j k}^{(0)} g_{j} g_{k}+\beta_{i j k l}^{(1)} g_{j} g_{k} g_{l}+\mathcal{O}\left(g_{i}^{4}\right)  \tag{5.3.11}\\
& \tilde{g}_{i}\left(g_{j}\right)=g_{i}+a_{i j k} g_{j} g_{k}+\mathcal{O}\left(g^{3}\right)
\end{align*}
$$

Proceeding as before, we get

$$
\begin{align*}
& \tilde{\beta}_{i j k}^{(0)}=\beta_{i j k}^{(0)} \\
& \tilde{\beta}_{i j k l}^{(1)}=\beta_{i j k l}^{(1)}+\frac{2}{3}\left(a_{i s l} \beta_{j k s}^{(0)}-a_{s k l} \beta_{i j s}^{(0)}+2 \text { perms. in }(j, k, l)\right) \tag{5.3.12}
\end{align*}
$$

The two-loop coefficients of the $\beta$-function for multiple couplings are in general schemedependent.

### 5.4 The Callan-Symanzik RG Equations

We have seen that at the quantum level we are necessarily led to introduce the sliding scale $\mu$. It is often useful to define (renormalize) at the same scale $\mu$ not only the coupling constant but also the fields themselves. For instance, as we will see in subsection 5.5, in renormalization schemes that make use of dimensional regularization, the scale $\mu$ shows up in a different way and affects all Green functions. In particular, the mass renormalized in those schemes does not coincide with the physical mass and the wave function renormalization factor $Z$ is no longer the one appearing in the LSZ reduction formulae. Changing the definition of $Z$, now a function of $\mu, Z=Z(\mu)$, will however make external legs contribute to physical processes, because the $Z$ entering the LSZ reduction formulae is the physical $\mu$-independent one. This is quite clear from eq.(2.3.10). Aside from irrelevant factors, the S-matrix reads

$$
\begin{equation*}
S_{m \rightarrow n}^{C} \sim \frac{1}{Z^{\frac{n+m}{2}}} G_{B}^{(n+m)}\left(p_{i},-q_{j}\right)=\frac{1}{Z^{\frac{n+m}{2}}} Z(\mu)^{\frac{n+m}{2}} G^{(n+m)}\left(p_{i},-q_{j}, \mu\right) \tag{5.4.1}
\end{equation*}
$$

where $G_{B}^{(n+m)}$ and $G^{(n+m)}$ are the bare and renormalized connected Green functions, respectively. Since $S$-matrix elements do not depend on $\mu$, we see that the bare, rather than the renormalized, Green functions are those that are independent of $\mu$. A similar reasoning applies of course also to the 1PI amplitudes. Due to the amputation of the external legs, one has the opposite power of $Z$ 's relating bare and renormalized 1PI amplitudes. Taking again the $\lambda \phi^{4}$ theory as our working example, we have

$$
\begin{equation*}
\Gamma_{B}^{(n)}=Z^{-\frac{n}{2}}(\mu) \Gamma(\mu), \tag{5.4.2}
\end{equation*}
$$

and hence

$$
\begin{equation*}
\mu \frac{d\left(Z^{-\frac{n}{2}} \Gamma^{(n)}\right)}{d \mu}=Z^{-\frac{n}{2}}\left[\mu \frac{\partial}{\partial \mu}+\beta\left(\lambda, \frac{m}{\mu}\right) \frac{\partial}{\partial \lambda}-n \gamma\left(\lambda, \frac{m}{\mu}\right)\right] \Gamma^{(n)}=0 \tag{5.4.3}
\end{equation*}
$$

where

$$
\begin{equation*}
\gamma\left(\lambda, \frac{m}{\mu}\right) \equiv \frac{1}{2} \mu \frac{d \log Z}{d \mu} \tag{5.4.4}
\end{equation*}
$$

is the anomalous dimension of the field $\phi$. Eqs.(5.4.3) are called the Callan-Symanzik (CS) equations. ${ }^{9}$ For simplicity, in eq.(5.4.3) we have assumed that $m$ is the physical $\mu$-independent mass, otherwise we would also get a term proportional to the $\beta$-function of the mass. We will consider the RG flow of the mass term, and more in general of dimensionful couplings, in section 5.9.

The CS equations (5.4.3) can be solved as follows in the UV regime where we can neglect the $m / \mu$ dependence of $\beta$ and $\gamma$. As a first step, we can get rid of the last term in eq. (5.4.3) by defining a new Green function $\hat{\Gamma}^{(n)}$ :
such that

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\beta(\lambda) \frac{\partial}{\partial \lambda}\right) \hat{\Gamma}^{(n)}\left(p_{i}, \lambda, m, \mu\right)=0 . \tag{5.4.6}
\end{equation*}
$$

The above equation is solved by introducing an auxiliary variable $t$ and $t$-dependent functions $\mu(t)$ and $\lambda(t)$ with

$$
\begin{equation*}
\mu(t)=e^{t} \mu, \quad \lambda(0)=\lambda \tag{5.4.7}
\end{equation*}
$$

We then demand that $\lambda(t)$ is such that

$$
\begin{equation*}
\frac{d}{d t} \hat{\Gamma}^{(n)}\left(p_{i}, \lambda(t), m, \mu(t)\right)=0 \tag{5.4.8}
\end{equation*}
$$

[^30]When eq. (5.4.8) is satisfied, $\hat{\Gamma}^{(n)}\left(p_{i}, \lambda(t), m, \mu(t)\right)$ is independent of $t$. Evaluating it at $t=0$, we see that it coincides with our original Green function $\hat{\Gamma}^{(n)}\left(p_{i}, \lambda, m, \mu\right)$. Using the chain rule of derivatives, eq. (5.4.8) equals

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\frac{d \lambda}{d t} \frac{\partial}{\partial \lambda}\right) \hat{\Gamma}^{(n)}\left(p_{i}, \lambda(t), m, \mu(t)\right)=0 \tag{5.4.9}
\end{equation*}
$$

Comparing eqs. (5.4.9) and (5.4.6), we see that the following equation should hold:

$$
\begin{equation*}
\frac{d \lambda(t)}{d t}=\beta(\lambda(t)) \tag{5.4.10}
\end{equation*}
$$

For any $t$, provided that $\lambda(t)$ satisfies eq. (5.4.10), we have

$$
\begin{equation*}
\hat{\Gamma}^{(n)}\left(p_{i}, \lambda, m, \mu\right)=\hat{\Gamma}^{(n)}\left(p_{i}, \lambda(t), m, e^{t} \mu\right) \tag{5.4.11}
\end{equation*}
$$

In terms of the original Green functions $\Gamma^{(n)}$ and changing variable $d \lambda \rightarrow d t$ inside the integral appearing in eq.(5.4.5), eq.(5.4.11) reads

$$
\begin{equation*}
e^{-n \int_{t_{0}}^{0} \gamma\left(t^{\prime}\right) d t^{\prime}} \Gamma^{(n)}\left(p_{i}, \lambda, m, \mu\right)=e^{-n \int_{t_{0}}^{t} \gamma\left(t^{\prime}\right) d t^{\prime}} \Gamma^{(n)}\left(p_{i}, \lambda(t), m, e^{t} \mu\right) \tag{5.4.12}
\end{equation*}
$$

Rescaling $p \rightarrow e^{t} p$, eq.(5.4.12) is rewritten as

$$
\begin{equation*}
\Gamma^{(n)}\left(e^{t} p_{i}, \lambda, m, \mu\right)=e^{-n \int_{0}^{t} \gamma\left(t^{\prime}\right) d t^{\prime}} \Gamma^{(n)}\left(e^{t} p_{i}, \lambda(t), m, e^{t} \mu\right) \tag{5.4.13}
\end{equation*}
$$

Recall that the classical dimension of $\Gamma^{(n)}$ is $4-n$, so that

$$
\begin{equation*}
\Gamma^{(n)}\left(e^{t} p_{i}, \lambda(t), m, e^{t} \mu\right)=e^{t(4-n)} \Gamma^{(n)}\left(p_{i}, \lambda(t), e^{-t} m, \mu\right) \tag{5.4.14}
\end{equation*}
$$

Combining eqs.(5.4.13) and (5.4.14) we finally get

$$
\begin{equation*}
\Gamma^{(n)}\left(e^{t} p_{i}, \lambda, m, \mu\right)=e^{t(4-n)-n \int_{0}^{t} \gamma\left(t^{\prime}\right) d t^{\prime}} \Gamma^{(n)}\left(p_{i}, \lambda(t), e^{-t} m, \mu\right) . \tag{5.4.15}
\end{equation*}
$$

Equation (5.4.15) tells us that a Green function at some energy scale $e^{t} p$, modulo an overall factor, equals the same Green function evaluated at the energy scale $p$, provided we replace the coupling $\lambda$ with its running counterpart $\lambda(t)$, solution of eq. (5.4.10). In the high energy regime $t \rightarrow \infty$, mass corrections in the Green functions are negligible, in agreement with our expectations for an (IR) relevant coupling. The anomalous dimension $\gamma$ affects the effective scaling dimension of the Green function, as expected.

The UV behaviour of $\Gamma^{(n)}$ is particularly simple for theories with an UV stable fixed point $\lambda^{*}$. In this case, $\int \gamma\left[\lambda\left(t^{\prime}\right)\right] d t^{\prime} \simeq \int \gamma\left(\lambda^{*}\right) d t^{\prime}=t \gamma^{*}$, with $\gamma^{*} \equiv \gamma\left(\lambda^{*}\right)$. The UV behaviour of $\Gamma^{(n)}$ reduces then to

$$
\begin{equation*}
\Gamma^{(n)}\left(e^{t} p_{i}, \lambda, m, \mu\right) \simeq e^{4 t-n t\left(1+\gamma^{*}\right)} \Gamma^{(n)}\left(p_{i}, \lambda^{*}, 0, \mu\right) . \tag{5.4.16}
\end{equation*}
$$

Another interesting situation arises for UV free theories (like QCD) where the (one-loop) running of the coupling is given by eq. (5.3.3) with $\beta_{0}<0$. Assuming that $\gamma(t)=\gamma_{0} g$, we have $\int \gamma\left[g\left(t^{\prime}\right)\right] d t^{\prime}=\int \gamma(g) / \beta\left(g^{\prime}\right) d g^{\prime}=\gamma_{0} / \beta_{0} \log g(t) / g$ and thus for large $t$

$$
\begin{equation*}
\Gamma^{(n)}\left(e^{t} p_{i}, g, m, \mu\right) \simeq e^{t(4-n)}\left(\frac{g}{g(t)}\right)^{\frac{n \gamma_{0}}{\beta_{0}}} \Gamma^{(n)}\left(p_{i}, 0, \mu\right) \tag{5.4.17}
\end{equation*}
$$

### 5.5 Minimal Subtraction

The sliding scale $\mu$ does not necessarily appear as the momenta where we renormalize the Green functions. In particular, in DR , it arises due to the departure from $d=4$ space-time dimensions. Let $g_{B}(d)$ be a dimensionless coupling constant in 4 space-time dimensions. In $d$ dimensions the coupling will acquire a mass dimension

$$
\begin{equation*}
\Delta(d)=(4-d) \rho, \tag{5.5.1}
\end{equation*}
$$

where $\rho$ is a coupling dependent coefficient. Since divergences appear as poles in $(d-4)$, a dimensionless renormalized coupling constant can be defined as

$$
\begin{equation*}
g_{B}(d) \mu^{-\Delta(d)} \equiv g(\mu, d)+\sum_{n=1}^{\infty}(4-d)^{-n} b_{n}(g(\mu, d)) . \tag{5.5.2}
\end{equation*}
$$

Let us now take a derivative with respect to $\mu d / d \mu$ of eq. (5.5.2). We get

$$
\begin{equation*}
-(4-d) \rho\left(g+\sum_{n=1}^{\infty}(4-d)^{-n} b_{n}(g)\right)=\beta_{d}(g)+\sum_{n=1}^{\infty}(4-d)^{-n} \beta_{d} \frac{\partial b_{n}}{\partial g} \tag{5.5.3}
\end{equation*}
$$

where for simplicity $g=g(\mu, d)$. Eq. (5.5.3) should apply to any $d$, and hence we get an independent relation for any power of $d-4$. At $\mathcal{O}(4-d)$ we have

$$
\begin{equation*}
\beta_{d}(g)=-(4-d) \rho g+\beta(g), \tag{5.5.4}
\end{equation*}
$$

where $\beta(g)$ does not vanish for $d \rightarrow 4$ and is the actual four-dimensional $\beta$-function. At $\mathcal{O}(d-4)^{0}$ we have

$$
\begin{equation*}
\beta(g)=-\rho b_{1}+\rho g \frac{\partial b_{1}}{\partial g} \tag{5.5.5}
\end{equation*}
$$

We see that the $\beta(g)$ is determined by looking at the simple poles in eq. (5.5.2). Such a scheme is denoted Minimal Subtraction (MS). Since $d-4$ poles arise typically in the combination

$$
\begin{equation*}
\Gamma\left(\frac{4-d}{2}\right)(4 \pi)^{\frac{4-d}{2}} \rightarrow_{d \rightarrow 4} \frac{2}{4-d}-\gamma_{E}+\log 4 \pi \tag{5.5.6}
\end{equation*}
$$

where $\gamma_{E} \simeq 0.577$ is the Euler-Mascheroni constant, sometimes it is convenient to subtract, together with the pole $1 /(d-4)$, the finite pieces given by $\gamma_{E} / 2-1 / 2 \log 4 \pi$. Such scheme is called modified Minimal Subtraction and is denoted as $\overline{\mathrm{MS}}$.

In contrast to the definition of $\beta$ given in eq. (5.2.8), the MS or $\overline{\mathrm{MS}} \beta$-functions do not depend on masses but only on coupling constants. A similar results applies to the field anomalous dimensions. For this reason, such schemes are called mass-independent. We will explain in next section why and how these two results are not in contradiction with each other.

### 5.6 Scheme Dependence

The detailed form of the $\beta$-function of couplings and anomalous dimensions $\gamma$ of fields has not per se an intrinsic physical meaning, since it depends on the renormalization scheme chosen. Only physical quantities are scheme-independent, and hence if distinct schemes give different expressions for physical amplitudes, necessarily the evolution of the couplings should compensate for the difference. For instance, as we have just seen in section 5.5, in MS or $\overline{\text { MS }}$ schemes, $\beta$ and $\gamma$ are mass-independent, as opposed to momentum subtraction used in section 5.2 (hereafter denoted by MOM) and more in general to all schemes where $\beta$ and $\gamma$ depend on the masses of the particles. Is it then meaningful to talk about running coupling? If so, how do we determine the "correct" running? In order to answer these questions, it is useful to focus on a concrete example and work it out in some detail. As usual, we take the $\phi^{4}$ theory and compare the 1PI 4-point function (5.2.5), evaluated using MOM, with the expression one obtains in $\overline{\mathrm{MS}}$. By denoting $d=4-\epsilon$, and recalling eqs. (4.2.3) and (5.5.2), one has

$$
\begin{equation*}
\lambda_{B}(d) \mu^{-\epsilon}=\lambda+\left(\frac{Z_{\lambda}}{Z^{2}}-1\right) \lambda \equiv \lambda+\delta_{\lambda} \tag{5.6.1}
\end{equation*}
$$

We compute

$$
\begin{align*}
& =\frac{\left(-i \lambda \mu^{\epsilon}\right)^{2}}{2} \int \frac{d^{d} k}{(2 \pi)^{d}} \int_{0}^{1} d x \frac{-i}{\left[k_{E}^{2}+m^{2}-s x(1-x)\right]^{2}} \\
& =\frac{i \lambda^{2} \mu^{2 \epsilon}}{2(4 \pi)^{d / 2}} \Gamma(\epsilon / 2) \int_{0}^{1} \frac{d x}{\left[m^{2}-s x(1-x)\right]^{\epsilon / 2}} \tag{5.6.2}
\end{align*}
$$

According to eq. (5.5.6), the $\overline{\mathrm{MS}}$ scheme is defined by adding the counter-term

$$
\begin{equation*}
\delta_{\lambda}=\frac{3 \lambda^{2}}{32 \pi^{2}}\left(\frac{2}{\epsilon}-\gamma_{E}+\log 4 \pi\right) . \tag{5.6.3}
\end{equation*}
$$

The counterterm above is related to the coefficient $b_{1}$ in eq.(5.5.2), $\delta_{\lambda}=b_{1} / \epsilon$. Summing over the $t$ and $u$ channel contributions, we get the finite, renormalized 4-point function

$$
\begin{equation*}
\Gamma \frac{(4)}{\mathrm{MS}}(s, t, u)=-\lambda(\mu)+\frac{\lambda^{2}(\mu)}{32 \pi^{2}} \int_{0}^{1} d x \log \frac{\mu^{2}}{m^{2}-s x(1-x)}+(s \rightarrow t)+(s \rightarrow u) \tag{5.6.4}
\end{equation*}
$$

Notice how the numerators of the log term in eqs. (5.2.5) and (5.6.4) differ, the latter being simply $\mu^{2}$. Demanding the $\mu$-independence of the amplitude, as in section 4.1, we get

$$
\begin{equation*}
\beta_{\overline{\mathrm{MS}}}=\frac{3 \lambda^{2}}{16 \pi^{2}} \tag{5.6.5}
\end{equation*}
$$

for any value of $\mu$, independently of $m$. Of course, the same result would have been obtained directly using eq. (5.5.5) with $\rho=1$ and $b_{1}$ given by eq. (5.6.3). Thus, $\beta_{\overline{\mathrm{MS}}}$ in eq. (5.6.5) differs from $\beta_{\text {MOM }}$ in eq. (5.2.8), although their UV behaviour is identical and given by eq. (5.2.10). This is a general result: the detailed form of $\beta$ is scheme-dependent, its UV behaviour, given by mass-independent coefficients, is scheme-independent, as we have shown in section 5.3. As far as the LL summation is concerned, both renormalization schemes are valid. At low energies, however, the two schemes lead to different behaviours: in MOM, $\lambda$ essentially stops running below the scale $m$, in $\overline{\mathrm{MS}}$ the running never stops and would lead to a free theory when $\mu \rightarrow 0$ ! In the IR the physical picture is best given by MOM. This can be seen by noticing that eq. (5.2.5) is regular when $\mu \rightarrow 0$, while eq. (5.6.4) seems IR divergent. This "fake" IR divergence and the corresponding large logs associated to it, are easily evaded by noticing that at arbitrarily low energies $s, t, u \ll m^{2}$, the large logs are avoided by taking $\mu^{2} \simeq m^{2}$. In other words, in the $\overline{\mathrm{MS}}$, the necessity of avoiding spurious large logs forces us to never use the IR evolution of the coupling, being $\mu \simeq m$ the correct sliding scale in the IR. We thus conclude that the "correct" IR running behaviour is the one given by the MOM scheme. However, provided that one keeps in mind that the IR running in mass-independent schemes is fake, the latter can reliably be used. The best way to automatically get rid of this non-decoupling of heavy particles in mass-independent schemes is provided by using them in an effective field theory approach (more on effective field theories in chapter 7), where one integrates out the heavy particle so that, for $\mu \ll m$, the heavy state is no longer in the spectrum and does not contribute to the running anymore.

Amplitudes are easier to compute in mass-independent rather than in mass-dependent schemes. However, physical couplings are typically defined by processes occurring at some energy scale and are directly related to the more physical mass-dependent schemes. It is important to understand how to match couplings in mass-independent schemes with the physical couplings. Again, this is best illustrated with the specific instance of the $\phi^{4}$ theory. First of all, let us find an approximation to the RG evolution of $\lambda$ in MOM which is more refined than eqs. (5.2.10) and (5.2.11). We proceed as follows. The solution of eq. (5.2.7), for $\mu \gg m$, is given by

$$
\begin{equation*}
\lambda^{-1}(\mu)=-\frac{3}{16 \pi^{2}} \log \mu+c \tag{5.6.6}
\end{equation*}
$$

where $c$ is an integration constant. This is fixed by matching eq. (5.6.6) with the exact flow implied by eq. (5.2.7):

$$
\begin{equation*}
\lambda(\mu)=\lambda\left(\mu_{0}\right)+\int_{\mu_{0}}^{\mu} \frac{\beta}{\mu^{\prime}} d \mu^{\prime} \tag{5.6.7}
\end{equation*}
$$

Let us take $\mu_{0}=0, \lambda(0) \equiv \lambda$ and $\mu \gtrsim m$, i.e. slightly larger than $\mu$, so that we can still neglect $m$ with respect to $\mu$, but small enough that no large $\log (\mu / m)$ term appears. In the range $\left[\mu_{0}, \mu\right], \lambda(\mu) \simeq \lambda$, and eq. (5.6.7) is well approximated by

$$
\begin{equation*}
\lambda(\mu) \simeq \lambda+\frac{3 \lambda^{2}}{32 \pi^{2}} \int_{0}^{1} d x \log \frac{\mu^{2} x(1-x)}{m^{2}}=\lambda+\frac{3 \lambda^{2}}{16 \pi^{2}}\left(\log \frac{\mu}{m}-1\right) \tag{5.6.8}
\end{equation*}
$$

from which $c=\lambda^{-1}+3 /\left(16 \pi^{2}\right)(1+\log m)$ and hence

$$
\begin{equation*}
\lambda^{-1}(\mu)=\lambda^{-1}-\frac{3}{16 \pi^{2}}\left(\log \frac{\mu}{m}-1\right) \theta(\mu-m) \quad(\mathrm{MOM}) \tag{5.6.9}
\end{equation*}
$$

Notice that $\lambda(\mu)$ in eq. (5.6.9) is discontinuous in $m$ :

$$
\begin{equation*}
\lim _{\mu \rightarrow m^{-}} \lambda^{-1}(\mu)=\lambda^{-1} \neq \lim _{\mu \rightarrow m^{+}} \lambda^{-1}(\mu)=\lambda^{-1}+3 /\left(16 \pi^{2}\right) \tag{5.6.10}
\end{equation*}
$$

This discontinuity essentially takes into account the mass term disturbance to the UV running for $\mu \gtrsim m$. The constant $3 /\left(16 \pi^{2}\right)$ is often called mass threshold effect. It can be verified that eq. (5.6.9) is an excellent approximation to the exact one-loop running given by eq. (5.2.7) far away from the threshold region $\mu \simeq m$. The running coupling in the $\overline{\mathrm{MS}}$ scheme is simply given by

$$
\begin{equation*}
\lambda^{-1}(\mu)=\lambda^{-1}-\frac{3}{16 \pi^{2}} \log \frac{\mu}{m} \theta(\mu-m) \quad(\overline{\mathrm{MS}}) \tag{5.6.11}
\end{equation*}
$$

where the step function $\theta(\mu-m)$ is put by hand, for the reasons explained before. Comparing eqs. (5.6.9) with (5.6.11), we get, for $\mu>m$,

$$
\begin{equation*}
\lambda_{\mathrm{MOM}}^{-1}(\mu)=\lambda_{\frac{-1}{\mathrm{MS}}}^{-1}(\mu)+\frac{3}{16 \pi^{2}} \tag{5.6.12}
\end{equation*}
$$

We can finally answer the previous questions: is it meaningful to talk about running coupling? If so, how do we determine the "correct" running? There is no notion of "correct" coupling. The running given by any sensible scheme is meaningful, provided we consistently associate it to expressions computed in that scheme. The simplest massindependent schemes and their associated simple running, can reliably be used in the UV, and then matched, by means of formulae analogous to eq. (5.6.12), to the physically defined coupling constants. Unless differently specified, in the rest of these lecture notes we will adopt the $\overline{\mathrm{MS}}$ renormalization scheme.

### 5.7 Leading Logs and Callan-Symanzik Equations

In section 5.2 we have seen that the $\beta$-function allows us to improve the perturbative expansion by resumming large logs. However, it was not clear from our derivation that the resummed expression (5.2.13) captures exactly all higher-loops leading logs. The Callan-Symanzik equations allow us to fill this gap. Although this result is effectively encoded in eq.(5.4.15), it is more transparent to work out the explicit form of the solution (5.4.15) in perturbation theory, which is what we will do in this section. As usual, for concreteness we consider $\Gamma^{(4)}$ in the $\lambda \phi^{4}$ theory, though the derivation is more general. At high energies, neglecting masses and taking $s=t=u=-E^{2} \gg m^{2}, \Gamma^{(4)}(E)$ is a function of $\lambda$ and $\mu / E$. In perturbation theory we can then write

$$
\begin{equation*}
\Gamma^{(4)}(E)=-\lambda \sum_{l=0}^{\infty} \lambda^{l} c_{l}(\mu / E) \tag{5.7.1}
\end{equation*}
$$

where $c_{l}$ are functions to be determined, and $c_{0}=1$. This expression should satisfy the Callan-Symanzik equation

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\beta \frac{\partial}{\partial \lambda}-4 \gamma\right) \Gamma^{(4)}(E)=0 \tag{5.7.2}
\end{equation*}
$$

In perturbation theory, $\beta=\beta_{0} \lambda^{2}+\beta_{1} \lambda^{3}+\mathcal{O}\left(\lambda^{4}\right), \gamma=\gamma_{0} \lambda+\gamma_{1} \lambda^{2}+\mathcal{O}\left(\lambda^{3}\right)$ and eq.(5.7.2) should be satisfied order by order. At the first non-trivial order $\mathcal{O}\left(\lambda^{2}\right)$ we have, recalling that $\gamma_{0}=0$,

$$
\begin{equation*}
\mu \partial_{\mu} c_{1}=-\beta_{0} \longrightarrow c_{1}=-\beta_{0} \log \frac{\mu}{E}+C_{1} \tag{5.7.3}
\end{equation*}
$$

where $C_{1}$ is an integration constant. At $\mathcal{O}\left(\lambda^{3}\right)$ we get

$$
\begin{equation*}
\mu \partial_{\mu} c_{2}=-2 \beta_{0} c_{1}-\beta_{1}+4 \gamma_{1}=0 \longrightarrow c_{2}=\beta_{0}^{2} \log ^{2} \frac{\mu}{E}+\left(4 \gamma_{1}-\beta_{1}-2 \beta_{0} C_{1}\right) \log \frac{\mu}{E}+C_{2} \tag{5.7.4}
\end{equation*}
$$

Proceeding in this way allows us to establish an important result. At loop order $l$ the Green function is a polynomial of degree $l$ in $\log \mu / E$. At high energies the most relevant terms at loop $l$ are the $\log ^{l}$ terms (the leading logs). These are entirely determined in terms of $\beta_{0}$ :

$$
\begin{equation*}
c_{l}=\left(-\beta_{0}\right)^{l} \log ^{l} \frac{\mu}{E}+\mathcal{O}\left(\log ^{l-1} \frac{\mu}{E}\right) . \tag{5.7.5}
\end{equation*}
$$

Resumming these terms give

$$
\begin{equation*}
\Gamma_{\mathrm{LL}}^{(4)}(E)=-\lambda \sum_{l=0}^{\infty}\left(-\beta_{0} \lambda\right)^{l} \log ^{l}(\mu / E)=\frac{-\lambda(\mu)}{1-\beta_{0} \lambda \log \frac{E}{\mu}}, \tag{5.7.6}
\end{equation*}
$$

If we take $\mu=m$ in eq.(5.7.6) and recall that $\beta_{0}=3 /\left(16 \pi^{2}\right)$ we precisely recover eq.(5.2.13). This proves that there are no extra leading logs contributions that are not
captured by the $\beta$-function. More in general, the leading log contributions of a correlation function also depends on the anomalous dimension coefficient $\gamma_{0}$.

This analysis can be repeated to sub-leading orders (next-to leading logs, etc.). For example, by knowing the two-loop coefficients $\beta_{1}, \gamma_{1}$ and the integration constant $C_{1}$ allows us to determine the next-to-leading logs, such as the log term in eq.(5.7.4), at any loop order. More in general, the series of logs of the form $\log ^{l-n}$ at loop order $l$ can be determined from the knowledge of $\beta_{0}, \ldots, \beta_{n}, \gamma_{0}, \ldots, \gamma_{n}, C_{1}, \ldots, C_{n}$.

## 5.8 "Irrelevant" RG Flow of Dimensionful Couplings

Depending on the energy scale, relevant and irrelevant operators parametrically either dominate the physics or are negligible. For instance, in the IR relevant operators, such as masses, are the dominant effect, while in the UV irrelevant operators make a theory ill-defined. We focus here on a perturbative situation in which the (ir)relevant operators can be seen as a small deformation in the theory. For instance, the effect of a mass term in the UV or the insertion of an irrelevant operator in the IR.

The concept of sliding scale and running coupling can be extended to relevant or irrelevant couplings, but care should be paid in this case to the scheme dependence of the results. Let us first consider irrelevant couplings, assuming that no relevant couplings are present or that their effect is negligible. In general, in presence of marginal and irrelevant couplings, the $\beta$-function coefficients are scheme-dependent even in the limit where mass effects are neglected. If we denote by $g_{i}$ and $\widetilde{g}_{i}$ two coupling constants with classical mass dimension $\Delta_{i}$ in two different schemes, we have

$$
\begin{equation*}
\widetilde{g}_{i}=g_{i}+a_{i j k}(\mu) g_{j} g_{k}+\mathcal{O}\left(g^{3}\right), \tag{5.8.1}
\end{equation*}
$$

where by dimensional analysis

$$
\begin{equation*}
a_{i j k}(\mu)=\mu^{\Delta_{i}-\Delta_{j}-\Delta_{k}} c_{i j k}, \tag{5.8.2}
\end{equation*}
$$

and $c_{i j k}$ constant coefficients. The perturbative expansion of the $\beta$-functions for the $g_{i}$ 's read

$$
\begin{equation*}
\beta_{i}=\mu \frac{d g_{i}}{d \mu}=\beta_{0}^{i j k}(\mu) g_{j} g_{k}+\mathcal{O}\left(g^{3}\right) \tag{5.8.3}
\end{equation*}
$$

where

$$
\begin{equation*}
\beta_{0}^{i j k}(\mu)=\mu^{\Delta_{i}-\Delta_{j}-\Delta_{k}} b_{0}^{i j k} \tag{5.8.4}
\end{equation*}
$$

On the other hand, in the other scheme we have

$$
\begin{equation*}
\widetilde{\beta}_{i}=\mu \frac{d \widetilde{g}_{i}}{d \mu}=\widetilde{\beta}_{0}^{i j k}(\mu) \widetilde{g}_{j} \widetilde{g}_{k}+\mathcal{O}\left(\widetilde{g}^{3}\right) \tag{5.8.5}
\end{equation*}
$$

where

$$
\begin{equation*}
\widetilde{\beta}_{0}^{i j k}(\mu)=\mu^{\Delta_{i}-\Delta_{j}-\Delta_{k}} \widetilde{b}_{0}^{i j k}, \quad \widetilde{b}_{0}^{i j k}=b_{0}^{i j k}+c_{i j k}\left(\Delta_{i}-\Delta_{j}-\Delta_{k}\right) . \tag{5.8.6}
\end{equation*}
$$

Universal coefficients arise only when $\Delta_{i}-\Delta_{j}-\Delta_{k}=0$. Renormalization schemes where classical dimensional analysis is preserved at the quantum level (i.e. where no powers of $\mu$ can be generated from quantum corrections), like MS or $\overline{\mathrm{MS}}$ in DR, give automatically $b_{0}^{i j k}=0$ when $\Delta_{i}-\Delta_{j}-\Delta_{k} \neq 0$, and only keep the scheme-independent coefficients. This is related to the property of dimensional regularization of setting to zero all powerlike divergences, keeping only the logarithmic ones. Logarithmic divergences are special, because they are the only ones not saturated by UV physics, and sample uniformly all energy scales, up to the IR. Since the IR physics should be insensitive to the details of the different renormalization schemes, it follows that the associated $\beta$-function coefficients should be scheme-independent.

For illustration and in order to be concrete, consider again the $\lambda \phi^{4}$ theory, but this time in five space-time dimensions, and let us compute the one-loop RG evolution of $\lambda$. In five dimensions the $\phi^{4}$ theory is non-renormalizable and $\lambda$ is an irrelevant coupling of mass dimension -1 . In presence of irrelevant couplings, dimensional regularization can no longer be considered a mass-independent scheme, since by dimensional analysis the $\beta$-functions can have a dependence on masses. If we neglect mass terms, however, dimensional analysis dictates that the irrelevant coupling with the smallest dimensions, $\lambda$ itself in this example, has a vanishing $\beta$-functions to all orders in perturbation theory:

$$
\begin{equation*}
\beta_{5 d}^{(\lambda)}=0 \tag{5.8.7}
\end{equation*}
$$

There is no analogue of the log resummation needed in treating marginal couplings and hence no need to improve the perturbative expansion.

Let us see what happens in a mass-dependent scheme. It is convenient to introduce a mass scale $M$ and write the interaction as $\lambda / M \phi^{4}$, with $\lambda$ dimensionless and of $\mathcal{O}(1)$. The scale $M$ is the scale below which this theory makes sense as an effective field theory. We assume that the size of all dimensionful operators is governed by this scale, that is any dimensionful coupling of mass dimension $-n$ can be written as a dimensionless coupling of order one times $1 / M^{n}$. Let us focus on the 1PI 4 -point function. This is linearly divergent in a cut-off regularization and can be renormalized by momentum subtraction using eq. (5.2.4), with $\lambda \rightarrow \lambda / M$. The finite, 1PI four-point function, the 5 d analogue of eq. (5.2.3), is

$$
\begin{align*}
\Gamma_{5 d}^{(4)}(s, t, u)= & -\frac{\lambda(\mu)}{M}-\frac{\lambda^{2}(\mu)}{32 \pi^{2} M^{2}} \int_{0}^{1} d x\left(\sqrt{m^{2}-s x(1-x)}-\sqrt{m^{2}+\mu^{2} x(1-x)}\right) \\
& +(s \rightarrow t)+(s \rightarrow u) \tag{5.8.8}
\end{align*}
$$

The logs are replaced by square roots in 5 d. In order to see that the sliding scale is not of great help, fix, say, $\mu=0$ and consider the high energy regime $-s \gg m^{2}$. Like the 4 d case (5.2.4), the one-loop term grows and can in principle be comparable to the tree-level term. This will occur at energies $\lambda \sqrt{|s|} /\left(32 \pi^{2} M\right) \sim 1$, which implies $|s|>M^{2}$. But this is beyond the regime in which the effective field theory makes sense, which is $|s| \ll M^{2}$ ! There is no energy regime in which the perturbative expansion can be improved by some RG resummation. The theory is simply weakly coupled for energies below $M$ and becomes strongly coupled for energies above $M$. Nevertheless, there is nothing intrinsically wrong in introducing the sliding scale $\mu$, so we can keep going and compute the $\beta$-function for $\lambda$ by using the CS equations (5.4.3). In this way we get the 5 d analogue of eq. (5.2.8):

$$
\begin{equation*}
\beta_{5 d}^{(\lambda)}=\frac{3 \lambda^{2}}{32 \pi^{2} M} \int_{0}^{1} d x \frac{\mu^{2} x(1-x)}{\sqrt{m^{2}+\mu^{2} x(1-x)}}+\mathcal{O}\left(\lambda^{3}\right) \tag{5.8.9}
\end{equation*}
$$

For $\mu \gg m$ this simplifies to

$$
\begin{equation*}
\beta_{5 d}^{(\lambda)} \simeq \frac{3 \lambda^{2} \sqrt{\mu^{2}}}{256 \pi M}, \quad \mu \gg m . \tag{5.8.10}
\end{equation*}
$$

The approximate solution of the RG flow is

$$
\begin{array}{ll}
\lambda_{U V}(\mu) \simeq \frac{\lambda\left(\mu_{0}\right)}{1-\frac{3 \lambda\left(\mu_{0}\right)}{256 \pi M}\left(\mu-\mu_{0}\right)}, & \mu_{0}, \mu \gg m \\
\lambda_{I R}(\mu) \simeq \mathrm{constant}, & \mu \ll m . \tag{5.8.12}
\end{array}
$$

Let us finally compare the $\Gamma_{5 d}^{(4)}$ one obtains with and without the use of the RG technique at high energies (the analogue of eqs. (5.2.12) and (5.2.13)) for $s=t=u \equiv-E^{2} \gg m^{2}$ :

$$
\begin{align*}
\Gamma_{5 d}^{(4)}(E) & \simeq-\frac{\lambda}{M}-\frac{3 \lambda^{2} E}{256 \pi M^{2}}+\mathcal{O}\left(\lambda^{2}\right),  \tag{5.8.13}\\
\Gamma_{5 d}^{(4)}(E) & \simeq-\frac{\lambda(E)}{M}=\frac{-\lambda / M}{1-\frac{3 \lambda E}{256 \pi M}}+\mathcal{O}\left(\lambda^{2}\right) . \quad(\mathrm{RG}) \tag{5.8.14}
\end{align*}
$$

Like in 4d, eq. (5.8.14) reproduces the one-loop result (5.8.13) and, in addition, encodes higher order terms. However, in contrast to the 4 d case, the higher order terms are scheme-dependent and are not "special" in the perturbative expansion. They are of order $\lambda^{l+1}(E / M)^{l}$, which is the order expected for a generic $l$-loop computation. There is no analogue of the logarithmic enhancement found in the 4 d case. Hence eq. (5.8.14) should only be trusted at $\mathcal{O}\left(\lambda^{2}\right)$, in which case it merely reproduces the perturbative result obtained with no RG technique.

The vanishing of $\beta_{5 d}^{(\lambda)}$ in DR at one-loop level is immediately seen by noting that the one-loop integral is proportional to $\Gamma(2-d / 2)$. No divergence then occurs because the
(analytic continuation of the) Gamma function is well-behaved for negative half-integers values of its argument.

Scheme-independent $\beta$-functions can occur in this example for higher dimensional couplings. Consider for instance the dimension 9 operator $\phi^{6}$, that will be generated at some order in perturbation theory due to the non-renormalizability of the theory. Its coupling $g$ has classical dimension -4 , so by dimensional analysis we expect in a mass-independent scheme

$$
\begin{equation*}
\beta_{5 d}^{(g)}=c \lambda^{4} \tag{5.8.15}
\end{equation*}
$$

where $c$ is a generally non-vanishing constant. ${ }^{10}$ Since $\lambda$ does not run, the solution to eq.(5.8.15) is simply

$$
\begin{equation*}
g(\mu)=g\left(\mu_{0}\right)+c \lambda^{4} \log \left(\frac{\mu}{\mu_{0}}\right) . \tag{5.8.16}
\end{equation*}
$$

We see that log terms can appear but in a rather dull way. Moreover, we should keep in mind that log's grow slowly and the range of energies we can explore is limited by the range of validity of the non-renormalizable effective theory.

A similar analysis can also be made for relevant couplings, with the obvious crucial difference that their effect decreases, rather than increases, in the UV. An example is once again provided by the $\phi^{4}$ theory, this time in three dimensions, where the theory is superrenormalizable. The coupling $\lambda$ has classical mass dimension +1 and in mass-independent schemes, and to all orders in perturbation theory, we have

$$
\begin{equation*}
\beta_{3 d}^{(\lambda)}=0, \tag{5.8.17}
\end{equation*}
$$

while in mass-dependent schemes $\beta_{3 d}^{(\lambda)} \neq 0 .{ }^{11}$

## 5.9 "Relevant" RG Flow of Dimensionful Couplings and Renormalization of Composite Operators

We have seen in section 5.8 that the use of RG techniques applied to relevant and irrelevant couplings is not always as useful as in the standard situation where we have marginal couplings only. We have provided an example, the $\phi^{4}$ theory in 5 space-time dimensions, where all couplings are irrelevant. In presence of marginal couplings, the situation changes.

[^31]In a theory where marginal and (ir)relevant couplings are present together, the former can lead to large logs effects on the latter that should be summed. In other words, the quantum RG flow of (ir)relevant operators, induced by marginal couplings, can and should be taken into account. This is what we mean by "relevant" RG flow in the title of this section.

If we denote by $c$ the coupling constant associated to the classically lowest dimensional irrelevant operator $\mathcal{O}_{c}$, assuming it to be unique for the moment, dimensional analysis requires that to all orders in perturbation theory ${ }^{12}$

$$
\begin{equation*}
\beta_{c}=f\left(\lambda_{i}\right) c, \tag{5.9.1}
\end{equation*}
$$

where we have denoted by $\lambda_{i}$ all the marginal couplings in the theory. Note that $\beta_{c}$ can have only a linear dependence on the coupling constant $c$, but an arbitrary dependence on the marginal couplings $\lambda_{i}$, parametrized in eq.(5.9.1) by the function $f$. The function $f$ corresponds to the anomalous dimension of the operator $\mathcal{O}_{c}$. Before showing this result, let us first introduce the concept of composite operator and its corresponding anomalous dimension.

Composite operators are obtained by taking products of elementary fields (and their derivatives) at the same space-time point. At the quantum level, when two fields approach the same space-time point, UV divergences arise. Naive product of fields does not hold, i.e. $\phi^{2}(x) \neq(\phi(x))^{2}$. Indeed, we know that $\left\langle\phi^{2}(x)\right\rangle \neq(\langle\phi(x)\rangle)^{2} .{ }^{13}$ For this reason, correlations functions involving composite operators are divergent, even in a renormalized theory where the infinities have been already buried in the counter-terms. Such additional divergences are removed by a renormalization of the composite operators themselves, generalization of the wave-function renormalization of the elementary fields. Composite operators can be relevant, marginal or irrelevant: $\phi^{2}, \phi^{4},(\partial \phi)^{2}, \phi^{6},(\partial \phi)^{4}, \partial_{\mu} \phi \partial_{\nu} \phi, \partial_{\mu} \partial_{\nu} \partial_{\rho} \phi^{4}$, etc. are all composite operators. They can in general carry Lorentz quantum numbers (spin), so we can have scalar operators, as well as tensor operators, like the last two illustrated above. If we denote by $\mathcal{O}$ a generic composite operator, in analogy to the elementary field case, we write the bare operator $\mathcal{O}^{B}=\mathcal{O}(\mu) Z^{\mathcal{O}}(\mu)$.

Let $G^{(n)}$ be a generic $n$-point connected Green function of $n$ operators $\mathcal{O}_{i}:{ }^{14}$

$$
\begin{equation*}
G_{B}^{(n)}\left(p_{1}, \ldots, p_{n}\right)=\left\langle\mathcal{O}_{1}^{B}\left(p_{1}\right) \ldots \mathcal{O}_{n}^{B}\left(p_{n}\right)\right\rangle=\prod_{i=1}^{n} Z^{\mathcal{O}_{i}}(\mu) G^{(n)}\left(p_{1}, \ldots, p_{n}\right) \tag{5.9.2}
\end{equation*}
$$

[^32]The straightforward generalization of the CS eqs. (5.4.3) - which are also valid for connected, rather than 1PI, Green functions - are

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\beta_{i}(\lambda) \frac{\partial}{\partial \lambda_{i}}+\sum_{i=1}^{n} \gamma^{\mathcal{O}_{i}}(\lambda)\right) G^{(n)}=0 \tag{5.9.3}
\end{equation*}
$$

where

$$
\begin{equation*}
\gamma^{\mathcal{O}_{i}}(\lambda)=\mu \frac{d \log Z^{\mathcal{O}_{i}}}{d \mu} \tag{5.9.4}
\end{equation*}
$$

is the anomalous dimension of the operator $\mathcal{O}_{i}$. In general, the situation is more complicated, because operators with the same quantum numbers typically mix under renormalization. ${ }^{15}$ The anomalous dimension becomes then a matrix of anomalous dimensions: $\mathcal{O}_{i}^{B}=Z_{i j}^{\mathcal{O}}(\mu) \mathcal{O}_{j}(\mu)$ and

$$
\begin{equation*}
\gamma_{i j}(\lambda)=\left(Z_{i k}^{\mathcal{O}}\right)^{-1} \mu \frac{d Z_{k j}^{\mathcal{O}}}{d \mu} \tag{5.9.5}
\end{equation*}
$$

We can now come back to our original problem of studying the RG evolution of an irrelevant coupling driven by marginal couplings, and show that the function $f$ appearing in eq.(5.9.1) can be identified with the anomalous dimension of its corresponding operator $\mathcal{O}_{c}$. Let us denote by $G^{(n)}$ a generic $n$-point Green function obtained from a Lagrangian containing, besides marginal interactions, the term $c \mathcal{O}_{c}$. We can bring down from the action the term $c \mathcal{O}_{c}$ so that (in momentum space)

$$
\begin{align*}
G^{(n)}\left(c, p_{1}, \ldots, p_{n}\right)=\left\langle\phi\left(p_{1}\right) \ldots \phi\left(p_{n}\right)\right\rangle_{c} & =\sum_{k=0}^{\infty} \frac{(i c)^{k}}{k!}\left\langle\phi\left(p_{1}\right) \ldots \phi\left(p_{n}\right) \mathcal{O}_{c}^{k}(0)\right\rangle_{0} \\
& \equiv \sum_{k=0}^{\infty} \frac{(i c)^{k}}{k!} G^{(n, k)}\left(0, p_{1}, \ldots, p_{n}\right) \tag{5.9.6}
\end{align*}
$$

The CS eqs. satisfied by the $G^{(n, k)}$ are

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\beta_{i}(\lambda) \frac{\partial}{\partial \lambda_{i}}+n \gamma+k \gamma^{\mathcal{O}_{c}}(\lambda)\right) G^{(n, k)}\left(0, p_{1}, \ldots, p_{n}\right)=0 \tag{5.9.7}
\end{equation*}
$$

We can multiply eq. (5.9.7) by $(i c)^{k} / k$ ! and sum over $k$ to write

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\beta_{i}(\lambda) \frac{\partial}{\partial \lambda_{i}}+n \gamma+c \gamma^{\mathcal{O}_{c}}(\lambda) \frac{\partial}{\partial c}\right) G^{(n)}\left(c, p_{1}, \ldots, p_{n}\right)=0 \tag{5.9.8}
\end{equation*}
$$

where we used that $c \partial G^{(n)} / \partial c=\sum_{k} k(i c)^{k} / k!G^{(n, k)}$. The $\beta$-function for $c$ can be identified by looking at the term multiplying $\partial / \partial c$. We then have

$$
\begin{equation*}
\beta_{c}=c \gamma^{\mathcal{O}}\left(\lambda_{i}\right) \tag{5.9.9}
\end{equation*}
$$

[^33]where the anomalous dimension $\gamma^{\mathcal{O}}\left(\lambda_{i}\right)$ is computed in the theory with $c=0$. Comparing eq.(5.9.9) with (5.9.1), we see that
\[

$$
\begin{equation*}
f=\gamma^{\mathcal{O}} \tag{5.9.10}
\end{equation*}
$$

\]

Let us solve the RG flow eqution (5.9.9) in the case in which we have only one marginal coupling $\lambda$, with $\gamma^{\mathcal{O}}(\lambda)=\gamma_{0} \lambda+\ldots, \beta_{\lambda}(\lambda)=b_{0} \lambda^{2}+\ldots$. In this case, using the chain rule, we can rewrite eq.(5.9.9) as

$$
\begin{equation*}
\frac{d c}{d \lambda} \beta_{\lambda}(\lambda)=c \gamma^{\mathcal{O}}(\lambda) \tag{5.9.11}
\end{equation*}
$$

that, at leading order in $\lambda$, admits the solution

$$
\begin{equation*}
c(\mu)=c\left(\mu_{0}\right)\left(\frac{\lambda(\mu)}{\lambda\left(\mu_{0}\right)}\right)^{\frac{\gamma_{0}}{b_{0}}} \tag{5.9.12}
\end{equation*}
$$

We see how the RG flow of the marginal coupling $\lambda(\mu)$ drives an RG flow for the irrelevant coupling $c$, with an effect that sensitively depends on the anomalous dimension of the operator $\mathcal{O}_{c}$ (determined by $\gamma_{0}$ ) and of course on the running of $\lambda$ (determined by $b_{0}$ ), that enters both explicitly and implicitly in defining $\lambda(\mu)$ in eq.(5.9.12).

More in general, if the lowest dimensional irrelevant operator is not unique, we can have a set of operators $c_{n} \mathcal{O}_{n}$ that mix under renormalization. Repeating the analysis above, we get

$$
\begin{equation*}
\beta_{c_{n}}=\gamma_{n m}(\lambda) c_{m} \tag{5.9.13}
\end{equation*}
$$

In many cases eq.(5.9.13) is enough to have a good description of the physics because, by definition, irrelevant couplings are small in the IR and the lowest dimensional ones capture the main effects we want to study. ${ }^{16}$

Irrelevant operators can also come from an underlying renormalizable theory when some degrees of freedom are integrated out. We will study this topic in chapter 7 , where we will also consider an explicit example of RG flow of an irrelevant coupling.

The above analysis can be repeated in the case in which $c$ is a relevant operator. We will not repeat the general analysis for relevant operators, but focus on the particular case of the mass term, which is a relevant operator. The physical mass, defined in a mass-dependent scheme as $\Gamma^{(2)}\left(p^{2}=m^{2}\right)=0$, cannot depend on the energy scale. But non-physical masses in mass-independent schemes do have such a dependence. This is particularly simple in a theory with no irrelevant couplings where, by dimensional analysis and to all orders in perturbation theory, we have

$$
\begin{equation*}
\beta_{m^{2}}=\gamma_{m}\left(\lambda_{i}\right) m^{2}, \tag{5.9.14}
\end{equation*}
$$

[^34]where $\gamma_{m}$ is the anomalous dimension of the mass operator. In the $\phi^{4}$ theory the composite operator associated to the mass is $\phi^{2}, \gamma_{m}=\gamma^{\phi^{2}}$. It is a useful exercise to determine $\gamma^{\phi^{2}}$ at one-loop level. This can be derived by studying the Green function $G^{(2,1)}=\left\langle\phi \phi \phi^{2}\right\rangle$. Since $Z=1$ at one-loop, we have $G^{(2,1)}=\left(Z^{\phi^{2}}\right)^{-1} G_{B}^{(2,1)}$. The divergent part of $G^{(2,1)}$ is readily computed by setting the external momentum to zero. We have
\[

$$
\begin{equation*}
G_{B, \operatorname{div}}^{(2,1)}(0)=-i \lambda \int \frac{d^{d} p}{(2 \pi)^{d}} \frac{i^{2}}{\left(p^{2}\right)^{2}}=-\frac{\lambda}{16 \pi^{2} \epsilon}, \tag{5.9.15}
\end{equation*}
$$

\]

and hence

$$
\begin{equation*}
Z^{\phi^{2}}=1-\frac{\lambda}{16 \pi^{2} \epsilon} . \tag{5.9.16}
\end{equation*}
$$

Recalling that $\beta(\lambda)=-\epsilon \lambda+\mathcal{O}(1)$ in $d$ dimensions, one gets

$$
\begin{equation*}
\gamma^{\phi^{2}}=\frac{d \log Z^{\phi^{2}}}{d \log \mu}=\frac{\lambda}{16 \pi^{2}} . \tag{5.9.17}
\end{equation*}
$$

Using eq. (5.9.14), we conclude that

$$
\begin{equation*}
\beta_{m^{2}}=m^{2} \frac{\lambda}{16 \pi^{2}} . \tag{5.9.18}
\end{equation*}
$$

We can also get eq.(5.9.18) directly, by-passing the above analysis. The one-loop mass correction in the $\phi^{4}$ theory is given by the tadpole graph in fig. 5.2. In DR, the one-loop 1PI 2-point function reads

$$
\begin{equation*}
i \Gamma^{(2)}(p)=\frac{-i \mu^{\epsilon} \lambda}{2} \int \frac{d^{d} k}{(2 \pi)^{d}} \frac{i}{k^{2}-m^{2}}-i \delta_{m}=-\frac{i \mu^{\epsilon} \lambda}{2(4 \pi)^{d / 2}} \frac{\Gamma(2-d / 2)}{(1-d / 2)} m^{d-2}-i \delta_{m} \tag{5.9.19}
\end{equation*}
$$

In the $\overline{\mathrm{MS}}$ scheme the counter-term is

$$
\begin{equation*}
\delta_{m}=\frac{m^{2} \lambda}{32 \pi^{2}}\left(\frac{2}{\epsilon}-\gamma_{E}+\log 4 \pi\right) \tag{5.9.20}
\end{equation*}
$$

so that the tree+1-loop 1PI 2-point function reads

$$
\begin{equation*}
\Gamma^{(2)}(p)=p^{2}-m \frac{2}{\mathrm{MS}}+\frac{m^{2} \lambda}{32 \pi^{2}}\left(1+\log \frac{\mu^{2}}{m^{2}}\right) \tag{5.9.21}
\end{equation*}
$$

In the physical scheme we simply have

$$
\begin{equation*}
\Gamma^{(2)}(p)=p^{2}-m^{2} \tag{5.9.22}
\end{equation*}
$$

with no finite one-loop corrections left, where $m$ is the physical, $\mu$-independent, mass. Matching eqs.(5.9.21) and (5.9.22) gives

$$
\begin{equation*}
m^{2}=m \frac{2}{\overline{\mathrm{MS}}}(\mu)-\frac{m^{2} \lambda}{32 \pi^{2}}\left(1+\log \frac{\mu^{2}}{m^{2}}\right), \tag{5.9.23}
\end{equation*}
$$



Figure 5.2: One-loop tadpole graph contributing to the renormalization of the mass in the $\phi^{4}$ theory.
and hence

$$
\begin{equation*}
0=\mu \frac{d m^{2}}{d \mu}=\beta_{m^{2}}-\frac{m^{2} \lambda}{16 \pi^{2}}+\mathcal{O}\left(\lambda^{2}\right) \Longrightarrow \beta_{m^{2}}=m^{2} \frac{\lambda}{16 \pi^{2}} \tag{5.9.24}
\end{equation*}
$$

which reproduces eq.(5.9.18), as expected.
The use of $\mu$-dependent masses is unavoidable when using mass-independent schemes such as MS or $\overline{\mathrm{MS}}$. There is nothing wrong in computing amplitudes in terms of unphysical masses (like the mass terms in the $\overline{\mathrm{MS}}$ scheme), in which case RG flow techniques should be used to relate mass terms at different energy scales. Eventually one relates these masses to the physical ones, using formulae like eq. (5.9.23). As a matter of fact, $\overline{\mathrm{MS}}$ masses are often used in the literature, being the $\overline{\mathrm{MS}}$ scheme one of the most popular schemes in perturbative computations. For particles like quarks that do not appear in asymptotic states and for which the direct physical mass definition is unavailable, the story is more complicated. The $\overline{\mathrm{MS}}$ quark mass, for example, can be defined to be $m^{2} \frac{2}{\mathrm{MS}}(M)$, evaluated at some indirectly derived "pole" mass $M$. For a heavy quark, for which $M \gg \Lambda_{Q C D}$, the mass $M$ can roughly be computed as the mass of the meson bound states $\bar{Q} Q$ divided by two. For light quarks, $M$ can be computed using chiral perturbation theory, as explained in section 8.7. For $M \simeq \Lambda_{Q C D}$, more complicated procedures are needed.

### 5.10 RG Improved Effective Potential

The RG technique is also useful in the context of effective potentials. The summation of LL leads to so called RG improved Effective Potentials. Our favorite $\phi^{4}$ theory is particularly instructive in this case, since it shows how RG improved potentials help us in correcting fake perturbative results. We have seen in section 4.2 that the CW potential for a massless $\phi^{4}$ theory reads (omitting the irrelevant constant term):

$$
\begin{equation*}
V_{e f f}(\phi)=\frac{\lambda}{4!} \phi^{4}+\frac{\lambda^{2}}{256 \pi^{2}} \phi^{4} \log \frac{\phi^{2}}{\phi_{0}^{2}} . \tag{5.10.1}
\end{equation*}
$$

where $\phi_{0}$ is an arbitrary fixed energy scale. Let us look for the extrema of $\phi$ :

$$
\begin{equation*}
0=\frac{d V_{e f f}(\phi)}{d \phi}=\frac{\phi^{3} \lambda}{384 \pi^{2}}\left(64 \pi^{2}+3 \lambda+6 \lambda \log \phi^{2} / \phi_{0}^{2}\right) \rightarrow \phi=0, \phi= \pm \bar{\phi}= \pm \phi_{0} e^{-\left(\frac{1}{4}+\frac{16 \pi^{2}}{3 \lambda}\right)} \tag{5.10.2}
\end{equation*}
$$

The classical minimum at $\phi=0$ turns into a maximum and two new symmetric minima arise at $\pm \bar{\phi}$. The one-loop truncation of the potential is reliable provided the tree-level term is greater than the one-loop term, namely for field values such that

$$
\begin{equation*}
\frac{3 \lambda}{16 \pi^{2}}|\log \phi| \ll 1 \tag{5.10.3}
\end{equation*}
$$

The condition (5.10.3) is manifestly violated at the minima $|\bar{\phi}|$, so we cannot trust the result we have found for such small values of $\phi$. This problem is easily solved by RG arguments. Let us define a running coupling $\lambda(\mu)$ by

$$
\begin{equation*}
V_{e f f}(\phi=\mu) \equiv \frac{\lambda(\mu)}{4!} \mu^{4} \tag{5.10.4}
\end{equation*}
$$

In terms of $\lambda(\mu)$ the potential reads

$$
\begin{equation*}
V_{e f f}(\phi)=\frac{\lambda(\mu)}{4!} \phi^{4}+\frac{\lambda^{2}(\mu)}{256 \pi^{2}} \phi^{4} \log \frac{\phi^{2}}{\mu^{2}} . \tag{5.10.5}
\end{equation*}
$$

We already know how $\lambda(\mu)$ flows with the energy scale, but it is instructive to see how the $\beta$-function for $\lambda$ can be computed by demanding the $\mu$-invariance of $V_{\text {eff }}$. Recalling that $\gamma_{\phi}=0$ up to one-loop level, we get

$$
\begin{equation*}
0=\mu \frac{d V_{e f f}}{d \mu}=\beta(\lambda) \frac{\phi^{4}}{4!}-\frac{\lambda^{2} \phi^{4}}{128 \pi^{2}}+\mathcal{O}\left(\lambda^{3}\right) \rightarrow \beta(\lambda)=\frac{3 \lambda^{2}}{16 \pi^{2}} \tag{5.10.6}
\end{equation*}
$$

By choosing $\mu=\phi$ in eq. (5.10.5), we can get rid of the log term and write

$$
\begin{equation*}
V_{e f f}(\phi)=\frac{\lambda(\phi)}{4!} \phi^{4}=\frac{\lambda_{0}}{1-\frac{3 \lambda_{0}}{16 \pi^{2}} \log \frac{\phi}{\phi_{0}}} \frac{\phi^{4}}{4!}, \tag{5.10.7}
\end{equation*}
$$

where $\phi_{0}$ is an arbitrary scale. The minima at $\phi= \pm \bar{\phi}$ have disappeared in the potential (5.10.7), which manifestly increases monotonically when $\phi$ increases. The potential (5.10.7) is the RG improved version of the effective potential (5.10.1). Expanding the log term in eq. (5.10.7), we recover eq. (5.10.1) plus all the LL terms summed by the RG technique. The origin of the fake result (5.10.2) should now be clear. The minima at $\phi= \pm \bar{\phi}$ were obtained by forgetting the large log's that appear for so small values of $\phi$ (small energies). Being the $\phi^{4}$ theory free in the IR, the effective coupling at such small energies becomes smaller and smaller and the actual minimum is in fact the tree-level one.

### 5.11 Anomalous Dimension of the Photon and QED $\beta$-function

The WT identities imply that the radiative corrections to the photon propagator are transverse to all orders in perturbation theory, as described by eq. (3.3.5). It is useful to explicitly check this result at one-loop order. In so doing, we will determine the one-loop counter-term $Z_{3}$ defined in eq.(3.3.11). Once $Z_{3}$ is known, we can compute the photon anomalous dimension and $\beta$-function for $e$. The only one-loop graph contributing to the photon propagator is the electron loop:


We can bring together the two electron propagators appearing in eq.(5.11.1) by introducing the Feynman parameter $x$ :

$$
\begin{equation*}
\frac{1}{a b}=\int_{0}^{1} d x[a x+b(1-x)]^{-2} \tag{5.11.2}
\end{equation*}
$$

Performing also the Dirac algebra, we have

$$
\begin{equation*}
i \Pi_{\mu \nu}^{(1)}=-d e^{2} \mu^{\epsilon} \int_{0}^{1} d x \int \frac{d^{d} k}{(2 \pi)^{d}} \frac{k_{\mu}(k+q)_{\nu}+k_{\nu}(k+q)_{\mu}+\left(m^{2}-k \cdot(q+k)\right) \eta_{\mu \nu}}{\left[\left(k^{2}-m^{2}\right)(1-x)+x(q+k)^{2}-x m^{2}\right]^{2}} . \tag{5.11.3}
\end{equation*}
$$

Let us now redefine $k \rightarrow k-q x$, so that the odd term $k \cdot q$ in the denominator is cancelled. By symmetry, after the momentum shift, the terms in the numerator odd in $k$ vanish. We have then

$$
\begin{equation*}
i \Pi_{\mu \nu}^{(1)}=-d e^{2} \mu^{\epsilon} \int_{0}^{1} d x \int \frac{d^{d} k}{(2 \pi)^{d}} \frac{2\left(k_{\mu} k_{\nu}-q_{\mu} q_{\nu} x(1-x)\right)+\eta_{\mu \nu}\left(m^{2}+q^{2} x(1-x)-k^{2}\right)}{\left(k^{2}-\Delta\right)^{2}} . \tag{5.11.4}
\end{equation*}
$$

where $\Delta \equiv m^{2}-q^{2} x(1-x)$. Recall that $k_{\mu} k_{\nu} f\left(k^{2} / d\right) \eta_{\mu \nu}$ and perform then the Wick rotation to euclidean signature: $k^{0} \rightarrow i k_{4}, k^{2} \rightarrow-k^{2}$. After these steps, eq.(5.11.4) is rewritten as
$i \Pi_{\mu \nu}^{(1)}=-i d e^{2} \mu^{\epsilon} \int_{0}^{1} d x \int \frac{d^{d} k}{(2 \pi)^{d}} \frac{(1-2 / d) k^{2} \eta_{\mu \nu}+\left(m^{2}+q^{2} x(1-x)\right) \eta_{\mu \nu}-2 q_{\mu} q_{\nu} x(1-x)}{\left(k^{2}+\Delta\right)^{2}}$.

Let us now focus on the divergent part of eq.(5.11.5), i.e. to the $1 / \epsilon$ pole. Using the tricks described in section 3.4, we immediately get

$$
\begin{align*}
i \Pi_{\mu \nu}^{(1) d i v .} & =-i d \frac{e^{2}}{8 \pi^{2} \epsilon} \int_{0}^{1} d x\left[\left(1-\frac{2}{d}\right) \eta_{\mu \nu}(-2 \Delta)+\left(m^{2}+q^{2} x(1-x)\right) \eta_{\mu \nu}-2 q_{\mu} q_{\nu} x(1-x)\right] \\
& =i\left(\eta_{\mu \nu} q^{2}-q_{\mu} q_{\nu}\right) \frac{e^{2}}{8 \pi^{2}}\left(-\frac{4}{3}\right) \frac{1}{\epsilon} \tag{5.11.6}
\end{align*}
$$

The one-loop correction $\Pi_{\mu \nu}^{(1)}$ is in the transverse form expected from the WT identity (3.3.5). The value of the counter-term $Z_{3}$ needed to cancel the divergence (5.11.6) is (see fig.3.2)

$$
\begin{equation*}
Z_{3}=1-\frac{e^{2}}{6 \pi^{2}} \frac{1}{\epsilon} \tag{5.11.7}
\end{equation*}
$$

where we used here the MS subtraction scheme. The photon anomalous dimension is

$$
\begin{equation*}
\gamma_{A}=\frac{1}{2} \frac{d \log Z_{3}}{d \log \mu}=-\frac{e \beta_{e}^{(d)}}{6 \pi^{2}} \frac{1}{\epsilon} \tag{5.11.8}
\end{equation*}
$$

where $\beta_{e}^{(d)}$ is the QED $\beta$-function in $d$ (and not 4) dimensions. This is easily determined from eq.(3.3.13), that in DR reads $e_{B}=Z_{3}^{-1 / 2} \mu^{\epsilon / 2} e$ (recall the identity $Z_{1}=Z_{2}$ ). Since the bare charge $e_{B}$ does not depend on $\mu$, we have

$$
\begin{equation*}
0=\mu \frac{d e_{B}}{d \mu}=Z_{3}^{-1 / 2}\left(-e \gamma_{A}+\frac{1}{2} \epsilon e+\beta_{e}^{(d)}\right) \Longrightarrow \beta_{e}^{(d)}=-\frac{1}{2} \epsilon e+e \gamma_{A} \tag{5.11.9}
\end{equation*}
$$

Plugging eq.(5.11.9) in eq.(5.11.8) gives, up to order $e^{2}$,

$$
\begin{equation*}
\gamma_{A}=\frac{e^{2}}{12 \pi^{2}} \tag{5.11.10}
\end{equation*}
$$

In turn, eq.(5.11.10) allows us to determine the actual 4D $\beta$-function $\beta_{e}$ :

$$
\begin{equation*}
\beta_{e}=e \gamma_{A}=\frac{e^{3}}{12 \pi^{2}} \tag{5.11.11}
\end{equation*}
$$

It is useful to write the RG behavior in terms of $\alpha \equiv e^{2} /(4 \pi)$ and solve for its inverse. One gets, for $\mu_{0}, \mu \gg m$,

$$
\begin{equation*}
\alpha^{-1}(\mu)=\alpha^{-1}\left(\mu_{0}\right)-\frac{2}{3 \pi} \log \frac{\mu}{\mu_{0}} \tag{5.11.12}
\end{equation*}
$$

For $\mu<m$, instead, like for the coupling $\lambda$ in the $\phi^{4}$ theory, $\alpha$ does not run and its value is approximately given by $\alpha(m)$. We are clearly assuming here that there is a single charged particle. In presence of more charged particles with different masses, $\alpha$ stops running at the scale given by the lowest charged particle. Notice that in the real world, where the electron is the lightest charged particle, the absence of massless charged particles is crucial to prevent the electric charge (and hence all electrodynamical interactions) to vanish in the far IR.

## Chapter 6

## Non-Abelian Gauge Theories

### 6.1 Introduction and Classical Analysis

Non-abelian gauge theories are at the base of our current understanding of particle physics. Both the strong and the electroweak interactions are described in terms of them. These theories are based on a generalization of the QED $U(1)$ gauge symmetry, where two transformations do not necessarily commute with each other (hence the name non-abelian). Before describing them, let us quickly review the role of the $U(1)$ symmetry in QED. We assume an invariance of the Lagrangian under local (i.e. space-time dependent) transformations parametrized by a function $\lambda(x)$, under which any field $\psi$ carrying charge $q$ transforms as

$$
\begin{equation*}
\psi(x) \rightarrow e^{i q \lambda(x)} \psi(x) \tag{6.1.1}
\end{equation*}
$$

Due to the space-time dependence of the transformation, the derivative of the field $\partial_{\mu} \psi$ does not transform covariantly:

$$
\begin{equation*}
\partial_{\mu} \psi(x) \rightarrow e^{i q \lambda(x)}\left(\partial_{\mu} \psi(x)+i q \psi(x) \partial_{\mu} \lambda(x)\right) . \tag{6.1.2}
\end{equation*}
$$

We then add a gauge field (photon) $A_{\mu}$ that transforms inhomogeneously:

$$
\begin{equation*}
A_{\mu}(x) \rightarrow A_{\mu}(x)+\partial_{\mu} \lambda(x), \tag{6.1.3}
\end{equation*}
$$

that allows us to define a new (covariant) derivative transforming covariantly under $U(1)$ gauge transformations:

$$
\begin{equation*}
D_{\mu} \psi(x) \equiv \partial_{\mu} \psi(x)-i q A_{\mu}(x) \psi(x) \rightarrow e^{i \lambda(x)} D_{\mu} \psi(x) \tag{6.1.4}
\end{equation*}
$$

Under infinitesimal $U(1)$ transformations parametrized by $\epsilon(x) \ll 1$, we have

$$
\begin{equation*}
\delta_{\epsilon} \psi(x)=i q \epsilon(x) \psi(x), \quad \delta_{\epsilon} A_{\mu}(x)=\partial_{\mu} \epsilon(x) \tag{6.1.5}
\end{equation*}
$$

The QED Lagrangian is constructed by forming gauge invariant combinations of $\psi, D_{\mu} \psi$ and of the field strength

$$
\begin{equation*}
F_{\mu \nu}=\partial_{\mu} A_{\nu}-\partial_{\nu} A_{\mu} \tag{6.1.6}
\end{equation*}
$$

Non-abelian gauge theories are constructed by generalizing the above construction to a set of fields (scalars or fermions) $\psi_{l}$, with $l$ labeling the different fields. We assume that under an infinitesimal gauge transformation, the $\psi_{l}$ are rotated among each other:

$$
\begin{equation*}
\delta_{\epsilon} \psi_{l}(x)=i \epsilon^{\alpha}\left(t_{\alpha}\right)_{l}^{m} \psi_{m}(x), \tag{6.1.7}
\end{equation*}
$$

where $t_{\alpha}$ are a set of constant matrices, labelled by the index $\alpha$, satisfying the relation

$$
\begin{equation*}
\left[t_{\alpha}, t_{\beta}\right]=i C_{\alpha \beta}^{\gamma} t_{\gamma} \tag{6.1.8}
\end{equation*}
$$

The coefficients $C_{\alpha \beta}^{\gamma}$ are a set of real parameters denoted the structure constants. They are manifestly antisymmetric in the two lower indices: $C_{\alpha \beta}^{\gamma}=-C_{\beta \alpha}^{\gamma}$. We have written the index $\gamma$ upstairs, because in principle there might be a non-trivial metric $g_{\alpha \beta}$ in group space, to raise and lower the group indices. In the physically most interesting cases, this metric can be chosen to be the identity. From now on, we will assume a trivial group metric. Correspondingly, the position of the indices $\alpha$, $\beta$, etc. will be irrelevant. In this case, it can be shown that the structure constants $C_{\alpha \beta \gamma}$ become antisymmetric in all three indices.

