

QFT II

0 The Free Field Case

To start, we will first consider one of the simplest QFTs we can imagine: a massive scalar field on a one-dimensional line. This theory in particular is the one that we arrive at if we take the limit as $N \rightarrow \infty$ of N identically coupled harmonic oscillators with an additional self-coupling term for each mass. The position-space Hamiltonian is given by

$$H_0 = \frac{1}{2} \int dx \left(\pi(x)^2 + \left(\frac{\partial \phi}{\partial x} \right)^2 + m_\phi^2 \phi(x)^2 \right), \quad (0.0.1)$$

where π is the conjugate momentum to the field ϕ . Performing a Fourier transform to new fields $\tilde{\phi}$ and $\tilde{\pi}$, this Hamiltonian is given by

$$H_0 = \frac{1}{2} \int \frac{dp}{2\pi} \left(|\tilde{\pi}(p)|^2 + (p^2 + m_\phi^2) |\tilde{\phi}(p)|^2 \right), \quad (0.0.2)$$

where the reality conditions of π and ϕ require $\tilde{\pi}^*(p) = \tilde{\pi}(-p)$ and $\tilde{\phi}^*(p) = \tilde{\phi}(-p)$, and we see that we have decoupled all of our momentum modes into single harmonic oscillator solutions with frequencies $\omega_p = \sqrt{p^2 + m_\phi^2}$.

When we move to the quantum theory, we diagonalize the Hamiltonian using the creation and annihilation operators

$$\tilde{\phi}(p) = \frac{1}{\sqrt{2\omega_p}} (a_p + a_{-p}^\dagger), \quad \tilde{\pi}(p) = -i\sqrt{\frac{\omega_p}{2}} (a_p - a_{-p}^\dagger), \quad (0.0.3)$$

which reduces the Hamiltonian to the form

$$H_0 = \int \frac{dp}{2\pi} \omega_p \left(a_p^\dagger a_p + \frac{1}{2} \right). \quad (0.0.4)$$

We can think of the field and momentum operators in Eq. (0.0.3) as measuring the value of the field and the time variation of the field, respectively, in some small momentum-space volume centered around p . Likewise, the spatial operators

$$\phi(x) = \int \frac{dp}{2\pi} \tilde{\phi}(p) e^{ipx}, \quad \pi(x) = \int \frac{dp}{2\pi} \tilde{\pi}(p) e^{ipx}, \quad (0.0.5)$$

give the similar values in a small spatial region centered at x .

The vacuum energy gives a divergent contribution, but it is constant, so it can always be subtracted off from the energy (recall that only differences in energy are physical things anyway). Eigenstates of the free-particle Hamiltonian are then simply given by¹

$$\bigotimes_p |n_p\rangle \Rightarrow H_0 \left(\bigotimes_p |n_p\rangle \right) = \int \frac{dp}{2\pi} n_p \omega_p \left(\bigotimes_p |n_p\rangle \right), \quad (0.0.6)$$

where the eigenvalue of the state is the total energy

$$E = \int \frac{dp}{2\pi} n_p \omega_p. \quad (0.0.7)$$

¹Note that these states are not the standard ones chosen with relativistic normalization, but for the purposes of our considerations, this point is not relevant.

0.1 “Pictures” of Quantum Mechanics

In the standard way of working with quantum mechanics, we start with an initial state $|\psi\rangle = |\Psi(t_0)\rangle$ which we then want to evolve in time. This time evolution is described by the Schrödinger equation

$$i\frac{\partial}{\partial t} |\Psi(t)\rangle = H |\Psi(t)\rangle, \quad (0.1.1)$$

where H is the Hamiltonian describing the theory. This equation has the solution

$$|\Psi(t)\rangle = U(t, t_0) |\Psi(t_0)\rangle, \quad U(t, t_0) = e^{-iH(t-t_0)}, \quad (0.1.2)$$

where $U(t, t_0)$ is known as the time-evolution operator. This way of calculating, where we choose to time-evolve the states of a system, is known as the “Schrödinger picture” of quantum mechanics.

We can instead work with a different picture by considering the following: take a time-independent operator, \mathcal{O}_S , in the Schrödinger picture and place it in a matrix element between two time-evolved states

$$\langle \chi_S(t) | \mathcal{O}_S | \psi_S(t) \rangle = \langle \chi_S(t_0) | e^{iH(t-t_0)} \mathcal{O}_S e^{-iH(t-t_0)} | \psi_S(t_0) \rangle \quad (0.1.3)$$

In some sense, this is the only reasonable thing to consider since any physical observation depends only on matrix elements of operators. But clearly, the exact same physics is described if we choose to time evolve *operators* instead of states using

$$\langle \chi_H | = \langle \chi_S(t_0) |, \quad |\psi_H\rangle = |\psi_S(t_0)\rangle, \quad \mathcal{O}_H(t) = e^{iH(t-t_0)} \mathcal{O}_S e^{-iH(t-t_0)} = e^{iH(t-t_0)} \mathcal{O}_H(t_0) e^{-iH(t-t_0)}. \quad (0.1.4)$$

It is a straightforward exercise to show that such an operator satisfies a modified version of the Schrödinger equation,

$$i\frac{\partial}{\partial t} \mathcal{O}_H(t) = [\mathcal{O}_H(t), H_H(t)], \quad (0.1.5)$$

where $H_H(t) = U^\dagger(t, t_0) H U(t, t_0) = H$. This picture, where we time evolve operators instead of the states, is known as the “Heisenberg picture.”

It turns out that working with the Heisenberg picture will come in handy a bit later, so this is what we will use. If we choose $\phi(x, t_0) = \phi(x)$ where $\phi(x)$ is the Schrödinger picture operator in Eq. (0.0.3), we find

$$\phi(x, t) = \int \frac{dp}{2\pi} \frac{1}{\sqrt{2\omega_p}} \left(U^\dagger(t, t_0) a_p U(t, t_0) e^{ipx} + U^\dagger(t, t_0) a_p^\dagger U(t, t_0) e^{-ipx} \right), \quad (0.1.6)$$

where, in the second term, we have taken $p \rightarrow -p$ in the integral so that the creation operator creates a positive-momentum mode. To see how the time-evolution operators act on the creation and annihilation operators, we recognize that these operators eat a state $|n_p\rangle$ and spit out a state $|(n+1)_p\rangle$ and $|(n-1)_p\rangle$, respectively, and produce the appropriate constants. This can be realized using the form

$$a_p = \sum_{n=0}^{\infty} \sqrt{n_p} |(n-1)_p\rangle \langle n_p|, \quad a_p^\dagger = \sum_{n=0}^{\infty} \sqrt{n_p+1} |(n+1)_p\rangle \langle n_p|. \quad (0.1.7)$$

Then, using the fact that the states, $|n_p\rangle$, are eigenstates of the Hamiltonian, we find

$$\begin{aligned} e^{iH(t-t_0)} a_p e^{-iH(t-t_0)} &= \sum_n \sqrt{n_p} |(n-1)_p\rangle \langle n_p| e^{i(n_p-1)\omega_p(t-t_0)} e^{-in_p\omega_p(t-t_0)} = a_p e^{-i\omega_p(t-t_0)}, \\ e^{iH(t-t_0)} a_p^\dagger e^{-iH(t-t_0)} &= \sum_n \sqrt{n_p+1} |(n+1)_p\rangle \langle n_p| e^{i(n_p+1)\omega_p(t-t_0)} e^{-in_p\omega_p(t-t_0)} = a_p^\dagger e^{i\omega_p(t-t_0)}. \end{aligned} \quad (0.1.8)$$

with this, we find the expression for the Heisenberg field operator

$$\phi(X) = \int \frac{dp}{2\pi} \frac{1}{\sqrt{2\omega_p}} \left(a_p e^{-i(\omega_p(t-t_0)-px)} + a_p^\dagger e^{i(\omega_p(t-t_0)-px)} \right) = \int \frac{dp}{2\pi} \frac{1}{\sqrt{2\omega_p}} \left(a_p e^{-iP \cdot X} + a_p^\dagger e^{iP \cdot X} \right), \quad (0.1.9)$$

where we denote spacetime vectors with capital letters and $P^\mu = (\omega_p, p)$ is the standard form of the two-momentum.

0.2 A first calculation

Now, we want to ask a “physical” question from this system. Suppose we start with some source which prepares a single-particle state at time $t = 0$ (really, we can think of this as a measurement of the vacuum state of the field which pulls a particle from the measurement device). We now set up a detector at position y and we want to know, after evolving the system for some time t , what is the probability of our detector seeing (and therefore annihilating) the particle. After we make our field measurement at y , the field should be left in a vacuum state so that we know that we did, indeed annihilate the particle instead of creating it. Furthermore, we will not consider the experiment as preparing the particle at a single position, but instead, we will consider the more realistic case where it is prepared in some region centered around x described by the distribution $f(x' - x)$.

Then, our question is equivalent to finding the square-magnitude of the matrix element

$$\begin{aligned}\mathcal{M}(t, x, y) &= \int dx' f(x' - x) \langle 0 | \phi(Y) \phi(X') | 0 \rangle \\ &= \int dx' \int \frac{dp_1}{2\pi} \int \frac{dp_2}{2\pi} f(x' - x) \frac{1}{2\sqrt{\omega_{p_1}\omega_{p_2}}} e^{-i(\omega t - p_2 y + p_1 x')} \langle 0 | a_{p_2} a_{p_1}^\dagger | 0 \rangle \\ &= \int dx' \int \frac{dp}{2\pi} \frac{1}{2\omega_p} f(x' - x) e^{-i(\omega_p t - p(y - x'))},\end{aligned}\tag{0.2.1}$$

where we defined $X'^\mu = (0, x')$, $Y^\mu = (t, y)$ and used the fact that the vacuum is invariant under time translations. In the second line, we passed to our Fourier transformed variables and in the third line, we used the commutation relations of the creation/annihilation operators and the fact that annihilation operators annihilate the zero-particle state, i.e. $a_p | 0 \rangle = 0$.

The momentum integral in Eq. (0.2.1) can be evaluated analytically considering three cases. Defining $z = y - x$, these cases are: (i) $t > |z|$ or the timelike case, (ii) $t < |z|$ the spacelike case, and (iii) $t = |z|$ or the null case. In all cases, it will be helpful to first redefine the variable $p = m \sinh \alpha$ which then gives $\omega_p = m \cosh \alpha$.

- Case (i): In the case where $t > |z|$, we can always perform a Lorentz transformation into a new, primed frame with relative velocity $v = z/t$ so that

$$t' = \gamma(t - vz) = \sqrt{t^2 - z^2}, \quad z' = \gamma(z - vt) = 0,\tag{0.2.2}$$

giving the momentum integral

$$\int_{-\infty}^{\infty} d\alpha e^{-imt' \cosh \alpha} = 2 \int_0^{\infty} d\alpha e^{-imt' \cosh \alpha} = -i\pi H_0^{(2)}(m\sqrt{t^2 - z^2}),\tag{0.2.3}$$

where $H_0^{(2)}(z)$ is the zeroth Hankel function of the second kind.

- Case (ii): Similar to case (i), we can boost to a frame where the separation is purely spacelike using velocity $v = t/z$. In this case, we find

$$\int_{-\infty}^{\infty} d\alpha e^{-imz' \sinh \alpha} = 2K_0(m\sqrt{z^2 - t^2}),\tag{0.2.4}$$

where $K_0(z)$ is the zeroth modified Bessel function of the second kind.

- Case(iii): The integral we wish to look at is actually divergent in the null case. However, we can find the behavior of this region by simply taking the limit as $t^2 - z^2 \rightarrow 0$ from above (limit of case (i)) or from below (limit of case (ii)). As we should expect, the limiting behavior of these two functions is the same for both of these considerations

$$\lim_{x \rightarrow 0^+} -i\pi H_0^{(2)}(x) = \lim_{x \rightarrow 0^-} 2K_0(x) = \lim_{x \rightarrow 0} \log \frac{1}{x^2}.\tag{0.2.5}$$

Note that, since this is a logarithmic divergence, the vanishing spacetime distance dominates the constant logarithm of the mass and the scaling near $|t^2 - z^2| \rightarrow 0$ becomes mass-independent. So, this is simply the same result we would get if we considered a massless particle, which is restricted to travel along its light cone, hence the pole in the result.

So, all in all, the expression for the state we wish to examine is given by

$$\mathcal{M}(t, x, y) = \int_{-\infty}^{\infty} dx' f(x' - x) \left(-\frac{i}{4} H_0^{(2)}(m\sqrt{t^2 - z'^2}) \Theta(t^2 - z'^2) + \frac{1}{2\pi} K_0(m\sqrt{z'^2 - t^2}) \Theta(z'^2 - t^2) \right), \quad (0.2.6)$$

where, of course, the null case arises from whether we choose the $+$ or $-$ for $\lim_{x \rightarrow 0^\pm} \Theta(x) = 1$ in the definition of our Heaviside function and we defined $z' = y - x'$.

The result in Eq. (0.2.6) is perfectly correct, but it has a feature that might bother us: it is non-zero outside of the particle's light cone. At its core, this has to do with correlation of any two points in the field, but we can think of it as we really asked a question that didn't necessarily require causality. In fact, our question only gives us the probability of the outcome of two measurements, it had nothing to do with if one measurement *caused* the other. So really, it only tells us how the two measurements are *correlated* with each other. This is why matrix elements of the type in Eq. (0.2.6) is often called "correlation functions."

If instead, we want to know whether or not the measurement at X' affects the measurement at Y , we need to subtract off the piece where we first make the measurement at Y and second at X' . Note that "first" here is purely the quantum mechanical sense, where the Y measurement acts on the state in the ordering of operators before X' and has nothing to do with the chronological ordering of events². If the two orderings are different, this difference is exactly the amount that one measurement can affect the other. The second ordering amounts to computing

$$\begin{aligned} \mathcal{M}'(t, x, y) &= \int dx' f(x' - x) \langle 0 | \phi(X') \phi(Y) | 0 \rangle \\ &= \int dx' \int \frac{dp}{2\pi} \frac{1}{2\omega_p} f(x' - x) e^{i(\omega_p t - p(y - x'))}. \end{aligned} \quad (0.2.7)$$

Using the same methods as before, it is a straightforward calculation to show that, if we subtract this contribution from the result in Eq. (0.2.6), we find that the transmission matrix element, $\mathcal{M}_R(t, x, y) = \mathcal{M}(t, x, y) - \mathcal{M}'(t, x, y)$ is perfectly causal!

Interestingly, it looks like the matrix element in Eq. (0.2.7) is exactly the same as the original matrix element, but instead of creating a particle with two-momentum P^μ , we instead create a particle with $-P^\mu$. Alternatively, if we look at the momentum-space amplitude

$$\mathcal{M}^{(\prime)}(t, x, y) = \int dx' \int \frac{dp}{2\pi} f(x' - x) \tilde{\mathcal{M}}^{(\prime)}(p), \quad (0.2.8)$$

we see that $\tilde{\mathcal{M}}^{(\prime)}(p) = \tilde{\mathcal{M}}^*(-p)$. So, this particle with reversed two-momentum can instead be thought of as the complex conjugated particle with spatial momentum reversed, but with all of the exact same properties of the standard field. This is what is known as an antiparticle (in fact this story is much easier to see when using complex fields, but this discussion will suffice for now). So, it seems that, in order for the result to be explicitly causal, we need to also consider the linear combinations with matrix elements including the creation of an antiparticle traveling in the opposite direction!

This is a very important note, but going forward, we will simply live with the fact that these correlations die off very rapidly outside the lightcone and knowing that they are not really a problem for causality, as long as we also allow for the creation of antiparticles as well as particles³.

²There is also an interesting point about Lorentz invariance hidden here. If the two measurements are space-like separated, we can always boost into a frame where we reverse the chronological ordering of the measurements. So, we can't really say which operators acts on the system first if they are causally disconnected.

³There is also the point that there is some uncertainty in our measurement, so the light-cone of a particle actually becomes "fuzzy," since we can't exactly know the time or location that the particle was created.

1 Adding Interactions

As we saw in the last section, we want to calculate the correlation functions of vacuum-to-vacuum transitions with some insertions of field operators. In other words, we want to consider starting with a vacuum state, creating some set of particles, letting these particles evolve (maybe creating some other particles in the process), then setting up a series of detectors to measure these particles.

This is a simple task for the non-interacting theory for one reason: the Hamiltonian is easily diagonalizable in terms of momentum-space harmonic oscillators. As soon as we add some interactions, i.e.

$$H = H_0 + H_I, \quad (1.0.1)$$

this decomposition into harmonic oscillators will no longer be possible in general. In fact, for a general interacting QFT, it is unknown how to even define eigenstates of the theory, or even if the mathematical object exists for that matter. Our lives as theoretical physicists can't be too easy!

1.1 Small couplings to the rescue!

Thankfully, in certain theories, we can get around the problem of not knowing how to calculate in interacting QFTs if they have the very nice property that (qualitatively) $H_I \ll H_0$.

This allows us to do many things, but first and foremost, it allows us to define a vacuum state. Once we introduce interactions, the vacuum (or zero-particle state) of the theory is *not necessarily the same* as the zero-particle state of the free theory. This is simply because interactions change the definition of a particle, because they can become “dressed” or altered by other fields that they interact with. However, if the interactions are very weak, as any particles propagate infinitely far away from each other, they will feel less and less of an effect from these interactions and get closer to their non-interacting counterparts. In particular, if we only consider the infinitely distant past and the infinitely far future, the two theories should coincide. This means that we can define

$$\lim_{t \rightarrow \pm\infty} a_{m,p} |\Omega\rangle = 0, \quad (1.1.1)$$

where we have added an additional flavor index, m , to the annihilation operators in the case where we have several fields⁴.

Furthermore, since we only insert the particles we create first and the ones we annihilate later, the operator insertions will always be in chronological order (remember that this will lead to some “non-causal” effects, but these are simply the correlations, they have nothing to do with transmission, and we can always obtain the purely causal pieces by considering antiparticle propagation). So our task is now to compute the vacuum-to-vacuum correlation functions of time-ordered products of operators. In other words, we want to compute matrix elements like

$$\langle \Omega | \mathbf{T} \{ \Phi_{m_n}(X_n) \Phi_{m_{n-1}}(X_{n-1}) \dots \Phi_{m_2}(X_2) \Phi_{m_1}(X_1) \} | \Omega \rangle \quad (1.1.2)$$

where again, m_i 's are flavor indices and \mathbf{T} denotes a time-ordered product so that operators are ordered chronologically right-to-left. From this point forward, we will simply assume time ordering and omit the \mathbf{T} .

To be even more general, we don't need to consider just field operator insertions, we could consider more general operator insertions, $\mathcal{O}_\alpha(\Phi_\ell, \Pi_\ell)$, which depend on both field and momentum operators. However, to keep the discussion simpler, we will just consider field operator insertions, as this is what we will actually want to calculate.

It is also important to keep in mind that this framework completely neglects all bound states of particles. So, when we make these assumptions, objects like atoms or nuclei can never appear in our theory explicitly (though they can sometimes be included in other clever ways).

1.2 Deriving the Path Integral

We can evaluate time ordered products like in Eq. (1.1.2) by expanding in creation/annihilation operators, reordering, and collapsing the resulting products using delta functions, but it turns out to be more economical to go about it in a different way. This section will pretty much exclusively follow Weinberg's textbook [?].

⁴In more general theories, we should also consider some spin index, σ .

The first step is to recognize that the Schrödinger-picture operators $\Phi_m(x)$ and $\Pi_m(x)$ are standard, albeit space-dependent, quantum mechanical operators, and therefore have complete sets of eigenstates and eigenvalues which span the Hilbert space

$$\Phi_m(x) |\varphi\rangle = \varphi_m(x) |\varphi\rangle, \quad \Pi_m(x) |\pi\rangle = \pi_m(x) |\pi\rangle. \quad (1.2.1)$$

Furthermore, since the classical versions of Φ and Π are canonical conjugates, their operator versions satisfy the canonical commutation relations

$$[\Phi_m(x), \Pi_n(x)] = i\delta_{mn}\delta(x-y), \quad (1.2.2)$$

and their corresponding eigenstates satisfy

$$\begin{aligned} \langle\varphi|\varphi'\rangle &= \prod_{m,x} \delta(\varphi_m(x) - \varphi'_m(x)) = \delta(\varphi - \varphi'), \\ \langle\pi|\pi'\rangle &= \prod_{m,x} \delta(\pi_m(x) - \pi'_m(x)) = \delta(\pi - \pi'), \\ \langle\varphi|\pi\rangle &= \prod_{m,x} \frac{1}{\sqrt{2\pi}} \exp(i\pi_m(x)\varphi_m(x)). \end{aligned} \quad (1.2.3)$$

The notation in Eq. (1.2.3) can be a bit confusing. The products are over all flavors and spacetime points, and the delta functions are meant to collapse sums over all possible configurations of fields e.g.

$$\int \left(\prod_{m,x} d\varphi_m(x) \right) f(\varphi) \delta(\varphi' - \varphi) = f(\varphi'), \quad (1.2.4)$$

and similar for π .

However, we want to consider insertions of Heisenberg operators, not Schrödinger operators. Upgrading to

$$\Phi_m(X) = e^{iHt}\Phi_m(x)e^{-iHt}, \quad \Pi_m(X) = e^{iHt}\Pi_m(x)e^{-iHt}, \quad (1.2.5)$$

these will have eigenstates

$$\Phi_m(X) |\varphi\rangle_t = \varphi_m(x) |\varphi\rangle_t, \quad \Pi_m(X) |\pi\rangle_t = \pi_m(x) |\pi\rangle_t, \quad (1.2.6)$$

which satisfy the same inner products as in Eq. (1.2.3). Notice that in order for these to work out properly with the Heisenberg operators, these states must be given by the “wrong-sign” evolution

$$|\varphi\rangle_t = e^{iHt} |\varphi\rangle, \quad |\pi\rangle_t = e^{iHt} |\pi\rangle. \quad (1.2.7)$$

Another major thing to note is that the eigenvalues in Eq. (1.2.6) are time-independent; they are the same eigenvalues as those in the Schrödinger picture.

Since we have complete sets of eigenstates for both incompatible operators, we are free to define the identity operator with respect to both bases, i.e.

$$\int \left(\prod_{m,x} d\varphi_m(x) \right) |\varphi\rangle_t \langle\varphi| = \int \left(\prod_{m,x} d\pi_m(x) \right) |\pi\rangle_t \langle\pi| = \mathbb{1}. \quad (1.2.8)$$

The final ingredient is to first recognize that the Hamiltonian commutes with itself, so the Schrödinger picture and Heisenberg picture Hamiltonians are equivalent, i.e. $H(\Phi(x), \Pi(x)) = H(\Phi(X), \Pi(X))$. Then we can always use the canonical commutation relations between Φ and Π to push Π to the right of Φ in all terms in the Hamiltonian.

