

Relativistic Quantum Mechanics described the state of a single particle either Klein-Gordon or Dirac. The wave function has only a single pair of conjugate variables.

In Quantum Field Theory we associate such a pair of conjugate variables with every such point in a field. This allows for the field to be described an infinite number of particles.

What do we mean by a field of conjugate points:

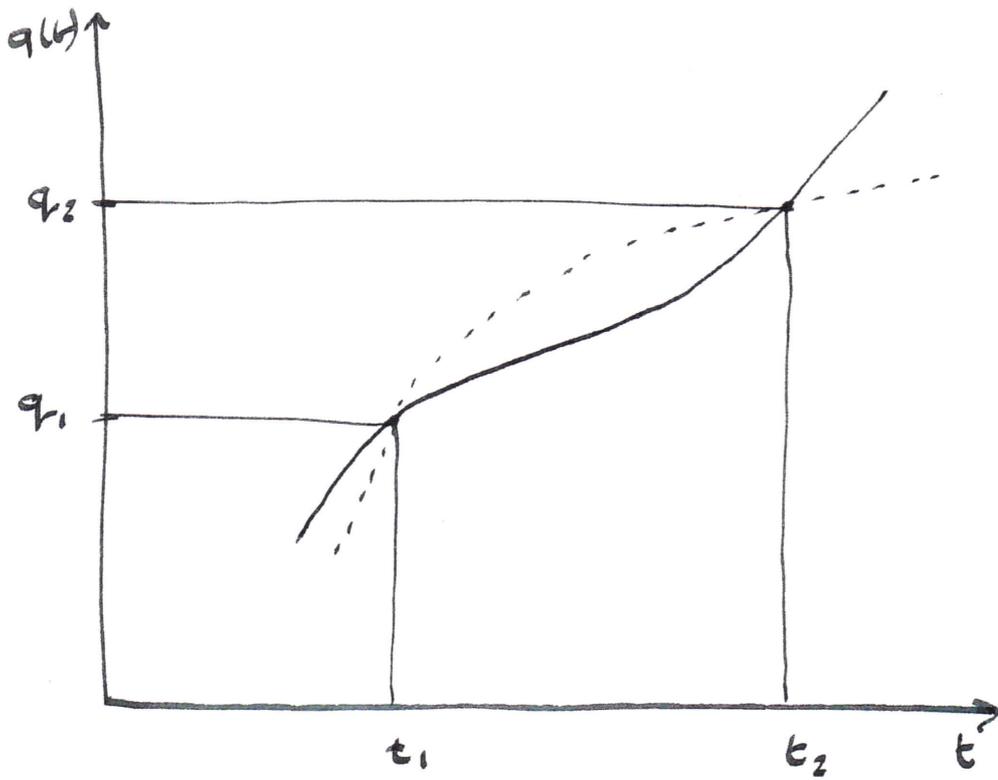
Imagine a string which can vibrate, each point along the string labelled by position x . There is a value $\phi(x)$ that can be attached to that string. It could be the transverse or longitudinal displacements or any other variable which can describe a wave.

One can then proceed to form a quantum mechanical description by using the Lagrangian and canonical quantization.

The Lagrangian and the principle of least

Action.

Supposing a system is described by $q(t)$ and t we can express the state of that system as a point on a $q(t)$ vs t graph.



The curve on this graph describes what the system does, as a function of $q(t)$.
 What describes or determines what route the system takes.

We can use the principle of least action to describe or find this curve. The curve it will take is the one for which the action $S(t_1, t_2)$ is a minimum.

For a conservative system the Action is described in terms of the Lagrangian

$$L = T - V \quad ; \quad S = \int_{t_1}^t L dt$$

We can use the calculus of variations to find the least action solutions.

We need to find the trajectory for which the integral of the Lagrangian (assumed only to depend on $(q(t), \dot{q}(t), t)$)

So

$$S = \int_{t_1}^{t_2} L(q(t), \dot{q}(t), t) dt$$

let $q(t)$ be the solution and

$$q(t) + \epsilon \phi(t)$$

a different path such that $\phi(t_1) = \phi(t_2) = 0$ and ϵ is a small quantity.

The condition that S is a minimum can

be expressed as $\frac{dS}{d\epsilon} = 0$.

(If $q(t)$ is for at t between t_1 and t_2 the minimum then $dS/d\epsilon = 0$ will be a minimum).