Matrices satisfying the condition (6.1.8) form a so-called Lie algebra, namely they define infinitesimal transformations of a so called Lie group. Differently from the QED case, the group of transformations (6.1.7) is multidimensional and this explains the origin of the index $\alpha$, that runs from 1 up to $\operatorname{dim} G$, the number of dimensions of the group. The matrices $t_{\alpha}$ are nothing else than a set of generators of the Lie group, meaning that any group transformation can be written in terms of these matrices, that are linearly independent from each other. Like any set of matrices, the $t_{\alpha}$ 's also satisfy the so called Jacobi identity

$$
\begin{equation*}
\left[\left[t_{\alpha}, t_{\beta}\right], t_{\gamma}\right]+\left[\left[t_{\gamma}, t_{\alpha}\right], t_{\beta}\right]+\left[\left[t_{\beta}, t_{\gamma}\right], t_{\alpha}\right]=0 . \tag{6.1.9}
\end{equation*}
$$

Using eq. (6.1.8), this identity implies the following constraints among the structure constants:

$$
\begin{equation*}
C_{\alpha \beta}^{\omega} C_{\omega \gamma}^{\delta}+C_{\gamma \alpha}^{\omega} C_{\omega \beta}^{\delta}+C_{\beta \gamma}^{\omega} C_{\omega \alpha}^{\delta}=0 \tag{6.1.10}
\end{equation*}
$$

Depending on how many fields $\psi_{l}$ we have, the matrices $t_{\alpha}$ are said to be in different representations of the Lie algebra. In general, there is an infinite set of matrices (of different size), satisfying eqs. (6.1.8) and (6.1.9). Among these, a special role is played by the "adjoint" representation. This is the representation in which the generators $t_{\alpha}^{A d j}$ are
$\operatorname{dim} G \times \operatorname{dim} G$ matrices, i.e. the indices $l, m$ coincide with the indices $\alpha, \beta$. An explicit form of this representation is given by

$$
\begin{equation*}
\left(t_{\alpha}^{A d j}\right)_{\gamma}^{\beta}=i C_{\alpha \gamma}^{\beta} . \tag{6.1.11}
\end{equation*}
$$

Indeed, it is straightforward to check that the Jacobi identity (6.1.10) can be rewritten as

$$
\begin{equation*}
\left[t_{\alpha}^{A d j}, t_{\beta}^{A d j}\right]=C_{\alpha \beta}^{\gamma} t_{\gamma}^{A d j} \tag{6.1.12}
\end{equation*}
$$

We do not enter here in any detail concerning the definition of Lie algebras, Lie groups, etc., because all this will be extensively treated in the group theory course. However, we need to introduce another relevant representation, the fundamental. It might be useful to consider a simple example of non-abelian Lie group, $G=S U(2)$. $S U(2)$ is defined by the set of $2 \times 2$ unitary matrices $U^{\dagger}=U^{-1}$ with unit determinant (that's why the name S (for special, with unit determinant) $\mathrm{U}($ for unitary $)(2))$. It is straightforward to check that this set is in fact a group and it is three dimensional. Any $S U(2)$ matrix $U$ can be written as

$$
\begin{equation*}
U=e^{i \omega_{\alpha} t_{\alpha}}, \quad \alpha=1,2,3 \tag{6.1.13}
\end{equation*}
$$

where $t_{\alpha}=\sigma_{\alpha} / 2$ and $\sigma_{\alpha}$ are the usual Pauli matrices

$$
\sigma_{1}=\left(\begin{array}{cc}
0 & 1  \tag{6.1.14}\\
1 & 0
\end{array}\right), \quad \sigma_{2}=\left(\begin{array}{cc}
0 & -i \\
i & 0
\end{array}\right), \quad \sigma_{3}=\left(\begin{array}{cc}
1 & 0 \\
0 & -1
\end{array}\right) .
$$

The structure constants are

$$
\begin{equation*}
C_{\alpha \beta \gamma}=\epsilon_{\alpha \beta \gamma}, \tag{6.1.15}
\end{equation*}
$$

where $\epsilon_{\alpha \beta \gamma}$ is the completely antisymmetric tensor, with $\epsilon_{123}=+1$. The above $2 \times 2$ matrices $t_{\alpha}$, that enter the definition of the group $S U(2)$, form the fundamental representation. The $3 \times 3$ matrices (6.1.11) define instead the adjoint representation. We can then have field "doublets" $D_{l}(l=1,2)$, that transforms as $D_{l} \rightarrow U_{l m} D_{m}$, with $U$ as in eq. (6.1.13) or field "triplets" $T_{l}(l=1,2,3)$, transforming as $T_{l} \rightarrow\left(U^{A d j}\right)_{l m} D_{m}$, where

$$
\begin{equation*}
U^{A d j}=e^{i \omega_{\alpha} t_{\alpha}^{A d j}}, \quad \alpha=1,2,3 . \tag{6.1.16}
\end{equation*}
$$

In addition to the fundamental and adjoint representations, there are an infinite number of other irreducible representations, labelled by the angular momentum $J$, of dimension $2 J+1$, with $J$ any positive integer or semi-integer number.

Similar considerations apply for more general groups, such as $S U(N)$, with $N>2$, or $S O(N)$, with $N>3 .{ }^{1}$ Instead of defining $\operatorname{dim} r \times \operatorname{dim} r$ matrices $U^{r}$, for each different

[^35]representation $r$, it is practically more convenient to write the transformation properties of $r$ in terms of, say, the matrices $U$ in the fundamental representation. All we need to know is how the representation $r$ in question appears in the tensor product of fundamentals. For instance, for $S U(N)$ groups, we have $\mathbf{N} \otimes \overline{\mathbf{N}}=\mathbf{N}^{2}-\mathbf{1} \oplus \mathbf{1}$, where $\mathbf{N}$ and $\overline{\mathbf{N}}$ are the fundamental and anti-fundamental (i.e. complex conjugate) representations, and $\mathbf{N}^{2}-\mathbf{1}$ is the adjoint one. A field $\psi$ in the adjoint representation of $S U(N)$ can correspondingly be written as an $N \times N$ matrix field $\psi_{i j}$, transforming as
\[

$$
\begin{equation*}
\psi \rightarrow U \psi U^{\dagger} \tag{6.1.17}
\end{equation*}
$$

\]

Coming back to physics, given the transformation (6.1.7), we add gauge fields $A_{\mu}^{\alpha}$, one for each independent direction in field space, so that we can form a covariant derivative. In analogy to the $U(1)$ case, the transformations of the gauge fields $A_{\mu}^{\alpha}$ must contain a term of the form $\partial_{\mu} \epsilon^{\alpha}$. Contrary to the $U(1)$ case, this cannot be the end of the story, since $\alpha$ is an index in the adjoint representation. The natural guess for the infinitesimal transformation of $A_{\mu}^{\alpha}$ is then

$$
\begin{equation*}
\delta A_{\mu}^{\alpha}=\partial_{\mu} \epsilon^{\alpha}+i \epsilon^{\beta}\left(t_{\beta}^{A d j}\right)_{\gamma}^{\alpha} A_{\mu}^{\gamma}=\partial_{\mu} \epsilon^{\alpha}+C_{\beta \gamma}^{\alpha} A_{\mu}^{\beta} \epsilon^{\gamma} . \tag{6.1.18}
\end{equation*}
$$

The covariant derivative is defined as

$$
\begin{equation*}
D_{\mu} \psi_{l} \equiv \partial_{\mu} \psi_{l}-i A_{\mu}^{\alpha}\left(t_{\alpha}\right)_{l}^{m} \psi_{m} \tag{6.1.19}
\end{equation*}
$$

It is straightforward to show that this guess is in fact correct and the covariant derivative transforms as it should:

$$
\begin{equation*}
\delta\left(D_{\mu} \psi\right)_{l}=i \epsilon^{\alpha}\left(t_{\alpha}\right)_{l}^{m}\left(D_{\mu} \psi\right)_{m} \tag{6.1.20}
\end{equation*}
$$

It is often convenient to write the components of the gauge fields $A_{\mu}^{\alpha}$ in matrix form by defining

$$
\begin{equation*}
A_{\mu}=A_{\mu}^{\alpha} t_{\alpha} . \tag{6.1.21}
\end{equation*}
$$

For simplicity of notation we have omitted, and from now on will be done most of the time, the gauge group indices in eq. (6.1.21). The finite form of eq. (6.1.18) is easily found by demanding that the gauge transformed connection $A \rightarrow A^{U}$, is such that

$$
\begin{equation*}
D_{\mu}(A) \psi \rightarrow U D_{\mu}(A) \psi \tag{6.1.22}
\end{equation*}
$$

for any field in a given representation $r$, transforming as $\psi \rightarrow \psi^{U}=U \psi$, where $U$ is defined as

$$
\begin{equation*}
U(x)=e^{i \Lambda_{\alpha}(x) t_{\alpha}} \tag{6.1.23}
\end{equation*}
$$

We have

$$
\begin{equation*}
\left(\partial_{\mu}-i A_{\mu}\right) \psi \rightarrow \partial_{\mu}(U \psi)-i A_{\mu}^{U} U \psi=U\left(\partial_{\mu} \psi-i A_{\mu} \psi\right)+\left(\partial_{\mu} U\right) \psi+\left(i U A_{\mu}-i A_{\mu}^{U} U\right) \psi \tag{6.1.24}
\end{equation*}
$$

Demanding that the last three terms in eq. (6.1.22) vanish uniquely fixes

$$
\begin{equation*}
A_{\mu}^{U}=U A_{\mu} U^{-1}-i\left(\partial_{\mu} U\right) U^{-1} \tag{6.1.25}
\end{equation*}
$$

The first term in eq. (6.1.25) is the one expected from a field in the adjoint representation (see eq. (6.1.17)), while the second is the inohomogenous one characterizing a gauge connection. For the $U(1)$ case the index $\alpha$ is trivial and eq. (6.1.25) reduces to eq. (6.1.3), with $\Lambda=\lambda$. For infinitesimal transformations, $\Lambda_{\alpha}=\epsilon_{\alpha}$, the transformation (6.1.25) correctly reproduces eq. (6.1.18), as it should.

The generalization of the $U(1)$ field strength $F_{\mu \nu}=\partial_{\mu} A_{\nu}-\partial_{\nu} A_{\mu}$ can be found by recalling that for any field $\psi$ with charge $q$, the field strength is proportional to the (commutator part of the) action of two covariant derivatives acting on the field itself. Taking $q=1$, we have

$$
\begin{equation*}
D_{\mu} D_{\nu} \psi=\left(\partial_{\mu}-i A_{\mu}\right)\left(\partial_{\nu}-i A_{\nu}\right) \psi=\left(\partial_{\mu} \partial_{\nu}-i \partial_{\mu} A_{\nu}-i A_{\nu} \partial_{\mu}-i A_{\mu} \partial_{\nu}-A_{\mu} A_{\nu}\right) \psi . \tag{6.1.26}
\end{equation*}
$$

Only the second term in the above equation survives when we take the antisymmetric combination in $\mu \leftrightarrow \nu$ :

$$
\begin{equation*}
\left[D_{\mu}, D_{\nu}\right] \psi=-i F_{\mu \nu} \psi \tag{6.1.27}
\end{equation*}
$$

The field strength in a non-abelian gauge theory can be defined by generalizing eq. (6.1.27). Denoting $\psi$ a field in an arbitrary representation of the gauge group, eq. (6.1.26) still applies, but now the last term does not vanish, since $A_{\mu}$ is a matrix. We have

$$
\begin{equation*}
\left[D_{\mu}, D_{\nu}\right] \psi=-i\left(\partial_{\mu} A_{\nu}-\partial_{\nu} A_{\mu}-i\left[A_{\mu}, A_{\nu}\right]\right) \psi \equiv-i F_{\mu \nu} \psi \tag{6.1.28}
\end{equation*}
$$

In components, $F_{\mu \nu}=F_{\mu \nu}^{\alpha} t_{\alpha}$, with

$$
\begin{equation*}
F_{\mu \nu}^{\alpha}=\partial_{\mu} A_{\nu}^{\alpha}-\partial_{\nu} A_{\mu}^{\alpha}+C_{\alpha \beta \gamma} A_{\mu}^{\beta} A_{\nu}^{\gamma} . \tag{6.1.29}
\end{equation*}
$$

Contrary to the abelian case, the field strength $F_{\mu \nu}$ is not gauge invariant. Its transformation properties can easily be found by considering the gauge transform of eq. (6.1.28):

$$
\begin{equation*}
\left[D_{\mu}, D_{\nu}\right] \psi \rightarrow U\left[D_{\mu}, D_{\nu}\right] \psi=-i U F_{\mu \nu} \psi=-i F_{\mu \nu}^{U} \psi^{U} \tag{6.1.30}
\end{equation*}
$$

from which we immediately get

$$
\begin{equation*}
F_{\mu \nu}^{U}=U F_{\mu \nu} U^{-1} \tag{6.1.31}
\end{equation*}
$$

In non-abelian gauge theories the gauge field strength transforms in the adjoint representation of the gauge group. The most general gauge-invariant Lagrangian can be written as a Lorentz invariant functional of matter fields and their covariant derivatives, and of the field strengths $F_{\mu \nu}$ and their covariant derivatives. At the level of dimension 4 (or less) operators, we have

$$
\begin{equation*}
\mathcal{L}=\mathcal{L}_{Y M}+\mathcal{L}_{M a t t e r} \tag{6.1.32}
\end{equation*}
$$

where

$$
\begin{align*}
\mathcal{L}_{Y M} & =-\frac{1}{2 g^{2}} \operatorname{Tr} F_{\mu \nu} F^{\mu \nu},  \tag{6.1.33}\\
\mathcal{L}_{\text {Matter }} & =\sum_{i} \bar{\psi}_{i}\left(\left\langle D>-m_{i}\right) \psi_{i}+\sum_{j}\left|D_{\mu} \phi_{j}\right|^{2}-V\left(\psi_{i}, \phi_{j}\right) .\right. \tag{6.1.34}
\end{align*}
$$

In eq. (6.1.33), we have introduced a dimensionless parameter $g$ that is identified as the gauge coupling constant of the non-abelian theory. By rescaling the gauge fields as

$$
\begin{equation*}
A_{\mu}^{\alpha} \rightarrow g A_{\mu}^{\alpha} \tag{6.1.35}
\end{equation*}
$$

we get canonical kinetic terms, having normalized the generators as

$$
\begin{equation*}
\operatorname{Tr} t^{\alpha} t^{\beta}=\frac{\delta^{\alpha \beta}}{2} \tag{6.1.36}
\end{equation*}
$$

in the fundamental representation. The coupling $g$ now appears in all covariant derivatives in the matter sector and in the self-coupling of the gauge fields in the Yang-Mills (YM) Lagrangian. Gauge invariance requires that the same coupling $g$ governs all these interactions. In eq. (6.1.34) $i$ and $j$ run over all fermions and scalars in the theory, $D_{\mu}$ are the covariant derivatives in the appropriate representations and $V$ encodes the scalar potential of the scalar fields $\phi_{j}$ and their Yukawa interactions with the fermions $\psi_{i}$. Gauge invariance requires that $\delta_{\epsilon} V=0$.

The carriers of the force associated to the non-abelian group, that we generally denote by "gluons", are themselves subject to the force they carry. The equations of motion (e.o.m.) for $A_{\mu}^{\alpha}$ deduced from the Lagrangian (6.1.32) are

$$
\begin{equation*}
\partial_{\mu} F_{\alpha}^{\mu \nu}=-g C_{\alpha \beta \gamma} A_{\mu}^{\beta} F_{\gamma}^{\mu \nu}-\frac{\delta \mathcal{L}_{\text {Matter }}}{\delta A_{\nu}^{\alpha}} \equiv-\mathcal{J}_{\alpha}^{\nu} \tag{6.1.37}
\end{equation*}
$$

where $\mathcal{J}_{\nu}^{\alpha}$ are the $\operatorname{dim} G$ conserved currents associated to the symmetry group $G$ :

$$
\begin{equation*}
\partial_{\mu} \mathcal{J}_{\alpha}^{\mu}=0 \tag{6.1.38}
\end{equation*}
$$

The e.o.m. (6.1.37), written in terms of $\mathcal{J}_{\mu}^{\alpha}$ are not covariant. It is convenient to shift the gluon contribution to the current, the first term in the second relation in eq. (6.1.37), to
the left-hand side of that equation. In doing so, the e.o.m. read

$$
\begin{equation*}
D_{\nu} F_{\alpha}^{\mu \nu}=-J_{\alpha}^{\mu} \tag{6.1.39}
\end{equation*}
$$

where

$$
\begin{align*}
D_{\rho} F_{\alpha}^{\mu \nu} & =\partial_{\rho} F_{\alpha}^{\mu \nu}-i g\left(t_{\beta}^{A d j}\right)_{\alpha \gamma} A_{\rho}^{\beta} F_{\gamma}^{\mu \nu} \\
J_{\alpha}^{\mu} & =\frac{\delta \mathcal{L}_{\text {Matter }}}{\delta A_{\mu}^{\alpha}}=\frac{\delta \mathcal{L}_{\text {Matter }}}{\delta D_{\mu} \Psi_{I}}\left(-i g t_{\alpha} \Psi_{I}\right) \tag{6.1.40}
\end{align*}
$$

In eq. (6.1.40) we have rewritten the form of the current $J_{\mu}^{\alpha}$ to make explicit its covariant properties, in contrast to to the conserved current $\mathcal{J}_{\mu}^{\alpha}$. The field $\Psi$ encodes both fermions and scalars and $I=(i, j)$. The current $J_{\mu}^{\alpha}$ is covariantly conserved, namely $D_{\mu} J_{\alpha}^{\mu}=0$. Indeed,

$$
\begin{equation*}
D_{\mu} J_{\alpha}^{\mu}=-D_{\nu} D_{\mu} F_{\alpha}^{\mu \nu}=\left[D_{\mu}, D_{\nu}\right] F_{\alpha}^{\mu \nu}=i F_{\nu \mu}^{\beta}\left(t_{\beta}^{A d j}\right)_{\alpha \gamma} F_{\gamma}^{\mu \nu}=C_{\alpha \beta \gamma} F_{\mu \nu}^{\beta} F_{\gamma}^{\mu \nu}=0 \tag{6.1.41}
\end{equation*}
$$

The field strength $F_{\mu \nu}$ satisfies another important relation, called Bianchi identity. It is a consistency relation, and can be derived starting from the Jacobi identity (6.1.9), applied to covariant derivatives:

$$
\begin{equation*}
\left[D_{\mu},\left[D_{\nu}, D_{\rho}\right]\right]+\left[D_{\rho},\left[D_{\mu}, D_{\nu}\right]\right]+\left[D_{\nu},\left[D_{\rho}, D_{\mu}\right]\right]=0 \tag{6.1.42}
\end{equation*}
$$

Using eq. (6.1.27), we can rewrite the above expression as

$$
\begin{equation*}
D_{\mu} F_{\nu \rho}+D_{\rho} F_{\mu \nu}+D_{\nu} F_{\rho \mu}=0 \tag{6.1.43}
\end{equation*}
$$

It is a straightforward exercise to show that eq. (6.1.43) is identically satisfied.

### 6.2 Quantum Treatment: the Faddeev-Popov Method

The quantization of gauge theories is non-trivial. The essential point is that Lorentz invariance forces us to describe helicity-one fields in terms of a four-vector field, but the latter has four components, and hence more degrees of freedom than necessary. The timecomponent of a vector field, in addition, is problematic, because it would lead to a kinetic term with an opposite sign with respect to its spatial components. Non-physical degrees of freedom are then expected. The way in which gauge theories solve the problem is to introduce a redundancy in the theory, gauge invariance, so that we can eliminate these extra unwanted and unphysical degrees of freedom.

Functional methods based on the path integral are by far the best way to quantize gauge theories. Before considering non-abelian theories, it is very useful to recall how

QED can be quantized using the path integral. Let's first see why a naive path integral quantization cannot work, by computing the photon propagator

$$
\begin{equation*}
\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle=\mathcal{N} \int \mathcal{D} A_{\mu} A_{\mu}(x) A_{\nu}(y) e^{i S(A)} \tag{6.2.1}
\end{equation*}
$$

where $\mathcal{N}$ is a normalization constant and

$$
\begin{equation*}
S(A)=-\frac{1}{4} \int d^{4} x F_{\mu \nu}^{2}=\frac{1}{2} \int \frac{d^{4} p}{(2 \pi)^{4}} A^{\mu}(-p) D_{\mu \nu}(p) A^{\nu}(p) \tag{6.2.2}
\end{equation*}
$$

is the usual free action, written both in configuration and momentum space. The tensor $D_{\mu \nu}$ equals

$$
\begin{equation*}
D_{\mu \nu}(p)=\left(-p^{2} \eta_{\mu \nu}+p_{\mu} p_{\nu}\right) \tag{6.2.3}
\end{equation*}
$$

The photon propagator is equal to the Fourier transform of the inverse of $D_{\mu \nu}(p)$ :

$$
\begin{equation*}
\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle=? \int \frac{d^{4} p}{(2 \pi)^{4}} i D_{\mu \nu}^{-1}(p) e^{i p \cdot(x-y)} \tag{6.2.4}
\end{equation*}
$$

However, $\operatorname{det} D_{\mu \nu}(p)=0$, no inverse exists and no propagator can be defined. Another way of looking at the problem is obtained by performing a shift of variables in the path integral: $A_{\mu}(x) \rightarrow A_{\mu}(x)+\partial_{\mu} \lambda(x)$. We get

$$
\begin{equation*}
\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle=\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle+\left(\partial_{\mu} \lambda \partial_{\nu} \lambda\right) \mathcal{N} \int \mathcal{D} A_{\mu} e^{i S(A)} \tag{6.2.5}
\end{equation*}
$$

whose only solution is $\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle=\infty$. The problem arises from the fact that the action is gauge invariant and we are integrating over all possible field configurations, including those that are related by a gauge transformation. All pure gauge configurations, such as $A_{\mu}(x)=\partial_{\mu} \lambda$, are not dumped by the action and lead to the above divergence. The problem is solved by restricting the integration to gauge inequivalent configurations only. In other words, we have to implement a gauge-fixing condition of the form $G(A)=0$, where $G(A)$ is a functional of the gauge field $A_{\mu}$. This can be imposed inside the path integral by means of a functional generalization of the Dirac delta-function. Care should be paid on possible Jacobian factors. These can arise by recalling the formula

$$
\begin{equation*}
\delta[f(x)]=\sum_{i}\left|\frac{d f}{d x}\right|_{x=x_{0}^{(i)}}^{-1} \delta\left(x-x_{0}^{(i)}\right) \tag{6.2.6}
\end{equation*}
$$

where $x_{0}^{(i)}$ are the values where the function $f$ vanishes. The $n$-dimensional integral generalization of eq. (6.2.6) is

$$
\begin{equation*}
1=\int \prod_{i} d x_{i} \delta^{(n)}\left(f_{j}\left(x_{i}\right)\right)\left|\operatorname{det} \frac{\partial f_{j}}{\partial x_{i}}\right| \tag{6.2.7}
\end{equation*}
$$

where $f_{j}$ are $n$ functions of the $n$ variables $x_{i}$ and we have assumed that they all vanish at a single point $x_{0, i}$. The further infinite dimensional generalization of eq. (6.2.7) is

$$
\begin{equation*}
1=\int \mathcal{D} \lambda \delta\left(G\left(A^{\lambda}\right)\right)\left|\operatorname{det} \frac{\delta G\left(A^{\lambda}\right)}{\delta \lambda(x)}\right| \tag{6.2.8}
\end{equation*}
$$

where $A_{\mu}^{\lambda}(x)=A_{\mu}(x)+\partial_{\mu} \lambda(x)$ and $G\left(A^{\lambda}\right)$ is an arbitrary functional, assumed to have a single function $\lambda_{0}$ where the functional vanishes. A simple choice for $G(A)$ is

$$
\begin{equation*}
G(A)=\partial_{\mu} A^{\mu} \Longrightarrow G\left(A^{\lambda}\right)=\partial_{\mu} A^{\mu}+\square \lambda \tag{6.2.9}
\end{equation*}
$$

Inserting eq. (6.2.8) inside the path integral gives

$$
\begin{align*}
& \mathcal{N} \int \mathcal{D} A_{\mu} e^{i S(A)} \int \mathcal{D} \lambda \delta\left(G\left(A^{\lambda}\right)\right)\left|\operatorname{det} \frac{\delta G\left(A^{\lambda}\right)}{\delta \lambda(x)}\right|=\mathcal{N}|\operatorname{det} \square| \int \mathcal{D} A_{\mu} e^{i S(A)} \int \mathcal{D} \lambda \delta\left(G\left(A^{\lambda}\right)\right) \\
& =\mathcal{N}^{\prime} \int \mathcal{D} \lambda \int \mathcal{D} A_{\mu} e^{i S(A)} \delta(G(A))=\mathcal{N}^{\prime \prime} \int \mathcal{D} A_{\mu} e^{i S(A)} \delta(G(A)) \tag{6.2.10}
\end{align*}
$$

where we have included into the normalization constants the gauge-field independent factor $\operatorname{det} \square$ and the integration over the gauge parameter $\lambda$. The functional delta into the last term of eq. (6.2.10) avoids to integrate over redundant field configurations. With a simple trick, we can also get rid of the functional delta and yet have a well-defined path integral. Instead of taking a single gauge fixing like the one in eq. (6.2.9) we can introduce a family of gauge fixing terms, parametrized by an arbitrary function $f$ :

$$
\begin{equation*}
G_{f}(A)=\partial_{\mu} A^{\mu}-f \tag{6.2.11}
\end{equation*}
$$

Since no physical observable can depend on the gauge fixing, we can average over the different gauge fixings by introducing a phase factor

$$
\begin{equation*}
\exp \left(-\frac{i}{2 \xi} \int d^{4} x f^{2}(x)\right) \tag{6.2.12}
\end{equation*}
$$

and integrating over $f(x)$. This is useful, since in so doing we can get rid in one step of the functional delta and of the integration over $f(x)$. In eq. (6.2.12) $\xi$ is a positive parameter. The path integral becomes now

$$
\begin{equation*}
\mathcal{N} \int \mathcal{D} A_{\mu} e^{i S(A)-\frac{i}{2 \xi} \int d^{4} x\left(\partial_{\mu} A^{\mu}\right)^{2}} \tag{6.2.13}
\end{equation*}
$$

The final outcome of these manipulations is the addition of a new gauge-variant term in the action, called gauge-fixing term. The latter is crucial to make sense of the photon propagator. Equation (6.2.1) is replaced by

$$
\begin{equation*}
\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle=\mathcal{N} \int \mathcal{D} A_{\mu} A_{\mu}(x) A_{\nu}(y) e^{i S(A)-\frac{i}{2 \xi} \int d^{4} x\left(\partial_{\mu} A^{\mu}\right)^{2}} \tag{6.2.14}
\end{equation*}
$$

The tensor (6.2.3) becomes

$$
\begin{equation*}
D_{\mu \nu}(p)=-p^{2} \eta_{\mu \nu}+p_{\mu} p_{\nu} \frac{(\xi-1)}{\xi} \tag{6.2.15}
\end{equation*}
$$

and it admits the inverse

$$
\begin{equation*}
D_{\mu \nu}^{-1}(p)=-\frac{1}{p^{2}}\left(\eta_{\mu \nu}-(1-\xi) \frac{p_{\mu} p_{\nu}}{p^{2}}\right) . \tag{6.2.16}
\end{equation*}
$$

Reintroducing the $i \epsilon$, we finally get

$$
\begin{equation*}
\left\langle A_{\mu}(x) A_{\nu}(y)\right\rangle=\int \frac{d^{4} p}{(2 \pi)^{4}} \frac{-i}{p^{2}+i \epsilon}\left(\eta_{\mu \nu}-(1-\xi) \frac{p_{\mu} p_{\nu}}{p^{2}}\right) e^{i p \cdot(x-y)} \tag{6.2.17}
\end{equation*}
$$

The photon propagator is not a direct physical observable and depends on $\xi$. Common choices for $\xi$ in explicit computations are $\xi=0$ (Landau gauge), in which case the tree-level propagator is transverse and $\xi=1$ (Feynman gauge), in which this simplifies considerably.

The quantization of non-abelian gauge theories proceeds along the same way as the abelian case, but it presents additional complications. The naive measure is

$$
\begin{equation*}
\mathcal{N} \int \mathcal{D} A_{\mu}^{\alpha} e^{i S(A)} \tag{6.2.18}
\end{equation*}
$$

It is gauge-invariant, but in a less trivial way than in abelian theories, since a gauge transformation rotates the fields. The Jacobian associated to the infinitesimal transformation (6.1.18) is

$$
\begin{equation*}
\operatorname{Jac}_{\mu \nu}^{\alpha \beta}(x, y)=\frac{\delta A_{\mu}^{\prime \alpha}(x)}{\delta A_{\nu}^{\beta}(y)}=\delta_{\mu}^{\nu} \delta(x-y)\left(\delta_{\alpha \beta}+\epsilon^{\gamma} C_{\alpha \beta \gamma}\right) . \tag{6.2.19}
\end{equation*}
$$

Since $\operatorname{Det}(1+\epsilon)=1+\operatorname{Tr} \epsilon+\mathcal{O}\left(\epsilon^{2}\right)$, we have

$$
\begin{equation*}
\operatorname{det} \operatorname{Jac}_{\mu \nu}^{\alpha \beta}(x, y)=\delta_{\mu}^{\mu} \delta(0)\left(\delta_{\alpha \alpha}+\epsilon^{\gamma} C_{\alpha \alpha \gamma}\right)=\delta_{\mu}^{\mu} \delta(0) \delta_{\alpha \alpha} \tag{6.2.20}
\end{equation*}
$$

that is the infinite dimensional generalization of the unit matrix. We then conclude that the measure $\mathcal{D} A_{\mu}^{\alpha}$ is gauge invariant. Let us now proceed like in the abelian case, introducing a delta-functional gauge-fixing in the path integral, like in eq. (6.2.8). Let us define

$$
\begin{equation*}
\Delta_{G}^{-1}(A)=\int \mathcal{D} U \delta\left(G\left(A^{U}(x)\right)\right) \tag{6.2.21}
\end{equation*}
$$

In eq. (6.2.21), $\mathcal{D} U$ is the so-called invariant measure of the group $G$, parametrized by group elements $U(\Lambda)$. It is not, like in the abelian case, simply the integration over the Lie algebra generators, $\prod_{\alpha} \mathcal{D} \Lambda^{\alpha}(x)$, but it includes a non trivial measure $\rho(\Lambda)$. It is called invariant measure because it satisfies the following properties:

$$
\begin{equation*}
\int \mathcal{D} U f(U)=\int \mathcal{D} U f\left(U^{-1}\right)=\int \mathcal{D} U f\left(U \cdot U_{0}\right)=\int \mathcal{D} U f\left(U_{0} \cdot U\right) \tag{6.2.22}
\end{equation*}
$$

where $U_{0}$ is a constant element of the group and $f(U)$ is an arbitrary function over the group.
** Although we will never need its explicit form in the following, it is worth to spend a few more words on $\rho(\Lambda)$. This can be defined starting from the metric in group space $g_{\alpha \beta}$ as

$$
\begin{equation*}
g_{\alpha \beta}=\operatorname{Tr}\left(U^{-1}(\Lambda)\left(\frac{\partial U(\Lambda)}{\partial \Lambda_{\alpha}}\right) U^{-1}(\Lambda)\left(\frac{\partial U(\Lambda)}{\partial \Lambda_{\beta}}\right)\right) . \tag{6.2.23}
\end{equation*}
$$

The invariant measure is

$$
\begin{equation*}
\rho(\Lambda)=\sqrt{\operatorname{det} g_{\alpha \beta}} . \tag{6.2.24}
\end{equation*}
$$

The parameters $\Lambda_{\alpha}$ in eq. (6.2.23) are space-time independent coordinates spanning the group $G$. They do not necessarily correspond to the $x$-independent version of the generators appearing in eq. (6.1.23). An example will clarify this point. The group $S U(2)$ is defined as the set of $2 \times 2$ unitary matrices $U$ of unit determinant. Any matrix $U$ can be written as

$$
U=\left(\begin{array}{cc}
z_{1} & z_{2}  \tag{6.2.25}\\
-z_{2}^{*} & z_{1}^{*}
\end{array}\right)
$$

with $\left|z_{1}\right|^{2}+\left|z_{2}\right|^{2}=1, z_{1,2} \in \mathbb{C}$. The group $S U(2)$ is isomorphic to the three-dimensional sphere $S^{3}$. Instead of using standard coordinates subject to a constraint, we might use radial coordinates. Denoting $z_{i}=x_{i}+i y_{i}, i=1,2$, we have

$$
\begin{array}{ll}
x_{1}=\sin \psi \sin \theta \cos \phi, & y_{1}=\sin \psi \sin \theta \sin \phi  \tag{6.2.26}\\
x_{2}=\sin \psi \cos \theta, & y_{1}=\cos \psi,
\end{array}
$$

with $0 \leq \phi<2 \pi, 0 \leq \theta<2 \pi, 0 \leq \psi \leq \pi$. The invariant $S U(2)$ measure coincides with the standard metric of $S^{3}: \rho(\psi, \theta, \phi)=\sin ^{2} \psi \sin \theta .{ }^{2}$ Given the explicit form of the metric (6.2.23) it is straightforward to prove the relations (6.2.22).

The functional $\Delta_{G}(A)$ is gauge-invariant: $\Delta_{G}\left(A^{U}\right)=\Delta_{G}(A)$, that immediately follows from eq. (6.2.22). The naive measure (6.2.18) can be rewritten as (omitting Lorentz and color indices in the gauge measure)

$$
\begin{equation*}
\mathcal{N} \int \mathcal{D} A \int \mathcal{D} U e^{i S(A)} \Delta_{G}(A) \delta\left(G\left(A^{U}\right)\right) \tag{6.2.27}
\end{equation*}
$$

We can change variables in the path integral by defining $A_{\mu}=A_{\mu}^{\prime U^{-1}}$. Since $\Delta_{G}(A)$, the measure and the action are all gauge invariant, we get (redefine $A^{\prime} \rightarrow A$ )

$$
\begin{equation*}
\mathcal{N} \int \mathcal{D} A \int \mathcal{D} U e^{i S(A)} \Delta_{G}(A) \delta(G(A))=\mathcal{N}^{\prime} \int \mathcal{D} A e^{i S(A)} \Delta_{G}(A) \delta(G(A)) \tag{6.2.28}
\end{equation*}
$$

[^36]and we can reabsorb the invariant group measure into the overall path integral normalization $\mathcal{N}$. Let us verify that any correlation function of gauge invariant operators does not depend on the gauge-fixing, namely does not depend on the specific choice of the functional $G(A)$. If $O(A)$ schematically represents some product of gauge invariant operators, given two arbitrary gauge fixing functionals $G$ and $F$, one has
\[

$$
\begin{align*}
\langle O(A)\rangle_{G} & =\mathcal{N} \int \mathcal{D} A e^{i S(A)} \Delta_{G}(A) \delta(G(A)) O(A)  \tag{6.2.29}\\
& =\mathcal{N} \int \mathcal{D} A e^{i S(A)} \Delta_{G}(A) \delta(G(A)) \int \mathcal{D} U \delta\left(F\left(A^{U}\right)\right) \Delta_{F}(A) O(A) \\
& =\mathcal{N} \int \mathcal{D} A \int \mathcal{D} U e^{i S(A)} \Delta_{G}(A) \delta\left(G\left(A^{U-1}\right)\right) \delta(F(A)) \Delta_{F}(A) O(A) \\
& =\mathcal{N} \int \mathcal{D} A \int \mathcal{D} U e^{i S(A)} \Delta_{G}(A) \delta\left(G\left(A^{U}\right)\right) \delta(F(A)) \Delta_{F}(A) O(A) \\
& =\mathcal{N} \int \mathcal{D} A e^{i S(A)} \delta(F(A)) \Delta_{F}(A) O(A)=\langle O(A)\rangle_{F}
\end{align*}
$$
\]

On the contrary, the correlation functions of gauge dependent quantities do depend on the choice of $G(A)$. Similarly to the abelian case, we can also take a family of gauge fixing functionals and integrate over them. The typical choice will be the non-abelian generalization of eq. (6.2.11):

$$
\begin{equation*}
G_{f}(A)=\partial_{\mu} A_{\alpha}^{\mu}-f_{\alpha} \tag{6.2.30}
\end{equation*}
$$

weighted by the phase factor

$$
\begin{equation*}
\exp \left(-\frac{i}{2 \xi} \int d^{4} x f_{\alpha}^{2}(x)\right) \tag{6.2.31}
\end{equation*}
$$

In this way, we get

$$
\begin{equation*}
\langle O(A)\rangle=\mathcal{N} \int \mathcal{D} A e^{i S(A)-\frac{i}{2 \xi} \int d^{4} x\left(\partial_{\mu} A_{\alpha}^{\mu}\right)^{2}} \Delta(A) O(A) \tag{6.2.32}
\end{equation*}
$$

where

$$
\begin{equation*}
\Delta^{-1}(A)=\int \mathcal{D} U \delta\left(\partial_{\mu}\left(A^{U}\right)_{\alpha}^{\mu}-\partial_{\mu} A_{\alpha}^{\mu}\right) \tag{6.2.33}
\end{equation*}
$$

The integrand in eq. (6.2.33) has a non-trivial support for values of $U$ infinitesimally close to the identity: $U=1+i \epsilon^{\alpha} t_{\alpha}$. We have

$$
\begin{equation*}
\delta\left(\partial_{\mu}\left(A^{U}\right)_{\alpha}^{\mu}-\partial_{\mu} A_{\alpha}^{\mu}\right)=\delta\left(\partial_{\mu} D^{\mu} \epsilon_{\alpha}\right)=\frac{\delta\left(\epsilon_{\alpha}\right)}{\left|\operatorname{det} \partial_{\mu} D^{\mu}\right|} . \tag{6.2.34}
\end{equation*}
$$

Modulo irrelevant constants, as usual absorbable in the path integral normalization $\mathcal{N}$, we get

$$
\begin{equation*}
\Delta(A)=\left|\operatorname{det} \partial_{\mu} D^{\mu}\right| \tag{6.2.35}
\end{equation*}
$$

This is nothing else than the non-abelian generalization of the |det $\square \mid$ term appearing in eq. (6.2.10). The crucial difference with the abelian case is its gauge field dependence by means of the covariant derivative. For this reason it should be kept inside the path integral. We can remove the absolute value, that would make the computation of $\Delta(A)$ quite complicated, because, loosely speaking,

$$
\begin{equation*}
\left|\operatorname{det} \partial_{\mu} D^{\mu}\right|=|\operatorname{det} \square||\operatorname{det}(1+\mathcal{O}(g))| . \tag{6.2.36}
\end{equation*}
$$

The first factor is irrelevant, while the second one is manifestly positive in perturbation theory. We can then remove the absolute value and write

$$
\begin{equation*}
\Delta(A)=\operatorname{det} \partial_{\mu} D^{\mu} \tag{6.2.37}
\end{equation*}
$$

It is not easy to directly compute determinants. It is more convenient to turn a determinant into a local action by inserting additional non-physical degrees of freedom and use the identity, valid for Grassmann variables $\omega$ and $\omega^{*},{ }^{3}$

$$
\begin{equation*}
\int \mathcal{D} \omega \mathcal{D} \omega^{*} e^{-i \int d^{4} x \omega^{*}(x) F(x) \omega(x)} \propto \operatorname{det} F \tag{6.2.38}
\end{equation*}
$$

where $F(x)$ is an arbitrary differential operator. Using eq. (6.2.38), we can write, modulo irrelevant constants,

$$
\begin{equation*}
\operatorname{det} \partial_{\mu} D^{\mu}=\int \mathcal{D} \omega_{\alpha} \mathcal{D} \omega_{\alpha}^{*} e^{i S_{g h o s t}} \tag{6.2.39}
\end{equation*}
$$

where

$$
\begin{equation*}
S_{g h o s t}=\int d^{4} x \partial_{\mu} \omega_{\alpha}^{*} D^{\mu} \omega_{\alpha}=\int d^{4} x \partial_{\mu} \omega_{\alpha}^{*}\left(\partial^{\mu} \omega_{\alpha}+g C_{\alpha \beta \gamma} A_{\beta}^{\mu} \omega_{\gamma}\right) \tag{6.2.40}
\end{equation*}
$$

$\omega_{\alpha}$ and $\omega_{\alpha}^{*}$ are dim-G scalar fields with fermion statistic, transforming in the adjoint representation of $G$. They are not associated to physical propagating particles and for this reason they are denoted ghost fields. They cannot appear as physical external states, but they can and do appear in loops as virtual particles, by means of their interaction with the gauge fields. Ghosts are crucial to restore unitarity in non-abelian gauge theories. Loosely speaking, they compensate for the (also unphysical) contribution of the longitudinal and time component of the gauge fields that, contrary to the abelian case, do not automatically decouple from scattering amplitudes.

Putting all together, the complete non-abelian Lagrangian density at the quantum level is the sum of three terms:

$$
\begin{equation*}
\mathcal{L}_{t o t}=\mathcal{L}+\mathcal{L}_{\text {g.f. }}+\mathcal{L}_{\text {ghost }}, \tag{6.2.41}
\end{equation*}
$$

[^37]where
\[

$$
\begin{align*}
\mathcal{L} & =-\frac{1}{4} F_{\mu \nu}^{\alpha} F_{\alpha}^{\mu \nu}+\sum_{i} \bar{\psi}_{i}\left(\left\langle\partial D-m_{i}\right) \psi_{i}+\sum_{j}\left|D_{\mu} \phi_{j}\right|^{2}-V\left(\psi_{i}, \phi_{j}\right)\right.  \tag{6.2.42}\\
\mathcal{L}_{\text {g.f. }} & =-\frac{1}{2 \xi}\left(\partial_{\mu} A_{\alpha}^{\mu}\right)^{2}  \tag{6.2.43}\\
\mathcal{L}_{\text {ghost }} & =\partial_{\mu} \omega_{\alpha}^{*} D^{\mu} \omega_{\alpha} . \tag{6.2.44}
\end{align*}
$$
\]

The propagator of gauge and ghost field is readily found by the quadratic term of the above Lagrangian density. We have

$$
\begin{align*}
\left\langle A_{\mu}^{\alpha}(x) A_{\nu}^{\beta}(y)\right\rangle & =\delta^{\alpha \beta} \int \frac{d^{4} p}{(2 \pi)^{4}} \frac{-i}{p^{2}+i \epsilon}\left(\eta_{\mu \nu}-(1-\xi) \frac{p_{\mu} p_{\nu}}{p^{2}}\right) e^{i p \cdot(x-y)}  \tag{6.2.45}\\
\left\langle\omega_{\alpha}^{*}(x) \omega_{\beta}(y)\right\rangle & =\delta_{\alpha \beta} \int \frac{d^{4} p}{(2 \pi)^{4}} \frac{i}{p^{2}+i \epsilon} e^{i p \cdot(x-y)}
\end{align*}
$$

The gauge field propagator is the trivial generalization of the photon propagator (6.2.16), while the ghost propagator coincides with that of a complex massless scalar field. Recall that the ghost Lagrangian (6.2.44) depends on the specific form of $\Delta(A)$, that in turn depends on the specific gauge fixing chosen. It is invariant under a $U(1)$ symmetry (ghost number) under which the ghost fields $\omega_{\alpha}$ and $\omega_{\alpha}^{*}$ have respectively charges +1 and -1 . For this reason the latter fields are commonly denoted anti-ghosts and this explains the notation $\omega_{\alpha}^{*}$. All the operators appearing in the Lagrangian (6.2.41) have dimensions less or equal to four, compatibly with a renormalizable theory. However, not all possible operators of dimensions $\Delta \leq 4$ appear in $\mathcal{L}_{\text {tot }}$ and hence the renormalizability of these theories is not obvious. The appearance of unphysical fields in the theory complicates also the notion of physical field in non-abelian gauge theories. All these issues are best addressed by introducing the BRST symmetry, subject of the next section.

Before going on we would like to comment on the assumption made below eq.(6.2.8), that was implicit in writing eq.(6.2.21), about the existence, at a given $A_{\mu}(x)$, of only one gauge element $U_{0}(x)$ where the functional $G\left(A^{U}\right)$ vanishes. This is equivalent to say that the equation

$$
\begin{equation*}
\partial^{\mu}\left(U^{\dagger}\left(\partial_{\mu} U-i\left[A_{\mu}, U\right]\right)\right)=0 \tag{6.2.46}
\end{equation*}
$$

has either no solutions or such solutions should be discarded in the path integral. In the abelian case, eq.(6.2.46) boils down to $\square \lambda=0$ that clearly admits plane wave solutions. These can be ruled out by demanding proper boundary conditions for $A_{\mu}$ in the path integral. This is particularly clear in the euclidean case, where we demand $A_{\mu}$ to vanish at infinity. The conditions $\lambda \rightarrow 0$ at infinity and $\square \lambda=0$ in euclidean space imply $\lambda=0$. In contrast, in the non-abelian case there exists in general non-trivial solutions of the
above equation that would force us to consider other gauge field configurations $A^{U}$, called Gribov copies. Their possible presence is important in a non-perturbative treatment of gauge theories, but will not affect the perturbative expansion around the free theory (like the absolute value in eq.(6.2.36)). The above assumption can then be justified only in a perturbative set-up.

### 6.3 BRST Symmetry

The Faddeev-Popov path integral quantization of gauge theories reviewed above requires to fix a gauge and hides the underlying gauge invariance of the theory. ${ }^{4}$ In other words, the Lagrangian $\mathcal{L}_{\text {tot }}$ cannot obviously be gauge invariant. On the other hand, it was found by Becchi, Rouet and Stora, and indepedendently by Tyutin, that $\mathcal{L}_{\text {tot }}$ is invariant under an additional symmetry, called BRST symmetry. It is useful to rewrite the gauge fixing term in a different fashion by "integrating in" an auxiliary field $H_{\alpha}(x)$ :

$$
\begin{equation*}
e^{-\frac{i}{2 \xi} \int d^{4} x f_{\alpha}^{2}(x)}=\int \mathcal{D} H_{\alpha} e^{\frac{i \xi}{2} \int d^{4} x H_{\alpha}^{2}-i \int d^{4} x f_{\alpha} H_{\alpha}}, \tag{6.3.1}
\end{equation*}
$$

so that

$$
\begin{equation*}
\mathcal{L}_{\text {tot }}=\mathcal{L}+\mathcal{L}_{\text {ghost }}+\frac{\xi}{2} H_{\alpha}^{2}-f_{\alpha} H_{\alpha} \tag{6.3.2}
\end{equation*}
$$

The total Lagrangian (6.3.2) is invariant under infinitesimal transformations parametrized by an anticommuting variable $\theta$ :

$$
\begin{align*}
\delta_{\theta} \Psi & =i t_{\alpha} \theta \omega_{\alpha} \Psi \\
\delta_{\theta} A_{\mu} & =\theta D_{\mu} \omega_{\alpha} \\
\delta_{\theta} \omega_{\alpha} & =-\frac{1}{2} \theta C_{\alpha \beta \gamma} \omega_{\beta} \omega_{\gamma}  \tag{6.3.3}\\
\delta_{\theta} \omega_{\alpha}^{*} & =-\theta H_{\alpha} \\
\delta_{\theta} H_{\alpha} & =0
\end{align*}
$$

These are the BRST transformations, written in the non-canonical basis for the gauge fields where the gauge coupling does not appear in the interactions. The field $\Psi$ represents any matter field, fermionic or bosonic. The BRST transformations are nilpotent, namely if we denote $\delta_{\theta} F \equiv \theta s F$, then $\delta_{\theta}(s F)=\theta s^{2} F=0$ for any functional of the fields $\Psi, A_{\mu}, \omega_{\alpha}, \omega_{\alpha}^{*}$ and $H_{\alpha}$. Let us check that $s^{2}$ vanishes when acting on single fields. From eq. (6.3.3) we have $s \Psi=i t_{\alpha} \omega_{\alpha} \Psi_{\alpha}$. Then

$$
\begin{align*}
\delta_{\theta}(s \Psi) & =i t_{\alpha}\left(\left(\delta_{\theta} \omega_{\alpha}\right) \Psi+\omega_{\alpha}\left(\delta_{\theta} \Psi\right)\right)=i t_{\alpha}\left(-\frac{1}{2} \theta C_{\alpha \beta \gamma} \omega_{\beta} \omega_{\gamma} \Psi+\omega_{\alpha} i t_{\beta} \theta \omega_{\beta} \Psi\right) \\
& =\theta \omega_{\alpha} \omega_{\beta} \Psi\left(-\frac{i}{2} C_{\alpha \beta \gamma} t_{\gamma}+\frac{1}{2}\left[t_{\alpha}, t_{\beta}\right]\right)=0 . \tag{6.3.4}
\end{align*}
$$

[^38]Consider now the BRST transformations of the gauge fields. We have $s A_{\mu \alpha}=D_{\mu} \omega_{\alpha}=$ $\partial_{\mu} \omega_{\alpha}+C_{\alpha \beta \gamma} A_{\mu}^{\beta} \omega_{\gamma}$. Then
$\delta_{\theta}\left(s A_{\mu}^{\alpha}\right)=\partial_{\mu}\left(-\frac{1}{2} \theta C_{\alpha \beta \gamma} \omega_{\beta} \omega_{\gamma}\right)+C_{\alpha \beta \gamma} \theta\left(\partial_{\mu} \omega_{\beta}+C_{\beta \rho \sigma} A_{\mu}^{\rho} \omega_{\sigma}\right) \omega_{\gamma}+C_{\alpha \beta \gamma} A_{\mu}^{\beta}\left(-\frac{1}{2} \theta C_{\gamma \rho \sigma} \omega_{\rho} \omega_{\sigma}\right)$.
Consider first the terms proportional to $A_{\mu}$ in eq. (6.3.5). Reshuffling indices, one has

$$
\begin{align*}
\left.\delta_{\theta}\left(s A_{\mu}^{\alpha}\right)\right|_{A} & =\theta A_{\mu}^{\beta} \omega_{\rho} \omega_{\sigma}\left(-\frac{1}{2} C_{\alpha \beta \gamma} C_{\gamma \rho \sigma}+C_{\alpha \rho \gamma} C_{\gamma \beta \sigma}\right)  \tag{6.3.6}\\
& =-\frac{1}{2} \theta A_{\mu}^{\beta} \omega_{\rho} \omega_{\sigma}\left(C_{\alpha \beta \gamma} C_{\gamma \rho \sigma}+C_{\rho \alpha \gamma} C_{\gamma \beta \sigma}+C_{\beta \rho \gamma} C_{\gamma \alpha \sigma}\right)=0
\end{align*}
$$

where the last equality follows from the Jacobi identity (6.1.10). The terms with no gauge fields give

$$
\begin{equation*}
\left.\delta_{\theta}\left(s A_{\mu}^{\alpha}\right)\right|_{n o A}=\theta C_{\alpha \beta \gamma}\left(-\frac{1}{2}\left(\partial_{\mu} \omega_{\beta}\right) \omega_{\gamma}-\frac{1}{2} \omega_{\beta}\left(\partial_{\mu} \omega_{\gamma}\right)+\left(\partial_{\mu} \omega_{\beta}\right) \omega_{\gamma}\right)=0 . \tag{6.3.7}
\end{equation*}
$$

Similarly, we have

$$
\begin{align*}
\delta_{\theta}\left(s \omega_{\alpha}\right) & =-\frac{1}{2} C_{\alpha \beta \gamma}\left(\left(-\frac{1}{2} \theta C_{\beta \rho \sigma} \omega_{\rho} \omega_{\sigma}\right) \omega_{\gamma}+\omega_{\beta}\left(-\frac{1}{2} \theta C_{\gamma \rho \sigma} \omega_{\rho} \omega_{\sigma}\right)\right) \\
& =\frac{1}{2} \theta \omega_{\rho} \omega_{\sigma} \omega_{\beta} C_{\rho \sigma \gamma} C_{\gamma \beta \alpha}=0, \tag{6.3.8}
\end{align*}
$$

where again the last equality follows from the Jacobi identity (6.1.10) and the antisymmetrization in the indices $\rho, \sigma$ and $\beta$. The nilpotency of the BRST transformations acting on $\omega_{\alpha}^{*}$ and $H_{\alpha}$ is trivial. We immediately have, from eq. (6.3.3),

$$
\begin{equation*}
\delta_{\theta}\left(s \omega_{\alpha}^{*}\right)=\delta_{\theta} s H_{\alpha}=0 \tag{6.3.9}
\end{equation*}
$$

Summarizing, we have shown that for any field $\Phi=A, \Psi, \omega, \omega^{*}, H$,

$$
\begin{equation*}
s^{2} \Phi=0 \tag{6.3.10}
\end{equation*}
$$

For two fields, we have

$$
\begin{equation*}
\delta_{\theta}\left(\Phi_{1} \Phi_{2}\right)=\theta\left(s \Phi_{1}\right) \Phi_{2}+\Phi_{1}\left(\theta s \Phi_{2}\right)=\theta\left(\left(s \Phi_{1}\right) \Phi_{2} \pm \Phi_{1}\left(s \Phi_{2}\right)\right) \tag{6.3.11}
\end{equation*}
$$

with + or - depending whether the field $\Phi_{1}$ is bosonic or fermionic. ${ }^{5}$ Hence

$$
\begin{equation*}
s\left(\Phi_{1} \Phi_{2}\right)=\left(\left(s \Phi_{1}\right) \Phi_{2} \pm \Phi_{1}\left(s \Phi_{2}\right)\right) \tag{6.3.12}
\end{equation*}
$$

Acting again with a BRST transformation gives

$$
\begin{align*}
\delta_{\theta}\left(s \Phi_{1} \Phi_{2}\right) & =\delta_{\theta}\left(s \Phi_{1}\right) \Phi_{2}+s \Phi_{1} \delta_{\theta} \Phi_{2} \pm\left(\delta_{\theta} \Phi_{1}\right) s \Phi_{2} \pm \Phi_{1} \delta_{\theta}\left(s \Phi_{2}\right) \\
& =\theta\left(\left(s^{2} \Phi_{1}\right) \Phi_{2} \mp\left(s \Phi_{1}\right)\left(s \Phi_{2}\right) \pm\left(s \Phi_{1}\right)\left(s \Phi_{2}\right)+\Phi_{1}\left(s^{2} \Phi_{2}\right)\right)=0 \tag{6.3.13}
\end{align*}
$$

[^39]if eq. (6.3.10) is satisfied, for $\Phi_{1}$ both bosonic and fermionic. Iterating the argument for more fields, we conclude that for any functional $F(\Phi)$,
\[

$$
\begin{equation*}
s^{2} F(\Phi)=0 \tag{6.3.14}
\end{equation*}
$$

\]

We now proceed to prove that the Lagrangian (6.3.2) is BRST-invariant. On the physical fields $\Psi$ and $A_{\mu}^{\alpha}$, the BRST transformations (6.3.3) can be seen as infinitesimal gauge transformations with parameter $\epsilon_{\alpha}(x)=\theta \omega_{\alpha}(x)$. The BRST-invariance of the gauge and matter Lagrangian term $\mathcal{L}$ immediately follows from the fact that $\mathcal{L}$ is gauge invariant. Let us now turn to the remaining three terms in eq. (6.3.2). We have

$$
\begin{equation*}
\delta_{\theta} f_{\alpha}=\delta_{\theta} \partial^{\mu} A_{\mu}^{\alpha}=\theta \partial_{\mu} D^{\mu} \omega_{\alpha}=\theta s f_{\alpha} \tag{6.3.15}
\end{equation*}
$$

Modulo total derivatives, we then get

$$
\begin{equation*}
\mathcal{L}_{\text {ghost }}=-\omega_{\alpha}^{*} s f_{\alpha} . \tag{6.3.16}
\end{equation*}
$$

Since $s \omega_{\alpha}^{*}=-H_{\alpha}$, the last three terms in eq. (6.3.2) can be rewritten as

$$
\begin{equation*}
\mathcal{L}_{\text {ghost }}+\frac{\xi}{2} H_{\alpha}^{2}-f_{\alpha} H_{\alpha}=s\left(f_{\alpha} \omega_{\alpha}^{*}-\frac{1}{2} \xi \omega_{\alpha}^{*} H_{\alpha}\right) . \tag{6.3.17}
\end{equation*}
$$

In this way, the BRST invariance of these terms is automatically ensured by the fact that $s^{2}=0$ for any functional. We conclude that the whole Lagrangian $\mathcal{L}_{t o t}$ is invariant under BRST transformations.

From an operatorial point of view, the BRST transformations (6.3.3) are generated by a Grassmann Hermitian operator $Q$. For any field $\Phi$, we have

$$
\begin{equation*}
\delta_{\theta} \Phi=i[\theta Q, \Phi]=i \theta[Q, \Phi]_{\mp}, \tag{6.3.18}
\end{equation*}
$$

where - and + denote commutator and anti-commutator, respectively, depending on whether the field $\Phi$ is bosonic or fermionic. We then have

$$
\begin{equation*}
[Q, \Phi]_{\mp}=-i s \Phi . \tag{6.3.19}
\end{equation*}
$$

The nilpotency of $s, s^{2}=0$, is equivalent to

$$
\begin{equation*}
(-i s)^{2} \Phi=\left[Q,[Q, \Phi]_{\mp}\right]_{ \pm}=\left[Q^{2}, \Phi\right]_{-}=0 \Longrightarrow Q^{2}=0 . \tag{6.3.20}
\end{equation*}
$$

The BRST operator $Q$ allows us to make the following partition of the Hilbert space in non-abelian gauge theories. Any state $|\phi\rangle$ in the Hilbert space falls in one of the following three categories:

$$
\begin{align*}
& Q\left|\phi_{1}\right\rangle \neq 0 \\
& Q\left|\phi_{2}\right\rangle=0, \quad \text { with }\left|\phi_{2}\right\rangle=Q\left|\phi_{1}\right\rangle  \tag{6.3.21}\\
& Q\left|\phi_{3}\right\rangle=0, \quad \text { but }\left|\phi_{3}\right\rangle \neq Q\left|\phi_{1}\right\rangle .
\end{align*}
$$

The fields in the second class are manifestly unphysical, since they have vanishing norm:

$$
\begin{equation*}
\left.\| \phi_{2}\right\rangle\left.\right|^{2}=\left\langle\phi_{2} \mid \phi_{2}\right\rangle=\left\langle\phi_{1}\right| Q^{2}\left|\phi_{1}\right\rangle=0 . \tag{6.3.22}
\end{equation*}
$$

We now show that gauge invariance implies that physical states should be annihilated by $Q$. More precisely, matrix elements between physical states $|\alpha\rangle$ and $|\beta\rangle$ should not depend on the choice of gauge fixing term. We have just seen that the total Lagrangian $\mathcal{L}_{\text {tot }}$ can be written as a physical gauge invariant term $\mathcal{L}$ plus a BRST variation of some functional $F: \mathcal{L}_{\text {tot }}=\mathcal{L}+s F(\Phi)$. The specific form of $F$ depends on the gauge-fixing chosen. If we infinitesimally deform the gauge fixing, the functional $F$ will also be deformed $F \rightarrow F+\delta F$. Demanding that

$$
\begin{equation*}
\langle\alpha \mid \beta\rangle_{F}=\langle\alpha \mid \beta\rangle_{F+\delta F} \tag{6.3.23}
\end{equation*}
$$

is equivalent to the condition

$$
\begin{equation*}
\langle\alpha|[Q, \delta F]_{+}|\beta\rangle_{F}=0, \quad \forall \alpha, \beta \in \text { physical } \tag{6.3.24}
\end{equation*}
$$

and for any sensible choice of gauge-fixing functional $\delta F$. We then conclude that $Q$ should annihilate physical states. These are then identified with the states in the third category in eq. (6.3.21). States that are annihilated by the operator $Q\left(\left|\phi_{3}\right\rangle\right)$ are said to be in the kernel of $Q$. The states $\left|\phi_{2}\right\rangle$ are said to be in the image of $Q$. The physical states are states in the kernel that are not in the image of $Q$. Such states are said to be in the cohomology of $Q$. It is clear that physical states are not uniquely defined. If $|\alpha\rangle$ is a given physical state, then any state of the form

$$
\begin{equation*}
|\widetilde{\alpha}\rangle=|\alpha\rangle+Q\left|\phi_{1}\right\rangle \tag{6.3.25}
\end{equation*}
$$

defines the same physical state, since

$$
\begin{equation*}
\langle\widetilde{\alpha} \mid \beta\rangle=\langle\alpha \mid \beta\rangle \tag{6.3.26}
\end{equation*}
$$

for any other physical state $|\beta\rangle$. Physical states $|p h y s\rangle$ correspond to equivalence classes within the class $\left|\phi_{3}\right\rangle$. This complicated structure of the Hilbert space in non-abelian gauge theories is a consequence of the redundancy introduced by the gauge symmetries. Like ghosts, any unphysical state, even if not present as an external line in a scattering process, can contribute as a virtual particle in loops. On the other hand, the optical theorem relates the imaginary part of a loop diagram with the square of scattering diagrams where the virtual particles become external on-shell states. Unitarity is then not obvious. BRST invariance is of great help to show us that, in fact, no problem arises because physical states are unitary by themselves, namely the contribution of unphysical states in loop
diagrams always cancels. Ghosts are crucial for this cancellation to occur. This is best seen by considering

$$
\begin{equation*}
\langle\alpha| S^{\dagger} S|\beta\rangle=\sum_{i=1,2,3}\langle\alpha| S^{\dagger}\left|\phi_{i}\right\rangle\left\langle\phi_{i}\right| S|\beta\rangle=\langle\alpha \mid \beta\rangle, \tag{6.3.27}
\end{equation*}
$$

where $|\alpha\rangle$ and $|\beta\rangle$ are two arbitrary physical states. In principle all states in the Hilbert space contribute in the completeness relation, but since the BRST operator $Q$ commutes with the $S$ matrix and $Q$ annihilates both $|\alpha\rangle$ and $|\beta\rangle$, the states $\phi_{i}$ should also be annihilated by $Q$. Hence the states $\left|\phi_{1}\right\rangle$ cannot enter eq. (6.3.27). The states $\left|\phi_{2}\right\rangle$ enter, but have vanishing inner product with states in both classes $\left|\phi_{2}\right\rangle$ and $\left|\phi_{3}\right\rangle$. We then conclude that effectively

$$
\begin{equation*}
\langle\alpha| S^{\dagger} S|\beta\rangle=\sum_{p h y s}\langle\alpha| S^{\dagger}|p h y s\rangle\langle p h y s| S|\beta\rangle=\langle\alpha \mid \beta\rangle, \tag{6.3.28}
\end{equation*}
$$

and hence unitarity is recovered.
The non-abelian generalizations of the QED WT identities, so called Slavnov-Taylor identities, provide constraints on the form of correlation functions. As an example, let consider the general relation (4.4.16), adapted to a BRST transformation in a non-abelian gauge theory. Let us take $J_{\psi}=J_{\omega}=0$ and keep only the currents for the non-abelian gauge field and the anti-ghosts. Explicitly, eq. (4.4.16) becomes

$$
\begin{align*}
& \int d^{4} x\left(\left\langle s A_{\alpha}^{\mu}(x)\right\rangle_{J} J_{A, \mu}^{\alpha}(x)+\left\langle s \omega_{\alpha}^{\star}(x)\right\rangle_{J} J_{\omega^{\star}}^{\alpha}(x)\right)= \\
& \int d^{4} x\left(\left\langle D^{\mu} \omega_{\alpha}(x)\right\rangle_{J} J_{A, \mu}^{\alpha}(x)-\frac{1}{\xi}\left\langle\partial_{\mu} A_{\alpha}^{\mu}(x)\right\rangle_{J} J_{\omega^{\star}}^{\alpha}(x)\right)=0, \tag{6.3.29}
\end{align*}
$$

where in the last equality we have integrated out the auxiliary field $H_{\alpha}=\partial_{\mu} A_{\alpha}^{\mu} / \xi$. Let us take two functional derivatives of eq. (6.3.29) with respect to the anti-ghost and gauge currents and then set all currents to zero. We get

$$
\begin{gather*}
\left.\int d^{4} x \frac{\delta^{2}}{\delta J_{A, \alpha}^{\nu}(y) \delta J_{\omega^{\star}}^{\beta}(z)}\left(\left\langle D^{\mu} \omega_{\gamma}(x)\right\rangle_{J} J_{A, \mu}^{\gamma}(x)-\frac{1}{\xi}\left\langle\partial_{\mu} A_{\gamma}^{\mu}(x)\right\rangle_{J} J_{\omega^{\star}}^{\gamma}(x)\right)\right|_{J_{A}=J_{\omega^{\star}}=0}=  \tag{6.3.30}\\
i\left\langle\omega_{\beta}^{\star}(z) D_{\nu} \omega_{\alpha}(y)\right\rangle-\frac{i}{\xi}\left\langle A_{\nu, \alpha}(y) \partial_{\mu} A_{\beta}^{\mu}(z)\right\rangle=0
\end{gather*}
$$

Multiply now expression (6.3.30) by $\partial / \partial y_{\nu}$. In so doing the first term becomes trivial, because the operator $\partial^{\nu} D_{\nu}$ is, by definition, the inverse of the exact ghost propagator:

$$
\begin{equation*}
\left\langle\omega_{\beta}^{\star}(z) \partial^{\nu} D_{\nu} \omega_{\alpha}(y)\right\rangle=-i \delta_{\alpha \beta} \delta^{(4)}(z-y) \tag{6.3.31}
\end{equation*}
$$

Plugging in eq. (6.3.30) gives

$$
\begin{equation*}
\left\langle\partial^{\nu} A_{\nu}^{\alpha}(y) \partial_{\mu} A_{\beta}^{\mu}(z)\right\rangle=-i \xi \delta_{\beta}^{\alpha} \delta^{(4)}(z-y) . \tag{6.3.32}
\end{equation*}
$$

The exact gluon propagator is trivial in the color indices and can be written as

$$
\begin{equation*}
G_{\mu \nu}^{\alpha \beta}(y, z) \equiv\left\langle A_{\nu}^{\alpha}(y) A_{\mu}^{\beta}(z)\right\rangle=\delta^{\alpha \beta} \int \frac{d^{4} p}{(2 \pi)^{4}} G_{\mu \nu}(p) e^{i p(y-z)}, \tag{6.3.33}
\end{equation*}
$$

with

$$
\begin{equation*}
p^{\mu} p^{\nu} G_{\mu \nu}(p)=-i \xi \tag{6.3.34}
\end{equation*}
$$

In perturbation theory, we can write $G_{\mu \nu}(p)=G_{\mu \nu}^{(0)}(p)+G_{\mu \nu}^{(q)}(p)$, where $G_{\mu \nu}^{(0)}(p)$ is the tree-level propagator we have computed in eq. (6.2.45) and $G_{\mu \nu}^{(q)}(p)$ encodes the quantum corrections. We notice that the classical propagator alone satisfies eq. (6.3.34), implying that the quantum corrections must be transverse: $p^{\mu} p^{\nu} G_{\mu \nu}^{(q)}(p)=0$. The gluon propagator to all orders in perturbation theory can be written as

$$
\begin{equation*}
G_{\mu \nu}(p)=\frac{-i}{p^{2}+i \epsilon}\left(\frac{\eta_{\mu \nu} p^{2}-p^{\mu} p^{\nu}}{p^{2}\left(1-\Pi\left(p^{2}\right)\right)}+\xi \frac{p_{\mu} p_{\nu}}{p^{2}}\right) \tag{6.3.35}
\end{equation*}
$$

where all quantum corrections are encoded in the scalar form factor $\Pi\left(p^{2}\right)$.
BRST invariance is very useful to also establish the renormalization properties of nonabelian gauge theories. Being a symmetry of the total action $S_{\text {tot }}=\int d^{4} x \mathcal{L}_{\text {tot }}$, one can show that the total quantum action $\Gamma_{t o t}$ is of the same form of $S_{t o t}$, with renormalized parameters. This derivation is discussed in detail in section 6.5.

### 6.4 The Background Field Method

The renormalization of non-abelian gauge theories can proceed in a perturbative way as it was explained in chapter 3 - and exemplified in the case of the abelian $U(1)$ gauge theory (QED), by singling out the leading divergences and the class of diagrams which are in need of renormalization. However, the corresponding analysis becomes quite challenging and demanding already for the lowest-order calculations, because of the rather large number of diagrams involved. In particular, in non-abelian gauge theories the identity (3.3.15) no longer holds. This important difference between abelian and non-abelian theories can be traced back to the different nature of the associated conserved currents in the two classes of theories. In abelian gauge theories we have a gauge-invariant divergence-less current $J^{\mu}$ and its associated well-defined conserved charge $Q$ given by the spatial integral of its time component. The notion of an absolute normalization of the charge $Q$ guarantees that $Z_{1}=Z_{2}$ for all charged particles. As we have seen below eq.(6.1.37), in non-abelian gauge theories the currents are either conserved but not covariant, or non-conserved but covariant. Hence there is no a well-defined standard gauge-invariant conserved charge $Q$ as in the QED case and correspondingly $Z_{1} \neq Z_{2}$, though the ratio $Z_{1} / Z_{2}$ is the same for
all charged fields. Contrary to QED, then, in addition to the gluon two-point function we should in general compute vertex corrections in order to derive the $\beta$-function for the gauge coupling. This can in principle be computed in different ways, using gauge, ghost or matter vertices. Gauge invariance guarantees that the same result is obtained independently of this choice.

As a matter of fact, also in non-abelian gauge theories the $\beta$-function can be deduced from the gluon two-point function using an alternative approach, denoted background field method, that will be discussed in this section. We present the principle of the background field method primarily as a tool to compute the renormalization constant $Z_{A}$ of the gauge fields and the associated $\beta$-function. It should be clear, however, that this method has a wider variety of applications.

### 6.4.1 The method

The basic idea behind the background field method consists in calculating the effective action in the presence of a classical background field gauge $A_{\mu}$, in total analogy to the scalar analogue (4.1.12):

$$
\begin{equation*}
e^{i \Gamma\left(A_{\mu}\right)}=\int_{1 P I} \mathcal{D} Q_{\mu} e^{i S\left(A_{\mu}+Q_{\mu}\right)} \delta(G(A, Q)) \Delta_{G}(A, Q) \tag{6.4.1}
\end{equation*}
$$

where we have denoted by $A_{\mu}$ and $Q_{\mu}$ the background and fluctuating fields, respectively. The key point is to choose a gauge fixing in the path integral such that $\Gamma(A)$ is gauge invariant under transformations of the background:

$$
\begin{align*}
\delta_{B} A_{\mu}^{\alpha} & =\partial_{\mu} \epsilon_{\alpha}+g C_{\alpha \beta \gamma} A_{\mu}^{\beta} \epsilon_{\gamma},  \tag{6.4.2}\\
\delta_{B} Q_{\mu}^{\alpha} & =g C_{\alpha \beta \gamma} Q_{\mu}^{\beta} \epsilon_{\gamma} .
\end{align*}
$$

It is important to distinguish the gauge transformations of the background (6.4.2) from the quantum gauge transformations under which ${ }^{6}$

$$
\begin{align*}
& \delta A_{\mu}^{\alpha}=0  \tag{6.4.3}\\
& \delta Q_{\mu}^{\alpha}=\partial_{\mu} \epsilon_{\alpha}+g C_{\alpha \beta \gamma}\left(A_{\mu}^{\beta}+Q_{\mu}^{\beta}\right) \epsilon_{\gamma}
\end{align*}
$$

The gauge fixing where $\Gamma(A)$ is gauge invariant is called background field gauge and is obtained by choosing the functional (6.2.30) as

$$
\begin{equation*}
G_{f}(A, Q)=D_{\mu}(A) Q_{\alpha}^{\mu}-f_{\alpha}=\partial_{\mu} Q_{\alpha}^{\mu}+g C_{\alpha \beta \gamma} A_{\mu}^{\beta} Q_{\gamma}^{\mu}-f_{\alpha} \tag{6.4.4}
\end{equation*}
$$

[^40]The ghost action associated to this gauge fixing is obtained, as usual, by taking a quantum gauge transformation of the gauge fixing functional:

$$
\begin{equation*}
\delta G(A, Q)=D_{\mu}(A) D^{\mu}(A+Q) \epsilon_{\alpha} \tag{6.4.5}
\end{equation*}
$$

and hence

$$
\begin{equation*}
\mathcal{L}_{\text {ghost }}=D_{\mu}(A) \omega_{\alpha}^{*}\left(D^{\mu}(A) \omega_{\alpha}+g C_{\alpha \beta \gamma} Q_{\beta}^{\mu} \omega_{\gamma}\right) . \tag{6.4.6}
\end{equation*}
$$

Weighting the path integral in eq. (6.4.1) with the phase factor (6.2.31) and integrating over the delta functional gives

$$
\begin{equation*}
e^{i \Gamma\left(A_{\mu}\right)}=\int_{1 P I} \mathcal{D} Q_{\mu} \mathcal{D} \omega \mathcal{D} \omega^{*} e^{i S\left(A_{\mu}+Q_{\mu}\right)+i S_{\text {ghost }}-\frac{i}{2 \xi} \int d^{4} x\left(D_{\mu}(A) Q_{\alpha}^{\mu}\right)^{2}} \tag{6.4.7}
\end{equation*}
$$

Under the background gauge transformations (6.4.2) the ghosts and anti-ghosts transform as fields in the adjoint representations:

$$
\begin{align*}
\delta_{B} \omega_{\alpha} & =g C_{\alpha \beta \gamma} \epsilon_{\beta} \omega_{\gamma},  \tag{6.4.8}\\
\delta_{B} \omega_{\alpha}^{*} & =g C_{\alpha \beta \gamma} \epsilon_{\beta} \omega_{\gamma}^{*} .
\end{align*}
$$

Notice that while the ghost and gauge fixing terms are clearly not gauge invariant under the quantum gauge transformations (6.4.3), they are invariant under the background gauge transformations (6.4.2) and (6.4.8). The matter action in $S(A+Q)$, in fact, is manifestly invariant. The gauge kinetic term is

$$
\begin{align*}
& -\frac{1}{4}\left(\partial_{\mu}\left(A_{\nu}^{\alpha}+Q_{\nu}^{\alpha}\right)-\partial_{\nu}\left(A_{\mu}^{\alpha}+Q_{\mu}^{\alpha}\right)+g C_{\alpha \beta \gamma}\left(A_{\mu}^{\beta}+Q_{\mu}^{\beta}\right)\left(A_{\nu}^{\gamma}+Q_{\nu}^{\gamma}\right)\right)^{2}=  \tag{6.4.9}\\
& -\frac{1}{4}\left(F_{\mu \nu}^{\alpha}(A)+D_{\mu}(A) Q_{\nu}^{\alpha}-D_{\nu}(A) Q_{\mu}^{\alpha}+g C_{\alpha \beta \gamma} Q_{\mu}^{\beta} Q_{\nu}^{\gamma}\right)^{2}
\end{align*}
$$

which is also invariant. We have then proved that $\delta_{B} \Gamma(A)=0$. The divergent part of $\Gamma(A)$ has to be a gauge-invariant polynomial of dimension 4 and therefore it is necessarily of the form

$$
\begin{equation*}
-\frac{Z_{A}}{4} \int d^{4} x F_{\mu \nu, \alpha}(A) F^{\mu \nu, \alpha}(A)+\text { finite parts } \tag{6.4.10}
\end{equation*}
$$

where $Z_{A}$ is a possibly divergent constant. Note that, as usual, $Z_{A}=1$ for $g \rightarrow 0$, and therefore $Z_{A}=1+\mathcal{O}\left(g^{2}\right)$. By comparing the renormalized action above to the bare one with coupling constant $g_{B}$ and bare fields $A_{B}$, it is possible to conclude that $A_{B}=Z_{A}^{1 / 2} A$ and that $g=g_{B} \mu^{-\epsilon / 2} Z_{A}^{1 / 2}$. As a consequence of the gauge invariance (which constraints the form of $\Gamma$ ) the renormalization of the coupling constant $g$ is fixed by the renormalization constant of the field $A$, as it was the case for QED. The $\beta$-function can then be calculated as in subsection 5.11.