With all this in place, we can return to the problem at hand: calculating the time-ordered product of vacuum-to-vacuum correlation functions. Consider the first step where we propagate from the vacuum to the first operator insertion $\Phi_{m_1}(X_1)$, i.e.

$$\Phi_{m_1}(X_1) |\Omega\rangle = \int \mathcal{D}\varphi_1 \Phi_{m_1}(X_1) |\varphi_1\rangle_{t_1} \langle\varphi_1|\Omega\rangle = \int \mathcal{D}\varphi_1 |\varphi_1\rangle_{t_1} \varphi_1(x_1)_{t_1} \langle\varphi_1|\Omega\rangle. \quad (1.2.9)$$

This is all well and good, but as we said previously, we can only really get a good definition of $|\Omega\rangle$ for time going to the infinite future or the infinite past. Clearly, in this case, we can't simply propagate from $t \rightarrow -\infty$ to t_1 in a single step. However, what we can do is consider many different smaller steps of size dt using

$$_{t+dt}\langle\varphi'|\varphi\rangle_t = {}_t\langle\varphi'|e^{-iH(\Phi_m(X),\Pi_m(x))dt}|\varphi\rangle_t = \int \left(\prod_{m,x} d\pi_m(x) \right) {}_t\langle\varphi'|e^{-iH(\Phi(X),\Pi(X))dt}|\pi\rangle_t {}_t\langle\pi|\varphi\rangle_t. \quad (1.2.10)$$

Since we have ordered our Hamiltonian so that momentum operators are always on the right, we can simply replace the operators in $H(\Phi_m(X),\Pi_m(X))$ with their corresponding *classical* field counterparts (or the eigenvalues) $\varphi_m(x)$ and $\pi_m(x)$ and use the inner product between momentum and field states to find

$$\begin{aligned} {}_{t+dt}\langle\varphi'|\varphi\rangle_t &= \int \left(\prod_{m,x} d\pi_m(x) \right) \exp(-iH(\varphi'_m(x),\pi_m(x))dt) {}_t\langle\varphi'|\pi\rangle_t {}_t\langle\pi|\varphi\rangle_t \\ &= \int \left(\prod_{m,x} \frac{d\pi_m(x)}{2\pi} \right) \exp \left(-iH(\varphi'_m(x),\pi_m(x))dt + i \sum_n \int dx [\varphi'_n(x) - \varphi_n(x)]\pi_m(x) \right) \\ &= \int \left(\prod_{m,x} \frac{d\pi_m(x)}{2\pi} \right) \exp \left[i dt \left(\sum_n \int dx \frac{[\varphi'_n(x) - \varphi_n(x)]}{dt} \pi_n(x) - H(\varphi'_m(x),\pi_m(x)) \right) \right] \end{aligned} \quad (1.2.11)$$

where we converted products over field flavor and space “indices” of exponentials into sums/integrals inside the argument of the exponential. Keep in mind that, in this inner product, $\varphi_m(x)$ and $\pi_m(x)$ are the classical field configurations at time t while $\varphi'_m(x)$ is the configuration at time $t+dt$. In the final step, we have simply taken a factor of dt out of all terms in the exponent.

Using this, we can now evaluate the inner product of any two time-ordered field eigenstates by splitting up the time interval into N parts each separated by dt using the completeness relations of the eigenstates

$$\begin{aligned} {}_{t_{N+1}}\langle\varphi^{N+1}|\varphi^0\rangle_{t_0} &= \int \left(\prod_{k=1}^N \prod_{m,x} d\varphi_m^k(x) \right) {}_{t_{N+1}}\langle\varphi^{N+1}|\varphi^N\rangle_{t_N} {}_{t_N}\langle\varphi^N|\varphi^{N-1}\rangle_{t_{N-1}} \times \cdots \times {}_{t_1}\langle\varphi^1|\varphi^0\rangle_{t_0} \\ &= \int \left(\prod_{k=1}^N \prod_{m,x} d\varphi_m^k(x) \right) \int \left(\prod_{k=1}^N \prod_{m,x} \frac{d\pi_m^k(x)}{2\pi} \right) \times \\ &\quad \times \exp \left[i \sum_{k=0}^{N+1} dt \left(\sum_n \int dx \frac{\varphi_n^{k+1}(x) - \varphi_n^k(x)}{dt} \pi_n^k(x) - H(\varphi_m^{k+1}(x),\pi_m^k(x)) \right) \right]. \end{aligned} \quad (1.2.12)$$

We now take $dt \rightarrow 0$ in which case, $\varphi_m^k(x) \rightarrow \varphi_m(X)$ ⁵ where the time component of X is t_k , and therefore the ratio

$$\frac{\varphi_m^{k+1}(x) - \varphi_m^k(x)}{dt} \stackrel{dt \rightarrow 0}{\equiv} \frac{\partial \varphi_m(X)}{\partial t} \equiv \dot{\varphi}_m(X), \quad (1.2.13)$$

and the discrete sum over time indices becomes a time integral. We may then write

$${}_{t_{N+1}}\langle\varphi^{N+1}|\varphi^0\rangle_{t_0} = \int_{t_0}^{t_{N+1}} \prod_m \mathcal{D}\varphi_m \mathcal{D}\pi_m \exp \left[i \int_{t_0}^{t_{N+1}} dt \left(\sum_n \int dx \dot{\varphi}_n(X) \pi_n(X) - H(\varphi_m(X),\pi_m(X)) \right) \right], \quad (1.2.14)$$

where $\mathcal{D}\varphi_m$ indicates the *path integral* over all possible field configurations between time t_0 and t_{N+1} with the endpoints fixed by φ^0 and φ^{N+1} , respectively.

Now, we need to approach the operator insertions. This is straightforward, especially if we only consider

⁵Really, we introduce a “smoothing function” which interpolates between the points at each time.

field operators. Then, we can always insert identity operators so that

$$\begin{aligned} \langle \Omega | \Phi_{m_n}(X_n) \dots \Phi_{m_1}(X_1) | \Omega \rangle &= \int \left(\prod_{m_{n+1}, x_{n+1}} d\varphi_{m_{n+1}}^{n+1}(x_{n+1}) \right) \times \dots \times \int \left(\prod_{m_0, x_n} d\varphi_{m_0}^0(x_0) \right) \times \\ &\times \langle \Omega | \varphi^{n+1} \rangle_{t_{n+1} t_{n+1}} \langle \varphi^{n+1} | \Phi_{m_n}(X_n) | \varphi^n \rangle_{t_n} \times \dots \times_{t_2} \langle \varphi^2 | \Phi_{m_1}(X_1) | \varphi^1 \rangle_{t_1 t_1} \langle \varphi^1 | \varphi^0 \rangle_{t_0 t_0} \langle \varphi^0 | \Omega \rangle, \end{aligned} \quad (1.2.15)$$

where we have included two additional insertions of the identity for reasons that will become clear soon. If we had more general momentum-dependent operators, we would simply order them in the opposite way of the Hamiltonian (momentum operators on the left) and insert complete momentum states on the left side of the operator insertions. Using the eigenstate conditions, the field operators are now turned into classical fields evaluated at spacetime position X_ℓ , and we find

$$\begin{aligned} \langle \Omega | \Phi_{m_n}(X_n) \dots \Phi_{m_1}(X_1) | \Omega \rangle &= \int \left(\prod_{m_{n+1}, x_{n+1}} d\varphi_{m_{n+1}}^{n+1}(x_{n+1}) \right) \times \dots \times \int \left(\prod_{m_0, x_n} d\varphi_{m_0}^0(x_0) \right) \times \\ &\times \langle \Omega | \varphi^{n+1} \rangle_{t_{n+1} t_{n+1}} \langle \varphi^{n+1} | \varphi^n \rangle_{t_n} \times \dots \times_{t_2} \langle \varphi^2 | \varphi^1 \rangle_{t_1 t_1} \langle \varphi^1 | \varphi^0 \rangle_{t_0 t_0} \langle \varphi^0 | \Omega \rangle \varphi_{m_n}(X_n) \times \dots \times \varphi_{m_1}(X_1). \end{aligned} \quad (1.2.16)$$

We know how to evaluate each inner product with Eq. (1.2.14) and the additional integrals over field configurations at the times of operator insertions is exactly an integral over the endpoints of each inner product. Defining, $t_0 = -T$ and $t_{n+1} = T$, we have

$$\begin{aligned} \langle \Omega | \Phi_{m_n}(X_n) \dots \Phi_{m_1}(X_1) | \Omega \rangle &= \int \prod_m \mathcal{D}\varphi_m \mathcal{D}\pi_m \exp \left[i \int_{-T}^T dt \left(\sum_n \int dx \dot{\varphi}_n(X) \pi_n(X) - H(\varphi_m(X), \pi_m(X)) \right) \right] \times \\ &\times \langle \Omega | \varphi \rangle_T \langle \varphi | \Omega \rangle_{(-T)} \varphi_{m_n}(X_n) \times \dots \varphi_{m_1}(X_1), \end{aligned} \quad (1.2.17)$$

where we note that the field configurations at $t = \pm T$ are *also* integrated over, denoted by not having bounds on the path integrals.

The final ingredient is to evaluate the inner products between the field eigenstates and the vacuum states at $t = \pm T$. This is where our weak-coupling assumption comes into play. If we take the limit where $T \rightarrow \infty$, the weakly interacting theory becomes arbitrarily close to the free theory. In that case, we can rearrange Eq. (0.0.3) (or the corresponding expressions for complex fields) to find an expression for the free-field annihilation operators

$$a_{m,p} \stackrel{X^0 \rightarrow \pm\infty}{=} e^{i\omega_p X^0} \int dx \left[\sqrt{\frac{\omega_p}{2}} \Phi_m(X) + i \frac{1}{\sqrt{2\omega_p}} \Pi_m(X) \right] e^{-ipx}. \quad (1.2.18)$$

We also know that this operator will annihilate the vacuum in these limits

$$a_{m,p} |\Omega\rangle \stackrel{T \rightarrow \pm\infty}{=} 0. \quad (1.2.19)$$

In field-space (as opposed to conjugate momentum space), the momentum operator acts on states like a field derivative exactly analogous to the P -operator in position space for single-particle systems, denoted

$$\Pi_m(X) |\varphi\rangle_t = -i \frac{\delta}{\delta\varphi(x)} |\varphi\rangle_t, \quad {}_t \langle \varphi | \Pi_m(X) = -i \frac{\delta}{\delta\varphi(x)} {}_t \langle \varphi |, \quad (1.2.20)$$

so, we may write

$$\lim_{T \rightarrow \pm\infty} {}_T \langle \varphi | a_{m,p} | \Omega \rangle = \lim_{X^0 \rightarrow \infty} \frac{e^{i\omega_p X^0}}{\sqrt{4\pi\omega_p}} \int dx \left[\frac{\delta}{\delta\varphi_m(x)} + \omega_p \varphi_m(x) \right] e^{-ipx} {}_{X^0} \langle \varphi | \Omega \rangle = 0. \quad (1.2.21)$$

The differential equation under the integral is solved by the Gaussian⁶

$$\lim_{T \rightarrow \pm\infty} {}_T\langle \varphi | \Omega \rangle = \mathcal{N} \exp \left[-\frac{1}{2} \sum_{mn} \int dx \int dy \mathcal{E}_{mn}(x, y) \varphi_m(x) \varphi_n(y) \right], \quad (1.2.22)$$

as long as the kernel $\mathcal{E}_{mn}(x, y)$ satisfies

$$\mathcal{E}_{mn}(x, y) = \delta_{mn} \int \frac{dp}{2\pi} e^{ip(x-y)} \omega_p, \quad (1.2.23)$$

and so is just the position-space representation of the energy of the system. The normalization \mathcal{N} is yet undetermined, but we will consider that momentarily. With this, the product of wavefunctions becomes

$$\lim_{T \rightarrow \infty} \langle \Omega | \varphi \rangle_T {}_{(-T)}\langle \varphi | \Omega \rangle = \lim_{T \rightarrow \infty} |\mathcal{N}|^2 \exp \left[-\frac{1}{2} \sum_{m,n} \int dx dy \mathcal{E}_{mn}(x, y) \left(\varphi_m(x, T) \varphi_n(y, T) + \varphi_m(x, -T) \varphi_n(y, -T) \right) \right]. \quad (1.2.24)$$

Finally, we use a trick (assuming smooth $f(t)$):

$$\lim_{T \rightarrow \infty} (f(t) + f(-t)) = \lim_{\epsilon \rightarrow 0^+} \epsilon \int_{-\infty}^{\infty} dt f(t) e^{-\epsilon|t|}, \quad (1.2.25)$$

to allow us to include these terms into the exponential factor of the path integral underneath the time integral, i.e.

$$\begin{aligned} \langle \Omega | \Phi_{m_n}(X_n) \dots \Phi_{m_1}(X_1) | \Omega \rangle &= |\mathcal{N}|^2 \lim_{\epsilon \rightarrow 0^+} \int \prod_m \mathcal{D}\varphi_m \mathcal{D}\pi_m \varphi_{m_1}(X_1) \times \dots \times \varphi_{m_n}(X_n) \times \\ &\times \exp \left[i \int_{-\infty}^{\infty} dt \left(\sum_n \int dx \dot{\varphi}_n(x, t) \pi_n(x, t) - H(\varphi_m(X), \pi_m(X)) \right. \right. \\ &\quad \left. \left. + \frac{i}{2} \epsilon \sum_{m,n} \int dx dy \mathcal{E}_{mn}(x, y) \varphi_m(x, t) \varphi_n(y, t) e^{-\epsilon|t|} \right) \right]. \end{aligned} \quad (1.2.26)$$

This expression is quite cumbersome, but it is the correct result. To make it a bit more wieldy, we will not always write the last term, but instead refer to it as “ $i\epsilon$ terms” and crucially remember that *it is quadratic in the fields*.

We can fix our normalization so that the vacuum is normalized, $\langle \Omega | \Omega \rangle = 1$, in which case we find that $1/|\mathcal{N}|^2$ is given by the path integral with no field insertions.

Eq. (1.2.26) is the final result for a general classical Hamiltonian, H . However, for many (and really, most interesting) physical systems, the Hamiltonian is at most quadratic in momenta, $\pi_m(X)$. In this case, we can write

$$H(\varphi_\ell(Z), \pi_\ell(Z)) = \frac{1}{2} \sum_{mn} \int dx \int dy M_{mn}(x, y) \pi_m(x, t) \pi_n(y, t) + \sum_m \int dx N_m(x) \pi_m(x, t) + C \quad (1.2.27)$$

where M_{mn} , N_m , and C can all depend on the fields $\varphi_m(x, t)$ in a complicated way. Then, the difference appearing in our correlation function can be written as

$$\begin{aligned} \sum_n \int dt \int dx \dot{\varphi}_n(x, t) \pi_n(x, t) - H(\varphi_\ell(Z), \pi_\ell(Z)) &= \\ -\frac{1}{2} \sum_{mn} \int d^2X d^2Y \mathcal{M}_{mn}(X, Y) \pi_m(X) \pi_n(Y) - \sum_m \int d^2X \mathcal{N}_m(X) \pi_m(X) - \mathcal{C}, \end{aligned} \quad (1.2.28)$$

⁶This will look slightly different in the case where we have complex fields or spinors, but at the end, what is important is that it is quadratic in fields.

with

$$\begin{aligned}\mathcal{M}_{mn}(X, Y) &= M_{mn}(X, Y)\delta(X^0 - Y^0), \\ \mathcal{N}_m(X) &= N_m(X) - \dot{\varphi}_m(X), \\ \mathcal{C} &= \int dt C,\end{aligned}\tag{1.2.29}$$

again remembering that M , N , and C all depend on $\varphi_\ell(Z)$. We have also defined the two-dimensional spacetime integrals, e.g. d^2X .

We won't prove it here (see Weinberg for such a proof), but Gaussian integrals of this form are proportional to the integrand with the integrated variable replaced by the stationary point of the exponent, i.e.

$$\begin{aligned}& \int \prod_m \mathcal{D}\pi_m \exp \left[-i \left(\frac{1}{2} \sum_{mn} \int d^2X d^2Y \mathcal{M}_{mn}(X, Y) \pi_m(X) \pi_n(Y) + \sum_m \int d^2X \mathcal{N}_m(X) \pi_m(X) + \mathcal{C} \right) \right] \\ & \propto \exp \left[-i \left(\frac{1}{2} \sum_{mn} \int d^2X d^2Y \mathcal{M}_{mn}(X, Y) \bar{\pi}_m(X) \bar{\pi}_n(Y) + \sum_m \int d^2X \mathcal{N}_m(X) \bar{\pi}_m(X) + \mathcal{C} \right) \right],\end{aligned}\tag{1.2.30}$$

where $\bar{\pi}_m(X)$ is momentum evaluated at the stationary point of the exponent, given by the solution to

$$\frac{\delta}{\delta \pi_r(W)} \left(\frac{1}{2} \sum_{mn} \int d^2X d^2Y \mathcal{M}_{mn}(X, Y) \pi_m(X) \pi_n(Y) + \sum_m \int d^2X \mathcal{N}_m(X) \pi_m(X) + \mathcal{C} \right) \Big|_{\pi=\bar{\pi}} = 0. \tag{1.2.31}$$

In this particular case, this integrand is just $\dot{\varphi}\pi$ minus the Hamiltonian (the $i\epsilon$ terms are π -independent), so $\bar{\pi}$ can be found as the solution to

$$\dot{\varphi}_m(X) = \frac{\delta H}{\delta \pi_m(X)} \Big|_{\pi=\bar{\pi}}. \tag{1.2.32}$$

Since we are only considering field operator insertions, we can *always* perform this integral as long as the Hamiltonian is quadratic in momenta, and the proportionality constant will always cancel between the main path integral and the normalization. Furthermore, once we replace the $\bar{\pi}$'s with corresponding $\dot{\varphi}$'s, we end up with a function known as the *Lagrangian*

$$L(\varphi_\ell(Z), \dot{\varphi}_\ell(Z)) = \sum_m \int dx \dot{\varphi}_m(X) \bar{\pi}_m(X) - H(\varphi_\ell(Z), \bar{\pi}_\ell(Z)). \tag{1.2.33}$$

and the time integral of the Lagrangian is given the special name of the *action*

$$S[\varphi_\ell(Z)] = \int dt L(\varphi_\ell(z, t), \dot{\varphi}_\ell(z, t)) = \int d^2X \mathcal{L}(\varphi_\ell(X), \dot{\varphi}_\ell(X)), \tag{1.2.34}$$

where in the second step, we defined the *Lagrangian density*, though we will often just refer to this as the Lagrangian.

With all of this, we can write our path integral in a more convenient form

$$\langle \Omega | \Phi_{m_n}(X_n) \dots \Phi_{m_1}(X_1) | \Omega \rangle = \mathcal{A} \int \prod_m \mathcal{D}\varphi_m \varphi_{m_1}(X_1) \times \dots \times \varphi_{m_n}(X_n) e^{i(S + i\epsilon \text{ terms})}, \tag{1.2.35}$$

where \mathcal{A} is the normalization with the momenta integrated out.

An important thing to note in evaluating the integral over momenta is that we assumed we could pull the proportionality constant in Eq. (1.2.30) outside of the φ integral for it to cancel the term in the normalization. It turns out that this is equivalent to assuming that \mathcal{M} was field-independent. This is going to be the case for many theories, but not all as it turns out, so we must be very cautious when actually performing this step.

1.3 Free Theories with the Path Integral

Before moving on to interacting theories, we should first familiarize ourselves with using the path integral in the context of the simpler case of the free theory. Examining Eq. (0.0.1), we see that the Hamiltonian is indeed quadratic in momentum $\pi(x)$, so we may freely move to the Lagrangian description. Varying the Hamiltonian, we find

$$\left. \frac{\delta H_0}{\delta \pi(X)} \right|_{\pi=\bar{\pi}} = \bar{\pi}(X) = \dot{\phi}(X). \quad (1.3.1)$$

Plugging this result in for $\bar{\pi}$, we find the result for the Lagrangian density

$$\mathcal{L} = \frac{1}{2} \left[\left(\frac{\partial \phi}{\partial t} \right)^2 - \left(\frac{\partial \phi}{\partial x} \right)^2 - m_\phi^2 \phi^2 \right] = \frac{1}{2} \partial_\mu \phi \partial^\mu \phi - \frac{m_\phi^2}{2} \phi^2, \quad (1.3.2)$$

where we have used the short-hand notation $\partial_\mu = \partial/\partial X^\mu$. Notice that the Lagrangian turns out to be naturally Lorentz-covariant, unlike the Hamiltonian. This makes it far more useful to work with than the Hamiltonian since it now becomes very simple to construct Lorentz-invariant theories!

We will next pass the action to its corresponding Fourier-transformed partner

$$S = \int d^2 X \mathcal{L} = -\frac{1}{2} \int d^2 X \int \frac{d^2 P_1}{(2\pi)^2} \int \frac{d^2 P_2}{(2\pi)^2} (P_1 \cdot P_2 + m_\phi^2) \tilde{\phi}(P_1) \tilde{\phi}(P_2) e^{-iX \cdot (P_1 + P_2)} \quad (1.3.3)$$

Evaluating the integral over X gives a two-momentum delta function $(2\pi)^2 \delta^{(2)}(P_1 + P_2)$,

$$S = \frac{1}{2} \int \frac{d^2 P_1}{(2\pi)^2} \int \frac{d^2 P_2}{(2\pi)^2} (P_1^2 - m_\phi^2) (2\pi)^2 \delta^{(2)}(P_1 + P_2) \tilde{\phi}(P_1) \tilde{\phi}(P_2), \quad (1.3.4)$$

where it will become evident as to why we chose not to collapse the delta function yet. We have also defined the short-hand notation $P^\mu P_\mu = P \cdot P = P^2$.

Now, let's revisit the two-point correlation function we calculated earlier. We know that this is given by

$$\int dx' f(x' - x) \langle \Omega | \phi(Y) \phi(X') | \Omega \rangle = \mathcal{A} \int dx' f(x' - x) \int \mathcal{D}\phi \phi(Y) \phi(X') e^{i(S + i\epsilon \text{ terms})} \quad (1.3.5)$$

Since the action and $i\epsilon$ terms are purely quadratic in the fields, we can evaluate this in a straightforward way using Gaussian integration techniques (again, see Weinberg for explicit details). In particular, we can evaluate a Gaussian integral of the form

$$\begin{aligned} & \int \prod_m \mathcal{D}\varphi_m \varphi_{m_1}(X_1) \varphi_{m_2}(X_2) \dots \exp \left(-\frac{i}{2} \sum_{mn} \int d^2 X \int d^2 Y \varphi_m(X) A_{mn}(X, Y) \varphi_n(Y) \right) \\ &= \frac{1}{\mathcal{A}} \sum_{\text{pairs}(\{(m_1, X_1), (m_2, X_2), \dots\})} \left(-i \prod_{\{(n_1, Y_1), (n_2, Y_2)\} \in \text{pairs}} A_{n_1, n_2}^{-1}(Y_1, Y_2) \right) \end{aligned} \quad (1.3.6)$$

where “pairs” refers to all possible ways of pairing the indices of the fields in the integrand of the path integral. Note, that the X_i 's and Y_i 's in Eq. (1.3.6) can either refer to spacetime points or two-momenta, depending on which space we are working in.

Examining Eq. (1.3.4), we see that the “matrix” we need to invert is given in momentum space by

$$\tilde{A}(P_1, P_2) = -(P_1^2 - m_\phi^2) (2\pi)^2 \delta^{(2)}(P_1 + P_2) + i\epsilon \text{ terms}. \quad (1.3.7)$$

The inverse of this must satisfy

$$\int \frac{d^2 Q}{(2\pi)^2} \tilde{A}^{-1}(P_1, Q) \tilde{A}(Q, P_2) = (2\pi)^2 \delta^{(2)}(P_1 - P_2), \quad (1.3.8)$$

which we can explicitly verify is satisfied by

$$\tilde{A}^{-1}(P_1, P_2) = -\frac{(2\pi)^2 \delta^{(2)}(P_1 + P_2)}{P_1^2 - m_\phi^2 + i\epsilon}. \quad (1.3.9)$$

or, in position space,

$$A^{-1}(X, Y) = - \int \frac{d^2 P}{(2\pi)^2} \frac{e^{-iP \cdot (X-Y)}}{P^2 - m_\phi^2 + i\epsilon}. \quad (1.3.10)$$

Using this, we can evaluate our two-point function by Fourier-transforming the fields, giving (dropping the $f(x' - x)$ piece for now)

$$\int \mathcal{D}\phi \int \frac{d^2 P_1}{(2\pi)^2} \int \frac{d^2 P_2}{(2\pi)^2} \tilde{\phi}(P_1) \tilde{\phi}(P_2) e^{-iP_1 \cdot X'} e^{-iP_2 \cdot Y} e^{i(S+i\epsilon \text{ terms})} = \int \frac{d^2 P}{(2\pi)^2} \frac{ie^{-iP \cdot (Y-X')}}{P^2 - m_\phi^2 + i\epsilon}. \quad (1.3.11)$$

The denominator has two poles in P^0 which we can see by splitting up $P^2 = (P^0)^2 - p^2$ and using

$$P^2 - m_\phi^2 + i\epsilon = (P^0 - \sqrt{p^2 + m_\phi^2} + i\epsilon)(P^0 + \sqrt{p^2 + m_\phi^2} - i\epsilon). \quad (1.3.12)$$

The poles therefore lie at $P_0 = \pm \sqrt{p^2 + m_\phi^2} \mp i\epsilon$. The integral over P^0 can then be evaluated using contour integration methods. In particular, since $Y^0 > X'^0$, we need to ensure that the imaginary part of the pole that is enclosed by our contour is negative, which is only satisfied by the positive-energy pole (i.e. $P^0 = \omega_p - i\epsilon$). We therefore find that the two-point function is given by (after taking $\epsilon \rightarrow 0$)

$$\int dx' f(x' - x) \langle \Omega | \phi(Y) \phi(X') | \Omega \rangle = \int dx' f(x' - x) \int \frac{dp}{2\pi} \frac{1}{2\omega_p} e^{-i(\omega_p(Y^0 - X'^0) - p(y - x'))}, \quad (1.3.13)$$

which is exactly what we found previously!

