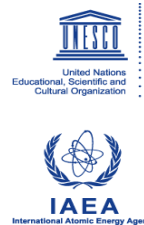




The Abdus Salam
**International Centre
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2442-5

Preparatory School to the Winter College on Optics

28 January - 1 February, 2013

Review of Quantum Mechanics and Semi-classical theory of Atom-field interaction

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1 REVIEW OF QUANTUM MECHANICS

1.1 CLASSICAL MECHANICS

Classical mechanics is based on the assumption that any physically interesting variables connected with a system/particle, such as its position, velocity or its energy can be measured with arbitrary precision and without mutual interference from any other such measurement. Laws of classical mechanics can be expressed in various mathematical forms,

1. Newtonian mechanics
2. Hamiltonian mechanics

Hamiltonian of a physical system gives its total energy

$$H = T + V$$

Quantum mechanics is based on the realization that the measuring process may affect the physical system. It is therefore impossible to measure simultaneously certain pairs of variables with precision. In quantum mechanics physical system is described by a state vector or wave function, and variables are represented by operators.

Quantum mechanics can be expressed by,

1. Wave mechanics
2. Dirac's notation

1.2 WAVE MECHANICS:

A quantum mechanical system (such as atoms, molecules, ions etc.) are given by its wave function $\psi(r, t)$. Itself $\psi(r, t)$ has no physical meaning but it allows to calculate the expectation values of all observables of interest.

Measurable quantities are called observables and are represented by hermitian operators \hat{O} . Expectation value is given by

$$\langle \hat{O} \rangle = \int d^3r \psi^*(r, t) \hat{O} \psi(r, t).$$

1.3 Probability:

The probability of finding the system in the volume element d^3r is

$$\psi^*(r, t) \hat{O} \psi(r, t).$$

As the system exist, its probability of being somewhere has to equal 1.

$$\begin{aligned} \int \psi^*(r, t) \psi(r, t) d^3r &= 1 \\ \int \psi_n^*(r, t) \psi_m(r, t) d^3r &= \delta_{nm}. \end{aligned}$$

The time development of wave function is determined by schrodinger equation,

$$i\hbar \frac{\partial}{\partial t} \psi(r, t) = H\psi(r, t),$$

where H is the Hamiltonian of the system. The energy of the unperturbed system for instance an atom not interacting with light is the sum of its potential and kinetic energies

$$H = \frac{p^2}{2m} + V(r).$$

1.4 Stationary state:

Stationary states of schrodinger equation are those for which space and time dependence are separated,

$$\psi_n(r, t) = U_n(\vec{r})e^{-i\omega_n t}.$$

Time independent equation,

$$HU_n(\vec{r}) = E_n U_n(\vec{r}) = \hbar\omega_n U_n(\vec{r}),$$

where $U_n(\vec{r})$ is an eigen function of H with eigen values $E_n = \hbar\omega_n$. The eigen values of hermitian operators are real numbers. The eigen functions of hermitian operators belonging to different eigen values are orthogonal, and eigen functions having same eigen values are normal.

$$\int U_n^*(r)U_m(r)d\vec{r} = \delta_{nm}$$

and complete

$$\sum_n U_n^*(r)U_n(r) = 1.$$

The completeness relation means that any function can be written as a linear combination of the $U_n(r, t)$. The wave function

$$\psi(r, t) = \sum_n \psi_n(r, t) = \sum_n C_n(t)U_n(\vec{r})e^{-i\omega_n t},$$

here $C_n(t)$ are the expansion coefficients.

$C_n(t)$ - constant for problems related to free part of Hamiltonian

$C_n(t)$ - change in time if we include the interaction part of Hamiltonian.

Putting the values of $\psi(r, t)$ in the normalization condition we get

$$\sum_n |C_n|^2 = 1,$$

gives the probability of finding the system in state n . The expectation value in terms of C_n

$$\langle \hat{O} \rangle = \sum_{n,m} C_n C_m^* \hat{O}_{nm} e^{-i\omega_{nm} t},$$

Where

$$\hat{O}_{nm} = \int d^3r U_m^*(r) \hat{O} U_n(r),$$

and

$$\omega_{nm} = \omega_n - \omega_m.$$

1.5 DIRAC NOTATION:

The wave function of wave mechanics corresponds to the state vector in Dirac's formulation of quantum mechanics. The relation between state vector and wave function is analogous to using vectors instead of coordinates. A vector \vec{V} can be expanded as,

$$\vec{V} = V_x \hat{x} + V_y \hat{y}.$$

In Dirac's notation

$$|V\rangle = V_x |x\rangle + V_y |y\rangle.$$

x-component of a vector is obtained by

$$\vec{V} \cdot \hat{x} = V_x,$$

in Dirac's notation

$$\langle x | V \rangle = V_x \quad \text{and} \quad \langle y | V \rangle = V_y.$$

Using these Eqn's. we can write

$$\begin{aligned} |V\rangle &= |x\rangle \langle x | V \rangle + |y\rangle \langle y | V \rangle \\ &= (|x\rangle \langle x| + |y\rangle \langle y|) |V\rangle \end{aligned}$$

The identity diadic (outer product of two vectors)

$$|x\rangle \langle x| + |y\rangle \langle y| = 1$$

for n dimensions

$$\begin{aligned} |V\rangle &= \sum_n |n\rangle \langle n | V \rangle \\ \sum_n |n\rangle \langle n| &= I \end{aligned}$$

where $\{ |n\rangle \}$ are complete set of vectors, i.e. a basis. The inner products $\langle n | V \rangle$ are the expansion coefficients of the vector $|V\rangle$ in this basis. Expansion coefficients are in general complex.

$$\langle k | V \rangle = \langle V | k \rangle^*$$

For continuous basis $\{ |r\rangle \}$

$$I = \int d^3r |r\rangle \langle r|$$

The wave vector

$$|\psi(t)\rangle = \int d^3r |r\rangle \langle r|\psi\rangle$$

Where the wave function

$$\begin{aligned}\psi(r) &= \langle r|\psi\rangle \\ \psi(x) &= \langle x|\psi\rangle\end{aligned}$$

The expectation value of the operator \hat{O} is given by,

$$\langle \hat{O} \rangle = \langle \psi(t)|\hat{O}|\psi(t)\rangle$$

Hermitian

$$\begin{aligned}\langle \psi(t)|\hat{O}|\psi(t)\rangle &= \left[\langle \psi(t)|\hat{O}^\dagger|\psi(t)\rangle \right]^* = \langle \psi(t)|\hat{O}|\psi(t)\rangle^* \\ \hat{O}^\dagger &= \hat{O}\end{aligned}$$

The set of eigen vectors of a hermitian operator is complete. This means that any arbitrary vector $|\psi(t)\rangle$ can be expressed as a sum of orthogonal eigen vectors.

$$|\psi\rangle = \sum_{n=0}^{\infty} C_n |x_n\rangle$$

Eigen vectors are orthonormal

$$\begin{aligned}\langle X_n|X_m\rangle &= \delta_{nm} \\ \delta_{nm} &= 1 \text{ for } n = m \\ &= 0 \text{ for } n \neq m\end{aligned}$$

Completeness relation for discrete case is

$$\sum_n |X_n\rangle \langle X_n| = I$$

The state vector $|\psi\rangle$ in terms of position eigen states (which are continuous).

$$\begin{aligned}|\psi(t)\rangle &= \int d\vec{x} |x\rangle \langle x|\psi\rangle \\ \int d\vec{x} |x\rangle \langle x| &= I\end{aligned}$$

The normalization of eigen vectors with a continuous set of eigen values must be normalized with the help of dirac delta function having properties,

$$\begin{aligned}\delta(x - x') &= 0 \quad \text{if } x \neq x' \\ \delta(x - x') &= \infty \quad x = x' \\ \langle x|x'\rangle &= \delta(x - x')\end{aligned}$$

State vectors obey the Schrodinger's equation

$$\begin{aligned} i\hbar \left| \dot{\psi} \right\rangle &= H |\psi\rangle, \\ |\psi\rangle &= \sum_n C_n e^{-i\omega_n t} |n\rangle \end{aligned}$$

Expectation value can be written as

$$|\psi\rangle = \sum_{n,m} C_n^* C_m e^{-i(\omega_n - \omega_m)t} |n\rangle \hat{O}_{nm}$$

Where

$$\hat{O}_{nm} = \langle m | \hat{O} | n \rangle$$

1.6 Two level system:

Wave function for two level system is

$$\psi(r, t) = C_a U_a(\vec{r}) e^{-i\omega_a t} + C_b U_b(\vec{r}) e^{-i\omega_b t}$$

State vector

$$|\psi\rangle = C_a e^{-i\omega_a t} |a\rangle + C_b e^{-i\omega_b t} |b\rangle$$

1.7 SCHRODINGER, INTERACTION AND HEISENBERG PICTURES:

1.8 SCHRODINGER PICTURE:

The interaction of radiation with matter involves a hamiltonian.

$$H = H_o + V$$

H_o - unperturbed energy

V - Interaction energy

The corresponding Schrodinger equation

$$\begin{aligned} \left| \dot{\psi}(t) \right\rangle &= \frac{-i}{\hbar} H |\psi(t)\rangle \\ \left| \dot{\psi}(t) \right\rangle &= \frac{-i}{\hbar} (H_o + V) |\psi(t)\rangle, \\ |\psi(t)\rangle &= e^{\frac{-iHt}{\hbar}} |\psi(0)\rangle \end{aligned}$$

Expectation value of an operator \hat{O} which represents the observables.

$$\langle \hat{O} \rangle = \langle \psi(t) | \hat{O}(0) | \psi(t) \rangle$$

Operator \hat{O} is independent of time, but $|\psi(t)\rangle$ is a function of time. This is the schrodinger picture way of writing the expectation value of an operator.