(a)

(b)

(c)

(d)

Figure 6.1: Trilinear vertices involving one background field $A$ within the background field method. Wiggly lines indicate the background field $A$, curly lines the quantum fluctuations $Q$, dashed directed lines the ghosts, while dotted directed lines the scalars. We list here only the trilinear vertices which, at the lowest order in perturbation theory, give a non-vanishing contribution to the two-point function of the background fields: (a) $A-\omega^{*}-\omega$ vertex $V^{A \omega^{*} \omega}$ in eq. (6.4.11), (b) $A-Q-Q$ vertex $V^{A Q Q}$ in eq. (6.4.17), (c) $A-\bar{\psi}-\psi$ vertex $V^{A \bar{\psi} \psi}$ in eq. (6.4.20), and (d) $A-\phi^{\dagger}-\phi$ vertex $V^{A \phi^{\dagger} \phi}$ in eq. (6.4.21).

### 6.4.2 Two-Point Function of the Background Field: Feynman Rules

The calculation of $Z_{A}$ requires to take into account for the divergences which arise when the integration over $Q, \omega$, and $\omega^{*}$ is performed in eq. (6.4.7). Given eq.(6.4.10), it is enough to consider the two-point function of the background field $A$. This is analogous to a standard propagator if were not for the fact that $A$ is considered as an external field. Let us report below the Feynman rules that are necessary to the computation of $Z_{A}$ at one-loop level.

From $\mathcal{L}_{\text {ghost }}$ in eq. (6.4.6) one gets a $A-\omega^{*}-\omega$-vertex (represented in fig. 6.1(a))

$$
\begin{equation*}
V_{\mu \alpha ; \beta ; \gamma}^{A \omega^{*} \omega} \equiv \frac{\delta i \mathcal{L}_{g h o s t}}{\delta A_{\alpha}^{\mu} \delta \omega_{\beta}(q) \delta \omega_{\gamma}^{*}(p)}=-g C_{\alpha \beta \gamma}(p+q)_{\mu} \tag{6.4.11}
\end{equation*}
$$

where $p_{\mu}$ and $q_{\mu}$ are the momenta of the ghosts. In eq.(6.4.11) we have omitted to write the dependence of the gauge field $A_{\mu}$ on its momentum, since the vertex does not depend on it. Analogue simplifications will be done below. The additional vertices involving $A$ which are generated by this term are of the form $A-A-\omega^{*}-\omega$ and $A-Q-\omega^{*}-\omega$. While the latter does not contribute at one-loop to the two-point correlation function of the background field, the former can. We have (see fig. 6.2(a))

$$
\begin{equation*}
V_{\mu \alpha ; \beta ; \gamma ; \delta}^{A A \omega^{*} \omega} \equiv \frac{\delta i \mathcal{L}_{\text {ghost }}}{\delta A_{\alpha}^{\mu} \delta A_{\beta}^{\nu} \delta \omega_{\gamma} \delta \omega_{\delta}^{*}}=i g^{2} \eta_{\mu \nu}\left(C_{\alpha \gamma \omega} C_{\beta \delta \omega}+C_{\alpha \delta \omega} C_{\beta \gamma \omega}\right) . \tag{6.4.12}
\end{equation*}
$$

The gauge-fixing term $\mathcal{L}_{\text {g.f. }}$. generates a $A-Q$ - $Q$-vertex

$$
\begin{equation*}
\widetilde{V}_{\mu \alpha ; \nu \beta ; \rho \gamma}^{A Q Q} \equiv \frac{\delta i \mathcal{L}_{g . f .}}{\delta A_{\alpha}^{\mu} \delta Q_{\beta}^{\nu}(q) \delta Q_{\gamma}^{\rho}(p)}=-\frac{g}{\xi}\left(\eta_{\mu \nu} C_{\alpha \beta \gamma} p_{\rho}+\eta_{\mu \rho} C_{\alpha \gamma \beta} q_{\nu}\right) \tag{6.4.13}
\end{equation*}
$$



Figure 6.2: Quartic vertices involving two background fields $A$ within the background field method. Wiggly lines indicate the background field $A$, curly lines the quantum fluctuations $Q$, dashed directed lines the ghosts, while dotted directed lines the scalars. We list here only the quartic vertices which, at the lowest order in perturbation theory, might give a non-vanishing contribution to the two-point function of the background fields: (a) $A$ -$A-\omega^{*}-\omega$ vertex $V^{A A \omega \omega^{*}}$ in eq. (6.4.12), (b) $A-A-Q-Q$ vertex $V^{A A Q Q}$ in eq. (6.4.18), (b) $A-A-\phi^{\dagger}-\phi$ vertex $V^{A A \phi^{\dagger} \phi}$ in eq. (6.4.22).
with the momenta of the fields oriented as in fig. 6.1(b). In addition to this interaction, $\mathcal{L}_{\text {g.f. }}$ generates a vertex $A-A-Q-Q$,

$$
\begin{equation*}
\widetilde{V}_{\mu \alpha ; \nu \beta ; \rho \gamma ; \sigma \delta}^{A A Q Q} \equiv \frac{\delta i \mathcal{L}_{g . f .}}{\delta A_{\alpha}^{\mu} \delta A_{\beta}^{\nu} \delta Q_{\gamma}^{\rho} \delta Q_{\delta}^{\sigma}}=-\frac{i g^{2}}{\xi}\left(\eta_{\mu \rho} \eta_{\nu \sigma} C_{\alpha \gamma \omega} C_{\beta \delta \omega}+\eta_{\mu \sigma} \eta_{\nu \rho} C_{\beta \gamma \omega} C_{\alpha \delta \omega}\right) . \tag{6.4.14}
\end{equation*}
$$

The vertex involving three gauge fields $A$ in the standard YM theory generates several vertices: $A-A-Q, A-Q-Q$, and $Q-Q-Q$ but only the second one is relevant for determining $Z_{A}$ at the lowest order. We have

$$
\begin{equation*}
\widehat{V}_{\mu \alpha ; \nu \beta ; \rho \gamma}^{A Q Q} \equiv \frac{\delta i \mathcal{L}}{\delta A_{\alpha}^{\mu}(k) \delta Q_{\beta}^{\nu}(q) \delta Q_{\gamma}^{\rho}(p)}=g C_{\alpha \beta \gamma}\left((k-q)_{\rho} \eta_{\mu \nu}+(q-p)_{\mu} \eta_{\nu \rho}+(p-k)_{\nu} \eta_{\rho \mu}\right) \tag{6.4.15}
\end{equation*}
$$

where all momenta are assumed to be outgoing from the vertex. The expression (6.4.15) coincides with the trilinear vertex computed in YM theories in standard gauges. The vertex involving four gauge fields $A$ in the standard YM theory generates several vertices: $A-A-A-A, A-A-A-Q, A-A-Q-Q, A-Q-Q-Q$, and $Q-Q-Q-Q$ but only the third one is relevant for determining $Z_{A}$ at the lowest order. The corresponding expression is again the standard one:

$$
\begin{align*}
& \widehat{V}_{\mu \alpha ; \nu \beta ; \rho \gamma ; \sigma \delta}^{A A Q Q} \equiv \frac{\delta i \mathcal{L}}{\delta A_{\alpha}^{\mu} \delta A_{\beta}^{\nu} \delta Q_{\gamma}^{\rho} \delta Q_{\delta}^{\sigma}}=-i g^{2}\left(C_{\alpha \beta \omega} C_{\gamma \delta \omega}\left(\eta^{\mu \rho} \eta^{\nu \sigma}-\eta_{\mu \sigma} \eta_{\nu \rho}\right)+\right.  \tag{6.4.16}\\
&\left.C_{\alpha \gamma \omega} C_{\beta \delta \omega}\left(\eta_{\mu \nu} \eta_{\rho \sigma}-\eta_{\mu \sigma} \eta_{\nu \rho}\right)+C_{\alpha \delta \omega} C_{\beta \gamma \omega}\left(\eta_{\mu \nu} \eta_{\rho \sigma}-\eta_{\mu \rho} \eta_{\nu \sigma}\right)\right) .
\end{align*}
$$

The total vertices $A-Q-Q$ and $A-A-Q-Q$ of this theory are given by the sum of the one stemming from the gauge-fixing term in eqs. (6.4.13) and (6.4.14) and the ones from $\mathcal{L}$ in
eqs. (6.4.15) and (6.4.16), see figs.6.1(b) and 6.2(b),

$$
\begin{align*}
& V_{\mu \alpha ; \nu \beta ; \rho \gamma}^{A Q Q}=g C_{\alpha \beta \gamma}\left(\left(k-q-\frac{1}{\xi} p\right)_{\rho} \eta_{\mu \nu}+(q-p)_{\mu} \eta_{\rho \nu}+\left(p-k+\frac{1}{\xi} q\right)_{\nu} \eta_{\rho \mu}\right),(6.4 .17  \tag{6.4.17}\\
& V_{\mu \alpha ; \nu \beta ; \rho \gamma ; \sigma \delta}^{A A Q Q}=-i g^{2}\left(C_{\alpha \beta \omega} C_{\gamma \delta \omega}\left(\eta_{\mu \rho} \eta_{\nu \sigma}-\eta_{\mu \sigma} \eta_{\nu \rho}\right)+C_{\alpha \gamma \omega} C_{\beta \delta \omega}\right.  \tag{6.4.18}\\
&\left.\left(\eta_{\mu \nu} \eta_{\rho \sigma}-\eta_{\mu \sigma} \eta_{\nu \rho}+\frac{1}{\xi} \eta_{\mu \rho} \eta_{\nu \sigma}\right)+C_{\alpha \delta \omega} C_{\beta \gamma \omega}\left(\eta_{\mu \nu} \eta_{\rho \sigma}-\eta_{\mu \rho} \eta_{\nu \sigma}+\frac{1}{\xi} \eta_{\mu \sigma} \eta_{\nu \rho}\right)\right)
\end{align*}
$$

where

$$
\begin{align*}
V_{\mu \alpha ; \nu \beta ; \rho \gamma}^{A Q Q} & \equiv \widetilde{V}_{\mu \alpha ; \nu \beta ; \rho \gamma}^{A Q Q}+\widehat{V}_{\mu \alpha ; \nu \beta ; \rho \gamma}^{A Q Q}, \\
V_{\mu \alpha ; \nu \beta ; \rho \gamma ; \sigma \delta}^{A A Q Q} & \equiv \widetilde{V}_{\mu \alpha ; \nu \beta ; \rho \gamma ; \sigma \delta}^{A A Q Q}+\widehat{V}_{\mu \alpha ; \nu \beta ; \rho \gamma ; \sigma \delta}^{A A Q Q} . \tag{6.4.19}
\end{align*}
$$

The interaction with the matter is ruled by the standard vertices of the YM theory. For fermions we have, see fig. 6.1(c),

$$
\begin{equation*}
V_{\mu \alpha ; i j}^{A \bar{\psi} \psi} \equiv \frac{\delta i \mathcal{L}}{\delta A_{\alpha}^{\mu} \delta \psi_{j} \delta \bar{\psi}_{i}}=i g \gamma_{\mu} t_{i j}^{\alpha} \tag{6.4.20}
\end{equation*}
$$

where $t^{\alpha}$ are the gauge generators in the appropriate representation. For scalars, we can have both the trilinear and quartic vertices, see figs. 6.1(d) and 6.2(c):

$$
\begin{align*}
V_{\mu \alpha ; i j}^{A \phi^{\dagger} \phi} & \equiv \frac{\delta i \mathcal{L}}{\delta A_{\alpha}^{\mu} \delta \phi_{j}(q) \delta \phi_{i}^{\dagger}(p)}=-i g t_{i j}^{\alpha}(p+q)_{\mu}  \tag{6.4.21}\\
V_{\mu \alpha ; \nu ; i j}^{A A \phi^{\dagger} \phi} & \equiv \frac{\delta i \mathcal{L}}{\delta A_{\alpha}^{\mu} \delta A_{\beta}^{\nu} \delta \phi_{j} \delta \phi_{i}^{\dagger}}=i g^{2} \eta_{\mu \nu}\left\{t^{\alpha}, t^{\beta}\right\}_{i j} \tag{6.4.22}
\end{align*}
$$

where again $t^{\alpha}$ are the gauge generators in the appropriate representation.
With all these vertices at hand, the radiative corrections to the two-point function of the background field can be computed. We leave to the reader to perform the computation of $Z_{A}$ at one-loop level and report here only the final result for the $\beta$-function.

In a $S U(N)$ gauge theory with $n_{f}$ Dirac fermions in representations $r_{i}^{(F)}$ and $n_{s}$ scalars in representations $r_{i}^{(S)}$ of $S U(N)$, one has at one-loop level

$$
\begin{equation*}
\beta(g)=-\frac{g^{3}}{(4 \pi)^{2}}\left(\frac{11 N}{3}-\frac{4}{3} \sum_{i=1}^{n_{f}} T_{2}\left(r_{i}^{(F)}\right)-\frac{1}{3} \sum_{i=1}^{n_{s}} T_{2}\left(r_{i}^{(S)}\right)\right) \tag{6.4.23}
\end{equation*}
$$

where

$$
\begin{equation*}
\operatorname{tr}_{r} t^{\alpha} t^{\beta} \equiv \delta^{\alpha \beta} T_{2}(r), \quad T_{2}(a d j .)=N \tag{6.4.24}
\end{equation*}
$$

In particular, for $n_{s}=0$ and $r_{i}^{(F)}$ equal to the fundamental representation for any fermion, we have:

$$
\begin{equation*}
\beta(g)=-\frac{g^{3}}{(4 \pi)^{2}}\left(\frac{11 N}{3}-\frac{2}{3} n_{f}\right) \tag{6.4.25}
\end{equation*}
$$

QCD corresponds to the choice $N=3$ and $n_{f}=6$ and is then an asymptotically free theory with $\beta(g)<0$, as we discuss in general in sec. 5.3.

### 6.5 Proof of the Renormalizability of Non-Abelian Gauge Theories**

The BRST symmetry provides a useful set-up to prove the renormalizability of non-abelian gauge theories. Before attaching the problem, we need however to develop a bit on the formalism and derive what is called the master equation. ${ }^{7}$

### 6.5.1 The Master Equation**

The relation (4.4.17), applied to the BRST transformations (6.3.3), give

$$
\begin{equation*}
\int d^{4} x\left\langle\Delta^{n}[\chi(x)]\right\rangle_{J(\chi)} \frac{\delta_{L} \Gamma}{\delta \chi^{n}(x)}=0 \tag{6.5.1}
\end{equation*}
$$

where $\chi^{n}$ is a compact notation to denote all the fields: $\chi^{n}=\left(A_{\mu}^{\alpha}, \omega_{\alpha}, \omega_{\alpha}^{*}, \psi, \phi, H_{\alpha}\right)$, and we have defined $\delta \chi^{n}=\theta s \chi^{n} \equiv \theta \Delta^{n}(\chi)$, Care has to be taken in taking functional derivatives in presence of Grassmann fields, so we distinguish with a subscript $L$ and $R$ derivatives defined as follows:

$$
\begin{equation*}
\frac{\delta_{L} F}{\delta \phi(x)}=\frac{\overrightarrow{\delta_{L}}}{\delta \phi(x)} F, \quad \frac{\delta_{R} F}{\delta \phi(x)}=F \frac{\overleftarrow{\delta_{R}}}{\delta \phi(x)} \tag{6.5.2}
\end{equation*}
$$

The BRST transformations are quadratic in the fields and hence $\left\langle\Delta^{n}[\chi(x)]\right\rangle_{J(\chi)} \neq \Delta^{n}[\chi(x)]$. In order to overcome this problem, we introduce, in addition to the source $J_{n}$ that couple to $\chi^{n}$, extra sources $K_{n}(x)$ that couple to $\Delta^{n}(\chi)$. Correspondingly we define

$$
\begin{equation*}
e^{i W[J, K]}=\int \mathcal{D} \chi^{n} e^{i S_{t o t}+i \int \Delta^{n} K_{n}+i \int \chi^{n} J_{n}} \tag{6.5.3}
\end{equation*}
$$

and the 1PI generator

$$
\begin{equation*}
\Gamma[\chi, K]=W[J, K]-\int d^{4} x \chi^{n}(x) J_{n}(x) \tag{6.5.4}
\end{equation*}
$$

where $J_{n}=J_{n}(\chi, K)$, obtained by inverting the relation

$$
\begin{equation*}
\frac{\delta_{R} W[J, K]}{\delta J_{n}(x)}=\chi^{n}(x) \tag{6.5.5}
\end{equation*}
$$

Under a BRST transformation $\delta \Delta^{n}=\theta s^{2} \chi^{n}=0$ and hence the 1 PI action $\Gamma[\chi, K]$ satisfies the analogue of eq.(6.5.1):

$$
\begin{equation*}
\int d^{4} x\left\langle\Delta^{n}[\chi(x)]\right\rangle_{J(\chi, K)} \frac{\delta_{L} \Gamma[\chi, K]}{\delta \chi^{n}(x)}=0 \tag{6.5.6}
\end{equation*}
$$

Taking a functional R-derivation with respect to $K_{n}(x)$ of eq.(6.5.4) gives

$$
\begin{align*}
& \frac{\delta_{R} \Gamma[\chi, K]}{\delta K_{n}(x)}=\frac{\delta_{R} W[J, K]}{\delta K_{n}(x)}+\int d^{4} y \frac{\delta_{R} W[J, K]}{\delta J_{m}(y)} \frac{\delta_{R} J_{m}(y)}{\delta K_{n}(x)}-\int d^{4} y \chi^{m}(y) \frac{\delta_{R} J_{m}(y)}{\delta K_{n}(x)}  \tag{6.5.7}\\
= & \frac{\delta_{R} W[J, K]}{\delta K_{n}(x)}=i\left\langle\Delta^{n}[\chi(x)]\right\rangle_{J(\chi, K)} .
\end{align*}
$$

[^41]We can then rewrite eq.(6.5.6) as

$$
\begin{equation*}
\int d^{4} x \frac{\delta_{R} \Gamma[\chi, K]}{\delta K_{n}(x)} \frac{\delta_{L} \Gamma[\chi, K]}{\delta \chi^{n}(x)}=0 . \tag{6.5.8}
\end{equation*}
$$

Since the sources $K_{n}$ have the same commuting properties as $\Delta^{n}$, which are opposite to those of $\chi^{n}, \delta_{R} \Gamma / \delta K_{n}$ and $\delta_{L} \Gamma / \delta J_{n}$ anti-commute with each other. We can then rewrite more compactly eq.(6.5.8) as

$$
\begin{equation*}
(\Gamma, \Gamma)=0, \tag{6.5.9}
\end{equation*}
$$

where we have defined, for any two functionals $F$ and $G$, the operator

$$
\begin{equation*}
(F, G) \equiv \int d^{4} x\left(\frac{\delta_{R} F}{\delta \chi^{n}(x)} \frac{\delta_{L} G}{\delta K_{n}(x)}-\frac{\delta_{R} F}{\delta K_{n}(x)} \frac{\delta_{L} G}{\delta \chi^{n}(x)}\right) \tag{6.5.10}
\end{equation*}
$$

Equation (6.5.9) is denoted the master equation and is the key relation that we will use in the next subsection to prove the renormalizability of non-abelian gauge theories.

### 6.5.2 Structure of divergences ${ }^{\star \star}$

In this subsection we will prove that all the divergences occurring at any loop in perturbation theory can be reabsorbed by a renormalization of the couplings appearing in the classical action, namely by a proper choice of counter-terms. We define the classical bare action including the extra $K$-source term:

$$
\begin{equation*}
S_{B}=S_{t o t}(\chi)+\int \Delta^{n} K_{n} \tag{6.5.11}
\end{equation*}
$$

As usual, we rewrite the action $S_{B}=S+S_{\infty}$, where $S$ is the action written in terms of physical fields and couplings and $S_{\infty}$ includes the counter-terms. The 1PI action $\Gamma[\chi, K]$ admits a loop expansion of the form

$$
\begin{equation*}
\Gamma[\chi, K]=\sum_{N=0}^{\infty} \Gamma_{N}[\chi, K] \tag{6.5.12}
\end{equation*}
$$

where $\Gamma_{N}[\chi, K]$ is the $N$-loop effective action. Clearly, we have $\Gamma_{0}=S$. Plugging eq.(6.5.12) in the master equation (6.5.9) gives an infinite set of equations, valid for any $N$ :

$$
\begin{equation*}
\sum_{N^{\prime}=0}^{N}\left(\Gamma_{N^{\prime}}, \Gamma_{N-N^{\prime}}\right)=0 \tag{6.5.13}
\end{equation*}
$$

We denote by $\Gamma_{N, \infty}$ the divergent part of the $N$-loop effective action. If the theory is renormalizable, by a proper choice of $S_{\infty}, \Gamma_{N}$ should be finite for each value of $N$, i.e. $\Gamma_{N, \infty}$ should all vanish. We will prove this statement by induction, assuming that this is
true for $M \leq N-1$, namely that all the divergences in $\Gamma_{M}$ have been reabsorbed in $S_{\infty}$. If this is true, divergences in eq.(6.5.13) can only arise from the terms with $N^{\prime}=0, N$, that give rise to the equation

$$
\begin{equation*}
\left(S, \Gamma_{N, \infty}\right)=0 \tag{6.5.14}
\end{equation*}
$$

Let us see the possible form of $\Gamma_{N, \infty}$. It should be a local functional, invariant under all the linearly realized symmetries of the classical action. By power counting renormalization, all operators should have dimension less or equal to 4 . Notice that

$$
\begin{align*}
& {\left[\Delta_{n}\right]=\left[\chi_{n}\right]+1, \quad\left[K_{n}\right]=3-\left[\chi_{n}\right],}  \tag{6.5.15}\\
& \operatorname{gh} .\left(\chi_{n}\right) \equiv \gamma_{n}, \quad \operatorname{gh.}\left(\Delta_{n}\right)=\gamma_{n}+1, \quad \operatorname{gh} .\left(K_{n}\right)=-\gamma_{n}-1,
\end{align*}
$$

where we have denoted by $\gamma_{n}$ the ghost number of the field $\chi^{n}$. Since $\Delta^{H}$ trivially vanishes, $K_{H}=0$. We see from eq.(6.5.15) that all $K$-currents have negative ghost charge, with the exception of $K_{\omega^{*}}$ that has ghost charge zero. Thus the only allowed operator quadratic in the $K$-currents is $K_{\omega^{*}}^{\alpha} K_{\omega^{*}}^{\alpha}$. This is already of dimension four, so no extra fields can be added to it. From eq.(6.5.7) we have

$$
\begin{equation*}
\frac{\delta_{R} \Gamma_{N, \infty}}{\delta K_{\omega^{*}}^{\alpha}(x)}=i\left\langle\Delta^{\omega^{*}}[\chi(x)]\right\rangle_{J(\chi, K)}=-i H_{\alpha}(x) \tag{6.5.16}
\end{equation*}
$$

where the last identity applies because the BRST transformation of $\omega_{\alpha}^{*}$ is linear. The right-hand side of eq.(6.5.16) is fully saturated by the classical action, and we conclude that $\Gamma_{N, \infty}$ does not depend on $K_{\omega^{*}}$ for $N>0$ and is at most linear in the $K$ 's for any $N$. We can write

$$
\begin{equation*}
\Gamma_{N, \infty}[\chi, K]=\Gamma_{N, \infty}[\chi, 0]+\int d^{4} x \mathcal{D}_{N}^{n}[\chi] K_{n}(x) \tag{6.5.17}
\end{equation*}
$$

Recall that

$$
\begin{equation*}
S[\chi, K]=S[\chi]+\int d^{4} x \Delta_{n}[\chi] K_{n}(x) \tag{6.5.18}
\end{equation*}
$$

Let us plug eqs.(6.5.17) and (6.5.18) in eq.(6.5.14). The terms with no $K$ 's give

$$
\begin{equation*}
\int d^{4} x\left(-\frac{\delta_{L} S[\chi]}{\delta \chi^{n}(x)} \mathcal{D}_{N}^{n}[\chi(x)]-\Delta_{n}[\chi(x)] \frac{\delta_{L} \Gamma_{N, \infty}}{\delta \chi^{n}(x)}\right)=0 \tag{6.5.19}
\end{equation*}
$$

In deriving eq.(6.5.19) we used the fact that $K_{n}$ and $\chi_{n}$ are bosonic (fermionic) and fermionic (bosonic) for any $n$, hence for any two arbitrary bosonic functionals $A$ and $B$ one has

$$
\begin{equation*}
\frac{\delta_{R} A}{\delta \chi^{n}} \frac{\delta_{L} B}{\delta K_{n}}=-\frac{\delta_{L} A}{\delta \chi^{n}} \frac{\delta_{R} B}{\delta K_{n}}=-\frac{\delta_{L} B}{\delta K_{n}} \frac{\delta_{L} A}{\delta \chi^{n}} . \tag{6.5.20}
\end{equation*}
$$

The terms linear in $K$ 's give

$$
\begin{equation*}
\int d^{4} x\left(\mathcal{D}_{N}^{n}[\chi(x)] \frac{\delta_{L} \Delta_{m}[\chi]}{\delta \chi^{n}(x)}+\Delta_{n}[\chi(x)] \frac{\delta_{L} \mathcal{D}_{N}^{m}[\chi]}{\delta \chi^{n}(x)}\right)=0 . \tag{6.5.21}
\end{equation*}
$$

Equation (6.5.19) can be rewritten as

$$
\begin{equation*}
\int d^{4} x \Delta_{N}^{(\epsilon) n} \frac{\delta_{L} \Gamma_{N}^{(\epsilon)}}{\delta \chi^{n}}=0 \tag{6.5.22}
\end{equation*}
$$

where we have defined

$$
\begin{align*}
\Gamma_{N}^{(\epsilon)}[\chi] & =S[\chi]+\epsilon \Gamma_{N, \infty}[\chi, 0] \\
\Delta_{N}^{(\epsilon) n} & =\Delta^{n}+\epsilon \mathcal{D}_{N}^{n} \tag{6.5.23}
\end{align*}
$$

with $\epsilon$ an infinitesimal parameter. The $\mathcal{O}\left(\epsilon^{0}\right)$ term vanishes because $S$ is BRST-invariant, while the $\mathcal{O}(\epsilon)$ terms reproduce eq.(6.5.19). Similarly, eq.(6.5.19) can be rewritten as the $\mathcal{O}(\epsilon)$ terms of the following relation:

$$
\begin{equation*}
\Delta_{N}^{(\epsilon) n} \frac{\delta_{L}}{\delta \chi^{n}}\left(\Delta_{N}^{(\epsilon) n} \frac{\delta_{L}}{\delta \chi^{n}}\right)=0 \tag{6.5.24}
\end{equation*}
$$

We conclude from eqs.(6.5.23) and (6.5.24) that the action $\Gamma_{N}^{(\epsilon)}$ is invariant under the transformations generated by $\Delta_{N}^{(\epsilon) n}$, which are nilpotent, like the original BRST transformations $\Delta^{n}$.

Comparing eqs.(6.5.17) and (6.5.18) we notice that $\mathcal{D}_{N}^{n}$ and $\Delta_{n}$ have the same ghost and Lorentz quantum numbers. The most general $\Delta_{N}^{(\epsilon) n}$ transformations is then parametrized by a simple generalization of the original BRST transformations (6.3.3):

$$
\begin{align*}
\delta_{\theta}^{(\epsilon) N} \Psi & =i \theta Q_{\alpha}^{N} \omega_{\alpha} \Psi \\
\delta_{\theta}^{(\epsilon) N} A_{\mu \alpha} & =\theta\left(B_{\alpha \beta}^{N} \partial_{\mu} \omega_{\beta}+D_{\alpha \beta \gamma}^{N} A_{\mu \beta} \omega_{\gamma}\right),  \tag{6.5.25}\\
\delta_{\theta}^{(\epsilon) N} \omega_{\alpha} & =-\frac{1}{2} \theta E_{\alpha \beta \gamma}^{N} \omega_{\beta} \omega_{\gamma} .
\end{align*}
$$

The transformations of $\omega_{\alpha}^{\star}$ and $H_{\alpha}$ are like the original BRST transformations (6.3.3) since $\mathcal{D}_{N}^{\omega^{\star}}=K_{H}=0$. For simplicity, from now on we omit the superscript $N$ on $\delta_{\theta}^{(\epsilon) N}$ and the coefficients $Q_{\alpha}^{N}, B_{\alpha \beta}^{N}, D_{\alpha \beta \gamma}^{N}$ and $E_{\alpha \beta \gamma}^{N}$. Nilpotency of the ghost transformation in eq.(6.5.25) requires that $E_{\alpha \beta \gamma}$ satisfies a Jacobi identity (compare with eq.(6.3.8)). Hence $E_{\alpha \beta \gamma}$ should be the structure constant of a Lie algebra. Since for $\epsilon \rightarrow 0$ we should recover the original BRST transformations, we conclude that $E_{\alpha \beta \gamma}$ must be the structure constant of the original Lie algebra, up to a constant:

$$
\begin{equation*}
E_{\alpha \beta \gamma}=Z C_{\alpha \beta \gamma} \tag{6.5.26}
\end{equation*}
$$

Consider now the gauge fields. We have $s^{(\epsilon)} A_{\mu \alpha}=B_{\alpha \beta} \partial_{\mu} \omega_{\beta}+D_{\alpha \beta \gamma} A_{\mu \beta} \omega_{\gamma}$. Then

$$
\begin{align*}
\delta_{\theta}^{(\epsilon)}\left(s^{(\epsilon)} A_{\mu}^{\alpha}\right)= & \partial_{\mu}\left(-\frac{1}{2} B_{\alpha \beta} \theta E_{\beta \gamma \delta} \omega_{\gamma} \omega_{\delta}\right)+D_{\alpha \beta \gamma} \theta\left(B_{\beta \sigma} \partial_{\mu} \omega_{\sigma}+D_{\beta \rho \sigma} A_{\mu}^{\rho} \omega_{\sigma}\right) \omega_{\gamma} \\
& +D_{\alpha \beta \gamma} A_{\mu}^{\beta}\left(-\frac{1}{2} \theta E_{\gamma \rho \sigma} \omega_{\rho} \omega_{\sigma}\right) . \tag{6.5.27}
\end{align*}
$$

Using eq.(6.5.26) and reshuffling indices, the terms proportional to $A_{\mu}$ in eq. (6.5.27) give

$$
\begin{equation*}
\left.\delta_{\theta}^{(\epsilon)}\left(s^{(\epsilon)} A_{\mu}^{\alpha}\right)\right|_{A}=-\frac{1}{2} \theta A_{\mu}^{\beta} \omega_{\rho} \omega_{\sigma}\left(D_{\alpha \gamma \rho} D_{\gamma \beta \sigma}-D_{\alpha \gamma \sigma} D_{\gamma \beta \rho}+Z C_{\gamma \rho \sigma} D_{\alpha \beta \gamma}\right) . \tag{6.5.28}
\end{equation*}
$$

This term vanishes, again due to the Jacobi identity, if

$$
\begin{equation*}
D_{\alpha \beta \gamma}=Z C_{\alpha \beta \gamma} \tag{6.5.29}
\end{equation*}
$$

The terms with no gauge fields give

$$
\begin{equation*}
\left.\delta_{\theta}^{(\epsilon)}\left(s^{(\epsilon)} A_{\mu}^{\alpha}\right)\right|_{n o A}=\theta \partial_{\mu} \omega_{\sigma} \omega_{\gamma} Z\left(C_{\alpha \beta \gamma} B_{\beta \sigma}-B_{\alpha \beta} C_{\beta \sigma \gamma}\right) \tag{6.5.30}
\end{equation*}
$$

and vanishes if

$$
\begin{equation*}
B_{\alpha \beta}=Z N \delta_{\alpha \beta} \tag{6.5.31}
\end{equation*}
$$

Finally, let's look at the fermion transformations.

$$
\begin{align*}
\delta_{\theta}^{(\epsilon)}\left(s^{(\epsilon)} \Psi\right) & =i Q_{\alpha}\left(\left(\delta_{\theta}^{(\epsilon)} \omega_{\alpha}\right) \Psi+\omega_{\alpha}\left(\delta_{\theta}^{(\epsilon)} \Psi\right)\right)=i Q_{\alpha}\left(-\frac{1}{2} Z \theta C_{\alpha \beta \gamma} \omega_{\beta} \omega_{\gamma} \Psi+\omega_{\alpha} i Q_{\beta} \theta \omega_{\beta} \Psi\right) \\
& =\theta \omega_{\alpha} \omega_{\beta} \Psi\left(-\frac{i}{2} Z C_{\alpha \beta \gamma} Q_{\gamma}+\frac{1}{2}\left[Q_{\alpha}, Q_{\beta}\right]\right) \tag{6.5.32}
\end{align*}
$$

This vanishes if

$$
\begin{equation*}
Q_{\alpha}=Z t_{\alpha} \tag{6.5.33}
\end{equation*}
$$

Summarizing, we find that the $\Delta_{N}^{(\epsilon) n}$ transformations (6.5.25) read

$$
\begin{align*}
\delta_{\theta}^{(\epsilon)} \Psi & =i \theta Z t_{\alpha} \omega_{\alpha} \Psi \\
\delta_{\theta}^{(\epsilon)} A_{\mu \alpha} & =\theta Z\left(Z^{\prime} \partial_{\mu} \omega_{\alpha}+C_{\alpha \beta \gamma} A_{\mu \beta} \omega_{\gamma}\right)  \tag{6.5.34}\\
\delta_{\theta}^{(\epsilon)} \omega_{\alpha} & =-\frac{1}{2} Z \theta C_{\alpha \beta \gamma} \omega_{\beta} \omega_{\gamma}
\end{align*}
$$

Modulo the two (N-dependent) constants $Z$ and $Z^{\prime}$, they are like the original BRST transformations (6.3.3). Let us now construct the possible form of the action $\Gamma_{N}^{(\epsilon)}[\chi]$. It is a local functional, has to contain up to dimension four operators, and has to be invariant under the transformations (6.5.34) and under all the linearly realized symmetries of the original total tree-level action (6.3.2). These are Lorentz invariance, ghost number conservation and the global gauge transformations

$$
\begin{align*}
\delta \Psi & =i t_{\alpha} \epsilon_{\alpha} \Psi \\
\delta A_{\mu} & =C_{\alpha \beta \gamma} A_{\mu \beta} \epsilon_{\gamma} \\
\delta \omega_{\alpha} & =C_{\alpha \beta \gamma} \omega_{\beta} \epsilon_{\gamma}  \tag{6.5.35}\\
\delta \omega_{\alpha}^{*} & =C_{\alpha \beta \gamma} \omega_{\beta}^{\star} \epsilon_{\gamma} \\
\delta H_{\alpha} & =C_{\alpha \beta \gamma} H_{\beta} \epsilon_{\gamma}
\end{align*}
$$

For the gauge fixing choice $f_{\alpha}=\partial_{\mu} A_{\alpha}$, we have an additional symmetry, invariance under anti-ghost shifts:

$$
\begin{equation*}
\delta_{T} \omega_{\alpha}^{\star}=c_{\alpha}, \tag{6.5.36}
\end{equation*}
$$

where $c_{\alpha}$ is an arbitrary Grassmann constant parameter. It is clear that the only dimension four or less terms containing ghosts and compatible with all the above symmetries are $\partial_{\mu} \omega_{\alpha}^{\star} \partial^{\mu} \omega_{\alpha}$ and $d_{\alpha \beta \gamma} \partial_{\mu} \omega_{\alpha}^{\star} A_{\gamma}^{\mu} \omega_{\beta}$. The only terms involving $H_{\alpha}$ are $H_{\alpha} H_{\alpha}, H_{\alpha} \partial_{\mu} A_{\alpha}^{\mu}$ and $e_{\alpha \beta \gamma} H_{\alpha} A_{\beta}^{\mu} A_{\mu \gamma}$. We have $\Gamma_{N}^{(\epsilon)}[\chi]=\int d^{4} x \mathcal{L}_{N}^{(\epsilon)}$, with

$$
\begin{align*}
\mathcal{L}_{N}^{(\epsilon)}= & \mathcal{L}^{\prime}(\psi, A)+\frac{1}{2} \xi^{\prime} H_{\alpha} H^{\alpha}+c H_{\alpha} \partial_{\mu} A_{\alpha}^{\mu}-e_{\alpha \beta \gamma} H_{\alpha} A_{\beta}^{\mu} A_{\mu \gamma}  \tag{6.5.37}\\
& -Z_{\omega} \partial_{\mu} \omega_{\alpha}^{\star} \partial^{\mu} \omega_{\alpha}-d_{\alpha \beta \gamma} \partial_{\mu} \omega_{\alpha}^{\star} A_{\gamma}^{\mu} \omega_{\beta} .
\end{align*}
$$

In eq.(6.5.37), $\mathcal{L}^{\prime}(\psi, A)$ is the part of the Lagrangian that depends only on the gauge and matter fields. Let us impose the modified BRST transformations (6.5.34) to $\mathcal{L}_{N}^{(\epsilon)}$. The terms linear in the $H_{\alpha}$ field give

$$
\begin{align*}
\left.\delta^{(\epsilon)} \mathcal{L}_{N}^{(\epsilon)}\right|_{H_{\alpha}}= & c H_{\alpha} \theta Z \partial^{\mu}\left(Z^{\prime} \partial_{\mu} \omega_{\alpha}+C_{\alpha \beta \gamma} A_{\mu \beta} \omega_{\gamma}\right)-2 e_{\alpha \beta \gamma} H_{\alpha} A_{\mu \gamma} \theta Z\left(Z^{\prime} \partial_{\mu} \omega_{\beta}+C_{\beta \rho \sigma} A_{\mu \rho} \omega_{\sigma}\right) \\
& +Z_{\omega} \theta \partial_{\mu} H_{\alpha} \partial^{\mu} \omega_{\alpha}+d_{\alpha \beta \gamma} \theta \partial_{\mu} H_{\alpha} A_{\gamma}^{\mu} \omega_{\beta} \\
= & \theta \partial_{\mu} H_{\alpha} \partial^{\mu} \omega_{\alpha}\left(-c Z Z^{\prime}+Z_{\omega}\right)+\theta \partial_{\mu} H_{\alpha} A_{\gamma}^{\mu} \omega_{\beta}\left(c Z C_{\alpha \beta \gamma}+d_{\alpha \beta \gamma}\right) \\
& -2 \theta H_{\alpha} A_{\mu}^{\gamma} A_{\mu \rho} \omega_{\sigma}\left(e_{\alpha \beta \gamma} Z C_{\beta \rho \sigma}\right)-2 \theta H_{\alpha} A_{\mu}^{\gamma} \partial^{\mu} \omega_{\beta}\left(e_{\alpha \beta \gamma} Z Z^{\prime}\right) . \tag{6.5.38}
\end{align*}
$$

This vanishes only if

$$
\begin{equation*}
e_{\alpha \beta \gamma}=0, \quad d_{\alpha \beta \gamma}=-c Z C_{\alpha \beta \gamma}, \quad c=\frac{Z_{\omega}}{Z Z^{\prime}} . \tag{6.5.39}
\end{equation*}
$$

The remaining terms are of the form $\partial \omega^{\star} \omega \partial \omega$ and $\partial \omega^{\star} A \omega^{2}$. They do not give additional constraints. Indeed,

$$
\begin{align*}
\delta^{(\epsilon)} \mathcal{L}_{N}^{(\epsilon)} \mid \partial \omega^{\star} \omega \partial \omega & =-Z_{\omega} \partial_{\mu} \omega_{\alpha}^{\star}\left(-\frac{1}{2} Z\right) C_{\alpha \beta \gamma} \theta \partial^{\mu}\left(\omega_{\beta} \omega_{\gamma}\right)-d_{\alpha \beta \gamma} \partial_{\mu} \omega_{\alpha}^{\star} \theta Z Z^{\prime} \partial^{\mu} \omega_{\gamma} \omega_{\beta}  \tag{6.5.40}\\
& =\theta \partial_{\mu} \omega_{\alpha}^{\star} \partial^{\mu} \omega_{\gamma} \omega_{\beta} Z\left(Z_{\omega} C_{\alpha \beta \gamma}+Z^{\prime} d_{\alpha \beta \gamma}\right)=0
\end{align*}
$$

and

$$
\begin{align*}
\left.\delta^{(\epsilon)} \mathcal{L}_{N}^{(\epsilon)}\right|_{\partial \omega^{\star} A \omega^{2}}= & -d_{\alpha \beta \gamma} \partial_{\mu} \omega_{\alpha}^{\star}\left(\theta C_{\gamma \rho \sigma} Z A_{\rho}^{\mu} \omega_{\sigma} \omega_{\beta}+A_{\gamma}^{\mu}\left(-\frac{1}{2} \theta Z\right) C_{\beta \rho \sigma} \omega_{\rho} \omega_{\sigma}\right) \\
& =\partial_{\mu} \omega_{\alpha}^{\star} \theta \omega_{\sigma} \omega_{\beta} A_{\rho}^{\mu}\left(\frac{Z_{\omega} Z}{2 Z^{\prime}}\right)\left(C_{\gamma \rho \sigma} C_{\gamma \alpha \beta}+C_{\gamma \rho \beta} C_{\gamma \sigma \alpha}+C_{\gamma \rho \alpha} C_{\gamma \beta \sigma}\right)=0 \tag{6.5.41}
\end{align*}
$$

Invariance of $\mathcal{L}^{\prime}(\psi, A)$ under the transformations (6.5.34) is ensured by demanding gauge invariance, with gauge parameter

$$
\begin{equation*}
\epsilon_{\alpha}(x)=Z Z^{\prime} \theta \omega_{\alpha}(x), \tag{6.5.42}
\end{equation*}
$$

in terms of rescaled Lie algebra generators and structure constants: $\tilde{t}_{\alpha}=t_{\alpha} / Z^{\prime}, \tilde{C}_{\alpha \beta \gamma}=$ $C_{\alpha \beta \gamma} / Z^{\prime}$. In the canonical basis, such rescalings correspond of course to a renormalization of the gauge coupling constant $\tilde{g}=g / Z^{\prime}$. Putting all together, we can finally write down the most general form of $\mathcal{L}_{N}^{(\epsilon)}$ :

$$
\begin{align*}
\mathcal{L}_{N}^{(\epsilon)}= & -Z_{A} \tilde{F}_{\mu \nu}^{\alpha} \tilde{F}_{\alpha}^{\mu \nu}-Z_{\psi} \bar{\psi} \gamma^{\mu}\left(\partial_{\mu}-i \tilde{t}_{\alpha} A_{\mu}^{\alpha}\right) \psi \\
& +\frac{1}{2} \xi^{\prime} H_{\alpha} H_{\alpha}+\left(\frac{Z_{\omega}}{Z Z^{\prime}}\right) H_{\alpha} \partial^{\mu} A_{\mu}^{\alpha}-Z_{\omega}\left(\partial_{\mu} \omega^{\star}\right)\left(\partial^{\mu} \omega_{\alpha}+\tilde{C}_{\alpha \beta \gamma} A_{\beta}^{\mu} \omega_{\gamma}\right), \tag{6.5.43}
\end{align*}
$$

where schematically $\tilde{F}=d A+\tilde{C} A^{2}$. We notice that this Lagrangian has the same form as the tree level one. Hence, by redefining at level $N$ the fields and couplings in the tree-level action, we can set $\mathcal{L}_{N}^{(\epsilon)}[\chi, 0]=S[\chi]$, namely $\Gamma_{N, \infty}[\chi, 0]=0$. This concludes the proof of the renormalizability of non-abelian gauge theories.

## Chapter 7

## Effective Field Theories

We have so far mostly considered renormalizable quantum field theories, since these theories have simple renormalization properties. Once a finite number of observables is fixed (by experiment), these theories allow for very accurate predictions. It is however clear that renormalizable theories such as the SM cannot be the ultimate theory of nature, since gravity is not included. From this perspective, renormalizable theories are special because they allow us to hide in a redefinition of a finite number of parameters all our ignorance of the UV physics which is beyond the model in consideration. Aside from an improved calculability, there is no conceptual reason to focus only on such a restricted class of theories.

Unless one is so ambitious to try to construct the ultimate, possibly, finite theory of everything, any quantum field theory should be seen as an effective field theory, namely a theory that is reasonably accurate in a given energy regime and is replaced by some other more complete theory at a given UV scale $M$. This happens all the time in physics and is the most efficient way to study a phenomenon keeping only the relevant degrees of freedom. Classical examples are the Fermi theory of electroweak interactions or the pion chiral Lagrangian. These theories are effective, being replaced, at sufficiently high energies, by the SM and QCD, respectively. It is obvious that if we are interested in processes happening at a given scale $E \ll M$ and involving light fields only, with masses much smaller than $M$, the heavy states with mass $\sim M$ cannot be produced, so that the latter can be integrated out. In a path integral approach, the complete 1PI generating functional of light fields is given by

$$
\begin{equation*}
e^{i \Gamma\left(\phi_{l}\right)}=\int_{1 P I} \mathcal{D} \eta_{l} e^{i S_{I R}\left(\phi_{l}+\eta_{l}\right)}, \quad e^{i S_{I R}\left(\phi_{l}\right)}=\int \mathcal{D} \phi_{h} e^{i S_{U V}\left(\phi_{l}, \phi_{h}\right)}, \tag{7.0.1}
\end{equation*}
$$

where $\phi_{l}$ and $\phi_{h}$ schematically denote light and heavy fields. Here $S_{U V}$ and $S_{I R}$ represent
the underlying UV and IR effective actions, respectively. At tree-level, we simply have

$$
\begin{equation*}
S_{I R}^{(0)}\left(\phi_{l}\right)=S_{U V}\left(\phi_{l}, \phi_{h}\left(\phi_{l}\right)\right) \tag{7.0.2}
\end{equation*}
$$

where $\phi_{h}\left(\phi_{l}\right)$ is the classical solution to the heavy field equations of motion. Being the integrated fields heavy, the action admits a well-defined momentum expansion in terms of local operators. We could proceed by computing the one-loop and higher loop effective actions $S_{I R}^{(l)}$ and reconstruct in this way the whole effective action $S_{I R}$. The key point of effective field theory is to replace this procedure by a simpler one, where we compute radiative effects involving the light fields starting from the action

$$
\begin{equation*}
S_{E F T}\left(\phi_{l}\right)=S_{I R}^{(0)}\left(\phi_{l}\right)+\Delta S\left(\phi_{l}\right) \tag{7.0.3}
\end{equation*}
$$

where $\Delta S$ encodes all higher dimensional local operators up to some order, depending on the accuracy we want to achieve, and compatible with the global symmetries of the theory. All the unknown couplings multiplying the higher dimensional local operators are then fixed by comparing given quantities as computed in the UV theory and in the EFT. In principle, one might want to directly compare $S$-matrix elements, but it is often easier to compare 1PI functions, generically off-shell. Note that 1PI functions can be compared at an unphysical point, since by analyticity we are ensured that their on-shell properties are equal. However, this is somewhat improper and too conservative, because physical observables correspond to on-shell $S$ matrix elements and not to the whole Green functions, and different off-shell Green functions can lead to the same on-shell results, as we will see in section 7.8 . We will nevertheless require the off-shell equality of 1PI functions in our examples for simplicity. Note also that in the UV theory 1PI refers to the light fields only. Graphs that can be splitted in two disconnected ones by cutting a propagator of a heavy field should be included.

The procedure of fixing the parameters of the IR theory using the UV theory is called "matching". Of course, matching requires to also perform a computation in the full theory but, as we will see, it is typically much easier to do computations in the EFT and then matching, rather than computing everything in the full theory.

### 7.1 1PI vs Wilsonian Actions

Before considering various examples of use of effective field theories techniques, we would like here to emphasize the difference between the various effective actions introduced so far. ${ }^{1}$ The effective action $\Gamma(\phi)$ defined in eq.(4.1.2) is the generator of all 1PI amplitudes

[^42]in the theory, namely those involving an arbitrary combination of heavy and light fields. In the context of EFT, we also have $\Gamma\left(\phi_{l}\right)$ appearing in the l.h.s. of the first equation in (7.0.1), which generates the 1PI amplitudes associated to the light fields. Another effective action is $S_{I R}$, defined in the l.h.s. of the second equation in (7.0.1). The effective actions $\Gamma(\phi)$ and $\Gamma\left(\phi_{l}\right)$ are shortly denoted 1PI actions, while $S_{I R}$ is called a Wilsonian effective action. The key difference between 1PI and Wilsonian actions is that in the latter the virtual particles are always massive. Correspondingly, while 1PI actions are intrinsically non-local, namely they cannot be written as the integral of some effective Lagrangian density (even with an infinite number of terms), the Wilsonian action always admits such rewriting. The key point can be understood already at tree-level. When we integrate out a scalar particle of mass $M$ at tree-level, roughly speaking we get an effective action term that reads in momentum space as
\[

$$
\begin{equation*}
S_{I R} \supset \int d^{4} p \Phi_{1}(-p) \frac{1}{p^{2}-M^{2}} \Phi_{2}(p), \tag{7.1.1}
\end{equation*}
$$

\]

where $\Phi_{1,2}$ are local functionals of the remaining light fields and $1 /\left(p^{2}-M^{2}\right)$ represents the propagator of the massive particle that has been integrated out. The term is manifestly non-local and cannot be written in configuration space as the integral of a Lagrangian density. At low energies, $p^{2} \ll M^{2}$, we can expand the propagator as

$$
\begin{equation*}
\frac{1}{p^{2}-M^{2}}=-\frac{1}{M^{2}}-\frac{1}{M^{2}} \frac{p^{2}}{M^{2}}+\ldots \tag{7.1.2}
\end{equation*}
$$

Plugging eq.(7.1.2) in eq.(7.1.1), however, gives a (generally infinite) sum of local terms of the form

$$
\begin{equation*}
S_{I R} \supset-\frac{1}{M^{2}} \int d^{4} x \Phi_{1}(x)\left(1+\frac{\square}{M^{2}}+\ldots\right) \Phi_{2}(x) \tag{7.1.3}
\end{equation*}
$$

On the contrary, when massless fields are present, the term in eq.(7.1.1) is intrinsically non-local and eq.(7.1.3) does not apply.

### 7.2 Two Scalars

Consider

$$
\begin{equation*}
\mathcal{L}_{U V}=\frac{1}{2}(\partial H)^{2}-\frac{1}{2} M^{2} H^{2}+\frac{1}{2}(\partial L)^{2}-\frac{1}{2} m^{2} L^{2}-\frac{g}{2} H L^{2}, \tag{7.2.1}
\end{equation*}
$$

where $m \ll M$ and the dimensionful coupling $g$ is assumed to be of $\mathcal{O}(M) .{ }^{2}$ The most general EFT for $L$ reads as follows:

$$
\begin{equation*}
\mathcal{L}_{I R}=\frac{1}{2} Z_{L}(\partial L)^{2}-\frac{1}{2} \tilde{m}^{2} L^{2}-\frac{\lambda}{4!} L^{4}+\text { h.d.o. }, \tag{7.2.2}
\end{equation*}
$$

[^43]where h.d.o. stands for higher dimensional operators. The tree-level integration of $H$ gives $H=-g L^{2} /\left(2 M^{2}\right)+\mathcal{O}\left(p^{2} / M^{4}\right)$. Plugged back in eq. (7.2.1) gives
\[

$$
\begin{equation*}
\mathcal{L}_{I R}^{(0)}=\frac{1}{2}(\partial L)^{2}-\frac{1}{2} m^{2} L^{2}-\frac{\lambda}{4!} L^{4}+\mathcal{O}\left(g^{2} M^{-4}\right) \tag{7.2.3}
\end{equation*}
$$

\]

so that tree-level matching gives

$$
\begin{equation*}
Z_{L}=1+\mathcal{O}\left(g^{2}\right), \quad \tilde{m}^{2}=m^{2}+\mathcal{O}\left(g^{2}\right), \quad \lambda=-\frac{3 g^{2}}{M^{2}}+\mathcal{O}\left(g^{4}\right) \tag{7.2.4}
\end{equation*}
$$

It is important to emphasize that the factors $Z_{L}, \tilde{m}^{2}$, etc. in eq. (7.2.4) are the analogues of $(1+\delta Z), m^{2}+\delta m^{2}$, etc. defined in eq. (5.1.5) in the Wilsonian RG flow. They are finite terms, which at the quantum level indicate how the EFT fields, masses etc. differ from their UV counterparts to compensate for the different UV behaviour of the two theories, so that low-energy observables match in the two descriptions. They should not be confused with the usual counter-terms needed to subtract divergences, which we will never explicitly write, always fixing them by an $\overline{\mathrm{MS}}$ subtraction scheme.

We can directly use $\mathcal{L}_{I R}^{(0)}$ as effective theory to compute Green functions for energies much smaller than $M$. For instance, the $\overline{\mathrm{MS}}$ renormalized four-point 1PI function reads (see eq. (5.6.4)):

$$
\begin{equation*}
\Gamma^{(4)}(s, t, u)=-\lambda(\mu)+\frac{\lambda^{2}(\mu)}{32 \pi^{2}} \int_{0}^{1} d x\left[\log \frac{m^{2}-s x(1-x)}{\mu^{2}}+(s \rightarrow t)+(s \rightarrow u)\right] . \tag{7.2.5}
\end{equation*}
$$

By taking $s \simeq t \simeq u \simeq \mu$ we get $\Gamma^{(4)}(\mu) \simeq-\lambda(\mu)$, where

$$
\begin{equation*}
\lambda^{-1}(\mu)=\lambda^{-1}\left(\mu_{0}\right)-\frac{3}{16 \pi^{2}} \log \left(\frac{\mu}{\mu_{0}}\right) \tag{7.2.6}
\end{equation*}
$$

is the usual one-loop running in the $\phi^{4}$ theory, assuming $\mu, \mu_{0}>m$. The natural scale where the effective theory parameter $\lambda$ should be matched, at one-loop level, with the underlying UV parameters $g$ and $M$ is at the scale $\mu_{0}=M$, since this is the energy scale boundary between the two theories. This is understood by comparing the 1PI four point amplitude $\Gamma^{(4)}$. The following diagrams contribute to $\Gamma^{(4)}$ in the UV theory (continuous and dashed lines correspond to light and heavy fields, respectively). At leading order in
an $1 / M$ expansion and for zero external momentum, we get

$$
\begin{align*}
+ \text { perms. } & =\left(\frac{1}{2} \times 3\right)(-i g)^{4} \mu^{2 \epsilon} \int \frac{d^{d} k}{(2 \pi)^{d}}\left(\frac{i}{k^{2}-m^{2}}\right)^{2}\left(\frac{-i}{M^{2}}\right)^{2} \\
& =\frac{3 i g^{4}}{32 \pi^{2} M^{4}} \log \frac{m^{2}}{\mu^{2}}, \\
+ \text { perms. } & =(6)(-i g)^{4} \mu^{2 \epsilon} \int \frac{d^{d} k}{(2 \pi)^{d}}\left(\frac{i}{k^{2}-m^{2}}\right)^{2} \frac{-i}{M^{2}} \frac{i}{k^{2}-M^{2}} \\
& =\frac{3 i g^{4}}{8 \pi^{2} M^{4}}\left(1+\log \frac{m^{2}}{M^{2}}\right),  \tag{7.2.7}\\
+ \text { perms. } & =(6)(-i g)^{4} \mu^{2 \epsilon} \int \frac{d^{d} k}{(2 \pi)^{d}}\left(\frac{i}{k^{2}-m^{2}}\right)^{2}\left(\frac{i}{k^{2}-M^{2}}\right)^{2} \\
& =\frac{6 i g^{4}}{8 \pi^{2} M^{4}}\left(1+\frac{1}{2} \log \frac{m^{2}}{M^{2}}\right),
\end{align*}
$$

where in parenthesis we have denoted the geometrical factor (permutations included) of the graph. Recall that $\Gamma^{(4)}$ is 1PI in the light fields only, and hence the first graph in eq.(7.2.7), reducible when cutting the propagator of the heavy field, should be included in the computation.

Including the tree-level term, we get

$$
\begin{equation*}
\Gamma^{(4)}(0)=\frac{3 g^{2}(\mu)}{M^{2}}+\frac{3 g^{4}(\mu)}{8 \pi^{2} M^{4}}\left(3+2 \log \frac{m^{2}}{M^{2}}+\frac{1}{4} \log \frac{m^{2}}{\mu^{2}}\right) . \tag{7.2.8}
\end{equation*}
$$

Comparing eq. (7.2.5) (at zero momentum) with eq. (7.2.8) and using the tree-level relation (7.2.4) for $\lambda$, we get

$$
\begin{equation*}
\lambda(\mu)=-\frac{3 g^{2}(\mu)}{M^{2}}-\frac{3 g^{4}}{8 \pi^{2} M^{4}}\left(3+2 \log \frac{\mu^{2}}{M^{2}}\right) . \tag{7.2.9}
\end{equation*}
$$

Notice how all the (potentially large) logs involving the light mass $m^{2}$ cancelled from eq. (7.2.9). This is not a coincidence. All the IR properties of the UV theory (as well as large IR effects) are reliably described by the EFT by construction, so that the matching equations are always regular in the IR. It is quite clear from eq. (7.2.9) that the best scale to match the two theories is $\mu=M$, in which case we have

$$
\begin{equation*}
\lambda(M)=-\frac{3 g^{2}(M)}{M^{2}}-\frac{9 g^{4}}{8 \pi^{2} M^{4}} . \tag{7.2.10}
\end{equation*}
$$

Taking $\mu_{0}=M$ in eq. (7.2.6) and using eq. (7.2.10), will allow us to take into account possible large logs of the form $\log (\mu / M)$. Notice that the $g^{4} / M^{4}$ term in eq. (7.2.10) can safely be neglected at one-loop level, starting to be relevant only at two-loop level, if NLL want to be resummed. In other words, no one-loop computation of the 1PI 4-point function
would have been needed in matching the theory at $\mu \simeq M$ and we would have been able to resum the leading $\log (\mu / M)$ terms without performing any radiative computation in the UV theory! In general, an $l$-loop computation in the effective theory requires an $l$ - 1 -loop matching computation in the UV theory. This example clearly shows the usefulness of effective theories even in such a simple set-up.

### 7.3 Yukawa Theory I: Heavy Scalar, Light Fermion

Consider a light fermion field coupled to a massive scalar by means of a Yukawa-like coupling. The UV and IR Lagrangians are

$$
\begin{align*}
\mathcal{L}_{U V} & =\frac{1}{2}(\partial \phi)^{2}-\frac{1}{2} M^{2} \phi^{2}+\bar{\psi}(i \not \partial-m) \psi-g \phi \bar{\psi} \psi  \tag{7.3.1}\\
\mathcal{L}_{I R} & =Z_{\psi} \bar{\psi} i \not \partial \psi-\tilde{m} \bar{\psi} \psi+\frac{\lambda}{2} \bar{\psi} \psi \bar{\psi} \psi+\text { h.d.o. } \tag{7.3.2}
\end{align*}
$$

The tree-level integration of $\phi$ gives $\phi=-g \bar{\psi} \psi / M^{2}+\mathcal{O}\left(p^{2} / M^{4}\right)$. Plugging back in eq. (7.3.1) gives

$$
\begin{equation*}
Z_{\psi}=1+\delta Z_{\psi}, \quad \tilde{m}=m+\delta_{m}, \quad \lambda=\frac{g^{2}}{M^{2}} \tag{7.3.3}
\end{equation*}
$$

Let us now compute the 1PI 2-point function $\Gamma^{(2)}(p)$ in both the UV and the IR EFT at one-loop level. In the UV theory the relevant one-loop diagram is

$$
\begin{aligned}
& =\frac{i g^{2}}{16 \pi^{2}} \int_{0}^{1} d x(\not p(1-x)+m) \log \left(\frac{\mu^{2}}{-p^{2} x(1-x)+m^{2} x+(1-x) M^{2}}\right) .
\end{aligned}
$$

Expanding for $M \gg m, p$, we get

$$
\begin{align*}
\Gamma^{(2)}(p) & =\not p-m-\frac{g^{2}}{16 \pi^{2}}\left[\not p\left(\frac{1}{2} \log \frac{M^{2}}{\mu^{2}}-\frac{1}{4}+\frac{m^{2}}{2 M^{2}}-\frac{p^{2}}{6 M^{2}}+\mathcal{O}\left(\frac{1}{M^{4}}\right)\right)\right. \\
& \left.+m\left(\log \frac{M^{2}}{\mu^{2}}-1-\frac{m^{2}}{M^{2}} \log \frac{m^{2}}{M^{2}}-\frac{p^{2}}{2 M^{2}}+\mathcal{O}\left(\frac{1}{M^{4}}\right)\right)\right] \tag{7.3.5}
\end{align*}
$$

In the effective theory, the relevant one-loop diagram is

$$
\begin{align*}
& \overbrace{p}^{k}=(-1)\left(i \lambda \mu^{\epsilon}\right) \int \frac{d^{d} k}{(2 \pi)^{d}}\left(\frac{i(\not k+m)}{k^{2}-m^{2}}+\frac{i \operatorname{Tr}(\not k+m)}{k^{2}-m^{2}}\right)  \tag{7.3.6}\\
= & \frac{-5 i \lambda m^{3}}{16 \pi^{2}}\left(1+\log \frac{\mu^{2}}{m^{2}}\right), \tag{7.3.7}
\end{align*}
$$

where the two terms in square brackets indicate the two different contractions among the four fermions in $\bar{\psi} \psi \bar{\psi} \psi$, and we have taken $\operatorname{Tr} I=4$ (rather than $d$ ). The 1PI 2-point function $\Gamma^{(2)}(p)$ in the IR reads then

$$
\begin{equation*}
\Gamma^{(2)}(p)=\left(1+\delta Z_{\psi}\right) \not p-\left(m+\delta_{m}\right)-\frac{5 \lambda m^{3}}{16 \pi^{2}}\left(1+\log \frac{\mu^{2}}{m^{2}}\right) \tag{7.3.8}
\end{equation*}
$$

Matching eqs. (7.3.5) and (7.3.8) gives

$$
\begin{align*}
\delta Z_{\psi}(\mu) & =-\frac{g^{2}}{16 \pi^{2}}\left(\frac{1}{2} \log \frac{M^{2}}{\mu^{2}}-\frac{1}{4}+\frac{m^{2}}{2 M^{2}}+\mathcal{O}\left(\frac{1}{M^{4}}\right)\right) \\
\delta_{m}(\mu) & =-\frac{5 \lambda m^{3}}{16 \pi^{2}}\left(1+\log \frac{\mu^{2}}{m^{2}}\right)+\frac{m g^{2}}{16 \pi^{2}}\left(\log \frac{M^{2}}{\mu^{2}}-1-\frac{m^{2}}{M^{2}} \log \frac{m^{2}}{M^{2}}+\mathcal{O}\left(\frac{1}{M^{4}}\right)\right) . \\
& =-\frac{m g^{2}}{16 \pi^{2}}\left(1+\log \frac{\mu^{2}}{M^{2}}\right)\left(1+\frac{m^{2}}{M^{2}}\right)-\frac{4 m^{3} g^{2}}{16 \pi^{2} M^{2}}\left(1+\log \frac{\mu^{2}}{m^{2}}\right), \tag{7.3.9}
\end{align*}
$$

where in the last line we have used the tree-level relation (7.3.3) for $\lambda$. It is clear from eq. (7.3.9) that the matching is best performed at $\mu=M$. Matching the $p^{2} / M^{2}$ terms in eq. (7.3.5) would require the addition of the higher dimensional operators of the form $\bar{\psi} \not \partial \square \psi$ or $\bar{\psi} \square \psi$. We neglect them, since their effect is small, suppressed by $p^{2} / M^{2}$ times a loop factor. The physical fermion mass is given, at one-loop level, by

$$
\begin{equation*}
m_{\text {phys }}=m(M)\left[1-\frac{g^{2}}{16 \pi^{2}}\left(\frac{5}{4}+\frac{9 m^{2}}{2 M^{2}}+\frac{4 m^{2}}{M^{2}} \log \frac{M^{2}}{m^{2}}\right)\right] \tag{7.3.10}
\end{equation*}
$$

### 7.4 Yukawa Theory II: Heavy Fermion, Light Scalar

Consider now the opposite case with an heavy fermion field coupled to a light scalar:

$$
\begin{align*}
\mathcal{L}_{U V} & =\frac{1}{2}(\partial \phi)^{2}-\frac{1}{2} m^{2} \phi^{2}+\bar{\psi}(i \not \partial-M) \psi-g \phi \bar{\psi} \psi  \tag{7.4.1}\\
\mathcal{L}_{I R} & =\frac{1}{2} Z_{\phi}(\partial \phi)^{2}-\frac{1}{2} \tilde{m}^{2} \phi^{2}+\frac{\lambda_{3}}{3!} \phi^{3}+\frac{\lambda_{4}}{4!} \phi^{4}+\text { h.d.o. } \tag{7.4.2}
\end{align*}
$$

The tree-level integration of $\psi$ gives $\psi=0$, so that

$$
\begin{equation*}
Z_{\phi}=1+\delta Z_{\phi}, \quad \tilde{m}^{2}=m^{2}+\delta_{m^{2}}, \quad \lambda_{3}=\mathcal{O}\left(g^{3}\right), \quad \lambda_{4}=\mathcal{O}\left(g^{4}\right) \tag{7.4.3}
\end{equation*}
$$

Let us again match the 1PI 2-point function in the UV and in the IR theory. In the UV we have

$$
\begin{align*}
& \text { ( }=-1)(-i g)^{2} \mu^{\epsilon} \operatorname{Tr} \int \frac{d^{d} k}{(2 \pi)^{d}} \frac{i(\not k+M)}{k^{2}-M^{2}} \frac{i(p p+\not k+M)}{(k+p)^{2}-M^{2}} \\
& =-\frac{4 i g^{2}}{16 \pi^{2}} \int_{0}^{1} d x\left(M^{2}-p^{2} x(1-x)\right)\left[3 \log \left(\frac{\mu^{2}}{-p^{2} x(1-x)+M^{2}}\right)-5\right], \tag{7.4.4}
\end{align*}
$$

taking $\operatorname{Tr} I=4$ (rather than $d$ ). Expanding for $M \gg m, p$, we get

$$
\begin{align*}
\Gamma^{(2)}(p) & =p^{2}-m^{2}-\frac{4 g^{2}}{16 \pi^{2}}\left[p^{2}\left(-\frac{1}{2} \log \frac{M^{2}}{\mu^{2}}+\frac{4}{3}-\frac{p^{2}}{20 M^{2}}+\mathcal{O}\left(\frac{1}{M^{4}}\right)\right)\right. \\
& \left.+M^{2}\left(-3 \log \frac{M^{2}}{\mu^{2}}-5+\mathcal{O}\left(\frac{1}{M^{4}}\right)\right)\right] . \tag{7.4.5}
\end{align*}
$$

In the EFT we simply have

$$
\begin{equation*}
\Gamma^{(2)}(p)=p^{2}-m^{2}+p^{2} \delta Z_{\phi}-\delta_{m^{2}} \tag{7.4.6}
\end{equation*}
$$

Matching eqs. (7.4.5) and (7.4.6) gives

$$
\begin{align*}
\delta Z_{\phi}(\mu) & =-\frac{4 g^{2}}{16 \pi^{2}}\left(-\frac{1}{2} \log \frac{M^{2}}{\mu^{2}}+\frac{4}{3}\right) \\
\delta_{m^{2}}(\mu) & =-\frac{4 g^{2} M^{2}}{16 \pi^{2}}\left(3 \log \frac{M^{2}}{\mu^{2}}+5\right) \tag{7.4.7}
\end{align*}
$$

The physical scalar mass is given by

$$
\begin{equation*}
m_{\text {phys }}^{2}=m^{2}(M)+\frac{g^{2}}{16 \pi^{2}} M^{2}\left(-20+\frac{16}{3} \frac{m^{2}}{M^{2}}\right) \tag{7.4.8}
\end{equation*}
$$

### 7.5 Naturalness and the Hierarchy Problem

The scalar mass correction in eq. (7.4.8) is of $\mathcal{O}\left(M^{2} g^{2} /\left(16 \pi^{2}\right)\right)$. For $M \gg m$, the ratio $(M / m)^{2}$ can overcome the one-loop suppression so that the radiative term is generically much bigger than the tree-level one. If we want to keep the physical scalar mass small, a fine-tuning between the $\overline{M S}$ mass term $m^{2}(M)$ and the one-loop correction is needed (of course this readjustment is needed at each order in perturbation theory). This is not a conceptual problem, and there is nothing wrong in principle to do that, but it is nevertheless unpleasant to have in the EFT a parameter that is so sensitive to the UV physics. Even a small change in the UV theory (say, $g \rightarrow g+\delta g$ in our toy example) will give rise to a large radiative correction to $m_{\text {phys. }}^{2}$ and to a new readjustment. This is the typical example of an "hierarchy problem" or "naturalness problem", namely the problem of explaining why a parameter in an EFT Lagrangian is much smaller than its expected value at the quantum level. ${ }^{3}$ Natural parameters are often defined as follows:

[^44]dimensionless couplings should be of order one and dimensionful couplings should be of order of the largest mass scale in the theory, to the appropriate power. Exceptions arise if a symmetry is restored when a coupling (dimensionless or not) vanishes, in which case it is natural to have that coupling arbitrarily small.

Let us check, using our results above, that this definition makes sense. In the scalar theory of section 7.2 , when $g=\mathcal{O}(M)$ (natural value), we get that the dimensionless coupling $\lambda$ in the EFT is of $\mathcal{O}\left(M^{2} / M^{2}\right)=\mathcal{O}(1)$. Contrary to the scalar case, the fermion mass correction in eq. (7.3.10) is proportional to the mass itself. This implies that, for $m \rightarrow 0$, all loop corrections vanish as well. Hence, light fermion masses are natural. According to our general definition of natural parameters, some symmetry should be restored when $m \rightarrow 0$. This is indeed what happens, the symmetry being the chiral symmetry $\psi \rightarrow \gamma_{5} \psi$ (combined with the discrete $\mathbf{Z}_{2}$ inversion symmetry $\phi \rightarrow-\phi$ in the UV theory). No symmetry is instead restored when the scalar mass term goes to zero, so that in this case we have a naturalness problem.

Despite hierarchy problems do not lead to real inconsistencies and we can live with them, they have been the main driving force in going beyond the SM, since the Higgs boson mass term is unnatural. Similarly, one of the main problems in theoretical physics today is provided by another hierarchy problem, which is why the vacuum energy we measure is so smaller than its natural value.