1.4 Small couplings to the rescue...again!

The free-field case is very nice because it results in purely Gaussian path integrals, which we know how to evaluate analytically. When we add interactions (even weak ones!) this is no longer the case: in fact, we *must* include non-Gaussian terms in order to introduce interaction terms which will include three or more fields meeting at the same spacetime point. We can, however use the fact that, as long as the interactions are suitably weak, we can split the action as $S = S_0 + S_I$ where

$$S_0 = \int d^2 X \mathcal{L}_{\text{free}}, \quad (1.4.1)$$

is the free action and S_I is the piece of the action which contains the interactions. In the case where $S_I \ll S_0$, we can use two very nice properties of the path integral formalism: first, the action is in terms of classical fields, not field operators, so we do not need to worry about commutation relations as long as we work under the path integral. Second, since the action is exponentiated in the path integral, as long as S_I is polynomial in the fields, we can expand the exponential of the action as

$$e^{i(S+i\epsilon \text{ terms})} = e^{i(S_0+i\epsilon \text{ terms})} \sum_{n=0}^{\infty} \frac{(iS_I)^n}{n!}, \quad (1.4.2)$$

which has the form of a polynomial of fields times a Gaussian, which we already saw how to solve exactly! Moreover, each term has more powers of the action, and therefore becomes smaller (assuming $S_I \ll 1$), as we expand to higher orders. This is exactly the perturbative expansion of a QFT.

2 A Simple Model

To start working with an interacting QFT, we will consider two fields, ϕ and θ , with only a single four-point interaction between the fields. This will be described by the Lagrangian

$$\mathcal{L} = \frac{1}{2} \left(\partial_\mu \phi \partial^\mu \phi + \partial_\mu \theta \partial^\mu \theta - m_\phi^2 \phi^2 - m_\theta^2 \theta^2 \right) - \frac{\lambda_{\phi\theta}}{4} \phi^2 \theta^2. \quad (2.0.1)$$

After integrating over X the term in parentheses will become S_0 while the term proportional to $\lambda_{\phi\theta}$ will give S_I . As it turns out, this situation will only be truly stable for $\lambda_{\phi\theta} \geq 0$, so we will restrict ourselves to this range (for negative values of $\lambda_{\phi\theta}$ and $m_\phi^2, m_\theta^2 > 0$, it will actually be metastable, but tunnelling effects will allow it to decay).

As we have seen, it is quite useful to work in momentum space instead of position space, so we will consider the momentum-space version of S_I , given by

$$S_I = -\frac{\lambda_{\phi\theta}}{4} \int \left(\prod_{n=1}^4 \frac{d^2 K_n}{(2\pi)^2} \right) \tilde{\phi}(K_1) \tilde{\phi}(K_2) \tilde{\theta}(K_3) \tilde{\theta}(K_4) (2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4), \quad (2.0.2)$$

where the delta function arises from the integral over X of $\exp[-iX \cdot (K_1 + K_2 + K_3 + K_4)]$. This delta function serves to conserve two-momentum for any interactions.

2.1 Two-Point Functions

As we will see, for any n -point correlation function, we will always end up with some subset consisting of the two-point functions

$$\Pi_\phi(X, Y) = \langle \Omega | \mathbf{T} \phi(Y) \phi(X) | \Omega \rangle, \quad \Pi_\theta(X, Y) = \langle \Omega | \mathbf{T} \theta(Y) \theta(X) | \Omega \rangle. \quad (2.1.1)$$

Since these will not only be extremely useful for future calculations, but are naively (we will see soon that they are not so simple after all!) the simplest n -point functions to compute, we will use them to demonstrate the techniques that will be useful for higher-point correlation functions. It will be useful to split up these correlation functions according to their powers of $\lambda_{\phi\theta}$, each power containing additional powers of S_I in the path integral,

$$\Pi_i(X, Y) = \Pi_i^{(0)}(X, Y) + \frac{\lambda_{\phi\theta}}{4\pi} \Pi_i^{(1)}(X, Y) + \left(\frac{\lambda_{\phi\theta}}{4\pi} \right)^2 \Pi_i^{(2)}(X, Y) + \dots \quad (2.1.2)$$

The extra factors of $1/4\pi$ will become apparent soon. We have already found the zero-order terms, which are exactly the free propagators

$$\Pi_i^{(0)}(X, Y) = \int \frac{d^2 P}{(2\pi)^2} \tilde{\Pi}_i^{(0)}(P^2) e^{-iP \cdot (Y-X)} = \int \frac{d^2 P}{(2\pi)^2} \frac{i e^{-iP \cdot (Y-X)}}{P^2 - m_i^2 + i\epsilon}. \quad (2.1.3)$$

Our next task will then be to compute the higher-order contributions to Π_i .

2.1.1 Leading Order

Although one can argue that the zero-order pieces are the leading-order contributions, it tends to make a bit more sense to count the number of couplings included in the computation, so we will use “leading order” to mean “single-coupling insertion.” Then, in momentum space, the leading-order two-point function for ϕ is given by

$$\begin{aligned} \Pi_\phi^{(1)}(X, Y) = & -\frac{i\lambda_{\phi\theta}}{4} \int \mathcal{D}\phi \mathcal{D}\theta \int \frac{d^2 P_1}{(2\pi)^2} \int \frac{d^2 P_2}{(2\pi)^2} \tilde{\phi}(P_1) \tilde{\phi}(P_2) e^{-iX \cdot P_1} e^{-iY \cdot P_2} \times \\ & \times \int \left(\prod_{n=1}^4 \frac{d^2 K_n}{(2\pi)^2} \right) \tilde{\phi}(K_1) \tilde{\phi}(K_2) \tilde{\theta}(K_3) \tilde{\theta}(K_4) (2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4) e^{i(S_0 + i\epsilon \text{ terms})}. \end{aligned} \quad (2.1.4)$$

Now, we simply use our rules for Gaussian integration and sum over all possible ways of connecting similar fields with the free momentum-space propagator, Eq. (1.3.9). Unfortunately, this is a combinatorial problem,

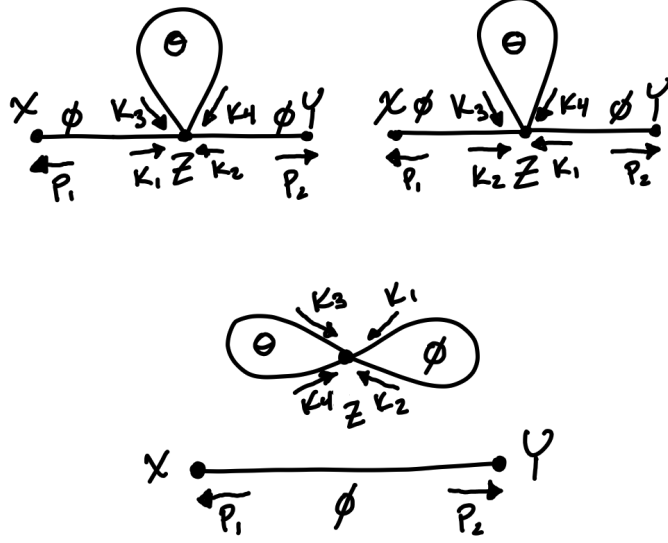


Figure 1: Feynman diagrams contributing to “leading-order” two-point correlator.

so the number of contractions grows quite fast. Luckily, in this case, we only get three terms

$$\begin{aligned}
 \Pi_{\phi}^{(1)}(X, Y) = & -\frac{i\lambda_{\phi\theta}}{4} \int \frac{d^2 P_1}{(2\pi)^2} \frac{d^2 P_2}{(2\pi)^2} \left(\prod_{n=1}^4 \frac{d^2 K_n}{(2\pi)^2} \right) \frac{i(2\pi)^2 \delta^{(2)}(K_3 + K_4)}{K_3^2 - m_{\theta}^2 + i\epsilon} \left[\right. \\
 & \frac{i(2\pi)^2 \delta^{(2)}(K_1 + P_1) i(2\pi)^2 \delta^{(2)}(K_2 + P_2)}{(K_1^2 - m_{\phi}^2 + i\epsilon)(K_2^2 - m_{\phi}^2 + i\epsilon)} + \frac{i(2\pi)^2 \delta^{(2)}(K_1 + P_2) i(2\pi)^2 \delta^{(2)}(K_2 + P_1)}{(K_1^2 - m_{\phi}^2 + i\epsilon)(K_2^2 - m_{\phi}^2 + i\epsilon)} \\
 & \left. + \frac{i(2\pi)^2 \delta^{(2)}(K_1 + K_2) i(2\pi)^2 \delta^{(2)}(P_1 + P_2)}{(K_1^2 - m_{\phi}^2 + i\epsilon)(P_1^2 - m_{\phi}^2 + i\epsilon)} \right] (2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4) e^{-i(P_1 \cdot X + P_2 \cdot Y)}.
 \end{aligned} \tag{2.1.5}$$

From this expression, we can see that it is possible to represent each of these terms with a diagram: we have fixed points at X and Y representing our external states, and a single internal vertex at some other point (call it Z) coming from the insertion of the interaction term. We then connect the lines from the vertex in all possible ways with corresponding lines (either from the external points or from other lines on the vertex itself). We can get the corresponding mathematical expression by replacing the vertex with $-i\lambda_{\phi\theta}/4(2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4)$, and each line with the free two-point function. Finally, we integrate over all momenta, including the extra Fourier exponentials for the external states. This method of generating all of such diagrams is known as the Feynman diagram expansion, and the corresponding rules for replacing each part of a diagram with mathematical expressions are known as the Feynman rules. For this theory,

then, the Feynman rules are given by⁷:

$$\begin{aligned}
\text{Two-point } (\phi(P_1), \phi(P_2)) : & \quad \frac{i(2\pi)^2 \delta^{(2)}(P_1 + P_2)}{P_1^2 - m_\phi^2 + i\epsilon} \\
\text{Two-point } (\theta(P_1), \theta(P_2)) : & \quad \frac{i(2\pi)^2 \delta^{(2)}(P_1 + P_2)}{P_1^2 - m_\phi^2 + i\epsilon} \\
\text{Four-point } (\phi(P_1), \phi(P_2), \theta(P_3), \theta(P_4)) : & \quad -\frac{i\lambda_{\theta\phi}}{4} (2\pi)^2 \delta^{(2)}(P_1 + P_2 + P_3 + P_4).
\end{aligned} \tag{2.1.6}$$

The diagrams for the three terms in Eq. (2.1.5) are shown in Fig. 1.

Examining Eq. (2.1.5), the second line can be simplified by the fact that the integrals are symmetric under $K_1 \leftrightarrow K_2$, so the two terms in the sum are actually the same, just giving a factor of two. Collapsing the delta functions, we arrive at the expression for the momentum-space two-point function

$$\begin{aligned}
\tilde{\Pi}_\phi^{(1)}(P) = & -\frac{\lambda_{\phi\theta}}{P^2 - m_\phi^2 + i\epsilon} \left[\frac{1}{2} \frac{1}{P^2 - m_\phi^2 + i\epsilon} \int \frac{d^2 K}{(2\pi)^2} \frac{1}{K^2 - m_\theta^2 + i\epsilon} \right. \\
& \left. + \frac{1}{4} \int \frac{d^2 K}{(2\pi)^2} \frac{d^2 L}{(2\pi)^2} \frac{(2\pi)^2 \delta^{(2)}(0)}{(K^2 - m_\theta^2 + i\epsilon)(L^2 - m_\phi^2 + i\epsilon)} \right].
\end{aligned} \tag{2.1.7}$$

Now, we have a problem: both of these terms are divergent. The second term features a far worse divergence than the first, proportional to $\delta^{(2)}(0)$, but it is also just proportional to the zero-order propagator, so we can simply view this as correction to the normalization of our constant vacuum energy, absorb it into \mathcal{A} , and ignore it entirely (again, remember that the vacuum energy is already a divergent quantity, but we can only really measure changes in energy, in which case this divergent piece will always cancel). While this is an important point to understand, and in some contexts can be physically important, these contributions are irrelevant to what we actually want to calculate, so we will ignore any diagrams that appear with such disconnected “vacuum bubbles.”

On the other hand, the first term, we cannot just get rid of and we have to treat with a bit of care because it actually affects the propagation (and other interactions) of the particle. We will discuss the treatment of this divergence more in the next section.

2.1.2 Renormalization

The divergence appearing in the first expression of Eq. (2.1.7) may be extremely disturbing at first glance, but as it turns out, it isn't as problematic as it seems. To see why, we need to again make sure we are asking physical questions. As a first taste of this, we will consider rescaling the fields in the theory by some field- and momentum-independent constant (the square roots are just convention)

$$\phi \rightarrow Z_\phi^{1/2} \phi, \quad \theta \rightarrow Z_\theta^{1/2} \theta. \tag{2.1.8}$$

Does this make any detectable change in the theory? Recall that the integrand of the path integral after doing a perturbative expansion in S_I will always be polynomial times a free-particle Gaussian. Then, to perform the integral, we connect any two like fields in the polynomial part with the inverse of the Gaussian operator (what we called A or \tilde{A}) in all possible ways. However, under the rescaling Eq. (2.1.8) each pair of fields in the polynomial part will come with a factor of Z_i , but the free action is also rescaled by a total factor of Z_i , since it is quadratic in the fields. This means that $A_i^{-1} \rightarrow 1/Z_i A^{-1}$ under such a rescaling and the Z_i 's cancel in each pair⁸! So, we find that *the overall normalization of the fields is an unphysical parameter*: any rescaling is guaranteed to drop out in calculations of correlation functions.

⁷Here, in the two-point function, we do not account for an additional factor of 4 which arises from the symmetry of the external legs, as is usually standard. This is because we wish to explicitly show how this arises from the pairing of indices in the path integral.

⁸It is important to note that the integration measures also change under such a rescaling, but since we divide by the zero-insertion path integral, the rescaling of the measures cancel between the numerator and denominator.

For other free parameters of the theory, like the couplings and the masses, we need to think about how we can actually measure these quantities. For couplings like $\lambda_{\phi\theta}$, we can only really measure such parameters by measuring the outcomes of certain experiments, such as $\phi\phi \rightarrow \theta\theta$ in this case. However, it is important to remember that, whether we like it or not, all orders of the perturbative expansion play out in nature. So, in such an experiment, we are not just measuring the leading-order, single- S_I insertion, so we see that the only physical thing we can extract from experiment is the combination of this single insertion (which is proportional to the classical parameter $\lambda_{\phi\theta}$) *plus all quantum corrections to the process*. So again, the $\lambda_{\phi\theta}$ that appears in the Lagrangian is not a physical parameter that we can unambiguously extract from experiment.

Finally, we will consider masses, which are perhaps the most concrete example of this. In this case, the physical thing that we can measure is the pole of the propagator, i.e. the mass of ϕ is when $P^2 = m_\phi^2$ (and similar for θ). However, again, we need to consider the quantum corrections to the free propagator as the particle travels from X to Y . In particular, we will consider the set of diagrams that cannot be cut along a single line to separate them, known as one-particle irreducible (1PI) diagrams. Then, the 1PI two-point function, $\tilde{\Pi}_i^{1\text{PI}}(P^2)$ is given by the sum of all such diagrams. We can then connect these 1PI two-point functions with free propagators to get the full two-point function

$$\begin{aligned} \tilde{\Pi}_i(P^2) = & \frac{i}{P^2 - m_i^2 + i\epsilon} + \frac{i}{P^2 - m_i^2 + i\epsilon} \tilde{\Pi}_i^{1\text{PI}}(P^2) \frac{i}{P^2 - m_i^2 + i\epsilon} \\ & + \frac{i}{P^2 - m_i^2 + i\epsilon} \tilde{\Pi}_i^{1\text{PI}}(P^2) \frac{i}{P^2 - m_i^2 + i\epsilon} \tilde{\Pi}_i^{1\text{PI}}(P^2) \frac{i}{P^2 - m_i^2 + i\epsilon} + \dots, \end{aligned} \quad (2.1.9)$$

or

$$\tilde{\Pi}_i(P^2) = \frac{i}{P^2 - m_i^2 + i\epsilon} \sum_{n=0}^{\infty} \left(\tilde{\Pi}_i^{1\text{PI}}(P^2) \frac{i}{P^2 - m_i^2 + i\epsilon} \right)^n = \frac{i}{P^2 - m_i^2 - i\tilde{\Pi}_i^{1\text{PI}}(P^2) + i\epsilon}, \quad (2.1.10)$$

where in the second step, we used the expression for the summation of a geometric series. From this, it is clear that the only physical thing we can measure is the *combination of m_i^2 and $\tilde{\Pi}_i^{1\text{PI}}(P^2)$* , which is again, the mass parameter appearing in the Lagrangian plus all additional quantum corrections.

So, we see that if we measure reasonably finite physical quantities (which again, are combinations of the Lagrangian parameters and the quantum corrections in e.g. $\tilde{\Pi}_i^{1\text{PI}}$), then any divergences that appear in the (unphysical) quantum corrections must cancel corresponding divergences in the (unphysical) Lagrangian parameters to yield a finite physical result. We stress that we are only playing around with unphysical quantities; other divergences can arise that have physical interpretations (and in fact, are not problematic at all), that we cannot get rid of.

A very nice methodology of book-keeping for all of these divergent quantities in a consistent way is known as *renormalization*. To renormalize the theory, we consider taking all of the unphysical parameters which appear in the original Lagrangian (denoted with a “(0)” superscript) and replace them with new, rescaled quantities

$$\phi^{(0)}(x) = \sqrt{Z_\phi} \phi(x), \quad \theta^{(0)}(x) = \sqrt{Z_\theta} \theta(x), \quad m_\phi^{(0)2} = Z_{m_\phi} m_\phi^2, \quad m_\theta^{(0)2} = Z_{m_\theta} m_\theta^2, \quad \lambda_{\phi\theta}^{(0)} = Z_\lambda \lambda_{\phi\theta}. \quad (2.1.11)$$

We now take into account the fact that we are doing a perturbative expansion in $\lambda_{\phi\theta}$, and quantum corrections are those that enter at higher orders in $\lambda_{\phi\theta}$. Another way of saying this is that the classical and quantum theories (and therefore the corresponding parameters) coincide if we neglect higher-order quantum corrections. To encapsulate this, we expand the rescaling factors as

$$Z_i = 1 + \sum_{n=1}^{\infty} \left(\frac{\lambda_{\phi\theta}}{4\pi} \right)^n \delta Z_i^{(n)} = 1 + \delta Z_i. \quad (2.1.12)$$

These δZ_i ’s are typically called “counterterms” because we choose them exactly to counteract the divergences in the quantum corrections. Then, the Lagrangian in Eq. (2.0.1) is given in terms of the rescaled parameters

as

$$\begin{aligned}
\mathcal{L} &= \frac{1 + \delta Z_\phi}{2} \left(\partial_\mu \phi \partial^\mu \phi - (1 + \delta Z_{m_\phi}) m_\phi^2 \phi^2 \right) + \frac{1 + \delta Z_\theta}{2} \left(\partial_\mu \theta \partial^\mu \theta - Z_{m_\theta} m_\theta^2 \theta^2 \right) \\
&\quad - (1 + \delta Z_\phi)(1 + \delta Z_\theta)(1 + \delta Z_\lambda) \frac{\lambda_{\phi\theta}}{4} \phi^2 \theta^2 \\
&= \frac{1}{2} \left(\partial_\mu \phi \partial^\mu \phi - m_\phi^2 \phi^2 \right) + \frac{1}{2} \left(\partial_\mu \theta \partial^\mu \theta - m_\theta^2 \theta^2 \right) - \frac{\lambda_{\phi\theta}}{4} \phi^2 \theta^2 \\
&\quad + \frac{\delta Z_\phi}{2} \partial_\mu \phi \partial^\mu \phi - (\delta Z_\phi + \delta Z_{m_\phi} + \delta Z_\phi \delta Z_{m_\phi}) \frac{m_\phi^2}{2} \phi^2 + \frac{\delta Z_\theta}{2} \partial_\mu \theta \partial^\mu \theta - (\delta Z_\theta + \delta Z_{m_\theta} + \delta Z_\theta \delta Z_{m_\theta}) \frac{m_\theta^2}{2} \theta^2 \\
&\quad - (\delta Z_\phi + \delta Z_\theta + \delta Z_\lambda + \delta Z_\phi \delta Z_\theta + \delta Z_\phi \delta Z_\lambda + \delta Z_\theta \delta Z_\lambda + \delta Z_\phi \delta Z_\theta \delta Z_\lambda) \frac{\lambda_{\phi\theta}}{4} \phi^2 \theta^2.
\end{aligned} \tag{2.1.13}$$

The first line of the second equality of Eq. (2.1.13) is just the original Lagrangian with classical parameters replaced with renormalized ones. The other terms we collect into a separate piece of the action that we denote S_{ct} , known as the “counterterm action.” Now, for the final trick: instead of splitting these into a quadratic piece in fields and an interaction piece, we will include *everything with δZ ’s into the interaction piece of the action*. This way, all propagators retain the same form as their unrenormalized counterparts, but we add an additional “counterterm action,” given in momentum space by

$$\begin{aligned}
S_{\text{ct}} &= \int \frac{d^2 K_1}{(2\pi)^2} \frac{d^2 K_2}{(2\pi)^2} \left(\frac{\delta Z_\phi}{2} K_1^2 - (\delta Z_\phi + \delta Z_{m_\phi} + \delta Z_\phi \delta Z_{m_\phi}) \frac{m_\phi^2}{2} \right) \tilde{\phi}(K_1) \tilde{\phi}(K_2) (2\pi)^2 \delta^{(2)}(K_1 + K_2) \\
&\quad + \int \frac{d^2 K_1}{(2\pi)^2} \frac{d^2 K_2}{(2\pi)^2} \left(\frac{\delta Z_\theta}{2} K_1^2 - (\delta Z_\theta + \delta Z_{m_\theta} + \delta Z_\theta \delta Z_{m_\theta}) \frac{m_\theta^2}{2} \right) \tilde{\theta}(K_1) \tilde{\theta}(K_2) (2\pi)^2 \delta^{(2)}(K_1 + K_2) \\
&\quad - \int \left(\prod_{n=1}^4 \frac{d^2 K_n}{(2\pi)^2} \right) (\delta Z_\phi + \delta Z_\theta + \delta Z_\lambda + \delta Z_\phi \delta Z_\theta + \delta Z_\phi \delta Z_\lambda + \delta Z_\theta \delta Z_\lambda + \delta Z_\phi \delta Z_\theta \delta Z_\lambda) \times \\
&\quad \times \frac{\lambda_{\phi\theta}}{4} \tilde{\phi}(K_1) \tilde{\phi}(K_2) \tilde{\theta}(K_3) \tilde{\theta}(K_4) (2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4).
\end{aligned} \tag{2.1.14}$$

This looks quite terrible, but remember that each counterterm we can treat as a perturbative expansion, so most of the terms won’t be relevant until higher orders anyway. Moreover, the counterterms must be momentum-independent (otherwise, we add derivatives to the terms in the action, and spoil the cancellations we need), so we can just treat these as constants, and they don’t actually add much more difficulty to the computation.