1.9 HEISENBERG PICTURE:

In Heisenberg picture total time dependence goes into operator, so state vector is independent of time, the expectation value of an operator in Schrodinger picture is,

$$\langle \hat{O} \rangle = \langle \psi(t) | \hat{O}(0) | \psi(t) \rangle$$

It can also be written as,

$$\langle \hat{O}(t) \rangle = \langle \psi(t) | e^{\frac{-iHt}{\hbar}} e^{\frac{+iHt}{\hbar}} \hat{O}(0) e^{\frac{-iHt}{\hbar}} e^{\frac{+iHt}{\hbar}} | \psi(t) \rangle$$

Where H is the total Hamiltonian. According to Schrodinger equation,

$$i\hbar \frac{\partial}{\partial t} |\psi(t)\rangle = H |\psi(t)\rangle$$

Integrating

$$\begin{aligned} |\psi(t)\rangle &= e^{\frac{-iHt}{\hbar}} |\psi(0)\rangle \\ e^{\frac{+iHt}{\hbar}} |\psi(t)\rangle &= |\psi(0)\rangle \end{aligned}$$

Using this we can write

$$\langle \hat{O}(t) \rangle = \langle \psi(0) | e^{\frac{+iHt}{\hbar}} \hat{O}(0) e^{\frac{-iHt}{\hbar}} | \psi(0) \rangle$$

Define

$$\hat{O}(t) = e^{\frac{+iHt}{\hbar}} \hat{O}(0) e^{\frac{-iHt}{\hbar}}$$

Then

$$\langle \hat{O}(t) \rangle = \langle \psi(0) | \hat{O}(t) | \psi(0) \rangle$$

Which is called Heisenberg picture. In this all time dependence lies in operator, while wave function is independent of time.

1.9.1 Why called Heisenberg picture?

$$\begin{aligned} \hat{O}(t) &= e^{\frac{+iHt}{\hbar}} \hat{O}(0) e^{\frac{-iHt}{\hbar}} \\ \dot{\hat{O}}(t) &= \frac{i}{\hbar} H \hat{O} + \frac{-i}{\hbar} \hat{O} H \\ \dot{\hat{O}}(t) &= \frac{i}{\hbar} [H, \hat{O}] \end{aligned}$$

Which is Heisenberg equation of motion. That is why we call it Heisenberg picture. In between two extremes of Schrodinger and Heisenberg picture. There is an intermediate picture called interaction picture.

1.10 INTERACTION PICTURE:

Consider again the equation,

$$\langle \hat{O}(t) \rangle = \langle \psi(0) | e^{\frac{+iHt}{\hbar}} \hat{O}(0) e^{\frac{-iHt}{\hbar}} | \psi(0) \rangle$$

As

$$H = H_0 + V$$

H_0 - free Hamiltonian,

V - interaction part of Hamiltonian,

If the time dependence created by the interaction energy is only assigned to the state vector and rest of time dependence goes to the operator, then expectation value is written as,

$$\begin{aligned} \langle \hat{O}(t) \rangle &= \langle \psi(0) e^{\frac{+iVt}{\hbar}} \left| e^{\frac{+iH_0t}{\hbar}} \hat{O}(0) e^{\frac{-iH_0t}{\hbar}} \right| e^{\frac{-iVt}{\hbar}} \psi(0) \rangle \\ \langle \hat{O}(t) \rangle &= \langle \psi^I(t) | \hat{O}(t) | \psi^I(t) \rangle \end{aligned}$$

Interaction picture state vector

$$|\psi^I(t)\rangle = e^{\frac{-iVt}{\hbar}} |\psi(0)\rangle$$

Equation of motion

$$\left| \dot{\psi}^I(t) \right\rangle = \frac{-iV}{\hbar} |\psi^I(t)\rangle$$

This equation is simpler than ordinary schrodinger equation, but requires the calculation of $\hat{O}^I(t)$, where

$$\hat{O}^I(t) = e^{\frac{+iH_0t}{\hbar}} \hat{O}(0) e^{\frac{-iH_0t}{\hbar}}$$

In interaction picture both state vector and operator are time dependent. The interaction picture state vector,

$$|\psi^I(t)\rangle = \sum_n C_n(t) |n\rangle$$

The schrodinger picture state vector is

$$|\psi(t)\rangle = \sum_n c_n(t) |n\rangle = \sum_n C_n(t) e^{\frac{-i\omega_n t}{\hbar}} |n\rangle$$

Where

$$c_n(t) = C_n(t) e^{\frac{-i\omega_n t}{\hbar}}$$

The complete time dependence is given by $c_n(t)$, but due to interaction energy is given by $C_n(t)$.

1.11 PAULI SPIN MATRIX:

Another method to describe two- level atom is a use of 2×2 matrix notation. The eigen function U_a and U_b or eigen vectors $|a\rangle$ and $|b\rangle$ can be represented by the column vectors

$$\begin{aligned}|a\rangle &= U_a \leftrightarrow \begin{pmatrix} 1 \\ 0 \end{pmatrix} \\ |b\rangle &= U_b \leftrightarrow \begin{pmatrix} 0 \\ 1 \end{pmatrix}\end{aligned}$$

And the wave function and wave vector by the column vectors

$$\psi(r, t) = \begin{bmatrix} C_a \\ C_b \end{bmatrix}$$

And

$$|\psi(r, t)\rangle = \begin{pmatrix} C_a \\ C_b \end{pmatrix}$$

The energy and electric- dipole operators are written in terms of the Pauli spin matrices as

$$\begin{aligned}\sigma_x &= \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \\ \sigma_y &= \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix} \\ \sigma_z &= \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}\end{aligned}$$

These matrices are hermitian, but the spin- flip operators

$$\begin{aligned}\sigma_+ &= \frac{1}{2}(\sigma_x + i\sigma_y) = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} \\ \text{and } \sigma_- &= \frac{1}{2}(\sigma_x - i\sigma_y) = \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix}\end{aligned}$$

are not hermitian, σ_- flips the system from upper- level to a lower- level

$$\sigma_- \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix} \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \begin{pmatrix} 0 \\ 1 \end{pmatrix}$$

While, σ_+ flips the system from lower- level to the upper- level

$$\sigma_+ \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \begin{pmatrix} 1 \\ 0 \end{pmatrix}$$

1.12 TWO LEVEL ATOMIC SYSTEM AND HAMILTONIAN IN TERMS OF MATRICES:

State vector of a system can be written as,

$$|\psi(r, t)\rangle = C_a e^{-i\omega_a t} |a\rangle + C_b e^{-i\omega_b t} |b\rangle,$$

Which corresponds to the wave function

$$\psi(r, t) = C_a U_a(\vec{r}) e^{-i\omega_a t} + C_b U_b(\vec{r}) e^{-i\omega_b t}$$

The matrix form from for the unit vectors (eigen vectors) $|a\rangle$ and $|b\rangle$ are

$$\psi(r, t) \leftrightarrow C_a e^{-i\omega_a t} \begin{pmatrix} 1 \\ 0 \end{pmatrix} + C_b e^{-i\omega_b t} \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \begin{pmatrix} C_a e^{-i\omega_a t} \\ C_b e^{-i\omega_b t} \end{pmatrix}$$

The two- level Hamiltonian of the semi-classical treatment is given as,

$$H = \hbar\omega_a |a\rangle \langle a| + V_{ab} |a\rangle \langle b| + V_{ba} |b\rangle \langle a| + \hbar\omega_b |b\rangle \langle b|$$

In matrix notation

$$H = \begin{pmatrix} H_{aa} & H_{ab} \\ H_{ba} & H_{bb} \end{pmatrix} = \begin{pmatrix} \hbar\omega_a & V_{ab} \\ V_{ba} & \hbar\omega_b \end{pmatrix}$$

Thus the matrix form of the schrodinger equation is

$$i\hbar \frac{d}{dt} \begin{pmatrix} C_a e^{-i\omega_a t} \\ C_b e^{-i\omega_b t} \end{pmatrix} = \begin{pmatrix} \hbar\omega_a & V_{ab} \\ V_{ba} & \hbar\omega_b \end{pmatrix} \begin{pmatrix} C_a e^{-i\omega_a t} \\ C_b e^{-i\omega_b t} \end{pmatrix}$$

1.13 EXPECTATION VALUE OF DIPOLE MOMENT OPERATOR FOR A TWO-LEVEL ATOM:

The expectation value of any operator is given by

$$\langle \psi | \hat{O} | \psi \rangle = C_a C_a^* \hat{O}_{aa} + C_b C_b^* \hat{O}_{bb} + \left\{ C_a C_b^* \hat{O}_{ab} e^{-i\omega_{ab}t} + c.c \right\}$$

The expectation value of “ er ” is

$$\langle \psi | er | \psi \rangle = e C_a C_a^* \langle a | r | a \rangle + e C_b C_b^* \langle b | r | b \rangle + \left\{ e C_a C_b^* e^{-i(\omega_a - \omega_b)t} \langle b | r | a \rangle + c.c \right\}$$

As the diagonal matrix element of “ er ” between eigen states of the Hamiltonian generally vanishes.

$$\begin{aligned} er_{aa} &= \langle a | er | a \rangle = e \int d^3r U_a^*(r) r U_a(r) = 0 \\ er_{bb} &= \langle b | er | b \rangle = e \int d^3r U_b^*(r) r U_b(r) = 0 \\ er_{ab} &= \langle a | er | b \rangle = e \int d^3r U_a^*(r) r U_b(r) \end{aligned}$$

In matrix form

$$\begin{aligned} \langle er \rangle &= e \begin{pmatrix} 0 & r_{ab} \\ r_{ba} & 0 \end{pmatrix} \\ \langle er \rangle &= e C_a C_b^* e^{-i(\omega_a - \omega_b)t} r_{ba} + c.c \end{aligned}$$

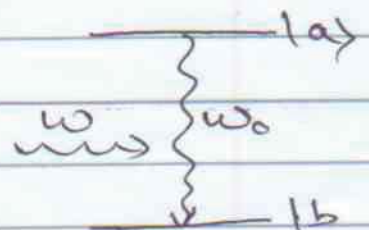
ATOM - FIELD INTERACTION SEMICLASSICAL THEORY

\Rightarrow Atom \rightarrow Quantum Mechanical

Field \rightarrow Classical

Two-Level atomic system.

$$|\psi(t)\rangle = c_a(t)|a\rangle + c_b(t)|b\rangle$$



Hamiltonian of the system

$$H = H_0 + H'$$

H_0 — free part of Hamiltonian

H' — perturbed part of Hamiltonian

Completeness relation for atomic system

$$|a\rangle\langle a| + |b\rangle\langle b| = 1$$

$$\Rightarrow H_0 = \underbrace{(|a\rangle\langle a| + |b\rangle\langle b|)}_1 H_0 (|a\rangle\langle a| + |b\rangle\langle b|)$$

$$\text{As } H_0|a\rangle = E_a|a\rangle, H_0|b\rangle = E_b|b\rangle$$

$$\Rightarrow H_0 = E_a|a\rangle\langle a| + E_b|b\rangle\langle b|$$

$$\Rightarrow H_0 = \hbar\omega_a|a\rangle\langle a| + \hbar\omega_b|b\rangle\langle b|$$