### 7.6 Non-Leptonic Decays

As a final example, we consider a commonly used EFT of the weak interactions, by means of four-fermion operators. The charged electroweak current in the SM contains the terms

$$
\begin{equation*}
\mathcal{L}_{S M} \supset \frac{g}{\sqrt{2}} W_{\mu}^{+}\left(V_{u d} \bar{u}_{L} \gamma^{\mu} d_{L}+V_{u s} \bar{u}_{L} \gamma_{\mu} s_{L}\right)+h . c . \tag{7.6.1}
\end{equation*}
$$

Integrating out the $W$ gives, among others, the flavour changing four-Fermi interaction term

$$
\begin{equation*}
-\frac{4 G_{F}}{\sqrt{2}} V_{u s} V_{d u}^{\star}\left(\bar{u}_{L} \gamma^{\mu} s_{L}\right)\left(\bar{d}_{L} \gamma_{\mu} u_{L}\right) \tag{7.6.2}
\end{equation*}
$$

with the matching relation

$$
\begin{equation*}
\frac{G_{F}}{\sqrt{2}}=\frac{g^{2}}{8 m_{W}^{2}} . \tag{7.6.3}
\end{equation*}
$$

The most general EFT dimension six interaction terms that can contribute to $K$ decays are

$$
\begin{equation*}
\mathcal{L}_{I R}=-c_{1} O_{1}-c_{2} O_{2}+h . c . \tag{7.6.4}
\end{equation*}
$$



Figure 7.1: One-loop graphs contributing to the renormalization of the four-Fermi operator (7.6.2). All external momenta are vanishing.
where

$$
\begin{equation*}
O_{1}=\left(\bar{u}_{L} \gamma^{\mu} s_{L}\right)\left(\bar{d}_{L} \gamma_{\mu} u_{L}\right), \quad O_{2}=\left(\bar{u}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{d}_{L} \gamma_{\mu} s_{L}\right) \tag{7.6.5}
\end{equation*}
$$

Tree-level matching gives

$$
\begin{equation*}
c_{1}\left(m_{W}\right)=\frac{g^{2}}{2 m_{W}^{2}} V_{u s} V_{d u}^{\star}\left(m_{W}\right), \quad c_{2}\left(m_{W}\right)=0 \tag{7.6.6}
\end{equation*}
$$

Although at tree-level only $O_{1}$ appears, symmetry arguments do not forbid $O_{2}$, which is expected to arise at the quantum level. Operators like (7.6.5) are relevant for the decays of the strange mesons, such as the $K$ 's. The natural scale where the operators $O_{1,2}$ should be renormalized is $\mathcal{O}\left(m_{K}\right)$, being this the scale of the decay process. It is important to compute the RG flow of the couplings $c_{i}=c_{i}(\mu), i=1,2$, and see its effect. The main contribution is given by the QCD corrections. In what follows we then compute the QCD effects on the RG flow of $c_{1}$ and $c_{2}$.

Recall that at the linear level, the RG evolution of the couplings $c_{1,2}$ is fixed by the anomalous dimensions of the corresponding operators (see eq. (5.9.13)):

$$
\begin{equation*}
\mu \frac{d c_{i}}{d \mu}=\beta_{i}=\gamma_{i j} c_{j} \tag{7.6.7}
\end{equation*}
$$

where $\gamma_{i j}$ is the $2 \times 2$ matrix of anomalous dimensions of the composite operators $O_{1}$ and $O_{2}$ which, as we will see, mix under renormalization. We have then to determine $\gamma_{i j}$. By definition,

$$
\begin{equation*}
O_{i}^{B}=Z_{i j}(\mu) O_{j}(\mu), \tag{7.6.8}
\end{equation*}
$$

where $O_{i}^{B}$ are the bare composite operators. In analogy to what done at the end of section 5.9 , where we computed the anomalous dimensions of the composite operator $\phi^{2}$ in the
$\lambda \phi^{4}$ theory by demanding the finiteness of the connected Green function $\left\langle\phi^{2} \phi \phi\right\rangle$ at zero momentum, we might determine $Z_{i j}$ by demanding that the connected Green functions $G_{i}^{(4,1)} \equiv\left\langle O_{i} \bar{u}_{L} u_{L} \bar{s}_{L} d_{L}\right\rangle$ are finite at zero momentum. The Green functions $G_{i}^{(4,1)}$ carry four spinor indices, one for each elementary quark field. In order to simplify the notation, we will multiply $G_{i}^{(4,1)}$ by the wave functions of the external fermion fields and denote them by the same symbol of the corresponding fermion field $u_{L}, d_{L}$ and $s_{L}$.

The relevant 1PI one-loop graphs renormalizing $O_{1}$ are depicted in fig. 7.1. A similar set of graphs, with $d \leftrightarrow u$ in the external lines, renormalizes $O_{2}$. Due to the conservation of the underlying electroweak currents, the diagrams (1) and (2) in fig. 7.1 precisely compensate for the non 1PI one-loop graphs correcting the external quark lines so that we do not need to compute them. This is best understood by noticing that $G_{i B}^{(4,1)}=Z_{i j} G_{j}^{(4,1)} / Z_{q}^{2}$, where $\sqrt{Z_{q}}$ is the wave function renormalization of the elementary quark fields, flavourindependent when only strong interactions are taken into account. The factors $Z_{q}^{2}$ compensate then for the divergences arising from the diagrams (1) and (2). In the following we neglect all quark masses and choose the Feynman gauge $\xi=1$ for the gluon propagator. Keeping only the divergent terms, diagram (3) gives

$$
\begin{align*}
(3) & =\left(i g_{c}\right)^{2} \mu^{\epsilon} \int \frac{d^{d} k}{(2 \pi)^{d}}\left(\bar{d}_{L} \gamma^{\mu} \frac{i \nmid k}{k^{2}} \gamma^{\nu} t^{a} u_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} \frac{-i \not k}{k^{2}} \gamma_{\nu} t^{a} s_{L}\right) \frac{-i}{k^{2}} \\
& =-\frac{g_{c}^{2}}{32 \pi^{2} \epsilon}\left(\bar{d}_{L} \gamma^{\mu} \gamma^{\rho} \gamma^{\nu} t^{a} u_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} \gamma_{\rho} \gamma_{\nu} t^{a} s_{L}\right)+\text { finite terms } . \tag{7.6.9}
\end{align*}
$$

Similarly, we get

$$
\begin{equation*}
(4)=(3), \quad(5)=(6)=\frac{g_{c}^{2}}{32 \pi^{2} \epsilon}\left(\bar{d}_{L} \gamma^{\mu} \gamma^{\rho} \gamma^{\nu} t^{a} u_{L}\right)\left(\bar{u}_{L} \gamma_{\nu} \gamma_{\rho} \gamma_{\mu} t^{a} s_{L}\right)+\text { finite terms } \tag{7.6.10}
\end{equation*}
$$

Eqs. (7.6.9) and (7.6.10) are best written using certain identities in both color and spinor indices, identities that are derived in detail in subsection 7.6.1. Using the results in that section, we have

$$
\begin{align*}
t_{i j}^{a} t_{k l}^{a} & =\frac{1}{2}\left(\delta_{i l} \delta_{j k}-\frac{1}{3} \delta_{i j} \delta_{k l}\right),  \tag{7.6.11}\\
\psi_{2 L} \bar{\psi}_{1 L} & =\frac{1}{2}\left(\bar{\psi}_{1 L} \gamma^{\mu} \psi_{2 L}\right) \gamma_{\mu} P_{R} \tag{7.6.12}
\end{align*}
$$

Starting from eq.(7.6.9), we first use eq.(7.6.12) to write

$$
\begin{align*}
K \equiv\left(\bar{d}_{L}^{i} \gamma^{\mu} \gamma^{\rho} \gamma^{\nu} t_{i j}^{a} u_{L}^{j}\right)\left(\bar{u}_{L}^{k} \gamma_{\mu} \gamma_{\rho} \gamma_{\nu} t_{k l}^{a} s_{L}^{l}\right) & =\frac{1}{2}\left(\bar{d}_{L}^{i} \gamma^{\mu} \gamma^{\rho} \gamma^{\nu} \gamma^{\alpha} P_{R} \gamma_{\mu} \gamma_{\rho} \gamma_{\nu} s_{L}^{l}\right) t_{a}^{k l} t_{a}^{i j} \\
\times\left(\bar{u}_{L}^{k} \gamma_{\alpha} u_{L}^{j}\right) & =-16\left(\bar{u}_{L}^{k} \gamma_{\alpha} u_{L}^{j}\right)\left(\bar{d}_{L}^{i} \gamma^{\alpha} s_{L}^{l}\right) t_{a}^{i j} t_{a}^{k l} \tag{7.6.13}
\end{align*}
$$

where in the second equality we have used the gamma matrix algebra relation

$$
\begin{equation*}
\gamma^{\mu} \gamma^{\rho} \gamma^{\nu} \gamma^{\alpha} \gamma_{\mu}=-2 \gamma^{\alpha} \gamma^{\nu} \gamma^{\rho} \tag{7.6.14}
\end{equation*}
$$

We can now apply eq.(7.6.11) to get

$$
\begin{align*}
K= & -16\left(\bar{u}_{L}^{k} \gamma^{\alpha} u_{L}^{j}\right)\left(\bar{d}_{L}^{i} \gamma_{\alpha} s_{L}^{l}\right) \frac{1}{2}\left(\delta_{i l} \delta_{j k}-\frac{1}{3} \delta_{i j} \delta_{k l}\right)=-8\left(\bar{u}_{L} \gamma^{\alpha} u_{L}\right)\left(\bar{d}_{L} \gamma_{\alpha} s_{L}\right) \\
& +\frac{8}{3}\left(\bar{u}_{L}^{k} \gamma^{\alpha} u_{L}^{i}\right)\left(\bar{d}_{L}^{i} \gamma_{\alpha} s_{L}^{k}\right) . \tag{7.6.15}
\end{align*}
$$

Applying again eq.(7.6.12) to the second term in eq.(7.6.15) gives

$$
\begin{align*}
K & =-8\left(\bar{u}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{d}_{L} \gamma_{\mu} s_{L}\right)+\frac{8}{3}\left(\frac{1}{2}\right)\left(\bar{d}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{u}_{L} \gamma^{\alpha} \gamma_{\mu} P_{R} \gamma_{\alpha} s_{L}\right) \\
& =-8\left(\bar{u}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{d}_{L} \gamma_{\mu} s_{L}\right)-\frac{8}{3}\left(\bar{d}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} s_{L}\right), \tag{7.6.16}
\end{align*}
$$

where we have the relation

$$
\begin{equation*}
\gamma^{\alpha} \gamma_{\mu} \gamma_{\alpha}=-2 \gamma_{\mu} \tag{7.6.17}
\end{equation*}
$$

The second row in eq.(7.6.18) is our final expression of $K$. As can be seen, we have started with the four fermion operator $O_{1}$ with the schematic form $(\bar{u} s)(\bar{d} u)$ and ended up with having generated also the form $(\bar{u} u)(\bar{d} s)$, corresponding to the operator $O_{2}$. Identical manipulations can be performed starting from the term (5) in eq.(7.6.10) and will not be reported. Summarizing, we have

$$
\begin{align*}
& \left(\bar{d}_{L} \gamma^{\mu} \gamma^{\rho} \gamma^{\nu} t^{a} u_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} \gamma_{\rho} \gamma_{\nu} t^{a} s_{L}\right)=-8\left(\bar{d}_{L} \gamma^{\mu} s_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} u_{L}\right)-\frac{8}{3}\left(\bar{d}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} s_{L}\right), \\
& \left(\bar{d}_{L} \gamma^{\mu} \gamma^{\rho} \gamma^{\nu} t^{a} u_{L}\right)\left(\bar{u}_{L} \gamma_{\nu} \gamma_{\rho} \gamma_{\mu} t^{a} s_{L}\right)=-2\left(\bar{d}_{L} \gamma^{\mu} s_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} u_{L}\right)-\frac{2}{3}\left(\bar{d}_{L} \gamma^{\mu} u_{L}\right)\left(\bar{u}_{L} \gamma_{\mu} s_{L}\right) \cdot(7 \tag{7.6.18}
\end{align*}
$$

Collecting all terms and taking into account that $O_{1}$ and $O_{2}$ are related by an odd exchange of fermion operators, we get $\left(\alpha_{c}=g_{c}^{2} /(4 \pi)\right)$

$$
\begin{equation*}
O_{1, B}=O_{1}(\mu)+\frac{\alpha_{c}}{2 \pi \epsilon}\left[O_{1}(\mu)-3 O_{2}(\mu)\right], \quad O_{2, B}=O_{2}(\mu)+\frac{\alpha_{c}}{2 \pi \epsilon}\left[O_{2}(\mu)-3 O_{1}(\mu)\right], \tag{7.6.19}
\end{equation*}
$$

from which we compute the wave-function renormalizaiton matrix $Z_{i j}$ :

$$
Z_{i j}=\delta_{i j}+\frac{\alpha_{c}}{2 \pi \epsilon}\left(\begin{array}{cc}
1 & -3  \tag{7.6.20}\\
-3 & 1
\end{array}\right)+\mathcal{O}\left(\alpha_{c}^{2}\right)
$$

The matrix of anomalous dimension is now easily computed from eq.(5.9.5):

$$
\gamma_{i j}=\frac{\alpha_{c}}{2 \pi}\left(\begin{array}{cc}
-1 & 3  \tag{7.6.21}\\
3 & -1
\end{array}\right)+\mathcal{O}\left(\alpha_{c}^{2}\right)
$$

where we have used the tree-level $d$-dimensional QCD $\beta$-function $\beta\left(\alpha_{c}\right)=-\epsilon \alpha_{c}+\mathcal{O}(1) .{ }^{4}$ The operator basis in which the anomalous dimension matrix is diagonal is given by the

[^45]operator combinations ${ }^{5}$
\[

$$
\begin{equation*}
O_{1 / 2}=\frac{O_{1}-O_{2}}{2}, \quad O_{3 / 2}=\frac{O_{1}+O_{2}}{2} \tag{7.6.22}
\end{equation*}
$$

\]

with eigenvalues

$$
\begin{equation*}
\gamma_{1 / 2}=-\frac{2 \alpha_{c}}{\pi}, \quad \gamma_{3 / 2}=\frac{\alpha_{c}}{\pi} . \tag{7.6.23}
\end{equation*}
$$

We then have $c_{i} O_{i}=c_{1 / 2} O_{1 / 2}+c_{3 / 2} O_{3 / 2}$, with $c_{1 / 2}=c_{1}-c_{2}, c_{3 / 2}=c_{1}+c_{2}$. In this new basis the RG flow for the couplings reads

$$
\begin{equation*}
\mu \frac{d c_{i}}{d \mu}=\gamma_{i} c_{i}, \quad i=1 / 2,3 / 2 \tag{7.6.24}
\end{equation*}
$$

The equations are easily integrated by writing $\mu d c_{i} / d \mu=d c_{i} / d \alpha_{c} \beta_{c}$, where $\beta_{c}$ is the (one-loop) QCD $\beta$-function

$$
\begin{equation*}
\beta_{c}=-\frac{\alpha_{c}^{2}}{2 \pi} b_{0} \quad b_{0}=11-\frac{2}{3} n_{f} . \tag{7.6.25}
\end{equation*}
$$

We get

$$
\begin{equation*}
c_{1 / 2}(\mu)=c_{1 / 2}\left(\mu_{0}\right)\left(\frac{\alpha_{c}(\mu)}{\alpha_{c}\left(\mu_{0}\right)}\right)^{\frac{4}{b_{0}}}, \quad c_{3 / 2}(\mu)=c_{3 / 2}\left(\mu_{0}\right)\left(\frac{\alpha_{c}(\mu)}{\alpha_{c}\left(\mu_{0}\right)}\right)^{-\frac{2}{b_{0}}} \tag{7.6.26}
\end{equation*}
$$

The integration constants $c_{i}\left(\mu_{0}\right)$ are fixed by taking $\mu_{0}=m_{W}$ and matching by using eq. (7.6.6): $c_{3 / 2}\left(m_{W}\right)=c_{1 / 2}\left(m_{W}\right)=g^{2} V_{u s} V_{d u}^{\star} /\left(2 m_{W}^{2}\right) \equiv c_{0}$. By taking $\mu \sim m_{K}$, we can estimate the effect of the QCD LL corrections. For $n_{f}=4,{ }^{6}$ one roughly gets $c_{1 / 2}\left(m_{K}\right) \simeq$ $2 c_{0}, c_{3 / 2}\left(m_{K}\right) \simeq 0.7 c_{0}$, with an enhancement by about a factor 3 of the isospin $1 / 2$ operator with respect to the $3 / 2$ one, showing that QCD corrections are far from being negligible.

### 7.6.1 Useful Color and Spinor identities

In this subsection we provide the details to derive eqs.(7.6.11) and (7.6.12). Let us first consider the color indices. When the combination of color generators $t_{i j}^{a} t_{k l}^{a}$ is summed over all generators, the indices $i, j, k, l$ can only be carried by the group space metric $\delta_{i j}$. Hence we must have

$$
\begin{equation*}
\sum_{a} t_{i j}^{a} t_{k l}^{a}=A \delta_{i l} \delta_{j k}+B \delta_{i j} \delta_{l k}, \tag{7.6.27}
\end{equation*}
$$

[^46]with $A$ and $B$ coefficients to be determined. Multiplying by $\delta_{i j}$ (or by $\delta_{k l}$ ) and by $\delta_{j k}$ (or by $\delta_{i l}$ ) we get, for $S U(N)$,
\[

$$
\begin{align*}
0 & =\delta_{l k}(A+N B) \\
C_{2}(r) \delta_{i l} \equiv \sum_{a}\left(t^{a} t^{a}\right)_{i l} & =\delta_{i l}(N A+B) \tag{7.6.28}
\end{align*}
$$
\]

where $C_{2}(r)$ is the quadratic Casimir group invariant, that depends on the representation of the generators $t^{a}$. For $S U(N)$ generators in the fundamental representation we have

$$
\begin{equation*}
C_{2}(\text { fund } .)=\frac{N^{2}-1}{2 N} \tag{7.6.29}
\end{equation*}
$$

The solution of eqs.(7.6.28) is

$$
\begin{equation*}
A=\frac{1}{2}, \quad B=-\frac{1}{2 N} \tag{7.6.30}
\end{equation*}
$$

For $S U(3)$ we have then

$$
\begin{equation*}
\sum_{a} t_{i j}^{a} t_{k l}^{a}=\frac{1}{2}\left(\delta_{i l} \delta_{j k}-\frac{1}{3} \delta_{i j} \delta_{k l}\right) \tag{7.6.31}
\end{equation*}
$$

that coincides with eq.(7.6.11). Let us now turn to the spinor indices. Here we look for identities relating products of two fermion bilinears expressed in different combinations. Schematically $\left(\bar{\psi}_{1} \Gamma \psi_{4}\right)\left(\bar{\psi}_{3} \Gamma \psi_{2}\right) \sim\left(\bar{\psi}_{1} \Gamma \psi_{2}\right)\left(\bar{\psi}_{3} \Gamma \psi_{4}\right)$, where $\Gamma$ are products of gamma matrices. Relations of this form are called Fierz identities. The latter are best derived by decomposing an open index fermion bilinear $\psi_{a} \bar{\psi}_{b}$ in terms of an independent basis in spinor space. A convenient independent basis is provided by the antisymmetric products of gamma matrices:

$$
\begin{equation*}
1_{a b}, \quad \gamma_{a b}^{\mu}, \quad \gamma_{a b}^{\mu \nu}, \quad \gamma_{a b}^{\mu \nu \rho}, \quad \gamma_{a b}^{\mu \nu \rho \sigma} \tag{7.6.32}
\end{equation*}
$$

where $\gamma^{\mu_{1} \ldots \mu_{n}}$ is the completely antisymmetrized product of gamma matrices:

$$
\begin{equation*}
\gamma^{\mu_{1} \ldots \mu_{n}} \equiv \frac{1}{n!}\left(\gamma^{\mu_{1}} \ldots \gamma^{\mu_{n}} \pm \text { perms } .\right) \tag{7.6.33}
\end{equation*}
$$

A spinor basis should consists of $4 \times 4=16$ matrices, with complex valued coefficients. The total number of matrices appearing in eq.(7.6.32) is indeed $1+4+6+4+1=16$. One also notices that the matrices (7.6.32) are orthogonal in spinor space, namely

$$
\begin{equation*}
\operatorname{Tr}\left(\gamma^{\mu_{1} \ldots \mu_{n_{1}}} \gamma^{\mu_{1} \ldots \mu_{n_{2}}}\right)=0, \quad \text { for } \quad n_{1} \neq n_{2} \tag{7.6.34}
\end{equation*}
$$

and hence they form a basis. ${ }^{7}$ Notice that the matrix $\gamma^{\mu \nu \rho \sigma}$ is equivalent to the chiral $\gamma_{5}$ matrix. Hence, an alternative, but simpler, basis is given by the set

$$
\begin{equation*}
1_{a b}, \quad \gamma_{a b}^{\mu}, \quad \gamma_{a b}^{\mu \nu}, \quad\left(\gamma^{\mu} \gamma_{5}\right)_{a b}, \quad \gamma_{a b}^{5} \tag{7.6.35}
\end{equation*}
$$

[^47]A generic fermion bilinear can then be written as follows

$$
\begin{equation*}
\psi_{2 a} \bar{\psi}_{1 b}=c_{1} \delta_{a b}+c_{2} \gamma_{a b}^{\mu}+c_{3} \gamma_{a b}^{\mu \nu}+c_{4}\left(\gamma^{\mu} \gamma_{5}\right)_{a b}+c_{5}\left(\gamma_{5}\right)_{a b} \tag{7.6.36}
\end{equation*}
$$

where $c_{i}(i=1, \ldots 5)$ are coefficients to be determined. Multiplying eq.(7.6.36) by $\delta_{a b}$ immediately allows us to fix $c_{1}$ :

$$
\begin{equation*}
c_{1}=\frac{\eta}{4}\left(\bar{\psi}_{1} \psi_{2}\right) \tag{7.6.37}
\end{equation*}
$$

where $\eta=+1$ for commuting spinors and $\eta=-1$ for anti-commuting spinors. Similarly, multiplying eq.(7.6.36) by the various gamma matrices allows us to fix the remaining coefficients. We get
$\psi_{2 a} \bar{\psi}_{1 b}=\frac{\eta}{4}\left(\left(\bar{\psi}_{1} \psi_{2}\right) \delta_{a b}+\left(\bar{\psi}_{1} \gamma_{\mu} \psi_{2}\right) \gamma_{a b}^{\mu}+\left(\bar{\psi}_{1} \gamma_{\mu \nu} \psi_{2}\right) \gamma_{a b}^{\mu \nu}+\left(\bar{\psi}_{1} \gamma_{\mu} \gamma_{5} \psi_{2}\right)\left(\gamma_{5} \gamma^{\mu}\right)_{a b}+\left(\bar{\psi}_{1} \gamma_{5} \psi_{2}\right)\left(\gamma_{5}\right)_{a b}\right)$.
Using eq.(7.6.38), it is clear how to relate products of bilinears of the form $\left(\bar{\psi}_{1} \Gamma \psi_{4}\right)\left(\bar{\psi}_{3} \Gamma \psi_{2}\right)$ to $\left(\bar{\psi}_{1} \Gamma \psi_{2}\right)\left(\bar{\psi}_{3} \Gamma \psi_{4}\right)$. The relation (7.6.38) is the master equation that, applied to different products of bilinears, gives rise to different sets of Fierz identities. In the situation discussed in section 7.6, eq.(7.6.9), we have to consider chiral fermions. This simplifies the analysis, because upon multiplying eq.(7.6.38) by $P_{L}$ from the left and by $P_{R}$ from the right, $\left(P_{L}=\left(1+\gamma_{5}\right) / 2, P_{R}=\left(1-\gamma_{5}\right) / 2\right)$, three out of five terms in the r.h.s. of that equation vanish. For two (commuting) spinor wave functions we take $\eta=1$, the remaining two terms combine to give

$$
\begin{equation*}
\psi_{2 L, a} \bar{\psi}_{1 L, b}=\frac{1}{2}\left(\bar{\psi}_{1 L} \gamma^{\mu} \psi_{2 L}\right)\left(\gamma_{\mu} P_{R}\right)_{a b} \tag{7.6.39}
\end{equation*}
$$

that is eq.(7.6.12) of section 7.6.

## 7.7 (Ir)relevance of Higher Dimensional Operators

We have so far neglected all higher dimensional operators appearing in the EFT, assuming that their effect is small. This assumption is actually correct, and this is the main reason why we study EFT after all, but it is not as trivial as what naively expected. At the quantum level the insertion of higher dimensional operators in Feynman amplitudes leads to bad divergences that can obscure the irrelevance of these operators at low energies. As a matter of fact, these operators are not always negligible at low-energies, their effect depending on the regularization and renormalization procedure used to deal with the divergences. As an illustrative example, let us go back to the two scalar theory in section 7.2 and let us add the dimension six operator

$$
\begin{equation*}
\frac{\tilde{\lambda}}{M^{2}} L^{3} \square L \tag{7.7.1}
\end{equation*}
$$

to the Lagrangian (7.2.2). ${ }^{8}$ The operator (7.7.1) is indeed generated, at tree-level, when we integrate $H$ out. Naively, we expect that (7.7.1) is suppressed at low-energies as $E^{2} / M^{2}$ with respect to the $\lambda$ coupling in eq. (7.2.2). However, at the quantum level eq. (7.7.1) leads to severe divergences. For instance, at one-loop level the 1PI 4-point function with one $\lambda$ and one $\tilde{\lambda}$ vertex is quadratically divergent, so that

$$
\begin{equation*}
\delta \Gamma^{(4)} \propto \lambda \frac{\tilde{\lambda}}{M^{2}} \int d^{4} k \frac{k^{2}}{\left(k^{2}+p^{2}\right)^{2}} \sim \lambda \frac{\tilde{\lambda}}{M^{2}}\left(\Lambda^{2}+p^{2} \log \frac{\Lambda^{2}}{p^{2}}+\ldots\right) \tag{7.7.2}
\end{equation*}
$$

where $p$ generically denotes external momenta or the mass $m$. Since the cut-off $\Lambda$ of the EFT is around $M$, we see that the divergent term can compensate for the manifest $1 / M^{2}$ suppression to give a contribution to $\Gamma^{(4)}$ of the same order of the leading $\lambda^{2}$ term.

This potential problem is actually scheme-dependent and is manifestly absent in DR with a mass-independent renormalization subtraction scheme, such as $\overline{\mathrm{MS}}$. In DR, divergences arise as poles in $1 / \epsilon$ and, when subtracted, leave a renormalized amplitude where the sliding scale $\mu$ appears in logs only. Since in the EFT the only other mass scales present are the light masses and the low-energy momenta, by dimensional analysis all higher dimensional operators are manifestly suppressed. In the example above, for instance, we get, once the poles are subtracted,

$$
\begin{equation*}
\delta \Gamma^{(4)} \propto \lambda \frac{\tilde{\lambda}}{M^{2}} \int d^{d} k \frac{k^{2}}{\left(k^{2}+p^{2}\right)^{2}} \sim \lambda \frac{\tilde{\lambda}}{M^{2}}\left(p^{2} \log \frac{\mu^{2}}{p^{2}}+\ldots\right), \tag{7.7.3}
\end{equation*}
$$

which is manifestly sub-leading with respect to the $\lambda^{2}$ term.
Of course by using any other regulator, after the divergence is subtracted, no depedence on the cut-off appears. However, in a generic regularization and renormalization prescription, power-like dependence on the sliding scale is induced, which can spoil the naive dimensional analysis. For instance, by renormalizing the operators $L^{4}$ and $L^{3} \square L$ at a scale $\mu$, one would get

$$
\begin{equation*}
\delta \Gamma^{(4)} \sim \lambda \frac{\tilde{\lambda}}{M^{2}}\left(p^{2} \log \frac{\mu^{2}}{p^{2}}+\mu^{2} \log \frac{\mu^{2}}{p^{2}}+\ldots\right) \tag{7.7.4}
\end{equation*}
$$

If the subtraction point $\mu \ll M$, we can still neglect higher dimensional operators, but in a sufficiently complicated set-up powers of $M$ might be induced by mixing with various operators. Moreover, we have seen that it is convenient in effective field theory to renormalize at the scale $\mu \sim M$, in which case dimensional analysis, when using eq.(7.7.4), would break down. These problems are avoided by using DR with a mass independent

[^48]subtraction. In other words, mass-independent schemes preserve dimensional analysis, an observation we already made in subsection 5.8 in the context of the RG flow of general dimensionful couplings. We now see that this property has the fundamental implication of validating a naive treatment of higher dimensional operators in the context of EFT.

### 7.8 Redundant Operators

We have discussed at length that in EFT one should write down a local effective action for the lights fields including all the higher dimensional operators up to the desired order, the coefficients of which will be then fixed by a matching procedure. It turns out, however, that there is a redundancy in this description, since the physical observables in general depend on only a subset of the EFT coupling constants (called physical in what follows). The remaining couplings are called redundant, because they can be expressed in terms of the physical ones. This redundancy can be traced back to the LSZ reduction formulas (2.3.12), for which on-shell amplitudes are associated to the residue of multiple poles in a Green function, not to the whole off-shell Green function. We can obtain different Green functions by changing, for instance, the asymptotic field $\phi$ that creates one-particle states. Any other quantum field $\phi^{\prime}$, as long as $\langle 0| \phi^{\prime}|p\rangle \neq 0$, is an equally good choice. The offshell Green functions associated to $\phi^{\prime}$ would generally differ from those associated to $\phi$, but both would lead to identical on-shell properties. In other words, different off-shell Green functions lead to the same on-shell physics.

The same redundancy can similarly be discussed in terms of operators, and not of their associated couplings, using the Schwinger-Dyson equations (4.4.3). Viewing $\delta S / \delta \phi$ as a particular composite operator in the theory, we can interpret eqs.(4.4.3) as the statement that such composite operator, when inserted in Green functions, gives only rise to contact terms with Green functions containing a lower number of fields. On-shell, this implies that some pole is lost and these operators should not contribute to physical processes (compare with the situation discussed in subsection 4.4.3 in the context of the WI in QED). These operators are called redundant [22]. Composite operators made of $\delta S / \delta \phi$ times other operators evaluated at the same space-time point (suitably renormalized) are also redundant. This can be shown as follows [22]. Let us define

$$
\begin{equation*}
\theta(x)=F(\phi(x)) \frac{\delta S}{\delta \phi(x)} \tag{7.8.1}
\end{equation*}
$$

a (suitably renormalized) composite operator, where $F$ is a local functional of $\phi$ and $\delta S / \delta \phi(x)$ are the equations of motion for the field $\phi$. We can introduce external sources $s$
and $J$ for $\phi$ and $\theta$, respectively, and consider the generating functional

$$
\begin{equation*}
e^{i W[J, s]}=\int \mathcal{D} \phi e^{i \int d^{d} x(\mathcal{L}+s \phi+J \theta)} \tag{7.8.2}
\end{equation*}
$$

We will be interested in correlation functions with at most one insertion of $\theta$, so we can expand in $J$ and keep only the terms up to linear level. Performing the path integral change of variables

$$
\begin{equation*}
\phi(x) \rightarrow \phi(x)-J(x) F(\phi(x)), \tag{7.8.3}
\end{equation*}
$$

we then get

$$
\begin{equation*}
e^{i W[J, s]}=\int \mathcal{D} \phi|\operatorname{det} \operatorname{Jac}(x, y)| e^{i \int d^{d} x(\mathcal{L}+s(\phi-J F))} \tag{7.8.4}
\end{equation*}
$$

where

$$
\begin{equation*}
\operatorname{Jac}(x, y)=\delta^{(d)}(x-y)-J(x) \frac{\delta F((\phi(x))}{\delta \phi(y)} \tag{7.8.5}
\end{equation*}
$$

is the Jacobian associated to the change of variables. By taking $n$ functional derivatives with respect to $s\left(x_{i}\right)$ and one with respect to $J(x)$, and setting $s=J=0$, we get the identity

$$
\begin{align*}
i\left\langle\phi\left(x_{1}\right) \ldots \phi\left(x_{n}\right) \theta(x)\right\rangle= & -\sum_{i=1}^{n} \delta^{(d)}\left(x-x_{i}\right)\left\langle\phi\left(x_{1}\right) \ldots F(\phi(x)) \ldots \phi\left(x_{n}\right)\right\rangle \\
& -\left\langle\phi\left(x_{1}\right) \ldots \phi\left(x_{n}\right) \frac{\delta F((\phi(x))}{\delta \phi(x)}\right\rangle \tag{7.8.6}
\end{align*}
$$

The last term in eq.(7.8.6) arises from the determinant of the Jacobian. For any local functional of fields $F(\phi)$, it is UV divergent and proportional to $\delta^{(d)}(0)$, so it vanishes in DR, where such divergences are set to zero. In this way eq.(7.8.6) is the generalization of the Schwinger-Dyson equations (4.4.3), which are reproduced by taking $F=1$. Equation (7.8.6) shows that any local composite operator $\theta$ of the form (7.8.1) does not contribute to physical processes.

In practice redundant operators can be removed in EFT either by using the equations of motion at the Lagrangian level or alternatively by field redefinitions. It is useful to see how this works in a simple example: an EFT featuring only a real scalar with $\mathbb{Z}_{2}$ symmetry $\phi \rightarrow-\phi$. Up to dimension six operators, the most general EFT reads as:

$$
\begin{equation*}
\mathcal{L}_{E F T}^{(d \leq 6)}=\frac{1}{2}(\partial \phi)^{2}-\lambda \phi^{4}+\frac{g_{1}}{M^{2}} \phi^{6}+\frac{g_{2}}{M^{2}}(\square \phi)^{2}+\frac{g_{3}}{M^{2}} \phi^{3} \square \phi+\mathcal{O}\left(\frac{1}{M^{4}}\right) \tag{7.8.7}
\end{equation*}
$$

where, for further simplification, we have neglected the mass term $m^{2} \phi^{2}$. One could write down other dimension six operators, but these are equal, up to total derivatives, to the operators in eq.(7.8.7). For example, integrating by parts, one easily gets

$$
\begin{equation*}
\phi^{2}(\partial \phi)^{2}=-\frac{1}{3} \phi^{3} \square \phi+\frac{1}{3} \partial_{\mu}\left(\phi^{3} \partial_{\mu} \phi\right), \tag{7.8.8}
\end{equation*}
$$

the last operator giving a vanishing contribution in the action. The equation of motion for $\phi$ is

$$
\begin{equation*}
-\square \phi-4 \lambda \phi^{3}+\mathcal{O}\left(\frac{1}{M^{2}}\right)=0 \tag{7.8.9}
\end{equation*}
$$

Plugging back eq.(7.8.9) in the dimension six terms in the Lagrangian (7.8.7) gives

$$
\begin{equation*}
\widetilde{\mathcal{L}}_{E F T}^{(d \leq 6)}=\frac{1}{2}(\partial \phi)^{2}-\lambda \phi^{4}+\frac{\tilde{g}_{1}}{M^{2}} \phi^{6}+\mathcal{O}\left(\frac{1}{M^{4}}\right), \tag{7.8.10}
\end{equation*}
$$

where

$$
\begin{equation*}
\tilde{g}_{1}=g_{1}+16 \lambda^{2} g_{2}-4 \lambda g_{3} \tag{7.8.11}
\end{equation*}
$$

We see that two out of the three dimension-six operators are redundant. This implies that, at this order, physical observables only depend on the two couplings $\lambda$ and $\tilde{g}_{1}$.

The same conclusion can be reached by making the most general field redefinition compatible with the symmetries at this order:

$$
\begin{equation*}
\phi \rightarrow \phi+\frac{a}{M^{2}} \square \phi+\frac{b}{M^{2}} \phi^{3}, \tag{7.8.12}
\end{equation*}
$$

where $a$ and $b$ are so far undetermined dimensionless parameters. In the new basis the Lagrangian (7.8.7) reads, up to total derivative terms, ${ }^{9}$

$$
\begin{equation*}
\mathcal{L}_{E F T}^{(d \leq 6)} \rightarrow \mathcal{L}_{E F T}^{(d \leq 6)}=\frac{1}{2}(\partial \phi)^{2}-\lambda \phi^{4}+\frac{\tilde{g}_{1}}{M^{2}} \phi^{6}+\frac{\tilde{g}_{2}}{M^{2}}(\square \phi)^{2}+\frac{\tilde{g}_{3}}{M^{2}} \phi^{3} \square \phi+\mathcal{O}\left(\frac{1}{M^{4}}\right), \tag{7.8.13}
\end{equation*}
$$

where

$$
\begin{align*}
& \tilde{g}_{1}=g_{1}-4 b \lambda  \tag{7.8.14}\\
& \tilde{g}_{2}=g_{2}-a  \tag{7.8.15}\\
& \tilde{g}_{3}=g_{3}-4 a \lambda-b . \tag{7.8.16}
\end{align*}
$$

We can choose $a$ and $b$ such that $\tilde{g}_{2}=\tilde{g}_{3}=0$, i.e. $a=g_{2}, b=g_{3}-4 \lambda g_{2}$. Plugging these values back in eq.(7.8.14) reproduces eq.(7.8.11) and gives rise to the same EFT (7.8.10). The presence of two redundant operators in the Lagrangian (7.8.7) can also be seen by taking the following basis for the dimension six operators:

$$
\begin{equation*}
\mathcal{L}_{E F T}^{(d \leq 6)}=\frac{1}{2}(\partial \phi)^{2}-\lambda \phi^{4}+\frac{\hat{g}_{1}}{M^{2}} \phi^{6}+\frac{\hat{g}_{2}}{M^{2}}\left(\square \phi+4 \lambda \phi^{3}\right)^{2}+\frac{\hat{g}_{3}}{M^{2}} \phi^{3}\left(\square \phi+4 \lambda \phi^{3}\right)+\mathcal{O}\left(\frac{1}{M^{4}}\right), \tag{7.8.17}
\end{equation*}
$$

where we immediately see that $\hat{g}_{2}$ and $\hat{g}_{3}$ are redundant couplings, and the operators $\hat{\mathcal{O}}_{2}$ and $\hat{\mathcal{O}}_{3}$ in eq.(7.8.17) are of the form (7.8.1), with $F_{2}=\square \phi+4 \lambda \phi^{3}, F_{3}=\phi^{3}$.

[^49]Removing redundant operators can lead to a significant simplification. It should however be noted that if one starts with a Lagrangian where redundant operators have been removed, like eq.(7.8.10), and considers off-shell Green functions, the cancellation of UV divergences will generally require the introduction of counterterms associated to the redundant operators. Such counterterms can then be rexpressed in terms of counterterms for the non-redundant operators using field redefinitions or the equations of motion.

## Chapter 8

## Spontaneously Broken Symmetries

Symmetries are a fundamental concept in QFT but they are often broken in Nature. If a physical system described by a Lagrangian $\mathcal{L}(\phi)$ is invariant under some symmetry, a possible breaking term is obtained by adding terms to $\mathcal{L}(\phi)$ which do not respect the symmetry (explicit breaking). If the corresponding breaking operators have dimensions less than four, the UV behaviour of the theory (and its divergences) will not be affected by the symmetry breaking terms. In this case we say that the symmetry is softly broken.

Another possibility occurs when $\mathcal{L}(\phi)$ is invariant under the symmetry but the ground state (the vacuum) is not. If $G$ is the operator in the Hilbert space parametrizing the action of the symmetry group, $G|0\rangle \neq|0\rangle$. In this case we have what is called a spontaneous symmetry breaking. Classical prototypical example is the breaking of the $S O(3)$ spatial rotations in a ferromagnet. The laws of Nature are all spatially symmetric, but the vacuum is not. The simplest example of a spontaneous symmetry breaking mechanism is provided by the $\lambda \phi^{4}$ theory with a negative squared mass term, that enjoys a $\mathbf{Z}_{2}$ symmetry under which $\phi \rightarrow-\phi$, with $\langle\phi\rangle= \pm \phi_{0} \neq 0$. Clearly, the $\mathbf{Z}_{2}$ symmetry exchanges the two degenerate vacua, $\mathbf{Z}_{2}\left| \pm \phi_{0}\right\rangle=\left|\mp \phi_{0}\right\rangle$. The two vacua are physically identical, so that we can choose any of the two and expand for small fluctuations around the selected one.

The story is however not so simple. We are tacitly assuming that the vacuum is one of the two vacua $\left| \pm \phi_{0}\right\rangle$, and not a linear combination of them. If the vacuum was
 broken (the vacua being mapped to themselves, up to a possible sign factor). This is in fact what happens in the quantum mechanical analogue of the double well potential: the hamiltonian eigenstates are given by $| \pm\rangle$, since tunneling effects induce non-trivial transitions of the form $\left\langle \pm \phi_{0} \mid \mp \phi_{0}\right\rangle$, and no spontaneous $\mathbf{Z}_{2}$ symmetry breaking occurs. So we start in the next section explaining why and how spontaneous symmetry breaking
takes place in QFT.

### 8.1 Why Spontaneous Symmetry Breaking?

As we mentioned, spontaneous symmetry breaking does not occur in quantum mechanics. ${ }^{1}$ More generally, it does not occur in systems with a finite number of degrees of freedom, like in QFT defined on a finite volume. The tunneling effects between different vacua in QFT are exponentially suppressed by the volume of space $V$. As $V \rightarrow \infty$, such transitions no longer occur and different degenerate vacua are equivalent and completely disconnected from each other: spontaneous symmetry breaking can then occur.

The absence of transitions between different degenerate vacua in QFT can be shown by looking at the equal-time commutators of two generic local operators $A$ and $B$. Assume that the vacua are discrete and take them orthonormal:

$$
\begin{equation*}
\langle u \mid v\rangle=\delta_{u, v} . \tag{8.1.1}
\end{equation*}
$$

Inserting a complete basis of states we have

$$
\begin{equation*}
\langle u| A(\vec{x}) B(0)|v\rangle=\sum_{w}\langle u| A(0)|w\rangle\langle w| B(0)|v\rangle+\int d^{3} \vec{p} e^{i \vec{p} \cdot \vec{x}} \rho(\vec{p}) \tag{8.1.2}
\end{equation*}
$$

where the first term is the sum over all the possible vacua $w$ present in the theory and

$$
\begin{equation*}
\rho(\vec{p}) \equiv \sum_{n}\langle u| A(0)\left|n, q_{n}\right\rangle\left\langle n, q_{n}\right| B(0)|v\rangle \delta^{(3)}\left(\vec{p}-\vec{q}_{n}\right) \tag{8.1.3}
\end{equation*}
$$

represents the sum over $n$-particles states with total momentum $\vec{q}_{n}$. If we assume that $\rho(\vec{p})$ is integrable, the Riemann-Lebesgue lemma applies and we have

$$
\begin{equation*}
\lim _{|\vec{x}| \rightarrow \infty} \int d^{3} \vec{p} e^{i \vec{p} \cdot \vec{x}} \rho(\vec{p})=0 \tag{8.1.4}
\end{equation*}
$$

This condition is automatically satisfied if the theory has a mass gap, namely if the one-particle states are all massive, in which case the Fourier transform of $\rho(\vec{p})$ vanishes exponentially in $|\vec{x}|$ at large distances. It ensures that operators inserted far away from each other have no correlation. This is cluster decomposition at the operator level, that will play an important role again in what follows. Since $A$ and $B$ are space-like separated, microcausality requires that

$$
\begin{equation*}
[A(\vec{x}), B(0)]=0 \tag{8.1.5}
\end{equation*}
$$

[^50]Eqs.(8.1.2), (8.1.4) and (8.1.5) imply that the matrix elements $\langle u| A(0)|v\rangle$ and $\langle u| B(0)|v\rangle$ should commute for any $u$ and $v$, and hence can be simultaneously diagonalizable. There exists then a basis of vacua where

$$
\begin{equation*}
\langle u| A(0)|v\rangle=a_{u} \delta_{u, v} \tag{8.1.6}
\end{equation*}
$$

Since $A$ is an arbitrary local operator, we can take it to be the hamiltonian density operator and thus we have proved that no transitions between degenerate vacua are possible in QFT. Notice how the infinite volume limit entered crucially in eq.(8.1.4).

However, this is not the end of the story, because we have still to rule out the possibility that the final vacuum is given by some linear combination $\sum_{u} c_{u}|u\rangle$, where $|u\rangle$ is the orthonormal basis defined by eq.(8.1.6), rather than by any of the $|u\rangle$. This possibility is ruled out by demanding cluster decomposition of the correlation function:

$$
\begin{equation*}
\lim _{|\vec{x}| \rightarrow \infty}\langle A(\vec{x}) B(0)\rangle_{c}=\lim _{|\vec{x}| \rightarrow \infty}(\langle A(\vec{x}) B(0)\rangle-\langle A(\vec{x})\rangle\langle B(0)\rangle)=0 \tag{8.1.7}
\end{equation*}
$$

For instance, for two vacua related by a $\mathbf{Z}_{2}$ symmetry, we can denote $| \pm\rangle$ the vacua in the diagonal basis (8.1.6) and by $|\alpha\rangle$ the linear combination

$$
\begin{equation*}
|\alpha\rangle=\cos \alpha|+\rangle+\sin \alpha|-\rangle \tag{8.1.8}
\end{equation*}
$$

with $0 \leq \alpha \leq \pi / 2$. In a $\alpha$-vacuum we have

$$
\begin{align*}
\lim _{|\vec{x}| \rightarrow \infty}\langle\phi(\vec{x}) \phi(0)\rangle_{\alpha} & =\langle\alpha| \phi(0)|+\rangle\langle+| \phi(0)|\alpha\rangle+\langle\alpha| \phi(0)|-\rangle\langle-| \phi(0)|\alpha\rangle  \tag{8.1.9}\\
& =v^{2}\left(\sin ^{2} \alpha+\cos ^{2} \alpha\right)=v^{2}
\end{align*}
$$

where $\pm v$ is the VEV of the field $\phi$ in the two vacua $| \pm\rangle$. On the other hand

$$
\begin{equation*}
\langle\phi(0)\rangle_{\alpha}=\langle\alpha| \phi(0)|\alpha\rangle=v\left(\cos ^{2} \alpha-\sin ^{2} \alpha\right)=v \cos 2 \alpha \tag{8.1.10}
\end{equation*}
$$

so that

$$
\begin{equation*}
\lim _{|\vec{x}| \rightarrow \infty}\langle\alpha| \phi(\vec{x}) \phi(0)|\alpha\rangle_{c}=v^{2}\left(1-\cos ^{2} 2 \alpha\right)=v^{2} \sin ^{2} 2 \alpha . \tag{8.1.11}
\end{equation*}
$$

The only $\alpha$-vacua that satisfy cluster decomposition are those with $\alpha=0, \pi / 2$, i.e. the states $|+\rangle$ and $|-\rangle$. Spontaneous symmetry breaking can then be seen as an obstruction in having a vacuum that at the same time is invariant under a symmetry and satisfies cluster decomposition.

The discussion above can be generalized to continuous symmetries where we have an infinite family of degenerate vacua. For continuous symmetries we also have an important theorem, the Goldstone theorem, relating the symmetry breaking pattern to the spectrum of the theory.

### 8.2 The Goldstone Theorem

Let $G$ be a group of global continuous symmetries with generators $t^{\alpha}, \alpha=1, \ldots, \operatorname{dim} G$, acting on some system. By Noether's theorem, we have $\operatorname{dim} G$ conserved currents $J_{\mu}^{\alpha}(x)$ : $\partial^{\mu} J_{\mu}^{\alpha}(x)=0$ and the associated charges $Q^{\alpha}=\int d^{3} x J_{0}^{\alpha}(x)$. The group $G$ is said to be spontaneously broken if, on the vacuum, the generators splits into two sets, labelled by $a$ and $i$, such that

$$
\begin{equation*}
Q^{a}|0\rangle \neq 0, \quad Q^{i}|0\rangle=0, \tag{8.2.1}
\end{equation*}
$$

with a non-empty set for $a$. The unbroken generators labelled by $t^{i}$ form a subgroup of $G$, denoted by $H$, so that we have $\operatorname{dim} G-\operatorname{dim} H$ broken generators $(a=1, \ldots, \operatorname{dim} G-$ $\operatorname{dim} H)$. Goldstone's theorem states that, independently of the specific pattern of breaking and physical system we are considering, in the spectrum there will appear one massless and spinless particle for each broken generator. The particle will be scalar or pseudoscalar, depending on the parity of the associated broken generator. These particles are called Goldstone or Nambu-Goldstone (NG) bosons. For simplicity, we will always refer to them in what follows as to the NG particles.

Proof of the theorem. Let $\phi_{n}$ be the set of fields responsible for the spontaneous symmetry breaking pattern $G \rightarrow H$. This implies that an infinitesimal action of the group $G$ on the fields $\phi_{n}$ do not leave them invariant, namely $\left\langle\phi_{n}^{\prime}(0)\right\rangle \neq\left\langle\phi_{n}(0)\right\rangle$, where $\phi^{\prime}$ are the fields one gets after the infinitesimal action. In other words, we must have

$$
\begin{equation*}
\left\langle\delta \phi_{n}(0)\right\rangle=\epsilon^{\alpha}\left\langle\left[Q^{\alpha}, \phi_{n}(0)\right]\right\rangle \neq 0, \tag{8.2.2}
\end{equation*}
$$

with $\left\langle\left[Q^{a}, \phi_{n}(0)\right]\right\rangle \equiv \delta \phi_{n}^{a} \neq 0$, for each $a$. Consider now the two-point function of a conserved current $J_{\mu}^{a}$ with the fields $\phi_{n}$. Since $\partial^{\mu} J_{\mu}^{a}(x)=0$, we have

$$
\begin{align*}
\partial_{x}^{\mu}\langle 0| T\left[J_{\mu}^{a}(x) \phi_{n}(0)\right]|0\rangle & =\partial_{x}^{\mu}\left(\theta\left(x^{0}\right)\left\langle J_{\mu}^{a}(x) \phi_{n}(0)\right\rangle+\theta\left(-x^{0}\right)\left\langle\phi_{n}(0) J_{\mu}^{a}(x)\right\rangle\right)  \tag{8.2.3}\\
& =\delta\left(x^{0}\right)\left\langle\left[J_{0}^{a}(x), \phi_{n}(0)\right]\right\rangle
\end{align*}
$$

Let us denote by $G_{\mu, n}^{a}$ the Fourier transform of the two-point function:

$$
\begin{equation*}
\langle 0| T\left[J_{\mu}^{a}(x) \phi_{n}(0)\right]|0\rangle=\int \frac{d^{4} p}{(2 \pi)^{4}} G_{\mu, n}^{a}(p) e^{-i p \cdot x} \tag{8.2.4}
\end{equation*}
$$

By Lorentz invariance, we have

$$
\begin{equation*}
G_{\mu, n}^{a}(p)=i p_{\mu} H_{n}^{a}\left(p^{2}\right) . \tag{8.2.5}
\end{equation*}
$$

Integrating eq. (8.2.3) over space-time gives

$$
\begin{equation*}
\int d^{4} x \partial_{x}^{\mu}\langle 0| T\left[J_{\mu}^{a}(x) \phi_{n}(0)\right]|0\rangle=\int d^{4} p \delta(p) p^{2} H_{n}^{a}\left(p^{2}\right)=\left\langle\left[Q^{a}, \phi_{n}(0)\right]\right\rangle \neq 0 \tag{8.2.6}
\end{equation*}
$$

The integral of a divergence does not vanish when long-range particles appear, and these are precisely the NG bosons. More precisely, we must have

$$
\begin{equation*}
H_{n}^{a}\left(p^{2}\right)=\frac{\delta \phi_{n}^{a}}{p^{2}}+\ldots \tag{8.2.7}
\end{equation*}
$$

where $\ldots$ represent terms that are regular when $p \rightarrow 0$. A pole in the two-point function of $J_{\mu}^{a}$ with $\phi_{n}$ is a signal of the presence of massless 1-particle states, one for each unbroken generator $a$. Since $\phi_{n}$ are spinless (otherwise in the vacuum we will break the Lorentz symmetries as well), such particles are necessarily of spin zero. Their intrinsic parity is then fixed by the parity of the associated current $J^{a}$. Q.e.d.

Denoting by $N G_{a}$ the one-NG particle states, we have

$$
\begin{align*}
\langle 0| J_{\mu}^{a}(x)\left|N G_{b}(p)\right\rangle & =\frac{i p_{\mu} F_{a b} e^{-i p \cdot x}}{\sqrt{2 p_{0}(2 \pi)^{3}}} \\
\left\langle N G_{b}(p)\right| \phi_{n}(x)|0\rangle & =\frac{Z_{n}^{b} e^{i p \cdot x}}{\sqrt{2 p_{0}(2 \pi)^{3}}} \tag{8.2.8}
\end{align*}
$$

where $F_{a b}$ and $Z_{n}^{a}$ are matrices (of mass dimension one and zero, respectively) related to $\delta \phi_{n}^{a}$ in eq.(8.2.7). This relation is found by noticing that $G_{\mu, n}^{a}$ above is associated to the spectral density $\rho_{\mu, n}^{a}$ of the two point function $\langle 0| J_{\mu}^{a}(x) \phi_{n}(0)|0\rangle$ (recall eq.(2.1.16)):

$$
\begin{equation*}
G_{\mu, n}^{a}(p)=\int_{0}^{\infty} d \sigma \rho_{\mu, n}^{a}(p, \sigma) \frac{i}{p^{2}-\sigma+i \epsilon} \tag{8.2.9}
\end{equation*}
$$

where by Lorentz invariance $\rho_{\mu, n}^{a}(p, \sigma)=p_{\mu} \rho_{n}^{a}(\sigma)$. The contribution of the one-particle NG bosons to $\rho_{n}^{a}$ is easily computed from eq.(8.2.8):

$$
\begin{equation*}
\rho_{n, N G}^{a}(\sigma)=i F_{a b} Z_{n}^{b} \delta(\sigma) \tag{8.2.10}
\end{equation*}
$$

Matching eqs.(8.2.7), (8.2.9) and (8.2.10) we then have

$$
\begin{equation*}
\delta \phi_{n}^{a}=i F_{a b} Z_{b}^{n} \tag{8.2.11}
\end{equation*}
$$

We can also define the NG fields $\pi^{a}(x)$ with canonical 1-point function with the NG particles:

$$
\begin{equation*}
\left\langle N G_{b}(p)\right| \pi^{a}(x)|0\rangle=\frac{\delta_{a b} e^{i p \cdot x}}{\sqrt{2 p_{0}(2 \pi)^{3}}} \tag{8.2.12}
\end{equation*}
$$

Putting together eqs. (8.2.8) and eq. (8.2.12), we see that

$$
\begin{equation*}
\phi_{n}(x)=Z_{n}^{a} \pi^{a}(x)+\ldots=(i F)_{a b}^{-1} \delta \phi_{n}^{b} \pi^{a}(x)+\ldots \tag{8.2.13}
\end{equation*}
$$

where ... denote field components not associated to the NG bosons. For linearly realized symmetries, in a basis where all the field components $\phi_{n}$ are real and all the generators $t^{\alpha}$ are purely imaginary, we simply have

$$
\begin{equation*}
\delta \phi_{n}^{a}=i t_{n m}^{a}\left\langle\phi_{m}\right\rangle . \tag{8.2.14}
\end{equation*}
$$

The appearance of one NG particle per broken generator is physically explained by noticing that the potential for the $\phi_{n}$ is, by assumption, invariant under the symmetry. Hence, around any vacuum $\left\langle\phi_{m}\right\rangle$, the broken generators do not leave the minimum invariant, but necessarily shift it in a new minimum with exactly the same energy. Hence, we have a ( $\operatorname{dim} G$ - $\operatorname{dim} H)$-dimensional space of flat directions in the potential. It costs no energy to fluctuate around a flat direction, and the small fluctuation is identified with the NG boson.

The above proof of the Goldstone's theorem does not rely on perturbation theory, indicating that NG particles are expected to arise whenever a global symmetry group is spontaneously broken, no matter whether the theory undergoing the symmetry breaking is weakly or strongly coupled. The fields $\phi_{n}$ responsible for the symmetry breaking are also not necessarily elementary fields appearing directly in the Lagrangian at some energy scale, but might be composite fields built with different fields. The most relevant example of this kind is the spontaneous breaking of the $S U(2)$ chiral symmetry in QCD, induced by effective scalar fields $\phi_{n}$ constructed out of quark bilinears. In this case, the three NG bosons are spin zero mesons that appear as bound states of the original quarks, the pions $\pi^{0}, \pi^{ \pm}$. We will come back to this relevant case in much more detail in the following.

We close this section by noticing that the Goldstone's theorem applies for internal symmetries only, namely for those symmetries whose generators commute with the ones of the Poincaré group. An example of non-internal symmetry is provided by the conformal group, that in four dimensions is given by $S O(4,2)$. A CFT is invariant under the conformal group. When the latter is broken spontaneously to the Poincaré group, one gets $15-10=5$ broken generators (special conformal transformations and dilatations), but actually only one massless NG boson, commonly denoted as the dilaton.

As we will shortly see, the Goldstone's theorem does not apply for local (i.e. space-time dependent) symmetries as well.

### 8.3 Vacuum Alignement and Pseudo-Goldstone Bosons

It is interesting to analyze theories where, in some approximation, a spontaneous symmetry breaking occurs and, in addition, sub-leading effects introduce explicit symmetry breaking
terms. ${ }^{2}$ Relevant examples include the above mentioned case of QCD, where the up and down quark masses can be seen as small explicit symmetry breaking terms of the $S U(2)$ chiral symmetry. We will consider in what follows linearly realized symmetries only, for which we know that quantum effects do not spoil the invariance of the action. By considering space-time independent field configurations $\phi_{n}$, the whole effective action boils down to the effective potential $V(\phi)$. The latter, by definition, can be written as

$$
\begin{equation*}
V(\phi)=V_{0}(\phi)+V_{1}(\phi) \tag{8.3.1}
\end{equation*}
$$

where $V_{0}$ is the invariant term, and $V_{1}$ is the breaking one. In the field region of interest, by assumption, we have $\left|V_{1}\right| \ll\left|V_{0}\right|$. In the basis (8.2.14), we have

$$
\begin{equation*}
\delta \phi_{n}=i \epsilon^{\alpha} t_{n m}^{\alpha} \phi_{m} \tag{8.3.2}
\end{equation*}
$$

Invariance of $V_{0}$ implies

$$
\begin{equation*}
\frac{\partial V_{0}}{\partial \phi_{n}} t_{n m}^{\alpha} \phi_{m}=0, \quad \forall \alpha \tag{8.3.3}
\end{equation*}
$$

Let $\phi_{0}$ be the minimum of $V_{0}$ and $\phi=\phi_{0}+\phi_{1}$ the minimum of the whole potential $V$ :

$$
\begin{equation*}
\left.\frac{\partial V}{\partial \phi_{n}}\right|_{\phi=\phi_{0}+\phi_{1}}=0 \tag{8.3.4}
\end{equation*}
$$

Since $\left|V_{1}\right| \ll\left|V_{0}\right|$, we also have $\phi_{1} \ll \phi_{0}$ and we can consistently expand eq. (8.3.4) for small $V_{1}$ and $\phi_{1}$. At first non-trivial order, we have

$$
\begin{equation*}
\left.\frac{\partial^{2} V_{0}}{\partial \phi_{n} \partial \phi_{m}}\right|_{\phi_{0}} \phi_{1, m}+\left.\frac{\partial V_{1}}{\partial \phi_{n}}\right|_{\phi_{0}}=0 \tag{8.3.5}
\end{equation*}
$$

Taking a derivative with respect to $\phi$ of eq. (8.3.3), and evaluating at $\phi_{0}$, we have

$$
\begin{equation*}
\left.\frac{\partial^{2} V_{0}}{\partial \phi_{n} \partial \phi_{m}}\right|_{\phi_{0}} t_{m p}^{\alpha} \phi_{0, p}=0 \tag{8.3.6}
\end{equation*}
$$

When the index $\alpha$ runs over the unbroken directions $i$, (8.3.6) trivially vanishes since $t_{m p}^{i} \phi_{0, p}=0$. For $\alpha=a,(8.3 .6)$ is non-trivial and indicates that there are $\operatorname{dim} G$-dim $H$ directions in field space with a massless eigenvector, that are the NG particles (this can be in fact seen as an alternative proof of the Goldstone's theorem). Multiplying eq. (8.3.5) by $t_{n p}^{a} \phi_{0, p}$ and using eq. (8.3.6), we finally have

$$
\begin{equation*}
\left.\frac{\partial V_{1}}{\partial \phi_{n}}\right|_{\phi_{0}} t_{n p}^{\alpha} \phi_{0, p}=0 \tag{8.3.7}
\end{equation*}
$$

[^51]In presence of a perturbation $V_{1}$, the vacuum $\phi_{0}$ is no longer arbitrary among the would-be degenerate vacua (in absence of the perturbation) but is restricted to satisfy eq. (8.3.7). Equation (8.3.7) is denoted vacuum alignment condition, because it shows that the symmetry breaking terms typically force the vacuum $\phi_{0}$ to be parallel to the direction of the explicit breaking term. This is easily seen in the particularly simple example of an $S O(N) \rightarrow S O(N-1)$ breaking pattern. We choose as explicit breaking term $V_{1}=u_{n} \phi_{n}$, where $u_{n}$ is a fixed vector, explicitly breaking $S O(N)$ down to the $S O(N-1)$ subgroup that leaves it fixed. The condition (8.3.7) gives $u_{n} t_{n p}^{a} \phi_{0, p}=0$, namely $\phi_{0, n} \propto u_{n}$ (the matrices $t^{a}$ being antisymmetric). The addition of the small perturbation to the invariant Lagrangian lifts the vacuum degeneracy and forces $\phi_{0}$ to be parallel to $u_{n}$. The final unbroken group is then $S O(N-1)$ and not $S O(N-2)$, intersection of the two would-be misaligned $S O(N-1)$ subgroups.

The explicit term $V_{1}$ is generally responsible for another important effect: they give masses to the NG bosons, which become what are sometimes denoted as pseudo-NG bosons. Their masses can be extracted from the potential by using eq. (8.2.13):

$$
\begin{equation*}
M_{a b}^{2}=\left.\frac{\partial^{2} V}{\partial \pi^{a} \partial \pi^{b}}\right|_{\phi}=\left.\left(i F_{a c}\right)^{-1} \delta \phi_{n}^{c}\left(i F_{b d}\right)^{-1} \delta \phi_{m}^{d} \frac{\partial^{2} V}{\partial \phi_{n} \partial \phi_{m}}\right|_{\phi} . \tag{8.3.8}
\end{equation*}
$$

Expanding for small $V_{1}$ and $\phi_{1}$, it is straightforward to see that the pseudo NG boson masses are linearly proportional to the explicit breaking term:

$$
\begin{equation*}
M_{a b}^{2} \propto V_{1} \tag{8.3.9}
\end{equation*}
$$

### 8.4 Spontaneously Broken Gauge Symmetries: the Higgs Mechanism

The Goldstone's theorem does not apply in the case in which the broken symmetry is local. ${ }^{3}$ It is impossible for gauge theories to keep at the same time Lorentz invariance and positivity of the Hilbert space, both conditions being important to establish the theorem. For local symmetries, no NG massless particles appear, but rather the would-be NG bosons are "eaten" by gauge fields that become massive.

The situation is easily illustrated by the abelian Higgs model (model that will be extensively analyzed in the final chapter in the special case $m=0$ ), whose Lagrangian is

$$
\begin{equation*}
\mathcal{L}=-\frac{1}{4} F_{\mu \nu} F^{\mu \nu}+\left|D_{\mu} \Phi\right|^{2}-V(\Phi), \quad V(\Phi)=-m^{2} \Phi^{\dagger} \Phi+\frac{\lambda}{2}\left(\Phi^{\dagger} \Phi\right)^{2} \tag{8.4.1}
\end{equation*}
$$

[^52]with $D_{\mu} \Phi=\partial_{\mu} \Phi-i e A_{\mu} \Phi$. For $m^{2}>0$, the minimum of the potential is at
\[

$$
\begin{equation*}
\left|\Phi_{0}\right|=\sqrt{m^{2} / \lambda} \equiv \frac{v}{\sqrt{2}} . \tag{8.4.2}
\end{equation*}
$$

\]

It is useful to choose radial coordinates for the field $\Phi$ and write

$$
\begin{equation*}
\Phi=\frac{v+\rho}{\sqrt{2}} e^{i \theta / v} \tag{8.4.3}
\end{equation*}
$$

where the factor $v$ in the exponential has been introduced to get canonical fields of dimension one. Expanding around small fluctuations, we quickly get at quadratic order

$$
\begin{equation*}
\mathcal{L}=-\frac{1}{4} F_{\mu \nu} F^{\mu \nu}+\frac{1}{2}\left(\partial_{\mu} \rho\right)^{2}+\frac{1}{2}\left(\partial_{\mu} \theta-e v A_{\mu}\right)^{2}-\frac{1}{2}\left(2 m^{2}\right) \rho^{2}+\ldots \tag{8.4.4}
\end{equation*}
$$

The radial field is massive, with $m_{\rho}=\sqrt{2} m$, while the angular field $\theta$ is massless. Actually $\theta$ mixes at quadratic level with the gauge field $A_{\mu}$ and it is not an eigenstate of the free Hamiltonian. We can easily get rid of the unwanted mixing term $e A^{\mu} \partial_{\mu} \theta$ by noticing that we have still to gauge-fix the $U(1)$ gauge symmetry. Under a $U(1)$ transformation with parameter $\alpha(x), \Phi(x) \rightarrow \exp (i \alpha(x)) \Phi(x)$. The radial field $\rho(x)$ is invariant, while the angular field $\theta(x)$ shifts as $\theta(x) \rightarrow \theta(x)+v \alpha(x)$. For any $v \neq 0$, we can take $\alpha=-\theta / v$ and set in this way $\theta(x)=0$. In this gauge, denoted unitary gauge, the third term in eq. (8.4.4) boils down to a mass term for the $U(1)$ gauge field, $m_{A}=e v$. We can say that the gauge field has "eaten" the field $\theta$ that becomes the longitudinal component of a massive gauge field. This mechanism is commonly denoted Higgs mechanism. The number of physical degrees of freedom (d.o.f.) in the process is unchanged. We started with 2 d.o.f. from the massless gauge field and 2 d.o.f. from the complex scalar field, for a total of 4 , and ended up in 1 d.o.f. from the neutral field $\rho$ and 3 d.o.f. from the massive gauge field, again for a total of 4 .

Notice that in the limit $e \rightarrow 0$ the $U(1)$ symmetry becomes global and the gauge field decouples. In this limit the Goldstone's theorem applies and we recover the massless NG boson $\theta$.

All the above analysis similarly applies to non-abelian symmetries. Let us consider the usual sets of real fields $\phi_{n}$ transforming as (8.3.2) under an infinitesimal local transformation of a group $G$. For simplicity, we assume the fields $\phi_{n}$ to be in some irreducible representation of $G$, although this is not strictly necessary. The Lagrangian of our system will include the terms

$$
\begin{equation*}
\mathcal{L} \supset \frac{1}{2}\left(\partial_{\mu} \phi_{n}-i g t_{n m}^{\alpha} A_{\mu}^{\alpha} \phi_{m}\right)^{2}-V(\phi), \tag{8.4.5}
\end{equation*}
$$

where we assume that the potential $V(\phi)$ has minima for $\phi_{m}=v_{m}$ such that, in the "ungauged" limit $g \rightarrow 0$, the global group $G$ is spontaneously broken to $H$. Expanding
around the non-trivial vacuum, $\phi_{m}=v_{m}+\phi_{m}^{\prime}$, the above covariant derivative gives rise to the following terms, up to quadratic order in the fluctuations,

$$
\begin{equation*}
\mathcal{L} \subset \frac{1}{2}\left(\partial_{\mu} \phi_{n}^{\prime}\right)^{2}-i g \partial_{\mu} \phi_{n}^{\prime} t_{n m}^{\alpha} v_{m} A_{\mu}^{\alpha}+\frac{g^{2}}{2}\left(i t_{n m}^{\alpha} v_{m}\right)\left(i t_{n p}^{\beta} v_{p}\right) A_{\mu}^{\alpha} A^{\mu, \beta} \tag{8.4.6}
\end{equation*}
$$

When $\alpha, \beta=i$, the second and third terms in eq. (8.4.6) vanish identically, since $t_{n m}^{i} v_{m}=$ 0 , while they are generically non-vanishing for $\alpha, \beta=a$. The second term in eq. (8.4.6) is the generalization of the $\partial_{\mu} \theta A^{\mu}$ mixing term in the $U(1)$ case. Recall that the directions in field space given by $t_{n m}^{a} v_{m}$ correspond to the NG boson directions. As we will show below, it is always possible to choose a gauge (denoted again unitary gauge) in which the fields $\phi_{n}$ do not contain NG boson fields, and hence

$$
\begin{equation*}
\phi_{n}^{\prime} t_{n m}^{\alpha} v_{m}=0 \tag{8.4.7}
\end{equation*}
$$

The third term in eq. (8.4.6) is a mass term for the gauge fields in the broken directions. The mass matrix

$$
\begin{equation*}
\mu_{\alpha \beta}^{2}=g^{2}\left(i t_{n m}^{\alpha} v_{m}\right)\left(i t_{n l}^{\beta} v_{l}\right) \tag{8.4.8}
\end{equation*}
$$

is symmetric, real and positive, showing that each gauge field $A_{\mu}^{a}$ along the broken directions gets a non-vanishing mass, by eating the corresponding would-be NG boson. In the basis in which $\mu^{2}$ is diagonal, the propagators $G_{\mu \nu}^{a}$ for the massive gauge fields read

$$
\begin{equation*}
G_{\mu \nu}^{a(U G)}(p)=\frac{-i}{p^{2}-\mu_{a}^{2}}\left(\eta_{\mu \nu}-\frac{p_{\mu} p_{\nu}}{\mu_{a}^{2}}\right), \tag{8.4.9}
\end{equation*}
$$

where $U G$ stands for unitary gauge and $\mu_{a}^{2}$ are the mass square eigenvalues. Notice that for large values of the momentum $p$, the gauge field propagator goes to a constant. The renormalizability properties of spontaneously broken gauge theories are correspondingly unclear. In order to fix this problem, it is sometimes more useful to use a more general class of gauge-fixing terms, denoted by $\xi$-gauges. The gauge fixing term $\mathcal{L}_{\text {g.f. }}$ for $\xi$-gauges is a generalization of the usual $1 /(2 \xi)\left(\partial_{\mu} A^{\mu}\right)^{2}$. It reads

$$
\begin{equation*}
\mathcal{L}_{\text {g.f. }}=-\frac{1}{2 \xi} f_{\alpha} f_{\alpha}, \quad f_{\alpha}=\partial_{\mu} A_{\alpha}^{\mu}+i \xi g \phi_{n}^{\prime} t_{n m}^{\alpha} v_{m} \tag{8.4.10}
\end{equation*}
$$

The associated ghost Lagrangian is obtained by taking an infinitesimal gauge variation of $f_{\alpha}$, with parameter $\omega^{\alpha}, \mathcal{L}_{g h .}=-g \omega_{\alpha}^{\star} \delta_{\omega} f_{\alpha}$. We get, modulo total derivatives, ${ }^{4}$

$$
\begin{equation*}
\mathcal{L}_{g h .}=\left(\partial^{\mu} \omega_{\alpha}^{\star}\right)\left(D_{\mu} \omega_{\alpha}\right)-\xi g^{2}\left(i t_{n m}^{\alpha} v_{m}\right)\left(i t_{n l}^{\beta} \phi_{l}\right) \omega_{\alpha}^{\star} \omega_{\beta} \tag{8.4.11}
\end{equation*}
$$

From eq. (8.4.11) we see that the ghost fields along the broken directions have a mass square that equals the gauge boson mass matrix times $\xi$. For any $\xi$, the term $i g \xi\left(\phi^{\prime} t v\right)_{\alpha} \partial_{\mu} A_{\alpha}^{\mu}$

[^53]appearing in eq. (8.4.10), upon an integration by parts, cancels the second unwanted term in eq. (8.4.6). The unitary gauge (8.4.7) corresponds to the limit $\xi \rightarrow \infty$ of the above class of gauges. In this limit, the term proportional to $\xi$ in $\mathcal{L}_{\text {g.f. }}$ oscillates very rapidly and averages to zero any field configuration for which this term is non-zero. We then effectively recover the unitary gauge (8.4.7). We see from eq. (8.4.11) that in the unitary gauge the ghost fields become infinitely massive and can be neglected.