A key thing to note: if our theory is to be finite, the only unphysical (called ultraviolet for reasons we will discuss) divergences we run into should be able to be absorbed into these five counterterms. This means that we must fix each one with *a single process*, and once fixed, *all other ultraviolet divergences at that order in perturbation theory must cancel when including these counterterm interactions*. Typically, for simplicity, we choose to fix the counterterms using the n -point correlation functions that have the same field content as the term in the action, i.e. we fix δZ_ϕ and δZ_{m_ϕ} from the two- ϕ function, δZ_θ and δZ_{m_θ} from the two- θ function, and δZ_λ from the four-point $\phi^2 \theta^2$ function.

The next thing we need to address is the fact that there is a certain amount of arbitrariness to how we choose these counterterms. Consider we break up the amplitude that we are calculating into a piece that depends on counterterms, \mathcal{A}_{ct} and a piece which does not, $\mathcal{A}_{\text{bare}}$. Both of these parts of the amplitude will be momentum-dependent in general, so we must require that

$$\mathcal{A}_{\text{total}}(P_\ell) = \mathcal{A}_{\text{ct}}(\delta Z_\ell, P_\ell) + \mathcal{A}_{\text{bare}}(P_\ell) = \text{UV finite}. \tag{2.1.15}$$

We can notice a couple of things about this. Recall that the counterterms cannot be momentum-dependent. Therefore, the momentum-dependence of the divergent pieces of \mathcal{A}_{ct} and $\mathcal{A}_{\text{bare}}$ must be the same in order for

us to define the divergent part of δZ_ℓ in a momentum-independent way. On the other hand, since the only requirement is that the final result is UV-finite, we are always free to add or subtract finite contributions from our counterterms; nothing fixes this contribution. In other words, we are always free to take finite pieces of $\mathcal{A}_{\text{bare}}$ and absorb them into \mathcal{A}_{ct} .

However, since the finite pieces of the sub-amplitudes are not fixed, there is no guarantee that the momentum-dependences of the *finite* parts of the two pieces of the total amplitude are the same. So, when we include some finite subtractions, we must define a particular value of the momentum that we do this rearranging at, known as the *subtraction point*. The way we do the subtractions (i.e. the finite pieces we absorb into the counterterms at the subtraction point) is known as the *renormalization scheme*.

One final thing before doing an actual calculation: in order to actually isolate the divergences in the integrals, we will need to regulate the integrals first, i.e. render them finite in a way where we can obtain the original result in some limit. Now, it turns out that whenever we do this, we will need to introduce some additional parameters which don't exist in the original theory. Of course, the way we choose to regulate the integral is not unique, so the physical results cannot depend on these parameters. Consider an unphysical parameter μ that we absorb into our counterterms. Then, since $Z = Z(\mu)$, we must cancel this contribution in the corresponding renormalized parameter so that the original, bare parameter (the one that shows up in our classical theory) does not depend on this parameter. We can concisely describe this in the so-called *renormalization group equations* (RGEs), which in a condensed way say that, if we have a bare parameter, $g^{(0)} = f(\mu)Z_g(\mu)g(\mu)$, then we must require (the logarithmic derivative is not necessary, but it will be convenient later)⁹

$$\frac{dg^{(0)}}{d\log\mu} = \frac{df}{d\log\mu}Z_g(\mu)g(\mu) + f(\mu)\frac{dZ_g}{d\log\mu}g(\mu) + f(\mu)Z_g(\mu)\frac{dg}{d\log\mu} = 0, \quad (2.1.16)$$

or

$$\frac{dg}{d\log\mu} = -\frac{g(\mu)}{f(\mu)}\frac{df}{d\log\mu} - \frac{g(\mu)}{Z_g(\mu)}\frac{dZ_g}{d\log\mu}. \quad (2.1.17)$$

Eq. (2.1.17) turns out to be critical for book-keeping: it is easy to lose track of the parametric dependence on μ in the computation, but the RGEs allow us to keep track of this in a very convenient way.

Again, this all may seem a bit sketchy, but remember that neither the sum of “bare” diagrams nor the Lagrangian parameters are physical; only their combination is physical. Furthermore, we know that any physical result *must* be UV-finite, so if the sum of bare diagrams is UV-divergent, the Lagrangian parameters must also be UV-divergent in exactly the opposite way.

Another way to think about this is that the divergences of the bare diagrams is telling us that we made a stupid assumption somewhere. Of course, it is easy to see where this assumption is: we assumed that as $T \rightarrow \pm\infty$, the theory coincides exactly with the free theory. However, as we see here, the particle *interacts with itself*, and so these interactions can never truly be switched off. We have to take this into account by adjusting the Lagrangian parameters from those in the classical theory that we tried to quantize. In this case, the deep-UV interactions are the most relevant, which result in divergent corrections to the parameters, but these divergences are exactly the divergences we get from the diagrams using the unrenormalized parameters. The net result is a finite physical result.

2.2 Back to the Two-Point Function

With all of this out of the way, we can now consider adding the terms in Eq. (2.1.14) to our perturbative expansion of the path integral. Keeping only the terms $\mathcal{O}(\lambda_{\phi\theta})$, it is straightforward to find that the total two-point function (neglecting the vacuum bubbles) is given by

$$\tilde{\Pi}_\phi^{(1)}(P^2) = -\frac{\lambda_{\phi\theta}}{(P^2 - m_\phi^2 + i\epsilon)^2} \left[\frac{i}{4\pi} \delta Z_\phi^{(1)} P^2 - \frac{i}{4\pi} \left(\delta Z_\phi^{(1)} + \delta Z_{m_\phi}^{(1)} \right) m_\phi^2 + \frac{1}{2} \int \frac{d^2 K}{(2\pi)^2} \frac{1}{K^2 - m_\theta^2 + i\epsilon} \right]. \quad (2.2.1)$$

⁹The additional function $f(\mu)$ is not strictly necessary and can be absorbed into Z_g , but as we have already defined the rescaling parameters to be expansions around one, this definition is more suited for future computations.

As we know, the integral, which we will call $\mathcal{I}_0(m_\theta^2)$, is divergent, so it will first be helpful to isolate where the problem is. To do this, we will first explicitly write this as the integral over two spacetime dimensions

$$\begin{aligned}\mathcal{I}_0(m_\theta^2) &= \int_{-\infty}^{\infty} \frac{dk^0}{2\pi} \int_{-\infty}^{\infty} \frac{dk}{2\pi} \frac{1}{(k^0)^2 - k^2 - m_\theta^2 + i\epsilon} \\ &= \int_{-\infty}^{\infty} \frac{dk^0}{2\pi} \int_{-\infty}^{\infty} \frac{dk}{2\pi} \frac{1}{(k^0 + \sqrt{k^2 + m_\theta^2 - i\epsilon})(k^0 - \sqrt{k^2 + m_\theta^2 + i\epsilon})}.\end{aligned}\quad (2.2.2)$$

Now, we will employ a particularly strange trick: notice that, in the integral over k^0 , we have two poles in the complex- k^0 plane: one above the real axis at negative-valued k^0 and one below the real axis for positive-valued k^0 . This means that, if we rotate our integration contour by 90 degrees counterclockwise, we will not need to pass through any poles. This is a form of analytic continuation, known as a Wick rotation, and it has a rigorous mathematical backing (though it isn't always guaranteed to work), but to not distract from the actual point, we will not go further into this and simply take it as something that does, in fact, work. So, we rotate our integral $k^0 \rightarrow ik_E^0$ to find

$$\mathcal{I}_0(m_\theta^2) = -i \int_{-\infty}^{\infty} \frac{dk_E^0}{2\pi} \int_{-\infty}^{\infty} \frac{dk}{2\pi} \frac{1}{(k_E^0)^2 + k^2 + m_\theta^2 - i\epsilon}.\quad (2.2.3)$$

Now, we will choose instead to work in polar coordinates where $k_E^0 = K \sin \phi$ and $k = K \cos \phi$

$$\mathcal{I}_0(m_\theta^2) = -i \int_0^\infty \frac{dK}{2\pi} \int_0^{2\pi} \frac{d\phi}{2\pi} \frac{K}{K^2 + m_\theta^2},\quad (2.2.4)$$

where we took the limit as $\epsilon \rightarrow 0$ since we no longer need to worry about the poles appearing on our axis of integration. Here, it is easy to identify the problem: consider the region of the integral where K is large. In this region, the K^2 piece of the denominator dominates and we find

$$\int_0^\infty \frac{dK}{K^2 + m_\theta^2} \sim \int_0^\infty \frac{dK}{K} \sim \lim_{K \rightarrow \infty} \log K.\quad (2.2.5)$$

So, the fact that the internal momentum of the Feynman diagram is allowed to go to infinity is what results in the divergence. This source of divergence is known as an *ultraviolet divergence*. There also exist infrared divergences which have a bit of a different resolution, but we won't worry about that for now, since all of our integrals in this theory are IR finite.

Now, there are many, many ways that we could regulate this integral to render it finite so that we can extract the divergence. However, one of the nicest in terms of its theoretical structure is known as *dimensional regularization*. The basic idea is that integrals of the type (2.2.4) can actually be neatly solved in an arbitrary number of dimensions. To see, this consider

$$\mathcal{I}_d^{(n)}(m^2) = \int \frac{d^d K}{(2\pi)^d} \frac{1}{(K^2 + m^2)^n} = \int \frac{d\Omega_d}{(2\pi)^d} \int_0^\infty \frac{dK}{K} \frac{K^{d-1}}{(K^2 + m^2)^n},\quad (2.2.6)$$

where in the second step, we chose to use d -dimensional spherical coordinates and $d\Omega_d$ is the differential solid angle of a $d-1$ sphere. The integral over this solid angle just gives the surface area of this $d-1$ -sphere, which has a very nice generalization to d -dimensions in terms of the Γ -function (basically the generalization of the factorial to continuous variables)

$$\int d\Omega_d = \frac{2\pi^{d/2}}{\Gamma(\frac{d}{2})}.\quad (2.2.7)$$

Next, we will rescale $K \rightarrow m\kappa$ and define the new variable $u = (\kappa^2 + 1)^{-1}$

$$\mathcal{I}_d(m^2) = \frac{1}{(2\pi)^d} \frac{2\pi^{d/2}}{\Gamma(\frac{d}{2})} m^{d-2n} \int_0^\infty \frac{d\kappa}{(\kappa^2 + 1)^n} \kappa^{d-1} = \frac{1}{(2\pi)^d} \frac{\pi^{d/2}}{\Gamma(\frac{d}{2})} m^{d-2n} \int_0^1 du u^{n-\frac{d}{2}-1} (1-u)^{\frac{d}{2}-1}.\quad (2.2.8)$$

This integral can actually be solved again for general n and d in terms of the Euler Beta Function, which can be written again in terms of Γ -functions as

$$\mathcal{I}_d^{(n)}(m^2) = \frac{m^{d-2n}}{(4\pi)^{d/2}} \frac{\Gamma(n - \frac{d}{2})\Gamma(\frac{d}{2})}{\Gamma(\frac{d}{2})\Gamma(n)}. \quad (2.2.9)$$

Now, the source of the poles is readily apparent: the Γ -function has poles at any integer value n for $n \leq 0$. So, in our case, when $n = 1$ and $d = 2$, the numerator will diverge! However, the Γ -function also admits a very nice property that it can be expanded around any one of its poles

$$\Gamma(\epsilon - n) = \frac{(-1)^n}{n!} \left(\frac{1}{\epsilon} - \gamma_E + \psi^{(0)}(n+1) + \frac{\epsilon\pi^2}{6} + \frac{\epsilon}{2} \left[\psi^{(0)}(n+1)^2 - \psi^{(1)}(n+1) \right] + \mathcal{O}(\epsilon^2) \right), \quad (2.2.10)$$

where $\gamma_E \approx 0.5772$ is the Euler constant and $\psi^{(m)}$ is the m th Polygamma function. For positive integer n , the zeroth Polygamma function is given by

$$\psi^{(0)}(n+1) = \sum_{m=1}^n \frac{1}{m}. \quad (2.2.11)$$

At this point, we are ready for the magic of dimensional regularization: since we are able to evaluate the integrals we need analytically in any number of spacetime dimensions, we will continue our spacetime away from $d = 2$ and instead work in $d = 2 - 2\epsilon$ dimensions. Then, once we have renormalized the theory and obtain a finite result, we can take $\epsilon \rightarrow 0$ and recover the original theory. However, we need to be a bit careful with dimensions. When we take our two-dimensional integrals and upgrade them to d -dimensions, the mass-dimensions of the problem will no longer work out. In particular, in two dimensions, the integral $\mathcal{I}_0(m_\theta^2)$ is dimensionless, but if we naively continue it to d -dimensions, it will have mass dimension $d - 2 = -2\epsilon$. So, to compensate, we will introduce an artificial mass parameter μ and take

$$\int \frac{d^2 K}{(2\pi)^2} \rightarrow \left(\frac{\bar{\mu}}{4\pi} e^{\gamma_E} \right)^{2\epsilon} \int \frac{d^d K}{(2\pi)^d} = \mu^{2\epsilon} \int \frac{d^d K}{(2\pi)^d}, \quad (2.2.12)$$

where the weird factors in the middle step are chosen to cancel corresponding factors that appear in the expansion of the other parts of the integral. More properly, we should view this as the coupling constant $\lambda_{\phi\theta}$ picking up an additional mass dimension which we cancel off by redefining $\lambda_{\phi\theta} \rightarrow \mu^{2\epsilon} \lambda_{\phi\theta}$.

With all of this together, we can finally evaluate our integral

$$\mathcal{I}_0(m_\theta^2) = -\frac{i}{(4\pi)^{1-\epsilon}} \left(\frac{\mu^2}{m_\theta^2} \right)^\epsilon \Gamma(\epsilon) = -\frac{i}{4\pi} \left(\frac{1}{\epsilon} + \log \frac{\bar{\mu}^2}{m_\theta^2} \right) + \mathcal{O}(\epsilon), \quad (2.2.13)$$

where we discard all terms order ϵ and higher because we want to take the limit as $\epsilon \rightarrow 0$ once the result is finite. Using this, we find

$$\tilde{\Pi}_\phi^{(1)}(P^2) = -\frac{i\lambda_{\phi\theta}}{4\pi} \frac{1}{(P^2 - m_\phi^2 + i\epsilon)^2} \left[\delta Z_\phi^{(1)} P^2 - \left(\delta Z_\phi^{(1)} + \delta Z_{m_\phi}^{(1)} \right) m_\phi^2 - \frac{1}{2\epsilon} - \frac{1}{2} \log \frac{\bar{\mu}^2}{m_\theta^2} \right]. \quad (2.2.14)$$

As we already stated, the divergent parts of the counterterms are entirely determined by matching the momentum-dependence of the poles since these divergences must cancel for all choices of momenta. Since the pole here is momentum-independent, $\delta Z_\phi^{(1,\text{div})} = 0$ and therefore

$$\delta Z_{m_\phi}^{(1,\text{div})} = -\frac{1}{\epsilon} \frac{1}{2m_\phi^2}. \quad (2.2.15)$$

The inverse mass dependence is to be expected since Z_{m_ϕ} is dimensionless, but $\lambda_{\phi\theta}$ has mass dimension two.

Finally, we need to fix our renormalization scheme. Of course, the simplest choice is to just wash our hands, be done and not absorb any more factors into the counterterms. This is known as the *minimal subtraction (MS)* (or $\overline{\text{MS}}$ since we use $\bar{\mu}$ instead of μ) scheme. However, we can make the mass parameters

a bit more physical if we require that the poles in the propagators (which if we recall, are the true, physical masses of the particles) coincide with m_ϕ^2 . In other words, we want to require that, when $P^2 = m_\phi^2$, the denominator of Eq. (2.1.10) vanishes (up to $i\epsilon$ pieces). This clearly means that we should require $\tilde{\Pi}_\phi(m_\phi^2) = 0$. With this, we choose our counterterms

$$\delta Z_\phi^{(1)} = 0, \quad \delta Z_{m_\phi} = -\frac{1}{2m_\phi^2} \left(\frac{1}{\epsilon} + \log \frac{\bar{\mu}^2}{m_\phi^2} \right). \quad (2.2.16)$$

This choice ensures that m_ϕ^2 (and m_θ^2) are the fixed parameters measured by experiment and therefore do not run via the RGEs. This is known as the *on-shell* (OS) scheme. Conveniently, we see that for this choice of renormalization scheme, the leading-order correction to the two-point correlator exactly vanishes for all momenta, $\tilde{\Pi}_\phi^{(1)}(P^2) = 0$! Of course, the story is the same for θ with masses swapped, and so

$$\delta Z_\theta^{(1)} = 0, \quad \delta Z_{m_\theta} = -\frac{1}{2m_\theta^2} \left(\frac{1}{\epsilon} + \log \frac{\bar{\mu}^2}{m_\theta^2} \right), \quad \tilde{\Pi}_\theta^{(1)}(P^2) = 0. \quad (2.2.17)$$

2.3 Two-to-two Scattering

With all of this in hand, we are now finally ready to tackle a true interacting process. The setup will be as follows: we will start with two ϕ particles at spacetime positions $X_1^\mu = (0, x_1)$ and $X_2^\mu = (0, x_2)$, with some spread determined by position-space distribution

$$f(x - x') = \int \frac{dp}{2\pi} \tilde{f}(p - p') e^{ip(x - x')}, \quad (2.3.1)$$

where we defined the corresponding momentum-space distribution since it will usually be easier to numerically evaluate the momentum-space integrals than the position-space ones. Then, after some time t , we will probe the system for a certain number of final state particles: $\{\phi(Y_1), \phi(Y_2)\}$ and $\{\theta(Y_1), \theta(Y_2)\}$. We can immediately see that we need an even number of the same particle flavor in the final state since we have two ϕ 's in the initial state, and our action has a \mathbb{Z}_2 symmetry under $\phi \rightarrow -\phi$ and/or $\theta \rightarrow -\theta$, so we can't start with two ϕ 's and end up with a final state like $\{\phi(Y_1), \theta(Y_2)\}$. Finally, we want just a probability to measure a single particle, without knowledge of other particles in the system, so we will integrate over y_2 to only get a distribution in terms of y_1 at time t .

2.3.1 Zeroth Order

To begin, we will consider the simplest (and indeed trivial) case of $\phi\phi \rightarrow \phi\phi$ at zeroth order. This is computed with

$$\begin{aligned} & \int dx'_1 dx'_2 f(x_1 - x'_1) f(x_2 - x'_2) \langle \Omega | \phi(Y_2) \phi(Y_1) \phi(X'_2) \phi(X'_1) | \Omega \rangle^{(0)} \\ &= \mathcal{A} \int \mathcal{D}\phi \mathcal{D}\theta \int \left(\prod_{i=1}^4 \frac{d^2 P_i}{(2\pi)^2} \right) \tilde{f}(p_1 - p'_1) \tilde{f}(p_2 - p'_2) \tilde{\phi}(P_1) \tilde{\phi}(P_2) \tilde{\phi}(P_3) \tilde{\phi}(P_4) e^{-i(P_1 \cdot X_1 + P_2 \cdot X_2 + P_3 \cdot Y_2 + P_4 \cdot Y_2)} e^{i(S_0 + i\epsilon \text{ terms})} \\ &= \int \left(\prod_{i=1}^4 \frac{d^2 P_i}{(2\pi)^2} \right) \tilde{f}(p_1 - p'_1) \tilde{f}(p_2 - p'_2) \left(\pi_\phi(P_1, P_2) \pi_\phi(P_3, P_4) \right. \\ & \quad \left. + \pi_\phi(P_1, P_3) \pi_\phi(P_2, P_4) + \pi_\phi(P_1, P_4) \pi_\phi(P_2, P_3) \right) e^{-i(P_1 \cdot X_1 + P_2 \cdot X_2 + P_3 \cdot Y_1 + P_4 \cdot Y_2)}, \end{aligned} \quad (2.3.2)$$

where we defined the momentum-space propagator

$$\pi_i(P_1, P_2) \equiv \frac{i(2\pi)^2 \delta^{(2)}(P_1 + P_2)}{P_1^2 - m_i^2 + i\epsilon}. \quad (2.3.3)$$

Notice that the first product in Eq. (2.3.2) connects causally disconnected points: it pairs the particles at X_1 and X_2 as well as those at Y_1 and Y_2 . Although this will give a non-zero result, it is a pure correlation and not really relevant to the question we are asking, so we will neglect it (otherwise, we would have to consider an infinite number of n -point functions that account for all pure correlations). This will be exponentially suppressed anyway, so neglecting these contributions will not make a big impact on the result.

Of course, the remaining two pieces of Eq. (2.3.2) are just products of two two-point functions. So our leading-order four-point function for $\phi\phi \rightarrow \phi\phi$ scattering is just

$$\begin{aligned} \Lambda_{\phi\phi}^{(0)}(X_1, X_2, Y_1, Y_2) = & \int \frac{d^2 P_1}{(2\pi)^2} \frac{i\tilde{f}(p_1 - p'_1)e^{-iP_1 \cdot (Y_1 - X_1)}}{P_1^2 - m_\phi^2 + i\epsilon} \int \frac{d^2 P_2}{(2\pi)^2} \frac{i\tilde{f}(p_2 - p'_2)e^{-iP_2 \cdot (Y_2 - X_2)}}{P_2^2 - m_\phi^2 + i\epsilon} \\ & + \int \frac{d^2 P_1}{(2\pi)^2} \frac{i\tilde{f}(p_1 - p'_1)e^{-iP_1 \cdot (Y_2 - X_1)}}{P_1^2 - m_\phi^2 + i\epsilon} \int \frac{d^2 P_2}{(2\pi)^2} \frac{i\tilde{f}(p_2 - p'_2)e^{-iP_2 \cdot (Y_1 - X_2)}}{P_2^2 - m_\phi^2 + i\epsilon}. \end{aligned} \quad (2.3.4)$$

Of course, if we insert any θ 's, we will simply get zero (aside from correlations) since ϕ cannot turn into a θ simply by propagation; it needs interactions to do so!

2.3.2 First Order

The story becomes a bit more interesting at leading order in $\lambda_{\phi\theta}$. However, the first process we consider is again a trivial (albeit necessary) one; we will again consider the $\phi\phi \rightarrow \phi\phi$ case to linear order in the interactions. Recalling that we also need to insert $\delta Z_{m_i}^{(1)}$ since it is linear in $\lambda_{\phi\theta}$ (we do not need to include any factors of $\delta Z_\lambda^{(1)}$ since this part of the action begins at $\mathcal{O}(\lambda_{\phi\theta}^2)$ due to the additional factor of the coupling in the bare term), neglecting pure correlations and vacuum bubbles, and skipping a couple of steps, we find

$$\begin{aligned} \Lambda_{\phi\phi}^{(1)}(X_1, X_2, Y_1, Y_2) = & -i\lambda_{\phi\theta} \int \left(\prod_{n=1}^4 \frac{d^2 P_n}{(2\pi)^2} \right) \tilde{f}_1 \tilde{f}_2 e^{-i(P_1 \cdot X_1 + P_2 \cdot X_2 + P_3 \cdot Y_1 + P_4 \cdot Y_2)} \left[\right. \\ & 2 \times \frac{1}{4} \int \left(\prod_{m=1}^4 \frac{d^2 K_m}{(2\pi)^2} \right) \pi_\theta(K_3, K_4) \left\{ \pi_\phi(K_1, P_1) \pi_\phi(K_2, P_3) \pi_\phi(P_2, P_4) + \pi_\phi(K_1, P_1) \pi_\phi(K_2, P_4) \pi_\phi(P_2, P_3) \right. \\ & \quad \left. + \pi_\phi(K_1, P_2) \pi_\phi(K_2, P_3) \pi_\phi(P_1, P_4) + \pi_\phi(K_1, P_2) \pi_\phi(K_2, P_4) \pi_\phi(P_1, P_3) \right\} (2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4) \\ & + 2 \times \frac{\delta Z_{m_\phi} m_\phi^2}{2} \int \left(\prod_{m=1}^2 \frac{d^2 K_m}{(2\pi)^2} \right) \left\{ \pi_\phi(K_1, P_1) \pi_\phi(K_2, P_3) \pi_\phi(P_2, P_4) + \pi_\phi(K_1, P_1) \pi_\phi(K_2, P_4) \pi_\phi(P_2, P_3) \right. \\ & \quad \left. + \pi_\phi(K_1, P_2) \pi_\phi(K_2, P_3) \pi_\phi(P_1, P_4) + \pi_\phi(K_1, P_2) \pi_\phi(K_2, P_4) \pi_\phi(P_1, P_3) \right\} (2\pi)^2 \delta^{(2)}(K_1 + K_2) \left. \right]. \end{aligned} \quad (2.3.5)$$

The factors of two that we kept explicit are coming from replacing $K_1 \leftrightarrow K_2$ in the integrals and getting the same overall factor. This is to essentially keep in mind that each diagram that we draw comes with a “symmetry factor” that we have to account for in order to count repeated diagrams.