Interaction part of Hamiltonian is written as.

$$H' = -e\mathbf{r} \cdot \mathbf{E}(r, t)$$

under dipole approximation

$$\mathbf{E}(R, t) = \mathbf{E}(r, t) \approx \mathbf{E}(t) \approx E_0 \cos \omega t$$

$$H' = -e\mathbf{r} \cdot E_0 \cos \omega t$$

$$H' = [|a\rangle\langle a| + |b\rangle\langle b|] H' [|a\rangle\langle a| + |b\rangle\langle b|]$$

$$= |a\rangle\langle a| \overset{\rightarrow 0}{H'} |a\rangle\langle a| + |a\rangle\langle a| \overset{\rightarrow 0}{H'} |b\rangle\langle b| + |b\rangle\langle b| \overset{\rightarrow 0}{H'} |a\rangle\langle a| + |b\rangle\langle b| \overset{\rightarrow 0}{H'} |b\rangle\langle b|$$

$$+ |a\rangle\langle a| H' |b\rangle\langle b| + |b\rangle\langle b| H' |a\rangle\langle a|$$

$$\text{As } H'_{aa} = 0 = H'_{bb}$$

$$H' = |a\rangle\langle b| H'_{ab} + |b\rangle\langle a| H'_{ba}$$

where $H'_{ab} = \langle a | H' | b \rangle$

$$= -\langle a | e r | b \rangle E_0 \cos \omega t$$

$$= -p_{ab} E_0 \cos \omega t$$

and

$$H'_{ba} = -p_{ba} E_0 \cos \omega t$$

$$\Rightarrow H' = - \left(p_{ab} |a\rangle\langle b| + p_{ba} |b\rangle\langle a| \right) E(t)$$

where

$p_{ab} = p_{ba}^* = e \langle a | r | b \rangle$ — matrix element of the electric dipole moment and $E(t)$ is the field at the atom.

The time-development of the system is given by Schrodinger equation.

$$i\hbar \frac{d}{dt} |\psi(t)\rangle = H |\psi(t)\rangle$$

Substitute values of state-vector and Hamiltonian we get.

$$\dot{c}_a |a\rangle + \dot{c}_b |b\rangle = -\frac{i}{\hbar} [c_a \hbar \omega_a |a\rangle + c_b \hbar \omega_b |b\rangle + H_{ab} c_b |a\rangle + H_{ba} c_a |b\rangle]$$

Multiplying with $\langle a|$ and using

$$\begin{aligned} \langle a|a\rangle &= 1 & ; & \langle a|b\rangle = 0 \\ \langle b|b\rangle &= 1 & ; & \langle b|a\rangle = 0 \end{aligned}$$

$$\dot{c}_a = -\frac{i}{\hbar} [\hbar \omega_a c_a + H_{ab} c_b]$$

and

$$\dot{c}_b = -\frac{i}{\hbar} [\hbar \omega_b c_b + H_{ba} c_a]$$

Putting values of interaction Hamiltonian

$$\dot{c}_a = -i \omega_a c_a + \frac{i p_{ab}}{\hbar} E_0 \cos \omega t c_b$$

and

$$\dot{c}_b = -i \omega_b c_b + \frac{i p_{ba}}{\hbar} E_0 \cos \omega t c_a$$

Define

$$\Omega_R = \frac{|p_{ba}|}{\hbar} E_0 - \text{Rabi frequency}$$

$$p_{ba} = |p_{ba}| e^{i\phi}$$

$$\Rightarrow p_{ab} = p_{ba}^* = |p_{ba}| e^{-i\phi}$$

where ϕ is the phase of the dipole matrix element

$$\Rightarrow \dot{c}_a = -i\omega_a c_a + i\Omega_R e^{-i\phi} \cos\omega t c_b$$

$$\dot{c}_b = -i\omega_b c_b + i\Omega_R e^{i\phi} \cos\omega t c_a$$

Transform c_a & c_b (Sch. picture amplitudes) into slowly varying interaction picture amplitudes

$$C_a = c_a e^{i\omega_a t}$$

$$C_b = c_b e^{i\omega_b t}$$

Differentiating above eqn's

$$\dot{C}_a = \dot{c}_a e^{i\omega_a t} + i\omega_a c_a e^{i\omega_a t}$$

$$\begin{aligned} \dot{C}_a &= (-i\omega_a c_a + i\Omega_R e^{-i\phi} \cos\omega t c_b) e^{i\omega_a t} + i\omega_a c_a e^{i\omega_a t} \\ &= i\Omega_R e^{-i\phi} \cos\omega t c_b e^{i\omega_a t} \end{aligned}$$

$$= i\Omega_R e^{-i\phi} \cos\omega t e^{i\omega_a t} C_b$$

$$\omega_{ab} = \omega_a - \omega_b$$

Similarly.

$$\dot{C}_b = i\Omega_R e^{i\phi} \cos \omega t e^{-i\omega_{ab}t} C_a$$

$\omega_{ab} = \omega_0$ - transition frequency.

Expanding $\cos \omega t$.

$$\dot{C}_a = \frac{i\Omega_R}{2} e^{-i\phi} [e^{i\omega t + i\omega_0 t} + e^{-i\omega t + i\omega_0 t}] C_b$$

Neglecting rapidly oscillating terms like $e^{i(\omega + \omega_0)t}$

$$\dot{C}_a = \frac{i\Omega_R}{2} C_b e^{i(\omega_0 - \omega)t} e^{-i\phi}$$

Similarly.

$$\dot{C}_b = \frac{i\Omega_R}{2} C_a e^{-i(\omega_0 - \omega)t} e^{+i\phi}$$

where $\Delta = \omega_0 - \omega$. - detuning.

→ Consider atom initially in the excited state $C_a(0) = 1$

$$C_b(0) = 0$$

Assume resonance $\Delta = \omega_0 - \omega = 0$

$$C_a(t) = \cos \frac{\Omega_R t}{2}$$

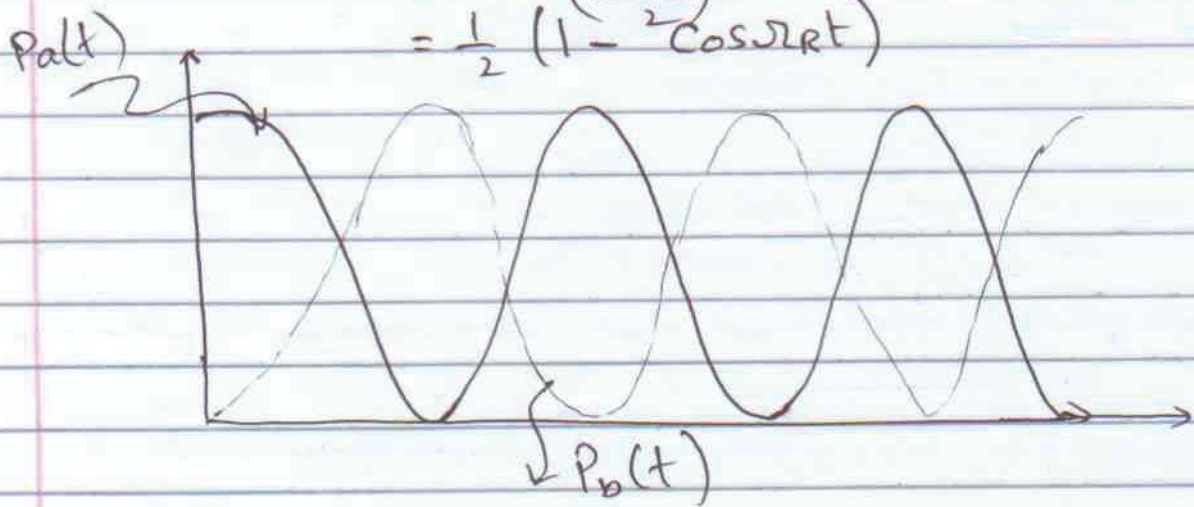
$$C_b(t) = i e^{i\phi} \sin \frac{\Omega_R t}{2}$$

Probability of atom in state $|a\rangle$
at time t

$$P_a(t) = |\langle \psi | a \rangle|^2 = |C_a(t)|^2 \\ = \cos^2\left(\frac{\Omega_R t}{2}\right) = \frac{1}{2} (1 + \cos \Omega_R t)$$

and

$$P_b(t) = |C_b(t)|^2 = \text{Probability of atom in state } |b\rangle \\ = \sin^2\left(\frac{\Omega_R t}{2}\right) \\ = \frac{1}{2} (1 - \cos \Omega_R t)$$



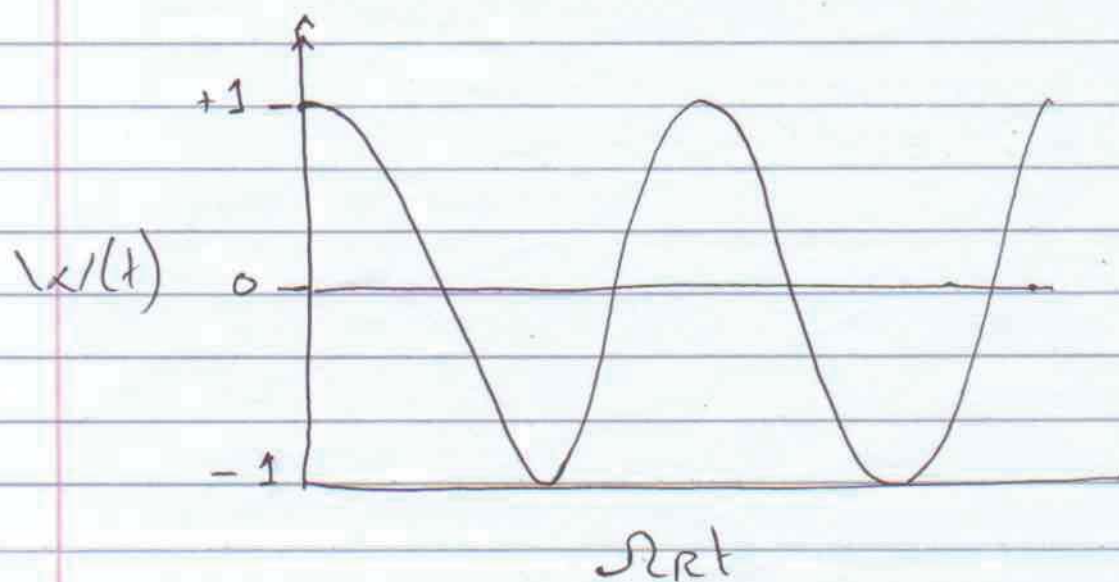
Atom oscillates with Rabi-freq.