Let us derive the propagators for gauge fields, ghosts and scalar (NG and not) fields in an arbitrary $\xi$-gauge. In momentum space, setting $\mu^{2}$ to diagonal form, the quadratic terms of the gauge bosons in the Lagrangian add to

$$
\begin{equation*}
\mathcal{L}_{\text {quad. }}(A)=\frac{1}{2} A_{\mu}^{\alpha}(-p) A_{\nu}^{\beta}(p) \delta_{\alpha \beta}\left(-p^{2} \eta^{\mu \nu}+p^{\mu} p^{\nu}\left(1-\frac{1}{\xi}\right)+\eta^{\mu \nu} \mu_{\alpha}^{2}\right) \tag{8.4.12}
\end{equation*}
$$

valid for both broken and unbroken generators (in the latter case $\mu_{i}^{2}=0$ ). By inverting the above quadratic terms, we easily get

$$
\begin{equation*}
G_{\mu \nu}^{\alpha \beta(\xi)}(p)=\frac{-i \delta_{\alpha \beta}}{p^{2}-\mu_{a}^{2}}\left(\eta_{\mu \nu}-\frac{(1-\xi) p_{\mu} p_{\nu}}{p^{2}-\xi \mu_{\alpha}^{2}}\right) . \tag{8.4.13}
\end{equation*}
$$

The ghost propagator $G_{\alpha \beta}^{\omega(\xi)}(p)$ is trivially obtained from eq. (8.4.11):

$$
\begin{equation*}
G_{\alpha \beta}^{\omega(\xi)}(p)=\frac{i \delta_{\alpha \beta}}{p^{2}-\xi \mu_{\alpha}^{2}} \tag{8.4.14}
\end{equation*}
$$

The scalar propagators are different for NG and non-NG bosons. The total mass matrix in the Lagrangian reads

$$
\begin{equation*}
M_{m n}^{2}=\left.\frac{\partial^{2} V}{\partial \phi_{m} \partial \phi_{n}}\right|_{\phi=v}+\xi g^{2}\left(i t_{m p}^{a} v_{p}\right)\left(i t_{n q}^{a} v_{q}\right) \tag{8.4.15}
\end{equation*}
$$

The mass matrix (8.4.15), acting on the NG boson directions $\left(t_{n s}^{b} v_{s}\right)$, gives

$$
\begin{equation*}
M_{m n}^{2}\left(t_{n s}^{b} v_{s}\right)=\xi g^{2}\left(i t_{m p}^{a} v_{p}\right)\left(i t_{n q}^{a} v_{q}\right)\left(t_{n s}^{b} v_{s}\right)=\xi \delta_{a b} \mu_{a}^{2}\left(t_{m p}^{a} v_{p}\right) \tag{8.4.16}
\end{equation*}
$$

where the first term in eq. (8.4.15) vanishes thanks to the relation (8.3.6). In a generic $\xi$-gauge, the would-be NG bosons have a mass which is $\sqrt{\xi}$ times the gauge boson mass. On the other hand, for the non-NG boson directions, defined by eq. (8.4.7), it is the second term in eq. (8.4.15) that vanishes, giving as masses the eigenvalues $m_{i}^{2}$ of $\partial^{2} V / \partial \phi_{m} \phi_{n}$. We then get as scalar propagators

$$
\begin{align*}
G_{i j}^{S}(p) & =\frac{i \delta_{i j}}{p^{2}-m_{i}^{2}}, \quad \text { (no NG bosons) }  \tag{8.4.17}\\
G_{a b}^{S(\xi)}(p) & =\frac{i \delta_{a b}}{p^{2}-\xi \mu_{a}^{2}}, \quad \text { (NG bosons) }
\end{align*}
$$

where the index $i$ runs over all, but the NG bosons, scalar directions.
In the unitary gauge $\xi \rightarrow \infty$, the NG bosons decouple and can be seen as eaten by the gauge fields. In any other gauge they should be kept. Notice how the gauge boson propagator (8.4.13) goes like $1 / p^{2}$ for large momenta for any finite values of $\xi$ and it is only in the unitary gauge that it behaves as $p^{0}$. This makes any finite $\xi$-gauge more suitable than the unitary gauge to address renormalizability properties of spontaneously broken gauge theories. Commom choices of $\xi$-gauges are the Landau gauge $\xi=0$, in which ghosts and NG bosons are massless, and the Feynman gauge $\xi=1$, in which the gauge propagator simplifies considerably.

### 8.5 The Goldstone Boson Equivalence Theorem ${ }^{\star \star}$

The Goldstone boson equivalence theorem, for short denoted equivalence theorem (ET) in what follows, is a theorem about the behaviour of the scattering amplitude of longitudinally polarized gauge bosons in a spontaneously broken gauge theory [24]. The theorem states that at high enough energies the scattering amplitude of longitudinally polarized gauge bosons is the same as the one obtained by replacing the gauge boson with its corresponding "eaten" Goldstone boson, considered as a physical state. The relevance of the ET relies in the fact that the scattering amplitude of massive gauge bosons is indeed dominated by the longitudinal polarizations at high energies.

More precisely, the ET states that for $E \gg m_{A}$

$$
\begin{equation*}
\epsilon_{L, \alpha_{1}}^{\mu_{1}}\left(p_{1}\right) \ldots \epsilon_{L, \alpha_{n}}^{\mu_{n}}\left(p_{n}\right) S_{\mu_{1} \ldots \mu_{n} \ldots}^{\alpha_{1} \ldots \alpha_{n} \ldots}\left(p_{1}, \ldots, p_{n}, \ldots\right)=S^{\alpha_{1} \ldots \alpha_{n} \ldots}\left(p_{1}, \ldots, p_{n}, \ldots\right)+\mathcal{O}\left(\frac{m_{A}}{E}\right) . \tag{8.5.1}
\end{equation*}
$$

In eq. (8.5.1), $S_{\mu_{1} \ldots \mu_{n} \ldots}^{\alpha_{1} \ldots \alpha_{n}}$ is the $S$-matrix element for the scattering of $n$ longitudinally polarized vector bosons with mass $m_{A}$ (taken equal for simplicity) with possibly other physical states, unspecified in eq. (8.5.1) and encoded in the second $\ldots$ in $S^{\mu_{1} \ldots \mu_{n} \ldots}$, while $S^{\alpha_{1} \ldots \alpha_{n} \ldots}$ is the scattering of their corresponding $n$ would-be Goldstone bosons $\phi$ (with the same physical states), treated as physical particles. In order to prove the ET we start by the generalization of eq. (6.3.24), which still applies in spontaneously broken gauge theories:

$$
\begin{equation*}
\langle\alpha|[Q, \mathcal{O}]|\beta\rangle=0, \tag{8.5.2}
\end{equation*}
$$

where $|\alpha\rangle$ and $|\beta\rangle$ are arbitrary physical states and $\mathcal{O}$ is any operator (not necessarily gauge invariant) of the theory. If we take $\mathcal{O}=\omega_{\alpha}^{\star}$ and work with the auxiliary field $H_{\alpha}$ (recall section 6.3), we get

$$
\begin{equation*}
\langle\alpha| H_{\alpha}|\beta\rangle=0 \tag{8.5.3}
\end{equation*}
$$

In the $R_{\xi}$-gauges defined in eq. (8.4.10), solving for $H_{\alpha}$ we have

$$
\begin{equation*}
H_{\alpha}=\frac{1}{\xi}\left(\partial_{\mu} A_{\alpha}^{\mu}+i \xi g \phi_{n}^{\prime} t_{n m}^{\alpha} v_{m}\right) \tag{8.5.4}
\end{equation*}
$$

For simplicity, let us consider the situation where all massive gauge fields have a common mass $m_{A}$ and let us define the would-be GB fields

$$
\begin{equation*}
m_{A} \pi^{\alpha} \equiv g \phi_{n}^{\prime} t_{n m}^{\alpha} v_{m} \tag{8.5.5}
\end{equation*}
$$

The relation (8.5.3) implies that any connected Green function of physical states involving $H_{\alpha}$ must vanish. Denoting collectively by $\Phi$ any physical field (transverse or longitudinal gauge field, physical scalar, matter field), we have

$$
\begin{equation*}
\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) H_{\alpha}(p)\right\rangle=0 \tag{8.5.6}
\end{equation*}
$$

S-matrix elements are given by the amputated Green-functions. In terms of these, eq. (8.5.6) becomes

$$
\begin{equation*}
p^{\mu} G_{\mu \nu}^{(\xi)}(p)\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) A^{\nu}(p)\right\rangle_{a m p}=-\xi m_{A} G^{S(\xi)}(p)\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) \pi(p)\right\rangle_{a m p} \tag{8.5.7}
\end{equation*}
$$

where the propagators are those defined in eqs. (8.4.13) and (8.4.17). We have

$$
\begin{equation*}
p^{\mu} G_{\mu \nu}^{(\xi)}(p)=\frac{-i p_{\nu}\left(p^{2}-\xi m_{A}^{2}-(1-\xi) p^{2}\right)}{\left(p^{2}-m_{A}^{2}\right)\left(p^{2}-\xi m_{A}^{2}\right)}=\frac{-i p_{\nu} \xi}{\left(p^{2}-\xi m_{A}^{2}\right)}=-p_{\nu} \xi G^{S(\xi)}(p) \tag{8.5.8}
\end{equation*}
$$

Plugging eq.(8.5.8) in eq.(8.5.7) gives

$$
\begin{equation*}
\frac{p^{\mu}}{m_{A}}\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) A_{\mu}(p)\right\rangle_{a m p}=\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) \pi(p)\right\rangle_{a m p} \tag{8.5.9}
\end{equation*}
$$

Notice that the $\xi$-dependence completely dropped. The longitudinal polarization for a vector with 4-momentum $p_{\mu}=(E, \vec{p})$ is given by

$$
\begin{equation*}
\epsilon_{\mu}^{L}(p)=\frac{1}{m_{A}}\left(|\vec{p}|, \frac{\vec{p}}{|\vec{p}|} E\right) \tag{8.5.10}
\end{equation*}
$$

For $E \gg m_{A},|\vec{p}|=E\left(1+\mathcal{O}\left(m_{A}^{2} / E^{2}\right)\right)$ and hence

$$
\begin{equation*}
\epsilon_{\mu}^{L}(p)=\frac{p_{\mu}}{m_{A}}+\mathcal{O}\left(\frac{m_{A}}{E}\right) \tag{8.5.11}
\end{equation*}
$$

We finally get

$$
\begin{equation*}
\epsilon_{L, \alpha}^{\mu}(p)\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) A_{\mu}(p)\right\rangle_{a m p}=\left\langle\Phi\left(q_{1}\right) \ldots \Phi\left(q_{n}\right) \pi_{\alpha}(p)\right\rangle_{a m p}+\mathcal{O}\left(\frac{m_{A}}{E}\right) \tag{8.5.12}
\end{equation*}
$$

The above derivation is easily generalized to multiple external longitudinal vector fields. If we take $\mathcal{O}=\omega_{\alpha}^{\star} \prod_{i=1}^{n-1} H_{\alpha_{i}}$ in eq. (8.5.2), we immediately get that the Green functions with an arbitrary number of insertions of $H_{\alpha}$ must vanish. Proceeding as above leads to

$$
\begin{equation*}
\left(\prod_{i=1}^{n} \epsilon_{L, \mu_{i}}^{\alpha_{i}}\left(p_{i}\right)\right)\left\langle A_{\mu_{1}}^{\alpha_{1}}\left(p_{1}\right) \ldots A_{\mu_{n}}^{\alpha_{n}}\left(p_{n}\right) \ldots\right\rangle_{a m p}=\left\langle\pi_{\alpha_{1}}\left(p_{1}\right) \ldots \pi_{\alpha_{n}}\left(p_{n}\right) \ldots\right\rangle_{a m p}+\mathcal{O}\left(\frac{m_{A}}{E}\right) \tag{8.5.13}
\end{equation*}
$$

that proves the relation (8.5.1). We conclude by noticing that the proof above applies only to tree-level amplitudes. Indeed, while the relation (8.5.2) and its generalizations are exact to all orders in perturbation theory, the form of the amputated Green functions crucially depends on the form of the propagator. In going from eq. (8.5.7) to eq. (8.5.9) we have used the tree-level form of the propagators. At the quantum level the relation (8.5.13) gets corrections, that however can be reabsorbed with a proper choice of the gauge fixing functional. We do not discuss these subtleties further.

### 8.6 Effective Field Theories for Broken Symmetries*

Massless NG bosons are a generic prediction of the Goldstone's theorem, no matter whether the underlying theory is weakly or strongly coupled. At sufficiently low energies we can then integrate out massive degrees of freedom and write down an EFT for the NG bosons, possibly interacting with other possibly present light fields, such as gauge fields. We will see in this section that, surprisingly enough, a lot can be said about this EFT without even knowing the underlying UV Lagrangian (both spectrum and interactions). All we need to know is the symmetry breaking pattern $G \rightarrow H$, namely the starting and final global groups. The key point is that the EFT should be not only invariant under the subgroup $H$, but under the entire group $G$ (this is precisely what we mean by spontaneous symmetry breaking!). However, while the symmetries for $H$ are manifest and linearly realized on the NG bosons, the ones for $G / H$ are non-linearly realized and in general complicated. Group theoretically considerations would allow us to find the proper way to repackage the NG bosons into fields that have relatively simple transformation properties so that a general EFT can be found.

Let us start by simply writing down the general commutation relations among the generators $t_{i} \in \mathcal{H}$ and $t_{a} \in \mathcal{G} / \mathcal{H}, \mathcal{G}$ and $\mathcal{H}$ being the Lie algebras of the corresponding groups. We have

$$
\begin{equation*}
\left[t_{i}, t_{j}\right]=i C_{i j k} t_{k}, \quad\left[t_{i}, t_{a}\right]=i C_{i a b} t_{b}, \quad\left[t_{a}, t_{b}\right]=i C_{a b c} t_{c}+i C_{a b i} t_{i} \tag{8.6.1}
\end{equation*}
$$

We take here the structure constants completely antisymmetric in its indices. Since $H$ is a subgroup of $G, C_{i j a}=-C_{i a j}=0$. Let $\phi_{m}$ be the fields responsible for the $G \rightarrow H$
spontaneous symmetry breaking pattern (again in a basis where all the fields are real). Since we eventually want to write down an EFT for the NG boson fields, integrating out the other model-dependent degrees of freedom present in the theory, we have to find a way to disentangle the NG bosons from the rest. In other words, we have to find a way to generalize for an arbitrary symmetry breaking pattern $G \rightarrow H$, the radial coordinates decomposition (8.4.3) that in the $U(1) \rightarrow \emptyset$ allows to disentangle the NG boson $(\theta)$ from the massive excitations $(\rho)$. For this purpose let us define fields $\tilde{\phi}$ that do not contain NG boson field directions:

$$
\begin{equation*}
\phi(x)=\gamma(x) \tilde{\phi}(x), \quad \gamma(x) \in G \tag{8.6.2}
\end{equation*}
$$

namely such that

$$
\begin{equation*}
\tilde{\phi}^{t}(x) t_{\alpha} v=0, \quad \forall \alpha . \tag{8.6.3}
\end{equation*}
$$

In eq.(8.6.2) and in what follows we omit for simplicity the matrix indices $m, n$. In eq.(8.6.3) we clearly have $t_{i} v=0$ and $t_{a} v \neq 0$, so this equation is non trivial only along the broken directions. In this basis the NG bosons all sit in the matrix field $\gamma(x)$. The definition of $\tilde{\phi}$ is not unique. If $\tilde{\phi}$ satisfies (8.6.3), so does $\tilde{\phi}^{\prime}=\tilde{\phi}+c_{i} t_{i} \tilde{\phi}$, for arbitrary coefficients $c_{i}$. Indeed,

$$
\begin{align*}
\tilde{\phi}^{\prime t} t_{\alpha} v & =\left(\tilde{\phi}^{t}+c_{i} \tilde{\phi}^{t} t_{i}^{t}\right) t_{\alpha} v=\left(\tilde{\phi}^{t}-c_{i} \tilde{\phi}^{t} t_{i}\right) t_{\alpha} v \\
& =-c_{i} \tilde{\phi}^{t} t_{i} t_{a} v=-c_{i} \tilde{\phi}^{t}\left[t_{i}, t_{a}\right] v=-i c_{i} C_{i a b} \tilde{\phi}^{t} t_{b} v=0 . \tag{8.6.4}
\end{align*}
$$

The field matrix $\gamma(x)$ is therefore defined only modulo $x$-dependent transformations under $H: \gamma(x) \sim \gamma(x) h(x)$, with $h(x) \in H$. Given this equivalence class, we can always choose as representative a $\gamma(x)$ of the form

$$
\begin{equation*}
\gamma(x)=e^{i \xi_{a}(x) t_{a}} \tag{8.6.5}
\end{equation*}
$$

where $\xi^{a}$ are essentially the NGB's fields. Under a global $G$ transformation, we have $\phi \rightarrow \phi^{\prime}=g \phi=g \gamma(x) \tilde{\phi}(x)$. The field $\phi^{\prime}$ can also be decomposed as in eq. (8.6.2). In general, we will have

$$
\begin{equation*}
\phi^{\prime}(x)=\gamma\left(\xi_{a}^{\prime}(x)\right) h\left(\xi_{a}(x), g\right) \tilde{\phi}(x), \tag{8.6.6}
\end{equation*}
$$

for some $h\left(\xi_{a}(x), g\right) \in H$, so that $\phi^{\prime}=\gamma\left(\xi_{a}^{\prime}\right) \tilde{\phi}^{\prime},{ }^{5}$ with

$$
\begin{align*}
\tilde{\phi}^{\prime} & =h\left(\xi_{a}, g\right) \tilde{\phi}  \tag{8.6.7}\\
\gamma\left(\xi_{a}^{\prime}\right) & =g \gamma\left(\xi_{a}\right) h^{-1}\left(\xi_{a}, g\right) \tag{8.6.8}
\end{align*}
$$

[^54]Notice that the field $\tilde{\phi}(x)$, under global $G$ transformations, transforms as effectively local transformations under $H$. The transformations of $\xi_{a}$ and $\tilde{\phi}$ simplify considerably when $g=h \in H$. The second commutation relations in eq. (8.6.1) shows that the broken generators $t_{a}$ transform linearly under transformations of $h$, namely

$$
\begin{equation*}
h t_{a} h^{-1}=t_{b} R_{b a}(h) \tag{8.6.9}
\end{equation*}
$$

where $R$ is some representation (in general reducible) of $H$. The very same commutation relations shows that when $g=h$, the factor $h(\xi(x), g)$ defined in eq. (8.6.6) reduces to the global $x$-independent element $h: h(\xi(x), g)=h$. We then have, using also eq. (8.6.9)

$$
\begin{equation*}
\gamma^{\prime} \equiv \gamma\left(\xi^{\prime}\right)=h \gamma(\xi) h^{-1}=\gamma(R(h) \xi) . \tag{8.6.10}
\end{equation*}
$$

Under $H$, then, the transformations of both $\xi$ and $\tilde{\phi}$ are simple and linear:

$$
\begin{align*}
\xi_{a}^{\prime}(x) & =R_{a b}(h) \xi_{b}(x) \\
\tilde{\phi}^{\prime}(x) & =h \tilde{\phi}(x) \tag{8.6.11}
\end{align*}
$$

On the other hand, for infinitesimal transformations along $G / H$, the NG fields $\xi_{a}$ transform as a shift, at leading order in the field fluctuations. This is immediately visible from eq.(8.6.8). For $g=1+i \epsilon_{a} t_{a}$, indeed, we have

$$
\begin{equation*}
\delta \xi_{a}=\epsilon_{a}+\mathcal{O}\left(\xi^{2}, \epsilon \xi\right) \tag{8.6.12}
\end{equation*}
$$

We now assume that the structure constants $C_{a b c}$ in the third commutator in eq. (8.6.1) vanish, in which case the coset $G / H$ is called symmetric and the formalism describing the interactions of NG bosons can be simplified. We refer the reader to subsection 8.9 where a more general formalism allows us to also consider non-symmetric cosets. With $C_{a b c}=0$, the Lie algebra defined by the relations (8.6.1) is invariant under the $\mathbf{Z}_{2}$ symmetry $R$ under which the broken generators change sign:

$$
\begin{equation*}
R\left(t_{i}\right)=t_{i}, \quad R\left(t_{a}\right)=-t_{a} \tag{8.6.13}
\end{equation*}
$$

We clearly have $R(\gamma)=\gamma^{-1}$. Under $G$ transformations

$$
\begin{equation*}
\gamma^{\prime}=g \gamma h^{-1} . \tag{8.6.14}
\end{equation*}
$$

On the other hand

$$
\begin{align*}
\gamma^{\prime-1}= & R\left(\gamma^{\prime}\right)=R(g) R(\gamma) R\left(h^{-1}\right)=R(g) \gamma^{-1} h^{-1} \\
& \Longrightarrow \gamma^{\prime}=h \gamma R(g)^{-1} \tag{8.6.15}
\end{align*}
$$

Using both eqs. (8.6.14) and (8.6.15), we see that

$$
\begin{equation*}
U^{\prime} \equiv \gamma^{\prime 2}=g \gamma h^{-1} h \gamma R(g)^{-1}=g U R(g)^{-1} . \tag{8.6.16}
\end{equation*}
$$

In other words, while $\gamma$ does not transform linearly under transformations of $G$, its square $U=\gamma^{2}$ does. This is a significant simplification, because it allows us to no longer worry of the complicated factor $h(x)$.

We have found that an EFT for the NG bosons, invariant under $G$, is best written when the NG bosons are repackaged in the matrix $U$. Let us see which kind of invariant terms we can write using $U$ and $U^{\dagger}$. Let us first focus on potential terms, i.e. no derivatives. Since $R(g) \neq g$, terms of the form $U^{n}$ or $\left(U^{\dagger}\right)^{n}$ are forbidden, the only allowed invariant terms being trace of operators proportional to $U^{\dagger} U$. But $U^{\dagger} U=1$, so no potential term is indeed allowed! This should not come as a surprise. We know that NG bosons must be massless. If a potential term would have been allowed, generally it would give a mass to them. The absence of any potential could have been anticipated by the shift symmetry (8.6.12), which is broken by any potential term. Non-trivial invariant operators can instead be written using derivatives. The leading, two-derivative, term involving the NG bosons and invariant under the symmetry (8.6.16) is

$$
\begin{equation*}
\mathcal{L}=\frac{f_{\pi}^{2}}{8} \operatorname{tr}\left(\partial_{\mu} U \partial^{\mu} U^{\dagger}\right)+\ldots \tag{8.6.17}
\end{equation*}
$$

where $f_{\pi}$ is a mass scale, related to the scale of spontaneous symmetry breaking, introduced for dimensional reasons, and ... stands for higher dimensional operators involving four or higher number of derivatives. The expansion of the the matrices $U$ and $U^{\dagger}$ in powers of $\xi^{a}$ in eq. (8.6.17) allows us to find the explicit form for an infinite number of leading interactions among the NG bosons, all fixed by a single term in the Lagrangian! The relation between the fields $\xi_{a}$ and the canonical NG fields $\pi_{a}$ defined in eq.(8.2.12) is obtained by computing the form of the broken currents $J_{\mu}^{a}$ from eq.(8.6.17). Using the trick of promoting a global symmetry to a local one, so that $\delta_{\epsilon} \mathcal{L}=-\partial_{\mu} \epsilon J^{\mu}$, we have,

$$
\begin{equation*}
\delta_{\epsilon^{a}} \mathcal{L}=f_{\pi}^{2} \partial_{\mu} \epsilon_{a} \partial^{\mu} \xi_{b} \operatorname{Tr} t^{a} t^{b}+\ldots=f_{\pi}^{2} \partial_{\mu} \epsilon^{a} \partial^{\mu} \xi_{a}+\ldots, \Rightarrow J_{\mu}^{a}=-f_{\pi}^{2} \partial_{\mu} \xi^{a} \tag{8.6.18}
\end{equation*}
$$

where we have used the leading order transformation (8.6.12) for the $\xi$ 's and taken $\operatorname{Tr} t_{a} t_{b}=$ $\delta_{a b}$. Matching eqs.(8.2.8) and (8.2.12) with eq.(8.6.18) and demanding that the kinetic term (8.6.17) is canonical in terms of the $\pi_{a}$ fields gives

$$
\begin{equation*}
\xi_{a}=\frac{F_{a b}}{f_{\pi}^{2}} \pi_{b} \tag{8.6.19}
\end{equation*}
$$

Since the NGB's transform linearly under $H$ transformations, the decay constant matrix $F_{a b}$ should correspondingly be invariant under such transformations. This is in general a strong constraint on the form of $F_{a b}$.

As we mentioned, all the interactions among the NG bosons, including their kinetic terms, are efficiently encoded in the single term (8.6.17). This is the first of an infinite series of higher derivative operators involving the matrices $U$. By expanding $U$ in terms of field fluctuations, we see that the effective coupling constant of all the interaction terms that can arise from eq. (8.6.17) is $E^{2} / f_{\pi}^{2}$. For sufficiently low-energies $E \ll f_{\pi}$, this term is the most important one, all the higher derivative interactions being suppressed by additional powers of $E / f_{\pi}$. The EFT Lagrangian (8.6.17) is of course non-renormalizable and becomes unreliable when $E \sim f_{\pi}$, in which case an alternative description is needed. More precisely, the energy scale where we expect our effective theory to break down is $E=\Lambda \simeq 4 \pi f_{\pi}$, including the $4 \pi$ factor coming from loops (analogue of the effective QED expansion parameter $\left.e^{2} /\left(16 \pi^{2}\right)\right)$. This is the energy scale where the non-NGBs degrees of freedom that we have (implicitly) integrated out start to matter and must be included to possibly embed the theory in a consistent UV complete model.

### 8.6.1 Adding Gauge Fields ${ }^{\star}$

Other light fields might appear in the EFT for the NG bosons, most notably gauge fields. Like the NG bosons, gauge particles can be naturally light or massless and their interactions are governed by gauge invariance. In fact, gauge interactions in this formalism do not pose any difficulty. Let us suppose that a subgroup $H_{g}$ of $G$, distinct from $H$, is gauged. A convenient way to proceed is to pretend that the whole group is gauged, namely take $H_{g}=G$. Since $U$ transforms linearly under $G$, the local symmetry is implemented in the usual way discussed in chapter 6 for matter fields; one introduces a gauge field $A_{\mu}$ and replace ordinary derivatives of $U$ with covariant ones. Under a gauge transformation $g$ we have

$$
\begin{equation*}
A_{\mu}^{g}=g A_{\mu} g^{-1}-i\left(\partial_{\mu} g\right) g^{-1}, \tag{8.6.20}
\end{equation*}
$$

where $A_{\mu}=A_{\mu}^{i} t_{i}+A_{\mu}^{a} t_{a}$ is the full (non-canonically normalized) gauge field. Given the transformation (8.6.16) and the above relation, the covariant derivative

$$
\begin{equation*}
D_{\mu} U=\partial_{\mu} U-i A_{\mu} U+i U A_{\mu}^{R} \tag{8.6.21}
\end{equation*}
$$

where $A_{\mu}^{R}=A_{\mu}^{i} t_{i}-A_{\mu}^{a} t_{a}$, transforms as $D_{\mu} U^{g}=g\left(D_{\mu} U\right) R(g)^{-1}$. Since the theory is now invariant under local transformations of $G$, by a proper choice of $g$ we might completely remove the NGB's, i.e. we can set $U=1$. This is nothing else than the unitary gauge in a non-linear realization of the symmetry. In this gauge the covariantization of eq.(8.6.17) leads to mass terms for the gauge fields $A_{\mu}-A_{\mu}^{R}=2 A_{\mu}^{a} t_{a}$, namely to those along the broken direction, as expected from the Higgs mechanism.

The generic case $H_{g} \subset G$ can be deduced from the case $H_{g}=G$ by simply switching off the gauge fields that are not in the direction of $H_{g}$. Denoting by $t^{\hat{\alpha}}$ the generators of $G$ that are gauged, in general we have

$$
\begin{equation*}
t^{\hat{\alpha}}=g_{\hat{\alpha} \alpha} t^{\alpha} \tag{8.6.22}
\end{equation*}
$$

where $g_{\hat{\alpha} \alpha}$ are the gauge coupling constants associated to the gauging, all different in the most general case. We now set

$$
\begin{equation*}
A_{\mu}^{\alpha}=g_{\hat{\alpha} \alpha} A_{\mu}^{\hat{\alpha}} \tag{8.6.23}
\end{equation*}
$$

where $A_{\mu}^{\hat{\alpha}}$ are the actual dynamical gauge fields. In the unitary gauge, as expected, only the gauge fields along the broken directions get a mass with the corresponding NGB's being eaten by them. Notice that gauging a subgroup of $H_{g}$ explicitly breaks the global symmetry $G$, since it determines a specific direction in field space. Taking into account the normalization (8.6.19) of the broken generators, the gauge field mass matrix reads

$$
\begin{equation*}
\mu_{\hat{\alpha} \hat{\beta}}^{2}=F_{a c} F_{a d} g_{\hat{\alpha} c} g_{\hat{\beta} d} \tag{8.6.24}
\end{equation*}
$$

This relation is the generalization of eq. (8.4.8). The latter only applies for weakly coupled descriptions of the spontaneous symmetry breaking in terms of free fields.

## 8.7 $S U(3)_{V} \times S U(3)_{A} \rightarrow S U(3)_{V}$ : Mesons in QCD ${ }^{\star}$

Consider QCD with three active quarks (say, the up, down and strange quark) in the limit in which they are all massless:

$$
\begin{equation*}
\mathcal{L}_{Q C D}=-\frac{1}{4} G_{\mu \nu}^{a} G_{a}^{\mu \nu}+\bar{Q} \not \partial D Q \tag{8.7.1}
\end{equation*}
$$

where $Q=\left(\begin{array}{ll}u d\end{array}\right)^{t} .{ }^{6}$ In addition to the $S U(3)_{c}$ gauge (color) symmetry, the Lagrangian (8.7.1) is invariant under an additional $S U(3)_{V} \times S U(3)_{A}$ global (flavour) symmety acting on $Q$ as

$$
\begin{equation*}
Q \rightarrow \exp \left(i \theta_{a}^{V} \lambda_{a}+i \theta_{a}^{A} \lambda_{a} \gamma_{5}\right) Q \tag{8.7.2}
\end{equation*}
$$

where $\lambda_{a}$ are the $S U(3)$ Gell-Mann matrices, normalized such that $\operatorname{tr} \lambda_{a} \lambda_{b}=\delta_{a b} / 2$ in the fundamental representation. At sufficiently low energies, when the QCD coupling constant becomes strong, a quark bilinear gets a non-vanishing vacuum expectation value:

$$
\begin{equation*}
\left\langle\bar{Q}_{i} Q_{j}\right\rangle=\hat{\Lambda}^{3} \delta_{i j} \tag{8.7.3}
\end{equation*}
$$

[^55]that breaks $G=S U(3)_{V} \times S U(3)_{A}$ down to $H=S U(3)_{V}$. Here $\hat{\Lambda}$ is the scale where the chiral symmetry breakdown occurs. The eight NG bosons are encoded in the matrix field
\[

$$
\begin{equation*}
\gamma(x)=e^{i \lambda^{a} \xi^{a}(x)} \tag{8.7.4}
\end{equation*}
$$

\]

The $S U(3)_{V}$ unbroken symmetry fixes the parameters $F_{a b}$ to be proportional to the identity. We take $F_{a b} \equiv \delta_{a b} f_{\pi}$ in eq.(8.6.19) and write the 8 NG bosons as

$$
\lambda_{a} \xi_{a}=\frac{1}{f_{\pi}}\left(\begin{array}{ccc}
\frac{1}{\sqrt{2}} \pi_{0}+\frac{1}{\sqrt{6}} \eta^{0} & \pi^{+} & K^{+}  \tag{8.7.5}\\
\pi^{-} & -\frac{1}{\sqrt{2}} \pi_{0}+\frac{1}{\sqrt{6}} \eta^{0} & K^{0} \\
\bar{K}^{-} & \bar{K}^{0} & -\sqrt{\frac{2}{3}} \eta^{0}
\end{array}\right)
$$

where $f_{\pi}$ is identified with the pion decay constant: $f_{\pi} \simeq 92 \mathrm{MeV}$. With this normalization the 8 spin zero mesons $\pi^{ \pm}, \pi^{0}$, etc. are all canonically normalized.

The commutator of two vector or two axial transformations is a vector transformation, while the commutator of a vector and axial transformation is an axial transformation. The schematic commutation relations for $S U(3)_{V} \times S U(3)_{A}$ are then $[V, V]=V,[V, A]=A$, $[A, A]=V$, where $V$ and $A$ schematically represent the $S U(3)_{V}$ and $S U(3)_{A}$ generators. This is a symmetric coset space, being this algebra invariant under the automorphism action $V \rightarrow V, A \rightarrow-A$. We can also construct L and R transformations defined as $V=(L+R) / 2, A=(L-R) / 2$, under which the above automorphism exchanges $L$ and $R: L \leftrightarrow R$. According to the general results of section 8.6, the matrix $U=\gamma^{2}$ transforms homogeneously under $g$. We have ${ }^{7}$

$$
\begin{equation*}
U \rightarrow R U L^{\dagger} \tag{8.7.6}
\end{equation*}
$$

with obvious notation. In other words, $U$ transforms as a $(\mathbf{3}, \overline{\mathbf{3}})$ representation of $S U(3)_{R} \times$ $S U(3)_{L}$. Correspondingly, the kinetic terms and leading interactions among the 8 mesons $\pi, K$ and $\eta$ are collected in the single term (8.6.17). As we have seen in the previous section, the effective field theory is expected to break down at $E=\Lambda \simeq 4 \pi f_{\pi} \sim 1 \mathrm{GeV}$. It is natural to assume that $\Lambda \simeq \hat{\Lambda}$, the scale where chiral symmetry breaking occurs. This scale is related to the scale $\Lambda_{Q C D}$ where quarks confine (confinement scale), but it is not the same scale, the chiral symmetry breaking scale being slightly higher than the confinement one. They are also conceptually different. In fact, there exist vacua in gauge (supersymmetric) theories where it is believed that quarks confine but no chiral symmetry breaking occurs.

[^56]In the real world, of course, the up, down and strange quarks are not massless and the spin zero mesons $\pi, K$ and $\eta$ are not exact NG bosons. The two things are related. The actual up, down and strange masses explicitly break the $S U(3)_{V} \times S U(3)_{A}$ symmetry, but are small enough, compared to $\Lambda$, to be considered as a perturbation of the QCD Lagrangian (8.7.1) (this would not be the case for the charm, bottom and top, that are all heavier than $\Lambda)$. We can get the relation between the quark and meson masses by using the following trick. We formally promote the quark mass term

$$
M=\left(\begin{array}{ccc}
m_{u} & 0 & 0  \tag{8.7.7}\\
0 & m_{d} & 0 \\
0 & 0 & m_{s}
\end{array}\right)
$$

to be an external source that transforms under $S U(3)_{L} \times S U(3)_{R}$ as $M \rightarrow L M R^{\dagger}$, so that the term $\bar{Q}_{L} M Q_{R}+h . c$. is made invariant. We have morally promoted the mass term to a so called spurion field, whose vacuum expectation value coincides with the mass matrix $M$ and whose dynamics is frozen. Spurions are often used in QFT to formally restore explicitly broken symmetries. In so doing, one proceeds in the computation demanding the full symmetry, and only at the end of the computation one plugs back the original form of the breaking term into the spurion field. This is a very efficient way to make manifest how breaking terms affect physical quantities. Thanks to this trick, at low-energies we can construct $S U(3)_{L} \times S U(3)_{R}$ invariant terms using both $U$ and $M$. If $M$ is small, we can consider terms linear in $M$ only. We simply have

$$
\begin{equation*}
\mathcal{L}_{m}=c f_{\pi}^{3} \operatorname{tr} U M+\text { h.c. }, \tag{8.7.8}
\end{equation*}
$$

where $c$ is an undetermined dimensionless coefficient. By plugging in eq. (8.7.8) the actual form of the NG fields (8.7.5) and of the mass term (8.7.7) we get, after some simple algebra:

$$
\begin{align*}
m_{\pi^{ \pm}}^{2} & =C\left(m_{u}+m_{d}\right) \\
m_{K^{ \pm}}^{2} & =C\left(m_{u}+m_{s}\right)  \tag{8.7.9}\\
m_{K^{0}}^{2} & =C\left(m_{d}+m_{s}\right),
\end{align*}
$$

where $C=2 c f_{\pi}$. The $\pi^{0}$ and $\eta^{0}$ mix with each other. Expanding for $m_{u, d} \ll m_{s}$ we get

$$
\begin{align*}
m_{\pi^{0}}^{2} & \simeq C\left(m_{u}+m_{d}\right) \\
m_{\eta}^{2} & \simeq C \frac{4 m_{s}+m_{u}+m_{d}}{3} . \tag{8.7.10}
\end{align*}
$$

It is surprising that using symmetries only we can fix the meson masses in terms of the quark masses and one unknown coefficient $C$. The above relations are a particular
example of eq. (8.3.9), showing that in general the masses of the pseudo NG bosons are linearly dependent on the source of the symmetry breaking violating term in the Lagrangian. In deriving eqs. (8.7.9) and (8.7.10) we have actually neglected another relevant source of explicit breaking of the $S U(3)_{L} \times S U(3)_{R}$ symmetry, given by the electric charge. As explained in section 8.6, in this EFT framework electromagnetic interactions are introduced by identifying a proper $U(1)$ subgroup and gauge it. This is a subgroup of the unbroken $S U(3)_{V}$, and its explicit form is found by looking at the infinitesimal transformations of the quark fields $Q$. We have $\delta_{\epsilon, E M} Q=i \epsilon Q_{e l} Q$, where

$$
Q_{e l}=\left(\begin{array}{ccc}
\frac{2}{3} & 0 & 0  \tag{8.7.11}\\
0 & -\frac{1}{3} & 0 \\
0 & 0 & -\frac{1}{3}
\end{array}\right)
$$

according to the actual electric charges of the up, down and strange quarks. We therefore get the leading order Lagrangian

$$
\begin{equation*}
\mathcal{L}_{K i n+\gamma}=\frac{f_{\pi}^{2}}{4} \operatorname{tr}\left(D_{\mu} U D^{\mu} U^{\dagger}\right)-\frac{1}{4} F_{\mu \nu} F^{\mu \nu}+\left(c f_{\pi}^{3} \operatorname{tr} U M+\text { h.c. }\right) \tag{8.7.12}
\end{equation*}
$$

with

$$
\begin{equation*}
D_{\mu} U=\partial_{\mu} U+i e Q_{e l} U A_{\mu}-i e U Q_{e l} A_{\mu} \tag{8.7.13}
\end{equation*}
$$

It is straightforward to check that eq. (8.7.12) contains the expected interactions of the QED form for the charged mesons $\pi^{ \pm}$and $K^{ \pm}$. These interactions, at the quantum level, contribute to the mass of the charged mesons by an amount $\Delta_{e l}$, which is the same, at leading order, for $\pi^{ \pm}$and $K^{ \pm}$, thanks to the $S U(3)_{V}$ global symmetry. Eventually, the effect of the QED interactions is to shift the charged meson masses in eq. (8.7.10) by this amount $\Delta_{e l}$ :

$$
\begin{align*}
m_{\pi^{ \pm}}^{2} & \rightarrow m_{\pi^{ \pm}}^{2}+\Delta_{e l}  \tag{8.7.14}\\
m_{K^{ \pm}}^{2} & \rightarrow m_{K^{ \pm}}^{2}+\Delta_{e l}
\end{align*}
$$

The neutral meson masses, at this order, are clearly unaffected by QED interactions. There are various ways in which the mass formulas (8.7.9), (8.7.10) and (8.7.14) can be used. For instance, the following relation holds, independently of $C, \Delta_{e l}$ and the quark masses:

$$
\begin{equation*}
3 m_{\eta}^{2}+2 m_{\pi^{ \pm}}^{2}-m_{\pi^{0}}^{2}=2 m_{K^{ \pm}}^{2}+2 m_{K^{0}}^{2} \tag{8.7.15}
\end{equation*}
$$

and it is experimentally very well reproduced. Alternatively, we might use the meson masses as experimental input parameters to compute the quark masses or, better, their
ratios. One has

$$
\begin{align*}
\frac{m_{u}}{m_{s}} & =\frac{2 m_{\pi^{0}}^{2}-m_{\pi^{ \pm}}^{2}+m_{K^{ \pm}}^{2}-m_{K^{0}}^{2}}{m_{K^{ \pm}}^{2}+m_{K^{0}}^{2}-m_{\pi^{ \pm}}^{2}} \simeq 0.027 \\
\frac{m_{d}}{m_{s}} & =\frac{m_{\pi^{ \pm}}^{2}+m_{K^{0}}^{2}-m_{K^{ \pm}}^{2}}{m_{K^{ \pm}}^{2}+m_{K^{0}}^{2}-m_{\pi^{ \pm}}^{2}} \simeq 0.050 \tag{8.7.16}
\end{align*}
$$

where we have inserted in the last relation the approximate masses for the spin zero mesons.

The actual global symmetry of the QCD Lagrangian (8.7.1) is $U(3)_{V} \times U(3)_{A}$, rather than $S U(3)_{V} \times S U(3)_{A}$, since the $U(1)_{V} \times U(1)_{A}$ transformations

$$
\begin{equation*}
Q \rightarrow \exp \left(i \theta^{V}+i \theta^{A} \gamma_{5}\right) Q \tag{8.7.17}
\end{equation*}
$$

are a symmetry of (8.7.1). The $U(1)_{V}$ symmetry is nothing else than the $U(1)$ baryon number and is unbroken, while $U(1)_{A}$ is manifestly broken by the quark condensate in eq. (8.7.3). Correspondingly, we should have a ninth pseudo NG boson, called $\eta^{\prime}$. It can be shown, by adding the $\eta^{\prime}$ to the matrix field (8.7.4), that its mass $m_{\eta^{\prime}}$ is bounded to be $m_{\eta^{\prime}} \leq \sqrt{3} m_{\pi}$. But no such particle has been observed within this range of energies. This puzzle, denoted the $U(1)_{A}$ problem, has been solved by noticing that the $U(1)_{A}$ symmetry is explicitly broken - by an amount larger than the explicit breaking of the quark masses - by instantons, non-perturbative effects that are outside the content of these lectures. The actual $\eta^{\prime}$ has indeed a much larger mass, $m_{\eta^{\prime}} \simeq 960 \mathrm{MeV}$.

## 8.8 $S O(5) \rightarrow S O(4):$ A Composite Higgs? ${ }^{\star \star}$

The example that follows is a bit more exotic and refers to the electroweak sector of the Standard Model (SM). We have briefly seen in section 7.5 that scalar masses are unnatural, in the sense that they are quadratically sensitive to UV physics that push their values to that scale. This problem clearly applies to elementary scalars, including the Higgs boson in the SM. But what if the Higgs is not elementary, but rather a composite particle, very much like the pion? We do not have a hierarchy problem for the pion mass, since we know that this particle is composite and at some energies of order $\Lambda_{Q C D}$, "diffuse" in its quark constituents. Moreover, pions are actually naturally lighter than $\Lambda_{Q C D}$ itself because, as we have extensively seen in the previous subsection, are pseudo NG bosons of an approximate global symmetry.

Before developing on this idea of considering the Higgs particle as a pseudo NG boson of some spontaneously broken global symmetry, it is useful to recall the symmetries of the
standard, elementary Higgs Lagrangian, in the SM. In the global limit in which we switch off the $S U(2)_{L} \times U(1)_{Y}$ gauge couplings, the SM Higgs sector is

$$
\begin{equation*}
\mathcal{L}_{H}^{S M}=\left(\partial_{\mu} H\right)^{\dagger}\left(\partial^{\mu} H\right)+m^{2} H^{\dagger} H-\frac{\lambda}{2}\left(H^{\dagger} H\right)^{2} \tag{8.8.1}
\end{equation*}
$$

where $H$ is an $S U(2)_{L}$ doublet of the form

$$
\begin{equation*}
H=\frac{1}{\sqrt{2}}\binom{h_{2}+i h_{1}}{h_{4}-i h_{3}} \tag{8.8.2}
\end{equation*}
$$

A closer look at eq. (8.8.1) reveals that the Higgs Lagrangian is invariant under an $S O(4)$, rather than $S U(2)_{L}$, global symmetry. This is manifest if we recast the four real Higgs components $h_{i}(i=1,2,3,4)$ in a 4-plet and notice that eq. (8.8.1) is invariant under the transformations $h_{i} \rightarrow O_{i j} h_{j}$, with $O \in S O(4)$. It is actually more convenient to exploit the local isomorphism of $S O(4)$ with $S U(2) \times S U(2)$. The fundamental 4 of $S O(4)$ becomes a bidoublet $(\mathbf{2}, \mathbf{2})$ of $S U(2) \times S U(2)$. One of the two $S U(2)$ is identified with the original $S U(2)_{L}$, and we denote by $S U(2)_{R}$ the other one. We can easily write the $S O(4)$ generators in the $S U(2)_{L} \times S U(2)_{R}$ basis. The 6 anti-symmetric hermitian generators of $S O(4)$ are proportional to

$$
\begin{equation*}
t_{i j}^{a b}=-t_{i j}^{b a}=\delta_{i}^{a} \delta_{j}^{b}-\delta_{i}^{b} \delta_{j}^{a}, \tag{8.8.3}
\end{equation*}
$$

where $a, b=1, \ldots, 4$ label the generators and $i, j$ their matrix components. A simple check of the algebra reveals that the combinations

$$
\begin{align*}
& t_{L}^{1}=-\frac{i}{2}\left(t^{23}+t^{14}\right), \quad t_{L}^{2}=-\frac{i}{2}\left(t^{31}+t^{24}\right), \quad t_{L}^{3}=-\frac{i}{2}\left(t^{12}+t^{34}\right), \\
& t_{R}^{1}=-\frac{i}{2}\left(t^{23}-t^{14}\right), \quad t_{R}^{2}=-\frac{i}{2}\left(t^{31}-t^{24}\right), \quad t_{R}^{3}=-\frac{i}{2}\left(t^{12}-t^{34}\right), \tag{8.8.4}
\end{align*}
$$

satisfy the commutation relations of the $S U(2)_{L}$ and $S U(2)_{R}$ algebras, respectively. The action of $S U(2)_{L} \times S U(2)_{R}$ is best seen by writing explicitly the Higgs field as a bidoublet using the $2 \times 2$ matrices $\sigma^{\mu}=\left(1, i \sigma_{k}\right)(k=1,2,3)$ :

$$
H_{B D}=\frac{h_{4} 1_{2}+i h_{k} \sigma_{k}}{2}=\frac{1}{2}\left(\begin{array}{cc}
h_{4}+i h_{3} & h_{2}+i h_{1}  \tag{8.8.5}\\
-h_{2}+i h_{1} & h_{4}-i h_{3}
\end{array}\right) .
$$

Under $S U(2)_{L} \times S U(2)_{R}$, the bidoublet (8.8.5) transforms as

$$
\begin{equation*}
H_{B D} \rightarrow g_{L} H_{B D} g_{R}^{\dagger} \tag{8.8.6}
\end{equation*}
$$

with $g_{L} \in S U(2)_{L}, g_{R} \in S U(2)_{R}$. It is manifest from eq. (8.8.6) that when the Higgs field develops a non-vanishing VEV, say $\left\langle h_{4}\right\rangle \equiv v \neq 0$, i.e. $\left\langle H_{B D}\right\rangle \propto 1_{2}$, the $S U(2)_{L} \times S U(2)_{R}$
global symmetry is broken to the diagonal $S U(2)_{c}$ symmetry, with $g_{L}=g_{R}$. This unbroken $S U(2)_{c}$ global symmetry is called custodial symmetry and plays an important role in establishing that the $W$ and $Z$ boson mass ratio at tree-level is

$$
\begin{equation*}
\frac{m_{Z}^{2}}{m_{W}^{2}}=\frac{g^{2}+g^{\prime 2}}{g^{2}} \tag{8.8.7}
\end{equation*}
$$

where $g$ and $g^{\prime}$ are the $S U(2)_{L}$ and $U(1)_{Y}$ gauge coupling constants. In this formalism, the gauge fields are introduced by gauging the whole $S U(2)_{L}$ global group and a subgroup of $S U(2)_{R}$, along the $\sigma_{3}$ generator, which is identified with the $U(1)_{Y}$ symmetry. The full, gauged, Higgs Lagrangian reads

$$
\begin{equation*}
\mathcal{L}_{H}^{S M}=\operatorname{tr}\left(D_{\mu} H_{B D}\right)^{\dagger}\left(D^{\mu} H_{B D}\right)+m^{2} \operatorname{tr} H_{B D}^{\dagger} H_{B D}-\frac{\lambda}{2} \operatorname{tr}\left(H_{B D}^{\dagger} H_{B D}\right)^{2} \tag{8.8.8}
\end{equation*}
$$

where

$$
\begin{equation*}
D_{\mu} H_{B D}=\partial_{\mu} H_{B D}-i g W_{\mu}^{L} H_{B D}+i g^{\prime} H_{B D} W_{\mu}^{R} \tag{8.8.9}
\end{equation*}
$$

and

$$
\begin{equation*}
W_{\mu}^{L}=\frac{1}{2} \sigma_{i} W_{\mu}^{i}, \quad W_{\mu}^{R}=\frac{1}{2} \sigma_{3} B_{\mu} . \tag{8.8.10}
\end{equation*}
$$

The custodial $S U(2)_{c}$ symmetry is broken by the hypercharge coupling $g^{\prime}$ only. ${ }^{8}$ If we set $g^{\prime}=0$, the Lagrangian (8.8.8) is invariant under $S U(2)_{L} \times S U(2)_{R}$ global transformations, provided we rotate $W_{\mu}^{L} \rightarrow g_{L} W_{\mu}^{L} g_{L}^{\dagger}$. The 3 would-be NG bosons associated to the $S U(2)_{L} \times S U(2)_{R} \rightarrow S U(2)_{c}$ breaking are all eaten by the $W^{L}$ 's, and $S U(2)_{c}$ ensures that the 3 vector bosons all have equal masses $m_{W}=g v / 2$. When $g^{\prime} \neq 0$, the custodial symmetry is broken and the 3 massive gauge fields no longer have equal masses. However, the breaking pattern fixes their tree-level masses. From eq. (8.8.9) we immediately see that the gauge bosons along the off-diagonal components of $S U(2)_{L}$ are unaffected by $g^{\prime}$ and retain their masses $m_{W}=g v / 2$. On the other hand, along the $U(1)_{L} \times U(1)_{R}$ gauged subgroups in the $\sigma_{3}$ directions, the gauge field along the unbroken symmetry remains massless (the photon) and the orthogonal direction (the $Z$ ) gets a mass equal to $m_{Z}=\sqrt{g^{2}+g^{\prime 2}} v / 2$. The importance of emphasizing the symmetries of the Higgs Lagrangian becomes clear when we replace the Higgs field by an unspecified sector responsible for the electroweak symmetry breaking. In this more general context, we would conclude that the $W$ and $Z$ boson masses are given by eq. (8.6.24), with $\hat{\alpha}=1 L, 2 L, 3 L, 3 R$. More explicitly, we have

$$
\begin{equation*}
g_{a L b}=g \delta_{a b}, \quad g_{a R b}=g^{\prime} \delta_{a b} \delta_{a 3}, \tag{8.8.11}
\end{equation*}
$$

[^57]where $a, b=1,2,3$ run over the broken $S U(2)$ generators. The crucial point is played by the custodial $S U(2)_{c}$ symmetry, that forces the $F_{a b}$ terms to be proportional to the identity (like $S U(3)_{V}$ in the QCD example before): $F_{a b}=\delta_{a b} F$. Putting all together, the gauge boson mass matrix is of the form
\[

\mu^{2}=F^{2}\left($$
\begin{array}{cccc}
g^{2} & 0 & 0 & 0  \tag{8.8.12}\\
0 & g^{2} & 0 & 0 \\
0 & 0 & g^{2} & g g^{\prime} \\
0 & 0 & g g^{\prime} & g^{\prime 2}
\end{array}
$$\right)
\]

We automatically recover the SM gauge boson masses, in particular the relation (8.8.7), with the identification $F=v / 2$. We conclude by stressing that any sector replacing the usual Higgs in the SM Lagrangian will give the correct leading order electroweak gauge boson masses, provided it includes an $S U(2)_{c}$ global symmetry.

After this long, but necessary, digression on the SM, let us come back to our idea of the Higgs field as a pseudo NG boson. Assuming this idea, the Higgs should be a composite of certain constituents, analogues of quarks and gluons in QCD, that appears at strong coupling at some energy scale. Being the analogue of the pion in QCD, the Higgs particle is expected to be the lightest resonance of the strongly coupled theory, that might include additional heavier resonances (in analogy with the hadron spectroscopy in QCD). Of course, in contrast to the QCD case, we do not know here what is the UV theory that becomes strongly coupled. But we have by now learned that a lot can be said about the dynamics of pseudo NG bosons, by only specifying the group theoretical structure of the spontaneous symmetry breaking pattern. As far as we are concerned, we have to assume that the UV theory, no matter what it is, has an approximate global symmetry group $G$, spontaneously broken to $H$, such that the NG bosons along the $G / H$ directions have the quantum numbers of the SM Higgs. Let us denote by $f_{H}$ the scale where this breaking occurs. The unbroken group $H$ should be large enough to accommodate an $S U(2)_{L} \times U(1)_{Y}$ subgroup, that we will gauge and identify with the SM electroweak gauge group. Finally, we might also demand that $H$ includes the custodial symmetry $S U(2)_{c}$ that ensures the correct tree-level mass ratio (8.8.7) between the SM gauge bosons. The minimal groups that give rise to the 4 Higgs NG bosons and nothing else are

$$
\begin{align*}
& G=S U(3) \rightarrow H=S U(2) \times U(1)  \tag{8.8.13}\\
& G=S O(5) \rightarrow H=S O(4) \cong S U(2)_{L} \times S U(2)_{R}
\end{align*}
$$

This is understood by looking at how the adjoint representations of $S U(3)$ and $S O(5)$
decompose under the breaking pattern above. We have

$$
\begin{align*}
\mathbf{8} & \rightarrow \mathbf{3}_{0} \oplus \mathbf{1}_{0} \oplus \mathbf{2}_{1 / 2} \oplus \overline{\mathbf{2}}_{-1 / 2}  \tag{8.8.14}\\
\mathbf{1 0} & \rightarrow \mathbf{6} \oplus \mathbf{4}=(\mathbf{1}, \mathbf{3}) \oplus(\mathbf{3}, \mathbf{1}) \oplus(\mathbf{2}, \mathbf{2}) .
\end{align*}
$$

The subscript in the first line of eq. (8.8.14) refers to the $U(1)$ charge and in the second line we have reported the decomposition in terms of $S O(4)$ and $S U(2)_{L} \times S U(2)_{R}$ representations, respectively. The broken generators in the $S U(3)$ case transform as the last two terms in the first line of eq. (8.8.14), and the corresponding NG bosons form a complex doublet of $S U(2)$. The broken generators in the $S O(5)$ case transform as the last term in the second line of eq. (8.8.14), and the corresponding NG bosons form a 4 -plet of $S O(4)$ or a bidoublet of $S U(2)_{L} \times S U(2)_{R}$. Both NG bosons have the quantum numbers of the ordinary Higgs field. However, the $S U(3) \rightarrow S U(2)$ case does not include the $S U(2)_{c}$ custodial symmetry, since the only unbroken $S U(2)$ must be the $S U(2)_{L}$ group. We then focus on the $S O(5) \rightarrow S O(4)$ case in the following.

The $S O(5)$ generators are as in eq. (8.8.3), with the indices $a, b, i, j$ running now from 1 to 5 , rather than from 1 to 4 . The $S O(4)$ subgroup can be taken to be the one generated by the matrices (8.8.4), with the understanding that they are now $5 \times 5$ matrices with an additional row and column of zeros. The remaining $S O(5) / S O(4)$ broken generators are given by

$$
\begin{equation*}
t^{a}=-\frac{i}{\sqrt{2}} t^{a 5}, \quad a=1,2,3,4 \tag{8.8.15}
\end{equation*}
$$

The 4 NG bosons $h_{a}$ are encoded as usual in the matrix

$$
\begin{equation*}
\gamma=e^{i \sqrt{2} \frac{h_{a} t_{a}}{f_{H}}} \tag{8.8.16}
\end{equation*}
$$

where the $\sqrt{2}$ factor arises from our choice of normalizations of the $S O(5)$ generators. It is straightforward to verify that $G / H$ is a symmetric coset space, invariant under the automorphism (8.6.13), with $t_{i} \in S O(4)(i=1, \ldots, 6)$ and $t_{a} \in S O(5) / S O(4)(a=$ $1, \ldots, 4)$. The matrix $U=\gamma^{2}$ transforms homogeneously under $S O(5)$ transformations, as in eq. (8.6.16). For a generic gauging, the covariant derivative for $U$ is given in eq.(8.6.21). In our case, the gauging is all within $H$, so that $A_{\mu}^{a}$ all vanish and $A_{R}=A$. The covariant derivative (8.6.21) reduces to (now for canonically normalized gauge fields, including gauge coupling constant factors)

$$
\begin{equation*}
D_{\mu} U=\partial_{\mu} U+i\left(g W_{\mu}^{L}+g^{\prime} W_{\mu}^{R}\right) U-i U\left(g W_{\mu}^{L}+g^{\prime} W_{\mu}^{R}\right) \tag{8.8.17}
\end{equation*}
$$

with

$$
\begin{equation*}
W_{\mu}^{L}=\sum_{\alpha=1}^{3} t_{L}^{\alpha} W_{\mu}^{\alpha}, \quad W_{\mu}^{R}=t_{R}^{3} B_{\mu} \tag{8.8.18}
\end{equation*}
$$

The Higgs Lagrangian associated to this "composite Higgs" scenario is

$$
\begin{equation*}
\mathcal{L}_{H}^{C H}=\frac{f_{H}^{2}}{16} \operatorname{tr}\left(D^{\mu} U\right)^{\dagger} D_{\mu} U \tag{8.8.19}
\end{equation*}
$$

The NG nature of the Higgs forbids a potential for $H$, in the limit in which the $S O(5)$ global symmetry is exact and only spontaneously broken to $S O(4)$. But like $S U(3)_{V}$ in the QCD case is broken by $U(1)_{E M}$, the $S O(5)$ symmetry is explicitly broken by the $S U(2)_{L} \times U(1)_{Y}$ gauging ${ }^{9}$. This implies that a potential for the Higgs field, even if not included at tree-level, will be generated by radiative effects. We will not elaborate more on this idea, but if suffices to say here that when fermions are also included and coupled to the Higgs matrix field $U$, electroweak symmetry breaking can in fact be induced at the quantum level. When $h_{4} \neq 0$, a straightforward computation shows that the Lagrangian (8.8.19) gives rise to the following mass terms for the SM gauge fields:

$$
\begin{equation*}
m_{W}^{2}=\frac{1}{4} g^{2} f_{H}^{2} \sin ^{2}\left(\frac{\left\langle h_{4}\right\rangle}{f_{H}}\right), \quad m_{Z}^{2}=\frac{1}{4}\left(g^{2}+g^{\prime 2}\right) f_{H}^{2} \sin ^{2}\left(\frac{\left\langle h_{4}\right\rangle}{f_{H}}\right), \quad m_{\gamma}^{2}=0 \tag{8.8.20}
\end{equation*}
$$

As expected, thanks to the $S U(2)_{c}$ custodial symmetry, the mass ratio (8.8.7) is reproduced. We see that for $\left\langle h_{4}\right\rangle / f_{H} \rightarrow 0$ we can expand the sin factor and recover the usual SM formula for the $W$ 's and $Z$, where we identify

$$
\begin{equation*}
\left\langle h_{4}\right\rangle=v \equiv \sqrt{\frac{\sqrt{2}}{2 G_{F}}} \simeq 246 \mathrm{GeV} \tag{8.8.21}
\end{equation*}
$$

This is the limit in which we push to very high energies the $S O(5) \rightarrow S O(4)$ breaking pattern, but in so doing we recover the hierarchy problem. On the other hand, various phenomenological bounds constrain the more natural limit $\left\langle h_{4}\right\rangle \simeq f_{H}$, so that a little tuning is needed to achieve a mild separation between $\left\langle h_{4}\right\rangle$ and $f_{H}$. Notice that away from the SM limit $f_{H} \rightarrow \infty,\left\langle h_{4}\right\rangle$ does not coincide with $v$, as defined in eq. (8.8.21). From eq. (8.8.20), for finite $f_{H}$, we get

$$
\begin{equation*}
v=f_{H} \sin \left(\frac{\left\langle h_{4}\right\rangle}{f_{H}}\right) . \tag{8.8.22}
\end{equation*}
$$

### 8.9 Effective Field Theories for Broken Symmetries: Generale Case*

In this section we consider the general construction of the effective field theory of NG bosons, discussed for the first time by Callan, Coleman, Wess and Zumino [25]. This

[^58]slightly more elaborated formalism allows to consider arbitrary symmetry breaking patterns.

The analysis made in section 8.6 continues to be valid up to eq.(8.6.12) included. If the coset is not symmetric, it is not particularly useful to consider the matrix $U=\gamma^{2}$ and a more elaborate analysis is needed. It is useful in this case to study the transformation properties of the field combination $\gamma^{-1} \partial_{\mu} \gamma$. By recalling the formula

$$
\begin{equation*}
e^{X} Y e^{-X}=Y+[X, Y]+\frac{1}{2!}[X,[X, Y]]+\frac{1}{3!}[X,[X,[X, Y]]]+\ldots \tag{8.9.1}
\end{equation*}
$$

valid for arbitrary matrices $X$ and $Y$, we see that the combination $\gamma^{-1} \partial_{\mu} \gamma$ is a field defined in the algebra $\mathcal{G}$ of the group $G$. As such, it can be decomposed as

$$
\begin{equation*}
\gamma^{-1} \partial_{\mu} \gamma=i t_{a} D_{\mu}^{a}(x)+i t_{i} E_{\mu}^{i}(x) \tag{8.9.2}
\end{equation*}
$$

where necessarily $D_{\mu}^{a}(x)=D^{a b}(\xi(x)) \partial_{\mu} \xi_{b}(x), E_{\mu}^{i}(x)=E^{i a}(\xi(x)) \partial_{\mu} \xi_{a}(x)$, for some fields $D^{a b}$ and $E^{i a}$. Under a global transformation $g \in G$ we clearly have

$$
\begin{equation*}
(g \gamma)^{-1} \partial_{\mu}(g \gamma)=\gamma^{-1} \partial_{\mu} \gamma \tag{8.9.3}
\end{equation*}
$$

On the other hand $g \gamma=\gamma^{\prime} h$ and hence, after multiplication by $h^{-1}$ to the right, we get

$$
\begin{equation*}
\gamma^{-1 \prime} \partial_{\mu} \gamma^{\prime}=h\left(\gamma^{-1} \partial_{\mu} \gamma\right) h^{-1}-\left(\partial_{\mu} h\right) h^{-1} \tag{8.9.4}
\end{equation*}
$$

Using the decomposition (8.9.2) for $\gamma^{\prime}$ and identifying both sides, we get the transformations of $D_{\mu}^{a}$ and $E_{\mu}^{i}$ under global $G$ transformations:

$$
\begin{align*}
t_{a} D_{\mu}^{a}\left(\xi^{\prime}\right) & =h t_{a} h^{-1} D_{\mu}^{a}(\xi)=t_{a} R_{a b}(h) D_{\mu}^{b}(\xi) \\
t_{i} E_{\mu}^{i}\left(\xi^{\prime}\right) & =h t_{i} h^{-1} E_{\mu}^{i}(\xi)+i\left(\partial_{\mu} h\right) h^{-1} \tag{8.9.5}
\end{align*}
$$

Since $H$ is a subgroup of $G$, we have

$$
\begin{equation*}
h t_{i} h^{-1}=R_{i j}(h) t_{j}, \quad\left(\partial_{\mu} h\right) h^{-1}=i t_{i} R_{i a}(h) \partial_{\mu} \xi_{a}(x), \tag{8.9.6}
\end{equation*}
$$

and thus

$$
\begin{align*}
D_{\mu}^{a}\left(\xi^{\prime}\right) & =R_{a b}(h) D_{\mu}^{b}(\xi)  \tag{8.9.7}\\
E_{\mu}^{i}\left(\xi^{\prime}\right) & =R_{i j}(h) E_{\mu}^{j}(\xi)-R_{i a}(h) \partial_{\mu} \xi_{a}(x)
\end{align*}
$$

Under global $G$ transformations, the factors $D_{\mu}^{a}$ transform linearly, while the factors $E_{\mu}^{i}$ transform non-linearly like a gauge field. We can in fact use $E_{\mu}^{i}$ to build covariant derivatives for $\tilde{\phi}$. From eqs. (8.6.7) and (8.6.8), we get that under $G$ transformations

$$
\begin{equation*}
\partial_{\mu} \tilde{\phi}^{\prime}=\left(\partial_{\mu} h\right) \tilde{\phi}+h\left(\partial_{\mu} \tilde{\phi}\right)=h\left(h^{-1}\left(\partial_{\mu} h\right) \tilde{\phi}+\partial_{\mu} \tilde{\phi}\right) . \tag{8.9.8}
\end{equation*}
$$

Defining

$$
\begin{equation*}
D_{\mu} \tilde{\phi} \equiv\left(\partial_{\mu} \tilde{\phi}+i t_{i} E_{\mu}^{i}\right) \tilde{\phi} \tag{8.9.9}
\end{equation*}
$$

we have

$$
\begin{equation*}
D_{\mu} \tilde{\phi}^{\prime}=h D_{\mu} \tilde{\phi} \tag{8.9.10}
\end{equation*}
$$

The symmetries constrain the NG bosons to appear in the low-energy Lagrangian only through the combinations $D_{\mu}^{a}$ and $E_{\mu}^{i}$ in derivative interactions. We conclude that any Lagrangian invariant under $H$ transformations and constructed with $D_{\mu}^{a}, \tilde{\phi}$ and their covariant derivatives, will automatically be invariant under the whole group $G$. Invariant terms can also be constructed using the "field strength"

$$
\begin{equation*}
F_{\mu \nu}(E)=\partial_{\mu} E_{\nu}-\partial_{\nu} E_{\mu}+\left[E_{\mu}, E_{\nu}\right] \tag{8.9.11}
\end{equation*}
$$

and derivatives thereof, keeping in mind that not all invariants constructed in this way are independent of each other.

The leading order term encoding the NGB kinetic terms is

$$
\begin{equation*}
\mathcal{L} \supset \frac{1}{2} F_{a b}^{2} D_{\mu}^{a} D_{\mu}^{b}, \tag{8.9.12}
\end{equation*}
$$

where $D_{\mu}^{a}=\partial_{\mu} \xi^{a}+\ldots$. As we have seen in section 8.6 , the linearly realized symmetry $H$ constrains the coefficients $F_{a b}^{2}$.

The above formalism can easily be extended when a subgroup $H_{g}$ is gauged. The physical considerations are clearly the same as discussed in section 8.6. Let us pretend that the whole group $G$ is gauged and only at the end recover the original theory by setting to zero the gauge fields not in $H_{g}$. One can check that all our previous results, until eq. (8.9.1) included, are formally still valid, provided the obvious understanding that now the $G$ transformations are $x$-dependent. The $G$ invariant field combination to consider now, analogue of $\gamma^{-1} \partial_{\mu} \gamma$, is ${ }^{10}$

$$
\begin{equation*}
\gamma^{-1} D_{\mu} \gamma \equiv \gamma^{-1}\left(\partial_{\mu}-i t^{\alpha} A_{\mu}^{\alpha}\right) \gamma \tag{8.9.13}
\end{equation*}
$$

so that the inhomogeneous term in the gauge field transformation cancels the analogue term coming from $\partial_{\mu} g$. The quantity $\gamma^{-1} D_{\mu} \gamma$ is also defined in the algebra $\mathcal{G}$, so we have

$$
\begin{equation*}
\gamma^{-1} D_{\mu} \gamma \equiv i t_{a} \hat{D}_{\mu}^{a}(x)+i t_{i} \hat{E}_{\mu}^{i}(x), \tag{8.9.14}
\end{equation*}
$$

where $\hat{D}_{\mu}^{a}$ and $\hat{E}_{\mu}^{i}$ are the gauged versions of the fields $D_{\mu}^{a}$ and $E_{\mu}^{i}$ defined in eq. (8.9.2), which now include the gauge fields $A_{\mu}^{\alpha}$. Following the same steps as above, we immediately

[^59]get
\[

$$
\begin{equation*}
\left(\gamma^{-1} D_{\mu} \gamma\right)^{\prime}=h(x)\left(\gamma^{-1} D_{\mu} \gamma\right) h^{-1}(x)-\partial_{\mu} h(x) h^{-1}(x) \tag{8.9.15}
\end{equation*}
$$

\]

Correspondingly, the relations (8.9.5) are valid for $\hat{D}_{\mu}^{a}$ and $\hat{E}_{\mu}^{i}$ as well. As before, any Lagrangian invariant under $H$ transformations and constructed with $\hat{D}_{\mu}^{a}, \tilde{\phi}$ and their covariant derivatives, will automatically be invariant under the whole gauge group $G$. In order to also construct $H$ invariant operators involving the field strength $F_{\mu \nu}=\partial_{\mu} A_{\nu}-$ $\partial_{\nu} A_{\mu}+\left[A_{\mu}, A_{\nu}\right]$, that transforms linearly under $G$ as $F_{\mu \nu} \rightarrow g F_{\mu \nu} g^{-1}$, we can define NG boson-dependent field strengths

$$
\begin{equation*}
f_{\mu \nu}=\gamma^{-1} F_{\mu \nu} \gamma \tag{8.9.16}
\end{equation*}
$$

The latter transform as $f_{\mu \nu} \rightarrow h \gamma^{-1} g^{-1} g F_{\mu \nu} g^{-1} g \gamma h^{-1}=h f_{\mu \nu} h^{-1}$ under $G$ and can be used together with $\hat{D}_{\mu}^{a}, \tilde{\phi}$ and their associated operators constructed acting with covariant derivatives. We can now set to zero the "spurionic" gauge field components that do not belong to $H_{g}$. The leading order term (8.9.12) in the low-energy effective Lagrangian becomes

$$
\begin{equation*}
\mathcal{L} \supset \frac{1}{2} F_{a b}^{2} \hat{D}_{\mu}^{a} \hat{D}_{\mu}^{b} \tag{8.9.17}
\end{equation*}
$$

where $\hat{D}_{\mu}^{a}=\partial_{\mu} \xi^{a}-g_{\hat{\alpha} a} A_{\mu}^{\hat{\alpha}}$, with $g_{\hat{\alpha} \alpha}$ as defined in eq.(8.6.22). As expected, the NGB's along the gauged directions are eaten by the gauge fields and their mass is given by eq.(8.6.24).

## Chapter 9

## Anomalies

There are different ways of regularizing a QFT. The best choice of regulator is the one which keeps the maximum number of symmetries of the classical action unbroken. Cut-off regularization, for instance, breaks gauge invariance and that's why we prefer to work in the somewhat more exotic DR , where instead gauge invariance is always manifestly unbroken. It is also possible that there exists no regulator that preserves a given classical symmetry. In this case we say that the symmetry is anomalous, namely the quantum theory necessarily breaks it, independently of the choice of regulator.

Roughly speaking, anomalies can affect global or local symmetries. The latter case is particularly important, because local symmetries are needed to decouple unphysical fields. Indeed, anomalies in linearly realized local gauge symmetries lead to inconsistent theories. Theories with anomalous global symmetries are instead consistent, yet the effect of the anomaly can have important effects on the theories. We have already seen an example of anomaly. In classically scale invariant theories, a scale dependence is typically generated by quantum effects by means of a non-vanishing $\beta$-function. This kind of anomaly will be briefly discussed in section 9.7. We will however mostly focus on global anomalies associated to chiral currents and their related anomalies in local symmetries. Historically, the first anomaly, discovered by Adler, Bell, and Jackiw [26, 27], was associated to the non-conservation of the axial current in QCD. Among other things, the axial anomaly resolved a puzzle related to the $\pi^{0} \rightarrow 2 \gamma$ decay rate, predicted by effective Lagrangian considerations to be about three orders of magnitude smaller than the observed one.

In the next section we will first study the basic anomaly associated to a global $U(1)$ chiral transformation using Feynman diagrams and then consider generalization to nonabelian groups and to local symmetries.