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$$S = \int_{t_1}^{t_2} L(q + \epsilon \phi, \dot{q} + \epsilon \dot{\phi}, t) dt$$

$$\frac{dS}{d\epsilon} = \int_{t_1}^{t_2} \left[\frac{\partial L}{\partial q} \phi + \frac{\partial L}{\partial \dot{q}} \dot{\phi} \right] dt = 0.$$

Integrating the second term by parts

$$\int_{t_1}^{t_2} \frac{\partial L}{\partial \dot{q}} \frac{d\phi}{dt} dt = \frac{\partial L}{\partial \dot{q}} \phi(t) \Big|_{t_1}^{t_2} - \int_{t_1}^{t_2} \phi(t) \frac{d}{dt} \frac{\partial L}{\partial \dot{q}} dt$$

"
0

as $\phi(t_1), \phi(t_2) = 0$.

$$\text{Thus } \frac{dS}{d\epsilon} = \int_{t_1}^{t_2} \left[\frac{\partial L}{\partial q} - \frac{d}{dt} \frac{\partial L}{\partial \dot{q}} \right] \phi(t) dt = 0.$$

This integral must be zero for arbitrary $\phi(t)$

Thus

$$\left(\frac{\partial L}{\partial q} - \frac{d}{dt} \frac{\partial L}{\partial \dot{q}} \right) = 0.$$

Euler - Lagrange Equation

For several independent coordinates

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$q_i (i = 1, 2, \dots, n)$ the generalization is simple.

$$S = \int_{t_1}^{t_2} L \{ q_1, q_2, \dots, q_n, \dot{q}_1, \dot{q}_2, \dots, \dot{q}_n, t \} dt.$$

but this simply gives an Euler Lagrange equation for each set of variables.

$$\frac{\partial L}{\partial q_i} - \frac{d}{dt} \frac{\partial L}{\partial \dot{q}_i} = 0.$$

Hamilton's Equations

Hamilton exchanged the dependence on the velocities with a dependence on the generalized momenta and is expressed as the sum of the $T+V$ is energy.

$$H = p\dot{q} - L(q, \dot{q}, t)$$

in
This term
is $2T$

$$= 2T - T - V$$

$$= T + V.$$

$$dH = p dq + \dot{q} dp - \frac{\partial L}{\partial q} dq - \frac{\partial L}{\partial \dot{q}} d\dot{q} - \frac{\partial L}{\partial t} dt$$

Since the first and 4th terms

$$p = \frac{\partial L}{\partial \dot{q}} \quad \text{cancel.}$$

$$\therefore dH = \dot{q} dp - \frac{\partial L}{\partial q} dq - \frac{\partial L}{\partial t} dt$$

But $\frac{\partial L}{\partial q} = \frac{d}{dt} \frac{\partial L}{\partial \dot{q}} = \frac{d}{dt} p = \dot{p}$

$$dH = \dot{q} dp - \dot{p} dq - \frac{\partial L}{\partial t} dt$$

$$\therefore \dot{q} = \frac{\partial H}{\partial p} \quad ; \quad \dot{p} = -\frac{\partial H}{\partial q} \quad ; \quad \frac{\partial H}{\partial t} = -\frac{\partial L}{\partial t}$$

$$\therefore \frac{dH}{dt} = \dot{q} \frac{dp}{dt} - \dot{p} \frac{dq}{dt} - \frac{\partial L}{\partial t}$$

$$\therefore \frac{dH}{dt} = \frac{\partial H}{\partial t}$$

Thus if t does not appear explicitly in the Hamiltonian it is a constant of time.

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So back to quantizing our simple string
where we are considering every point on the
string a set of independent variables

$$L(q, \dot{q}) = T - V$$

$$p_i \equiv \frac{\partial L}{\partial \dot{q}_i}$$

$$H(p, q) \equiv \sum_i p_i \dot{q}_i - L$$

We use the Heisenberg picture where
state vectors stay constant and all time
dependencies are carried by the operator.

Then we introduce equal-time commutation
relations.

$$[q_i(t), p_j(t)] = i \delta_{ij} \quad ; \quad [q_i(t), q_j(t)] = 0 \\ [p_i(t), p_j(t)] = 0.$$

In the Heisenberg picture $O(t)$

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$$-i\dot{O}(t) = [H, O]$$

The matrix element is still defined the same way,

$\langle \phi_1 | O(t) | \phi_2 \rangle$ except the states are independent of time and that is transferred to $O(t)$.