Population inversion

$$W(t) = P_a(t) - P_b(t) \quad \text{at } \Delta = 0$$

$$W(t) = \cos^2\left(\frac{\Omega_R t}{2}\right) - \sin^2\left(\frac{\Omega_R t}{2}\right) = \cos \Omega_R t$$

It oscillates between -1 and +1



There are three frequency involved

- 1, $\omega_0 = \omega_a - \omega_b = \frac{E_a - E_b}{\hbar}$ — Transition frequency
- 2, ω frequency of the field
- 3, Rabi frequency $\Omega_R = \frac{|P_{ab}| \mathcal{E}}{\hbar}$

1.14 DENSITY OPERATOR

For a given physical system there exist a state vector $|\psi\rangle$ which contain all possible information about the system. If we want to extract a piece of information about the system we must calculate the expectation value of the corresponding operator \hat{O}

$$\langle \hat{O} \rangle_{Q.M} = \langle \psi | \hat{O} | \psi \rangle$$

In many situations we do not know $|\psi\rangle$ but we know P_ψ , probability of finding the system in $|\psi\rangle$. For such a situation we not only need to take quantum mechanical average but also the ensemble average over many identical systems that have been similarly prepared

$$\left\langle \langle \hat{O} \rangle_{Q.M} \right\rangle_{ensemble} = \sum_{\psi} P_{\psi} \langle \psi | \hat{O} | \psi \rangle$$

It is called a quantum statistical system.

$$\begin{aligned} \sum_n |n\rangle \langle n| &= 1 \\ \left\langle \langle \hat{O} \rangle_{Q.M} \right\rangle_{ensemble} &= \sum_n \sum_{\psi} P_{\psi} \langle \psi | n \rangle \langle n | \hat{O} | \psi \rangle \\ &= \sum_n \sum_{\psi} P_{\psi} \langle n | \hat{O} | \psi \rangle \langle \psi | n \rangle \\ &= \sum_n \langle n | \hat{O} \left[\sum_{\psi} P_{\psi} | \psi \rangle \langle \psi | \right] | n \rangle \\ &= \sum_n \langle n | \hat{O} \rho | n \rangle = \sum_n (\hat{O} \rho)_{nn} \end{aligned}$$

where

$$\rho = \sum_{\psi} P_{\psi} | \psi \rangle \langle \psi |$$

The sum of diagonal elements gives

$$\left\langle \langle \hat{O} \rangle_{Q.M} \right\rangle_{ensemble} = Tr(\hat{O} \rho) = Tr(\rho \hat{O})$$

where

$$\rho = \sum_{\psi} P_{\psi} | \psi \rangle \langle \psi |$$

is called density operator. In a particular case where all P_{ψ} are zero except the one for a state $|\psi_0\rangle$ then

$$\rho = |\psi_o\rangle \langle \psi_o|$$

and state is called a Pure state. It follows from the conservation of probability that $Tr(\rho) = 1$. For a pure state

$$Tr(\rho^2) = 1$$

For mixed case

$$Tr(\rho^2) < 1$$

1.15 EQUATION OF MOTION FOR DENSITY OPERATOR

$$\begin{aligned}\rho &= \sum_{\psi} P_{\psi} |\psi\rangle \langle \psi| \\ \dot{\rho} &= \sum_{\psi} P_{\psi} \left[\left| \dot{\psi} \right\rangle \langle \psi| + |\psi\rangle \left\langle \dot{\psi} \right| \right]\end{aligned}$$

From Schrodinger equation, we know that

$$\left| \dot{\psi} \right\rangle = \frac{1}{i\hbar} H |\psi\rangle$$

Also

$$\langle \dot{\psi} | = -\frac{1}{i\hbar} \langle \psi | H$$

Using these values, we get

$$\begin{aligned}\dot{\rho} &= \frac{1}{i\hbar} \sum_{\psi} P_{\psi} [H |\psi\rangle \langle \psi| - |\psi\rangle \langle \psi| H] \\ &= \frac{1}{i\hbar} [H\rho - \rho H]\end{aligned}$$

where ρ and H are operators.

$$\dot{\rho} = \frac{1}{i\hbar} [H, \rho]$$

It is called “Liouville equation” or “von Neumann equation” and it is equivalent to Schrodinger , but more general (because it has quantum mechanical as well as statistical aspect). In the above equation we have not included the decay of atomic levels due spontaneous emission. The excited atomic levels can also decay because of collision and other phenomena. The finite life time of the atomic level can be dicribed by adding phenominalogical decay terms to the density operator.

The decay rates can be incorporated in equation by a relaxation matrix Γ , which is defined by the equation

$$\Gamma_{nm} = \langle n | \Gamma | m \rangle = \gamma_n \delta_{nm}$$

$$\dot{\rho} = -\frac{i}{\hbar} [H, \rho] - \frac{1}{2} \{\Gamma, \rho\}; \quad \{\Gamma \rho + \rho \Gamma\} = \{\Gamma, \rho\}$$

Equation of motion for the density matrix elements is

$$\dot{\rho}_{ij} = \frac{1}{i\hbar} \sum_k [H_{ik} \rho_{kj} - \rho_{ik} H_{kj}] - \frac{1}{2} \sum_k [\Gamma_{ik} \rho_{kj} + \rho_{ik} \Gamma_{kj}]$$

1.16 TWO LEVEL ATOM

$$\begin{aligned} |\psi\rangle &= c_a(t) |a\rangle + c_b(t) |b\rangle \\ \rho &= |\psi\rangle \langle\psi| \\ \langle\psi| &= c_a^*(t) \langle a| + c_b^*(t) \langle b| \\ \rho &= |\psi\rangle \langle\psi| = |c_a(t)|^2 |a\rangle \langle a| + c_a(t) c_b^*(t) |a\rangle \langle b| + c_a^*(t) c_b(t) |b\rangle \langle a| + |c_b(t)|^2 |b\rangle \langle b| \\ \rho_{aa} &= \langle a | \rho | a \rangle = |c_a|^2 \quad \text{probability of upper state} \\ \rho_{bb} &= \langle b | \rho | b \rangle = |c_b|^2 \quad \text{probability of lower state} \\ \rho_{ab} &= c_a c_b^* \quad \text{is propotional to dipole moment} \\ \rho_{ba} &= c_b c_a^* = \rho_{ab}^* \end{aligned}$$

In matrix form

$$\rho = \begin{pmatrix} |c_a|^2 & c_a c_b^* \\ c_a^* c_b & |c_b|^2 \end{pmatrix}$$

$$P(z, t) = c_a c_b^* p_{ba} + c.c = \rho_{ab}(z, t) p_{ba} + c.c$$

In spiner notation,

$$\begin{aligned} |\psi\rangle &= \begin{pmatrix} c_a \\ c_b \end{pmatrix} & \langle\psi| &= \begin{pmatrix} c_a^* & c_b^* \end{pmatrix} \\ \rho &= \begin{pmatrix} c_a \\ c_b \end{pmatrix} \begin{pmatrix} c_a^* & c_b^* \end{pmatrix} = \begin{pmatrix} |c_a|^2 & c_a c_b^* \\ c_a^* c_b & |c_b|^2 \end{pmatrix} \end{aligned}$$

1.17 EQUATION OF MOTION FOR DENSITY MATIRX ELEMENTS

Consider $i = a, j = a$

$$\dot{\rho}_{aa} = \frac{1}{i\hbar} \sum_{k=a,b} [H_{ak} \rho_{ka} - \rho_{ak} H_{ka}] - \frac{1}{2} \sum_{k=a,b} [\Gamma_{ak} \rho_{ka} + \rho_{ak} \Gamma_{ka}]$$

using this

$$\begin{aligned}
\Gamma_{nm} &= \langle n|\Gamma|m\rangle = \gamma_n \delta_{nm} \quad \text{and} \\
H &= H_\circ + \gamma \\
\dot{\rho}_{aa} &= \frac{1}{i\hbar} \sum_{k=a,b} [(H_\circ)_{ak} \rho_{ka} - \rho_{ak} (H_\circ)_{ka}] - \frac{1}{2} [\gamma_a \rho_{aa} + \rho_{aa} \gamma_a] \\
&\quad + \frac{1}{i\hbar} \sum_{k=a,b} [V_{ak} \rho_{ka} - \rho_{ak} V_{ka}] \\
(H_\circ)_{ab} &= \langle a|H_\circ|b\rangle = E_b \langle a|b\rangle = 0 \\
(H_\circ)_{aa} &= \langle a|H_\circ|a\rangle = E_a \quad \text{and} \quad V_{aa}=V_{bb}=0 \\
\dot{\rho}_{aa} &= -\gamma_a \delta_{aa} + \frac{1}{i\hbar} \sum_{k=a,b} [V_{ab} \rho_{ba} - \rho_{ab} V_{ba}] \\
\dot{\rho}_{bb} &= -\gamma_b \delta_{bb} + \frac{1}{i\hbar} [V_{ba} \rho_{ab} - \rho_{ba} V_{ab}] \\
\dot{\rho}_{ab} &= -(i\omega_\circ + \gamma_{ab}) \delta_{ab} + \frac{1}{i\hbar} [V_{ab} \rho_{bb} - \rho_{aa} V_{ab}] \\
&\quad \text{where} \\
\gamma_{ab} &= \frac{1}{2}(\gamma_a + \gamma_b) \quad \text{and} \quad \omega_\circ = \frac{1}{\hbar}(E_a - E_b)
\end{aligned}$$

The population of excited level decays in time because of spontaneous emission. In soome cases the upper level decays to ground state lower level then

$$\dot{\rho}_{aa} = -\Gamma \rho_{aa} + \frac{1}{i\hbar} [V_{ab} \rho_{ba} - \rho_{ab} V_{ba}]$$

and

$$\dot{\rho}_{bb} = \Gamma \rho_{bb} + \frac{1}{i\hbar} [V_{ba} \rho_{ab} - \rho_{ba} V_{ab}]$$

Where Γ is the upper-to-lower level decay constant.