Eq. (2.3.5) looks very obscure, but when we represent it in the language of Feynman diagrams, it is clear what is happening: we are simply computing the corrected propagation between each of the points from the first-order two-point functions! However, if we choose the on-shell renormalization scheme, we saw that, since the θ loops do not feature an explicit dependence on momentum, the first order two-point function exactly vanishes. So, we have the very nice result

$$\Lambda_{\phi\phi}^{(1)}(X_1, X_2, Y_1, Y_2) = 0. \quad (2.3.6)$$

Note that, in another renormalization scheme, such as $\overline{\text{MS}}$, this would not be the case. However, in this scheme, the masses would be running parameters, and would pick up a $\log \bar{\mu}$ -dependence, which would cancel

between the zeroth order (the propagators would depend on $m_\phi(\bar{\mu})$) and this leading-order result (see the Appendix for an explicit calculation).

Next, we need to find an expression for $\phi\phi \rightarrow \theta\theta$ at this order. This expression turns out to be a bit simpler once we neglect correlations and bubbles:

$$\Lambda_{\theta\theta}(X_1, X_2, Y_1, Y_2) = -i \frac{\lambda_{\phi\theta}}{4} \int \left(\prod_{n=1}^4 \frac{d^2 P_n}{(2\pi)^2} \right) \tilde{f}_1 \tilde{f}_2 e^{-i(P_1 \cdot X_1 + P_2 \cdot X_2 + P_3 \cdot Y_1 + P_4 \cdot Y_2)} \left[\right. \\ \left. 4 \times \int \left(\prod_{m=1}^4 \frac{d^2 K_m}{(2\pi)^2} \right) \pi_\phi(K_1, P_1) \pi_\phi(K_2, P_2) \pi_\theta(K_3, P_3) \pi_\theta(K_4, P_4) (2\pi)^2 \delta^{(2)}(K_1 + K_2 + K_3 + K_4) \right], \quad (2.3.7)$$

where again, the symmetry factor arises due to the fact that there are two ways of connecting ϕ 's and two ways of connecting θ 's that give the same integral under renaming of integration variables. The K_i integrals can all be easily evaluated using the delta functions in π_i to give

$$\Lambda_{\theta\theta}^{(0)}(X_1, X_2, Y_1, Y_2) = \\ -i \lambda_{\phi\theta} \int \left(\prod_{n=1}^4 \frac{d^2 P_n}{(2\pi)^2} \right) \left[\frac{i \tilde{f}_1 e^{-i P_1 \cdot X_1}}{P_1^2 - m_\phi^2 + i\epsilon} \frac{i \tilde{f}_2 e^{-i P_2 \cdot X_2}}{P_2^2 - m_\phi^2 + i\epsilon} \frac{i e^{-i P_3 \cdot Y_1}}{P_3^2 - m_\theta^2 + i\epsilon} \frac{i e^{-i P_4 \cdot Y_2}}{P_4^2 - m_\theta^2 + i\epsilon} \right] (2\pi)^2 \delta^{(2)}(P_1 + P_2 + P_3 + P_4). \quad (2.3.8)$$

This integral is much more challenging than the previous ones we considered, and it only gets worse from here!

One can try to evaluate Eq. (2.3.8) directly using e.g. the residue theorem, but after collapsing the delta function and integrating over all the P_i^0 's, one finds non-analytic poles in the remaining integration variables which one cannot evaluate using the residue theorem. The other problem is that the integral is highly oscillatory, making numerical evaluation (even after using an e.g. Feynman reparameterization) extremely difficult. In particular, since this is a high-dimensional integral, one would want to use Monte-Carlo integration techniques, but this will not converge with such a rapidly oscillating integrand.

3 What We Really Should Calculate

At the end of the day, the sorts of calculations shown in the last section are not what we do in particle physics; this is partly because the analytical integrals that we saw are very intractable to actually compute, especially when one goes beyond leading order. They are highly oscillatory integrals with non-trivial pole structures, which makes them very difficult to compute either analytically or even numerically. However, the question that we asked is still not really aligned with what is studied in an experiment. Typically, an experimentalist will not delicately prepare two single particles in a particular momentum/position distribution and measure a small number of particles that comes out. Instead, usually beams of many particles are accelerated and focused into a particular area to interact with each other at a very small interaction point. Then, once the particles have interacted, they are seen in detectors far from the interaction point.

3.1 The S -Matrix

This story is reminiscent of what we used to derive the path integral: when we start the particles very far from each other and wait until they are far away from each other after interacting, the two sets of particles behave like the free-field particles (aside from self-interaction effects, which are accounted for via renormalization). We can therefore simply consider “in” and “out” states which are sufficiently non-interacting so that we can treat them as living at times $T \rightarrow \pm\infty$. We then define the elements of the S -matrix

$$S_{\alpha\beta} = \langle \beta, \text{out} | \alpha, \text{in} \rangle. \quad (3.1.1)$$

The indices α and β are really short-hand for a tensor product of all single-particle states that live in the initial and final systems. The S -matrix will give, in general, two contributions: the interacting piece and the

non-interacting piece. The latter will simply be given by a delta function between states α and β , while the former will give the interesting matrix elements, $M_{\beta\alpha}$. Since this is a Lorentz-invariant theory, we can then write

$$S_{\alpha\beta} = \delta(\alpha - \beta) + (2\pi)^d i \delta^{(d)}(P_\alpha - P_\beta) M_{\beta\alpha}, \quad (3.1.2)$$

where we assumed we are working in d -spacetime dimensions, and the factor of $2\pi i$ is a normalization.

We again emphasize that the states making up the S -matrix live in the infinitely distant past and infinitely far future, a good approximation when most interactions occur at very small length scales. However, we note that the free-field theory will never have bound states in it, so if our interactions do end up producing bound states which are inseparable even in the limit as $T \rightarrow \pm\infty$, the S -matrix method will fail¹⁰.

3.2 LSZ Reduction

We now wish to get an actual form for the interaction matrix elements, $M_{\beta\alpha}$ in terms of the time-ordered correlation functions that we already know how to compute. The key assumption is that, after scattering, all of our particles have changed state, even if just by a little. In the language of Feynman diagrams, this corresponds to the subset of all *connected* Feynman diagrams, simply meaning that we never have any particles propagating directly from the initial to the final state; they must always participate in some interactions. This means that no particles that live in α live in β . Next, we use the fact that $\Pi(x) = \partial_0 \Phi(x)$, as we found from our Legendre transformation to re-write the distant-past creation operator

$$a_{\mathbf{p},m;\text{in}}^\dagger = \lim_{x^0 \rightarrow -\infty} \int d^{d-1} \mathbf{x} e^{-i(\omega_{\mathbf{p}} x^0 - \mathbf{p} \cdot \mathbf{x})} \left[\sqrt{\frac{\omega_{\mathbf{p}}}{2}} \Phi_m(x) - \frac{i}{\sqrt{2\omega_{\mathbf{p}}}} \partial_0 \Phi_m(x) \right], \quad (3.2.1)$$

where we now represent spacetime vectors as un-bolded letters and spatial vectors as bolded. Introducing the notation $f \overleftrightarrow{\partial}_0 g = f \partial_0 g - g \partial_0 f$, this can be re-written as

$$a_{\mathbf{p},m;\text{in}}^\dagger = \lim_{x^0 \rightarrow -\infty} \int d^{d-1} \mathbf{x} \frac{-i e^{-i(\omega_{\mathbf{p}} x^0 - \mathbf{p} \cdot \mathbf{x})}}{\sqrt{2\omega_{\mathbf{p}}}} \overleftrightarrow{\partial}_0 \Phi_m(x). \quad (3.2.2)$$

Similarly, for the “out” annihilation operator, we have

$$a_{\mathbf{p},m;\text{out}} = \lim_{x^0 \rightarrow \infty} \int d^{d-1} \mathbf{x} \frac{i e^{i(\omega_{\mathbf{p}} x^0 - \mathbf{p} \cdot \mathbf{x})}}{\sqrt{2\omega_{\mathbf{p}}}} \overleftrightarrow{\partial}_0 \Phi_m(x). \quad (3.2.3)$$

It should be noted that, with this definition, the single-particle states have the relativistic normalization $|\mathbf{p}\rangle = \sqrt{2\omega_{\mathbf{p}}} a_{\mathbf{p}}^\dagger |0\rangle$. Then, by removing one particle from the “in” state, our matrix element in question is given by

$$M_{\beta\alpha} = \sqrt{2\omega_{\mathbf{p}}} \langle \beta, \text{out} | a_{\mathbf{p}_1, m_1; \text{in}}^\dagger | \alpha', \text{in} \rangle = \sqrt{2\omega_{\mathbf{p}}} \langle \beta, \text{out} | a_{\mathbf{p}_1, m_1; \text{in}}^\dagger | \alpha', \text{in} \rangle - \sqrt{2\omega_{\mathbf{p}}} \langle \beta, \text{out} | a_{\mathbf{p}_1, m_1; \text{out}}^\dagger | \alpha', \text{in} \rangle, \quad (3.2.4)$$

where we used the fact that the $\mathbf{p}_1; m_1$ state doesn't exist in β , and therefore $\langle \beta, \text{out} | a_{\mathbf{p}_1, m_1; \text{out}}^\dagger = 0$ and the prime denotes the fact that we have removed a single particle from the in state. Using the explicit expressions for the creation operators, we find

$$M_{\beta\alpha} = -i \left(\lim_{x_1^0 \rightarrow -\infty} - \lim_{x_1^0 \rightarrow \infty} \right) \int d^{d-1} \mathbf{x}_1 e^{-i(\omega_{\mathbf{p}_1} x_1^0 - \mathbf{p}_1 \cdot \mathbf{x}_1)} \overleftrightarrow{\partial}_{x_1^0} \langle \beta, \text{out} | \Phi_{m_1}(x_1) | \alpha', \text{in} \rangle. \quad (3.2.5)$$

By the fundamental theorem of calculus (or the one-dimensional version of Stokes' theorem),

$$\left(\lim_{t \rightarrow \infty} - \lim_{t \rightarrow -\infty} \right) f(t) = \int_{-\infty}^{\infty} dt \frac{\partial f}{\partial t}, \quad (3.2.6)$$

¹⁰Interestingly, this happens in nature with Quantum Chromodynamics (QCD). It turns out that interactions in this theory actually get *stronger* at large separation, so as $T \rightarrow \pm\infty$, particles charged under QCD will always form bound states, called hadrons. This means that one must use a different technique other than computing S -matrix elements to calculate QCD correlation functions.

giving

$$\begin{aligned}
M_{\beta\alpha} &= i \int dx_1^0 \partial_{x_1^0} \int d^{d-1} \mathbf{x}_1 e^{-i(\omega_{\mathbf{p}_1} x_1^0 - \mathbf{p}_1 \cdot \mathbf{x}_1)} \overleftrightarrow{\partial}_{x_1^0} \langle \beta, \text{out} | \Phi_{m_1}(x_1) | \alpha, \text{in} \rangle \\
&= i \int d^d x_1 \left[e^{-i(\omega_{\mathbf{p}_1} x_1^0 - \mathbf{p}_1 \cdot \mathbf{x}_1)} \partial_{x_1^0}^2 \langle \beta, \text{out} | \Phi_{m_1}(x_1) | \alpha, \text{in} \rangle - \langle \beta, \text{out} | \Phi_{m_1}(x_1) | \alpha, \text{in} \rangle \partial_{x_1^0}^2 e^{-i(\omega_{\mathbf{p}_1} x_1^0 - \mathbf{p}_1 \cdot \mathbf{x}_1)} \right].
\end{aligned} \tag{3.2.7}$$

The complex exponential is just the plane-wave solution to the Klein-Gordon equation, i.e.

$$(\partial_t^2 - \partial_{\mathbf{x}} \cdot \partial_{\mathbf{x}} + m^2) e^{-i(\omega_{\mathbf{p}} t - \mathbf{p} \cdot \mathbf{x})} = 0, \tag{3.2.8}$$

where $\omega_{\mathbf{p}} = \sqrt{|\mathbf{p}|^2 + m^2}$. Using this and integrating by parts, assuming that the integrand is suitably well-behaved¹¹ as $|\mathbf{x}| \rightarrow \infty$, we find the matrix element

$$M_{\beta\alpha} = i \int d^d x_1 e^{-i\bar{p}_1 \cdot x_1} (\partial_{x_1, \mu} \partial_{x_1}^\mu + m_{m_1}^2) \langle \beta, \text{out} | \Phi_{m_1}(x_1) | \alpha', \text{in} \rangle, \tag{3.2.9}$$

where $\bar{p}^\mu = (\omega_{\mathbf{p}}, \mathbf{p})$ is the on-shell d -momentum.

Following similar steps, we will now remove a single particle from the “out” state, in which case we find the interesting result

$$\begin{aligned}
\langle \beta, \text{out} | \Phi_{m_1}(x_1) | \alpha', \text{in} \rangle &= \sqrt{2\omega_{\mathbf{k}_1}} \langle \beta', \text{out} | a_{\mathbf{k}_1, n_1; \text{out}} \Phi_{m_1}(x_1) | \alpha', \text{in} \rangle \\
&= \sqrt{2\omega_{\mathbf{k}_1}} \langle \beta', \text{out} | a_{\mathbf{k}_1, n_1; \text{out}} \Phi_{m_1}(x_1) | \alpha', \text{in} \rangle - \sqrt{2\omega_{\mathbf{k}_1}} \langle \beta', \text{out} | \Phi_{m_1}(x_1) a_{\mathbf{k}_1, n_1; \text{in}} | \alpha', \text{in} \rangle \\
&= i \int d^d y_1 e^{i\bar{k}_1 \cdot y_1} (\partial_{y_1, \mu} \partial_{y_1}^\mu + m_{n_1}^2) \langle \beta', \text{out} | \mathbf{T}\{\Phi_{n_1}(y_1) \Phi_{m_1}(x_1)\} | \alpha', \text{in} \rangle.
\end{aligned} \tag{3.2.10}$$

Repeating this process, we find the *Lehmann-Symanzik-Zimmermann (LSZ) reduction formula*

$$\begin{aligned}
M_{\beta\alpha} &= \prod_{\beta=1}^N \left(i \int d^d y_\beta e^{i\bar{k}_\beta \cdot y_\beta} (\partial_{y_\beta, \mu} \partial_{y_\beta}^\mu + m_{n_\beta}^2) \right) \prod_{\alpha=1}^M \left(i \int d^d x_\alpha e^{-i\bar{p}_\alpha \cdot x_\alpha} (\partial_{x_\alpha, \mu} \partial_{x_\alpha}^\mu + m_{m_\alpha}^2) \right) \\
&\quad \times \langle \Omega | \mathbf{T}\{\Phi_{n_N}(y_N) \Phi_{n_{N-1}}(y_{N-1}) \times \cdots \times \Phi_{n_1}(y_1) \Phi_{m_1}(x_1) \Phi_{m_2}(x_2) \times \cdots \times \Phi_{m_M}(x_M)\} | \Omega \rangle.
\end{aligned} \tag{3.2.11}$$

This formula is even more illuminating if we pass to Fourier variables, defining

$$\begin{aligned}
&\langle \Omega | \mathbf{T}\{\Phi_{n_N}(y_N) \times \cdots \times \Phi_{n_1}(y_1) \Phi_{m_1}(x_1) \times \cdots \times \Phi_{m_M}(x_M)\} | \Omega \rangle = \\
&\prod_{\beta=1}^N \left(\int \frac{d^d k_\beta}{(2\pi)^d} e^{i k_\beta \cdot y_\beta} \right) \prod_{\alpha=1}^M \left(\int \frac{d^d p_\alpha}{(2\pi)^d} e^{i p_\alpha \cdot x_\alpha} \right) \Gamma_{m_1, \dots, m_M; n_1, \dots, n_N}(p_1, \dots, p_M; k_1, \dots, k_N),
\end{aligned} \tag{3.2.12}$$

where, after shifting the integration variables $k \rightarrow -k$, and performing all x and y integrals, we find

$$\begin{aligned}
M_{\beta\alpha} &= \prod_{\beta=1}^N \left(-i \int d^d k_\beta \delta^{(d)}(k_\beta - \bar{k}_\beta) (k_\beta^2 - m_{n_\beta}^2) \right) \prod_{\alpha=1}^M \left(-i \int d^d p_\alpha \delta^{(d)}(p_\alpha - \bar{p}_\alpha) (p_\alpha^2 - m_{m_\alpha}^2) \right) \\
&\quad \times \Gamma_{m_1, \dots, m_M; n_1, \dots, n_N}(p_1, \dots, p_M; -k_1, \dots, -k_N).
\end{aligned} \tag{3.2.13}$$

This formula is actually quite amazing: it tells us that the interesting part of our S -matrix is given exclusively by the residues of the poles in the momentum-space time-ordered correlation function, setting external momenta on-shell (also note that ingoing versus outgoing momenta are already taken care of in the formula).

¹¹Really, here we should replace the plane-wave exponential with a well-behaved wave packet that is a linear combination of plane-waves to guarantee the proper behavior as $|\mathbf{x}| \rightarrow \infty$. However, this wave-packet will still be a solution to the Klein-Gordon equation, so the argument is unchanged.

This actually makes our lives substantially easier due to the fact that the big difficulty in calculating correlation functions comes from the complicated pole structure arising from the propagators on each external leg.

One last comment is in order. Even when we let $T \rightarrow \pm\infty$, although interactions between particles will be turned off, self-interactions of a single particle will never disappear. We must therefore not only take into account the renormalization of the fields, but also the fact that these self-interactions will shift the physical mass. The latter is simple to account for: we simply take m_i in Eq. (3.2.13) and replace it with m_i^{pole} , which is the mass of the particle measured as the pole in the propagator (we already saw that this accounts for all self-interactions). The former is a bit trickier, but can be accounted for by recognizing that after renormalization, the free one-particle states are normalized so that

$$|\langle \Omega | \Phi_m(0) | \lambda_{0,m} \rangle|^2 = Z_m, \quad (3.2.14)$$

where $|\lambda_{0,m}\rangle$ is the zero-momentum, single-particle state with flavor m . This immediately implies that the properly normalized states are then given by

$$|\mathbf{p}; m\rangle = \frac{\sqrt{2\omega_{\mathbf{p};m}}}{\sqrt{Z_m}} a_{\mathbf{p},m}^\dagger |\Omega\rangle, \quad (3.2.15)$$

in which case, the renormalized LSZ-reduction formula is given by

$$\begin{aligned} M_{\beta\alpha} = & \prod_{\beta=1}^N \left(-i \int d^d k_\beta \delta^{(d)}(k_\beta - \bar{k}_\beta) \frac{(k_\beta^2 - m_{n_\beta}^{\text{pole}2})}{\sqrt{Z_{n_\beta}}} \right) \prod_{\alpha=1}^N \left(-i \int d^d p_\alpha \delta^{(d)}(p_\alpha - \bar{p}_\alpha) \frac{(p_\alpha^2 - m_{m_\alpha}^{\text{pole}2})}{\sqrt{Z_{m_\alpha}}} \right) \\ & \times \Gamma_{m_1, \dots, m_M; n_1, \dots, n_N}^{\text{ren}}(p_1, \dots, p_M; -k_1, \dots, -k_N). \end{aligned} \quad (3.2.16)$$

Qualitatively, this doesn't change much. In fact, Eq. (3.2.16) just tells us that the renormalization of the external legs is automatically accounted for when computing S -matrix elements. In other words, we don't need to consider self-energy insertions on external legs, and we can just consider “amputated” diagrams where the external legs are automatically put on-shell.

3.3 Scattering Processes

Now that we know how to actually compute S -matrix elements, we should probably figure out how to use them in calculations of physical observables. We know that the S -matrix tells us how probable it is for some initial state α to scatter to final state β , so the simplest thing that we can think of is to compute *rates* at which a particular process will occur. To do this, we will pretty much exactly follow the steps of Sec. 3.4 of Weinberg.

We will start by putting the system we care about into a hyper-toroidal “box” (i.e. a hyper-cube with all opposite ends identified with each other) with volume V , in which case the momenta of the particles are quantized. Then, the spatial delta functions become

$$\delta^{(d-1)}(\mathbf{p}' - \mathbf{p}) \rightarrow V \delta_{\mathbf{p}', \mathbf{p}} \equiv \delta_V^{(d-1)}(\mathbf{p}' - \mathbf{p}), \quad (3.3.1)$$

where $\delta_{\mathbf{p}', \mathbf{p}}$ is the ordinary Kronecker delta function. Since we have restricted ourselves to a finite volume, if we allow interactions to run for an infinite amount of time, all possible interactions will occur an infinite number of times, so we won't get anything meaningful for scattering rates. In that case, we have to also consider a “time box” where we turn on interactions at time $-T/2$ and turn them off at later time $T/2$. In this case, the energy delta functions are given by

$$\delta(p'^0 - p^0) \rightarrow \frac{1}{2\pi} \int_{-T/2}^{T/2} dt e^{i(p'^0 - p^0)t} \equiv \delta_T(p'^0 - p^0). \quad (3.3.2)$$

Recall that in the infinite-volume case, the initial- and final-state wavefunctions are normalized so that

$$\langle \Psi_\beta | \Psi_\alpha \rangle = \delta(\alpha - \beta) \prod_{n=1}^{N_\alpha} 2E_n, \quad (3.3.3)$$

which is just shorthand for saying that if the particles are unchanged between the initial and final state, they should have the same flavors, energies, and momenta. The additional normalization comes from the standard relativistic normalization of states and N_α is the number of particles in state Ψ_α . However, the spatial momenta delta functions in the box pick up an additional factor of V for each initial/final state particle. Therefore, the initial/final state wavefunctions pick up an extra factor when put in a box

$$|\Psi_\alpha\rangle = V^{N_\alpha/2} \prod_{n=1}^{N_\alpha} \sqrt{2E_n} |\Psi_\alpha^{\text{box}}\rangle \Rightarrow \langle \Psi_\beta^{\text{box}} | \Psi_\alpha^{\text{box}} \rangle = \delta_{\alpha\beta}. \quad (3.3.4)$$

With this, the “box” S -matrix is given by

$$S_{\beta\alpha}^{\text{box}} = \left(\frac{1}{V}\right)^{\frac{N_\alpha + N_\beta}{2}} \prod_{m=1}^{N_\alpha} \frac{1}{\sqrt{2E_m}} \prod_{n=1}^{N_\beta} \frac{1}{\sqrt{2E_n}} S_{\beta\alpha}. \quad (3.3.5)$$

This picks out a specific final state β with a single fixed momentum state, but often times, we want to know the transition of particles with fixed initial momentum state and *any* final momentum state, or we want to know how the transition depends on the values of the final-state momentum. We therefore need to consider a properly-normalized volume in momentum space, which we will take to be the small momentum-space box, $d^{d-1}\mathbf{p}/(2\pi)^{d-1}$ for each particle. Then, the total number of states corresponding to this momentum-space box will correspond to the combined volume of this momentum-space box and the spatial box, i.e. $V d^{d-1}\mathbf{p}/(2\pi)^{d-1}$. Finally, if we do this for all final states of the system, we get a total number of states in this phase-space volume

$$d\mathcal{N}_\beta = V^{N_\beta} \prod_{n=1}^{N_\beta} \frac{d^{d-1}\mathbf{p}_n}{(2\pi)^{d-1}}. \quad (3.3.6)$$

The probability of measuring a final state within this phase-space volume is given by the probability of transitioning from state α to a state β within this volume, times the total number of states within the volume. The latter is given by $d\mathcal{N}_\beta$, and the former is just given by $|S_{\beta\alpha}^{\text{box}}|^2$

$$dP(\alpha \rightarrow \beta) = V^{-N_\alpha} |S_{\beta\alpha}|^2 \prod_{m=1}^{N_\alpha} \frac{1}{2E_m} \prod_{n=1}^{N_\beta} \frac{d^{d-1}\mathbf{p}_n}{(2\pi)^{d-1} 2E_n}. \quad (3.3.7)$$

Again, we will only really be interested in the case where we have a true interaction, so we will only consider connected matrix elements. Therefore, $\delta(\alpha - \beta) = 0$ and we can use

$$S_{\beta\alpha} = i(2\pi)^d \delta_T(p_\beta^0 - p_\alpha^0) \delta_V^{(d-1)}(\mathbf{p}_\beta - \mathbf{p}_\alpha) M_{\beta\alpha}, \quad (3.3.8)$$

where now that we work in the finite-volume limit, we can take the square of the “delta functions” to find

$$(\delta_T(p_\beta^0 - p_\alpha^0) \delta_V^{(d-1)}(\mathbf{p}_\beta - \mathbf{p}_\alpha))^2 \stackrel{T, V \rightarrow \infty}{=} \frac{T V}{(2\pi)^d} \delta^{(d)}(p_\beta - p_\alpha), \quad (3.3.9)$$

where we used the fact that, as T and V become very large, the finite delta functions become closer and closer to the true Dirac delta functions. Then, in the large volume limit, the differential transition probability is given by

$$dP(\alpha \rightarrow \beta) = (2\pi)^d T V^{1-N_\alpha} \delta^{(d)}(p_\beta - p_\alpha) |M_{\beta\alpha}|^2 \prod_{m=1}^{N_\alpha} \frac{1}{2E_m} \prod_{n=1}^{N_\beta} \frac{d^{d-1}\mathbf{p}_n}{(2\pi)^{d-1} 2E_n}. \quad (3.3.10)$$

For any system of particles, the differential probability is just proportional to T . We can then define the rate to be the probability per unit time

$$dR(\alpha \rightarrow \beta) = \frac{V^{1-N_\alpha}}{\prod_{m=1}^{N_\alpha} 2E_m} |M_{\beta\alpha}|^2 d\Phi_n(p_\alpha; p_\beta), \quad (3.3.11)$$

where we defined the n -body Lorentz-invariant phase-space factor

$$d\Phi_n(p_\alpha; p_\beta) = (2\pi)^d \delta^{(d)}(p_\alpha - p_\beta) \prod_{n=1}^{N_\beta} \frac{d^{d-1}\mathbf{p}_n}{(2\pi)^{d-1} 2E_n}, \quad (3.3.12)$$

with the implicit understanding that all external momenta are on-shell, i.e. $E_i = \sqrt{|\mathbf{p}_i|^2 + m_i^2}$. When all is said and done, Eq. (3.3.11) is really what we are after when we want to make the prediction for the outcome of a particle physics experiment. There are really two very special cases of $N_\alpha = 1$ and $N_\alpha = 2$ which are by far the most common to discuss in particle physics contexts.