Thus we can define $\phi(x)$ as the displacement or as $\phi_i(x_i)$ for the i th molecule in the string.

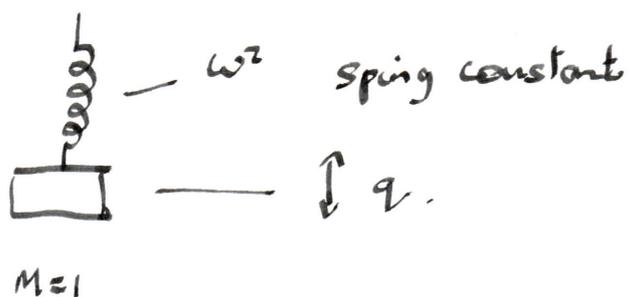
or we can choose normal modes.

$$\phi(x) = \sum_p c_p e^{ipx} + \sum_p c_p^* e^{-ipx}$$

These normal modes would just be those harmonics one would expect from the boundary conditions and taking the Fourier transform.

So it becomes natural to take the amplitudes of the normal modes as the general coordinates. The energy in each normal mode can be thought of as due to a number of particles.

Harmonic Oscillator



$$L(q, \dot{q}) = T - V = \frac{1}{2} (\dot{q}^2 - \omega^2 q^2)$$

The Euler-Lagrange equation becomes

$$\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}} \right) = \frac{\partial L}{\partial q} \quad \rightarrow \quad \dot{q} = -\omega^2 q$$

$\underbrace{\hspace{1.5cm}}_{\dot{q}} \qquad \underbrace{\hspace{1.5cm}}_{-\omega^2 q}$

$$\therefore q(t) \propto e^{\pm i\omega t}$$

$$q(t) = c e^{-i\omega t} + c^* e^{i\omega t}$$

\uparrow arbitrary complex constant.

$$p \equiv \frac{\partial L}{\partial \dot{q}} = \dot{q}$$

$$\begin{aligned} H(p, q) &\equiv p\dot{q} - \frac{1}{2} (\dot{q}^2 - \omega^2 q^2) \\ &= \dot{q}^2 - \frac{1}{2} (\dot{q}^2 - \omega^2 q^2) \Rightarrow \\ &= \frac{1}{2} (p^2 + \omega^2 q^2) \end{aligned}$$

if we impose the equal time
commutation relations

$$[q(t), p(t)] = i \quad [q(t), q(t)] = 0$$

$$[p(t), p(t)] = 0$$

$$-i\dot{q} = [H, q]$$

$$-i\dot{q} = [H, q] = \frac{1}{2} [p^2 + \omega^2 q^2, q] = \frac{1}{2} [p^2, q]$$

Now,

$$[AB, C] = A[B, C] + [A, C]B.$$

easily proved

$$= -i\dot{p} \quad \therefore \dot{q} = \dot{p}$$

$$-i\dot{p} = [H, p] = \frac{1}{2} [p^2 + \omega^2 q^2, p]$$

$$= \frac{1}{2} [\omega^2 q^2, p]$$

$$= i\omega^2 q$$

$$\therefore \dot{p} = -\omega^2 q$$

Define

$$a(t) = \frac{1}{\sqrt{2}} \left(\sqrt{\omega} q + i \frac{p}{\sqrt{\omega}} \right)$$

$$a^\dagger(t) = \frac{1}{\sqrt{2}} \left(\sqrt{\omega} q - i \frac{p}{\sqrt{\omega}} \right)$$

note $a(t)$ is not hermitian.

$$\begin{aligned}
 [a(t), a^\dagger(t)] &= \frac{1}{2} \left[\sqrt{\omega} q + i \frac{p}{\sqrt{\omega}}, \sqrt{\omega} q - i \frac{p}{\sqrt{\omega}} \right] \\
 &= \frac{1}{2} \left(\underbrace{[p, q]}_{-i} - i \underbrace{[q, p]}_i \right) = 1
 \end{aligned}$$