1 Quantization of the Free Electromagnetic Field

Dirac combined the wave and particle like aspects of light. Wave nature shows all the interference phenomena. Particle nature shows the excitation of a specific atom absorbing one photon of energy.

Classical field fails to explain

1. Spontaneous emission
2. Atomic decay
3. Lamb shift
4. Photon statistics

An interesting consequences of the quantization of the radiation is the fluctuations associated with the zero-point energy or so called vacuum fluctuations. These fluctuations have no classical analog and are responsible for many interacting phenomena in quantum optics.

1.1 Spontaneous Emission and Atomic Decay

A phenomena which we described phenomenologically in our treatment of semi-classical theory requires a quantum field. Spontaneous emission is often said to be the result of stimulating the atom by vacuum fluctuations.

1.2 Lamb Shift

According to the classical description of the field (*Dirac theory*) the $2S_{\frac{1}{2}}$ and $2P_{\frac{1}{2}}$ states in the hydrogen atom should have equal energies. Experimentally the two levels differ by approximately 1057 MHz. a fully quantized treatment of the field and atomic systems gives impressive agreement with the experimentally observed shift, because of the radiative correction due to the interaction between the atomic electron and the vacuum shift the $2S_{\frac{1}{2}}$ level higher in energy by around 1057 MHz relative to the $2P_{\frac{1}{2}}$ level.

1.3 Photon Statistics

In order to explain the photon statistics the concept of a particle, the photon is either necessary or convenient. For the quantization of the electromagnetic field in free space, it is convenient to begin with the classical description of the field based on Maxwell's equations. In MKS system

$$\begin{aligned}\nabla \cdot D &= 0 \\ \nabla \cdot B &= 0 \\ \nabla \times E &= -\frac{\partial B}{\partial t} \\ \nabla \times H &= \frac{\partial D}{\partial t}\end{aligned}$$

where in free space

$$\begin{aligned} D &= \epsilon_0 E \\ B &= \mu_0 H \end{aligned}$$

here ϵ_0 and μ_0 are the free space permittivity and permeability respectively and

$$\mu_0 \epsilon_0 = \frac{1}{c^2}$$

where c is the speed of light in vacuum. Using these Maxwell's equations we know that $E(r, t)$ and also $B(r, t)$ satisfies the wave equation

$$\nabla^2 E - \frac{1}{c^2} \frac{\partial^2 E}{\partial t^2} = 0$$

Following the Dirac approach we associate each mode of the radiation field with a quantized simple Harmonic oscillator. Energy of the Harmonic oscillator (classically) is given by the hamiltonian

$$H = \frac{1}{2} m \omega^2 x^2 + \frac{p^2}{2m}$$

and quantum mechanically it is written as

$$H = \frac{1}{2} m \omega^2 \hat{x}^2 + \frac{\hat{p}^2}{2m}$$

2 Mode Expansion of The Field

2.1 Quantization of Field Inside the Cavity of Length L

Electric field is linearly polarized in the x direction. Expanding the field in the normal modes of the cavity

$$E_x(z, t) = \sum_j A_j q_j(t) \sin(k_j z)$$

j corresponds to different modes such that

$$L = j \frac{\lambda}{2}; \lambda = \frac{2\pi}{k} \text{ and } L = \frac{j\pi}{k_j}, \text{ where } j = 1, 2, 3, \dots$$

Where q_j is the normal mode amplitude with the dimensions of length (position) and

$$\begin{aligned} A_j &= \left(\frac{2\omega_j^2 m_j}{V \epsilon_0} \right)^{1/2} \\ \text{where } \omega_j &= ck_j = \frac{j\pi c}{L} \end{aligned}$$

is the cavity eigen frequency. $V = LA$ is the volume (A is the transverse area of the optical resonator) m_j is a constant with the dimensions of mass, included to make an analogy with SHO nothing to do with mass of photon. The E.M.F is assumed to be transverse with electric field polarized in the x-direction. Such field satisfies

$$\nabla \cdot E = 0$$

The nonvanishing component of the magnetic field in the cavity is obtained by using Maxwell's 4th equation i.e,

$$\nabla \times H = \frac{\partial D}{\partial t} = \epsilon_0 \frac{\partial E}{\partial t}$$

$$\begin{aligned} \nabla \times H &= \begin{pmatrix} \hat{i} & \hat{j} & \hat{k} \\ \frac{\partial}{\partial x} & \frac{\partial}{\partial y} & \frac{\partial}{\partial z} \\ H_x & H_y & H_z \end{pmatrix} \\ &= \hat{i} \left(\frac{\partial H_z}{\partial y} - \frac{\partial H_y}{\partial z} \right) + \hat{j} \left(\frac{\partial H_x}{\partial z} - \frac{\partial H_z}{\partial x} \right) + \hat{k} \left(\frac{\partial H_y}{\partial x} - \frac{\partial H_x}{\partial y} \right) \end{aligned}$$

x-component of $(\nabla \times H)_x$ is written as

$$\left(\frac{\partial H_z}{\partial y} - \frac{\partial H_y}{\partial z} \right)_x = \epsilon_0 \frac{\partial E_x}{\partial t}$$

$H_z = 0$, z is the direction of propagation

$$\begin{aligned} -\frac{\partial H_y}{\partial z} &= \epsilon_0 \frac{\partial}{\partial t} \left(\sum_j A_j q_j(t) \sin(k_j z) \right) \\ &= \epsilon_0 \sum_j A_j \dot{q}_j(t) \sin(k_j z). \end{aligned}$$

As

$$\sin(k_j z) = -\frac{1}{k_j} \frac{\partial}{\partial z} \cos(k_j z),$$

putting the value of $\sin(k_j z)$ in the above equation we can write

$$\begin{aligned} -\frac{\partial H_y}{\partial z} &= -\frac{\partial}{\partial z} \left(\sum_j A_j \left(\frac{\dot{q}_j(t) \epsilon_0}{k_j} \right) \cos(k_j z) \right), \\ H_y &= \sum_j A_j \left(\frac{\dot{q}_j(t) \epsilon_0}{k_j} \right) \cos(k_j z). \end{aligned}$$

The classical Hamiltonian of the field i.e, the (total energy of the field) is

$$H = \frac{1}{2} \int_v d\tau (\epsilon_0 E_x^2 + \mu_0 H_y^2)$$

where the integration is over the volume of the cavity. Substituting the values of E_x and H_y in the above equation and performing the integration we get,

$$\begin{aligned} H &= \frac{1}{2} \sum_j \left(m_j \omega_j^2 q_j^2 + m_j \dot{q}_j^2 \right) \\ &= \frac{1}{2} \sum_j \left(m_j \omega_j^2 q_j^2 + \frac{p_j^2}{m_j} \right) \end{aligned}$$

where $p_j = m_j \dot{q}_j$ is the canonical momentum of the j th mode. The above equation expresses the hamiltonian of the radiation field as a sum of independent oscillator energies. Each mode of the field is dynamically equivalent to a mechanical harmonic oscillator.

3 Quantization

The present dynamical problem can be quantized by identifying q_j and p_j as operators, which obey the commutation relations

$$\begin{aligned} [\hat{q}_j, \hat{p}_{j'}] &= i\hbar \delta_{jj'} \\ [\hat{q}_j, \hat{q}_{j'}] &= [\hat{p}_j, \hat{p}_{j'}] = 0. \end{aligned}$$

It can be transformed as

$$\hat{a}_j = \frac{1}{\sqrt{2m_j\hbar\omega_j}} (m_j\omega_j\hat{q}_j + i\hat{p}_j) \exp(i\omega_j t)$$

and

$$\hat{a}_j^\dagger = \frac{1}{\sqrt{2m_j\hbar\omega_j}} (m_j\omega_j\hat{q}_j - i\hat{p}_j) \exp(-i\omega_j t)$$

$$\hat{q}_j = \left(\hat{a}_j \exp(-i\omega_j t) + \hat{a}_j^\dagger \exp(i\omega_j t) \right) \sqrt{\frac{\hbar}{2m_j\omega_j}}$$

$$\hat{p}_j = -i\sqrt{\frac{m_j\omega_j\hbar}{2}} \left(\hat{a}_j \exp(-i\omega_j t) - \hat{a}_j^\dagger \exp(i\omega_j t) \right)$$

The commutation relations between \hat{a}_j and \hat{a}_j^\dagger follow from those between \hat{q}_j and \hat{p}_j ,

$$\begin{aligned} [\hat{a}_j, \hat{a}_j^\dagger] &= \frac{1}{2m_j\hbar\omega_j} \left[-im_j\omega_j [\hat{q}_j, \hat{p}_j] + im_j\omega_j [\hat{p}_j, \hat{q}_j] \right] \\ &= \frac{1}{2m_j\hbar\omega_j} [-im_j\omega_j (i\hbar) + im_j\omega_j (-i\hbar)] \\ &= 1. \end{aligned}$$

Similarly

$$\begin{aligned}\left[\hat{a}_j, \hat{a}_{j'}\right] &= \left[\hat{a}_j^\dagger, \hat{a}_{j'}^\dagger\right] = 0 \\ \left[\hat{a}_j, \hat{a}_{j'}^\dagger\right] &= \delta_{jj'}\end{aligned}$$

The operators \hat{a} and \hat{a}^\dagger are referred to as the destruction and creation operators, they are not hermitian. Substituting the value of \hat{q}_j and \hat{p}_j in the equation for Hamiltonian we get

$$\begin{aligned}H &= \sum_j \frac{1}{2} m_j \omega_j^2 \left(\frac{\hbar}{2m_j \omega_j} \right) \left(a_j^2 \exp(-2i\omega_j t) + a_j^{\dagger 2} \exp(2i\omega_j t) + \hat{a}_j \hat{a}_j^\dagger + \hat{a}_j^\dagger \hat{a}_j \right) \\ &\quad + \frac{1}{2m_j} \left(-\frac{m_j \hbar \omega_j}{2} \right) \left(a_j^2 \exp(-2i\omega_j t) + a_j^{\dagger 2} \exp(2i\omega_j t) - \hat{a}_j \hat{a}_j^\dagger - \hat{a}_j^\dagger \hat{a}_j \right) \\ &= \sum_j \left(\frac{\hbar \omega_j}{2} \right) \left(\hat{a}_j \hat{a}_j^\dagger + \hat{a}_j^\dagger \hat{a}_j \right)\end{aligned}$$

