

$$S[\Psi, A] = \int d^d x \bar{\Psi} (\not{D} + m) \Psi$$

$$\not{D} = \gamma^\mu (\partial_\mu + ie A_\mu)$$

↑ canonically normalised

We are expanding $e^{-S[\Psi, A]}$

$$\rightarrow (-ie)^2 \int \text{tr} \left(\gamma^\mu \frac{1}{i\not{D} + m} \gamma^\nu \frac{1}{i\not{D} + m} \right) \frac{d^d p}{(2\pi)^d}$$

↑ fermion around the loop

There are two (equivalent) ways to see where this extra minus sign comes from.

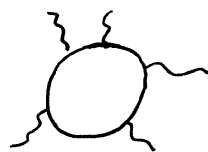
① Suppose $\langle \Psi(x) \bar{\Psi}(y) \rangle = S(x, y)$ is the Dirac propagator in position space

Then $\langle \Psi(x) \bar{\Psi}(y) \rangle = -\langle \bar{\Psi}(y) \Psi(x) \rangle$ because $\bar{\Psi}, \Psi$ anticommute.

Therefore in position space we'd get a term $(-ie)^2 \int d^d x \bar{\Psi}(x) \not{A} \Psi(x) \int d^d y \bar{\Psi}(y) \not{A} \Psi(y)$ from expanding the fermion interactions. Joining the electron fields with propagators and one of these terms is out of order \Rightarrow we get a minus sign.

More generally

$$\bar{\Psi}(x_1) \not{A} \Psi(x_1) \dots \bar{\Psi}(x_2) \not{A} \Psi(x_2) \dots \bar{\Psi}(x_n) \not{A} \Psi(x_n)$$



② Alternatively, if we perform the path integral over electrons, we get

$$\int D A D \bar{\Psi} D \Psi e^{-S[A, \bar{\Psi}, \Psi]} = \int D A \det(\not{D} + m) e^{-\frac{1}{4} \int F^2 dx}$$

if bosons, would get \det instead

$$= \int D A e^{-S_{\text{eff}}[A]} \quad \text{where } S_{\text{eff}}[A] = \frac{1}{4} \int F^2 dx - \ln \det(\not{D} + m)$$

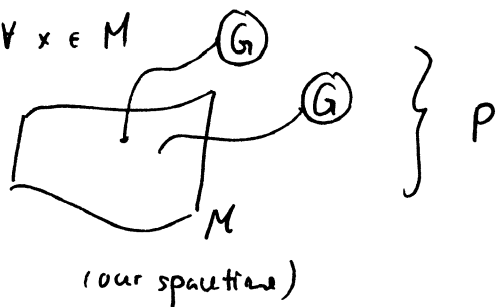
$$= \frac{1}{4} \int F^2 dx - \text{tr} \ln(\not{D} + m)$$

and perturbatively, we get $\ln(\not{D} + m) = \ln(\not{D} + m + ie \not{A})$

$$= \ln(\not{D} + m) + \sum_{n=1}^{\infty} \frac{1}{n} (ie \not{A} \frac{1}{\not{D} + m} ie \not{A} \frac{1}{\not{D} + m} \dots \frac{1}{\not{D} + m})$$

Non-Abelian Gauge Theories

Non-Abelian gauge theories are based on saying our world is a principal G bundle. This is a manifold P together with a projection $\pi: P \rightarrow M$ for some other manifold M , where $\pi^{-1}(x) \cong G$ (a lie group) for $\forall x \in M$

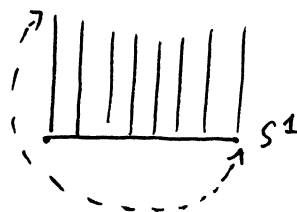
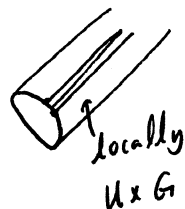


G is often called the structure group or the gauge group.