The Hamiltonian can be written in terms of these terms $a(t)$ $a^\dagger(t)$ as follows:-

$$\begin{aligned}
 & a(t) a^\dagger(t) + a^\dagger(t) a(t) \\
 &= \frac{1}{2} \left\{ \left(\sqrt{\omega} q + i \frac{p}{\sqrt{\omega}} \right) \left(\sqrt{\omega} q - i \frac{p}{\sqrt{\omega}} \right) + \left(\sqrt{\omega} q - i \frac{p}{\sqrt{\omega}} \right) \left(\sqrt{\omega} q + i \frac{p}{\sqrt{\omega}} \right) \right\} \\
 &= \frac{1}{2} \left(2\omega q^2 + 2 \frac{p^2}{\omega} \right) = \frac{1}{\omega} \underbrace{(\omega^2 q^2 + p^2)}_{2H}
 \end{aligned}$$

Thus

$$H = \frac{\omega}{2} \underbrace{(a(t) a^\dagger(t) + a^\dagger(t) a(t))}_{1 + a^\dagger(t) a(t)}$$

$$H = \omega \left(a^\dagger(t) a(t) + \frac{1}{2} \right)$$

The time dependence of $a(t)$

$$\begin{aligned}
 +i \dot{a}(t) &= [a(t), H] = [a(t), a^\dagger(t) a(t)] \\
 &= \omega \underbrace{[a^\dagger(t), a(t)]}_{-1} a(t)
 \end{aligned}$$

$$\therefore \dot{a}(t) = -i\omega a(t)$$

This has general solution

$$a(t) = a e^{-i\omega t} \quad \text{with } a \equiv a(0).$$

The symbol $a \equiv a(0)$

We can define the commutation relations set in terms of a, a^\dagger .

$$[a, a^\dagger] = 1$$

$$H = \omega \left(N + \frac{1}{2} \right) \quad N \equiv a^\dagger a.$$

where N is called the NUMBER OPERATOR.

The Number Operator

$$[A, Bc] = [A, B]c + B[A, c]$$

$$\begin{aligned} [a^\dagger, N] &= [a^\dagger, a^\dagger a] = \underbrace{[a^\dagger, a^\dagger]}_0 a + a^\dagger \underbrace{[a^\dagger, a]}_{-1} \\ &= -a^\dagger \end{aligned}$$

$$\therefore Na^\dagger = a^\dagger(N+1)$$

(In other words if you operate by N^\dagger first than you get a $N+1$ state.)

$$[a, N] = [a, a^\dagger a] = a^\dagger \underbrace{[a, a]}_0 + \underbrace{[a, a^\dagger]}_{+1} a = a.$$

$\therefore Na = a(N-1)$

So what does this imply?

Let $|n\rangle$ be an eigenstate of N eigenvalue n .

$\therefore N|n\rangle = n|n\rangle$

Note that since N is hermitian n is real.
 $(N = a^\dagger a \quad N^\dagger = (a^\dagger a)^\dagger = a^\dagger a = N)$

So $N a^\dagger |n\rangle = a^\dagger (N+1) |n\rangle$
 $\Rightarrow N (a^\dagger |n\rangle) = (n+1) (a^\dagger |n\rangle)$ } Thus a^\dagger can be thought of as a creation operator

$N a |n\rangle = a (N-1) |n\rangle$
 $N (a |n\rangle) = (n-1) (a |n\rangle)$ } a an annihilator operator.

$$\underline{n > 0}$$

$$\text{Let } \langle n | N | n \rangle = \langle n | a^\dagger a | n \rangle = |a|n\rangle|^2 \geq 0. \quad (17)$$

$$\text{But } \langle n | N | n \rangle = n \underbrace{\langle n | n \rangle}_{> 0}$$

$$\text{Thus } n \underbrace{\langle n | n \rangle}_{> 0} = |a|n\rangle|^2 \geq 0$$

$$\text{Thus } n > 0.$$

Thus the number of states n must be positive.

$n = \text{integer}$

Now if n were a fractional value for a given eigenvalue. It would be possible to apply $a^k |n\rangle$ until we reached a state where $n < 0$.

The only way to avoid this is for every n to be integer. At some point $a^k |n\rangle$

$$= |0\rangle$$

So what happens if you attempt to go past that point

Consider $a|0\rangle$

$$|a|0\rangle|^2 = \langle 0|a^\dagger a|0\rangle = 0 \cdot \langle 0|0\rangle = 0.$$

This implies $a|0\rangle = 0$ and so if the original state $|n\rangle$ is integer applying a will always result in $a|0\rangle = 0$ and the negative states are not accessible.