3.3.1 Particle Decays ($N_\alpha = 1$)

In the case where we only have a single particle in the initial state, we are looking at a particle which decays into a set of N_β other particles. In this case, Eq. (3.3.11) reduces to the *differential decay width*

$$d\Gamma((p_0; m_0) \rightarrow \beta) = \frac{|M_{\beta\alpha}|^2}{2E_0} d\Phi(\mathbf{p}_0; \mathbf{p}_\beta). \quad (3.3.13)$$

This will be simplest to evaluate in the initial-state particle's rest frame, where $E_0 = m_0$ and $\mathbf{p}_0 = 0$.

There is a slight issue with this case. Recall that, when we increase the size of the box, we also take the time which we allow the particles to interact for to also be very large, $T \rightarrow \infty$. However, for the particle to definitively decay, the interactions should only take place during a time period much smaller than the particle's lifetime, τ_0 . Therefore, we have to require that $T < \tau_0$, and we end up with a “delta function” with non-zero width $\Delta E \gtrsim 1/\tau_0$. This is just a statement of the uncertainty principle: if we are measuring the energies of the system, we should have at least good enough energy resolution to be able to say that we actually had a single particle in the initial state, but the uncertainty principle requires that $\Delta E \Delta t \geq 1$. Therefore, for us to even be able to say that there was a single particle in the system, we need the particle to survive long enough for us to actually resolve it!

3.3.2 Two-to- N Scattering ($N_\alpha = 2$)

The second special case occurs in most particle colliders, where we smash together two particles to produce some final states in the end. In this case, the total rate is proportional to $1/V$. It is therefore much more convenient to talk about the rate of reaction for some stream of particles going through the volume. This is described by the flux of particles $\mathcal{F}_\alpha = u_\alpha/V$, where u_α is the relative velocity of the two particles in the initial state, defined as

$$u_\alpha = \frac{\sqrt{(p_1 \cdot p_2)^2 - m_1^2 m_2^2}}{E_1 E_2}. \quad (3.3.14)$$

We then define the rate per unit flux, or *cross-section* as

$$d\sigma((p_1; m_1), (p_2; m_2) \rightarrow \beta) = \frac{|M_{\beta\alpha}|^2}{4\sqrt{(p_1 \cdot p_2)^2 - m_1^2 m_2^2}} d\Phi(\mathbf{p}_1, \mathbf{p}_2; \mathbf{p}_\beta). \quad (3.3.15)$$

The cross-section is a very general observable, independent of our given choice of experiment. However, if we do know the details of the experiment, e.g. the beam widths, the initial energies of the particles, and the interaction volume, it is straightforward to get back to the interaction rate from the cross section.

Note that with this definition, $d\sigma$ is a Lorentz invariant and can be defined irrespective of the frame of the experiment.

3.4 An Interesting Example

To showcase how a “real” (perhaps a better term is “practical”) quantum field theory calculation looks, we will consider a similar theory as before, but now upgraded to two spatial dimensions to take a bit of the triviality out of the problem. This theory then has the renormalized momentum-space action

$$S = S_0 + S_I + S_{\text{c.t.}}, \quad (3.4.1)$$

where the free and interacting actions are (we neglect the tildes, since it is assumed that these are momentum-space fields)

$$S_0 = \frac{1}{2} \int \left(\frac{d^3 k_1}{(2\pi)^3} \frac{d^3 k_2}{(2\pi)^3} \right) \left[(k_1^2 - m_\phi^2) \phi(k_1) \phi(k_2) + (k_1^2 - m_\theta^2) \theta(k_1) \theta(k_2) \right] (2\pi)^3 \delta^{(3)}(k_1 + k_2),$$

$$S_I = -\frac{\lambda_{\phi\theta}}{4} \int \left(\prod_{n=1}^4 \frac{d^3 k_n}{(2\pi)^3} \right) \phi(k_1) \phi(k_2) \theta(k_3) \theta(k_4) (2\pi)^3 \delta^{(3)}(k_1 + k_2 + k_3 + k_4),$$
(3.4.2)

respectively. The counterterm piece of the action can be found in the same way as before, but it turns out to not be relevant to what we will consider, so we ignore it for now.

Now that we have a model and the interesting experimental question, we will design a similar experiment as before: we will build a collider to smash together initial-state ϕ particles with “fixed” momentum (fixed within the uncertainty of our experiment). The particles will be provided by two beams with some width σ_b , and will cross at some fixed angle, ψ . This will form an interaction “volume” of size

$$V = \frac{\sigma_b^2}{\sin \psi}.$$
(3.4.3)

Finally, we just need to figure out the rates at which we expect certain processes to occur, which we can simply compute from scattering cross sections. We will consider two cases up to $\lambda_{\phi\theta}^2$ in perturbation theory.

3.4.1 $\phi(p_1)\phi(p_2) \rightarrow \theta(q_1)\theta(q_2)$ at Leading Order

This is by far the simplest case and is computed from all connected, amputated Feynman diagrams with a single insertion of S_I . The calculation is almost identical to what we did previously for this reaction, except now, we discard all external leg propagators to find

$$iM_{\phi^2 \rightarrow \theta^2}^{(0)} = -i\lambda_{\phi\theta}.$$
(3.4.4)

The situation simplifies considerably if we work in the center of mass frame, with $\mathbf{p}_1 = -\mathbf{p}_2 = \mathbf{p}$

$$\sqrt{(p_1 \cdot p_2)^2 - m_\phi^4} = 2|\mathbf{p}|E_0,$$
(3.4.5)

where $E_0 = \sqrt{|\mathbf{p}|^2 + m_\phi^2}$. Defining

$$\mathcal{C}_{\phi^2 \rightarrow \theta^2} = \frac{\lambda_{\phi\theta}^2}{64\pi|\mathbf{p}|E_0},$$
(3.4.6)

the differential scattering cross-section is given by

$$d^4\sigma(\phi^2 \rightarrow \theta^2) = \frac{1}{2} \mathcal{C}_{\phi^2 \rightarrow \theta^2} \frac{d^2 \mathbf{q}_1 d^2 \mathbf{q}_2}{\sqrt{|\mathbf{q}_1|^2 + m_\theta^2} \sqrt{|\mathbf{q}_2|^2 + m_\theta^2}}$$

$$\times \delta\left(2E_0 - \sqrt{|\mathbf{q}_1|^2 + m_\theta^2} - \sqrt{|\mathbf{q}_2|^2 + m_\theta^2}\right) \delta^{(2)}(\mathbf{q}_1 + \mathbf{q}_2),$$
(3.4.7)

Note the funny factor of 1/2 in the above equation: this arises due to the fact that our two θ particles in the final state are indistinguishable, and therefore there is no way to differentiate the case where one is labelled with p_1 and the other with p_2 or vice versa. Because of this redundancy, we have to get rid of the double-counted phase-space we will end up integrating over.

We can immediately evaluate the integral over \mathbf{q}_2 by collapsing the two-momentum delta function to give

$$d^2\sigma(\phi^2 \rightarrow \theta^2) = \frac{1}{2} \mathcal{C}_{\phi^2 \rightarrow \theta^2} \frac{d^2 \mathbf{q}_1}{|\mathbf{q}_1|^2 + m_\theta^2} \delta\left(2E_0 - 2\sqrt{|\mathbf{q}_1|^2 + m_\theta^2}\right).$$
(3.4.8)

We will now switch to polar coordinates where $\mathbf{q}_1 = (q \cos \xi, q \sin \xi)$ where ξ is the angle of \mathbf{q}_1 as measured from \mathbf{p} . We will then want to collapse the integral over q using the energy-conserving delta function. To do this, we can use the identity

$$\delta(f(x)) = \sum_i \frac{\delta(x - x_i)}{|f'(x_i)|}, \quad (3.4.9)$$

where x_i are the roots of the function $f(x)$. This gives

$$\delta(2E_0 - 2\sqrt{q^2 + m_\theta^2}) = \frac{E_0}{2\sqrt{E_0^2 - m_\theta^2}} \delta\left(q - \sqrt{E_0^2 - m_\theta^2}\right), \quad (3.4.10)$$

giving the differential scattering cross section

$$\frac{d\sigma_{\phi^2 \rightarrow \theta^2}}{d\xi} = \frac{\lambda_{\phi\theta}^2}{256\pi|\mathbf{p}|E_0^2}. \quad (3.4.11)$$

However, this isn't exactly what we want. Remember that, in our lab setup, we need to have the beams cross at angle ψ , which means that we will not be taking measurements in the center of mass frame. To make things simple, we will still assume that the two beams have the same energy, meaning $|\mathbf{p}'_1| = |\mathbf{p}'_2| = |\mathbf{p}'|$, where the prime denotes the true experimental frame. In this frame, we have

$$p_1^\mu = \begin{pmatrix} E'_0 \\ |\mathbf{p}'| \cos \frac{\psi}{2} \\ |\mathbf{p}'| \sin \frac{\psi}{2} \end{pmatrix}, \quad p_2^\mu = \begin{pmatrix} E'_0 \\ -|\mathbf{p}'| \cos \frac{\psi}{2} \\ |\mathbf{p}'| \sin \frac{\psi}{2} \end{pmatrix}, \quad (3.4.12)$$

where $E'_0 = \sqrt{|\mathbf{p}'|^2 + m_\phi^2}$. We want to find the boost which takes us from the center of mass frame to this frame, where in the center of mass frame, we have

$$p_1^\mu = \begin{pmatrix} E_0 \\ |\mathbf{p}| \\ 0 \end{pmatrix}, \quad p_2^\mu = \begin{pmatrix} E_0 \\ -|\mathbf{p}| \\ 0 \end{pmatrix}. \quad (3.4.13)$$

Clearly, we can achieve this by only a boost in the x^2 -direction so that $|\mathbf{p}| = |\mathbf{p}'| \cos \psi/2$. Then, we simply need to find β so that

$$\sqrt{|\mathbf{p}'|^2 + m_\phi^2} = \sqrt{\frac{|\mathbf{p}'|^2 \cos^2 \psi/2 + m_\phi^2}{1 - \beta^2}}, \quad |\mathbf{p}'| \sin \frac{\psi}{2} = -\beta \sqrt{\frac{|\mathbf{p}'|^2 \cos^2 \psi/2 + m_\phi^2}{1 - \beta^2}}. \quad (3.4.14)$$

It is straightforward to show that both of these conditions are satisfied by

$$\beta = -\frac{|\mathbf{p}'| \sin \frac{\psi}{2}}{\sqrt{|\mathbf{p}'|^2 + m_\phi^2}}. \quad (3.4.15)$$

We can then apply this Lorentz transformation to the final-state three-momenta in the center of mass frame

$$\begin{aligned} q_1^\mu = \begin{pmatrix} E_0 \\ q \cos \xi \\ q \sin \xi \end{pmatrix} &\rightarrow q_1'^\mu = \begin{pmatrix} \frac{1}{\sqrt{1-\beta^2}}(E_0 - \beta q \sin \xi) \\ q \cos \xi \\ \frac{1}{\sqrt{1-\beta^2}}(q \sin \xi - \beta E_0) \end{pmatrix} = \begin{pmatrix} E'_0 + \frac{|\mathbf{p}'| q \sin \xi \sin \frac{\psi}{2}}{E_0} \\ q \cos \xi \\ \frac{E'_0}{E_0} q \sin \xi + |\mathbf{p}'| \sin \frac{\psi}{2} \end{pmatrix} \\ q_2^\mu = \begin{pmatrix} E_0 \\ -q \cos \xi \\ -q \sin \xi \end{pmatrix} &\rightarrow q_2'^\mu = \begin{pmatrix} \frac{1}{\sqrt{1-\beta^2}}(E_0 + \beta q \sin \xi) \\ -q \cos \xi \\ \frac{1}{\sqrt{1-\beta^2}}(-q \sin \xi + \beta E_0) \end{pmatrix} = \begin{pmatrix} E'_0 - \frac{|\mathbf{p}'| q \sin \xi \sin \frac{\psi}{2}}{E_0} \\ -q \cos \xi \\ -\frac{E'_0}{E_0} q \sin \xi + |\mathbf{p}'| \sin \frac{\psi}{2} \end{pmatrix}. \end{aligned} \quad (3.4.16)$$

As a check, after some work, one can indeed verify that

$$(q_1'^0)^2 = |\mathbf{p}'|^2 + m_\phi^2 \left(|\mathbf{p}'|^2 - \frac{m_\theta^2 |\mathbf{p}'|^2}{E_0} \right) \sin^2 \frac{\psi}{2} \sin^2 \xi + \frac{2E'_0}{E_0} q |\mathbf{p}'| \sin \frac{\psi}{2} \sin \xi = |\mathbf{q}'_1|^2 + m_\theta^2 \quad (3.4.17)$$

and similar for q_2^μ . At this point, one can define a new variable ξ_1 such that

$$q_1'^\mu = \begin{pmatrix} q_1'^0 \\ |\mathbf{q}_1'| \cos \xi_1 \\ |\mathbf{q}_1'| \sin \xi_1 \end{pmatrix}, \quad (3.4.18)$$

giving the relation between ξ and ξ_1

$$\tan \xi_1 = \frac{E_0' q \sin \xi + E_0 |\mathbf{p}'| \sin \frac{\psi}{2}}{E_0 q \cos \xi} = \frac{E_0'}{E_0} \tan \xi + \frac{|\mathbf{p}'| \sin \frac{\psi}{2}}{q} \sec \xi. \quad (3.4.19)$$

Using this, along with the chain rule $d/d\xi = d\xi_1/d\xi d/d\xi_1$, one can get a differential scattering cross-section in terms of ξ' . However, the resulting expression is unnecessarily complicated, so we will not do this. Instead, we know that this gives the angle between the momentum of final-state particle 1 and the x^1 -axis. Similarly, we can find the corresponding angle ξ_2 for final-state particle 2 with

$$q_2'^\mu = \begin{pmatrix} q_2'^0 \\ |\mathbf{q}_2'| \cos \xi_2 \\ |\mathbf{q}_2'| \sin \xi_2 \end{pmatrix}, \quad (3.4.20)$$

which gives

$$\tan \xi_2 = \frac{|\mathbf{p}'| \sin \frac{\psi}{2}}{q} \sec \xi - \frac{E_0'}{E_0} \tan \xi. \quad (3.4.21)$$

Of course, these two angles are not independent, and one can check explicitly that the relations found here satisfy conservation of momentum.

Using Eq. (3.4.11), we can perform the trivial ξ integral to find the total scattering cross section in terms of the laboratory variables

$$\sigma_{\phi^2 \rightarrow \theta^2} = \frac{\lambda_{\phi\theta}^2}{128 |\mathbf{p}'| \sin \frac{\psi}{2} (|\mathbf{p}'|^2 \cos^2 \frac{\psi}{2} + m_\phi^2)}. \quad (3.4.22)$$

We can notice some interesting things about this equation. First and foremost, if we make the spatial momentum too large, the cross-section begins to decrease. This is actually a very necessary feature of a cross-section, and in fact, a scattering cross section that increases with momentum is a sign that the theory is incomplete. Remember that the cross section is related to the probability for an interaction to occur, so if it increases with momentum instead of decreases, we will run into probabilities greater than one at some energy! Historically, this was the case when considering scattering processes involving the heavy electroweak gauge bosons (W^\pm and Z^0), and was used as a justification for the existence of the Higgs boson!

Next, we note that this cross-section can help us to build our experiment (assuming that this is the process we want to measure): we know that, in order to produce the final-state particles, we need at least enough energy in the initial-state beams to produce two θ particles on-shell, i.e.

$$2E_0 \geq 2m_\theta \quad \Rightarrow \quad |\mathbf{p}'| \geq \max \left\{ \sqrt{\frac{m_\theta^2 - m_\phi^2}{\cos^2 \frac{\psi}{2}}}, 0 \right\}. \quad (3.4.23)$$

Since our cross-section is maximized for small ψ and momenta, we need to try to balance having reactions localized enough so that when we see them in a detector, we can pinpoint where they came from, having small enough momentum to not kill the reaction rate totally, and having enough momentum to keep our ϕ particles localized, i.e. behaving like particles instead of spreading out. We can help ourselves with the latter two points by considering just the case where $m_\theta \ll m_\phi$, so we are forced to have a high momentum in order for the process to occur. We still can't produce θ 's with zero momentum, though since we need them to still propagate to our detector, preferably relativistically so we don't need to account for the smearing effect of non-relativistic massive particles.

Alternatively, we may want to instead tune our experiment in order to *reduce* the amount of θ 's we pair-produce in order to see rarer processes like the ones we consider next.

3.4.2 $\phi(p_1)\phi(p_2) \rightarrow \phi(q_1)\phi(q_2)$

This process arises only at next-to-leading order in $\lambda_{\phi\theta}$ from two interaction insertions. Pairing fields in the same way as before and contracting delta-functions, we find the matrix element

$$iM_{\phi^2 \rightarrow \phi^2}^{(1)} = -\frac{\lambda_{\phi\theta}^2}{2} \int \frac{d^3k}{(2\pi)^3} \frac{i}{k^2 - m_\theta^2 + i\epsilon} \left(\frac{i}{(k+p_s)^2 - m_\theta^2 + i\epsilon} + \frac{i}{(k+p_t)^2 - m_\theta^2 + i\epsilon} + \frac{i}{(k+p_u)^2 - m_\theta^2 + i\epsilon} \right), \quad (3.4.24)$$

where $p_s = p_1 + p_2 = q_1 + q_2$, $p_t = p_1 - q_1 = q_2 - p_2$, and $p_u = p_1 - q_2 = q_1 - p_2$. Here, we are lucky: the integral in the matrix element turns out to be finite. If it were divergent, we would run into the problem that we have no term in the counterterm action that would subtract such a divergence. To solve this, we would simply add a new term to the action corresponding to a ϕ^4 interaction. While this may seem ad-hoc, it actually makes sense when we recall that, after renormalization, couplings run with the energy scale of the interaction according to the renormalization group equations. This means that, even if we set the coupling to zero at one scale, it will just be generated at a different scale, so we *must* include the interaction in the theory!

Returning to the matrix element at hand, we will need to evaluate the integral

$$\mathcal{I}_2(p^2, m^2) = \int \frac{d^3k}{(2\pi)^3} \frac{1}{(k^2 - m^2 + i\epsilon)[(k+p)^2 - m^2 + i\epsilon]}. \quad (3.4.25)$$

For this, we can use the convenient trick

$$\frac{1}{AB} = \int_0^1 dx \frac{1}{[Ax + B(1-x)]^2}, \Rightarrow \mathcal{I}_2(p^2, m^2) = \int_0^1 dx \int \frac{d^3k}{(2\pi)^3} \frac{1}{[k^2 + x(2k \cdot p + p^2) - m^2 + i\epsilon]^2}. \quad (3.4.26)$$

We can now complete the square in the denominator and shift $k \rightarrow k - xp$ to give

$$\mathcal{I}_2(p^2, m^2) = \int_0^1 dx \int \frac{d^3k}{(2\pi)^3} \frac{1}{(k^2 - \Delta + i\epsilon)^2}, \quad (3.4.27)$$

where $\Delta = x(x-1)p^2 + m^2$. With this, we can use our usual trick of Wick rotating and directly evaluate the integral over k to find

$$\mathcal{I}_2(p^2, m^2) = \frac{i}{8\pi} \int_0^1 dx \frac{1}{\sqrt{x(x-1)p^2 + m^2}} = \frac{i}{8\pi} \frac{1}{\sqrt{p^2}} \log \left(\frac{2m + \sqrt{p^2}}{2m - \sqrt{p^2}} \right), \quad (3.4.28)$$

giving the matrix element (defining the *Mandelstam variables* $p_s^2 = s$, $p_t^2 = t$, and $p_u^2 = u$)

$$iM_{\phi^2 \rightarrow \phi^2}^{(1)} = \frac{i\lambda_{\phi\theta}^2}{16\pi} [L(s) + L(t) + L(u)], \quad (3.4.29)$$

where

$$L(r) = \frac{1}{\sqrt{r}} \log \left(\frac{2m_\theta + \sqrt{r}}{2m_\theta - \sqrt{r}} \right). \quad (3.4.30)$$

Here, we see the magic of computing rates with the S -matrix. The complicated phase space integration that we did in the previous section to collapse the three-momentum-conserving delta functions was completely independent of the details of the matrix element. In this sense, computations with the S -matrix separate out the calculation of the interactions (the matrix elements) and the calculation of the kinematics (conservation of energy/momentum). We can therefore immediately write (sticking to the center of mass frame)

$$\frac{d\sigma_{\phi^2 \rightarrow \phi^2}}{d\xi} = \frac{|M_{\phi^2 \rightarrow \phi^2}|^2}{256\pi |\mathbf{p}| E_0}, \quad (3.4.31)$$

where the associated final-state three-momenta are given by

$$q_1^\mu = \begin{pmatrix} E_0 \\ |\mathbf{p}| \cos \xi \sin \xi \end{pmatrix}, \quad q_2^\mu = \begin{pmatrix} E_0 \\ -|\mathbf{p}| \cos \xi - |\mathbf{p}| \sin \xi \end{pmatrix}, \quad (3.4.32)$$

by conservation of energy and momentum, where we used the fact that the initial- and final-state masses are the same. Noting that the Mandelstam variables are Lorentz-invariants, we are free to evaluate them in any frame we wish, so in the center of mass frame, we find

$$s = 4E_0^2, \quad t = -2|\mathbf{p}|^2(1 - \cos \xi), \quad u = -2|\mathbf{p}|^2(1 + \cos \xi). \quad (3.4.33)$$

The resulting matrix element in terms of ξ is straightforward to find, but it is quite ugly. Moreover, the integral over ξ is very challenging analytically, but it is well-behaved enough that a numerical evaluation is not particularly difficult.