### 9.1 The $U(1)_{A}$ Chiral Anomaly from One-Loop Graphs*

The Lagrangian for a free Dirac fermion $\psi$ is

$$
\begin{equation*}
\mathcal{L}=i \bar{\psi} \not \partial \psi-m \bar{\psi} \psi \tag{9.1.1}
\end{equation*}
$$

and is invariant under the vectorial $U(1)_{V}$ symmetry under which $\psi \rightarrow e^{i \alpha} \psi, \bar{\psi} \rightarrow \bar{\psi} e^{-i \alpha}$. The associated conserved $U(1)_{V}$ current is

$$
\begin{equation*}
J^{\mu}=\bar{\psi} \gamma^{\mu} \psi \tag{9.1.2}
\end{equation*}
$$

If $m=0$, as we assume from now on, the Lagrangian (9.1.1) is also invariant under the axial $U(1)_{A}$ symmetry under which

$$
\begin{equation*}
\psi \rightarrow e^{i \alpha \gamma_{5}} \psi, \quad \bar{\psi} \rightarrow \bar{\psi} e^{i \alpha \gamma_{5}} \tag{9.1.3}
\end{equation*}
$$

The associated conserved $U(1)_{A}$ current is

$$
\begin{equation*}
J_{5}^{\mu}=\bar{\psi} \gamma^{\mu} \gamma_{5} \psi \tag{9.1.4}
\end{equation*}
$$

While the classical $U(1)_{V}$ symmetry is quantum mechanically preserved and leads to Ward identities as explained in section 4.4.2, it turns out that there is no way to keep the $U(1)_{A}$ chiral symmetry at the quantum level. This is clear from a functional integral point of view. An invariant theory not only requires an invariant action, but also an invariant measure. And the functional measure is not invariant under the chiral transformation (9.1.3) (see section 9.4 below for a detailed analysis). The $U(1)_{A}$ current is quantum mechanically not conserved,

$$
\begin{equation*}
\left\langle\partial_{\mu} J_{5}^{\mu}\right\rangle=\mathcal{A}(x) \neq 0 \tag{9.1.5}
\end{equation*}
$$

and we say that there is a $U(1)_{A}$ chiral anomaly given by $\mathcal{A}(x)$. We can probe the anomaly by adding to the free fermion theory external spin one sources $K^{\mu}$ and $K_{5}^{\mu}$ that couple to $J^{\mu}$ and $J_{5}^{\mu}$, respectively:

$$
\begin{equation*}
Z\left[K^{\mu}, K_{5}^{\mu}\right]=\int \mathcal{D} \psi \mathcal{D} \bar{\psi} e^{i \int d^{4} x\left[\mathcal{L}+K_{\mu} J^{\mu}+K_{\mu 5} J_{5}^{\mu}\right]} \tag{9.1.6}
\end{equation*}
$$

Anomalies were originally discovered by evaluating three-point functions between external sources at one-loop level. They arise from loops of the internal fermion lines from current correlations that involve the chiral matrix $\gamma_{5}$. Despite these one-loop graphs are divergent and would require a choice of regularization, their divergence (i.e. the anomaly) is finite. Since we are only interested to the anomaly, we can proceed with the computation without the need of introducing a regulator.

We are then led to compute the three point functions between two vector currents $J^{\mu}$ and one axial current $J_{5}^{\mu}$ :

$$
\begin{equation*}
\Gamma_{\mu \nu \rho}(x, y, z)=\left.\frac{i \delta Z}{\delta K_{5}^{\mu}(x) \delta K^{\nu}(y) \delta K^{\rho}(z)}\right|_{K=K_{5}=0}=\left\langle J_{5}^{\mu}(x) J^{\nu}(y) J^{\rho}(z)\right\rangle \tag{9.1.7}
\end{equation*}
$$

which in momentum space reads as

$$
\begin{equation*}
\Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right) \equiv\left\langle J_{\mu, 5}\left(-k_{1}-k_{2}\right) J_{\nu}\left(k_{1}\right) J_{\rho}\left(k_{2}\right)\right\rangle \tag{9.1.8}
\end{equation*}
$$

where, with some abuse of notation, we denote the Fourier transform with the same symbol of the corresponding term. The would-be conservation of the two currents, $\partial_{\mu} J^{\mu}=\partial_{\mu} J_{5}^{\mu}=$ 0 , would imply at the quantum level the relations

$$
\begin{equation*}
\left(k_{1}+k_{2}\right)^{\mu} \Gamma_{\mu \nu \rho}=k_{1}^{\nu} \Gamma_{\mu \nu \rho}=k_{2}^{\rho} \Gamma_{\mu \nu \rho}=0 \tag{9.1.9}
\end{equation*}
$$

No contact terms arise in eqs.(9.1.9), because the axial and vector currents are classically invariant under the $U(1)_{V} \times U(1)_{A}$ transformations. As we will see, due to the anomaly, it turns out to be impossible to impose all three conditions (9.1.9) simultaneously.

The two diagrams in fig. 9.1 contribute at one-loop level to $\Gamma_{\mu \nu \rho}$. They give

$$
\begin{array}{r}
\Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)=-\int \frac{d^{4} p}{(2 \pi)^{4}} \operatorname{tr}\left(\gamma_{\mu} \gamma_{5} S(p+a+k) \gamma_{\rho} S\left(p+a+k_{1}\right) \gamma_{\nu} S(p+a)+\right. \\
\left.\gamma_{\mu} \gamma_{5} S(p+b+k) \gamma_{\nu} S\left(p+b+k_{2}\right) \gamma_{\rho} S(p+b)\right) \tag{9.1.10}
\end{array}
$$

where $k=k_{1}+k_{2}$ and $S(p)=\not p / p^{2}$ is the free fermion propagator for a massless fermion. In eq. (9.1.10) $a$ and $b$ are arbitrary constant vectors, whose relevance will be clear in the following. Let us first compute the divergence of the axial current, $k^{\mu} \Gamma_{\mu \nu \rho}$. It is convenient to write

$$
\begin{equation*}
S(p+a) \not k S(p+a+k)=S(p+a)(p p+\not p+\not k b-\not p-\not p) S(p+a+k)=S(p+a)-S(p+a+k), \tag{9.1.11}
\end{equation*}
$$

so that
$k^{\mu} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)=\int \frac{d^{4} p}{(2 \pi)^{4}}\left[f_{\rho \nu}\left(p, b+k_{2}, b+k\right)-f_{\rho \nu}\left(p, a, a+k_{1}\right)+\left(\rho \leftrightarrow \nu, a \leftrightarrow b, k_{1} \leftrightarrow k_{2}\right)\right]$
where

$$
\begin{equation*}
f_{\rho \nu}(p, a, b) \equiv \operatorname{tr}\left(S(p+a) \gamma_{\rho} S(p+b) \gamma_{\nu} \gamma_{5}\right)=4 i \epsilon_{\gamma \rho \omega \nu} \frac{p_{\omega}(a-b)_{\gamma}+a_{\gamma} b_{\omega}}{(p+a)^{2}(p+b)^{2}} \tag{9.1.13}
\end{equation*}
$$

Noticing that $f_{\rho \nu}(p+c, a, b)=f_{\rho \nu}(p, a+c, b+c)$, the second term in eq. (9.1.12) becomes


Figure 9.1: One-loop graphs contributing to the anomaly. All external momenta are incoming. The wavy and dashed lines represent the external vector and axial sources, respectively.
identical to the first one by shifting the virtual momentum $p \rightarrow p+b-a+k_{2}$, giving naively $k^{\mu} \Gamma_{\mu \nu \rho}=0$. However, these expressions are divergent and momentum shift is not allowed. Indeed, for a generic function $f(x)$, we have

$$
\begin{equation*}
\int_{-\infty}^{\infty} d x[f(x+a)-f(x)]=\int_{-\infty}^{\infty} d x\left[a f^{\prime}(x)+\ldots\right]=a[f(\infty)-f(-\infty)]+\ldots \tag{9.1.14}
\end{equation*}
$$

Equation (9.1.14) vanishes if $f( \pm \infty)=f^{\prime}( \pm \infty)=\ldots=0$, conditions automatically satisfied for convergent integrals, but not for divergent ones. In the case at hand, the analogue of $f$ is $p^{3} f_{\rho \nu}$ and $f_{\rho \nu} \sim 1 / p^{3}$, so $f( \pm \infty) \neq 0$. The integral is governed by the asymptotic behaviour of $f$ and thus it is easy to evaluate. Rotating the virtual momentum $p$ to euclidean values and applying Stokes theorem, we have

$$
\begin{align*}
& i \int \frac{d^{4} p_{E}}{(2 \pi)^{4}}\left[f_{\rho \nu}\left(p_{E}+k, a, b\right)-f_{\rho \nu}\left(p_{E}, a, b\right)\right]=\frac{i}{(2 \pi)^{4}} \lim _{p_{E} \rightarrow \infty} \int k^{\mu} p_{E, \mu} p_{E}^{2} f_{\rho \nu}\left(p_{E}\right) d \Omega_{4} \\
= & \frac{i}{(2 \pi)^{4}} \int k_{\mu} p_{E}^{\mu} p_{E}^{2} \frac{4 i \epsilon_{\gamma \rho \omega \nu} k^{\omega}(a-b)^{\gamma} p_{E}^{\omega}}{\left(p_{E}^{2}\right)^{2}} d \Omega_{4}=\frac{\Omega_{4}}{(2 \pi)^{4}} \epsilon_{\gamma \rho \omega \nu} k^{\omega}(a-b)^{\gamma}, \tag{9.1.15}
\end{align*}
$$

where $\Omega_{4}=2 \pi^{2}$ is the volume of the unit four-sphere and in the last relation we have used $S O(4)$ invariance to write

$$
\begin{equation*}
\int p_{E}^{\mu} p_{E}^{\omega} f\left(p_{E}^{2}\right) d \Omega_{4}=-\frac{\delta^{\mu \omega}}{4} \Omega_{4} p_{E}^{2} f\left(p_{E}^{2}\right) \tag{9.1.16}
\end{equation*}
$$

Using eq. (9.1.15), the derivative of the axial current is easily computed:

$$
\begin{align*}
k^{\mu} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right) & =\frac{1}{8 \pi^{2}} \epsilon_{\gamma \rho \omega \nu}\left[\left(k_{2}+b-a\right)^{\omega}\left(-k_{1}\right)^{\gamma}+\left(k_{1}+a-b\right)^{\omega} k_{2}^{\gamma}\right] \\
& =\frac{1}{8 \pi^{2}} \epsilon_{\gamma \rho \omega \nu}\left(k_{1}-k_{2}+\Delta\right)^{\omega}\left(k_{1}+k_{2}\right)^{\gamma}, \tag{9.1.17}
\end{align*}
$$

with $\Delta \equiv a-b$. It is now clear why we have introduced the otherwise redundant parameters $a$ and $b$. In a convergent expression, they would trivially be reabsorbed in a shift of the virtual momentum $p$, but in the case at hand the final expression turns out to depend on their difference $\Delta$. Being $\Delta$ arbitrary, one could choose $\Delta=k_{2}-k_{1}$, so that the "anomalous" term (9.1.17) would cancel. However, care has to be paid on the divergence of the two vector currents $J_{\nu}$ and $J_{\rho}$. Proceeding exactly as before, we can compute

$$
\begin{align*}
k_{1}^{\nu} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right) & =\frac{1}{8 \pi^{2}} \epsilon_{\gamma \mu \omega \rho}\left[\left(k_{1}+k_{2}\right)^{\gamma}(b-a)^{\omega}-\left(b-a-k_{1}\right)^{\omega} k_{2}^{\gamma}\right] \\
& =\frac{1}{8 \pi^{2}} \epsilon_{\gamma \mu \omega \rho} k_{1}^{\omega}\left(\Delta+k_{2}\right)^{\gamma},  \tag{9.1.18}\\
k_{2}^{\rho} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right) & =\frac{1}{8 \pi^{2}} \epsilon_{\gamma \mu \omega \nu}\left[\left(k_{1}+k_{2}\right)^{\gamma}(a-b)^{\omega}-\left(a-b-k_{2}\right)^{\omega} k_{1}^{\gamma}\right] \\
& =\frac{1}{8 \pi^{2}} \epsilon_{\gamma \mu \omega \nu} k_{2}^{\gamma}\left(\Delta-k_{1}\right)^{\omega} . \tag{9.1.19}
\end{align*}
$$

The choice $\Delta=k_{2}-k_{1}$ would then lead to the non-conservation of the two vector currents $J_{\nu}$ and $J_{\rho}$. Since three-point functions between 3 vector currents do not lead to any anomaly, we insist in having vanishing divergence for the vector currents. This uniquely fixes $\Delta=k_{1}-k_{2}$, opposite to the choice leading to the conservation of the axial current. We can shift the anomalous term from one current to another, but what is important is that there is no choice of $\Delta$ for which all three currents are conserved. Plugging $\Delta=k_{1}-k_{2}$ in eq. (9.1.17) gives

$$
\begin{equation*}
k^{\mu} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)=\frac{1}{2 \pi^{2}} \epsilon_{\gamma \rho \omega \nu} k_{1}^{\omega} k_{2}^{\gamma} . \tag{9.1.20}
\end{equation*}
$$

This is the final form of the chiral anomaly in momentum space. When $K_{\mu}$ are actual external sources, the anomaly term in eq.(9.1.20) corresponds to a local counterterm for these external sources and effectively do not lead to any non-conservation of currents in the quantum theory. However, we can equally couple the free fermion theory to dynamical gauge fields, gauging the $U(1)_{V}$ symmetry, in which case the vector currents $K_{\mu}$ are nothing else than the $U(1)_{V}$ gauge fields $A_{\mu}$ (modulo a coupling constant factor, depending on the normalization chosen). The gauging of $U(1)_{V}$ does not break the $U(1)_{A}$ symmetry, that still remains at the classical level. In this case the wavy lines depicted in fig.9.1 are actual photons and it is not difficult to reconstruct the local anomaly functional $\mathcal{A}(x)$ entering in eq.(9.1.5). This should be such that, when taking two functional derivatives with respect to the gauge fields $A_{\mu}$ gives, in momentum space, the term in the right hand side of eq.(9.1.20). For canonically normalized gauge fields we get, in configuration space,

$$
\begin{equation*}
\left\langle\partial_{\mu} J_{5}^{\mu}\right\rangle=-\frac{g^{2}}{16 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu} F_{\rho \sigma} \tag{9.1.21}
\end{equation*}
$$

This is the standard form of the chiral anomaly. We see that when gauge fields are coupled to the fermion field the chiral current is not conserved.

It is not difficult to generalize the above analysis to non-abelian global symmetries. The Lagrangian for $N$ massless free Dirac fields $\psi_{i}(i=1, \ldots, N)$ has an $S U(N)_{V} \times$ $S U(N)_{A} \times U(1)_{A} \times U(1)_{V}$ global symmetry. The vector and axial currents read

$$
\begin{equation*}
J_{\mu}^{\alpha}=\bar{\psi} \gamma_{\mu} T^{\alpha} \psi, \quad J_{\mu 5}^{\alpha}=\bar{\psi} \gamma_{\mu} \gamma_{5} T^{\alpha} \psi \tag{9.1.22}
\end{equation*}
$$

where $T^{\alpha}$ the $U(N)$ generators, including the $U(1)$ factor. We then evaluate the 3-point function at one-loop level between two vector and 1 axial currents:

$$
\begin{equation*}
\Gamma_{\mu \nu \rho}^{\alpha \beta \gamma}\left(k_{1}, k_{2}\right) \equiv\left\langle J_{\mu 5}^{\alpha}\left(-k_{1}-k_{2}\right) J_{\nu}^{\beta}\left(k_{1}\right) J_{\rho}^{\gamma}\left(k_{2}\right)\right\rangle . \tag{9.1.23}
\end{equation*}
$$

Similarly to eq. (9.1.10), we have

$$
\begin{align*}
\Gamma_{\mu \nu \rho}^{\alpha \beta \gamma}\left(k_{1}, k_{2}\right)=- & \int \frac{d^{4} p}{(2 \pi)^{4}}\left[\operatorname { t r } ( T ^ { \alpha } T ^ { \gamma } T ^ { \beta } ) \operatorname { t r } \left(\gamma_{\mu} \gamma_{5} S(p+a+k) \gamma_{\rho} S\left(p+a+k_{1}\right) \gamma_{\nu} S(p+a)\right.\right. \\
& \left.\left.+\operatorname{tr}\left(T^{\alpha} T^{\beta} T^{\gamma}\right) \gamma_{\mu} \gamma_{5} S(p+b+k) \gamma_{\nu} S\left(p+b+k_{2}\right) \gamma_{\rho} S(p+b)\right)\right] . \tag{9.1.24}
\end{align*}
$$

Since the currents transform under a symmetry transformation, their classical conservation turns in non-trivial WT identities, see eq.(4.4.13):

$$
\begin{equation*}
i \partial_{x_{1}}^{\mu}\left\langle J_{\mu 5}^{\alpha}\left(x_{1}\right) J_{\nu}^{\beta}\left(x_{2}\right) J_{\rho}^{\gamma}\left(x_{3}\right)\right\rangle=\delta\left(x_{1}-x_{2}\right)\left\langle\delta_{\alpha} J_{\nu}^{\beta}\left(x_{2}\right) J_{\rho}^{\gamma}\left(x_{3}\right)\right\rangle+\delta\left(x_{1}-x_{3}\right)\left\langle J_{\nu}^{\beta}\left(x_{2}\right) \delta_{\alpha} J_{\rho}^{\gamma}\left(x_{3}\right)\right\rangle, \tag{9.1.25}
\end{equation*}
$$

where, under an infinitesimal chiral transformation, $\delta_{\alpha} J_{\mu}^{\beta}=C_{\beta \alpha \gamma} J_{\mu 5}^{\gamma}$. By parity invariance the two-point function between a vector and an axial current vanishes, the contact terms in eq.(9.1.25) then vanish and in momentum space we simply have ${ }^{1}$

$$
\begin{equation*}
\left(k_{1}+k_{2}\right)^{\mu} \Gamma_{\mu \nu \rho}^{\alpha \beta \gamma}=0 . \tag{9.1.26}
\end{equation*}
$$

It is convenient to write

$$
\begin{equation*}
\operatorname{tr}\left(T^{\alpha} T^{\beta} T^{\gamma}\right)=\frac{1}{2} \operatorname{tr}\left(\left\{T^{\alpha}, T^{\beta}\right\} T^{\gamma}\right)+\frac{1}{2} \operatorname{tr}\left(\left[T^{\alpha}, T^{\beta}\right] T^{\gamma}\right)=D^{\alpha \beta \gamma}+\frac{i}{2} C^{\alpha \beta \gamma} \tag{9.1.27}
\end{equation*}
$$

where in eq.(9.1.27) we have introduced the factor

$$
\begin{equation*}
D^{\alpha \beta \gamma} \equiv \frac{1}{2} \operatorname{tr}\left(\left\{T^{\alpha}, T^{\beta}\right\} T^{\gamma}\right) \tag{9.1.28}
\end{equation*}
$$

symmetric under permutation of its indices. It can be shown that the terms proportional to $C^{\alpha \beta \gamma}$ do not give rise to anomalies, which are found in the terms proportional to $D^{\alpha \beta \gamma}$.

[^60]These can be manipulated exactly as before and eventually lead to the same form of the anomaly:

$$
\begin{equation*}
\left.\Gamma_{\mu \nu \rho}^{\alpha \beta \gamma}\left(k_{1}, k_{2}\right)\right|_{\text {ano. }}=\left.D^{\alpha \beta \gamma} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)\right|_{\text {ano. }} . \tag{9.1.29}
\end{equation*}
$$

As before, a non-conservation of currents in the quantum theory arises only when we gauge some of the global symmetries. With non-abelian symmetries we have to pay attention in the gauging procedure, because it can lead to an explicit violation of some of the global symmetries. For example, gauging the whole $U(N)_{V}$ factor would break $S U(N)_{A}$, because the gauge coupling interaction is not $S U(N)_{A}$ invariant. The $U(1)_{A}$ symmetry is however classical preserved, so we might consider the case in which the axial current is given by $U(1)_{A}$, while the vector currents are gauged and correspond to gauge fields. In this case $T^{\alpha}$ can be taken equal to unity and we get

$$
\begin{equation*}
\left\langle\partial_{\mu} J_{5}^{\mu}\right\rangle=-\frac{g^{2}}{4 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} \partial_{\mu} A_{\nu}^{\beta} \partial_{\rho} A_{\sigma}^{\gamma} \operatorname{tr}\left(T^{\beta} T^{\gamma}\right) \tag{9.1.30}
\end{equation*}
$$

where $D_{\alpha \beta \gamma} \rightarrow \operatorname{tr}\left(T^{\beta} T^{\gamma}\right)$. In the non-abelian case, additional one-loop (square and pentagon) diagrams contribute to the anomaly and should be considered. When they are summed to eq. (9.1.30), the whole non-abelian form of the field strength is reconstructed and we get the final result

$$
\begin{equation*}
\left\langle\partial_{\mu} J_{5}^{\mu}\right\rangle=-\frac{g^{2}}{16 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu}^{\beta} F_{\rho \sigma}^{\gamma} \operatorname{tr}\left(T^{\beta} T^{\gamma}\right) . \tag{9.1.31}
\end{equation*}
$$

Eq.(9.1.31) is the generalization of eq.(9.1.21) for non-abelian gauge fields and, as expected, boils down to eq.(9.1.21) in the abelian case with $T^{\beta}=T^{\gamma}=1$.

In any given regularization, where all amplitudes are made finite and the shift in the virtual momentum allowed, the anomaly would appear differently. For instance, in DR, the anomaly arises from subtleties related to the definition of $\gamma_{5}$ in $d \neq 4$ dimensions (see, e.g., ref. [1] for a computation of the anomaly in DR). In Pauli-Villars regularization, where a heavy ( PV ) fermion is added, the anomaly is related to the explicit breaking of the axial symmetry given by the PV fermion mass term. There is no regulator that preserves at the same time the vector and axial symmetries and as a result an anomaly always appears.

### 9.2 Gauge Anomalies^

As we have already mentioned, anomalies can also affect local symmetries, in which case we refer to them as "gauge anomalies". The latter can arise in chiral, parity non-invariant, gauge theories, where left and right-moving fermions transform in representations of the
gauge group that are not complex conjugate with each other. The simplest gauge anomaly can be described by starting from the Lagrangian for a free chiral fermion $\psi_{L}$ :

$$
\begin{equation*}
\mathcal{L}=i \bar{\psi}_{L} \not \partial \psi_{L} \tag{9.2.1}
\end{equation*}
$$

invariant under $U(1)_{L}$ transformations $\psi_{L} \rightarrow e^{i \alpha} \psi_{L}, \bar{\psi}_{L} \rightarrow \bar{\psi}_{L} e^{-i \alpha}$, with $U(1)_{L}$ current $J^{\mu}=\bar{\psi}_{L} \gamma^{\mu} \psi_{L}=\bar{\psi} \gamma^{\mu} P_{L} \psi$, with $P_{L}=\left(1+\gamma_{5}\right) / 2$ the chirality projector. The 3 -point function at one-loop level between three currents is

$$
\begin{align*}
\Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)=- & \int \frac{d^{4} p}{(2 \pi)^{4}}\left[\operatorname { t r } \left(\gamma_{\mu} P_{L} S(p+a+k) \gamma_{\rho} P_{L} S\left(p+a+k_{1}\right) \gamma_{\nu} P_{L} S(p+a)\right.\right. \\
& \left.\left.+\gamma_{\mu} P_{L} S(p+b+k) \gamma_{\nu} P_{L} S\left(p+b+k_{2}\right) \gamma_{\rho} P_{L} S(p+b)\right)\right] \tag{9.2.2}
\end{align*}
$$

where $\Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right) \equiv\left\langle J_{\mu}\left(-k_{1}-k_{2}\right) J_{\nu}\left(k_{1}\right) J_{\rho}\left(k_{2}\right)\right\rangle$. Since $P_{L}^{2}=P_{L}$, eq.(9.2.2) turns into a sum of a non-anomalous 3 -point correlators with no $\gamma_{5}$ factors and an anomalous one with one insertion of $\gamma_{5}$. The last term, modulo a factor $1 / 2$ coming from $P_{L}$, equals eq.(9.1.10). We then get

$$
\begin{equation*}
k^{\mu} \Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)=\frac{1}{2} \frac{1}{8 \pi^{2}} \epsilon_{\gamma \rho \omega \nu}\left(k_{1}-k_{2}+\Delta\right)^{\omega}\left(k_{1}+k_{2}\right)^{\gamma}, \tag{9.2.3}
\end{equation*}
$$

and similarly $1 / 2$ of the divergences (9.1.18) and (9.1.19) for the $\mathcal{J}^{\nu}$ and $\mathcal{J}^{\rho}$ currents. Gauging the chiral theory (9.2.1) amounts to couple the above three currents to a $U(1)$ gauge field. Since the theory is chiral, there are no vector and axial currents, but only left currents to couple to the gauge field. As a result, there is no choice of the factor $\Delta$ that allows us to conserve all the three vector currents at the same time. Because of permutation symmetry of the three identical external currents, we should choose $\Delta$ so that the anomaly is symmetric under the exchange of any of the three currents, as required by Bose symmetry. This is achieved by taking

$$
\begin{equation*}
\Delta=-\frac{1}{3}\left(k_{1}-k_{2}\right) \tag{9.2.4}
\end{equation*}
$$

which gives an anomaly one-third smaller than the anomaly (9.1.21):

$$
\begin{equation*}
\left\langle\partial_{\mu} J^{\mu}\right\rangle=-\frac{1}{2} \frac{g^{2}}{48 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu} F_{\rho \sigma} \tag{9.2.5}
\end{equation*}
$$

This is the simplest form of a gauge anomaly. Such an anomaly is deadly for a theory. The WI derived in section 4.4.3 no longer hold. In particular, the violation of eq.(4.4.44) does not allow the decoupling of unphysical states. In scattering amplitudes the factor $\Gamma_{\mu \nu \rho}\left(k_{1}, k_{2}\right)$ is multiplied either by polarizations of the external photons or by internal
photon propagators when the triangular graph is embedded within a more complicated Feynman graph. For external photons the computation of the divergence of one of the gauge currents is equivalent to replace the photon polarization vector by its momentum, i.e. to consider longitudinal photons. A non-vanishing result implies that these states do not decouple. If the line is internal, the longitudinal photon runs into the loop, yet we have a problem, because unitarity (the optical theorem) would imply their non-decoupling and a violation of unitarity (at any scale). In other words, minimally coupling a $U(1)$ gauge field to a chiral fermion is quantum mechanically inconsistent.

Similar considerations can be generalized to multiple fermions and to non-abelian gauge theories. We will not enter into the details of non-abelian gauge anomalies. It is sufficient here to say that the key coefficient governing the anomaly is the factor $D^{\alpha \beta \gamma}$ defined in eq.(9.1.28). When we have several fermions transforming in different representations $r$ of a gauge group $G$, the trace in $D^{\alpha \beta \gamma}$ should be taken over all fermions, each in the corresponding representation $r$. Crucially, left-handed and right-handed fermions contribute with an opposite sign to $D^{\alpha \beta \gamma}$, due to the chiral nature of the gauge anomaly. A gauge theory is non-anomalous if and only if in total $D^{\alpha \beta \gamma}=0$. The explicit form of the non-abelian anomaly can be significantly constrained by demanding consistency with the structure of the gauge group (see section 9.5 for some other detail) and cannot be written in a covariant way in terms of field strengths only.

So far, we have only considered massless fermions. Fermion mass terms explicitly break the axial symmetry (9.1.3), but it can easily be shown that they do not change the form of the axial anomaly (9.1.30). In fact, left-handed and right-handed fermions contribute the same to the axial anomaly, because the sign change due to the different chirality of the fermions compensate their opposite charge with respect to the axial symmetry. This explains why we got a non-vanishing axial anomaly starting from vector-like Dirac fermions. The situation is different for gauge anomalies, that can receive a non-vanishing contribution only from massless fermions. This is proved as follows. It is convenient to consider all fermions as left-handed by defining, for each right-handed component $\psi_{R}$ its left-handed counterpart

$$
\begin{equation*}
\psi_{L}^{c}=C \psi_{R}^{\star} \tag{9.2.6}
\end{equation*}
$$

$C$ being the matrix of charge conjugation in spinor space. If $\psi_{R}$ transforms under a representation $r_{R}$ of a gauge group $G, \psi_{L}^{c}$ will transform as $r_{R}^{\star}$. In the basis of left-handed two-components spinors $\chi_{i}$, a generic mass term reads

$$
\begin{equation*}
\mathcal{L}_{m}=\chi_{i}^{t} \sigma^{2} \hat{M}_{i j} \chi_{j}+h . c . \tag{9.2.7}
\end{equation*}
$$

where $i$ run over all the fermions in the theory and $\hat{M}$ is a symmetric mass matrix. Let
us focus on a subset of terms in eq. (9.2.7) coupling two left-handed fermion multiplets $\chi_{1}$ and $\chi_{2}$ in irreducible representations $r_{1}$ and $r_{2}$ of the group $G$, with $\operatorname{dim} r_{1}=\operatorname{dim} r_{2}$ :

$$
\begin{equation*}
\mathcal{L}_{m} \supset \chi_{1}^{t} \sigma^{2} M \chi_{2}+h . c . \tag{9.2.8}
\end{equation*}
$$

where $M$ is a non-singular mass matrix. The term (9.2.8) is gauge-invariant if

$$
\begin{equation*}
-T_{1}^{t} M=M T_{2} \tag{9.2.9}
\end{equation*}
$$

where $T_{1}$ and $T_{2}$ are the generators in the representations $r_{1}$ and $r_{2}$. Eq. (9.2.9) implies that $-T_{1}^{t}$ and $T_{2}$ are related by a similarity transformation, since $-T_{1}^{t}=M T_{2} M^{-1}$. Then

$$
\begin{equation*}
D_{\alpha \beta \gamma}^{1}=\frac{1}{2} \operatorname{tr}\left\{T_{1 \alpha}, T_{1 \beta}\right\} T_{1 \gamma}=(-1)^{3} \frac{1}{2} \operatorname{tr}\left\{T_{2 \alpha}^{t}, T_{2 \beta}^{t}\right\} T_{2 \gamma}^{t}=-D_{\alpha \beta \gamma}^{2} \tag{9.2.10}
\end{equation*}
$$

The contribution to the gauge anomaly given by $\chi_{1}$ is then exactly cancelled by that of $\chi_{2}$. Let us clarify this result with a couple of simple examples. Consider a $U(1)$ gauge theory with a Dirac fermion with charge $q$. In terms of left-handed fields, the Dirac fermion consists of one left-handed fermion $\chi_{1}$ with charge $q$ and its conjugate $\chi_{2}$ with charge $-q$ and admits the mass term $m \chi_{1} \sigma^{2} \chi_{2}+h . c$. The total gauge anomaly is proportional to

$$
\begin{equation*}
+q^{3}+(-q)^{3}=0 \tag{9.2.11}
\end{equation*}
$$

Similarly, Dirac fermions in any representation $T_{r}$ of a gauge group consist of one lefthanded fermion multiplet in the representation $T_{r}$ and its conjugate in the complex conjugate representation $T_{\bar{r}}=-T_{r}^{t}$ and thus do not lead to anomalies. In manifestly parityinvariant theories the absence of anomalies is obvious since all currents are manifestly vector-like (i.e no $\gamma_{5}$ appears in a 4 -component Dirac notation). Non-trivial anomalies can only arise in non-parity-invariant theories, namely in so called chiral gauge theories, where a fermion and its complex conjugate transform in representations of the gauge group that are not complex conjugate of each other. The absence of anomalies in chiral gauge theories requires a non-trivial cancellation between different fermion multiplets. The most important example of a theory of this sort is the SM, where gauge anomalies cancel between quarks and leptons, as we will see in the next section.

We can now come back to the fate of the $U(1)_{A}$ symmetry in QCD, mentioned at the end of section 8.7. QCD with $n_{f}$ massless quarks has actually an $S U\left(n_{f}\right)_{V} \times S U\left(n_{f}\right)_{A} \times$ $U(1)_{V} \times U(1)_{A}$ global symmetry, spontaneously broken to $S U\left(n_{f}\right)_{V} \times U(1)_{V}$ by the quark condensate. Let us compute the possible anomalies that we can have by looking at the $D_{\alpha \beta \gamma}$ coefficients associated to the various groups. Since quarks are vector-like, the gauge $S U(3)^{3}$ anomalies all cancel. Anomalies with one $S U(N)$ non-abelian current of any kind
also manifestly vanish, because $\operatorname{tr} T=0$. The only possible non-vanishing anomalies are of the form $S U(3)^{2} \times U(1)_{V}$ and $S U(3)^{2} \times U(1)_{A}$. In terms of left-handed quarks, a quark and its conjugate have opposite charges with respect to $U(1)_{V}$ and equal charges with respect to $U(1)_{A}$, thus the $S U(3)^{2} \times U(1)_{V}$ anomaly vanishes while $S U(3)^{2} \times U(1)_{A}$ does not. This is nothing else than the anomaly considered before in eq.(9.1.31), with $F_{\mu \nu}^{\alpha}$ identified with the gluon field strengths. In QCD, then, the $U(1)_{A}$ symmetry, in addition of being spontaneously broken by the condensate, is also explicitly broken by the anomaly. The latter breaking is large when the theory becomes strongly coupled (due to instantons) and is responsible for the absence of a light $\eta^{\prime}$ in the QCD scalar meson spectrum.

Summarizing, we can have three qualitatively different scenarios:

- An anomaly involving three global currents $\rightarrow$ The theory is consistent, no real effect, the global symmetry is preserved.
- An anomaly involving one global and two local currents $\rightarrow$ The theory is consistent, the global symmetry is broken.
- An anomaly involving three local currents $\rightarrow$ The theory is inconsistent.


### 9.3 A Relevant Example: Cancellation of Gauge Anomalies in the $\mathrm{SM}^{\star}$

The SM gauge group is $G_{S M}=S U(3) \times S U(2) \times U(1)$. Its fermion content, in terms of lefthanded fields, is composed by three copies (generations) of the following representations:

$$
\begin{equation*}
(\mathbf{3}, \mathbf{2})_{\frac{1}{6}}+(\overline{\mathbf{3}}, \mathbf{1})_{-\frac{2}{3}}+(\overline{\mathbf{3}}, \mathbf{1})_{\frac{1}{3}}+(\mathbf{1}, \mathbf{2})_{-\frac{1}{2}}+(\mathbf{1}, \mathbf{1})_{1} \tag{9.3.1}
\end{equation*}
$$

corresponding to the quark doublet, up quark singlet, down quark singlet, lepton doublet and charged lepton singlet, respectively. In principle there can be ten possible kinds of gauge anomalies, associated to all possible combinations of $S U(3) S U(2)$ and $U(1)$ currents in the trangular graph. Five of them, where a non-abelian group factor $(S U(3)$ or $S U(2))$ appears only once, $S U(3)^{2} \times S U(2), S U(3) \times S U(2)^{2}, S U(3) \times S U(2) \times U(1), S U(3) \times U(1)^{2}$ and $S U(2) \times U(1)^{2}$ are trivially vanishing, since for $S U(n)$ groups the generators are traceless: $\operatorname{tr} T=0$. The $S U(2)^{3}$ anomaly is also manifestly vanishing because for $S U(2)$ the symmetric factor $D^{\alpha \beta \gamma}$ vanishes. In the case at hand, with doublets only, this is easily seen: $D^{\alpha \beta \gamma}=1 / 2 \operatorname{tr}\left\{t^{\alpha}, t^{\beta}\right\} t^{\gamma}=\delta^{\alpha \beta} \operatorname{tr} t^{\gamma}=0$. In the general case, $D^{\alpha \beta \gamma}$ vanishes because all $S U(2)$ representations are equivalent to their complex conjugates, namely there exists a matrix $A$ such that $T^{t}=T^{\star}=-A T A^{-1}$. Using this relation, one immediately sees that $D^{\alpha \beta \gamma}=0$ for any representation.

The remaining combinations $S U(3)^{3}, S U(3)^{2} \times U(1), S U(2)^{2} \times U(1)$ and $U(1)^{3}$ have to be checked. Let us then compute the values of the symmetric coefficients $D^{\alpha \beta \gamma}$ in each of the above 4 cases and show that they always vanish. In order to distinguish the different coefficients, we denote by a subscript $c, w$ and $Y$ the $S U(3), S U(2)$ and $U(1)$ factors, respectively. It is enough to consider a single generation of quarks and doublets, because the cancellation occurs generation per generation. Let us start with the $S U(3)^{3}$ anomaly. Only quarks contribute to it. We get

$$
\begin{equation*}
D_{c c c}^{\alpha \beta \gamma}=2 D_{3}^{\alpha \beta \gamma}+D_{\overline{3}}^{\alpha \beta \gamma}+D_{\overline{3}}^{\alpha \beta \gamma}=2 D_{\mathbf{3}}^{\alpha \beta \gamma}-D_{3}^{\alpha \beta \gamma}-D_{3}^{\alpha \beta \gamma}=0 \tag{9.3.2}
\end{equation*}
$$

using that $D_{\overline{3}}^{\alpha \beta \gamma}=-D_{3}^{\alpha \beta \gamma}$. For $S U(3)^{2} \times U(1)$ we have

$$
\begin{equation*}
D_{c c Y}^{\alpha \beta}=2 \operatorname{tr}_{\mathbf{3}} t^{\alpha} t^{\beta} \times \frac{1}{6}+\operatorname{tr}_{\overline{3}} t^{\alpha} t^{\beta} \times\left(-\frac{2}{3}\right)+\operatorname{tr}_{\mathbf{3}} t^{\alpha} t^{\beta} \times \frac{1}{3}=\operatorname{tr}_{\mathbf{3}} t^{\alpha} t^{\beta}\left(\frac{1}{3}-\frac{2}{3}+\frac{1}{3}\right)=0 \tag{9.3.3}
\end{equation*}
$$

with $\operatorname{tr}_{\mathbf{3}} t^{\alpha} t^{\beta}=\operatorname{tr}_{\mathbf{3}} t^{\alpha} t^{\beta}$. For $S U(2)^{2} \times U(1)$, only doublets contribute. We get

$$
\begin{equation*}
D_{w w Y}^{\alpha \beta}=3 \operatorname{tr}_{2} t^{\alpha} t^{\beta} \times \frac{1}{6}+\operatorname{tr}_{2} t^{\alpha} t^{\beta} \times\left(-\frac{1}{2}\right)=\operatorname{tr}_{2} t^{\alpha} t^{\beta}\left(\frac{1}{2}-\frac{1}{2}\right)=0 \tag{9.3.4}
\end{equation*}
$$

For $U(1)^{3}$ anomalies all quarks and leptons contribute and one gets

$$
\begin{equation*}
D_{Y Y Y}=3 \times 2 \times\left(\frac{1}{6}\right)^{3}+3 \times\left(-\frac{2}{3}\right)^{3}+3 \times\left(\frac{1}{3}\right)^{3}+2 \times\left(-\frac{1}{2}\right)^{3}+(1)^{3}=0 \tag{9.3.5}
\end{equation*}
$$

There is actually a fifth non-trivial anomaly to check, that arises when we couple the SM to gravity. It is an anomaly involving a $U(1)$ current and two energy-momentum tensors, and it is called a mixed $U(1)$-gravitational anomaly. ${ }^{2}$ This anomaly is proportional to $\sum_{n} q_{n}$, where $n$ runs over all fermions with charges $q_{n}$. In the SM, the mixed $U(1)$-gravitational anomaly is proportional to

$$
\begin{equation*}
3 \times 2 \times \frac{1}{6}+3 \times\left(-\frac{2}{3}\right)+3 \times \frac{1}{3}+2 \times\left(-\frac{1}{2}\right)+1=0 . \tag{9.3.6}
\end{equation*}
$$

Notice how the fermion charges nicely combine to give a vanishing result for all the anomalies, in particular the $U(1)^{3}$ and the mixed $U(1)$-gravitational ones.

The SM is then a chiral, anomaly-free gauge theory. More precisely, we have shown that the SM is anomaly-free in the unbroken phase where all the gauge group is linearly realized (i.e. no Higgs mechanism is at work). Since the anomaly does not depend on the Higgs field, our results automatically imply that the SM is anomaly-free in the broken phase as well. In particular, the unbroken gauge group $S U(3) \times U(1)_{E M}$ is manifestly anomaly-free being all currents vector-like.

[^61]
### 9.4 Path Integral Derivation of the Chiral Anomaly**

In the path-integral formulation of field theory, anomalies arise from the transformation of the measure used to define the fermion path integral. We will consider in what follows path-integral in euclidean space-time.

Let $\psi_{A}(x)$ be a massless Dirac fermion defined on $R^{4}$, in an arbitrary representation $\mathcal{R}$ of a gauge group $G(A=1, \ldots, \operatorname{dim} \mathcal{R})$. The minimal coupling of the fermion to the gauge fields is described by the Lagrangian

$$
\begin{equation*}
\mathcal{L}=\bar{\psi}(x)_{A} i \gamma^{\mu} D_{\mu B}^{A} \psi^{B}(x) . \tag{9.4.1}
\end{equation*}
$$

The covariant derivative is given by

$$
\begin{equation*}
D_{\mu B}^{A}=\partial_{\mu} \delta_{B}^{A}+A_{\mu}^{\alpha} T_{\alpha B}^{A}, \tag{9.4.2}
\end{equation*}
$$

in terms of the gauge connection $A_{\mu}^{\alpha}$ and the anti-Hermitian generators $T_{\alpha B}^{A}$ of the group $G$ in the representation $\mathcal{R}(\alpha=1, \ldots, \operatorname{dim} G)$, with $\gamma^{\mu}$ satisfying the anticommutation relations $\left\{\gamma_{\mu}, \gamma_{\nu}\right\}=2 \delta_{\mu \nu}(\mu, \nu=1,2,3,4)$.

The classical Lagrangian (9.4.1) is invariant under the global chiral transformation

$$
\begin{equation*}
\psi \rightarrow e^{i \alpha \gamma_{5}} \psi \tag{9.4.3}
\end{equation*}
$$

where $\gamma_{5}=\prod_{\mu=1}^{4} \gamma_{\mu}$, normalized so that $\gamma_{5}^{2}=I$, and $\alpha$ is a constant parameter. The chiral current

$$
\begin{equation*}
J_{5}^{\mu}=\bar{\psi}_{A} \gamma^{\mu} \gamma_{5} \psi^{A} \tag{9.4.4}
\end{equation*}
$$

is classically conserved. At the quantum level, however, this conservation law can be violated and turns into an anomalous WT identity. To derive it, we consider the quantum effective action $\Gamma$ defined by

$$
\begin{equation*}
e^{-\Gamma(A)}=\int \mathcal{D} \psi \mathcal{D} \bar{\psi} e^{-\int d^{4} x \mathcal{L}} \tag{9.4.5}
\end{equation*}
$$

and study its behavior under an infinitesimal chiral transformation of the fermions, with a space-time-dependent parameter $\alpha(x)$, given by ${ }^{3}$

$$
\begin{equation*}
\delta_{\alpha} \psi=i \alpha \gamma_{5} \psi, \quad \delta_{\alpha} \bar{\psi}=i \alpha \bar{\psi} \gamma_{5} \tag{9.4.6}
\end{equation*}
$$

Since the external gauge fields $A$ are inert, the transformation (9.4.6) represents a redefinition of dummy integration variables, and should not affect the effective action: $\delta_{\alpha} \Gamma=0$.

[^62]This statement carries however a non-trivial piece of information, since neither the classical action nor the integration measure is invariant under (9.4.6). The variation of the classical action under (9.4.6) is non-vanishing only for non-constant $\alpha$, and has the form $\delta_{\alpha} \int \mathcal{L}=\int J_{5}^{\mu} \partial_{\mu} \alpha$. The variation of the measure is instead always non-vanishing, because the transformation (9.4.6) leads to a non-trivial Jacobian factor, which has the form $\delta_{\alpha}[\mathcal{D} \psi \mathcal{D} \bar{\psi}]=\exp \left\{i \int \alpha \mathcal{A}\right\} \mathcal{D} \psi \mathcal{D} \bar{\psi}$, as we will see below. In total, the effective action therefore transforms as

$$
\begin{equation*}
\delta_{\alpha} \Gamma=\int d^{4} x \alpha(x)\left[i \mathcal{A}(x)+\left\langle\partial_{\mu} J_{5}^{\mu}(x)\right\rangle\right] . \tag{9.4.7}
\end{equation*}
$$

The condition $\delta_{\alpha} \Gamma=0$ then implies the anomalous WT identity:

$$
\begin{equation*}
\left\langle\partial_{\mu} J_{5}^{\mu}\right\rangle=-i \mathcal{A} \tag{9.4.8}
\end{equation*}
$$

In order to compute the anomaly $\mathcal{A}$, we need to define the integration measure more precisely. This is best done by considering the eigenfunctions of the Dirac operator $\langle\boldsymbol{} \boldsymbol{Z} \equiv$ $i \gamma^{\mu} D_{\mu}$. Since the latter is Hermitian, the set of its eigenfunctions $\psi_{k}(x)$ with eigenvalues $\lambda_{k}$, defined by $\backslash \supset \nmid \psi_{n}=\lambda_{n} \psi_{n}$, form an orthonormal and complete basis of spinor modes:

$$
\begin{equation*}
\int d^{4} x \psi_{k}^{\dagger a}(x) \psi_{l}^{a}(x)=\delta_{k, l}, \quad \sum_{k} \psi_{k}^{\dagger a}(x) \psi_{k}^{b}(y)=\delta^{a b} \delta^{(4)}(x-y), \tag{9.4.9}
\end{equation*}
$$

where we have made explicit the spinor indices $a$ and $b$. The fermion fields $\psi$ and $\bar{\psi}$, which are independent from each other in Euclidean space, can be decomposed as

$$
\begin{equation*}
\psi=\sum_{k} a_{k} \psi_{k}, \quad \bar{\psi}=\sum_{k} \bar{b}_{k} \psi_{k}^{\dagger} \tag{9.4.10}
\end{equation*}
$$

so that the measure becomes

$$
\begin{equation*}
\mathcal{D} \psi \mathcal{D} \bar{\psi}=\prod_{k, l} d a_{k} d \bar{b}_{l} \tag{9.4.11}
\end{equation*}
$$

Under the chiral transformation (9.4.6), we have

$$
\begin{equation*}
\delta_{\alpha} a_{k}=i \int d^{4} x \sum_{l} \psi_{k}^{\dagger} \alpha \gamma_{5} \psi_{l} a_{l}, \quad \delta_{\alpha} \bar{b}_{k}=i \int d^{4} x \sum_{l} \bar{b}_{l} \psi_{l}^{\dagger} \alpha \gamma_{5} \psi_{k} \tag{9.4.12}
\end{equation*}
$$

and the measure (9.4.11) transforms as

$$
\begin{align*}
\delta_{\alpha}[\mathcal{D} \psi \mathcal{D} \bar{\psi}] & =\mathcal{D} \psi \mathcal{D} \bar{\psi} \operatorname{det}\left(\delta_{k l}+i \int d^{4} x \psi_{k}^{\dagger} \alpha \gamma_{5} \psi_{l}\right)^{-2} \\
& =\mathcal{D} \psi \mathcal{D} \bar{\psi} \exp \left\{-2 i \sum_{k} \int d^{4} x \psi_{k}^{\dagger} \alpha \gamma_{5} \psi_{k}+\mathcal{O}\left(\alpha^{2}\right)\right\} . \tag{9.4.13}
\end{align*}
$$

For simplicity we can take $\alpha$ to be constant. This formal expression is ill-defined as it stands, since it decomposes into a vanishing trace over spinor indices $\left(\operatorname{tr} \gamma_{5}=0\right)$ times an infinite sum over the modes $\left(\sum_{k} 1=\infty\right)$. A convenient way of regularizing it is to introduce a gauge-invariant Gaussian cut-off. The anomaly $\mathcal{A}$ can then be defined as

$$
\begin{equation*}
\mathcal{A}=-2 \lim _{\beta \rightarrow 0} \sum_{k} \psi_{k}^{\dagger} \gamma_{5} e^{-\beta(\not \supset)^{2}} \psi_{k} . \tag{9.4.14}
\end{equation*}
$$

Using the completeness relation in eq. (9.4.9), we can write

$$
\begin{equation*}
\mathcal{A}=-2 \lim _{\beta \rightarrow 0} \lim _{y \rightarrow x} \operatorname{Tr}\left[\gamma_{5} e^{-\beta(\not \supset \not)^{2}}\right] \delta^{(4)}(x-y)=-2 \lim _{\beta \rightarrow 0} \int \frac{d^{4} k}{(2 \pi)^{4}} \operatorname{Tr}\left[\gamma_{5} e^{-\beta\left(k \mid y+\lfloor\not \supset)^{2}\right.}\right], \tag{9.4.15}
\end{equation*}
$$

where the trace is over the spinor and the gauge indices. By using the commutator properties of the $\gamma$ matrices, we can rewrite

$$
\begin{equation*}
(\not k t+i D)^{2}=(k+i D)^{2}-\frac{1}{4} F^{\mu \nu}\left[\gamma_{\mu}, \gamma_{\nu}\right] . \tag{9.4.16}
\end{equation*}
$$

Rescaling the momentum $k \rightarrow k / \sqrt{\beta}$, we get

$$
\begin{equation*}
\mathcal{A}=-2 \lim _{\beta \rightarrow 0} \frac{1}{\beta^{2}} \int \frac{d^{4} k}{(2 \pi)^{4}} \operatorname{Tr} \gamma_{5} e^{-\left(k^{2}+2 i \sqrt{\beta} D \cdot k-\beta D^{2}-\beta F\right)}, \tag{9.4.17}
\end{equation*}
$$

where $F \equiv F^{\mu \nu}\left[\gamma_{\mu}, \gamma_{\nu}\right] / 4$. The trace over spinor indices is vanishing unless at least two factors of $F$ (i.e. $4 \gamma^{\prime}$ s) are included in the trace. In this way, we get two powers of $\beta$ that compensate for the overall factor $1 / \beta^{2}$ in eq. (9.4.17). Hence, in the limit $\beta \rightarrow 0$, we can safely neglect the terms proportional to $D \cdot k$ and $D^{2}$ in the exponential. In this way, we finally get

$$
\begin{equation*}
\mathcal{A}=-\frac{1}{16 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu}^{\alpha} F_{\rho \sigma}^{\beta} \operatorname{tr}\left(T_{\alpha} T_{\beta}\right) \tag{9.4.18}
\end{equation*}
$$

In Minkowski space, with $\partial_{4}=-i \partial_{0}$, and in terms of hermitian (rather than antihermitian) generators and canonically normalized gauge fields $(A T \rightarrow-i g A T)$ the nonconservation of the axial current takes the form

$$
\begin{equation*}
\left\langle\partial_{\mu} J_{5}^{\mu}\right\rangle=-\frac{g^{2}}{16 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu}^{\alpha} F_{\rho \sigma}^{\beta} \operatorname{tr}\left(T_{\alpha} T_{\beta}\right) \tag{9.4.19}
\end{equation*}
$$

Anomalies in the chiral transformation (9.4.3) occur in any number of even spacetime dimensions (in odd-dimensional space-times, chirality is not defined, since there is no analogue of the matrix $\gamma_{5}$ ). The derivation of the anomaly reviewed here, due to Fujikawa, is easily generalized to any number of dimensions. In $2 n$ dimensions, we define the chiral matrix $\gamma_{2 n+1}$ as $\gamma_{2 n+1}=i^{n} \prod_{\mu=1}^{2 n} \gamma_{\mu}$ so that $\gamma_{2 n+1}^{2}=I$ for any $n$. All the steps are essentially identical to the 4 d case and we get the generalization of eq. (9.4.17):

$$
\begin{equation*}
\mathcal{A}=2 \lim _{\beta \rightarrow 0} \frac{1}{\beta^{n}} \int \frac{d^{2 n} k}{(2 \pi)^{2 n}} \operatorname{Tr} \gamma_{2 n+1} e^{-\left(k^{2}+2 i \sqrt{\beta} D \cdot k-\beta D^{2}-i \beta F\right)} . \tag{9.4.20}
\end{equation*}
$$

In $2 n$ dimensions, spinors have $2^{n}$ components, so we get for the anomaly

$$
\begin{equation*}
\mathcal{A}=-\frac{2}{(4 \pi)^{n} n!} \epsilon^{\mu_{1} \ldots \mu_{2 n}} F_{\mu_{1} \mu_{2}}^{\alpha_{1}} \ldots F_{\mu_{2 n-1} \mu_{2 n}}^{\alpha_{n}} \operatorname{tr}\left(T_{\alpha_{1}} \ldots T_{\alpha_{n}}\right) \tag{9.4.21}
\end{equation*}
$$

The anomaly (9.4.21) can more elegantly be rewritten in a compact form in terms of differential two-forms. Let us introduce auxiliary anticommuting variables $\phi^{1}, \ldots, \phi^{2 n}$ and define

$$
\begin{equation*}
F=\frac{1}{2} F_{\mu \nu} \phi^{\mu} \wedge \phi^{\nu} \tag{9.4.22}
\end{equation*}
$$

Then eq. (9.4.21) becomes simply

$$
\begin{equation*}
\mathcal{A}=-2 \int d^{2 n} \phi \operatorname{tr} e^{F /(2 \pi)}, \tag{9.4.23}
\end{equation*}
$$

where the integration over the auxiliary fermion variables $\phi^{\mu}$ automatically selects the correct number of field strengths $F$. In evaluating the integrated form of the anomaly, $\int d^{2 n} x \mathcal{A}$, we can replace the auxiliary fermion variables $\phi$ by the differentials $d x^{\mu}$, so that we have

$$
\begin{equation*}
\int d^{2 n} x \mathcal{A}=-2 \int d^{2 n} x \operatorname{ch}(F), \tag{9.4.24}
\end{equation*}
$$

where

$$
\begin{equation*}
\operatorname{ch}(F)=\operatorname{tr} e^{F /(2 \pi)} \tag{9.4.25}
\end{equation*}
$$

is the so-called Chern character of the gauge connection. There is a deep connection between anomalies and certain mathematical results that will not be discussed here.

### 9.5 The Wess-Zumino Consistency Conditions**

In this section we will consider some more formal developments about anomalies. It is convenient in this context to consider the effective action $\Gamma(A)$ arising upon integration of the fermions:

$$
\begin{equation*}
e^{-\Gamma(A)}=\int \mathcal{D} \psi \mathcal{D} \bar{\psi} e^{-S(\psi, \bar{\psi}, A)} \tag{9.5.1}
\end{equation*}
$$

In presence of an anomaly, $\Gamma(A)$ is not gauge invariant. Indeed, under an infinitesimal gauge transformation $A \rightarrow A+D \epsilon_{1}$

$$
\begin{equation*}
\delta_{\epsilon_{1}} \Gamma(A)=\int d^{4} x \frac{\delta \Gamma(A)}{\delta A_{\mu}^{\alpha}} D_{\mu} \epsilon_{1}^{\alpha}=-\int d^{4} x D_{\mu} J_{\alpha}^{\mu} \epsilon_{1}^{\alpha} \equiv-\int d^{4} x \mathcal{A}_{\alpha} \epsilon_{1}^{\alpha}=-\int d^{4} x \epsilon_{1}^{\alpha} \mathcal{F}_{\alpha} \Gamma(A) \tag{9.5.2}
\end{equation*}
$$

where the third equality defines the anomaly term $\mathcal{A}_{\alpha}$ and we have defined the functional operator

$$
\begin{equation*}
\mathcal{F}_{\alpha}=\partial_{\mu} \frac{\delta}{\delta A_{\mu}^{\alpha}}+C_{\alpha \beta \gamma} A_{\mu}^{\beta} \frac{\delta}{\delta A_{\mu}^{\gamma}} . \tag{9.5.3}
\end{equation*}
$$

Under a further infinitesimal transformations, we have

$$
\begin{equation*}
\delta_{\epsilon_{2}} \delta_{\epsilon_{1}} \Gamma(A)=\int d^{4} x \int d^{4} y \epsilon_{2}^{\beta}(y) \epsilon_{1}^{\alpha}(x) \mathcal{F}_{\beta}(y) \mathcal{F}_{\alpha}(x) \Gamma(A) \tag{9.5.4}
\end{equation*}
$$

Due to the group structure, performing the commutators of the two infinitesimal transformations parametrized by $\epsilon_{1}$ and $\epsilon_{2}$ should be equivalent to perform a single infinitesimal transformation with parameter $\left[\epsilon_{1}, \epsilon_{2}\right]$, where $\epsilon_{1,2}=\epsilon_{1,2}^{\alpha} T_{\alpha}$. In other words, we should have

$$
\begin{equation*}
\left[\delta_{\epsilon_{1}}, \delta_{\epsilon_{2}}\right] \Gamma(A)=\delta_{\left[\epsilon_{1}, \epsilon_{2}\right]} \Gamma(A), \tag{9.5.5}
\end{equation*}
$$

identity that can easily be verified to hold in general. The relation (9.5.5) implies a non-trivial condition for the anomaly, known as the Wess-Zumino consistency conditions:

$$
\begin{equation*}
\mathcal{F}_{\alpha}(x) \mathcal{A}_{\beta}(y)-\mathcal{F}_{\beta}(y) \mathcal{A}_{\alpha}(x)=-C_{\alpha \beta \gamma} \mathcal{A}_{\gamma}(x) \delta(x-y) \tag{9.5.6}
\end{equation*}
$$

These conditions can be more conveniently expressed in terms of BRST transformations by defining

$$
\begin{equation*}
G(\omega, A) \equiv \int d^{4} x \omega_{\alpha}(x) \mathcal{A}_{\alpha}(x) \tag{9.5.7}
\end{equation*}
$$

Equations (9.5.6) imply that $G$ is BRST-invariant, namely

$$
\begin{equation*}
s G(\omega, A)=0 \tag{9.5.8}
\end{equation*}
$$

This is easily shown by recalling the BRST transformations (6.3.3) of $\omega$ and $A$, according to which

$$
\begin{align*}
s G(\omega, A)= & \int d^{4} x\left(-\frac{1}{2} C_{\alpha \beta \gamma} \omega_{\beta}(x) \omega_{\gamma}(x) \mathcal{A}_{\alpha}(x)\right. \\
& -\omega_{\alpha}(x) \int d^{4} y \frac{\delta \mathcal{A}_{\alpha}(x)}{\delta A_{\mu}^{\beta}(y)}\left(\partial_{\mu} \omega_{\beta}(y)+C_{\beta \gamma \delta} A_{\mu}^{\gamma}(y) \omega_{\delta}(y)\right)  \tag{9.5.9}\\
= & \int d^{4} x d^{4} y\left(-\frac{1}{2} \omega_{\alpha}(x) \omega_{\beta}(y)\right)\left[\delta(x-y) C_{\alpha \beta \gamma} \mathcal{A}_{\gamma}(x)+\mathcal{F}_{\alpha}(x) \mathcal{A}_{\beta}(y)-\mathcal{F}_{\beta}(x) \mathcal{A}_{\alpha}(y)\right],
\end{align*}
$$

which vanishes if eq. (9.5.6) is satisfied.
We have extensively seen in the previous sections that there is some arbitrariness in computing anomalies, related to the fact that we can shift the latter from one current to another. From an effective action point of view, this shift corresponds to the possibility of adding local, non-gauge invariant, counter-terms to $\Gamma(A)$ that change $\delta_{\epsilon} \Gamma(A)$ and hence the anomaly. The impossibility of keeping all currents conserved corresponds to the im-
possibility of finding a local ${ }^{4}$ counter-term such that $\delta_{\epsilon} \Gamma(A)=0$. We can now be more precise exploiting eq. (9.5.8). If there existed a local functional of the gauge fields $F(A)$, such that $G(\omega, A)=s F(A)$, namely suppose that $G(\omega, A)$ would be BRST-exact, then the anomaly would be cancelled by adding to $\Gamma(A)$ the counter-term $-F(A)$. Anomalies then form an equivalence class. Two anomalies related by the addition of a local functional $F(A)$ are equivalent and belong to the same cohomology of the BRST operator $s$. Thanks to the Wess-Zumino consistency conditions, it is possible to reconstruct the whole form of the non-abelian anomaly by only knowing the terms quadratic in the gauge fields. The procedure does not uniquely fix the anomaly since, as we have just said, the latter can be changed by adding local counter-terms to the action.

We finally mention about the existence of an elegant formalism, known as the StoraZumino descent equations, that allows, in any number of even dimensions, to get anomaly functionals $\mathcal{A}_{\alpha}$ that automatically satisfy the conditions (9.5.6). The descent equations make also manifest the close relationship between chiral anomalies in $2 n+2$ dimensions and gauge anomalies in $2 n$ dimensions.

## 9.6 't Hooft Anomaly Matching and the Wess-Zumino-Witten Term*

Asymptotically free gauge theories are strongly coupled in the IR and can give rise to confinement of its fermion constituents, like quarks in QCD. At low energies the propagating degrees of freedom are bound states of the elementary high energy (UV) states, such as mesons and hadrons in QCD. What is the fate at low energies of possible global anomalies coming from the elementary constituents at high energy? t'Hooft has answered to this question by arguing that anomalies must arise in the low energy effective theory as well. For concreteness, let us consider a gauge theory coupled to fermions with local symmetry $G_{s}$ and global symmetry $G_{f}$ and denote schematically by $D_{a b c}$ the associated anomaly coefficients, where $a, b, c=s, f$. We assume that the theory in the IR becomes strongly coupled. For the theory to be consistent $D_{s s s}=0$ so that no gauge anomalies arise. We also have $D_{f s s}=0$, which means that we keep in $G_{f}$ only the quantum global symmetries of the theory, not the classical ones. For instance, in QCD we should not include $U(1)_{A}$ in $G_{f}$. We can however have $D_{f f f} \neq 0$ since, as we have seen, no real effect occurs for anomalies involving purely global currents. Although these anomalies do not lead to any

[^63]\[

$$
\begin{equation*}
-F(A)=\frac{g^{2}}{96 \pi^{2}} \frac{1}{\square} \partial^{\mu} A_{\mu} \epsilon^{\alpha \beta \gamma \delta} F_{\alpha \beta} F_{\gamma \delta} \tag{9.5.10}
\end{equation*}
$$

\]

effect, 't Hooft has shown that they should appear in the IR description of the theory and might then be useful to put constraints on the IR theory. Such anomalies are sometimes called 't Hooft anomalies.

The 't Hooft anomaly matching condition states that if in the UV theory $D_{f f f} \neq 0$ then the same anomaly should appear in the IR theory. While the UV anomaly is induced by the elementary UV fermion fields, the IR one should be given by the appearance of massless fermion bound states with suitable quantum numbers to have the same $D_{f f f} \neq 0$. If $G_{f}$ is spontaneously broken (like chiral symmetries in QCD), the NG bosons can replace the role of the massless fermion bound states that are no longer required. 't Hooft's argument is very simple. Let us assume of (weakly) gauging the group $G_{f}$ (or a subgroup of it) of the UV theory. In this way the innocuous global anomalies turn into deadly gauge anomalies making the theory inconsistent. We might cancel the gauge anomaly $D_{f f f}$ by adding suitable additional massless "spectator" fermions that give a contribution $D_{f f f}^{\text {spect. }}=-D_{f f f}$ to the anomaly. Crucially, these spectators can be taken neutral under $G_{s}$ and charged only under the weakly gauged symmetries $G_{f}$. At low energies, when the strong gauge group $G_{s}$ confines, the spectrum of the theory will include IR bound states plus the "spectator" fermions that, being neutral under $G_{s}$, are unaffected by the condensation of $G_{s}$. The UV theory with the spectators is, by construction, consistent and it has to remain so for all values of the gauge coupling constant. Hence, at low energies, there must be an anomaly contribution equal to $-D_{f f f}^{\mathrm{spect}}=D_{f f f}$, canceling that of the fermion spectators. The argument is valid for an arbitrarily weak gauging and thus has to also apply in the original theory where $G_{f}$ is a global symmetry.

Given the quantum numbers of the elementary constituents, the possible quantum numbers of the massless fermion bound states can be argued. It might happen that there is no choice of quantum numbers for the bound states to reproduce the anomaly of the constituents. In this case, the 't Hooft anomaly matching conditions can be used to prove that our original assumption of unbroken global symmetries is necessarily violated, and some global symmetries are spontaneously broken. Notably, one can show in this way that in QCD the $S U(3)_{V} \times S U(3)_{A} \times U(1)_{V}$ global symmetry must be spontaneously broken, see e.g. section 22.5 of ref.[2]. ${ }^{5}$

As we mentioned, when $G_{f}$ is spontaneously broken the anomaly of the UV fermions can be reproduced by the effective action of the (pseudo) Goldstone bosons. The latter situation is actually realized in Nature in QCD, in which case the weak gauging used in 't Hooft argument can be identified with the electromagnetic interactions. Consider

[^64]for simplicity $n_{f}=2$ (up and down quarks only). The quantum global symmetry group is $G_{f}=S U(2)_{V} \times S U(2)_{A} \times U(1)_{V}$. This is explicitly broken by the electromagnetic interactions to $H_{f}=U(1)_{V} \times U(1)_{E M} \times \widetilde{U}(1)_{A}$, where $\widetilde{U}(1)_{A} \subset S U_{A}(2)$ is the abelian subgroup generated by $t^{3}$. Consider now the $U(1)_{E M}^{2} \times \widetilde{U}(1)_{A}$ axial anomaly. We get
\[

$$
\begin{align*}
\partial_{\mu} J_{A}^{\mu} & =-\frac{N_{c}}{16 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu} F_{\rho \sigma}\left[\left(\frac{2 e}{3}\right)^{2} \times 1+\left(\frac{-e}{3}\right)^{2} \times(-1)\right] \\
& =-\frac{N_{c} e^{2}}{48 \pi^{2}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu} F_{\rho \sigma} \tag{9.6.1}
\end{align*}
$$
\]

Anomaly matching requires that a similar anomaly should show up in the IR. Since no massless fermion bound states occur Nature, this should appear in the low-energy effective chiral Lagrangian $\mathcal{L}_{\pi}$ describing the dynamics of the $\pi$ mesons interacting with photons. Under the $\widetilde{U}(1)_{A}$ chiral transformation above, $\delta_{\epsilon} \mathcal{L}_{\pi}$ should then not vanish, but rather reproduce the axial anomaly (9.6.1). Considering that $\delta_{\epsilon} \pi^{0}=\epsilon f_{\pi}, \mathcal{L}_{\pi}$ should include the coupling

$$
\begin{equation*}
\mathcal{L}_{\pi} \supset-\frac{N_{c} e^{2}}{48 \pi^{2} f_{\pi}} \epsilon^{\mu \nu \rho \sigma} F_{\mu \nu} F_{\rho \sigma} \pi^{0} \tag{9.6.2}
\end{equation*}
$$

The axial anomaly (9.6.1) has allowed to resolve the puzzle of the $\pi^{0} \rightarrow 2 \gamma$ decay. In absence of any anomaly, the term (9.6.2) would still appear in the chiral Lagrangian $\mathcal{L}_{\pi}$ but with a much more suppressed coupling. On the contrary, anomaly considerations uniquely fix its coefficient and it turns out that the experimental rate $\Gamma\left(\pi^{0} \rightarrow 2 \gamma\right)$ is successfully reproduced with $N_{c}=3$. We have seen in section 8.7 that the chiral Lagrangian $\mathcal{L}_{\pi}$ should be described in terms of the matrix of fields $U=\exp \left(2 i \pi^{a} t^{a} / f_{\pi}\right)$, rather than by the $\pi$ 's mesons themselves. The anomalous term (9.6.2) should then be rewritten in terms of the $U$ 's. This rewriting is not totally straightforward and will not be done here. The ending result goes under the name of the gauged version of the Wess-Zumino-Witten term. The latter includes many other couplings, including the term (9.6.2).

### 9.7 Anomalous Breaking of Scale Invariance ${ }^{\star}$

As we have mentioned at the beginning of this chapter, in most QFT's the symmetry under scale transformations, if present at tree level, is broken by quantum effects due to the energy dependence of the couplings. ${ }^{6}$ Let us formalize a bit better this observation.