(e.g. Maxwell theory has $G = U(1)$, while SM has $G = SU(3) \times SU(2) \times U(1)$)

There's a right action of $G: P \rightarrow P$ which preserves the fibres, i.e. given $g \in G$, we have $s: P \rightarrow Pg$ where $\pi(pg) = \pi(p)$

The simplest examples are to take $M = S^1$ and $G = \mathbb{R}$. Then we have two possible \mathbb{R} -bundles: the cylinder and the moebius strip

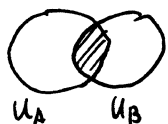


For an open set $U \subset M$, we have a local trivialisation Φ

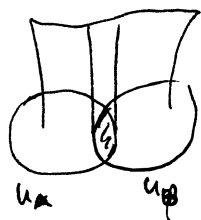
$$\Phi: P|_{\pi^{-1}(U)} \cong U \times G$$

Suppose $\{U_\alpha\}$ are a collection of open sets and we are given trivialisations Φ_α on each U_α . We have that $\Phi_\alpha: p \mapsto (x, \underbrace{\Phi_\alpha(p)}_{\text{group element}})$ and similarly

$\Phi_\beta: p' \mapsto (x', \Phi_\beta(p'))$ Now suppose U_α and U_β overlap



then on the intersection we can compare the two trivialisations.



Since $\phi_\alpha(p), \phi_\beta(p) \in G$, we must have

$$\phi_\beta(p) = \phi_\alpha(p) t_{\alpha\beta} \text{ for some } t_{\alpha\beta} \in G.$$

If we want to compare these trivialisations throughout the overlapping set, $t_{\alpha\beta}$ may need to vary with $x \in U_\alpha \cap U_\beta$. We define a transition fⁿ

$$T_{\alpha\beta}: U_\alpha \cap U_\beta \rightarrow G \quad \text{that obeys } t_{\alpha\beta}^{-1}(x) = t_{\beta\alpha}(x) \quad \forall x \in U_\alpha \cap U_\beta$$

$$x \mapsto t_{\alpha\beta}(x)$$

$$\text{and also } t_{\alpha\beta}(x) t_{\beta\gamma}(x) = t_{\alpha\gamma}(x) \quad \forall x \in U_\alpha \cap U_\beta \cap U_\gamma$$



In physics, the most application is when $M \cong \mathbb{R}^n$ and all the U_α are the whole space, so we are just comparing different trivialisations of $\pi^{-1}(\mathbb{R}^n) \cong \mathbb{R}^n \times G$

- For example, in EM, $M \cong \mathbb{R}^{3,1}$, $G = U(1)$, so our local trivialisations are a choice of gauge and transition fⁿs are gauge transformations

$$t_{\alpha\beta}(x) = e^{i\lambda_{\alpha\beta}(x)} \quad (\psi(x) \rightarrow e^{i\lambda_{\alpha\beta}(x)} \psi(x))$$

- In GR, we take $G = GL(n, \mathbb{R})$ if $\dim M = n$. Then transition fⁿs allow us to change coordinates $t_{\alpha\beta}(x) = \frac{\partial y^\mu(x)}{\partial x^\mu(x)}$

Vector bundles

To a Lie group G , we often want to study its representations. A representation is a map $\rho: G \rightarrow \text{Mat}(r, \mathbb{C})$ where $\rho(gh) = \rho(g)\rho(h)$

Picking a repn for a principle G bundle gives a vector bundle of rank r .

This is $E \xrightarrow{\pi} M$ st. we have local trivialisation $\Phi: E|_{U_\alpha} \rightarrow U_\alpha \times \mathbb{C}^r$

and transition fⁿ $T_{\alpha\beta}: U_\alpha \cap U_\beta \rightarrow \text{Mat}(r, \mathbb{C})$

If the original P has structure group $G \subset GL(r, \mathbb{C})$ then the transition fⁿs preserve some extra structure.

If $U(r) = G$, then the transition f^n preserve the inner product

$$\langle z_1, z_2 \rangle = \sum_i \bar{z}_1^i z_2^i$$

If $G = SU(r)$, then the transition f^n also have unit det.