There is an interesting feature of the matrix element, particularly in $L(s)$. Examining Eq. (3.4.30), we see that $L(r)$ diverges at $r = 4m_\theta^2$. This is precisely when the center-of-mass energy of the particles is exactly $E_0 = m_\theta$. However, this is exactly the case where we have just the right energy to produce two *on-shell* θ particles at rest at the interaction point. If these particles are produced exactly at rest, they will not be able to propagate away from each other and they will eventually recombine in a $\theta\theta \rightarrow \phi\phi$ scattering process. This effect is known as *rescattering*. We may be worried that such a pole will cause physical divergences in scattering rates, but recall that we can never exactly tune the center of mass energy to sit at this pole due to the uncertainty principle. Furthermore, higher-order corrections will “smear” the peak, giving it some finite height.

We can note that, even though the $\phi\phi \rightarrow \phi\phi$ matrix element is suppressed by an additional factor of $\lambda_{\phi\theta}/8\pi$ as compared to the $\phi\phi \rightarrow \theta\theta$ matrix element, this rescattering effect can enhance the rate enough so that the two processes become competitive.

3.4.3 $\phi(p_1)\phi(p_2) \rightarrow \theta(q_1)\theta(q_2)$ at Next-to-Leading Order

We now need to be consistent to which order we calculate; we computed the $\phi\phi \rightarrow \phi\phi$ matrix element to $\mathcal{O}(\lambda_{\phi\theta}^2)$, so in principle, we should do the same for $\phi\phi \rightarrow \theta\theta$. The procedure is nearly identical to the previous, except for the fact that there is no s -channel diagram (and there is an additional factor of two from symmetry). Therefore, we get the next-to-leading order matrix element

$$iM_{\phi^2 \rightarrow \theta^2}^{(1)} = \lambda_{\phi\theta}^2 \int \frac{d^3k}{(2\pi)^3} \frac{1}{k^2 - m_\phi^2 + i\epsilon} \left[\frac{1}{(k + p_t)^2 - m_\theta^2 + i\epsilon} + \frac{1}{(k + p_u)^2 - m_\theta^2 + i\epsilon} \right]. \quad (3.4.34)$$

Note that now, we must perform the integral with two different masses. The procedure is the same, but with

$$\Delta = x(x-1)p^2 + m_\phi^2 + x(m_\theta^2 - m_\phi^2). \quad (3.4.35)$$

After integrating over the Feynman parameter, x , we find

$$\mathcal{I}_2(p^2, m_1^2, m_2^2) = \frac{i}{8\pi} \frac{1}{\sqrt{p^2}} \log \left(\frac{m_1 + m_2 + \sqrt{p^2}}{m_1 + m_2 - \sqrt{p^2}} \right), \quad (3.4.36)$$

resulting in the matrix element

$$M_{\phi^2 \rightarrow \theta^2}^{(1)} = \frac{i\lambda_{\phi\theta}^2}{8\pi} [L_2(t) + L_2(u)], \quad (3.4.37)$$

where, similar to before, we defined

$$L_2(r) = \frac{1}{\sqrt{r}} \log \left(\frac{m_\theta + m_\phi + \sqrt{r}}{m_\theta + m_\phi - \sqrt{r}} \right). \quad (3.4.38)$$

It's at this point, we see a potential problem. Recall that the rate depends on the modulus-squared matrix element, not the matrix element itself. If we then expand the two matrix elements that we have considered as

$$iM_{\phi^2 \rightarrow \theta^2} = \lambda_{\phi\theta} \sum_{n=1}^{\infty} \left(\frac{\lambda_{\phi\theta}}{8\pi} \right)^n A_\theta^{(n)}, \quad iM_{\phi^2 \rightarrow \phi^2} = \lambda_{\phi\theta} \sum_{n=1}^{\infty} \left(\frac{\lambda_{\phi\theta}}{8\pi} \right)^n A_\phi^{(n)}, \quad (3.4.39)$$

it is clear that, due to the vanishing of $A_\phi^{(0)}$, the leading piece of the $\phi\phi \rightarrow \phi\phi$ cross-section will occur at $\mathcal{O}(\lambda_{\phi\theta}^4/(8\pi)^2)$. The potential problem arises when we consider also calculating $A_\theta^{(2)}$, in which case the squared matrix element is

$$|M_{\phi^2 \rightarrow \theta^2}|^2 = \lambda_{\phi\theta}^2 |A_\theta^{(0)}|^2 + \frac{\lambda_{\phi\theta}^3}{8\pi} \text{Re}(A_\theta^{(0)*} A_\theta^{(1)}) + \frac{\lambda_{\phi\theta}^4}{(8\pi)^2} |A_\theta^{(1)}|^2 + \frac{2\lambda_{\phi\theta}^4}{(8\pi)^2} \text{Re}(A_\theta^{(0)*} A_\theta^{(2)}) + \mathcal{O}(\lambda_{\phi\theta}^5), \quad (3.4.40)$$

and so it is clear that, in order to be truly consistent, we need to consider the interference term between the leading order and the next-to-next-to leading order $\phi\phi \rightarrow \theta\theta$ scattering. However, $A_\theta^{(2)}$ is much more difficult to calculate. To circumvent this, we note that the higher-order contribution to the scattering rate of $\phi\phi \rightarrow \theta\theta$ will still be a small correction to this rate, so what our detector actually sees is dominated by the leading piece of the cross section. In principle, if our experiment is precise enough, it will be able to detect such differences, but this could easily be rectified by going through the pain of calculating $A_\theta^{(2)}$.

That said, we still aren't quite done. If we suppose that we are working in the case where $m_\theta \gg m_\phi$, then it does not cost us much additional energy to produce more ϕ 's in the final state. Moreover, if we really crank up the energies of experiments, we could even produce four θ 's in the final state. If we are looking at processes at $\mathcal{O}(\lambda_{\phi\theta}^2)$ at the matrix-element level in perturbation theory, then we have to consider two more processes.

3.4.4 $\phi(p_1)\phi(p_2) \rightarrow \theta(q_1)\theta(q_2)\phi(q_3)\phi(q_4)$

Luckily, this interaction's leading contribution comes with two insertions of S_I , so we will not need to worry about any loop integrals. That said, there are six total diagrams that contribute to this process, with six kinematic variables

$$\begin{aligned} s_1 &= (p_1 + p_2 - q_1)^2 = (q_2 + q_3 + q_4)^2, & s_2 &= (p_1 + p_2 - q_2)^2 = (q_1 + q_3 + q_4)^2 \\ t_1 &= (p_1 - q_1 - q_3)^2 = (p_2 - q_2 - q_4)^2, & t_2 &= (p_1 - q_1 - q_4)^2 = (p_2 - q_2 - q_3)^2 \\ u_1 &= (p_1 - q_2 - q_3)^2 = (p_2 - q_1 - q_4)^2, & u_2 &= (p_1 - q_2 - q_4)^2 = (p_2 - q_1 - q_3)^2. \end{aligned} \quad (3.4.41)$$

Summing the resulting diagrams gives the leading-order matrix element

$$iM_{\phi^2 \rightarrow \theta^2 \phi^2}^{(0)} = -i\lambda_{\phi\theta}^2 \sum_{i=1,2} \left[\frac{1}{s_i - m_\theta^2} + \frac{1}{t_i - m_\theta^2} + \frac{1}{u_i - m_\theta^2} \right]. \quad (3.4.42)$$

This doesn't seem so bad, but the real challenge comes from performing the four-body phase space integral. In fact, even just collapsing the energy-momentum-conserving delta function is in general quite challenging. To ease the calculation, we can utilize the following trick: any n -body phase space factor can be factorized using

$$d\Phi_n(p_\alpha; q_1, \dots, q_n) = \frac{1}{2\pi} d\Phi_j(r; q_1, \dots, q_j) d\Phi_{n-j+1}(p_\alpha; r, q_{j+1}, \dots, q_n) dr^2, \quad (3.4.43)$$

where

$$r^\mu = \sum_{i=1}^j p_i^\mu \quad \Rightarrow \quad r^2 = \left(\sum_{i=1}^j E_i \right)^2 - \left| \sum_{i=1}^j \mathbf{p}_i \right|^2. \quad (3.4.44)$$

In our particular case of the four-body phase space, we can apply this formula twice to reduce it to a product of two-body phase space integrals. Defining $r_{ij}^\mu = q_i^\mu + q_j^\mu$, this gives

$$d\Phi_n(p_1 + p_2; q_1, q_2, q_3, q_4) = \frac{dr_{12}^2 dr_{34}^2}{(2\pi)^2} d\Phi_2(p_1 + p_2; r_{12}, r_{34}) d\Phi_2(r_{12}; q_1, q_2) d\Phi_2(r_{34}; q_3, q_4). \quad (3.4.45)$$

Each independent two-body phase space is much easier to analyze. In fact, since these are each independently Lorentz invariant, we can choose which frame to evaluate each phase-space factor in separately. Of course,

the simplest choice is the initial-state center of mass frame, in which case, a general two-body phase space factor is given by (in three dimensions)

$$d\Phi_2(p; q_1, q_2) = \frac{1}{8\pi} \frac{d^2\mathbf{q}_1 d^2\mathbf{q}_2}{\sqrt{|\mathbf{q}_1|^2 + m_1^2} \sqrt{|\mathbf{q}_2|^2 + m_2^2}} \delta(m_p - \sqrt{|\mathbf{q}_1|^2 + m_1^2} - \sqrt{|\mathbf{q}_2|^2 + m_2^2}) \delta^{(2)}(\mathbf{q}_1 + \mathbf{q}_2), \quad (3.4.46)$$

where here, the masses denote the invariant mass of a given three-momentum, $m_i = \sqrt{p_i^2}$, and don't necessarily correspond to masses in the traditional sense. Like before, the spatial momentum delta function can be immediately evaluated yielding $\mathbf{q}_1 = -\mathbf{q}_2$. Then, we can move to polar coordinates defining ξ_{12} to be the angle of \mathbf{q}_1 with respect to some reference axis, where we evaluate the energy-conserving delta function to give

$$d\Phi_2(p; q_1, q_2) = \frac{d\xi_1}{8\pi \sqrt{p^2}}. \quad (3.4.47)$$

We also wish to note that the magnitude of each three-momentum is fixed solely by kinematics to be

$$|\mathbf{q}_1| = |\mathbf{q}_2| = \frac{1}{2m_p} \sqrt{(m_p^2 - (m_1 + m_2)^2)(m_p^2 - (m_1 - m_2)^2)} = \lambda(m_p, m_1, m_2). \quad (3.4.48)$$

It is also good to note that the ranges of the variables are

$$r_{12}^2 \in [4m_\theta^2, (\sqrt{s} - \sqrt{r_{34}^2})^2], \quad r_{34}^2 \in [4m_\phi^2, (\sqrt{s} - 2m_\theta)^2], \quad (3.4.49)$$

Using this, it is simple to define the fifth-differential cross-section in the center of mass frame (there is an extra factor of $1/4$ due to the two pairs of indistinguishable particles)

$$\frac{d^5\sigma_{\phi^2 \rightarrow \theta^2 \phi^2}}{dr_{12}^2 dr_{34}^2 d\xi_{12} d\xi_1 d\xi_3} = \frac{|M_{\phi^2 \rightarrow \theta^2 \phi^2}|^2}{8|\mathbf{p}|E_0(2\pi)^2(8\pi)^3 \sqrt{sr_{12}^2 r_{34}^2}}. \quad (3.4.50)$$

Notice the huge suppression by an additional factor of $4096\pi^5$ in the denominator. Despite the fact that this is just $\mathcal{O}(\lambda_{\phi\theta}^2)$ in the matrix element, this process will be extremely rare just simply due to the fact that there is not a lot of room in phase space for this decay to occur!

We will choose the angles as follows: ξ_{12} will be the angle of $\mathbf{r}_{12} = \mathbf{q}_1 + \mathbf{q}_2$ in the center of mass frame with respect to the beam axis (which we choose to be the x^1 -axis). Then, ξ_1 will be the angle between the momentum vector of particle 1 in the $1-2$ center of mass frame and the $1-2$ center of mass momentum vector in the overall center of mass frame, and ξ_3 will be similar for particle 3 in the $3-4$ system.

We now need to do two things: first, we need our six kinematical variables in terms of the kinematical variables that we integrate over. Then, we need to know how to relate the integration variables to the angles and energies that we would measure in our experiment. Both of these can be accomplished by first finding the momenta in terms of the integration variables in the overall center of mass frame, which can be done using standard Lorentz transformations.

We can find a boost that takes us into the $i-j$ rest frame using the boost parameter, β_{ij} , which solves

$$0 = \frac{1}{\sqrt{1 + |\beta_{ij}|^2}} (\mathbf{r}_{ij} - \beta_{ij} \sqrt{|\mathbf{r}_{ij}|^2 + r_{ij}^2}) \Rightarrow \beta_{ij} = \frac{\mathbf{r}_{ij}}{\sqrt{|\mathbf{r}_{ij}|^2 + r_{ij}^2}}, \quad (3.4.51)$$

where the two spatial vectors are given by conservation of momentum as well as Eq. (3.4.48)

$$\mathbf{r}_{12} = -\mathbf{r}_{34} = \lambda(\sqrt{s}, \sqrt{r_{12}^2}, \sqrt{r_{34}^2}) \begin{pmatrix} \cos \xi_{12} \\ \sin \xi_{12} \end{pmatrix}. \quad (3.4.52)$$

In the $1-2$ and $3-4$ rest frames, the three-momenta of the final-state particles are given by

$$\begin{aligned} q_1^{(12)\mu} &= \begin{pmatrix} \sqrt{\lambda(\sqrt{r_{12}^2}, m_\theta, m_\theta)^2 + m_\theta^2} \\ \lambda(\sqrt{r_{12}^2}, m_\theta, m_\theta) \cos \xi_1 \\ \lambda(\sqrt{r_{12}^2}, m_\theta, m_\theta) \sin \xi_1 \end{pmatrix}, & q_2^{(12)\mu} &= \begin{pmatrix} \sqrt{\lambda(\sqrt{r_{12}^2}, m_\theta, m_\theta)^2 + m_\theta^2} \\ -\lambda(\sqrt{r_{12}^2}, m_\theta, m_\theta) \cos \xi_1 \\ -\lambda(\sqrt{r_{12}^2}, m_\theta, m_\theta) \sin \xi_1 \end{pmatrix} \\ q_3^{(34)\mu} &= \begin{pmatrix} \sqrt{\lambda(\sqrt{r_{34}^2}, m_\phi, m_\phi)^2 + m_\phi^2} \\ \lambda(\sqrt{r_{34}^2}, m_\phi, m_\phi) \cos \xi_1 \\ \lambda(\sqrt{r_{34}^2}, m_\phi, m_\phi) \sin \xi_1 \end{pmatrix}, & q_4^{(34)\mu} &= \begin{pmatrix} \sqrt{\lambda(\sqrt{r_{34}^2}, m_\phi, m_\phi)^2 + m_\phi^2} \\ -\lambda(\sqrt{r_{34}^2}, m_\phi, m_\phi) \cos \xi_1 \\ -\lambda(\sqrt{r_{34}^2}, m_\phi, m_\phi) \sin \xi_1 \end{pmatrix}. \end{aligned} \quad (3.4.53)$$

Then, after performing inverse Lorentz transformations as well as a rotation, we find the momenta in the overall center of mass frame. The result is quite ugly, so we won't display it, but suffice to say that the center of mass momenta all take the form

$$q_i^\mu = \begin{pmatrix} \sqrt{|\mathbf{q}_i|^2 + m_i^2} \\ |\mathbf{q}_i| \cos \xi_i^{\text{cm}} \\ |\mathbf{q}_i| \sin \xi_i^{\text{cm}} \end{pmatrix}, \quad (3.4.54)$$

which we can then give in the experiment's frame using the boost we found before as

$$q_i'^\mu = \begin{pmatrix} \frac{E'_0}{E_0} q_i^0 + \frac{|\mathbf{p}'| \sin \psi/2}{E'_0} q_i^2 \\ q_i^1 \\ \frac{E'_0}{E_0} q_i^2 + \frac{|\mathbf{p}'| \sin \psi/2}{E'_0} q_i^0 \end{pmatrix}, \quad (3.4.55)$$

where, like before $|\mathbf{p}'|$ is the momentum of the particles in the beam, ψ is the relative angle of the beams, $E_0 = \sqrt{|\mathbf{p}'|^2 \cos^2 \psi/2 + m_\phi^2}$ is the center of mass energy of one particle in the collision, and $E'_0 = \sqrt{|\mathbf{p}'|^2 + m_\phi^2}$ is the total energy of a single particle. With this, one can find the magnitude and angle of the particle hitting the detector.

We can now evaluate our kinematic variables in the center of mass frame. Again, the results are quite ugly, so we don't give them here, but the method to obtain them is straightforward.

3.4.5 $\phi(p_1)\phi(p_2) \rightarrow \theta(q_1)\theta(q_2)\theta(q_3)\theta(q_4)$

This case is nearly identical to the $\phi\phi \rightarrow \theta\theta\phi\phi$ case, except for some minor alterations. First and foremost, it is forbidden to have any “s-like” diagrams, so the six kinematic variables that we have are t_1, t_2, u_1, u_2 , along with two new variables

$$t_3 = (p_1 - q_1 - q_2)^2 = (p_2 - q_3 - q_4)^2, \quad u_3 = (p_1 - q_3 - q_4)^2 = (p_2 - q_1 - q_2)^2, \quad (3.4.56)$$

giving the matrix element

$$iM_{\phi^2 \rightarrow \theta^4} = -i\lambda_{\phi\theta}^2 \sum_{i=1}^3 \left(\frac{1}{t_i - m_\phi^2} + \frac{1}{u_i - m_\phi^2} \right). \quad (3.4.57)$$

Other than this change and a replacement of the factor of $1/4 \rightarrow 1/4!$ due to the symmetry of the final state, the phase-space integration is identical with $m_\phi \rightarrow m_\theta$ in the final state.

3.5 Converting a Cross-Section to a Rate

Now that we have differential cross-sections for all of our processes we wish to consider, we want to convert this into a rate, which is something we actually measure at an experiment. We will assume that both beamlines in our experiment are the same, just moving in opposite directions.

To understand how to do this, we first can think of the cross-section as the “size” of the particles when viewed from the other's perspective. With this, it is clear that the rate of interaction will depend on the density of particles in both beams, n ; if we increase the density of just one beam, it should increase linearly. If we allow the particles to interact in a larger volume, the rate should also increase. Finally, if we increase the number of particles going through the interaction volume in a given amount of time, the rate should also increase. We then arrive at an expression for the *instantaneous luminosity*

$$\mathcal{L} = n^2 V u_\alpha, \quad (3.5.1)$$

where like before, u_α is the relative velocity of our initial-state particles, and V is the interaction volume. Then, the rate for a particular process to occur is given by

$$R_{\alpha \rightarrow \beta} = \mathcal{L} \sigma_{\alpha \rightarrow \beta}, \quad (3.5.2)$$

for a given cross section, $\sigma_{\alpha \rightarrow \beta}$. Notice that, since the cross section is Lorentz invariant, the rate is *not* Lorentz invariant due to the factor of the volume and relative velocity in the definition of the luminosity.

This is, however, to be expected since the rate is the number of interactions in a given time, but the clocks run at different speeds between any two reference frames.

In the frame of our experiment, we have already computed the volume, and a straightforward computation of the relative velocity gives

$$u'_\alpha = \frac{2|\mathbf{p}'|E_0 \cos \frac{\psi}{2}}{E_0'^2} \Rightarrow \mathcal{L} = \frac{2n^2 \sigma_b^2 |\mathbf{p}'|E_0 \cos \frac{\psi}{2}}{E_0'^2 \sin \psi}. \quad (3.5.3)$$

As a final comment, this isn't actually how real particle experiments work. It turns out that having a continuous beam of particles is not particularly useful since it leads to a huge mess of constant interactions. Instead, what is done at e.g. the Large Hadron Collider at CERN is that the particles are lumped into "packets" with a fixed number of particles, N , crossing at a frequency f . In this case, the luminosity can be given by

$$\mathcal{L} = \frac{N^2 f}{A}, \quad (3.5.4)$$

where A is the cross-sectional area of the beam (in our case, this is just σ_b).

4 Unstable Particles and Resonances

The cases considered in the previous sections are certainly interesting, but the \mathbb{Z}_2 symmetry that we have imposed in the theory prevents an important physical phenomenon: particle decay. This is particularly important when trying to understand how real-world experiments study short-lived particles which decay before they reach the detector.

To examine the consequences of unstable particles, consider we include an additional scalar into our theory, ψ , such that the (bare) Lagrangian is given by

$$\mathcal{L}_\psi = \frac{1}{2} \left(\partial_\mu \psi \partial^\mu \psi - m_\psi^2 \psi^2 - g_\phi \psi \phi^2 - g_\theta \psi \theta^2 \right). \quad (4.0.1)$$

This introduces three new (momentum-space) Feynman rules to our theory

$$\psi \text{ Propagator : } \frac{i}{p^2 - m_\psi^2 + i\epsilon}, \quad \psi - \phi - \phi \text{ Vertex : } -ig_\phi, \quad \psi - \theta - \theta \text{ Vertex : } -ig_\theta, \quad (4.0.2)$$

where the momentum-space conserving delta functions are implied in the vertices.

4.1 ψ Decay

Assuming that $m_\psi > 2m_\phi > 2m_\theta$, this now opens the possibility that ψ can decay into two ϕ particles or two θ particles¹². Using the notation $\varphi_i \in \{\phi, \theta\}$, the matrix elements for such a decay are easy to find

$$iM_{\psi \rightarrow \varphi_i \varphi_i} = -ig_i. \quad (4.1.1)$$

We already computed the two-body phase space integrals in the case of scattering, which we can use (replacing $2E_0 \rightarrow m_\psi$) along with Eq. (3.3.13) to find

$$\frac{d\Gamma_{\psi \rightarrow \varphi_i \varphi_i}}{d\xi} = \frac{1}{2} \frac{|M_{\psi \rightarrow \varphi_i \varphi_i}|^2}{16\pi m_\psi^2}, \quad (4.1.2)$$

where, again, the factor of 1/2 comes from the fact that we have two identical particles in the final state. Integrating over ξ and summing the two leading contributions in perturbation theory, we find the total decay rate

$$\Gamma_\psi = \frac{1}{\tau_\psi} \approx \frac{g_\phi^2 + g_\theta^2}{16m_\psi^2}, \quad (4.1.3)$$

where τ_ψ is the half-life of the particle, ψ . We can now change how long-lived these ψ particles are by turning the knobs on g_ϕ , g_θ , and m_ψ . However, it is important to note that m_ψ must always be greater than $2m_\theta$ and $2m_\phi$ for the decays to occur, otherwise we can never satisfy the energy-conserving delta functions in the definition of the decay width.