We know that the energy-momentum tensor is the conserved tensor associated to the translational symmetries. In general, under an infinitesimal symmetry transformation of the action, the Lagrangian density is not necessarily invariant, but can change by a total

[^65]derivative: $\mathcal{L} \rightarrow \mathcal{L}+\epsilon \partial_{\mu} \mathcal{J}^{\mu}$, with $\mathcal{J}^{\mu}$ a given tensor. ${ }^{7}$ The associated conserved Noether current in this case is given by
\[

$$
\begin{equation*}
J^{\mu}=\frac{\delta \mathcal{L}}{\delta \partial_{\mu} \phi} \delta \phi-\mathcal{J}^{\mu} \tag{9.7.1}
\end{equation*}
$$

\]

Under an infinitesimal translation $x^{\mu} \rightarrow x^{\mu}+\epsilon^{\mu}$, the Lagrangian, being a scalar, transforms as $\mathcal{L} \rightarrow \mathcal{L}+\epsilon^{\mu} \partial_{\mu} \mathcal{L}$. The Noether energy-momentum tensor reads

$$
\begin{equation*}
T^{\mu \nu}=\frac{\delta \mathcal{L}}{\delta \partial_{\mu} \phi} \partial^{\nu} \phi-\eta^{\mu \nu} \mathcal{L} \tag{9.7.2}
\end{equation*}
$$

and is conserved: $\partial_{\mu} T^{\mu \nu}=0$. In general, the Noether energy-momentum tensor $T^{\mu \nu}$ is not symmetric, $T^{\mu \nu} \neq T^{\nu \mu}$. A new, symmetric, energy-momentum tensor $\theta^{\mu \nu}$ can be defined by adding a trivially conserved term to $T^{\mu \nu}$ :

$$
\begin{equation*}
\theta^{\mu \nu}=T^{\mu \nu}+\partial_{\rho} M^{\mu \nu \rho} \tag{9.7.3}
\end{equation*}
$$

where $M^{\mu \nu \rho}$ is a tensor antisymmetric in $\mu$ and $\rho$, so that $\partial_{\mu} \partial_{\rho} M^{\mu \nu \rho}$ automatically vanishes. The symmetric energy-momentum tensor $\theta^{\mu \nu}$ can directly be derived by coupling the QFT to gravity and noticing that $\theta^{\mu \nu}$ minimally couples to gravity. We have

$$
\begin{equation*}
\theta^{\mu \nu}=\left.\frac{2}{\sqrt{-g}} \frac{\delta S}{\delta g_{\mu \nu}}\right|_{g_{\mu \nu}=\eta_{\mu \nu}} \tag{9.7.4}
\end{equation*}
$$

where $g_{\mu \nu}$ is the metric and $g=\operatorname{det} g_{\mu \nu}$.
Let us see this how this works for the massless $\phi^{4}$ theory. In this case we have

$$
\begin{equation*}
T^{\mu \nu}=\theta^{\mu \nu}=\partial^{\mu} \phi \partial^{\nu} \phi-\eta^{\mu \nu} \mathcal{L} \tag{9.7.5}
\end{equation*}
$$

Coupling to gravity we have

$$
\begin{equation*}
S=\int d^{4} x \sqrt{-g}\left(\frac{1}{2} g_{\mu \nu} \partial^{\mu} \phi \partial^{\nu} \phi-\frac{\lambda}{4!} \phi^{4}\right) \tag{9.7.6}
\end{equation*}
$$

Noticing that

$$
\begin{equation*}
\frac{\delta \sqrt{-g}}{\delta g_{\mu \nu}}=\frac{\delta}{\delta g_{\mu \nu}} e^{1 / 2 \operatorname{Tr} \log (-g)_{\mu \nu}}=-\frac{1}{2} g^{\mu \nu} \sqrt{-g} \tag{9.7.7}
\end{equation*}
$$

we immediately recover eq. (9.7.5).
A QFT is (classically) scale invariant if its Lagrangian density $\mathcal{L}(x) \rightarrow e^{-4 \sigma} \mathcal{L}(x)$ when $x^{\mu} \rightarrow e^{\sigma} x^{\mu}$, with $\sigma$ a real constant parameter, so that the action $S=\int d^{4} x \mathcal{L}$ remains invariant. The transformations of fields are dictated by their classical dimension $\Delta$ :

$$
\begin{equation*}
\phi(x) \rightarrow e^{-\sigma \Delta} \phi(x) . \tag{9.7.8}
\end{equation*}
$$

[^66]The classical scaling dimension $\Delta$ coincides with the power counting dimension of a field, $\Delta=+1$ for scalars and vectors, and $\Delta=3 / 2$ for fermions in four space-time dimensions. Given the action on the coordinates, each derivative acting on a field increases by one the scaling dimension of the operator:

$$
\begin{equation*}
\partial_{\mu_{1}} \ldots \partial_{\mu_{n}} \phi(x) \rightarrow e^{-\sigma(\Delta+n)} \partial_{\mu_{1}} \ldots \partial_{\mu_{n}} \phi(x) . \tag{9.7.9}
\end{equation*}
$$

A QFT is classically scale invariant if only marginal operators appear in its Lagrangian density. In particular, any mass term explicitly breaks the symmetry. The current associated to scaling transformations is called the dilatation current $D_{\mu}$. Finding its explicit form with the Noether method is non-trivial, because the Lagrangian density is not invariant, even up to total derivatives. On the other hand, it is easy to find $D_{\mu}$ by using the same trick used before of coupling the QFT to gravity. In this way, we can reinterpret the scaling transformation $x^{\mu} \rightarrow e^{\sigma} x^{\mu}$ as a rescaling of the metric $g_{\mu \nu} \rightarrow e^{-2 \sigma} g_{\mu \nu}$. Multiplying eq. (9.7.4) by $\delta g_{\mu \nu}=-2 \epsilon g_{\mu \nu}$, we get that the action variation vanishes if and only if

$$
\begin{equation*}
\theta_{\mu}^{\mu}=0 \tag{9.7.10}
\end{equation*}
$$

The dilatation current is

$$
\begin{equation*}
D^{\mu}=x_{\nu} \theta^{\mu \nu} \tag{9.7.11}
\end{equation*}
$$

and is in fact conserved if $\theta^{\mu \nu}$ is traceless. Notice that $\theta^{\mu \nu}$ entering eq. (9.7.11) does not in general coincide with the $\theta^{\mu \nu}$ in eq. (9.7.4), but is related to it by a total derivative of the form (9.7.3). This is the case even in the simplest theories. For instance, in the massless $\phi^{4}$ theory,

$$
\begin{equation*}
\theta_{\mu}^{\mu}=\left(\partial_{\mu} \phi\right)^{2}-4 \mathcal{L}=-\left(\partial_{\mu} \phi\right)^{2}+\frac{\lambda}{3!} \phi^{4} \neq 0 \tag{9.7.12}
\end{equation*}
$$

We can fix this problem by redefining the energy-momentum tensor. By adding a manifestly conserved tensor, we have for the $\lambda \phi^{4}$ theory

$$
\begin{equation*}
\theta^{\mu \nu} \rightarrow \theta^{\mu \nu}-\frac{1}{6}\left(\partial^{\mu} \partial^{\nu}-\eta^{\mu \nu} \square\right) \phi^{2} \tag{9.7.13}
\end{equation*}
$$

The trace of the redefined energy-momentum tensor reads

$$
\begin{equation*}
\theta_{\mu}^{\mu}=-\left(\partial_{\mu} \phi\right)^{2}+\frac{\lambda}{3!} \phi^{4}+\left(\partial_{\mu} \phi\right)^{2}+\phi \square \phi=\phi\left(\square \phi+\frac{\lambda}{3!} \phi^{3}\right)=0, \tag{9.7.14}
\end{equation*}
$$

where in the last equality we used the equation of motion of $\phi$. One can easily show that by adding a mass term to the $\phi^{4}$ theory, one gets

$$
\begin{equation*}
\theta_{\mu}^{\mu}=m^{2} \phi^{2} \tag{9.7.15}
\end{equation*}
$$

No redefinition of $\theta_{\nu}^{\mu}$ can get rid of the mass term, which is in fact a genuine source of explicit violation of the scale symmetry. Alternatively, we can keep the energy-momentum tensor as defined in eq. (9.7.4) and change the definition of the dilatation current as

$$
\begin{equation*}
D^{\mu}=x_{\nu} \theta^{\mu \nu}+V^{\mu} \tag{9.7.16}
\end{equation*}
$$

The field $V^{\mu}$ is denoted the field virial. For the $\phi^{4}$ theory $V^{\mu}=\phi \partial^{\mu} \phi$.
At the quantum level, scale invariance is typically broken by quantum effects. The "would-be" WT identities implied by scale invariance are

$$
\begin{equation*}
\left\langle\theta_{\mu}^{\mu}(q) \phi_{1}\left(p_{1}\right) \ldots \phi_{n}\left(p_{n}\right)\right\rangle=-\sum_{r=1}^{n} \Delta_{r}\left\langle\phi_{1}\left(p_{1}\right) \ldots \phi_{r}\left(p_{r}+q\right) \ldots \phi_{n}\left(p_{n}\right)\right\rangle \tag{9.7.17}
\end{equation*}
$$

where $\Delta_{r}$ is the scaling dimension of the field $\phi_{r}$. Quantum corrections responsible of the anomalous dimensions of the field do not violate the WT identities (9.7.17), that continue to hold with a redefinition of $\Delta_{r}$ that includes the anomalous dimension contribution. The actual breaking arises from the energy-dependence of couplings, namely from their $\beta$-function. Invariance of the quantum action under scaling transformations would imply that

$$
\begin{align*}
& \int d^{4} x \mathcal{L}\left(\phi(x), \partial_{\mu} \phi(x), \lambda(x)\right)=\int d^{4} x \mathcal{L}\left(e^{-\sigma \Delta} \phi\left(e^{-\sigma} x\right), e^{-\sigma(\Delta+1)} \partial_{\mu} \phi\left(e^{-\sigma} x\right), \lambda(x)\right)  \tag{9.7.18}\\
& =\int d^{4} x e^{4 \sigma} \mathcal{L}\left(e^{-\sigma \Delta} \phi(x), e^{-\sigma(\Delta+1)} \partial_{\mu} \phi(x), \lambda\left(e^{\sigma} x\right)\right)
\end{align*}
$$

where $\lambda$ denotes a coupling constant and in the last equality we have redefined the coordinates $x^{\mu} \rightarrow e^{\sigma} x^{\mu}$. For a classically scale invariant theory $\mathcal{L}\left(e^{-\sigma \Delta} \phi(x), e^{-\sigma(\Delta+1)} \partial_{\mu} \phi(x), \lambda\right)=$ $e^{-4 \sigma} \mathcal{L}\left(\phi(x), \partial_{\mu} \phi(x), \lambda\right)$. Under an infinitesimal transformation,

$$
\begin{equation*}
\lambda(x+\epsilon x)=\lambda(x)+\epsilon \beta(\lambda), \tag{9.7.19}
\end{equation*}
$$

and the violation of the scaling symmetry is given by

$$
\begin{equation*}
\delta S=\int d^{4} x\left(\mathcal{L}\left(\phi(x), \partial_{\mu} \phi(x), \lambda(x+\epsilon x)\right)-\mathcal{L}\left(\phi(x), \partial_{\mu} \phi(x), \lambda(x)\right)\right)=\epsilon \int d^{4} x \frac{\partial \mathcal{L}}{\partial \lambda} \beta(\lambda) \tag{9.7.20}
\end{equation*}
$$

Correspondingly, the divergence of the dilatation current is anomalous, with the anomaly given by $\beta$ :

$$
\begin{equation*}
\partial_{\mu} D^{\mu}=\beta(\lambda) \frac{\partial \mathcal{L}}{\partial \lambda} \tag{9.7.21}
\end{equation*}
$$

Let us see how this explicitly works for the usual $\phi^{4}$ theory. We work in DR with $d=4-\epsilon$. First of all, since we are no longer in 4 dimensions, we have to reconsider how to redefine $\theta^{\mu \nu}$, like in eq. (9.7.13). Let

$$
\begin{equation*}
\theta_{B}^{\mu \nu} \rightarrow \theta_{B}^{\mu \nu}+\alpha\left(\partial^{\mu} \partial^{\nu}-\eta^{\mu \nu} \square\right) \phi_{B}^{2} \tag{9.7.22}
\end{equation*}
$$

where the subscript $B$ stands for bare quantities. On-shell, the trace of the redefined energy-momentum tensor reads

$$
\begin{equation*}
\theta_{B \mu}^{\mu}=\left(1-\frac{d}{2}+2 \alpha(1-d)\right)\left(\partial_{\mu} \phi_{B}\right)^{2}-(d+8 \alpha(1-d)) \frac{\lambda_{B}}{4!} \phi_{B}^{4} \tag{9.7.23}
\end{equation*}
$$

Let us choose $\alpha$ so that the term proportional to $\left(\partial_{\mu} \phi_{B}\right)^{2}$ vanishes:

$$
\begin{equation*}
\alpha=\frac{\left(1-\frac{d}{2}\right)}{2(d-1)} \tag{9.7.24}
\end{equation*}
$$

We get

$$
\begin{equation*}
\theta_{B \mu}^{\mu}=(d-4) \frac{\lambda_{B}}{4!} \phi_{B}^{4} . \tag{9.7.25}
\end{equation*}
$$

We now have

$$
\begin{equation*}
\lambda_{B} \phi_{B}^{4}=Z_{\lambda} \lambda \phi^{4} \tag{9.7.26}
\end{equation*}
$$

with

$$
\begin{equation*}
Z_{\lambda}=1+\frac{3 \lambda}{16 \pi^{2}} \frac{1}{\epsilon} \tag{9.7.27}
\end{equation*}
$$

at one-loop level. Substituting in eq. (9.7.25) gives

$$
\begin{equation*}
\theta_{\mu}^{\mu}=\theta_{B \mu}^{\mu}=-\frac{3 \lambda^{2}}{16 \pi^{2}} \frac{1}{4!} \phi^{4}=\beta(\lambda) \frac{\partial \mathcal{L}}{\partial \lambda}, \tag{9.7.28}
\end{equation*}
$$

which is the expected result. Notice that $\theta_{\mu}^{\mu}$ eventually is finite with no need of any wave function renormalization, in agreement to the general result that conserved currents do not renormalize, $Z_{J}=1$, and correspondingly have vanishing anomalous dimensions.

### 9.8 The Strong CP Problem and a Possible Solution: Axions

Strong interactions seem to respect to a high degree of accuracy invariance under parity P , charge conjugation C and time reversal T . On the other hand, the QCD Lagrangian might contain a gauge and Lorentz invariant dimension four operator of the form

$$
\begin{equation*}
\frac{\theta}{64 \pi^{2}} \int d^{4} x \epsilon_{\mu \nu \rho \sigma} G_{a}^{\mu \nu}(x) G_{a}^{\rho \sigma}(x) \tag{9.8.1}
\end{equation*}
$$

where $G_{a}^{\mu \nu}$ are the (non-canonically normalized) gluon field-strengths and $\theta$ is a real parameter. The parameter $\theta$ in eq.(9.8.1) is actually not physical. Under a chiral rotation of the quark fields

$$
\begin{equation*}
Q_{f} \rightarrow e^{i \alpha_{f} \gamma_{5}} Q_{f} \tag{9.8.2}
\end{equation*}
$$

where $f$ is a flavour index, the QCD action is not invariant due to the global anomaly (9.1.31). The non-invariance of the action corresponds to a redefinition of the parameter $\theta$ :

$$
\begin{equation*}
\theta \rightarrow \theta+2 \sum_{f} \alpha_{f} \tag{9.8.3}
\end{equation*}
$$

The chiral transformation (9.8.2) affects also the quark masses:

$$
\begin{equation*}
m_{f} \rightarrow e^{2 i \alpha_{f}} m_{f} \tag{9.8.4}
\end{equation*}
$$

Hence the invariant parameter $\bar{\theta}$ is the combination

$$
\begin{equation*}
\bar{\theta}=\theta+i \log \prod_{f} m_{f} \tag{9.8.5}
\end{equation*}
$$

If any of the quark masses was zero, the parameter $\theta$ would not be observable, since we will always have the freedom to set it to zero by rotating the massless quark. If $m_{f} \neq 0$, as it appears to be in the real world, $\bar{\theta}$ is a genuine physical source of parity and time reversal violation (and hence also of CP by the CPT theorem) in the strong interactions.

The term (9.8.1) can be shown to be a total derivative and is completely irrelevant at the perturbative level. ${ }^{8}$ Yet, at low energies, when QCD confines, it is expected to give rise to CP violating operators. A particularly relevant CP odd operators is the neutron electric dipole moment operator,

$$
\begin{equation*}
i d_{N} \bar{N} \sigma^{\mu \nu} \gamma_{5} N F_{\mu \nu} \tag{9.8.6}
\end{equation*}
$$

where $N$ is the neutron fermion field. Needless to say, computing $d_{N}$ from first principles is an hard task, considering also that the neutron has a mass of order $\Lambda_{Q C D}$ and cannot be straightforwardly included in the meson low energy action described in section 8.7. A rough simple estimate of the order of magnitude of $d_{N}$ is however possible. Since $d_{N}$ has to vanish if any quark mass is zero, eq.(8.7.9) and dimensional analysis suggest that

$$
\begin{equation*}
\left|d_{N}\right| \sim e|\bar{\theta}| \frac{m_{\pi}^{2}}{m_{N}^{3}} \tag{9.8.7}
\end{equation*}
$$

The current experimental upper bounds on the neutron electric dipole moment is $\left|d_{N}\right| \leq$ $2.9 \times 10^{-26} \mathrm{e} \mathrm{cm}$. Using the theoretical guess (9.8.7) gives the bound

$$
\begin{equation*}
|\bar{\theta}| \lesssim 10^{-10} \tag{9.8.8}
\end{equation*}
$$

More refined estimates confirm the order of magnitude bound given in eq.(9.8.8).
According to the naturalness criterium reviewed in section 7.5, a dimensionless coupling can be very small if a symmetry is restored in the limit it vanishes. This would indeed be the case for $\bar{\theta}$ in QCD, in isolation. In the real world, however, the electroweak interactions break both P and CP and the naturalness of a small $\bar{\theta}$ is not guaranteed. We are left with

[^67]the question: why is $\bar{\theta}$ so small? ${ }^{9}$ This naturalness problem is often denoted strong CP problem.

An elegant solution is achieved by introducing a new field $a(x)$, called the axion, that is supposed to be the NBG of a spontaneously broken $U(1)$ symmetry (at energies much larger than $\left.\Lambda_{Q C D}\right)$. The original formulation of this idea was due to Peccei and Quinn and for this reason this $U(1)$ symmetry is often denoted Peccei-Quinn (PQ) symmetry. Under a $U(1)_{P Q}$ transformation parametrized by $\omega$, the axion shifts

$$
\begin{equation*}
a(x) \rightarrow a(x)+f_{a} \omega \tag{9.8.9}
\end{equation*}
$$

where $f_{a}$ is the axion decay constant. According to the analysis made in sections 8.6 and 8.9, the non-linear realization of the $U(1)_{P Q}$ symmetry requires the axion to have only derivative interactions. The only allowed non-derivative term is an interaction of the form

$$
\begin{equation*}
\frac{1}{64 \pi^{2} f_{a}} \int d^{4} x a(x) \epsilon_{\mu \nu \rho \sigma} G_{a}^{\mu \nu}(x) G_{a}^{\rho \sigma}(x) \tag{9.8.10}
\end{equation*}
$$

From the discussion above, it is clear that the coupling (9.8.10) breaks $U(1)_{P Q}$ only at the non-perturbative level. In presence of this interaction, the $\theta$ term (9.8.1) can trivially be reabsorbed in eq. $(9.8 .10)$ by the $U(1)_{P Q}$ symmetry (9.8.9). We have essentially promoted $\theta$ to be a field, so the problem of the value of $\theta$ becomes now a dynamical one. We have to compute the low-energy effective potential for the axion $a(x)$ and find its minima. The best way to study the axion potential is by exploiting the QCD NGB Lagrangian already introduced in section 8.7. For simplicity consider $n_{f}=2$, keeping only the up and down quarks. By a chiral rotation, we can eliminate the VEV of $a(x)$ from eq.(9.8.10) and reabsorb it in, say, the up quark mass term. As we have seen, quark mass terms are sources of explicit violation of the chiral symmetry, responsible for the pion masses and, more generally, for a pion potential. The leading terms are given in eq.(8.7.8). When the axion VEV is reabsorbed into $M$, this term can be seen as the leading axion-pion effective potential term:

$$
\begin{equation*}
V(a, \pi)=-c f_{\pi}^{3} \operatorname{tr}\left(U M+M^{\dagger} U^{\dagger}\right) \tag{9.8.11}
\end{equation*}
$$

where $U=\exp \left(i \sigma^{a} \pi^{a} / f_{\pi}\right)$ and

$$
M=\left(\begin{array}{cc}
m_{u} e^{-i \frac{a}{f_{a}}} & 0  \tag{9.8.12}\\
0 & m_{d}
\end{array}\right)
$$

[^68]In eq.(9.8.12), $a$ is the axion VEV. By setting $\pi^{ \pm}=0$, the explicit form of $V$ reads

$$
\begin{equation*}
V\left(a, \pi_{0}\right)=-2 c f_{\pi}^{3}\left(m_{d} \cos \left(\frac{\pi_{0}}{f_{\pi}}\right)+m_{u} \cos \left(\frac{a}{f_{a}}-\frac{\pi_{0}}{f_{\pi}}\right)\right) . \tag{9.8.13}
\end{equation*}
$$

The extrema of $V$ are found at $a=0 \bmod \pi f_{a}$ and $\pi_{0}=0 \bmod \pi f_{\pi}$. A straightforward computation shows that the minimum is at

$$
\begin{equation*}
a=\pi_{0}=0 \tag{9.8.14}
\end{equation*}
$$

Hence the $\theta$ angle dynamically vanishes, providing a solution to the strong CP problem. From the potential (9.8.13) we can easily compute the axion mass. Expanding for $f_{a} \gg$ $f_{\pi}, m_{u}, m_{d}$ we get (assuming a canonical kinetic term for $a$ )

$$
\begin{equation*}
m_{a}^{2}=\frac{f_{\pi}^{2}}{f_{a}^{2}} \frac{m_{u} m_{d}}{\left(m_{u}+m_{d}\right)^{2}} m_{\pi}^{2} \tag{9.8.15}
\end{equation*}
$$

Once again, we see the power of an effective description of NGB, that has allowed us to compute the axion mass in a few simple steps. The axion decay constant $f_{a}$ governs all the derivative interactions that the axion can possibly have in the UV theory with quarks and gauge fields, and with mesons, baryons and photons in the IR. Astrophysical constraints coming from red giant cooling puts a lower bound on $f_{a}$ :

$$
\begin{equation*}
f_{a}>10^{9} \mathrm{GeV} \tag{9.8.16}
\end{equation*}
$$

Given eq.(9.8.15) and the above bound, we conclude that the axion must be an extremely light particle:

$$
\begin{equation*}
m_{a} \lesssim 10^{-3} \mathrm{eV} \tag{9.8.17}
\end{equation*}
$$

The axion has the added virtue of being also a viable dark matter candidate. Upper bounds, coming mainly from cosmological considerations, also exist. They depend on some details about the early evolution of the universe, and will not be discussed here.

## Chapter 10

## Some Formal Developments ${ }^{\star}$

### 10.1 Asymptotic Nature of Perturbation Theory*

Most of the considerations made in these notes are based on perturbation theory. It is natural at this stage to ask what are the convergent properties of the associated series. Perturbative expansions in QFT give rise to power series in some coupling constant $\lambda$ or to $\hbar$, if we consider the analogous loopwise expansion. It is useful to recall here some basic mathematical properties of power series of holomorphic functions. If $f(\lambda)$ is analytic at a point $\lambda_{0}$, in a small disc around $\lambda_{0}$ the function is given by the power series

$$
\begin{equation*}
f(\lambda)=\sum_{n=0}^{\infty} f_{n}(\lambda), \quad f_{n}(\lambda)=c_{n}\left(\lambda-\lambda_{0}\right)^{n} \tag{10.1.1}
\end{equation*}
$$

The radius of convergence $R$ of the series (10.1.1) is given, e.g., by

$$
\begin{equation*}
R=\lim _{n \rightarrow \infty}\left|\frac{c_{n}}{c_{n+1}}\right| \tag{10.1.2}
\end{equation*}
$$

If $\lambda_{0}$ is a regular point of $f(\lambda), R$ is non-vanishing and is given by the distance of $\lambda_{0}$ from the closest singularity of $f(\lambda)$. Viceversa, if $R$ is non-vanishing, necessarily $\lambda_{0}$ is a regular point of the function $f$. The power series (10.1.1) is uniformly convergent for any $|\lambda|<R$ and divergent for $|\lambda|>R$. The convergence at $|\lambda|=R$ depends on the particular cases, but necessarily there is at least a point where the series diverges, corresponding to the singular point of $f(\lambda)$ closest to $\lambda_{0}$.
F. Dyson in 1952 presented an argument that led to the conclusion that the perturbative series expansion in QED has zero radius of convergence. It is worth to sketch here Dyson's original simple and brilliant argument. Recall that the QED expansion is a power series in $\alpha \propto e^{2}$ and not in the charge $e$. Suppose now that the QED expansion has a non-vanishing radius of convergence. This would imply that the point $\alpha=0$ should be a
regular point of a physical observable $O(\alpha)$, seen as an analytic function of $\alpha$. As we recalled, this would imply that $O(\alpha)$ has a finite radius of convergence where the function is analytic, including regions where $\alpha<0$. A world with $\alpha<0$ is however unstable, because electrons and positrons would repel each other destroying any vacuum in an uncontrolled production of electron-positron pairs. ${ }^{1}$ We conclude that sensible physical observables cannot be analytic for $\alpha<0$ and thus the point $\alpha=0$ itself cannot be analytic. In turn, this implies that the series around $\alpha=0$ has zero radius of convergence.

A modern and more general version of Dyson's argument can be obtained by considering the euclidean path integral formulation in QFT. Consider a generic $n$-point function

$$
\begin{equation*}
G^{(n)}\left(x_{1}, x_{2}, \ldots x_{n} ; \hbar\right)=\int \mathcal{D} \phi \phi\left(x_{1}\right) \phi\left(x_{2}\right) \ldots \phi\left(x_{n}\right) e^{-S(\phi) / \hbar}, \tag{10.1.3}
\end{equation*}
$$

where we denote collectively all the fields in the action by $\phi .{ }^{2}$ In euclidean space the action is positive definite and the path integral converges. ${ }^{3}$ Consider now the loopwise expansion, that is the expansion of $G^{(n)}$ in powers of $\hbar$. The point $\hbar=0$ is necessarily non-analytic because for any value of $\hbar<0$ the Green function $G^{(n)}$ blows up. We conclude that loopwise perturbative expansions in QFT have generically zero radius of convergence and are divergent. A similar conclusion applies for the coupling constant expansion, that upon rescaling of the fields is equivalent to the loopwise expansion. Considering for simplicity a single coupling constant $\lambda$, we write the action as $S=S_{0}+\lambda \Delta S$, where $S_{0}$ is the free theory and $\Delta S$ the interaction term. Simplifying the notation, we have

$$
\begin{align*}
G^{(n)} & =\int \mathcal{D} \phi \phi_{1} \ldots \phi_{n} \sum_{p=0}^{\infty} \frac{(-\Delta S)^{p} \lambda^{p}}{p!} e^{-S_{0}} \stackrel{?}{=} \sum_{p=0}^{\infty} \lambda^{p} \int \mathcal{D} \phi \phi_{1} \ldots \phi_{n} \frac{(-\Delta S)^{p}}{p!} e^{-S_{0}}  \tag{10.1.4}\\
& =\sum_{p=0}^{\infty} \lambda^{p} G_{p}^{(n)} .
\end{align*}
$$

What we typically compute in perturbation theory are the correction terms $G_{p}^{(n)}$ to the exact Green function $G^{(n)}$. But since the power series in $\lambda$ is never uniformly convergent, having zero radius of convergence, we are not allowed to exchange the order of sum and integration in the second identity above. Despite we know this is an improper step, this is often the only thing we can do to make concrete computations and is indeed what we

[^69]have tacitly done all the times so far. The divergent power series appearing in the second row of eq.(10.1.4) is generally an asymptotic series. For sufficiently small values of the coupling such series reproduces quite accurately the exact result $G^{(n)}$ and this explains why perturbation theory in QFT is useful, despite being divergent.

### 10.1.1 Asymptotic Series and Optimal Truncation*

A series expansion associated to a function $Z(\lambda)$ is asymptotic if, at any fixed order $N$,

$$
\begin{equation*}
Z(\lambda)-\sum_{n=0}^{N-1} Z_{n} \lambda^{n}=\mathcal{O}\left(\lambda^{N}\right), \quad \text { as } \quad \lambda \rightarrow 0 \tag{10.1.5}
\end{equation*}
$$

Notice the crucial difference with respect to convergent series where for $N \rightarrow \infty$ the sum $\sum_{n=0}^{N} Z_{n} \lambda^{n}$ approaches $Z(\lambda)$ for any $\lambda$ within the domain of convergence. For convergent series different functions lead to different series. This is not the case for asymptotic expansions, where different functions can have the same asymptotic expansion. Indeed, if $Z(\lambda)$ has the asymptotic expansion (10.1.5), any other function of the form, say,

$$
\begin{equation*}
\widetilde{Z}(\lambda)=Z(\lambda)+e^{-1 / \lambda} a(\lambda) \tag{10.1.6}
\end{equation*}
$$

with $a(\lambda)$ sufficiently regular, will have exactly the same expansion as $Z(\lambda)$. Asymptotic series are anyhow useful because they approximate the true result with an accuracy that depends on the value of the coupling $\lambda$ and on the behaviour of the series coefficients $Z_{n}$ for $n \gg 1$. Contrary to convergent series, where the more terms are added in the series and the more accurate is the result, in asymptotic series there is an optimal number of terms one should keep, after which adding more terms results in worse and worse accuracy. This is called optimal truncation. Suppose that for $n \gg 1$

$$
\begin{equation*}
Z_{n} \sim n!a^{n} n^{c} \tag{10.1.7}
\end{equation*}
$$

for some real parameters $a$ and $c .^{4}$ The best accuracy for $Z(\lambda)$ is obtained by finding the value $N=N_{\text {Best }}$ that minimizes the error, estimated as the value of the last term not included in the expansion (10.1.5):

$$
\begin{equation*}
\Delta_{Z} \sim Z_{N} \lambda^{N} \tag{10.1.8}
\end{equation*}
$$

Recalling Stirling approximation

$$
\begin{equation*}
n!=\sqrt{2 \pi n} n^{n} e^{-n}\left(1+\mathcal{O}\left(\frac{1}{n}\right)\right) \tag{10.1.9}
\end{equation*}
$$

[^70]one has
\[

$$
\begin{equation*}
Z_{N} \lambda_{N} \sim e^{f(N)}, \quad f(N)=-N+N \log (a \lambda N)+\left(c+\frac{1}{2}\right) \log N \tag{10.1.10}
\end{equation*}
$$

\]

Since by assumption $N \gg 1$, the minimal error will be given by the minimum of the function $f(N)$ :

$$
\begin{equation*}
\partial_{N} f(N)=\log (a \lambda N)+\left(c+\frac{1}{2}\right) \frac{1}{N}=0 \Longrightarrow N_{\text {Best }}=\frac{1}{a \lambda}\left(1+\mathcal{O}\left(\frac{1}{N}\right)\right) . \tag{10.1.11}
\end{equation*}
$$

Plugging back in eq.(10.1.8) we find

$$
\begin{equation*}
\Delta_{Z} \sim e^{-\frac{1}{a \lambda}}, \tag{10.1.12}
\end{equation*}
$$

independently of $c$ at leading order. We see that the smaller is the coupling and the smaller (exponentially) is the error, but no matter how many terms we compute in perturbation theory, asymptotic series fail to reproduce the exact function by (at best) exponentially suppressed terms. This is consistent with the intrinsic ambiguity related to asymptotic series shown in eq.(10.1.6). Keeping more than $N_{\text {Best }}$ terms in the asymptotic series would lead to an increase in the error.

It has been shown that generally the coefficients of the perturbative expansion in QFT behave at parametrically large order as in eq.(10.1.7), due to the exponentially growing number of Feynman diagrams as the number of loops increases.

### 10.1.2 Borel Summation*

Borel summation is a summation method for asymptotic series. Suppose that a function $Z(\lambda)$ admits an asymptotic expansion of the form (10.1.5) with coefficients $Z_{n}$ that at large order goes like in eq.(10.1.7). We define the Borel transform as the function obtained by dividing the original series by a factorially growing factor:

$$
\begin{equation*}
\mathcal{B} Z(t) \equiv \sum_{n=0}^{\infty} \frac{Z_{n}}{n!} t^{n} \tag{10.1.13}
\end{equation*}
$$

Thanks to the division by $n$ !, the series in eq.(10.1.13) has a non-zero radius of convergence where it defines an analytic function $\mathcal{B} Z(t)$. If the analytic continuation of $\mathcal{B} Z(t)$ over the complex $t$ plane is free of singularities for $t>0$, the integral

$$
\begin{equation*}
Z_{B}(\lambda)=\int_{0}^{\infty} d t e^{-t} \mathcal{B} Z(t \lambda) \tag{10.1.14}
\end{equation*}
$$

defines a function of $\lambda$ that is said to be the Borel resum of the original asymptotic series. Recalling the definition of Gamma function

$$
\begin{equation*}
\Gamma(x)=\int_{0}^{\infty} d t e^{-t} t^{x-1} \tag{10.1.15}
\end{equation*}
$$

it is immediate to verify that $Z_{B}(\lambda)$ has the same asymptotic expansion as $Z(\lambda)$. On the other hand, it is also clear that if we approximate the series defining $\mathcal{B Z}(t)$ in eq.(10.1.13) with its truncated version, $Z_{B}(\lambda)$ boils down to the original asymptotic expansion of $Z(\lambda)$ and no progress is achieved. The Borel resummation method works when we can resum the whole series (10.1.13) or estimate the function $\mathcal{B Z}(t)$ by some other means. In general the Borel resummed function $Z_{B}(\lambda)$ is not guaranteed to coincide with the original function $Z(\lambda)$. Indeed, if different functions can admit the same asymptotic series, it is clear that manipulating the latter cannot be enough to uniquely fix $Z(\lambda)$. The uniqueness (and hence the condition $\left.Z_{B}(\lambda)=Z(\lambda)\right)$ is ensured by a theorem provided that certain analyticity properties of $Z(\lambda)$ near the origin are assumed. When $\mathcal{B} Z(t)$ is free of singularities and gives rise to a well-defined $Z_{B}(\lambda)$ we say that the associated asymptotic series is Borel resummable. When $Z_{B}(\lambda)=Z(\lambda)$ we say that the asymtptotic series is Borel resummable to the exact result.

We can get some intuition on Borel functions by working out the Borel transform of an asymptotic series with coefficients as in eq.(10.1.7), with $c=0$. The Borel series in this case collapses to a simple power series and gives

$$
\begin{equation*}
\mathcal{B} Z(t)=\frac{1}{1-a t}, \tag{10.1.16}
\end{equation*}
$$

that has a simple pole at $t=1 / a$. The radius of convergence of the Borel series is given by $R=1 /|a|$, but the function can be analytically continued over the whole $t$-plane. Borel summability depends on the sign of $a$. If $a>0$ (same sign series) the singularity is on the positive real $t$ axis, the integral (10.1.14) is divergent and the series is not Borel resummable. If $a<0$ (alternating series), the singularity is over the negative real $t$ axis, the integral (10.1.14) is finite and the Borel resummed series is given by $Z_{B}(\lambda)$. More in general, eq.(10.1.7) gives only the asymptotic form of the coefficients of the series, while the precise form of the latter might be unavailable. When the exact Borel function $\mathcal{B Z}(t)$ is not known, some information on its analytic structure can still be deduced, because the large order behaviour of the asymptotic series determines the position of the singularity closest to the origin. If $a<0$ the series is no longer guaranteed to be Borel resummable, because further singularities on the positive real axis might occur, depending on the next to leading large order behaviour of the series coefficients. On the other hand, for $a>0$ the series is certainly not Borel resummable.

The order of sum and integration in the Borel function cannot be inverted because the Borel series expansion has generally a finite radius of convergence while the integral is taken over the whole positive $t$ axis. Indeed, if we erroneously interchange the two
operations, we get back the asymptotic divergent series we started with:

$$
\begin{equation*}
Z_{B}(\lambda)=\int_{0}^{\infty} d t e^{-t} \mathcal{B} Z(\lambda t) \neq \sum_{n=0}^{\infty} \frac{Z_{n}}{n!} \lambda^{n} \int_{0}^{\infty} d t t^{n} e^{-t}=\sum_{n=0}^{\infty} Z_{n} \lambda^{n} \tag{10.1.17}
\end{equation*}
$$

Unfortunately the asymptotic series in the coupling in four dimensional gauge theories in four space-time dimensional are generally not Borel resummable. Examples of Borel resummable series include the $\lambda \phi^{4}$ theories in two and three space-time dimensions, important in the study of critical phenomena in statistical physics.

### 10.2 Vacuum Decay in the Presence of External Fields ${ }^{\star}$

In this section we discuss the issue of the stability of the quantum vacuum in the presence of external static and constant electric and magnetic fields. Heuristically, the various fields in the quantum vacuum fluctuate around their (vanishing) mean values and these fluctuations are responsible for some measurable effects, which are particularly important in QED: (i) the Lamb shift in atomic physics, due to the quantum corrections to the vacuum polarization, (ii) the Casimir effect due to the spatial confinement of fluctuations imposed (at the semiclassical level) by conducting bodies, and (iii) pair production due to applied external electric fields. Here we focus on this last case, also known as Schwinger effect. In particular, one can think of this phenomenon as the analogous of the ionization of a neutral atom: in vacuum the virtual pairs of particles (of mass $m$ ) form "bound" pairs of energy $\mathcal{E}_{b}=2 m$ and are confined within a potential well of a typical extension set by the Compton length $\lambda_{c}=1 / m$. In order for an electric field of strength $E$ to unbound this pair and make the constituent particle "real", it is necessary that it makes the bound pair overcome the energy barrier, such that $\lambda_{c} e E \geq \mathcal{E}_{b}$, i.e., $E$ has to exceed the critical value $E_{\text {cr }}=\mathcal{E}_{b} /\left(\lambda_{c} e\right)=2 m^{2} / e$ which, for electrons, turn out to be $E_{\text {cr }} \simeq 2 \times 10^{17} \mathrm{~V} / \mathrm{m}$. In a sense, this phenomenon requires a tunneling of that potential barrier and therefore it is expected that in the expression for the occurrence probability of such a tunneling the field strength $E$ will appear in the denominator of an exponential law, as it happens for thermally activated processes. Still on this heuristic intuition, one expects the phenomenon to be controlled by $E / E_{\text {cr }}$. Note that this pair production implies the decay of the vacuum, which we are going to study in the rest of the section. However, before discussing in details this issue in Sec. 10.2.2, we consider first a seemingly unrelated problem.

### 10.2.1 Landau Levels by Path Integral*

Consider a charged particle of mass $m$ in the presence of a magnetic field. Its classical Lagrangian is

$$
\begin{equation*}
L=\frac{m}{2} \dot{\vec{q}}^{2}+e \dot{\vec{q}} \cdot \vec{A}(\vec{q}) \tag{10.2.1}
\end{equation*}
$$

where $\vec{q}(t)$ is the particle coordinate and $\dot{\vec{q}}=d \vec{q} / d t$. The corresponding Hamiltonian $H$ is given by

$$
\begin{equation*}
H \equiv \vec{p} \cdot \dot{\vec{q}}-L=\frac{(\vec{p}-e \vec{A}(\vec{q}))^{2}}{2 m} \tag{10.2.2}
\end{equation*}
$$

where $H$ has been expressed in terms of the momentum $\vec{p} \equiv \delta L / \delta \dot{\vec{q}}=m \dot{\vec{q}}+e \vec{A}(\vec{q})$.
A classical problem in quantum mechanics is the determination of the energy levels of a quantum particle when it is confined in the $x$ - $y$-plane while being subject to a constant magnetic field $\vec{B}=B \vec{e}_{z}$ (where $\vec{e}_{i}$ is the unit vector along direction $i$ ). Given that $\vec{B}=\vec{\nabla} \times \vec{A}$, one can choose $\vec{A}=B x \vec{e}_{y}$, such that eq. (10.2.1) in terms of operators becomes

$$
\begin{equation*}
\hat{H}=\frac{\hat{p}_{x}^{2}}{2 m}+\frac{\left(\hat{p}_{y}-e B \hat{x}\right)^{2}}{2 m} . \tag{10.2.3}
\end{equation*}
$$

Translation invariance of $\hat{H}$ along $\vec{e}_{y}$ implies $\left[\hat{H}, \hat{p}_{y}\right]=0$ and therefore one can look for the eigenstates of $\hat{H}$ in the form of plane waves with definite momentum $p_{y}=\hbar k_{y}$, for which

$$
\begin{equation*}
\hat{H}=\frac{\hat{p}_{x}^{2}}{2 m}+\frac{\left(\hbar k_{y}-e B \hat{x}\right)^{2}}{2 m}=\frac{\hat{p}_{x}^{2}}{2 m}+\frac{m \omega_{c}^{2}}{2}\left(\hat{x}-x_{0}\right)^{2}, \tag{10.2.4}
\end{equation*}
$$

where

$$
\begin{equation*}
\omega_{c}=\frac{e B}{m} \tag{10.2.5}
\end{equation*}
$$

is the cyclotron frequency and the harmonic oscillator is centered around $x_{0}=\hbar k_{y} /(e B)$. For a given $k_{y}$, the energy levels of this harmonic oscillator are given by $E_{n}=\hbar \omega_{c}(n+1 / 2)$ with $n=$ $0,1, \ldots$ (Landau levels). Each of these levels has a degeneracy $g$ determined by the corresponding possible values of $k_{y}$. In order to determine $g$ consider a two-dimensional rectangular area with edges of length $L_{x}$ and $L_{y}$. Assuming periodic boundary conditions, one finds that $k_{y}=2 \pi n_{y} / L_{y}$, with $n_{y}=0,1, \ldots$, each corresponding to an harmonic oscillator centered at $x_{0}=\hbar 2 \pi n_{y} /\left(e B L_{y}\right)$. Requiring that $x_{0}$ is within the rectangle $0 \leq x_{0} \leq L_{x}$, one finds that $n_{y} \leq e B L_{x} L_{y} /(2 \pi \hbar)$ and therefore $g=e B \mathcal{A} /(2 \pi \hbar)$, where $\mathcal{A}=L_{x} L_{y}$ is the area of the surface across which $\vec{B}$ flows. The fact that $g$ is an integer implies a magnetic flux quantization in multiple integers of the quantum of flux $2 \pi \hbar / e$. Below we recover this result via a path-integral approach.

Consider a gas of non-interacting particles with Hamiltonian given by eq. (10.2.2) and at temperature $\beta^{-1}$. The corresponding single-particle partition function is given by

$$
\begin{equation*}
Z(\beta)=\operatorname{Tr} e^{-\beta H}=\int d \vec{q}_{a}\left\langle\vec{q}_{a}\right| e^{-\beta H}\left|\vec{q}_{a}\right\rangle . \tag{10.2.6}
\end{equation*}
$$

The matrix element $\left\langle\vec{q}_{a}\right| e^{-\beta H}\left|\vec{q}_{a}\right\rangle$ can be obtained via a Wick's rotation of the standard propagator of quantum mechanics $\left\langle\vec{q}_{b}\right| e^{-i H\left(t_{b}-t_{a}\right)}\left|\vec{q}_{a}\right\rangle$ which quantifies the probability amplitude for a particle with Hamiltonian $H$ to propagate from the position $\vec{q}_{a}$ at time $t_{a}$ to the position $\vec{q}_{b}$ at time $t_{b}$. This propagator can be represented as a Feynman pathintegral with an exponential weight determined by the action associated to Lagrangian $L$ in eq. (10.2.1):

$$
\begin{equation*}
\left\langle\vec{q}_{b}\right| e^{-i H\left(t_{b}-t_{a}\right)}\left|\vec{q}_{a}\right\rangle=\mathcal{N} \int_{\substack{\vec{q}\left(t_{a}\right)=\vec{q}_{a} \\ \vec{q}\left(t_{b}\right)=\vec{q}_{b}}} \mathcal{D} \vec{q}(t) e^{i \int_{t_{a}}^{t_{b}} d t L} \tag{10.2.7}
\end{equation*}
$$

By setting $t_{a}=0$ and by continuing $t_{b} \rightarrow-i \tau$, with $\dot{\vec{q}} \rightarrow i d \vec{q} / d \tau$ one gets

$$
\begin{equation*}
Z=\operatorname{Tr} e^{-\beta H}=\mathcal{N} \int d \vec{q}_{a} \int_{\vec{q}(\beta)=\vec{q}(0)=\vec{q}_{a}} \mathcal{D} \vec{q}(\tau) e^{-\int_{0}^{\beta} d \tau L_{E}} \tag{10.2.8}
\end{equation*}
$$

where the Euclidean Lagrangian $L_{E}$ is given by

$$
\begin{equation*}
L_{E}=-\left.L\right|_{\vec{q} \mapsto i d \vec{q} / d \tau}=\frac{m}{2}\left(\frac{d \vec{q}}{d \tau}\right)^{2}-i e \frac{d \vec{q}}{d \tau} \cdot \vec{A}(\vec{q}(\tau)) \tag{10.2.9}
\end{equation*}
$$

The path integral in eq. (10.2.8) is calculated by summing over all closed paths originating in a certain point $\vec{q}_{a}=\vec{q}(0)=\vec{q}(\beta)$, over which one eventually integrates. These paths are in a one-to-one correspondence with periodic functions of period $\beta$ and therefore can be expressed in terms of their Fourier transforms:

$$
\begin{equation*}
\vec{q}(\tau)=\vec{q}_{0}+\sum_{n=1}^{\infty} \sqrt{\frac{2}{\beta}}\left(\vec{c}_{n} \cos \frac{2 \pi n \tau}{\beta}+\vec{s}_{n} \sin \frac{2 \pi n \tau}{\beta}\right) \tag{10.2.10}
\end{equation*}
$$

where $\vec{s}_{n}=\left(s_{x, n}, s_{y, n}, s_{z, n}\right)$ and $\vec{c}_{n}=\left(c_{x, n}, c_{y, n}, c_{z, n}\right)$ are the vector coefficients of the transform which can be obtained by inverting it, i.e.,

$$
\left.\vec{q}_{0}=\frac{1}{\beta} \int_{0}^{\beta} d \tau \vec{q}(\tau), \quad \begin{array}{c}
\vec{c}_{n}  \tag{10.2.11}\\
\vec{s}_{n}
\end{array}\right\}=\int_{0}^{\beta} d \tau \vec{q}(\tau) \sqrt{\frac{2}{\beta}}\left\{\begin{array}{l}
\cos (2 \pi n \tau / \beta) \\
\sin (2 \pi n \tau / \beta)
\end{array}\right.
$$

Note that because the Fourier transform is a unitary transformation, one has

$$
\begin{equation*}
\int_{\vec{q}(0)=\vec{q}(\beta)} \mathcal{D} \vec{q}(\tau)=\int_{-\infty}^{+\infty} d \vec{q}_{0} \prod_{n=1}^{\infty} d \vec{c}_{n} d \vec{s}_{n} \tag{10.2.12}
\end{equation*}
$$

Assume now that this gas of particles is constrained to move on the $x-y$-plane with $q_{z}(\tau)=0$ (and therefore $c_{z, n}=s_{z, n}=0$ ) subject to a constant magnetic field $\vec{B}=B \vec{e}_{z}$, as in the case of the Landau levels discussed above. As before, we can choose $\vec{A}=B x \vec{e}_{y}$ and, by representing the trajectories as in eq. (10.2.10), one can easily express the (Euclidean)
action $S_{E}=\int_{0}^{\beta} d \tau L_{E}$ associated with the (Euclidean) Lagrangian in eq. (10.2.8) in terms of the coefficients $\vec{c}_{n}$ and $\vec{s}_{n}$. In particular, it is useful to note that

$$
\begin{align*}
& \int_{0}^{\beta} d \tau \frac{d \vec{q}}{d \tau} \cdot \vec{A}(\vec{q})=B \int_{0}^{\beta} d \tau \frac{d q_{y}}{d \tau} q_{x}=B \sum_{n=1}^{\infty} \frac{2 \pi n}{\beta}\left(-c_{x, n} s_{y, n}+s_{x, n} c_{y, n}\right),  \tag{10.2.13}\\
& \int_{0}^{\beta} d \tau\left(\frac{d \vec{q}}{d \tau}\right)^{2}=\sum_{n=1}^{\infty}\left(\frac{2 \pi n}{\beta}\right)^{2}\left(c_{x, n}^{2}+c_{y, n}^{2}+s_{x, n}^{2}+s_{y, n}^{2}\right) \tag{10.2.14}
\end{align*}
$$

and therefore

$$
\begin{align*}
S_{E}= & \sum_{n=1}^{\infty}\left\{\frac{m}{2}\left(\frac{2 \pi n}{\beta}\right)^{2}\left(c_{x, n}^{2}+c_{y, n}^{2}+s_{x, n}^{2}+s_{y, n}^{2}\right)-i e B \frac{2 \pi n}{\beta}\left(s_{x, n} c_{y, n}-c_{x, n} s_{y, n}\right)\right\} \\
= & \sum_{n=1}^{\infty}\left\{\frac{m}{2}\left(\frac{2 \pi n}{\beta}\right)^{2}\left(c_{y, n}^{2}+s_{x, n}^{2}\right)-i e B \frac{2 \pi n}{\beta} s_{x, n} c_{y, n}\right\} \\
& +\sum_{n=1}^{\infty}\left\{\frac{m}{2}\left(\frac{2 \pi n}{\beta}\right)^{2}\left(c_{x, n}^{2}+s_{y, n}^{2}\right)+i e B \frac{2 \pi n}{\beta} c_{x, n} s_{y, n}\right\} \\
= & \sum_{n=1}^{\infty} \frac{1}{2}\left(c_{y, n}, s_{x, n}\right) M_{n}^{*}\binom{c_{y, n}}{s_{x, n}}+\sum_{n=1}^{\infty} \frac{1}{2}\left(c_{x, n}, s_{y, n}\right) M_{n}\binom{c_{x, n}}{s_{y, n}} \tag{10.2.15}
\end{align*}
$$

where

$$
M_{n} \equiv\left(\begin{array}{cc}
m(2 \pi n / \beta)^{2} & i e B(2 \pi n / \beta)  \tag{10.2.16}\\
i e B(2 \pi n / \beta) & m(2 \pi n / \beta)^{2}
\end{array}\right)
$$

As expected, $S_{E}$ is independent of $\vec{q}_{0}$. Accordingly, eq. (10.2.8) becomes

$$
\begin{align*}
Z & =\mathcal{N} \int_{-\infty}^{+\infty} d q_{x, 0} d q_{y, 0} \prod_{n=1}^{\infty} d c_{x, n} d c_{y, n} d s_{x, n} d s_{y, n} e^{-S_{E}} \\
& =\mathcal{N} \mathcal{A} \prod_{n=1}^{\infty} \frac{\pi}{\sqrt{\operatorname{det} M_{n}}} \frac{\pi}{\sqrt{\operatorname{det} M_{n}^{*}}}=\mathcal{N}^{\prime} \mathcal{A} \prod_{n=1}^{\infty}\left[1+\left(\frac{\beta \omega_{c}}{2 \pi n}\right)^{2}\right]^{-1} \tag{10.2.17}
\end{align*}
$$

where $\omega_{c}$ is defined in eq.(10.2.5), we used the fact that the integral over $\vec{q}_{0}$ renders the area $\mathcal{A}$, and

$$
\begin{equation*}
\operatorname{det} M_{n}=m^{2}\left(\frac{2 \pi n}{\beta}\right)^{4}\left[1+\left(\frac{\beta \omega_{c}}{2 \pi n}\right)^{2}\right] \tag{10.2.18}
\end{equation*}
$$

(see eq. (10.2.5)), while we included all the factors which do not depend on $B$ in the overall normalization $\mathcal{N}^{\prime}$ of the path integral. In turn, this constant can be easily calculated by considering the same problem for $B=0$, i.e., by calculating the partition function $Z_{0}(\beta)$
of a free quantum particle with Hamiltonian $H_{0}=\vec{p}^{2} /(2 m)$ which is confined in a $d$ dimensional space with volume $\mathcal{A}$ :

$$
\begin{equation*}
Z_{0}(\beta)=\int \frac{d \vec{q} d \vec{p}}{(2 \pi \hbar)^{d}} e^{-\beta H_{0}}=\mathcal{A}\left(\frac{m}{2 \pi \hbar \beta}\right)^{d / 2} \tag{10.2.19}
\end{equation*}
$$

By requiring that $Z$ in eq. (10.2.5) reduces to $Z_{0}$ for $\omega_{c}=0$ (with $d=2$ ), one fixes the value of $\mathcal{N}^{\prime}$ and finds that

$$
\begin{equation*}
Z(\beta)=\mathcal{A} \frac{m}{2 \pi \hbar \beta} \frac{\beta \omega_{c} / 2}{\sinh \left(\beta \omega_{c} / 2\right)}=\frac{\mathcal{A} e B}{2 \pi \hbar} \sum_{n=0}^{\infty} e^{-\beta \omega_{c}(n+1 / 2)} \tag{10.2.20}
\end{equation*}
$$

where we used the identity

$$
\begin{equation*}
\prod_{n=1}^{\infty}\left(1+\frac{x^{2}}{\pi^{2} n^{2}}\right)=\frac{\sinh x}{x} \tag{10.2.21}
\end{equation*}
$$

and the expansion of sinh in terms of exponentials. Recalling the definition of $Z(\beta)$ in eq. (10.2.6) one can alternatively calculate it in the basis of the eigenstate of the energy $H$ (with eigenvalues $E_{n}$ and degeneracy $g_{n}$ ), finding that $Z(\beta)=\sum_{n} g_{n} e^{-\beta E_{n}}$. By comparing this general expansion with eq. (10.2.20) it is possible to identify $E_{n}=\hbar \omega_{c}(n+1 / 2)$ and $g_{n}=\mathcal{A} e B /(2 \pi \hbar)$, which are readily recognized to be the Landau levels that we discussed before. Note that, due to the fact that $g_{n}$ has to be an integer, one concludes that the flux $\mathcal{A} B$ of a magnetic field trough a certain surface of area $\mathcal{A}$ has to be a multiple integer of the elementary flux $2 \pi / e$. This is nothing but the Dirac quantization condition for a flux of $B$ through a (large) close surface enclosing a magnetic monopole.

### 10.2.2 Vacuum Instability for a Constant Electric Field ${ }^{\star}$

In order to address the issue of the vacuum instability in the presence of external fields, we consider first the case of a scalar field, i.e., the scalar QED with Lagrangian

$$
\begin{equation*}
L=-\frac{1}{4} F_{\mu \nu} F^{\mu \nu}+\left(D_{\mu} \phi\right)^{\dagger} D^{\mu} \phi-m^{2} \phi^{2} \tag{10.2.22}
\end{equation*}
$$

where $D_{\mu}=\partial_{\mu}-i e A_{\mu}$ is the covariant derivative and the field $A_{\mu}$ is assumed to be assigned from the outset. (We neglect here the self-interaction of the scalar field as it only provides a small correction to the phenomenon of vacuum instability.) Denoting by $|0\rangle$ the vacuum of the system (of large volume $V$ ), its decay rate in the presence of the field can be determined by looking at the effective action $S_{Q}[A]$ obtained by integrating out the fluctuations of $\phi$, i.e.,

$$
\begin{equation*}
e^{i S_{Q}[A]}=\langle 0| e^{-i H T}|0\rangle=\int \mathcal{D} \phi \mathcal{D} \phi^{*} e^{i \int d^{4} x L} . \tag{10.2.23}
\end{equation*}
$$

In fact, as long as the vacuum is stable, the probability amplitude $\langle 0| e^{-i H T}|0\rangle$ has unit modulus and therefore

$$
\begin{equation*}
S_{Q}[A]=-\frac{1}{4} \int d^{4} x F_{\mu \nu} F^{\mu \nu}+\Delta S_{Q}[A] \tag{10.2.24}
\end{equation*}
$$

is real, whereas if $\Delta S_{Q}$ acquires an imaginary part, the vacuum decays with a rate per unit volume given by

$$
\begin{equation*}
\Gamma \equiv \frac{2 \operatorname{Im} \Delta S_{Q}[A]}{V T} \tag{10.2.25}
\end{equation*}
$$

the same relationship applies to the case of QED, discussed further below. Let us mention that according to the heuristic picture of the Schwinger effect provided at the beginning of this section, the rate $\Gamma$ can be interpreted as the rate (per volume) of production of real pairs from the virtual ones of the vacuum.

Most of the results presented hereafter were originally derived by J. Schwinger in ref. [28].

### 10.2.3 Instability of a Scalar Field Vacuum ${ }^{\star}$

The one-loop contribution $\Delta S_{Q}[A]$ to the quantum action $S_{Q}[A]$ due to the interaction of the complex scalar field with the background field $A_{\mu}$ is given, according to eqs. (10.2.22) and (10.2.23), by

$$
\begin{equation*}
i \Delta S_{Q}[A]=\ln \left\{\mathcal{N} \int \mathcal{D} \phi \mathcal{D} \phi^{*} e^{i \int d^{4} x \phi^{*}\left(-D^{2}-m^{2}\right) \phi}\right\} \tag{10.2.26}
\end{equation*}
$$

In order to calculate the Gaussian integral on the r.h.s. it is convenient to perform a Wick's rotation by introducing a coordinate $x_{4}$ such that $x_{0}=-i x_{4}$ and correspondingly by replacing the temporal component $A_{0}$ of the vector potential by $i A_{4}$. As a result, $D_{0}=\partial_{0}-i e A_{0}=i D_{4}$ with $D_{4} \equiv \partial_{4}-i e A_{4}$ and $D^{2}=D_{0}^{2}-\vec{D}^{2}=-\left(D_{4}^{2}+\vec{D}^{2}\right) \equiv$ $-D_{E}^{2}$, where we define $D_{E, \mu}=\partial_{\mu}-i e A_{E, \mu}$ with $\mu=1,2,3,4$, and Euclidean metric. Correspondingly, one defines $\vec{A}_{E}=\vec{A}$ and $A_{E, 4}=-i A_{0}$. As a result, $i \int d^{4} x \phi^{*}\left(-D^{2}-\right.$ $\left.m^{2}\right) \phi=-\int d^{4} x_{E} \phi^{*}\left(-D_{E}^{2}+m^{2}\right) \phi$ and the Gaussian integral can be easily calculated, leading to

$$
\begin{equation*}
i \Delta S_{Q}[A]=-\operatorname{Tr} \ln \frac{-D_{E}^{2}+m^{2}}{\mu^{2}}+\ldots \tag{10.2.27}
\end{equation*}
$$

where henceforth we neglect irrelevant constants and we introduce a convenient momentum scale $\mu$. Note that the operator $D_{E}^{2}$ cannot be trivially diagonalized in momentum space, due to the spatial dependence of $A_{\mu}$. However, the problem of calculating $\operatorname{Tr} \ln (\cdots)$ can be conveniently related to the one we discussed in Sec. 10.2.1, i.e., to the calculation of
$\operatorname{Tr} e^{-(\cdots)}$, by noticing that

$$
\begin{equation*}
\ln \frac{a}{b}=-\int_{0}^{\infty} \frac{d t}{t}\left(e^{-a t}-e^{-b t}\right) \tag{10.2.28}
\end{equation*}
$$

By using eq. (10.2.28) we get

$$
\begin{equation*}
i \Delta S_{Q}[A]=\int_{0}^{\infty} \frac{d \beta}{\beta} e^{-\beta m^{2}} \operatorname{Tr} e^{-\beta\left(-D_{E}^{2}\right)}+\text { const. . } \tag{10.2.29}
\end{equation*}
$$

The constant term in eq. (10.2.29) is independent of $A$. It simply provides a regularization of the first term and can be neglected in what follows. The connection with the problem discussed in Sec. 10.2.1 is apparent by noticing that

$$
\begin{equation*}
-D_{E}^{2}=\left(-i \vec{\partial}-e \vec{A}_{E}\right)^{2} \tag{10.2.30}
\end{equation*}
$$

formally coincides with the Hamiltonian reported in eq. (10.2.2) of a quantum particle in 4 spatial dimensions, with momentum $\vec{p}=-i \vec{\partial}$, mass $m=1 / 2$, in the presence of an external field $\vec{A}_{E}$. Accordingly, $\operatorname{Tr} e^{-\beta\left(-D_{E}^{2}\right)}$ can be identified with the partition function $Z(\beta)$ of this particle and it can be calculated as explained in Sec. 10.2.1 (see eqs. (10.2.8) and (10.2.9)), where now the position vector $\vec{q}(\tau)$ of the fictitious particle has 4 components instead of 2 . This fact introduces a slight complication in the calculation, which we discuss here. In order to carry out the calculation we need to specify the form of the vector potential $\vec{A}_{E}$. Assuming that the corresponding Euclidean field $F_{E, \mu \nu}=\partial_{\mu} A_{E, \nu}-$ $\partial_{\nu} A_{E, \mu}=-F_{E, \nu \mu}$ is constant in space and time, a possible choice of the vector potential is

$$
\begin{equation*}
A_{E, \mu}(\vec{q})=\frac{1}{2} F_{E, \mu \nu} q_{\nu} \tag{10.2.31}
\end{equation*}
$$

such that the Euclidean Lagrangian (see eq. (10.2.9)) of this particle becomes

$$
\begin{equation*}
L_{E}=\frac{m}{2}\left(\frac{d \vec{q}}{d \tau}\right)^{2}-i e\left(\frac{d \vec{q}}{d \tau}\right)^{T} \frac{F_{E}}{2} \vec{q}(\tau) \tag{10.2.32}
\end{equation*}
$$

where $F_{E}$ is the (antisymmetric) matrix of the field and $T$ indicates the transposition. With this $L_{E}$, the path integral in eq. (10.2.8) can be calculated by decomposing $\vec{q}$ as in eq. (10.2.10) and by expressing the corresponding Euclidean action $S_{E}$ in terms of the coefficients $\vec{s}_{n}$ and $\vec{c}_{n}$, taking into account the natural generalizations of eqs. (10.2.13) and (10.2.14), i.e.,

$$
\begin{align*}
& \int_{0}^{\beta} d \tau \frac{d q_{\mu}}{d \tau} q_{\nu}(\tau)=\sum_{n=1}^{\infty} \frac{2 \pi n}{\beta}\left(-c_{\mu, n} s_{\nu, n}+s_{\mu, n} c_{\nu, n}\right)  \tag{10.2.33}\\
& \int_{0}^{\beta} d \tau\left(\frac{d q_{\mu}}{d \tau}\right)^{2}=\sum_{n=1}^{\infty}\left(\frac{2 \pi n}{\beta}\right)^{2}\left(c_{\mu, n}^{2}+s_{\mu, n}^{2}\right) \tag{10.2.34}
\end{align*}
$$

As the term on the l.h.s. of the first equation is contracted in $S_{E}$ with the antisymmetric matrix $F_{E}$, the two terms on the r.h.s. eventually give the same contribution and therefore

$$
\begin{equation*}
S_{E}=\sum_{n=1}^{\infty}\left\{\frac{m}{2}\left(\frac{2 \pi n}{\beta}\right)^{2} \sum_{\mu=1}^{4}\left(c_{\mu, n}^{2}+s_{\mu, n}^{2}\right)-i e\left(\frac{2 \pi n}{\beta}\right) \sum_{\mu, \nu=1}^{4} s_{\mu, n} F_{E, \mu \nu} c_{\nu, n}\right\} . \tag{10.2.35}
\end{equation*}
$$

Accordingly, the partition function can be written as

$$
\begin{equation*}
Z=\mathcal{N} \int \prod_{\mu=1}^{4}\left(d q_{\mu, 0} \prod_{n=1}^{\infty} d c_{\mu, n} d s_{\mu, n}\right) e^{-S_{E}} \tag{10.2.36}
\end{equation*}
$$

where the integral over $\vec{q}_{0}$ renders the Euclidean space-time volume $i T V$, while the integral over each single $s_{\mu, n}$ is a Gaussian integral in the presence of a linear term of the form

$$
\begin{equation*}
\int d s_{\mu, n} e^{-\alpha s_{\mu, n}^{2}+i \gamma s_{\mu, n}}=\sqrt{\frac{\pi}{\alpha}} e^{-\gamma^{2} /(4 \alpha)}, \tag{10.2.37}
\end{equation*}
$$

with $\alpha=(\pi n / \beta)^{2}$ (with $\left.m=1 / 2\right)$, and $\gamma=e(2 \pi n / \beta) F_{E, \nu \mu} c_{\mu, n}$. After this first integration, one finds (up to irrelevant constants which are absorbed in the definition of $\mathcal{N}^{\prime}$ and then $\mathcal{N}^{\prime \prime}$ )

$$
\begin{align*}
Z & =\mathcal{N}^{\prime} i T V \int \prod_{n=1}^{\infty}\left(\prod_{\mu=1}^{4} d c_{\mu, n} \exp \left\{-(\pi n / \beta)^{2} c_{\mu, n} M_{\mu \nu} c_{\nu, n}\right\}\right)  \tag{10.2.38}\\
& =\mathcal{N}^{\prime \prime} i T V \prod_{n=1}^{\infty}\left(\operatorname{det} M_{n}\right)^{-1 / 2}
\end{align*}
$$

where we introduced the $4 \times 4$ matrix

$$
\begin{equation*}
M_{n, \mu \nu} \equiv \delta_{\mu \nu}-\left(\frac{e \beta}{\pi n}\right)^{2}\left(F_{E}^{2}\right)_{\mu \nu} \tag{10.2.39}
\end{equation*}
$$

As in the case discussed in Sec. 10.2.1, the normalization constant $\mathcal{N}^{\prime \prime}$ can be fixed by comparing this result with the one in the absence of the coupling to the field $A_{E, \mu}$, i.e., for $e=0$, which reduces to the partition function of a free particle (with mass $m=1 / 2$ ) calculated in eq. (10.2.19) for generic dimension $d$, with an area $\mathcal{A}$ given by the Euclidean volume $i V T$. Accordingly,

$$
\begin{equation*}
Z(\beta)=i T V(4 \pi \beta)^{-2} \prod_{n=1}^{\infty}\left(\operatorname{det} M_{n}\right)^{-1 / 2} \tag{10.2.40}
\end{equation*}
$$

In order to proceed further we have to specify the fields $\vec{E}$ and $\vec{B}$, which enter the tensor $F_{\mu \nu}$ as $F_{0 j}=E_{j}$ with $j=1,2$, and 3 , while $F_{j k}=\epsilon_{j k l} B_{l}$. Accordingly, the components
of the Euclidean tensor are given by $F_{E, 4 j}=-i F_{0 j}=-i E_{j}$ and $F_{E, j k}=F_{j k}=\epsilon_{j k l} B_{l}$. Without loss of generality, we consider here the case in which the electric field $\vec{E}=E \vec{e}_{x}$ is directed along the $x$-axis, while the magnetic field $\vec{B}=B_{x} \vec{e}_{x}+B_{y} \vec{e}_{y}$ is on the $x$ - $y$-plane, which result in the Euclidean field

$$
F_{E}=\left(\begin{array}{cccc}
0 & 0 & -B_{y} & i E  \tag{10.2.41}\\
0 & 0 & B_{x} & 0 \\
B_{y} & -B_{x} & 0 & 0 \\
-i E & 0 & 0 & 0
\end{array}\right)
$$

with

$$
F_{E}^{2}=\left(\begin{array}{cccc}
-B_{y}^{2}+E^{2} & B_{x} B_{y} & 0 & 0  \tag{10.2.42}\\
B_{x} B_{y} & -B_{x}^{2} & 0 & 0 \\
0 & 0 & -B_{y}^{2}-B_{x}^{2} & i B_{y} E \\
0 & 0 & i B_{y} E & E^{2}
\end{array}\right)
$$

Due to the block-diagonal structure of $F_{E}^{2}$, the 4 eigenvalues of this matrix are given by the 2 eigenvalues of the upper left $2 \times 2$ matrix and by those of the lower right $2 \times 2$ matrix. However it is easy to see that these two $2 \times 2$ matrices have the same eigenvalues, which therefore have multiplicity two and which are solutions of the equation

$$
\begin{equation*}
\lambda^{2}+\left(\vec{B}^{2}-\vec{E}^{2}\right) \lambda-(\vec{B} \cdot \vec{E})^{2}=0 \tag{10.2.43}
\end{equation*}
$$

i.e.,

$$
\begin{equation*}
\lambda_{ \pm}=\frac{1}{2}\left[\vec{E}^{2}-\vec{B}^{2} \pm \sqrt{\left(\vec{B}^{2}-\vec{E}^{2}\right)^{2}+4(\vec{B} \cdot \vec{E})^{2}}\right] \equiv \pm a_{ \pm}^{2} \tag{10.2.44}
\end{equation*}
$$

Accordingly

$$
\begin{equation*}
\operatorname{det} M_{n}=\left(1-\frac{e^{2} \beta^{2}}{\pi^{2} n^{2}} a_{+}^{2}\right)^{2}\left(1+\frac{e^{2} \beta^{2}}{\pi^{2} n^{2}} a_{-}^{2}\right)^{2} \tag{10.2.45}
\end{equation*}
$$

and $Z$ in eq. (10.2.40) can be calculated by using eq. (10.2.21):

$$
\begin{equation*}
Z(\beta)=\operatorname{Tr} e^{-\beta\left(-D_{E}^{2}\right)}=\frac{i T V}{(4 \pi \beta)^{2}} \frac{e \beta a_{+}}{\sin \left(e \beta a_{+}\right)} \frac{e \beta a_{-}}{\sinh \left(e \beta a_{-}\right)} \tag{10.2.46}
\end{equation*}
$$

Note that

$$
\begin{equation*}
Z(\beta \rightarrow 0)=\frac{i T V}{(4 \pi \beta)^{2}}\left[1+\frac{e^{2} \beta^{2}}{6}\left(\vec{E}^{2}-\vec{B}^{2}\right)+O\left(\beta^{4}\right)\right] \tag{10.2.47}
\end{equation*}
$$

and therefore the integral over $\beta$ in eq. (10.2.29) (which defines $i \Delta S_{Q}[A]$ ) displays a leading divergence $\sim \beta^{-2}$ for $\beta \rightarrow 0$ which is independent of the field and therefore of the coupling constant $e$ and which is cured by the regularizing term that was omitted in that equation. The next, subleading logarithmic divergence is proportional to $E^{2}-\vec{B}^{2}$, i.e., it depends on the fields and therefore it cannot be taken care of by the additive counter-terms mentioned
above. In order to understand the origin of this divergence it is convenient to generalize the present analysis to generic dimensionality $d$ close to 4 . The primary effect of having $d \neq 4$ is to change the determination of the normalization constant $\mathcal{N}^{\prime \prime}$ by comparison with the case of a free particle. Going back to eq. (10.2.19) one concludes that this generalization amounts at multiplying the r.h.s. of eq. (10.2.46) by the factor $(4 \pi \beta)^{2-d / 2}$, while $V T$ has to be understood as the volume of space-time in the corresponding dimensionality. Accordingly, from eq. (10.2.29), one finds

$$
\begin{equation*}
\Delta S_{Q}[A]=V T \int_{0}^{\infty} \frac{d \beta}{\beta} e^{-\beta m^{2}}(4 \pi \beta)^{-d / 2}\left[1+\frac{e^{2} \beta^{2}}{6}\left(\vec{E}^{2}-\vec{B}^{2}\right)+O\left(\beta^{4}\right)\right]+\text { const. } \tag{10.2.48}
\end{equation*}
$$

which contains a dimensional pole for $\epsilon=4-d \rightarrow 0$, as expected. In particular, the part proportional to the fields turn out to be

$$
\begin{equation*}
\Delta S_{Q}[A]=V T \frac{e^{2}}{6(4 \pi)^{2}}\left(\vec{E}^{2}-\vec{B}^{2}\right) \times \frac{2}{\epsilon}+\ldots \tag{10.2.49}
\end{equation*}
$$

and we expect this dimensional pole to be cured by a suitable renormalization of the field amplitude. In fact, by introducing the field-strength renormalization constant $Z_{3}$, the quantum action in eq. (10.2.24) for spatially and temporally constant fields can be written as

$$
\begin{equation*}
S_{Q}[A]=V T \frac{\vec{E}^{2}-\vec{B}^{2}}{2} Z_{3}+V T \frac{e^{2}}{6(4 \pi)^{2}}\left(\vec{E}^{2}-\vec{B}^{2}\right) \times \frac{2}{\epsilon}+\ldots \tag{10.2.50}
\end{equation*}
$$

where we used the fact that $F_{\mu \nu} F^{\mu \nu}=-F_{\mu \nu} F^{\nu \mu}=-F_{E, \mu \nu} F_{E, \nu \mu}=-\operatorname{Tr} F_{E}^{2}=2\left(\vec{B}^{2}-\vec{E}^{2}\right)$ (see eq. (10.2.42)) on the field configuration that we are discussing. Accordingly, within the minimal subtraction scheme, one can remove the divergence by fixing

$$
\begin{equation*}
Z_{3}=1-\frac{e^{2}}{(4 \pi)^{2}} \frac{2}{3 \epsilon}+O\left(e^{4}\right) \tag{10.2.51}
\end{equation*}
$$

which is indeed the field-strength renormalization constant that one determines by a direct renormalization of the photon propagator in scalar QED. Accordingly, the expression of $S_{Q}[A]$ which follows from eqs. (10.2.29) and (10.2.46) can be made finite with suitable renormalizations. Having established this, let us focus on the one-loop contribution directly in $d=4$ :

$$
\begin{equation*}
\Delta S_{Q}[A]=\frac{V T}{16 \pi^{2}} \int_{0}^{\infty} \frac{d u}{u} e^{-u} \frac{e a_{+}}{\sin \left(e u a_{+} / m^{2}\right)} \frac{e a_{-}}{\sinh \left(e u a_{-} / m^{2}\right)} \tag{10.2.52}
\end{equation*}
$$

where the integral has been expressed in terms of the variable $u=\beta m^{2}$. Even after renormalization, a close inspection of this integral reveals that it does not converge because of the presence of zeros of the denominator also for $u>0$ and corresponding to
$u=u_{n} \equiv m^{2} \pi n /\left(e a_{+}\right)$. However, one should remember of the Feynman prescription for regularizing the path integral (as when calculating propagators), which effectively amounts at substituting $m^{2} \rightarrow m^{2}-i \varepsilon$ with $\varepsilon \rightarrow 0^{+}$. With this prescription, the singularities just mentioned move out of the real axis and the integral can be safely calculated. However - as it is the case when studying the decay of unstable particles - this also implies the emergence of an imaginary part in $\Delta S_{Q}[A]$, which is eventually responsible for the vacuum decay. In particular $1 / \sin x \simeq(-1)^{n} /(x-\pi n)$ for $x \rightarrow \pi n$ and taking into account eq. (2.1.26) one concludes that

$$
\begin{equation*}
\operatorname{Im} \frac{1}{\sin (x+i \varepsilon)}=-\pi \sum_{n=0}^{\infty}(-1)^{n} \delta(x-\pi n) \tag{10.2.53}
\end{equation*}
$$

Accordingly,

$$
\begin{equation*}
\operatorname{Im} \Delta S_{Q}[A]=\frac{V T}{(4 \pi)^{2}} e^{2} a_{+} a_{-} \sum_{n=1}^{\infty} \frac{(-1)^{n+1}}{n} \frac{e^{-n \pi m^{2} /\left(e a_{+}\right)}}{\sinh \left(n \pi a_{-} / a_{+}\right)}, \tag{10.2.54}
\end{equation*}
$$

(where the contribution due to $n=0$ is discarded, being cancelled by the renormalization discussed above) and, according to eq. (10.2.25), we find that the decay rate of the vacuum is

$$
\begin{equation*}
\Gamma=\frac{e^{2} a_{+} a_{-}}{8 \pi^{2}} \sum_{n=1}^{\infty} \frac{(-1)^{n+1}}{n} \frac{e^{-n \pi m^{2} /\left(e a_{+}\right)}}{\sinh \left(n \pi a_{-} / a_{+}\right)} \tag{10.2.55}
\end{equation*}
$$

Note that this expression is non-perturbative in the coupling constant $e$, due to the essential singularity associated with the exponential factor. Consider now the two relevant cases in which either field (a) $\vec{E}$ or (b) $\vec{B}$ is present: from eq. (10.2.44) one concludes that $a_{+}=|\vec{E}|$ and $a_{-}=0$ in the former case, whereas $a_{+}=0$ and $a_{-}=|\vec{B}|$ in the latter. As a consequence of the vanishing of $a_{+}$, the decay rate $\Gamma$ of the vacuum vanishes in the presence of $\vec{B}$ alone, as it could have been expected on the basis of the fact that a magnetic field does not do any work on a charged particle and therefore it cannot provide the necessary energy to transform virtual pairs into real ones. On the other hand, an electric field $\vec{E}$ induces a decay of the vacuum even in the absence of a magnetic field $\vec{B}$ and, in fact, eq. (10.2.55) becomes

$$
\begin{equation*}
\Gamma=\frac{e^{2}|\vec{E}|^{2}}{8 \pi^{3}} \sum_{n=1}^{\infty} \frac{(-1)^{n+1}}{n^{2}} e^{-n \pi m^{2} /(e|\vec{E}|)} \tag{10.2.56}
\end{equation*}
$$

for $\vec{B}=0$. As expected from the heuristic discussion presented in the introduction to the issue of vacuum instability, $\Gamma / m^{4}$ in this equation is a function $|\vec{E}| / E_{\text {cr }}$ with $E_{\text {cr }}=2 m^{2} / e$. Figure 10.1, presented at the end of next section, provides a plot of $\Gamma /(e|\vec{E}|)^{2}$ as a function of $|\vec{E}| / E_{\text {cr }}$ and compares it with the case of QED which we discuss next.