As

$$\begin{aligned}\left[\hat{a}_j, \hat{a}_j^\dagger\right] &= 1 \\ \left(\hat{a}_j \hat{a}_j^\dagger - \hat{a}_j^\dagger \hat{a}_j\right) &= 1 \\ \hat{a}_j \hat{a}_j^\dagger &= \hat{a}_j^\dagger \hat{a}_j + 1\end{aligned}$$

thus we can write

$$H = \sum_j \hbar \omega_j \left(\hat{a}_j^\dagger \hat{a}_j + \frac{1}{2} \right)$$

In terms of \hat{a}_j and \hat{a}_j^\dagger , the electric and magnetic fields can be written as

$$\begin{aligned}E_x(z, t) &= \sum_j \varepsilon_j \left(\hat{a}_j \exp(-i\omega_j t) + \hat{a}_j^\dagger \exp(i\omega_j t) \right) \sin(k_j z) \\ H_y(z, t) &= -i\epsilon_0 c \sum_j \varepsilon_j \left(\hat{a}_j \exp(-i\omega_j t) - \hat{a}_j^\dagger \exp(i\omega_j t) \right) \cos(k_j z)\end{aligned}$$

where the quantity

$$\varepsilon_j = \left(\frac{\hbar \omega_j}{\epsilon_0 V} \right)^{1/2}$$

has the dimensions of an electric field.

3.1 Quantization of Field Inside a Large Cavity of Finite Length L

Consider the field in a large but finite cubic cavity of side L . We consider the running wave solutions instead of the standing wave solutions. The classical electric and magnetic field can be expanded in terms of plane waves.

$$E(r, t) = \sum_k \hat{\epsilon}_k \varepsilon_k \alpha_k \exp(-i\omega_k t + ik \cdot r) + c.c$$

using Maxwell's equation i.e,

$$\nabla \times H = \frac{\partial D}{\partial t}$$

we get

$$H(r, t) = \frac{1}{\mu_0} \sum_k \frac{k \times \hat{\epsilon}_k}{\omega_k} \varepsilon_k \alpha_k \exp(-i\omega_k t + ik \cdot r) + c.c$$

where the summation is taken over an infinite discrete set of values of wave vector $k = (k_x, k_y, k_z)$, ϵ_k is a unit polarization vector, α_k is a dimensionless amplitude and

$$\varepsilon_k = \left(\frac{\hbar \omega_k}{2\epsilon_0 V} \right)^{1/2}$$

Periodic boundary conditions require that

$$k_x = \frac{2\pi n_x}{L}, k_y = \frac{2\pi n_y}{L}, k_z = \frac{2\pi n_z}{L}$$

where n_x, n_y, n_z are integers $0, \pm 1, \pm 2, \dots$. A set of numbers (n_x, n_y, n_z) defines a mode of electromagnetic field. For transverse field

$$\nabla \cdot D = 0$$

which requires

$$\vec{k} \cdot \hat{\epsilon}_k = 0$$

There are two independent polarization directions of $\hat{\epsilon}_k$ for each k . Changing from a discrete distribution of modes to a continuous distribution i.e,

$$\sum_k \Rightarrow 2 \left(\frac{L}{2\pi} \right)^3 \int d^3k$$

where factor of 2 accounts for two possible states of polarization. The number of modes available in a cavity is infinite, however the number of modes whose frequency lies between ω and $\omega + d\omega$ is finite. This is the same number of field modes having the magnitude of k , between k and $k + dk$. Making

transformation from rectangular coordinates (k_x, k_y, k_z) to the polar coordinates $(k \sin \theta \cos \phi, k \sin \theta \sin \phi, k \cos \theta)$, the volume element in k -space is written as

$$\begin{aligned} d^3k &= k^2 dk \sin \theta d\theta d\phi \\ &= \frac{\omega^2}{c^3} d\omega \sin \theta d\theta d\phi. \end{aligned}$$

The total number of modes in the volume L^3 in the range between ω and $\omega + d\omega$ is given by

$$\begin{aligned} dN &= 2 \left(\frac{L}{2\pi} \right)^3 \frac{\omega^2 d\omega}{c^3} \int_0^\pi \sin \theta d\theta \int_0^{2\pi} d\phi \\ &= \frac{L^3 \omega^2}{\pi^2 c^3} d\omega \end{aligned}$$

Radiation field is quantized by identifying α_k and α_k^* by the harmonic oscillator operators \hat{a}_k and \hat{a}_k^\dagger respectively, which satisfy the commutation relation

$$\left[\hat{a}_k, \hat{a}_k^\dagger \right] = 1$$

The quantized electric and magnetic fields takes the form

$$\begin{aligned} E(r, t) &= \sum_k \hat{\epsilon}_k \epsilon_k \hat{a}_k \exp(-i\omega_k t + ik \cdot r) + H.C \\ H(r, t) &= \frac{1}{\mu_0} \sum_k \frac{k \times \hat{\epsilon}_k}{\omega_k} \epsilon_k \hat{a}_k \exp(-i\omega_k t + ik \cdot r) + H.C \end{aligned}$$

where H.C is Hermetian conjugate. Separating positive and negative frequency parts of these field operators

$$\begin{aligned} E^+(r, t) &= \sum_k \hat{\epsilon}_k \epsilon_k \hat{a}_k \exp(-i\omega_k t + ik \cdot r) \\ E^-(r, t) &= \sum_k \hat{\epsilon}_k \epsilon_k \hat{a}_k^\dagger \exp(i\omega_k t - ik \cdot r) \end{aligned}$$

where $E^+(r, t)$ is the annihilation operator and $E^-(r, t)$ is the creation operator.

4 Fock or Number States of Radiation Field

Consider a single mode of the field of frequency ω having creation and annihilation operators \hat{a}^\dagger and \hat{a} respectively. Let $|n\rangle$ be the energy eigen state corresponding to the energy eigen value E_n , i.e.

$$\begin{aligned} H |n\rangle &= \hbar\omega \left(a^\dagger a + \frac{1}{2} \right) |n\rangle \\ &= E_n |n\rangle \end{aligned} \quad (1)$$

applying operator \hat{a} from the left of the eigenstates we have

$$\begin{aligned} Ha |n\rangle &= \hbar\omega \left(a^\dagger a + \frac{1}{2} \right) a |n\rangle \\ [a, a^\dagger] &= aa^\dagger - a^\dagger a = 1 \\ \implies aa^\dagger - 1 &= a^\dagger a \end{aligned} \quad (2)$$

Putting in Eq. (2) we get

$$\begin{aligned} Ha |n\rangle &= \hbar\omega \left(aa^\dagger - 1 + \frac{1}{2} \right) a |n\rangle \\ &= \hbar\omega \left(aa^\dagger a - a + \frac{a}{2} \right) |n\rangle \\ &= a\hbar\omega \left(a^\dagger a + \frac{1}{2} - 1 \right) |n\rangle \\ &= a(\hbar\omega(a^\dagger a + \frac{1}{2}) - \hbar\omega) |n\rangle \\ &= a(E_n - \hbar\omega) |n\rangle \\ &= (E_n - \hbar\omega)a |n\rangle \end{aligned}$$

where $a |n\rangle$ is an energy eigen state with eigen value $(E_n - \hbar\omega)$. Operator a lowers the energy and therefore it is called annihilation, destruction or absorption operator.

$$\implies |n-1\rangle = \frac{a}{\alpha_n} |n\rangle,$$

is an energy eigen state but with the reduced eigen value i.e.

$$\begin{aligned} E_{n-1} &= (E_n - \hbar\omega) \\ H |n-1\rangle &= E_{n-1} |n-1\rangle, \end{aligned}$$

and α_n is a constant which will be determined from the normalization condition,

$$\langle n-1 | n-1 \rangle = 1.$$

If we repeat this procedure n times we move down the energy ladder in steps of $\hbar\omega$ until we obtain

$$Ha |0\rangle = (E_0 - \hbar\omega) a |0\rangle$$

E_0 is the ground state energy . $E_n - \hbar\omega$ is smaller than E_0 i.e, $E_n - \hbar\omega$ is negative. Since energy eigen value cannot be negative

$$a |0\rangle = 0$$

The state $|0\rangle$ is called the vacuum state. (in which no photon is excited).

$$\begin{aligned} H |0\rangle &= \hbar\omega \left(a^\dagger a + \frac{1}{2} \right) |0\rangle \\ &= \frac{1}{2} \hbar\omega |0\rangle \\ \implies E_0 &= \frac{1}{2} \hbar\omega \end{aligned}$$

is the energy of the ground state. Now we go step by step up as

$$\begin{aligned} E_{n-1} &= E_n - \hbar\omega \\ E_n &= E_{n-1} + \hbar\omega \end{aligned}$$

For $n = 1$ we can write

$$\begin{aligned} E_1 &= E_0 + \hbar\omega \\ &= \frac{1}{2} \hbar\omega + \hbar\omega = \frac{3}{2} \hbar\omega \end{aligned}$$

Similarly

$$\begin{aligned} E_2 &= E_1 + \hbar\omega \\ &= \frac{3}{2} \hbar\omega + \hbar\omega \\ &= \frac{5}{2} \hbar\omega \end{aligned}$$

It can also be written as

$$\begin{aligned} E_2 &= \left(2 + \frac{1}{2}\right) \hbar\omega, \\ &\vdots \\ E_n &= \left(n + \frac{1}{2}\right) \hbar\omega \\ \implies H |n\rangle &= E_n |n\rangle \end{aligned}$$

$$\begin{aligned} \hbar\omega \left(a^\dagger a + \frac{1}{2} \right) |n\rangle &= \left(n + \frac{1}{2} \right) \hbar\omega |n\rangle \\ \implies a^\dagger a |n\rangle &= n |n\rangle \end{aligned}$$