All eigenstates can be formed by:-

$$|n\rangle \equiv \frac{1}{\sqrt{n!}} a^{\dagger n} |0\rangle$$

Proof:

We have shown that

$$N(a^\dagger |n\rangle) = (n+1) \underbrace{(a^\dagger |n\rangle)}_{\propto |n+1\rangle}$$

$$\therefore a^\dagger |n\rangle = \alpha |n+1\rangle, \quad \alpha = \text{arbitrary factor}$$

But $|a^\dagger |n\rangle|^2 = \langle n| a a^\dagger |n\rangle$ $[a, a^\dagger] = 1$
 $= \langle n| 1 + N |n\rangle$ $a a^\dagger = 1 + N$
 $= 1 + n \langle n|n\rangle = 1 + n$
 $\therefore |a^\dagger |n\rangle|^2 = 1 + n$

$$\therefore \boxed{a^\dagger |n\rangle = \sqrt{n+1} |n+1\rangle}$$

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Also

$$N(a|n\rangle) = (n-1) \underbrace{(a|n\rangle)}_{|n+1\rangle}$$

$$\therefore a|n\rangle = \alpha |n+1\rangle$$

But $|a|n\rangle|^2 = \langle n|a^\dagger a|n\rangle$
 $= n \langle n|n\rangle$

$$\therefore \boxed{a|n\rangle = \sqrt{n} |n-1\rangle}$$

So if we consider starting from $|0\rangle$

$$a^\dagger |0\rangle = \sqrt{n+1} |n+1\rangle$$

$$= \sqrt{1} |1\rangle$$

$$\sqrt{1} a^\dagger |1\rangle = \sqrt{2} |2\rangle$$

$$\sqrt{1}\sqrt{2} a^\dagger |2\rangle = \sqrt{3} |3\rangle$$

$$\therefore \sqrt{n!} a^\dagger |n\rangle = \sqrt{n+1} |n+1\rangle$$

a

\therefore if we want a normalized state

$$|n\rangle = \frac{1}{\sqrt{n!}} a^{\dagger n} |0\rangle$$

Fermionic Oscillator

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We know that for fermions the state can only be occupied by one or zero particles. This is achieved by replacing the commutator relationships with anticommutator relations.

$$\{a, a^\dagger\} = 1 \quad \text{instead of } [a, a^\dagger] = 1.$$

$$\{a, a\} = \{a^\dagger, a^\dagger\} = 0 \quad \text{instead of } [a, a] = [a^\dagger, a^\dagger] = 0.$$

$$\text{But } \{a, a\} = aa + aa = 2aa = 0.$$

$$\{a^\dagger, a^\dagger\} = a^\dagger a^\dagger + a^\dagger a^\dagger = 2a^\dagger a^\dagger = 0$$

$$\therefore a^2 = a^{\dagger 2} = 0.$$

Definition of $N = a^\dagger a$ as before

$$N^2 = a^\dagger a \underbrace{a^\dagger a}_{1 - aa^\dagger} = a^\dagger a (1 - aa^\dagger)$$

$$= a^\dagger a - \underbrace{a^\dagger a a a^\dagger}_0$$

$$= N$$

$$\therefore N(N-1) = 0 \quad \Rightarrow \quad N = 0, 1. \quad \text{only.}$$

If we define

$$\langle 0|0\rangle = 1.$$

Applying the creation operator a^\dagger

$$N a^\dagger = \underbrace{a^\dagger a}_{1 - a a^\dagger} a^\dagger = a^\dagger - \underbrace{a a^\dagger}_0 = a^\dagger.$$

Thus $N(a^\dagger|0\rangle) = (a^\dagger|0\rangle)$

$$\therefore N(a^\dagger|0\rangle) = 1 \cdot (a^\dagger|0\rangle).$$

$$\therefore |1\rangle = a^\dagger|0\rangle$$

$$\langle 1|1\rangle = \langle 0| \underbrace{a a^\dagger}_{1 - a a^\dagger} |0\rangle = \langle 0| \underbrace{1 - N}_{1} |0\rangle = \langle 0|0\rangle = 1$$

The annihilation operator

$$a|1\rangle = a a^\dagger|0\rangle = (1 - N)|0\rangle = |0\rangle.$$

Summary

$$a^\dagger|1\rangle = a^{\dagger 2}|0\rangle = 0.$$

$$a|0\rangle = a^2|0\rangle = 0.$$

The hamiltonian is proportional to the number of particles up to a constant which can be ignored.

$$H = \omega a^\dagger a$$