¹²it can also decay into other states, but these will be suppressed further by higher powers of g_i or $\lambda_{\theta\phi}$

4.2 ψ Two-Point Function

Now, let us consider the amputated (i.e. no external propagators) two-point function with $\psi \rightarrow \psi$. This matrix element occurs only at one loop, and obtains two contributions: one from ϕ and one from θ , giving

$$iM_{\psi \rightarrow \psi} = \sum_i \frac{(-ig_i)^2}{2} \int \frac{d^3k}{(2\pi)^3} \frac{i}{k^2 - m_i^2 + i\epsilon} \frac{i}{(k+p)^2 - m_i^2 + i\epsilon}, \quad (4.2.1)$$

with p^μ the three-momentum of the external particle(s). Again, we end up with a symmetry factor, but this time it comes from un-cancelled factors of $1/2$ in the Feynman rules when we expand the path integral and perform the Gaussian integration. The integral is the same one we performed for $\phi\phi \rightarrow \phi\phi$ scattering, giving

$$iM_{\psi \rightarrow \psi} = \sum_i \frac{ig_i^2}{16\pi} \frac{1}{\sqrt{p^2}} \log \left(\frac{2m_i + \sqrt{p^2}}{2m_i - \sqrt{p^2}} \right), \quad (4.2.2)$$

where it will be useful to note that

$$\log x = \begin{cases} \log x, & x > 0 \\ \log(-x) + i\pi, & x < 0 \end{cases}. \quad (4.2.3)$$

Note that Eq. (4.2.2) is finite, so no renormalization is strictly necessary. However, if we want to work in the on-shell scheme, we should include a finite counterterm, which we will choose to only subtract off the real part of M —as we will see, the imaginary part actually contains physical information, and introducing an imaginary part to the mass can mess with the spectrum of the theory

$$\delta Z_{m_\psi} = - \sum_i \frac{g_i^2}{16\pi m_\psi^3} \log \left(\frac{m_\psi + 2m_i}{m_\psi - 2m_i} \right), \quad (4.2.4)$$

to the given order we care about in perturbation theory.

4.3 The Optical Theorem

We now wish to prove a very interesting property of the S -matrix in general. Consider the quantum operator, S , whose matrix elements are the S -matrix elements, i.e.

$$\langle \beta | S | \alpha \rangle = \delta(\alpha - \beta) + (2\pi)^d i \delta^{(d)}(p_\alpha - p_\beta) M_{\alpha\beta}. \quad (4.3.1)$$

We can then split this operator into an “identity” piece and a transition piece, $S = 1 + iT$. In order to preserve the unitarity of the theory (we know quantum mechanics is a unitarity theory), we must require $S^\dagger S = 1$. Then,

$$(1 - iT^\dagger)(1 + iT) = 1 + i(T - T^\dagger) + T^\dagger T = 1 \quad \Rightarrow \quad T^\dagger T = -i(T - T^\dagger). \quad (4.3.2)$$

We now insert both sides of the latter equation into a matrix element between α and β , and insert a complete set of intermediate states to find

$$-i(2\pi)^d \delta^{(d)}(p_\alpha - p_\beta) (M_{\alpha \rightarrow \beta} - M_{\beta \rightarrow \alpha}^*) = \sum_\gamma M_{\beta \rightarrow \gamma}^* M_{\alpha \rightarrow \gamma} (2\pi)^d \delta^{(d)}(p_\alpha - p_\gamma) (2\pi)^d \delta^{(d)}(p_\gamma - p_\beta), \quad (4.3.3)$$

where we used

$$\langle a | O^\dagger | b \rangle = \langle b | O | a \rangle^*. \quad (4.3.4)$$

Now, the sum over states, γ , is entirely schematic. Really, we should sum over all sets of particles in the states as well as the particles’ phase spaces. Denoting a particular set of particles X_γ , we have

$$-i(M_{\alpha \rightarrow \beta} - M_{\beta \rightarrow \alpha}^*) = \sum_{X_\gamma} \int \left(\prod_{n \in X_\gamma} \frac{d^{d-1} \mathbf{p}_n}{(2\pi)^{d-1} 2E_n} \right) M_{\beta \rightarrow X_\gamma}^* M_{\alpha \rightarrow X_\gamma} (2\pi)^d \delta^{(d)} \left(p_\alpha - \sum_{n \in X_\gamma} p_n \right). \quad (4.3.5)$$

Eq. (4.3.5) is known as the *optical theorem*, and is particularly interesting in the case where $\alpha = \beta$:

$$2\text{Im}M_{\alpha \rightarrow \alpha} = \sum_{X_\gamma} \frac{1}{S_{X_\gamma}} \int \left(\prod_{n \in X_\gamma} \frac{d^{d-1}\mathbf{p}_n}{(2\pi)^{d-1}2E_n} \right) |M_{\alpha \rightarrow X_\gamma}|^2 (2\pi)^d \delta^{(d)} \left(p_\alpha - \sum_{n \in X_\gamma} p_n \right), \quad (4.3.6)$$

where S_{X_γ} is the symmetry factor associated to redundant phase-space of final-state X_γ . Note that the d -momentum conserving delta function restricts γ to contain only states which are on-shell. Eq. (4.3.6) tells us that physical processes (i.e. total scattering rates and decay rates) can be computed by calculating loop-level $n \rightarrow n$ -point functions at the matrix-element level. In other words, we can always exchange integrals over phase space for loop integrals in so-called inclusive processes.

We can see an explicit example of the optical theorem using our ψ particle. Using the fact that $E = \sqrt{p^2}$ in a particle's rest frame (we do not assume in this case that the ψ particle is on-shell), we see that

$$2\text{Im}M_{\psi \rightarrow \psi}(p^2) = 2\sqrt{p^2}\Gamma_\psi(p^2), \quad (4.3.7)$$

where $\Gamma_\psi(p^2)$ indicates that the ψ in the decay with is treated as off-shell. Using Eqs. (4.2.2) and (4.2.3), we find

$$\text{Im}M_{\psi \rightarrow \psi}(p^2) = \frac{g_\phi^2 + g_\theta^2}{16\sqrt{p^2}} + \dots, \quad (4.3.8)$$

which is exactly $\sqrt{p^2}$ times the total decay rate in Eq. (4.1.3) with $m_\psi^2 \rightarrow p^2$ (the “...” denote terms that are higher order in perturbation theory).

4.4 Resonances

At this point, one may be wondering how the interactions with this new particle may affect the scattering cross sections we calculated previously. This is straightforward enough to check: all we need to do is calculate the matrix elements of the previous processes to leading order in g_ϕ and g_θ .

4.4.1 Two-to-two scattering

The leading correction to the matrix element from the ψ to the $\phi\phi \rightarrow \theta\theta$ process only arises in the s -channel, giving

$$iM_{\phi\phi \rightarrow \theta\theta}^{g^2} = -i \frac{g_\phi g_\theta}{s - m_\psi^2}. \quad (4.4.1)$$

The contribution to $\phi\phi \rightarrow \phi\phi$ is slightly more complicated due to the fact that it contributes in all three channels, but is still relatively straightforward to calculate

$$iM_{\phi\phi \rightarrow \phi\phi}^{g^2} = -ig_\phi^2 \left(\frac{1}{s - m_\psi^2} + \frac{1}{t - m_\psi^2} + \frac{1}{u - m_\psi^2} \right). \quad (4.4.2)$$

4.4.2 Two-to-four scattering

The tree-level contributions to the $\phi\phi \rightarrow \theta\theta\theta\theta$ and $\phi\phi \rightarrow \phi\phi\theta\theta$ processes are not in principle more challenging, but the calculation is much more tedious since both processes have 24 contributing Feynman diagrams.

After evaluating all diagrams, we obtain the results

$$\begin{aligned} iM_{\phi^2 \rightarrow \theta^4}^{g^2} = & -i\lambda_{\theta\phi} \left[\frac{1}{r_{12} - m_\psi^2} \left(\frac{g_\theta^2}{s_3 - m_\theta^2} + \frac{g_\theta^2}{s_4 - m_\theta^2} + \frac{g_\theta g_\phi}{t_3 - m_\phi^2} + \frac{g_\theta g_\phi}{u_3 - m_\phi^2} \right) \right. \\ & + \frac{1}{r_{13} - m_\psi^2} \left(\frac{g_\theta^2}{s_2 - m_\theta^2} + \frac{g_\theta^2}{s_3 - m_\theta^2} + \frac{g_\theta g_\phi}{t_1 - m_\phi^2} + \frac{g_\theta g_\phi}{u_2 - m_\phi^2} \right) \\ & \left. + \frac{1}{r_{14} - m_\psi^2} \left(\frac{g_\theta^2}{s_2 - m_\theta^2} + \frac{g_\theta^2}{s_4 - m_\theta^2} + \frac{g_\theta g_\phi}{t_2 - m_\phi^2} + \frac{g_\theta g_\phi}{u_1 - m_\phi^2} \right) \right] \end{aligned}$$

$$\begin{aligned}
& + \frac{1}{r_{23} - m_\psi^2} \left(\frac{g_\theta^2}{s_1 - m_\theta^2} + \frac{g_\theta^2}{s_3 - m_\theta^2} + \frac{g_\theta g_\phi}{t_2 - m_\phi^2} + \frac{g_\theta g_\phi}{u_1 - m_\phi^2} \right) \\
& + \frac{1}{r_{24} - m_\psi^2} \left(\frac{g_\theta^2}{s_1 - m_\theta^2} + \frac{g_\theta^2}{s_4 - m_\theta^2} + \frac{g_\theta g_\phi}{t_1 - m_\phi^2} + \frac{g_\theta g_\phi}{u_2 - m_\phi^2} \right) \\
& + \frac{1}{r_{34} - m_\psi^2} \left(\frac{g_\theta^2}{s_1 - m_\theta^2} + \frac{g_\theta^2}{s_2 - m_\theta^2} + \frac{g_\theta g_\phi}{t_3 - m_\phi^2} + \frac{g_\theta g_\phi}{u_3 - m_\phi^2} \right) \Bigg], \\
iM_{\phi^2 \rightarrow 2\phi^2\theta^2}^{g^2} = & -i\lambda_{\theta\phi} \left[\frac{1}{s - m_\psi^2} \left(\frac{g_\phi^2}{s_1 - m_\phi^2} + \frac{g_\phi^2}{s_2 - m_\phi^2} + \frac{g_\theta g_\phi}{s_3 - m_\theta^2} + \frac{g_\theta g_\phi}{s_4 - m_\theta^2} \right) \right. \\
& + \frac{1}{r_{12} - m_\psi^2} \left(\frac{g_\phi^2}{t_3 - m_\phi^2} + \frac{g_\phi^2}{u_3 - m_\phi^2} + \frac{g_\theta g_\phi}{s_3 - m_\theta^2} + \frac{g_\theta g_\phi}{s_4 - m_\theta^2} \right) \\
& + \frac{1}{d_{11} - m_\psi^2} \left(\frac{g_\phi^2}{s_1 - m_\phi^2} + \frac{g_\phi^2}{t_3 - m_\phi^2} + \frac{g_\theta g_\phi}{t_1 - m_\theta^2} + \frac{g_\theta g_\phi}{t_2 - m_\theta^2} \right) \\
& + \frac{1}{d_{12} - m_\psi^2} \left(\frac{g_\phi^2}{s_2 - m_\phi^2} + \frac{g_\phi^2}{t_3 - m_\phi^2} + \frac{g_\theta g_\phi}{u_1 - m_\theta^2} + \frac{g_\theta g_\phi}{u_2 - m_\theta^2} \right) \\
& + \frac{1}{d_{21} - m_\psi^2} \left(\frac{g_\phi^2}{s_1 - m_\phi^2} + \frac{g_\phi^2}{u_3 - m_\phi^2} + \frac{g_\theta g_\phi}{u_1 - m_\theta^2} + \frac{g_\theta g_\phi}{u_2 - m_\theta^2} \right) \\
& \left. + \frac{1}{d_{22} - m_\psi^2} \left(\frac{g_\phi^2}{s_2 - m_\phi^2} + \frac{g_\phi^2}{u_3 - m_\phi^2} + \frac{g_\theta g_\phi}{t_1 - m_\theta^2} + \frac{g_\theta g_\phi}{t_2 - m_\theta^2} \right) \right],
\end{aligned}$$

where, along with the previously-defined kinematic variables, we defined

$$\begin{aligned}
s_3 = (p_1 + p_2 - q_4)^2 = (q_1 + q_2 + q_3)^2, \quad s_4 = (p_1 + p_2 - q_3)^2 = (q_1 + q_2 + q_4)^2, \\
r_{ij} = (q_i + q_j)^2, \quad d_{ij} = (p_i - q_j)^2.
\end{aligned} \tag{4.4.3}$$

4.4.3 More Divergences?

Now, the forms of the above matrix elements may be a bit distressing, as there are many poles where the expression diverges, all corresponding to the propagator of the internal ψ particle¹³. Unlike the UV poles which appeared in our discussion on renormalization, these poles are associated with kinematics – they correspond to specific choices of kinematical variables and cannot be absorbed into the definitions of parameters of the theory.

However, whenever we have a divergence which appears in our theory, it typically means we are making some kind of mistake somewhere along the way in our calculation. In this particular case, the source is a bit obscure, but we can get a first idea from the fact that the problem arises from the ψ propagator of the form

$$\Pi_\psi^{(0)}(p^2) = \frac{i}{p^2 - m_\psi^2}. \tag{4.4.4}$$

Clearly, this diverges when the ψ goes on-shell, i.e. $p^2 = m_\psi^2$. However, this is the propagator for a free field, which is of course *stable*. Here, we see the mistake we made: in our calculation, we used the propagator for a stable particle for ψ , but we are calculating the rate of scattering into only ϕ 's and θ 's. So, when we tune our initial state to produce a ψ exactly on-shell, if it is stable, it will just continue being a ψ instead of re-interacting to produce the intended particles in our final state. In other words, at this exact kinematical

¹³Although there are apparent poles corresponding to ϕ and θ propagators, these turn out to be unreachable kinematically.

point, we are actually calculating the wrong thing! Such a divergence in a physical process is known as a *resonance*.

Of course, the assumption of stability of ψ is a bad one since we know that it decays. To incorporate these results, we need to consider the corrections to the stable-particle propagator which arise from interactions. However, we already saw in the section on renormalization that the fully interacting two-point function is actually given by

$$\Pi_\psi(p^2) = \frac{i}{p^2 - m_\psi^2 - i\Pi_\psi^{\text{1PI}}(p^2)}, \quad (4.4.5)$$

where $\Pi_\psi^{\text{1PI}}(p^2)$ is the sum of one-particle-irreducible $\psi \rightarrow \psi$ diagrams. But this is going to be given exactly by $iM_{\psi \rightarrow \psi}$! Breaking this matrix element into real and imaginary parts as $M_{\psi \rightarrow \psi} = M_{\psi \rightarrow \psi}^R + iM_{\psi \rightarrow \psi}^I$ and including the explicit momentum dependence, we have

$$\Pi_\psi(p^2) = \frac{i}{p^2 - m_\psi^2 + M_{\psi \rightarrow \psi}^R(p^2) + iM_{\psi \rightarrow \psi}^I(p^2)}. \quad (4.4.6)$$

As we saw previously, though, the imaginary part of this matrix element is given exactly in terms of the off-shell decay rate, $\Gamma_\psi(p^2)$, giving

$$\Pi_\psi(p^2) = \frac{i}{p^2 - m_\psi^2 + M_{\psi \rightarrow \psi}^R(p^2) + i\sqrt{p^2}\Gamma_\psi(p^2)}. \quad (4.4.7)$$

For our last trick, we will note that the 1PI diagrams arise at higher orders in perturbation theory, so if we are working in a weakly coupled theory, either p^2 or m_ψ^2 will dominate the denominator for most choices of p^2 . This means that the majority of the contribution of the matrix element terms in the propagator will come from the region where $p^2 \approx m_\psi^2$. When $p^2 = m_\psi^2$, the real part of the two-point function vanishes since we are working in the on-shell scheme and the off-shell decay rate reduces to the *on-shell* or full decay rate, giving the corrected propagator

$$\Pi_\psi(p^2) \approx \frac{i}{p^2 - m_\psi^2 + im_\psi\Gamma_\psi}. \quad (4.4.8)$$

The parameterization of a resonance in Eq. (4.4.8) is known as a *Breit-Wigner resonance*. Notice that the additional term in the propagator has shifted the pole off of the real- p^2 axis so that the two-point function no longer diverges when the ψ particle goes on-shell.

We can think of this more intuitively in the following way: since the ψ has some finite lifetime, τ_ψ , we need at least a time resolution of $\Delta t = \tau_\psi$ in order to be able to say that the ψ was actually produced in our reaction. Of course, if we have a finite resolution in time, we must also have a corresponding resolution in our energy, $\Delta E \geq 1/\tau_\psi$. This means that we must “smear out” the energies required to produce the real ψ intermediate state by an amount $1/\tau_\psi = \Gamma_\psi$. So, while the central value of the energy needed to produce this state stays m_ψ , there is some induced width to the distribution of the energy due to this uncertainty. This distribution is exactly parameterized by Eq. (4.4.8). Furthermore, this is the reason that a total decay rate is often referred to as the *decay width* of a particle.

A $\overline{\text{MS}}$ Renormalization of m_i^2

An interesting case to consider is the one where we instead choose to renormalize our masses using $\overline{\text{MS}}$ instead of the on-shell scheme. Recall that any physical matrix element must be independent of the particular choice of renormalization scheme chosen, so we should be able to recover the same results in $\overline{\text{MS}}$ as in the on-shell scheme. However, it may not be immediately obvious how this works when we use a completely different subtraction method. In particular, we may worry that the end results are no longer independent of the unphysical parameter, μ .

To reconcile this, we recall that the renormalization group equations (RGEs) imply that the unphysical parameters, m_ϕ^2 in general pick up a μ -dependence, designated by (noting that $m_i^{(0)2} = Z_{m_i} m_i^2$)

$$\frac{dm_i^2}{d \log \mu} = -\frac{m_i^2}{Z_{m_i}} \frac{dZ_{m_i}}{d \log \mu}. \quad (\text{A.0.1})$$

We also have an analogous equation for $\lambda_{\phi\theta}$, which is related to its bare coupling as $\lambda_{\phi\theta}^{(0)} = \mu^{2\epsilon} Z_\lambda \lambda_{\phi\theta}$

$$\frac{d\lambda_{\phi\theta}}{d\log\mu} = -2\epsilon\lambda_{\phi\theta} - \frac{\lambda_{\phi\theta}}{Z_\lambda} \frac{dZ_\lambda}{d\log\mu}, \quad (\text{A.0.2})$$

where in the first term, we used $d/d\log\mu = \mu d/d\mu$. Now, since we are just using the $\overline{\text{MS}}$ scheme, our counterterms will have no explicit dependence on μ , only implicit dependence through $m_i^2(\mu)$ and $\lambda_{\phi\theta}(\mu)$. Therefore,

$$\frac{dZ_{m_i}}{d\log\mu} = \frac{\partial Z_{m_i}}{\partial \lambda_{\phi\theta}} \frac{d\lambda_{\phi\theta}}{d\log\mu} + \sum_j \frac{\partial Z_{m_i}}{\partial m_j^2} \frac{dm_j^2}{d\log\mu}, \quad (\text{A.0.3})$$

and a similar expression for $dZ_\lambda/d\log\mu$. Next, we are only considering up to $\mathcal{O}(\lambda_{\phi\theta})$ in our counterterms, and in this particular case,

$$\delta Z_{m_i}^{(1)} = -\frac{1}{2m_i^2\epsilon} \quad \delta Z_\lambda^{(1)} = 0. \quad (\text{A.0.4})$$

So, our coupled equations simplify dramatically

$$\frac{dm_i^2}{d\log\mu} = -\frac{m_i^2}{Z_{m_i}} \left(-2\epsilon \frac{\lambda_{\phi\theta}}{4\pi} \delta Z_{m_i}^{(1)} + \frac{\lambda_{\phi\theta}}{4\pi} \frac{\partial \delta Z_{m_i}^{(1)}}{\partial m_i^2} \frac{dm_i^2}{d\log\mu} \right). \quad (\text{A.0.5})$$

Combining the logarithmic derivative terms and multiplying by a factor of Z_{m_i} , we find

$$\left(1 + \frac{\lambda_{\phi\theta}}{4\pi} \delta Z_{m_i}^{(1)} + \frac{\lambda_{\phi\theta}}{4\pi} m_i^2 \frac{\partial \delta Z_{m_i}^{(1)}}{\partial m_i^2} \right) \frac{dm_i^2}{d\log\mu} = 2m_i^2 \epsilon \frac{\lambda_{\phi\theta}}{4\pi} \delta Z_{m_i}^{(1)}. \quad (\text{A.0.6})$$

From the explicit form of $\delta Z_{m_i}^{(1)}$, we find

$$m_i^2 \frac{\partial \delta Z_{m_i}^{(1)}}{\partial m_i^2} = -\delta Z_{m_i}^{(1)}, \quad (\text{A.0.7})$$

so the two divergent pieces on the left-hand side of Eq. (A.0.6) exactly cancel to give

$$\frac{dm_i^2}{d\log\mu} = -\frac{\lambda_{\phi\theta}}{4\pi} + \mathcal{O}\left(\left(\frac{\lambda_{\phi\theta}}{4\pi}\right)^2\right). \quad (\text{A.0.8})$$

This equation is solved by simple integration as long as we specify an initial condition at scale μ_0

$$m_i^2(\mu) = m_i^2(\mu_0) - \frac{\lambda_{\phi\theta}}{8\pi} \log\left(\frac{\mu^2}{\mu_0^2}\right). \quad (\text{A.0.9})$$

Here, μ_0 defines our subtraction point: one way of thinking of it is that it is the chosen value of the scale at which we extract the experimental value of the parameter.

Note that m_i^2 is the parameter that appears in the quadratic term of our Lagrangian and is therefore the mass in the propagator. In other words, the leading-order two-point function at renormalization scale μ is given by

$$\Pi_i^{(0)}(P^2, \mu) = \frac{i}{p^2 - m_i^2(\mu) + i\epsilon} = \frac{i}{P^2 - m_i^2(\mu_0) + i\epsilon} \left[1 - \frac{\lambda_{\phi\theta}}{8\pi} \frac{1}{P^2 - m_i^2(\mu_0) + i\epsilon} \log\left(\frac{\mu^2}{\mu_0^2}\right) \right]. \quad (\text{A.0.10})$$

In the $\overline{\text{MS}}$ scheme, since we have only subtracted off the $1/\epsilon$ divergence, the one-loop two-point function becomes

$$\Pi_i^{(1)}(P^2, \mu) = \frac{i\lambda_{\phi\theta}}{8\pi} \frac{1}{(P^2 - m_i^2(\mu_0) + i\epsilon)^2} \log\left(\frac{\mu^2}{m_i^2(\mu_0)}\right), \quad (\text{A.0.11})$$

where $j \neq i$. Note that we have evaluated the masses in this expression at μ_0 , since the difference between $m_i^2(\mu)$ and $m_i^2(\mu_0)$ is higher-order than we consider. It is clear that, at this order in perturbation theory, the μ -dependence of the result vanishes and one finds

$$\Pi_i^{(0)}(P^2, \mu) + \Pi_i^{(1)}(P^2, \mu) = \frac{i}{P^2 - m_i^2(\mu_0) + i\epsilon} \left[1 + \frac{\lambda_{\theta\phi}}{8\pi} \frac{1}{P^2 - m_i^2(\mu_0) + i\epsilon} \log \left(\frac{\mu_0^2}{m_j^2(\mu_0)} \right) \right]. \quad (\text{A.0.12})$$

One may not be totally convinced of the μ -independence of Eq. (A.0.12); after all, the result still depends on μ_0 . However, these dependencies are extremely different: recall the μ was *an entirely arbitrary, unfixed energy scale* while μ_0 was *the single energy scale where we fixed the initial condition*, where the two values are connected via the renormalization group equations. The distinction is that, without the RGEs, correlation functions naively depend on μ , which is a completely arbitrary, unphysical parameter, meaning we cannot hope to get sensible results from the theory. What the RGEs have accomplished by only leaving a dependence on the initial condition is that we can now simply *fix* μ_0 to be a single, arbitrary scale, often chosen to be the energy scale of some experiment. We can then extract the parameters such as $m_i^2(\mu_0)$ from the experiment operating at that chosen scale, and then use that measured value in subsequent correlation functions.