$|n\rangle$ is also an energy eigen state of the number operator

$$n = a^\dagger a$$

The normalization constant can be now calculated

$$\langle n-1 | n-1 \rangle = 1$$

as

$$\begin{aligned} |n-1\rangle &= \frac{\hat{a}}{\alpha_n} |n\rangle \\ \langle n-1| &= \frac{\langle n| a^\dagger}{\alpha_n^*} \\ \langle n-1 | n-1 \rangle &= \frac{\langle n| a^\dagger a |n\rangle}{\alpha_n^* \alpha_n} \\ &= \frac{1}{|\alpha_n|^2} \langle n| a^\dagger a |n\rangle \\ 1 &= \frac{n}{|\alpha_n|^2} \langle n | n \rangle = |\alpha_n|^2 = n \\ \alpha &= \sqrt{n} e^{i\phi} \end{aligned}$$

If we take the phase of the normalization constant α_n to be zero then $\alpha_n = \sqrt{n}$

$$\begin{aligned} a |n\rangle &= \alpha_n |n-1\rangle \\ &= \sqrt{n} |n-1\rangle \end{aligned}$$

now for operator a^\dagger

$$\begin{aligned} H a^\dagger |n\rangle &= \hbar\omega \left(a^\dagger a + \frac{1}{2} \right) a^\dagger |n\rangle \\ &= \hbar\omega \left(a^\dagger a a^\dagger + \frac{a^\dagger}{2} \right) |n\rangle \end{aligned}$$

using $aa^\dagger = a^\dagger a + 1$,

$$\begin{aligned} H a^\dagger |n\rangle &= \hbar\omega \left(a^\dagger a^\dagger a + a^\dagger + \frac{a^\dagger}{2} \right) |n\rangle \\ &= a^\dagger (E_n + \hbar\omega) |n\rangle \\ (\hbar\omega \left(a^\dagger a + \frac{1}{2} \right) a^\dagger |n\rangle) &= (E_n + \hbar\omega) a^\dagger |n\rangle \end{aligned}$$

Thus $a^\dagger |n\rangle$ is also an energy eigen state of the field with eigen value $E_n + \hbar\omega$.
We define

$$\begin{aligned} |n+1\rangle &= \frac{\hat{a}^\dagger}{\beta_n} |n\rangle \\ E_{n+1} &= E_n + \hbar\omega \\ \implies H |n+1\rangle &= E_{n+1} |n+1\rangle \end{aligned}$$

using the same procedure we get

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

A repeated use of the above equation gives,

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

The energy eigen states $|n\rangle$ are called fock states or photon number states. They form a complete set of state i.e.

$$\sum_{n=0}^{\infty} |n\rangle \langle n| = 1$$

The energy eigen value are discrete in contrast to classical electromagnetic theory where energy can have any value. State vector is written as the superposition of energy eigen states. i.e

$$|\Psi\rangle = \sum_n c_n |n\rangle$$

where c_n are complex coefficients. The energy $E_0 = \frac{1}{2}\hbar\omega$ is called zero-point energy. The energy levels for Q.M oscillations associated with the electromagnetic field are given as

$$\begin{array}{rcl} \text{-----} \frac{a^\dagger \uparrow}{a^\dagger \uparrow} \text{-----} E_{n+1} & = & \left(n + \frac{3}{2} \right) \hbar\omega \\ \text{-----} \frac{a \downarrow}{a \downarrow} \text{-----} E_n & = & \left(n + \frac{1}{2} \right) \hbar\omega \\ \text{-----} E_{n-1} & = & \left(n - \frac{1}{2} \right) \hbar\omega \\ & & \vdots \\ \text{-----} E_2 & = & \frac{5}{2} \hbar\omega \\ \text{-----} E_1 & = & \frac{3}{2} \hbar\omega \\ \text{-----} E_0 & = & \frac{1}{2} \hbar\omega \end{array}$$

The operators a and a^\dagger are not hermitian but some of the combinations are Hermitian such as,

$$\begin{aligned} a_1 &= (a + a^\dagger)/2, \\ a_2 &= (a - a^\dagger)/2i. \end{aligned}$$

Different energy eigen states of the field are orthogonal. The only non-vanishing matrix elements of a and a^\dagger are of the types;

$$\begin{aligned}\langle n-1 \mid a \mid n \rangle &= \sqrt{n} \\ \langle n+1 \mid a^\dagger \mid n \rangle &= \sqrt{n+1}\end{aligned}$$

An important property of $|n\rangle$ is that the expectation value of the single mode linearly polarized field operator vanishes. Using

$$E_x(z, t) = \varepsilon(ae^{-i\omega t} + a^\dagger e^{i\omega t}) \sin kz,$$

or

$$E(r, t) = \varepsilon a e^{-i\omega t + ik.r} + \varepsilon^* a^\dagger e^{i\omega t - ik.r}$$

$$\langle n \mid E(r, t) \mid n \rangle = \varepsilon \langle n \mid a \mid n \rangle e^{-i\omega t + ik.r} + \varepsilon^* \langle n \mid a^\dagger \mid n \rangle e^{i\omega t - ik.r} = 0$$

Now in order to find the average value of $\langle E^2 \rangle$, we write

$$\begin{aligned}\langle n \mid E^2 \mid n \rangle &= |\varepsilon|^2 \langle n \mid aa^\dagger + a^\dagger a \mid n \rangle + \varepsilon^2 e^{-2i\omega t + 2ik.r} \langle n \mid a^2 \mid n \rangle \\ &+ \varepsilon^2 e^{2i\omega t - 2ik.r} \langle n \mid a^{\dagger 2} \mid n \rangle.\end{aligned}$$

$$\begin{aligned}\langle E^2 \rangle &= (2n+1) |\varepsilon|^2 = 2 \left(n + \frac{1}{2} \right) |\varepsilon|^2 \\ \Delta E^2 &= \langle E^2 \rangle - \langle E \rangle^2 \\ &= 2 \left(n + \frac{1}{2} \right) |\varepsilon|^2,\end{aligned}$$

as $\langle E \rangle^2 = 0$. For $n = 0$ i.e. in vacuum

$$\Delta E^2 \neq 0,$$

but is equal to

$$\Delta E^2 = |\varepsilon|^2$$

From these equations we conclude that the mean value is zero but fluctuations are present. These fluctuations are considered to be responsible for spontaneous emission, Lamb shift etc.

1 The Coherent Photon States

The single-mode states of physical importance are not the individual number states $|n\rangle$ (because the electromagnetic wave generated by practical light source do not have definite numbers of photons), but the linear superposition of states $|n\rangle$. There is a wide variety of possible superposition states.

A superposition state can be constructed for which uncertainties in the expectation values of the phase operators $\hat{\cos}\phi$ and $\hat{\sin}\phi$ are both equal to zero. Such states have $\Delta n = \infty$. They cannot be excited in any real experiment. Another kind is the coherent state. A coherent state has equal amount of uncertainties in amplitude and phase. A field in coherent state is in a minimum uncertainty state. For coherent state an electric field variation approaches that of classical wave of stable amplitude and fixed phase, in the limit of high excitation. They are important because, they are the closest quantum mechanical approach to a classical electromagnetic wave. A single mode laser operated well above threshold generates a coherent state excitation.

The coherent state $|\alpha\rangle$ is the eigen state of the positive frequency part of the electric field operator or the eigen state of the destruction operator of the field.

$$\hat{a} |\alpha\rangle = \alpha |\alpha\rangle,$$

where α is complex, $|\alpha\rangle$ in terms of linear superposition of number state $|n\rangle$ is given by

$$\begin{aligned} |\alpha\rangle &= \sum_{n=0}^{\infty} C_n |n\rangle \\ \hat{a} |\alpha\rangle &= \sum_{n=0}^{\infty} C_n \hat{a} |n\rangle \\ &= \sum_{n=0}^{\infty} C_n \sqrt{n} |n-1\rangle \\ &= 0 + C_1 \sqrt{1} |0\rangle + C_2 \sqrt{2} |1\rangle + \dots \\ &= \sum_{n=0}^{\infty} C_{n+1} \sqrt{n+1} |n\rangle \end{aligned} \tag{1}$$

From eqn(1) multiplying with α we can write

$$\begin{aligned} \alpha |\alpha\rangle &= \sum_{n=0}^{\infty} C_n \alpha |n\rangle \\ \hat{a} |\alpha\rangle &= \sum_{n=0}^{\infty} C_n \alpha |n\rangle \end{aligned} \tag{b}$$

comparing eqn(a) and (b)

$$\begin{aligned}
C_{n+1}\sqrt{n+1} &= C_n\alpha \\
C_n\sqrt{n} &= C_{n-1}\alpha \\
C_n &= \frac{\alpha}{\sqrt{n}}C_{n-1} \\
C_n &= \frac{\alpha}{\sqrt{n}}\frac{\alpha}{\sqrt{n-1}}\frac{\alpha}{\sqrt{n-2}}\dots\frac{\alpha}{\sqrt{1}}C_0 \\
&= \frac{\alpha^n}{\sqrt{n!}}C_0
\end{aligned}$$

putting in Eqn(1),

$$|\alpha\rangle = C_0 \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} |n\rangle$$

the constant C_0 can be found by normalization

$$\begin{aligned}
\langle\alpha|\alpha\rangle &= C_0^*C_0 \sum_{n=0}^{\infty} \sum_{m=0}^{\infty} \frac{\alpha^{*n}}{\sqrt{n!}} \frac{\alpha^m}{\sqrt{m!}} \langle n|m\rangle \\
\langle\alpha|\alpha\rangle &= |C_0|^2 \sum_{n=0}^{\infty} \frac{|\alpha|^{2n}}{n!}
\end{aligned}$$

using

$$\sum_{n=0}^{\infty} \frac{x^n}{n!} = e^x$$

we can write

$$\begin{aligned}
1 &= |C_0|^2 e^{|\alpha|^2} \\
|C_0|^2 &= e^{-|\alpha|^2} \\
C_0 &= e^{-\frac{|\alpha|^2}{2}}
\end{aligned}$$

Therefore

$$|\alpha\rangle = e^{-\frac{|\alpha|^2}{2}} \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} |n\rangle$$