### 10.2.4 Instability of a Fermion Field Vacuum ${ }^{\star \star}$

Consider now the case in which the vacuum is that one of fermions, i.e., the case of QED with fermions of mass $m$. Following the same line of argument as in the previous subsection, the decay rate is determined by the effective quantum action

$$
\begin{align*}
i \Delta S_{Q}[A] & =\ln \left\{\mathcal{N} \int \mathcal{D} \psi \mathcal{D} \bar{\psi} \exp \left[i \int d^{4} x \bar{\psi}(I D D-m) \psi\right]\right\} \\
& =\operatorname{Tr} \ln \left(\frac{\text { ID }-m}{\mu}\right)+\ldots \tag{10.2.57}
\end{align*}
$$

Here we note that in taking the trace $\operatorname{Tr}_{D}$ over Dirac matrices of a generic scalar function $f$ of $\not D$ one has that $\operatorname{Tr}_{D} f(i \not D)=\operatorname{Tr}_{D} f(-i \not D)$ because $\operatorname{Tr}_{D}(i \not D)^{n}=\operatorname{Tr}_{D} \gamma_{5}(i \not D)^{n} \gamma_{5}=$ $\operatorname{Tr}_{D}\left(\gamma_{5}\left\langle\boldsymbol{D} \gamma_{5}\right)^{n}=\operatorname{Tr}_{D}(-i \not D)^{n}\left(\right.\right.$ being $\left(\gamma_{5}\right)^{2}=\mathbb{\square}$ and $\left.\gamma_{5}, \gamma_{\mu}=0\right)$. Accordingly,

$$
\begin{equation*}
\operatorname{Tr} \ln \left(\frac{\ddot{D} \mid \mathrm{D}-m}{\mu}\right)=\frac{1}{2} \operatorname{Tr} \ln \left(\frac{\ddot{D D}-m}{\mu} \frac{-\ddot{D}-m}{\mu}\right)=\frac{1}{2} \operatorname{Tr} \ln \left(\frac{\not D^{2}+m^{2}}{\mu^{2}}\right) \tag{10.2.58}
\end{equation*}
$$

which renders an expression quite similar to eq. (10.2.27). Using the facts that $\left\{\gamma_{\mu}, \gamma_{\nu}\right\}=$ $2 \eta_{\mu \nu},\left[\gamma_{\mu}, \gamma_{\nu}\right] \equiv-2 i \sigma_{\mu \nu}$, and that the convariant derivative satisfies $\left[D_{\mu}, D_{\nu}\right]=i e F_{\mu \nu}$, we can write

$$
\begin{equation*}
\not D^{2}=\frac{1}{2}\left(\left\{\gamma_{\mu}, \gamma_{\nu}\right\}+\left[\gamma_{\mu}, \gamma_{\nu}\right]\right) D^{\mu} D^{\nu}=D^{2}+\frac{e}{2} \sigma_{\mu \nu} F^{\mu \nu} \tag{10.2.59}
\end{equation*}
$$

As in the case discussed in Sec. 10.2.3, it is convenient to perform a Wick's rotation by introducing, in addition to the Euclidean coordinates and fields discussed right before eq. (10.2.27), also a set of Euclidean $\gamma$ matrices $\gamma_{E}$ such that $\gamma_{i}=\gamma_{E, i}$ with $i=1,2$, 3, while $\gamma_{0}=i \gamma_{E, 4}$. As a result, $\not \partial=\gamma_{0} \partial_{0}-\vec{\gamma} \cdot \vec{\partial}=-\left(\gamma_{E, 4} \partial_{4}+\vec{\gamma} \cdot \vec{\partial}\right) \equiv-\not \phi_{E}$ with Euclidean metric. In addition, we introduce $\left[\gamma_{E, \mu}, \gamma_{E, \nu}\right] \equiv-2 i \sigma_{E, \mu \nu}$ with $\sigma_{E, 4 k}=-i \sigma_{0 k}$ and $\sigma_{E, l k}=\sigma_{l k}$, with $l, k \in\{1,2,3\}$, such that $F_{\mu \nu} \sigma^{\mu \nu}=F_{E, \mu \nu} \sigma_{E, \mu \nu}$. Using the chiral representation for the $\gamma$ matrices one has, in terms of Pauli matrices $\left\{\sigma_{l}\right\}_{l=1,2,3}$,

$$
\gamma_{0}=\left(\begin{array}{ll}
0 & 1  \tag{10.2.60}\\
1 & 0
\end{array}\right), \quad \gamma_{l}=\left(\begin{array}{cc}
0 & \sigma_{l} \\
-\sigma_{l} & 0
\end{array}\right), \quad \sigma_{0 l}=i\left(\begin{array}{cc}
-\sigma_{l} & 0 \\
0 & \sigma_{l}
\end{array}\right), \quad \sigma_{i j}=\epsilon_{i j k}\left(\begin{array}{cc}
\sigma_{k} & 0 \\
0 & \sigma_{k}
\end{array}\right),
$$

and therefore

$$
F_{E, \mu \nu} \sigma_{E, \mu \nu}=2\left(\begin{array}{cc}
\vec{\sigma} \cdot(-i \vec{E}+\vec{B}) & 0  \tag{10.2.61}\\
0 & \vec{\sigma} \cdot(i \vec{E}+\vec{B})
\end{array}\right)
$$

Now it is possible to express the quantum action $\Delta S_{Q}$ with the help of eqs. (10.2.57), (10.2.58), (10.2.59), and (10.2.28), as

$$
\begin{align*}
i \Delta S_{Q}[A] & =\frac{1}{2} \operatorname{Tr} \ln \left(\frac{-D_{E}^{2}+(e / 2) F_{E} \sigma_{E}+m^{2}}{\mu^{2}}\right) \\
& =-\frac{1}{2} \int_{0}^{\infty} \frac{d \beta}{\beta} e^{-\beta m^{2}} \operatorname{Tr}\left[e^{-\beta\left(-D_{E}^{2}\right)} e^{-(\beta e / 2) F_{E} \sigma_{E}}\right]+\text { const. } \tag{10.2.62}
\end{align*}
$$

The trace in the integrand involves a differential operator $e^{-\beta\left(-D_{E}^{2}\right)}$ acting on the coordinate space which is the same as the one discussed in Sec. 10.2.3 and which does not affect the spinorial structure; the operator $e^{-(\beta e / 2) F_{E} \sigma_{E}}$, instead, acts only on the spinorial structure while it does not affect the coordinate space. As a result, the trace factorizes into the traces over these two different spaces and the only new element in the analysis compared to the case of scalar QED is the calculation of $\operatorname{Tr}_{\mathrm{D}} e^{-(\beta e / 2) F_{E} \sigma_{E}}$. This can be easily done because of the simple structure of $F_{E} \sigma_{E}$ (see eq. (10.2.61)), which yields

$$
\operatorname{Tr} \exp \left(\begin{array}{cc}
\vec{\sigma} \cdot \vec{x} & 0  \tag{10.2.63}\\
0 & \vec{\sigma} \cdot \vec{y}
\end{array}\right)=\operatorname{Tr}\left(\begin{array}{cc}
e^{\vec{\sigma} \cdot \vec{x}} & 0 \\
0 & e^{\vec{\sigma} \cdot \vec{y}}
\end{array}\right)=2(\cosh |\vec{x}|+\cosh |\vec{y}|)
$$

where we used the facts that $e^{\vec{\sigma} \cdot \vec{x}}=\rrbracket \cosh |\vec{x}|-(\vec{x} \cdot \vec{\sigma} /|\vec{x}|) \sinh |\vec{x}|$ and $\operatorname{Tr}_{\text {Pauli }} \sigma_{l}=0$, with $\operatorname{Tr}_{\text {Pauli }} \rrbracket=2$. (In the previous relation $|\vec{x}|=\sqrt{\vec{x} \cdot \vec{x}}$.) Accordingly, using eq. (10.2.63) in eq. (10.2.61), we find

$$
\begin{align*}
& \operatorname{Tr}_{D} e^{-(\beta e / 2) F_{E} \sigma_{E}}=2[\cosh (\beta e|-i \vec{E}+\vec{B}|)+\cosh (\beta e|i \vec{E}+\vec{B}|)] \\
& \quad=4 \cosh \left(\beta e \frac{|-i \vec{E}+\vec{B}|+|i \vec{E}+\vec{B}|}{2}\right) \cosh \left(\beta e \frac{|-i \vec{E}+\vec{B}|-|i \vec{E}+\vec{B}|}{2}\right)  \tag{10.2.64}\\
& \quad=4 \cos \left(\beta e a_{+}\right) \cosh \left(\beta e a_{-}\right)
\end{align*}
$$

where $a_{ \pm}$were introduced in eq. (10.2.44). ${ }^{5}$
Accordingly, on the basis of eqs. (10.2.46) and (10.2.64), one can express eq. (10.2.62) as

$$
\begin{equation*}
\Delta S_{Q}[A]=-\frac{V T}{8 \pi^{2}} \int_{0}^{\infty} \frac{d \beta}{\beta} \frac{1}{(4 \pi \beta)^{d / 2-2}} e^{-\beta m^{2}} \frac{e^{2} a_{+} a_{-}}{\tanh \left(\beta e a_{-}\right) \tan \left(\beta e a_{+}\right)} \tag{10.2.65}
\end{equation*}
$$

As in the case of the analogous expression for the scalar vacuum, the last factor in the integrand behaves as $\simeq \beta^{-2}+\left(\vec{B}^{2}-\vec{E}^{2}\right) / 3+O\left(\beta^{2}\right)$ for $\beta \rightarrow 0$, the leading field-independent term of which is regularized by additional additive terms that we have omitted; the next term of the expansion, instead, is responsible for the emergence of a field-dependent logarithmic divergence (equivalently, of a dimensional pole in DR ) which can be cancelled by a suitable renormalization of the field. Repeating the analysis presented in Sec. 10.2.3 for the scalar case it is not difficult to recover in this way the field-strength renormalization constant of QED

$$
\begin{equation*}
Z_{3}=1-\frac{e^{2}}{(4 \pi)^{2}} \frac{8}{3 \epsilon}+O\left(e^{4}\right) \tag{10.2.66}
\end{equation*}
$$

[^71]

Figure 10.1: Vacuum decay rate $\Gamma$ per unit volume as a function of the strength $E$ of the electric field (in the absence of a magnetic field), for scalar QED (sQED, red lines) and QED (blue lines). The solid curves correspond to eqs. (10.2.56) and (10.2.68) for sQED and QED, respectively, while the dashed lines indicate the corresponding asymptotic values for $E \gg E_{\text {cr }}$, with $E_{\text {cr }}=2 m^{2} / e$.
as the reader can easily verify. Once these divergent terms have been properly subtracted from eq. (10.2.65), its renormalized expression can be used in order to calculate the vacuum decay rate $\Gamma$ (via eqs. (10.2.25) and (10.2.53)), which turns out to be given by the expression in eq. (10.2.55) corrected by an overall factor -2 (compare eq. (10.2.65) for $d=4$ to eq. (10.2.52)) times the residue of the factors $\cos \left(e u a_{+} / m^{2}\right) \cosh \left(e u a_{-} / m^{2}\right)$ at $u=u_{n}$ (see after eq. (10.2.52)) introduced by the trace over the spinorial structure; accordingly,

$$
\begin{equation*}
\Gamma=\frac{e^{2} a_{+} a_{-}}{4 \pi^{2}} \sum_{n=1}^{\infty} \frac{1}{n} \frac{e^{-n \pi m^{2} /\left(e a_{+}\right)}}{\tanh \left(n \pi a_{-} / a_{+}\right)} . \tag{10.2.67}
\end{equation*}
$$

As in the case of $\Gamma$ in eq. (10.2.55) one can easily verify that the presence of a sole magnetic field $\vec{B}$ is not sufficient to give a non-vanishing $\Gamma$. On the other hand a finite electric field $\vec{E}$, with $\vec{B}=0$, is responsible for a decay rate $\Gamma$ which can be easily determined according to the discussion reported after eq. (10.2.55) and which is given by

$$
\begin{equation*}
\Gamma=\frac{e^{2}|\vec{E}|^{2}}{4 \pi^{3}} \sum_{n=1}^{\infty} \frac{e^{-n \pi m^{2} /(e|\vec{E}|)}}{n^{2}} \tag{10.2.68}
\end{equation*}
$$

Figure 10.1 compares, as a function of $E / E_{\text {cr }}$ the vacuum decay rate $\Gamma$ per unit volume (normalized as $\Gamma /(e E)^{2}$ ) with $\vec{E} \neq 0$ and $\vec{B}=0$ of QED (blue lines, given by the equation above) with that one of scalar QED (red lines, see eq. (10.2.56)). Due to its non-perturbative nature, the Schwinger effect is practically absent as long as $E$ does not
exceed $\simeq E_{\text {cr }}$ (as it is clearly shown by the behavior of the solid curves in Fig. 10.1 for $E \lesssim E_{\text {cr }}$ ), which makes its observation a long-standing experimental challenge. Consider, for example, the typical electric field within an atom, i.e., $E_{a}=e / a_{0}^{2} \simeq 6 \cdot 10^{11} \mathrm{~V} / \mathrm{m}$ (where $a_{0} \simeq 5 \cdot 10^{-11} \mathrm{~m}$ is Bohr's radius): its ratio to the critical field $E_{\text {cr }}$ is $E_{a} / E_{\text {cr }} \simeq 3 \cdot 10^{-6}$ and eq. (10.2.68) would render the $\Gamma / m^{4} \simeq 10^{-10^{6}}$ : the pair creation rate $\gamma$ from the vacuum surrounding an atom of volume $a_{0}^{3}$ would therefore be $\gamma=\Gamma a_{0}^{3} \simeq 10^{-10^{6}} \mathrm{~s}^{-1}$ which results in a number of produced pairs that is surely negligible even for an observation time of the order of the age of the universe ( $\simeq 4 \cdot 10^{17} \mathrm{~s}$ ).

## Chapter 11

## Final Project: The Abelian Higgs Model ${ }^{\star}$

In this last chapter we study the abelian Higgs model, along the lines of ref. [10]. This chapter should be seen as a sort of long exercise in which many of the notions and techniques introduced in these notes (effective potential, background field method, ghosts, gauge-fixing, CS equations, $\beta$-functions and anomalous dimensions) are considered together.

The Lagrangian is

$$
\begin{equation*}
\mathcal{L}=-\frac{1}{4} F_{\mu \nu} F^{\mu \nu}+\left|D_{\mu} \Phi\right|^{2}-\frac{\lambda}{6}\left(\Phi^{\dagger} \Phi\right)^{2} \tag{11.0.1}
\end{equation*}
$$

where $D_{\mu} \Phi=\partial_{\mu} \Phi-i e A_{\mu} \Phi$. Our aim will be to understand the vacuum of this theory, namely whether the $U(1)$ gauge symmetry is broken or not, and the RG flows of the two couplings $e$ and $\lambda$. First of all, we have to gauge fix the theory. Since we want to study the effective potential as a function of the VEV of $\Phi$, it is convenient to use a generalized $\xi$-gauge which is valid for any value of $\langle\Phi\rangle$ :

$$
\begin{equation*}
\mathcal{L}_{\text {g.f. }}=-\frac{1}{2 \xi}\left[\partial_{\mu} A^{\mu}+i e \xi\left(\phi^{\dagger} \phi_{0}-\phi \phi_{0}^{\dagger}\right)\right]^{2} \tag{11.0.2}
\end{equation*}
$$

where $\Phi=\phi_{0}+\phi$, with $\phi_{0}$ the VEV of $\Phi$ and $\phi$ its quantum fluctuation. It is straightforward to verify that the quadratic mixing terms between $\phi$ and $A_{\mu}$ vanish when $\mathcal{L}_{\text {g.f. }}$ is added to the Lagrangian (11.0.1). Even if the local symmetry is abelian, the ghosts do not decouple in the $\xi$-gauge we have chosen. The ghost Lagrangian, as usual, is derived by taking the infinitesimal variation of $\mathcal{L}_{\text {g.f. }}$ with respect to a $U(1)$ transformation. One gets

$$
\begin{equation*}
\mathcal{L}_{\text {ghosts }}=\partial_{\mu} \omega^{\star} \partial^{\mu} \omega-e^{2} \xi \omega^{\star} \omega\left[2\left|\phi_{0}\right|^{2}+\left(\phi^{\dagger} \phi_{0}+\phi_{0}^{\dagger} \phi\right)\right] . \tag{11.0.3}
\end{equation*}
$$

In the Landau gauge $\partial_{\mu} A^{\mu}=0$, reached for $\xi \rightarrow 0$, the ghosts are free and decouple, whereas in the unitary gauge $\xi \rightarrow \infty$ they are infinitely massive and decouple again. We will not fix in the following a specific value of $\xi$, so that the ghosts should be taken into account. The total Lagrangian is

$$
\begin{equation*}
\mathcal{L}_{t o t}=\mathcal{L}+\mathcal{L}_{\text {g.f. }}+\mathcal{L}_{\text {ghosts }} . \tag{11.0.4}
\end{equation*}
$$

### 11.1 One-loop Effective Potential*

The 1-loop effective potential is completely determined by the terms in $\mathcal{L}_{\text {tot }}$ quadratic in the field fluctuations. In momentum space, we have

$$
\begin{equation*}
\mathcal{L}_{t o t, q u a d}(p)=-\frac{1}{2} A^{\mu}(-p) A^{\nu}(p) \mathcal{L}_{\mu \nu}^{(A)}(p)+\omega^{\star}(-p) \mathcal{L}^{(\omega)}(p) \omega(p)+\frac{1}{2} \phi_{i}(-p) \mathcal{L}_{i j}^{(\phi)}(p) \phi_{j}(p) \tag{11.1.1}
\end{equation*}
$$

where $\phi=\left(\phi_{1}+i \phi_{2}\right) / \sqrt{2}$, and
$\mathcal{L}_{\mu \nu}^{(A)}(p)=\eta_{\mu \nu}\left(p^{2}-2 e^{2}\left|\phi_{0}\right|^{2}\right)-\left(1-\frac{1}{\xi}\right) p_{\mu} p_{\nu}$,
$\mathcal{L}^{(\omega)}(p)=p^{2}-2 \xi e^{2}\left|\phi_{0}\right|^{2}$,
$\mathcal{L}_{i j}^{(\phi)}(p)=\left(\begin{array}{cc}p^{2}-\left|\phi_{0}\right|^{2}\left(e^{2} \xi+\frac{2 \lambda}{3}\right)+\left(\phi_{0}^{2}+\phi_{0}^{\star 2}\right)\left(\frac{e^{2} \xi}{2}-\frac{\lambda}{6}\right) & i\left(\phi_{0}^{2}-\phi_{0}^{\star 2}\right)\left(\frac{e^{2} \xi}{2}-\frac{\lambda}{6}\right) \\ i\left(\phi_{0}^{2}-\phi_{0}^{\star 2}\right)\left(\frac{e^{2} \xi}{2}-\frac{\lambda}{6}\right) & p^{2}-\left|\phi_{0}\right|^{2}\left(e^{2} \xi+\frac{2 \lambda}{3}\right)-\left(\phi_{0}^{2}+\phi_{0}^{\star 2}\right)\left(\frac{e^{2} \xi}{2}-\frac{\lambda}{6}\right)\end{array}\right)$.
Modulo irrelevant factors,

$$
\begin{align*}
\operatorname{det} \mathcal{L}_{\mu \nu}^{(A)}(p) & =\left(p^{2}-2 e^{2}\left|\phi_{0}\right|^{2}\right)^{3}\left(p^{2}-2 e^{2} \xi\left|\phi_{0}\right|^{2}\right) \\
\operatorname{det} \mathcal{L}_{i j}^{(\phi)}(p) & =\left(p^{2}-\lambda\left|\phi_{0}\right|^{2}\right)\left[p^{2}-\left|\phi_{0}\right|^{2}\left(2 e^{2} \xi+\frac{\lambda}{3}\right)\right] \tag{11.1.3}
\end{align*}
$$

Summing over all contributions (gauge, ghosts and scalar fields), we get

$$
\begin{align*}
V_{1-\text { loop }}(\rho)= & \frac{1}{2} \int \frac{d^{4} p_{E}}{(2 \pi)^{4}}\left[3 \log \left(p_{E}^{2}+e^{2} \rho^{2}\right)-\log \left(p_{E}^{2}+\xi e^{2} \rho^{2}\right)\right. \\
& \left.+\log \left(p_{E}^{2}+\frac{\lambda}{2} \rho^{2}\right)+\log \left(p_{E}^{2}+\left(\xi e^{2}+\frac{\lambda}{6}\right) \rho^{2}\right)\right] \tag{11.1.4}
\end{align*}
$$

where $\rho^{2}=2\left|\phi_{0}\right|^{2}$. We renormalize $V_{1-\text { loop }}$ by demanding that

$$
\begin{equation*}
\left.\frac{d^{2} V_{1-\text { loop }}}{d \rho^{2}}\right|_{\rho=0}=0,\left.\quad \frac{d^{4} V_{1-l o o p}}{d \rho^{4}}\right|_{\rho=\mu}=0 \tag{11.1.5}
\end{equation*}
$$

After some simple algebra, we obtain

$$
\begin{equation*}
V_{\text {eff }}(\rho)=V_{\text {tree }}(\rho)+V_{1-\text { loop }}(\rho)=\frac{\lambda}{4!} \rho^{4}+\frac{\rho^{4}}{64 \pi^{2}}\left(3 e^{4}+\frac{5}{18} \lambda^{2}+\frac{1}{3} \xi \lambda e^{2}\right)\left(\log \frac{\rho^{2}}{\mu^{2}}-\frac{25}{6}\right) . \tag{11.1.6}
\end{equation*}
$$



Figure 11.1: One-loop graph leading to a $\lambda \rho^{4}$ coupling.

Let us study the minima of $V_{e f f}$, assuming that $\lambda \sim e^{4}$, so that at leading order we can neglect the $\lambda^{2}$ and $\xi \lambda e^{2}$ terms in eq. (11.1.6). It is important to emphasize here that at tree-level (or, alternatively, at any given energy scale) we can assume any relation we like of the form $\lambda \sim e^{n}$ for any $n$, but such relations at the quantum level cannot generally hold at any energy scale for any $n$. Indeed, radiative corrections will nevertheless generate the $\lambda \rho^{4}$ coupling in the theory. The leading one-loop correction arises from virtual photons as illustrated in fig. 11.1. Being this correction of $\mathcal{O}\left(e^{4}\right)$, we see that $\lambda \sim e^{n}$, with $n \leq 4$, are the only radiatively stable assumptions we can make. The extrema of $V_{e f f}$ are

$$
\begin{equation*}
\frac{d V_{e f f}}{d \rho}=\rho^{3}\left(\frac{\lambda}{6}+\frac{e^{4}}{16 \pi^{2}}\left(3 \log \frac{\rho^{2}}{\mu^{2}}-11\right)\right)=0 \tag{11.1.7}
\end{equation*}
$$

Taking $\mu=\langle\rho\rangle$ in eq. (11.1.7), we get $\rho=0$ and

$$
\begin{equation*}
\lambda(\langle\rho\rangle)=\frac{33}{8 \pi^{2}} e^{4}(\langle\rho\rangle) \tag{11.1.8}
\end{equation*}
$$

Equation (11.1.8) is an instance of dimensional transmutation: we have traded the VEV of $\rho$ for the coupling $\lambda$. Plugging back in $V_{\text {eff }}$ gives

$$
\begin{equation*}
V_{e f f} \simeq \frac{3 e^{4}}{64 \pi^{2}} \rho^{4}\left(\log \frac{\rho^{2}}{\langle\rho\rangle^{2}}-\frac{1}{2}\right) \tag{11.1.9}
\end{equation*}
$$

The extremum (11.1.8) is a minimum. The photon and scalar masses are

$$
\begin{equation*}
m_{\gamma}^{2}=e^{2}\langle\rho\rangle^{2}, \quad m_{\rho}^{2}=\frac{3 e^{4}}{8 \pi^{2}}\langle\rho\rangle^{4} \tag{11.1.10}
\end{equation*}
$$

We conclude that in this theory a dynamical spontaneous symmetry breaking of the $U(1)$ gauge symmetry can occur. In order to firmly establish that, we have to compute the RG evolutions of $\lambda$ and of the charge $e$ to check the existence of an energy scale $\langle\rho\rangle$ where eq. (11.1.8) is valid. This computation will be the subjects of the following sections.

### 11.2 The Quantum Effective Action ${ }^{\star}$

An instructive way of computing the $\beta$-functions of $\lambda$ and $e$, as well as the anomalous dimensions of $A_{\mu}$ and $\phi$, makes use of a functional form of the CS equations (5.4.3), which reads:

$$
\begin{equation*}
\mu \frac{d \Gamma}{d \mu}=\left(\mu \frac{\partial}{\partial \mu}+\beta \frac{\partial}{\partial \lambda}-\gamma \int d^{4} x \phi(x) \frac{\delta}{\delta \phi(x)}\right) \Gamma(\phi)=0, \tag{11.2.1}
\end{equation*}
$$

where $\Gamma(\phi)$ is the quantum effective action. For a single real scalar field, keeping up to two derivative terms, the latter reads

$$
\begin{equation*}
\Gamma(\phi)=\int d^{4} x\left(\frac{1}{2} Z(\phi)\left(\partial_{\mu} \phi\right)^{2}-V_{e f f}(\phi)\right), \tag{11.2.2}
\end{equation*}
$$

where $V_{\text {eff }}$ is the Coleman-Weinberg potential and $Z$ is the radiative correction to the kinetic term. Once $\Gamma$ is known, eq. (11.2.1) gives us $\beta$ and $\gamma$. It is straightforward to see that eq. (11.2.1) encodes all eqs. (5.4.3) for any $n$, by recalling that

$$
\begin{equation*}
\Gamma(\phi)=\sum_{n=0}^{\infty} \frac{1}{n!} \int d^{4} x_{1} \ldots d^{4} x_{n} \Gamma^{(n)}\left(x_{1}, \ldots, x_{n}\right) \phi\left(x_{1}\right) \ldots \phi\left(x_{n}\right) . \tag{11.2.3}
\end{equation*}
$$

In the case at hand, with two fields and two couplings, eq. (11.2.1) generalizes to

$$
\begin{align*}
& \left(\mu \frac{\partial}{\partial \mu}+\beta_{\lambda} \frac{\partial}{\partial \lambda}+\beta_{e} \frac{\partial}{\partial e}-\gamma_{A} \int d^{4} x A_{\mu, 0}(x) \frac{\delta}{\delta A_{\mu, 0}(x)}\right. \\
& \left.-\gamma_{\phi} \int d^{4} x\left(\phi_{0}(x) \frac{\delta}{\delta \phi_{0}(x)}+\phi_{0}^{\dagger}(x) \frac{\delta}{\delta \phi_{0}^{\dagger}(x)}\right)\right) \Gamma\left(\phi_{0}, A_{\mu, 0}\right)=0 \tag{11.2.4}
\end{align*}
$$

where $\phi_{0}, \phi_{0}^{\dagger}$ and $A_{\mu, 0}$ are the background field configurations. Invariance under the background $U(1)$ gauge invariance implies that the lowest dimensional operators appearing in $\Gamma$ are of the form

$$
\begin{equation*}
\Gamma\left(\phi_{0}, A_{\mu, 0}\right)=\int d^{4} x\left[-\frac{1}{4} H(\rho) F_{\mu \nu, 0}^{2}+Z(\rho)\left|D_{\mu} \phi_{0}\right|^{2}-V_{e f f}(\rho)\right] . \tag{11.2.5}
\end{equation*}
$$

The effective potential $V_{e f f}$ has been already computed and is given by eq. (11.1.6). We have then to determine $H$ and $Z$ only. This can be done by decomposing the gauge field as well in terms of background and fluctuation fields: $A_{\mu} \rightarrow A_{\mu}^{0}+A_{\mu}$.

Let us start by computing $H$, in which case we can take $\phi_{0}=$ constant. The relevant interaction terms are

$$
\begin{equation*}
\mathcal{L} \supset i e A_{0}^{\mu}\left(\phi^{\dagger} \partial_{\mu} \phi-\partial_{\mu} \phi^{\dagger} \phi\right), \tag{11.2.6}
\end{equation*}
$$

and $H$ is determined by the contraction of the two scalar currents. The contractions of the form $\langle\phi \phi\rangle$ or $\left\langle\phi^{\dagger} \phi^{\dagger}\right\rangle$, although non-vanishing for $\phi_{0} \neq 0$ due to mass terms of the form $\phi^{2}$
and $\left(\phi^{\dagger}\right)^{2}$, give only rise to irrelevant divergent contact terms. In DR, for instance, they all trivially vanish. The only relevant contraction is the usual one, of the form $\left\langle\phi^{\dagger} \phi\right\rangle$. The computation essentially boils down to the one-loop photon vacuum polarization in scalar QED. Two diagrams contribute. The first reads

where, according to eq. (11.1.2),

$$
\begin{equation*}
m^{2}=m^{2}(\rho)=\rho^{2}\left(\frac{2 \lambda}{3}+\frac{e^{2} \xi}{2}\right) \tag{11.2.8}
\end{equation*}
$$

By performing the usual manipulations (introduce the Feynman parameter $x$, shift $q \rightarrow$ $q-x p$ and Wick rotate to euclidean momentum), we get

$$
\begin{align*}
i \Pi_{\mu \nu}^{1}(p) & =(i e)^{2} \mu^{\epsilon} \int \frac{d^{d} q}{(2 \pi)^{d}} \frac{i^{2}(p+2 q)_{\mu}(p+2 q)_{\nu}}{\left(q^{2}-m^{2}\right)\left[(p+q)^{2}-m^{2}\right]}=\frac{i e^{2} \mu^{\epsilon}}{(4 \pi)^{d / 2}} \int_{0}^{1} d x\left[m^{2}-p^{2} x(1-x)\right]^{\frac{d-4}{2}} \\
& \times \Gamma\left(\frac{4-d}{2}\right)\left(p_{\mu} p_{\nu}(1-2 x)^{2}-\frac{4}{2-d} \eta_{\mu \nu}\left[m^{2}-p^{2} x(1-x)\right]\right) \tag{11.2.9}
\end{align*}
$$

The second diagram is a tadpole, that does not depend on the external momentum $p$. It can be cast in a form close to eq. (11.2.9) by multiplying and dividing it by $(p+q)^{2}-m^{2}$. Using the same manipulations as before, we get


$$
\times \int_{0}^{1} d x\left[m^{2}-p^{2} x(1-x)\right]^{\frac{d-4}{2}} \Gamma\left(\frac{4-d}{2}\right)\left((1-x)^{2} p^{2}-m^{2}-\frac{d}{2-d}\left[m^{2}-p^{2} x(1-x)\right]\right) .
$$

Summing the two contributions, we have

$$
\begin{align*}
i \Pi_{\mu \nu}=i \Pi_{\mu \nu}^{1}+i \Pi_{\mu \nu}^{2} & =\frac{i e^{2}}{24 \pi^{2}}\left(p_{\mu} p_{\nu}-\eta_{\mu \nu} p^{2}\right)\left(\frac{1}{\epsilon}+\text { const. }\right)-\frac{i e^{2}}{16 \pi^{2}} \int_{0}^{1} d x \log \frac{m^{2}-p^{2} x(1-x)}{\mu^{2}} \\
& \times\left(p_{\mu} p_{\nu}(1-2 x)^{2}-\eta_{\mu \nu} p^{2}\left(4 x^{2}-6 x+2\right)\right) \tag{11.2.11}
\end{align*}
$$

We are interested in computing $H$, which is the coefficient of the $F^{2}$ term, quadratic in the external momentum $p$. The last term in eq. (11.2.11) is already quadratic in $p$, so we
can safely neglect the $p^{2}$ term inside the log and keep only the mass term $m^{2}$. In this way we get

$$
\begin{equation*}
i \Pi_{\mu \nu}=\frac{i e^{2}}{24 \pi^{2}}\left(p_{\mu} p_{\nu}-\eta_{\mu \nu} p^{2}\right)\left(\frac{1}{\epsilon}+\text { const. }-\log (\rho / \mu)\right) \tag{11.2.12}
\end{equation*}
$$

where we have reabsorbed in the arbitrary constant the $\rho$-independent factors coming from eq. (11.2.8). As expected, $\Pi_{\mu \nu}$ is transverse. We now perform a non-minimal subtraction, i.e. we add the counter-term ${ }^{1} i(Z-1)\left(p_{\mu} p_{\nu}-\eta_{\mu \nu} p^{2}\right)$ and require that

$$
\begin{equation*}
i \Pi_{\mu \nu}+i(Z-1)\left(p_{\mu} p_{\nu}-\eta_{\mu \nu} p^{2}\right)=0, \text { at } \rho=\mu \tag{11.2.13}
\end{equation*}
$$

The renormalized one-loop photon vacuum polarization is then

$$
\begin{equation*}
i \Pi_{\mu \nu}^{R}=\frac{i e^{2}}{24 \pi^{2}}\left(\eta_{\mu \nu} p^{2}-p_{\mu} p_{\nu}\right) \log \frac{\rho}{\mu} \equiv i A p^{2} \mathcal{P}_{\mu \nu} \tag{11.2.14}
\end{equation*}
$$

where $A=e^{2} /\left(24 \pi^{2}\right) \log \phi / \mu$ and we have defined the projector

$$
\begin{equation*}
\mathcal{P}_{\mu \nu}=\eta_{\mu \nu}-\frac{p_{\mu} p_{\nu}}{p^{2}} \tag{11.2.15}
\end{equation*}
$$

The tree-level photon propagator is

$$
\begin{equation*}
G_{\mu \nu}(p)=-\frac{i}{p^{2}}\left(\mathcal{P}_{\mu \nu}+\xi \frac{p_{\mu} p_{\nu}}{p^{2}}\right) . \tag{11.2.16}
\end{equation*}
$$

Iterating the one-loop correction (11.2.14) in the tree-level expression (11.2.16) we get

$$
\begin{equation*}
G_{\mu \nu}(p) \rightarrow \frac{-i \mathcal{P}_{\mu \nu}}{p^{2}(1-A)}-\frac{i}{p^{2}} \frac{\xi p_{\mu} p_{\nu}}{p^{2}} \tag{11.2.17}
\end{equation*}
$$

from which we see that

$$
\begin{equation*}
-\frac{1}{4} F_{\mu \nu, 0}^{2} \rightarrow-\frac{1}{4}(1-A) F_{\mu \nu, 0}^{2} \tag{11.2.18}
\end{equation*}
$$

finally giving the desired $H(\rho)$ :

$$
\begin{equation*}
H(\rho)=1-A(\rho)=1-\frac{e^{2}}{24 \pi^{2}} \log \frac{\rho}{\mu} . \tag{11.2.19}
\end{equation*}
$$

Let us now determine $Z$, the wave-function renormalization of $\phi_{0}$. In this case we can set $A_{\mu, 0}=0$ but of course $\phi_{0}$ can no longer be taken constant. The tadpole graph given by the quartic scalar interaction induces only a mass renormalization, so we need only to consider the contribution induced by the gauge interactions. These are of the form $A_{\mu} \phi \phi_{0}$, given by

$$
\begin{equation*}
\mathcal{L} \supset 2 i e A^{\mu}\left(\phi^{\dagger} \partial_{\mu} \phi_{0}-\partial_{\mu} \phi_{0}^{\dagger} \phi\right) \tag{11.2.20}
\end{equation*}
$$

[^72]We have

$$
\begin{equation*}
\rightarrow \underset{p+q}{\rightarrow} \quad \overbrace{0}^{\sim}=i \Sigma(p)=(2 i e)^{2} \mu^{\epsilon} p^{\mu} p^{\nu} \int \frac{d^{d} q}{(2 \pi)^{d}} G_{\mu \nu}(q) \frac{i}{(p+q)^{2}-m^{2}}, \tag{11.2.21}
\end{equation*}
$$

where $G_{\mu \nu}(q)$ is the photon propagator in the generically broken phase, as given by in eq.(8.4.13), with $\mu_{\gamma}^{2}=e^{2} \rho^{2} .{ }^{2}$ For $\xi \neq 1$, up to three denominators appear in eq. (11.2.21) and the computation is a bit involved. We can greatly simplify it by noticing that i) the $\log \mu$ (and hence the $\log \rho$ ) terms we are looking for can be detected by finding the residues of the appropriate $1 / q^{4}$ poles in the integrand, as explained in section 3.4 and ii) we can set $p=0$ in the scalar propagator since we already have two powers of the external momentum coming from the vertices. After Wick rotating to euclidean momenta and using $S O(d)$ invariance of the integrand, we get

$$
\begin{align*}
i \Sigma(p) & =-4 i e^{2} \mu^{\epsilon} p^{2} \int \frac{d^{d} q_{E}}{(2 \pi)^{d}}\left(\frac{1}{q_{E}^{2}+\mu_{\gamma}^{2}}-\frac{(1-\xi) q_{E}^{2}}{d\left(q_{E}^{2}+\xi \mu_{\gamma}^{2}\right)\left(q_{E}^{2}+\mu_{\gamma}^{2}\right)}\right) \frac{1}{q_{E}^{2}+m^{2}} \\
& =\left(-4 i e^{2}\right) p^{2} \frac{1}{8 \pi^{2}}\left(\frac{1}{\epsilon}+\log \mu+\ldots\right) \frac{\xi+3}{4} \tag{11.2.22}
\end{align*}
$$

where the $\ldots$ include the $\rho$-dependent terms we are looking for plus additional finite pieces. The renormalized scalar two-point function reads then

$$
\begin{equation*}
i \Sigma^{R}(p)=\frac{i e^{2}}{8 \pi^{2}}(3+\xi) p^{2} \log \frac{\rho}{\mu} \equiv i p^{2} B \tag{11.2.23}
\end{equation*}
$$

where, as before, we demand that $\Sigma^{R}(p)=0$ when $\rho=\mu$. Iterating the one-loop correction in the tree-level scalar propagator, as in the photon case, gives

$$
\begin{equation*}
\frac{i}{p^{2}} \rightarrow \frac{i}{p^{2}(1+B)}, \tag{11.2.24}
\end{equation*}
$$

from which we extract $Z(\rho)$ :

$$
\begin{equation*}
Z(\rho)=1+B(\rho)=1+\frac{e^{2}}{8 \pi^{2}}(3+\xi) \log \frac{\rho}{\mu} \tag{11.2.25}
\end{equation*}
$$

The form of the effective action (11.2.5) is finally determined and we can proceed to use it to compute the RG evolution of the couplings $e$ and $\lambda$. It is worth to emphasize that the effective action (11.2.5) should be gauge invariant, but on the contrary it seems gaugedependent, since $\xi$ enters in both $Z(\rho)$ and in $V_{\text {eff }}(\rho)$. This apparent paradox is explained by noticing that $\phi_{0}$ is not a canonically normalized field and the rescaling

$$
\begin{equation*}
\phi_{0} \rightarrow \frac{1}{\sqrt{Z(\rho)}} \phi_{0} \tag{11.2.26}
\end{equation*}
$$

[^73]is needed. The potential (11.1.6) is then rescaled by a factor $1 / Z(\rho)^{2}$ and we have
\[

$$
\begin{align*}
\frac{1}{Z(\rho)^{2}} V_{e f f}(\rho) & =\frac{\lambda}{4!} \rho^{4}\left(1-\frac{e^{2}}{4 \pi^{2}}(3+\xi) \log \frac{\rho}{\mu}\right)+\frac{\rho^{4}}{32 \pi^{2}}\left(3 e^{4}+\frac{5}{18} \lambda^{2}+\frac{1}{3} \xi \lambda e^{2}\right) \log \frac{\rho}{\mu} \\
& =\frac{\lambda}{4!} \rho^{4}+\frac{\rho^{4}}{32 \pi^{2}}\left(3 e^{4}+\frac{5}{18} \lambda^{2}-e^{2} \lambda\right) \log \frac{\rho}{\mu} \tag{11.2.27}
\end{align*}
$$
\]

and the $\xi$-dependence cancels. In eq. (11.2.27) we have focused on the log terms only, since the constant term is scheme-dependent and we have not been careful in systematically using a given scheme.

### 11.3 RG equations and Their Solutions ${ }^{\star}$

For a single real scalar field, the functional RG equation (11.2.1) splits into two ordinary differential equations for $Z$ and $V$ :

$$
\begin{align*}
\left(\mu \frac{\partial}{\partial \mu}+\beta \frac{\partial}{\partial \lambda}-\gamma \phi \frac{\partial}{\partial \phi}-2 \gamma\right) Z & =0 \\
\left(\mu \frac{\partial}{\partial \mu}+\beta \frac{\partial}{\partial \lambda}-\gamma \phi \frac{\partial}{\partial \phi}\right) V_{e f f} & =0 \tag{11.3.1}
\end{align*}
$$

In dealing with quartic potentials, like in our case, it is actually convenient to write an RG equation for $V^{(4)} \equiv \partial^{4} V_{\text {eff }} / \partial \phi^{4}$ rather than $V_{\text {eff }}$ itself. Given that $\partial_{\phi}^{4}\left(\gamma \phi \partial_{\phi}\right)=$ $4 \gamma \partial_{\phi}^{4}+\gamma \phi \partial_{\phi} \partial_{\phi}^{4}$, we have

$$
\begin{equation*}
\left(\mu \frac{\partial}{\partial \mu}+\beta \frac{\partial}{\partial \lambda}-4 \gamma-\gamma \phi \frac{\partial}{\partial \phi}\right) V^{(4)}=0 \tag{11.3.2}
\end{equation*}
$$

The RG equation (11.3.2) can further be simplified by noting that $V^{(4)}$ depends on $\phi$ only through the dimensionless combination $\phi / \mu$. Then $\phi \partial_{\phi} V^{(4)}=-\mu \partial_{\mu} V^{(4)}=\partial_{t} V^{(4)}$, where $t=\log \phi / \mu$ and thus we get

$$
\begin{equation*}
\left(-\frac{\partial}{\partial t}+\bar{\beta} \frac{\partial}{\partial \lambda}-4 \bar{\gamma}\right) V^{(4)}=0 \tag{11.3.3}
\end{equation*}
$$

where

$$
\begin{equation*}
\bar{\beta}=\frac{\beta}{1+\gamma}, \quad \bar{\gamma}=\frac{\gamma}{1+\gamma} \tag{11.3.4}
\end{equation*}
$$

Using the same manipulations, the RG equation for $Z$ becomes

$$
\begin{equation*}
\left(-\frac{\partial}{\partial t}+\bar{\beta} \frac{\partial}{\partial \lambda}-2 \bar{\gamma}\right) Z=0 \tag{11.3.5}
\end{equation*}
$$

In the situation at hand, the functional RG equation is given by eq. (11.2.4). It is clear that there can be no cancellations between the three terms (11.2.5) appearing in $\Gamma$, so
eq. (11.2.4) splits into three independent equations. Let us first focus on the scalar kinetic term $\left|D \phi_{0}\right|^{2}$ :

$$
\begin{align*}
& \left|D_{\mu} \phi_{0}\right|^{2}\left(-\frac{\partial}{\partial t}+\beta_{e} \frac{\partial}{\partial e}-\gamma_{\phi} \rho \partial_{\rho}\right) Z+Z\left(\beta_{e} \frac{\partial}{\partial e}\left|D_{\mu} \phi_{0}\right|^{2}-\gamma_{A} \int d^{4} x A_{\nu, 0}(x) \frac{\delta\left|D_{\mu} \phi_{0}\right|^{2}}{\delta A_{\nu, 0}(x)}\right. \\
& \left.-\gamma_{\phi} \int d^{4} x\left(\phi_{0}(x) \frac{\delta}{\delta \phi_{0}(x)}+\phi_{0}^{\dagger}(x) \frac{\delta}{\delta \phi_{0}^{\dagger}(x)}\right)\right)\left|D_{\mu} \phi_{0}\right|^{2}=0 \tag{11.3.6}
\end{align*}
$$

where we have used the fact that $Z$ depends on constant $\rho$ only, with $\phi \partial_{\phi}+\phi^{\dagger} \partial_{\phi^{\dagger}}=\rho \partial_{\rho}$. The term in the second line of eq. (11.3.6) equals $-2 \gamma_{\phi}\left|D_{\mu} \phi_{0}\right|^{2}$, while the last two terms in the first line of eq. (11.3.6) gives rise to a different operator. As such, two independent equations arise from eq. (11.3.6). Requiring the vanishing of the coefficient proportional to $\left|D_{\mu} \phi_{0}\right|^{2}$ gives

$$
\begin{equation*}
\left(-\frac{\partial}{\partial t}+\bar{\beta}_{e} \frac{\partial}{\partial e}-2 \bar{\gamma}_{\phi}\right) Z=0 \tag{11.3.7}
\end{equation*}
$$

with

$$
\begin{equation*}
\bar{\beta}_{e}=\frac{\beta_{e}}{1+\gamma_{\phi}}, \quad \bar{\gamma}_{\phi}=\frac{\gamma_{\phi}}{1+\gamma_{\phi}} \tag{11.3.8}
\end{equation*}
$$

It is straightforward to see that

$$
\begin{equation*}
\int d^{4} x A_{\nu, 0}(x) \frac{\delta\left|D_{\mu} \phi_{0}\right|^{2}}{\delta A_{\nu, 0}(x)}=e \frac{\partial}{\partial e}\left|D_{\mu} \phi_{0}\right|^{2} \tag{11.3.9}
\end{equation*}
$$

and hence the vanishing of the coefficient multiplying $\partial\left|D_{\mu} \phi_{0}\right|^{2} / \partial e$ gives

$$
\begin{equation*}
\beta_{e}=e \gamma_{A} \tag{11.3.10}
\end{equation*}
$$

Consider now the gauge kinetic term $F_{\mu \nu, 0}^{2}$. We get

$$
\begin{equation*}
-\frac{1}{4} F_{\mu \nu, 0}^{2}\left(-\frac{\partial}{\partial t}+\beta_{e} \frac{\partial}{\partial e}-\gamma_{\phi} \rho \partial_{\rho}\right) H-2 \gamma_{A} H\left(-\frac{1}{4} F_{\mu \nu, 0}^{2}\right)=0, \tag{11.3.11}
\end{equation*}
$$

that gives rise to

$$
\begin{equation*}
\left(-\frac{\partial}{\partial t}+\bar{\beta}_{e} \frac{\partial}{\partial e}-2 \bar{\gamma}_{A}\right) H=0, \quad \bar{\gamma}_{A}=\frac{\gamma_{A}}{1+\gamma_{\phi}} \tag{11.3.12}
\end{equation*}
$$

Finally we have the potential term. As explained before, we write an RG equation for $V^{(4)}$ rather than $V_{e f f}$, which is the obvious generalization of eq. (11.3.3):

$$
\begin{equation*}
\left(-\frac{\partial}{\partial t}+\bar{\beta}_{\lambda} \frac{\partial}{\partial \lambda}+\bar{\beta}_{e} \frac{\partial}{\partial e}-4 \bar{\gamma}_{\phi}\right) V^{(4)}=0 . \tag{11.3.13}
\end{equation*}
$$

Equations (11.3.7), (11.3.10), (11.3.12) and (11.3.13) are enough to determine $\beta_{e}, \beta_{\lambda}, \gamma_{A}$ and $\gamma_{\phi}$. Let us recall below the explicit form of $Z, H$ and $V^{(4)}$, the latter computed from
eq. (11.1.6):

$$
\begin{align*}
Z(t) & =1+\frac{e^{2}}{8 \pi^{2}}(\xi+3) t \\
H(t) & =1-\frac{e^{2}}{24 \pi^{2}} t  \tag{11.3.14}\\
V^{(4)}(t) & =\lambda+\frac{1}{4 \pi^{2}}\left(9 e^{4}+\frac{5}{6} \lambda^{2}+\xi \lambda e^{2}\right) t
\end{align*}
$$

One immediately gets from eqs. (11.3.7), (11.3.10) and (11.3.12)

$$
\begin{align*}
\gamma_{\phi} & =\bar{\gamma}_{\phi}+\mathcal{O}\left(e^{4}\right)=-\frac{e^{2}}{16 \pi^{2}}(\xi+3)+\mathcal{O}\left(e^{4}\right) \\
\gamma_{A} & =\bar{\gamma}_{A}+\mathcal{O}\left(e^{4}\right)=\frac{e^{2}}{48 \pi^{2}}+\mathcal{O}\left(e^{4}\right) \\
\beta_{e} & =\bar{\beta}_{e}+\mathcal{O}\left(e^{5}\right)=\frac{e^{3}}{48 \pi^{2}}+\mathcal{O}\left(e^{5}\right) \tag{11.3.15}
\end{align*}
$$

Plugging the values (11.3.15) in eq. (11.3.13) allows us to determine $\beta_{\lambda}$ :

$$
\begin{equation*}
\beta_{\lambda}=\bar{\beta}_{\lambda}+\mathcal{O}\left(e^{6}, e^{4} \lambda, e^{2} \lambda^{2}\right)=\frac{1}{4 \pi^{2}}\left(9 e^{4}+\frac{5}{6} \lambda^{2}-3 \lambda e^{2}\right)+\mathcal{O}\left(e^{6}, e^{4} \lambda, e^{2} \lambda^{2}\right) \tag{11.3.16}
\end{equation*}
$$

Notice how all $\xi$-dependent factors have cancelled in $\beta_{\lambda}$ as it should be, being the latter gauge invariant, like $\beta_{e}$ (and $\gamma_{A}$ ). The scalar field anomalous dimension $\gamma_{\phi}$, instead, does depend on $\xi$. This is expected since $\phi$ changes by a phase under a gauge transformation and at the quantum level there is no gauge invariant notion of $\gamma_{\phi}$.

The RG flow of $e$ is easily computed from $\beta_{e}$. We get

$$
\begin{equation*}
e^{2}(t)=\frac{e_{0}^{2}}{1-\frac{e_{0}^{2}}{24 \pi^{2}} t} \tag{11.3.17}
\end{equation*}
$$

The RG flow of $\lambda$ requires some more work. It is convenient to define $R(t)=\lambda(t) / e^{2}(t)$ and write an RG equation for $R$. One gets

$$
\begin{equation*}
e^{2}(t) \dot{R}(t)=\frac{e^{4}(t)}{4 \pi^{2}}\left(\frac{5}{6} R^{2}(t)-\frac{19}{6} R(t)+9\right) \tag{11.3.18}
\end{equation*}
$$

which is further simplified by considering $R=R\left(e^{2}\right)$, so that

$$
\begin{equation*}
\dot{R}=\frac{d R}{d e^{2}} 2 e \dot{e}=\frac{e^{4}}{24 \pi^{2}} \frac{d R}{d e^{2}} . \tag{11.3.19}
\end{equation*}
$$

Our desired final equation reads

$$
\begin{equation*}
e^{2} \frac{d R\left(e^{2}\right)}{d e^{2}}=5 R^{2}\left(e^{2}\right)-19 R\left(e^{2}\right)+54 \tag{11.3.20}
\end{equation*}
$$



Figure 11.2: Comparison between the RG behaviour of $e^{2}(t)$ (red line) and $\lambda(t)$ (blue line) over 100 orders of magnitude. We have taken $e^{2}(0)=\lambda(0)=1 / 10$.
whose solution is

$$
\begin{equation*}
R\left(e^{2}\right)=\frac{1}{10}\left(19+\sqrt{719} \tan \left(\frac{1}{2} \sqrt{719} \log e^{2}+\theta\right)\right) \tag{11.3.21}
\end{equation*}
$$

giving

$$
\begin{equation*}
\lambda(t)=\frac{e^{2}(t)}{10}\left(19+\sqrt{719} \tan \left(\frac{1}{2} \sqrt{719} \log e^{2}(t)+\theta\right)\right) \tag{11.3.22}
\end{equation*}
$$

where $\theta$ is an integration constant. Both $e$ and $\lambda$ grow in the UV but, as explicitly shown in fig. 11.2, the quartic coupling $\lambda$ varies significantly over a range in which the electric charge remains essentially constant. We can then conclude that for a wide range of initial conditions for $e$ and $\lambda$ there exists an energy scale where $\lambda \sim e^{4}$, and in particular eq. (11.1.8) is valid.

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[^0]:    ${ }^{1}$ Note that the normalizations of the creation/annihilation operators and of the single- and multi-particle states introduced here differ from the corresponding ones of ref. [1].

[^1]:    ${ }^{2}$ The analysis in this section closely follows section 4.3 of ref. [2].
    ${ }^{3}$ Strictly speaking this statement is not correct, since quantum entanglement among states provide correlations that do not decrease as the distance increases (this was at the base of the so called Einstein-Podolski-Rosen paradox claiming possible violations of causality in quantum mechanics). Such states are however fragile and short-lived, since decoherence effects given by arbitrarily small external perturbations will destroy the entanglement among the states asymptotically. A more precise statement would be that stable long-lived states should satisfy cluster decomposition. In general, cluster decomposition applies when the vacuum is a pure state, and it fails when it is a mixed state. The phenomenon of spontaneous symmetry breaking is also related to cluster decomposition, see section 8.1.

[^2]:    ${ }^{4}$ Note that in some textbooks, e.g., ref. [1] the factors $h_{\vec{p}}$ are actually absorbed in the definition of the multiparticle states and therefore they are absent in the formula above.

[^3]:    ${ }^{5}$ This expression differ from the one presented, e.g., in eq. (4.79) of ref. [1] because of the different conventions for the normalization of the states.

[^4]:    ${ }^{6}$ Clearly, the distinction is dictated by the time scale of the process. Particles with a time decay much longer than the interacting time of a process are effectively stable and can be effectively described by asymptotic states.

[^5]:    ${ }^{7}$ Unless $a_{\text {inc }}(\omega)$ has zeros that induce poles in $f(\omega)$, but these are avoidable by simply changing the (quite arbitrary) incident wave packet.

[^6]:    ${ }^{8}$ If the same particle can be exchanged in the process, we have $\mu=m$, but this is not necessary.

[^7]:    ${ }^{9}$ We briefly discussed crossing symmetry as a "simple" consequence of the LSZ formula (2.3.12) below that equation. In fact, this property is far from trivial because the crossed amplitudes are functions defined in disconnected regions in the Mandelstam variables with respect to the original amplitude function. Proving crossing symmetry is equivalent to prove that the crossed amplitude is the analytic continuation in the crossed region of the original amplitude.

[^8]:    ${ }^{10}$ Note that the possible branch-cut singularities around $p=0\left(E=-V_{0}\right)$ eventually cancel between each other, so that $T(E)$ is single-valued around $p=0$.

[^9]:    ${ }^{1}$ The analysis in this chapter closely follows chapter 12 of ref. [2], with a straightforward generalization to space-time dimensions different than 4.
    ${ }^{2}$ The renormalization of spontaneously broken gauge theories will not be discussed in detail. However, see the end of section 8.4 for a brief comment.
    ${ }^{3}$ This is called $D$ in ref. [2]. We call it $\delta$, because we will denote with the symbol $D$ the number of space-time dimensions in what follows.

[^10]:    ${ }^{4}$ It is understood here and in the following that these integrals have been "regularized" (i.e. made finite). In this way derivatives commute with the integration and can act directly on the integrand.

[^11]:    ${ }^{5}$ By interactions we here mean also terms quadratic in the fields.

[^12]:    ${ }^{6}$ As we will see, there are symmetries that are not respected by any regulator. When this happens, the symmetry is called anomalous, since it is no longer there at the quantum level. Anomalies will be the subject of chapter 9 .

[^13]:    ${ }^{7}$ The identity (3.3.5) is proved in section 4.4.3. Its non-abelian generalization is proved in section 6.3.

[^14]:    ${ }^{8}$ Identities such as eq. (3.3.9) are shown to hold in section 4.4.2.
    ${ }^{9}$ It is easy to understand why gauge invariance forbids the presence of the factor $K$ in eq.(3.3.8) or $C_{1}$ in eq.(3.3.4). If present, these divergences would require a counter-term in the Lagrangian of the form $\left(A_{\mu} A^{\mu}\right)^{2}$ or $A_{\mu} A^{\mu}$ respectively, which obviously does not respect gauge invariance.

[^15]:    ${ }^{10}$ As we discuss below, the analytic continuation of integrals such as eq. (3.4.1) can be done without invoking $\Gamma$ functions and their properties.

[^16]:    ${ }^{11} \Omega_{d}$ can be determined by calculating $I_{d}$ in eq. (3.4.3) with $f(|\vec{p}|)=\exp \left\{-\sum_{i=1}^{d} p_{i}^{2}\right\}$ both in cartesian and in spherical coordinates (see eq. (3.4.4)), which give $I_{d}=\pi^{d / 2}$ and $I_{d}=\Omega_{d} \Gamma(d / 2) / 2$, respectively, and therefore eq. (3.4.5).

[^17]:    ${ }^{1}$ Recall that a 1PI amplitude is defined as the amplitude whose defining graphs cannot be divided in two by cutting one internal line.
    ${ }^{2}$ The analysis in this section closely follows section 16.1 of ref. [2].
    ${ }^{3}$ It is important to keep in mind that $\Gamma$, contrary to $S$, is not in general a local functional of the fields: $\Gamma[\Phi] \neq \int d^{4} x \mathcal{L}_{e f f}(\Phi(x))$, see eq.(4.1.11).

[^18]:    ${ }^{4}$ The analysis in this section closely follows section 16.2 of ref. [2].

[^19]:    ${ }^{5}$ Strictly speaking, the coupling constant $\lambda$ appearing in eq. (4.2.18) is not the same as that appearing in eq. (4.2.17) because of a shift induced by the one-loop term. We neglect these finite corrections, possibly relevant when a careful definition of the coupling $\lambda$ is provided. We postpone to chapter 5 a more careful discussion of the definition of coupling constants.

[^20]:    ${ }^{6}$ The usual purely linear transformations of the fields are obtained from eq.(4.4.20) by taking $s(x)=0$ and $t(x, y)=t \delta(x-y)$.

[^21]:    ${ }^{7}$ The derivation of $\mathcal{L}_{g . f \text {. from a functional point of view, for both abelian and non-abelian theories, is }}^{\text {f }}$ discussed in section 6.2.
    ${ }^{8}$ Pay attention to the anticommuting nature of the Grassmann fields to get the signs right!

[^22]:    ${ }^{1}$ The analysis in this section closely follows section 12.1 of ref. [1].

[^23]:    ${ }^{2}$ Strictly speaking, the actions (5.1.3) should be written in momentum space using eq. (5.1.1) and recalling the overall bound $|k|<\Lambda$. This results in unnecessarily long expressions that are avoided in the rough, but more brief form (5.1.3).

[^24]:    ${ }^{3}$ The couplings $\delta Z, \delta m^{2}$, etc. appearing in eq. (5.1.5), after the redefinition $k \rightarrow b k$, can depend on $b$, as we will explicitly see in what follows. However, in order to avoid confusion with the couplings defined in eqs. (5.1.8), we have omitted to write this $b$-dependence.

[^25]:    ${ }^{4}$ On the contrary, when the mass of the field is sufficiently large, all frequency modes, i.e. the whole field, can and should be integrated out. In this case the Wilsonian picture is the way to go and results in a simplification of the physical system. More on this in chapter 7.

[^26]:    ${ }^{5}$ Since $s+t+u=4 m^{2}$, the Mandelstam variables will necessarily have, at high energies, different signs, so that the argument of some of the logs in eq. (5.2.3) will be negative with branch cuts singularities. This is of course expected, since by the optical theorem $\Gamma^{(4)}$ should have an imaginary part. In order to avoid branch cut-singularities and imaginary amplitudes, that will not change the discussion that follows, we consider off-shell amplitudes at euclidean values of the momenta, where $s, t$ and $u$ are all negative.

[^27]:    ${ }^{6}$ Notice that $\beta_{U V}$ in eq. (5.2.9) coincides with $b d \lambda(b) / d b$ in eq. (5.1.12), as expected, being the same thing.

[^28]:    ${ }^{7}$ The scale dependence of $\mu_{L}$ on $\mu_{0}$ is only apparent. One can check that $d \mu_{L} / d \mu_{0}=0$.

[^29]:    ${ }^{8}$ There can be exceptions to this theorem when the particle mass arises from a spontaneous breaking of a symmetry, in which case it is governed by a coupling.

[^30]:    ${ }^{9}$ Actually, the original Callan [17] and Symanzik [18] equations were a bit different. There was no sliding scale $\mu$, the physical mass playing essentially the role of $\mu$, and the equations were not homogeneous in $\Gamma^{(n)}$. The more modern version of the Callan Symanzik equations (5.4.3) in terms of $\mu$ were developed shortly later, see refs.[19] and [20].

[^31]:    ${ }^{10}$ By dimensional analysis other terms could appear in eq.(5.8.15), proportional to couplings associated to operators of dimension 7 and 8 . These operators are however redundant and by the equations of motion (or equivalently a field redefinition) can be removed. We will discuss redundant operators in section 7.8.
    ${ }^{11} \mathrm{RG}$ techniques can be used in this theory in another way, by introducing the so called $\epsilon$-expansion [21]. This is a very interesting and useful technique, which will not be discussed here due to lack of time.

[^32]:    ${ }^{12}$ Again, this is only true in a mass-independent scheme. From now on, the use of a mass-independent scheme will be assumed.
    ${ }^{13}$ In the operatorial formalism, where $\phi$ is an operator in the Hilbert space, some (in general not all) divergences of composite fields are removed by the normal ordering procedure.
    ${ }^{14}$ Composite or elementary. In the latter case $Z^{\phi}(\mu)=\sqrt{Z(\mu)}$.

[^33]:    ${ }^{15}$ The degree of mixing depends on the renormalization scheme. In particular, operators with different classical dimensions cannot mix in mass-independent schemes in theories with marginal couplings only.

[^34]:    ${ }^{16}$ Unless selection rules require to consider higher dimensional operators. Moreover, this is only true in a perturbative context. At strong coupling everything can possibly happen, e.g. a classically irrelevant operator might turn into a marginal or even a relevant operator.

[^35]:    ${ }^{1}$ The group $S O(3)$ is locally isomorphic to $S U(2)$ and is defined by the same 3 generators $t_{\alpha}^{\text {Adj. introduced }}$ above for $S U(2)$. The group $S O(2)$ is isomorphic to $U(1)$.

[^36]:    ${ }^{2}$ Notice the similarity of the definition (6.2.24) with the definition of diffeomorphism-invariant measure in general relativity.

[^37]:    ${ }^{3}$ Recall that inside the path integral $\omega$ and $\omega^{*}$ are two independent variables. In particular, $\omega^{*}$ is not the complex conjugate of $\omega$.

[^38]:    ${ }^{4}$ The analysis in this section closely follows section 15.7 of ref. [2].

[^39]:    ${ }^{5}$ Namely with Fermi statistics. In particular, ghosts and anti-ghosts are fermionic.

[^40]:    ${ }^{6}$ Note that the sum $A+Q$ transforms in the same way (i.e., as in eq. (6.1.18)) under both transformations (6.4.2) or (6.4.3), as expected from the full gauge field. The different splitting distinguishes however the background from the quantum gauge transformations.

[^41]:    ${ }^{7}$ The analysis in this section closely follows sections $16.4,17.1$ and 17.2 of ref. [2].

[^42]:    ${ }^{1}$ See 4.3 for a further subtlety.

[^43]:    ${ }^{2}$ This theory is not realistic, since the cubic potential $H L^{2}$ is not positive definite, but the considerations we want to make are insensitive to this stability problem.

[^44]:    ${ }^{3}$ The hierarchy problem is often defined as due to quadratically (or higher) divergent graphs in regularizations with some cut-off $\Lambda$. Although this argument is essentially correct, it relies on the use of $\Lambda$ and, taken as it is, does not apply to regularizations like DR , where no $\Lambda$ is introduced and there are no real quadratic divergences. It would then seem that the hiearchy problem is "scheme dependent", whereas of course it is not, as we have just shown.

[^45]:    ${ }^{4}$ Recall that the limit $d \rightarrow 4$ should always be performed at the end of the computation.

[^46]:    ${ }^{5}$ The label $1 / 2$ and $3 / 2$ refers to the quantum numbers of such operators under an approximate global $S U(2)_{L}$ symmetry under which $u_{L}$ and $d_{L}$ form a doublet and $s_{L}$ is a singlet. This is an abuse of language, because while $O_{1 / 2}$ is in fact a component of an operator doublet with isospin $1 / 2$, the operator $O_{3 / 2}$ is a component of a mixture of operators with isospin $1 / 2$ and $3 / 2$. We follow here the notation of ref.[1].
    ${ }^{6}$ This is, of course, a simplification, since we should take into account the $b$ quark for energies above its mass and integrate out the $c$ quark below its mass, by performing a more refined matching.

[^47]:    ${ }^{7}$ Notice that this true in any number of even dimensions $d$. In particular, the total number of matrices equal $(1+1)^{d}=2^{d}=2^{d / 2} \times 2^{d / 2}$, where $2^{d / 2}$ is the dimension of the spinor space in $d$ even dimensions.

[^48]:    ${ }^{8}$ As we will see in the next section, this is a so called redundant operator, but this is inessential for the argument we want to make here.

[^49]:    ${ }^{9}$ We have seen that no contribution in DR arises from the Jacobian associated to the change of variables (7.8.12), which can then be neglected.

[^50]:    ${ }^{1}$ The analysis in this section closely follows section 19.1 of ref. [2]. I also found useful chapter 2.2 of Parisi's book [23], that contains several interesting observations that are often omitted in QFT textbooks, in particular about the role of cluster decomposition.

[^51]:    ${ }^{2}$ The analysis in this section closely follows section 19.3 of ref. [2].

[^52]:    ${ }^{3}$ This often used terminology is actually misleading. Gauge symmetries, being merely redundancies of the system, cannot be broken. A more proper term would be gauge theories in a non-linearly realized, or Higgs, phase.

[^53]:    ${ }^{4}$ We take $\delta v_{m}=0$, so that $\delta \phi_{m}^{\prime}=\omega^{\alpha} i t_{m}^{\alpha} \phi_{n}$.

[^54]:    ${ }^{5}$ From now on we omit, for simplicity, the partial or total dependence on $x, \xi_{a}(x)$, etc. of the various fields.

[^55]:    ${ }^{6}$ The analysis in this section closely follows section 19.7 of ref. [2].

[^56]:    ${ }^{7}$ There is actually no way to distinguish an $S U(3)_{L}$ from an $S U(3)_{R}$ transformation when acting on a $3 \times 3$ matrix such as $U$. Strictly speaking, we should write $U$ as a $6 \times 6$ matrix, containing $U$ and $U^{\dagger}$, to properly distinguish the two $S U(3)$ 's. In doing that, one gets eq. (8.7.6).

[^57]:    ${ }^{8}$ The fermion Yukawa couplings also break $S U(2)_{c}$, but here for simplicity we are focusing on the bosonic sector of the SM, neglecting fermions altogether.

[^58]:    ${ }^{9} S O(5)$ might also be broken by additional terms, such as the analogues of the quarks masses in QCD. Without knowing the UV theory, we might for simplicity assume that there are no such terms.

[^59]:    ${ }^{10}$ In order not to confuse the group element $g$ with the gauge coupling constant $g$, we adopt here a non-canonical normalization for the gauge fields $A_{\mu}^{\alpha}$.

[^60]:    ${ }^{1}$ Contact terms will not vanish when taking the divergence of currents in other correlators, such as three vector currents. These terms are always proportional to the structure constants $C_{\beta \alpha \gamma}$ and can be shown to not give rise to anomalies.

[^61]:    ${ }^{2}$ There also exist pure gravitational anomalies. These are vanishing in 4 space-time dimensions, but can occur in $4 n+2$ dimensions ( $n$ a non-negative integer). We will not discuss these anomalies, that are beyond our course.

[^62]:    ${ }^{3}$ For simplicity of the notation, we omit the gauge index $A$ in the following equations. It will be reintroduced later on in this section.

[^63]:    ${ }^{4}$ Locality is crucial. By using non-local functionals, any anomaly can be cancelled. For instance, a gauge $U(1)^{3}$ anomaly would be cancelled by adding to $\Gamma(A)$ the non-local functional

[^64]:    ${ }^{5}$ The argument requires $n_{f}=3$. Solutions for the bound states quantum numbers can instead be found when $n_{f}=2$.

[^65]:    ${ }^{6}$ The analysis in this section closely follows section 19.5 of ref. [1].

[^66]:    ${ }^{7}$ In fact, this is not the most general possibility. A conterexample is provided by the transformation of the Lagrangian density under scale transformations considered below.

[^67]:    ${ }^{8}$ This term plays an important role when studying non-perturbative configurations in QCD, instantons. We will not consider such effects in these lectures.

[^68]:    ${ }^{9}$ Notice that this naturalness problem is qualitatively different from the one affecting the Higgs mass (or more generally, relevant operators). There, a delicate cancellation (or fine-tuning) has to occur to get a small number out of big ones, induced by quantum corrections. Here $\bar{\theta}$ is radiatively stable and we have "simply" to understand the smallness of its classical value.

[^69]:    ${ }^{1}$ A similar event occurs for $\alpha>0$ in presence of a constant electric field. We will study in some quite detail this phenomenon in the subsequent sections of this chapter.
    ${ }^{2}$ This argument works also in presence of fermions. We integrate them out and identify the resulting effective action with $S$ in eq.(10.1.3).
    ${ }^{3}$ With an appropriate measure and upon renormalization. Strictly speaking a full non-perturbative definition of the path integral requires a lattice discretization of space-time.

[^70]:    ${ }^{4}$ The analysis that follows can easily be generalized for large-order behaviours of the kind $Z_{n} \sim$ $(n!)^{k} a^{n} n^{c}$.

[^71]:    ${ }^{5}$ Indeed, indicating by $q_{ \pm}=[|-i \vec{E}+\vec{B}| \pm|i \vec{E}+\vec{B}|] / 2=\left\{\left[|\vec{B}|^{2}-|\vec{E}|^{2}-2 i \vec{E} \cdot \vec{B}\right]^{1 / 2} \pm\left[|\vec{B}|^{2}-|\vec{E}|^{2}+\right.\right.$ $\left.2 i \vec{E} \cdot \vec{B}]^{1 / 2}\right\} / 2$ the factors in the arguments of cosh on the second line of eq. (10.2.64), the last equality in that equation follows from the fact that a comparison with eq. (10.2.44) yields the equality $q_{ \pm}^{2}=\mp a_{ \pm}^{2}$ and therefore $q_{+}=i a_{+}$and $q_{-}=a_{-}$(a possible ambiguity in the overall sign does not affect the final result).

[^72]:    ${ }^{1}$ Of course, the counter-term $Z$ should not be confused with the finite $Z(\rho)$ appearing in the effective action (11.2.5).

[^73]:    ${ }^{2}$ Pay attention in not confusing the photon mass $\mu_{\gamma}$ with the sliding scale $\mu$ !