Another way of proving the above relation which interpret $|\alpha\rangle$ as a superposition of number state is,

$$|\alpha\rangle = \sum_n |n\rangle \langle n|\alpha\rangle \quad (1)$$

where $\sum_n |n\rangle \langle n| = 1$ is the completeness relation for number state. As

$$|n\rangle = \frac{(a^\dagger)^n}{\sqrt{n!}} |0\rangle \quad (2)$$

$$\langle n | = \langle 0 | \frac{(a)^n}{\sqrt{n!}} \quad (3)$$

Putting in Eqn(1) we can write as

$$| \alpha \rangle = \sum_n | n \rangle \langle 0 | \frac{(a)^n}{\sqrt{n!}} | \alpha \rangle$$

we know that

$$\begin{aligned} \hat{a} | \alpha \rangle &= \alpha | \alpha \rangle \\ (\hat{a})^n | \alpha \rangle &= \alpha^n | \alpha \rangle \\ \Rightarrow | \alpha \rangle &= \sum_n | n \rangle \frac{(\alpha)^n}{\sqrt{n!}} \langle 0 | \alpha \rangle \end{aligned} \quad (4)$$

The value of $\langle 0 | \alpha \rangle$ is obtained by normalization i.e

$$\langle \alpha | \alpha \rangle = 1$$

$$\begin{aligned} \langle \alpha | \alpha \rangle &= \sum_n \sum_m \frac{(\alpha^*)^m}{\sqrt{m!}} \frac{(\alpha)^n}{\sqrt{n!}} \langle m | n \rangle |\langle 0 | \alpha \rangle|^2 \\ &= \sum_n \frac{(\alpha^* \alpha)^n}{n!} |\langle 0 | \alpha \rangle|^2 \\ &= \sum_n \frac{(|\alpha|^2)^n}{n!} |\langle 0 | \alpha \rangle|^2 \\ \Rightarrow 1 &= e^{|\alpha|^2} |\langle 0 | \alpha \rangle|^2 \\ |\langle 0 | \alpha \rangle|^2 &= e^{-|\alpha|^2} \\ \langle 0 | \alpha \rangle &= e^{-\frac{|\alpha|^2}{2}} \end{aligned}$$

Putting in Eqn(4) we get

$$| \alpha \rangle = e^{-\frac{|\alpha|^2}{2}} \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} | n \rangle$$

hence proved.

Some other representations of the coherent state

$$\begin{aligned}
|\alpha\rangle &= e^{-\frac{|\alpha|^2}{2}} \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} |n\rangle \\
|n\rangle &= \frac{(a^\dagger)^n}{\sqrt{n!}} |0\rangle \\
|\alpha\rangle &= e^{-\frac{|\alpha|^2}{2}} \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} \frac{(a^\dagger)^n}{\sqrt{n!}} |0\rangle \\
|\alpha\rangle &= e^{-\frac{|\alpha|^2}{2}} e^{\alpha a^\dagger} |0\rangle
\end{aligned}$$

since we know that $e^{-\alpha^* a} |0\rangle = |0\rangle$

$$\begin{aligned}
\Rightarrow |\alpha\rangle &= e^{-\frac{|\alpha|^2}{2}} e^{\alpha a^\dagger} e^{-\alpha^* a} |0\rangle \\
D(\alpha) &= e^{-\frac{|\alpha|^2}{2}} e^{\alpha a^\dagger} e^{-\alpha^* a}
\end{aligned}$$

where $D(\alpha)$ is called the displacement operator.

$$|\alpha\rangle = D(\alpha) |0\rangle$$

1.1 Baker-Hausdorff identity

If

$$[[A, B], A] = [[A, B], B] = 0$$

then

$$e^{A+B} = e^{-[A,B]/2} e^A e^B$$

Let we have $A = \alpha a^\dagger$ and $B = -\alpha^* a$

$$\begin{aligned}
e^{\alpha a^\dagger - \alpha^* a} &= e^{-\frac{1}{2}[-\alpha a^\dagger \alpha^* a + \alpha^* a \alpha a^\dagger]} e^{\alpha a^\dagger} e^{-\alpha^* a} \\
&= e^{-\frac{1}{2}|\alpha|^2[-a^\dagger a + a a^\dagger]} e^{\alpha a^\dagger} e^{-\alpha^* a} \\
&= e^{-\frac{|\alpha|^2}{2}} e^{\alpha a^\dagger - \alpha^* a}
\end{aligned}$$

$$D(\alpha) = e^{\alpha a^\dagger - \alpha^* a}$$

another definition

$$\Rightarrow |\alpha\rangle = e^{\alpha a^\dagger - \alpha^* a} |0\rangle$$

The other equivalent antinormal form of $D(\alpha)$ is obtained by using $A = -\alpha^* a$ and $B = \alpha a^\dagger$, then we get

$$D(\alpha) = e^{\frac{|\alpha|^2}{2}} e^{-\alpha^* a} e^{\alpha a^\dagger}$$

or by using

$$\begin{aligned} e^{A+B} &= e^{\frac{1}{2}[A,B]} e^B e^A \\ D(\alpha) &= e^{\frac{|\alpha|^2}{2}} e^{-\alpha^* a} e^{\alpha a^\dagger} \end{aligned}$$

The operator $D(\alpha)$ is a unitary operator. i.e.

$$D^\dagger(\alpha) = D(-\alpha) = D^{-1}(\alpha)$$

It acts as a displacement operator upon the amplitudes a and a^\dagger i.e.

$$\begin{aligned} D^{-1}(\alpha) \hat{a} D(\alpha) &= a + \alpha \\ D^{-1}(\alpha) \hat{a}^\dagger D(\alpha) &= a^\dagger + \alpha^* \end{aligned}$$

This can be proved by writing

$$\begin{aligned} D(\alpha) &= e^{\alpha a^\dagger - \alpha^* a} \\ D^\dagger(\alpha) &= e^{\alpha^* a - \alpha a^\dagger} = D^{-1}(\alpha) \end{aligned}$$

using these equations we get

$$D^{-1}(\alpha) \hat{a} D(\alpha) = e^{\alpha^* a} e^{-\alpha a^\dagger} \hat{a} e^{\alpha a^\dagger} e^{-\alpha^* a}$$

For any operators A and B we have

$$e^{-\alpha A} B e^{\alpha A} = B - \alpha [A, B] + \frac{\alpha^2}{2!} [A, [A, B]] + \dots$$

For $A = \hat{a}^\dagger$ and $B = \hat{a}$

$$\begin{aligned} e^{-\alpha \hat{a}^\dagger} \hat{a} e^{\alpha \hat{a}^\dagger} &= a + \alpha \\ D^{-1}(\alpha) \hat{a} D(\alpha) &= a + \alpha \end{aligned}$$

Similarly for

$$D^{-1}(\alpha) \hat{a}^\dagger D(\alpha) = e^{\alpha^* a} \hat{a}^\dagger e^{-\alpha^* a}$$

here $A = \hat{a}$ and $B = \hat{a}^\dagger$

$$\implies D^{-1}(\alpha) \hat{a}^\dagger D(\alpha) = a^\dagger + \alpha^*$$

Prove

$$\begin{aligned} e^{-\alpha A} B e^{\alpha A} &= B - \alpha [A, B] + \frac{\alpha^2}{2!} [A, [A, B]] + \dots \\ \left(1 - \alpha A + \frac{\alpha^2 A^2}{2!} + \dots \right) B \left(1 + \alpha A + \frac{\alpha^2 A^2}{2!} + \dots \right) \\ &= B - \alpha (AB - BA) + \frac{\alpha^2}{2!} (\dots) \\ e^{-\alpha a^\dagger} a e^{\alpha a^\dagger} &= a - \alpha [a^\dagger, a] + \frac{\alpha^2}{2!} [a^\dagger, [a^\dagger, a]] + \dots \\ &= a^\dagger + \alpha \end{aligned}$$

1.2 Properties of coherent states

Properties of a cavity mode excited to a coherent state $|\alpha\rangle$ can be determined by the method applied to the number state $|n\rangle$.

1- The mean number of photon in the coherent state $|\alpha\rangle$ is given by

$$\begin{aligned}\langle n \rangle &= \langle \alpha | \hat{n} | \alpha \rangle = \langle \alpha | a^\dagger a | \alpha \rangle = |\alpha|^2 \\ | \alpha \rangle &= e^{-\frac{|\alpha|^2}{2}} \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} | n \rangle \\ \langle \alpha | &= e^{-\frac{|\alpha|^2}{2}} \sum_{m=0}^{\infty} \frac{(\alpha^*)^m}{\sqrt{m!}} \langle m | \end{aligned}$$

therefore

$$\begin{aligned}\langle n \rangle &= e^{-|\alpha|^2} \sum_{n=0}^{\infty} \sum_{m=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} \frac{(\alpha^*)^m}{\sqrt{m!}} \langle m | a^\dagger a | n \rangle \\ &= e^{-|\alpha|^2} \sum_n \frac{(\alpha^* \alpha)^n}{n!} n \\ &= e^{-|\alpha|^2} \sum_n \frac{(|\alpha|^2)^n}{n!} n \end{aligned}$$

Let $x = |\alpha|^2$, and also

$$\begin{aligned}x \frac{\partial}{\partial x} \sum_n \frac{x^n}{n!} &= \sum_n \frac{x^n}{n!} n \\ x \frac{\partial}{\partial x} e^x &= \sum_n \frac{x^n}{n!} n \end{aligned}$$

therefore we can write

$$\begin{aligned}\langle n \rangle &= e^{-|\alpha|^2} |\alpha|^2 \frac{\partial}{\partial |\alpha|^2} e^{|\alpha|^2} \\ &= |\alpha|^2 \end{aligned}$$

Find

$$\begin{aligned}\langle \alpha | \hat{n}^2 | \alpha \rangle &= e^{-|\alpha|^2} \sum_n \frac{(\alpha^* \alpha)^n}{n!} n^2 \\ &= e^{-|\alpha|^2} \sum_n \frac{|\alpha|^{2n}}{n!} \{n(n-1) + n\} \end{aligned}$$

Let again $x = |\alpha|^2$, by the definition

$$\begin{aligned}\sum_n \frac{x^n}{n!} n(n-1) &= x^2 \frac{\partial^2}{\partial x^2} \sum_n \frac{x^n}{n!} \\ &= x^2 \frac{\partial^2}{\partial x^2} e^x \end{aligned}$$

so we can write

$$\langle \alpha | \hat{n}^2 | \alpha \rangle = e^{-|\alpha|^2} e^{|\alpha|^2} \left(|\alpha|^4 + |\alpha|^2 \right)$$

Root-mean-square deviation is

$$\Delta n = \sqrt{\langle n^2 \rangle - \langle n \rangle^2} = \sqrt{|\alpha|^2}$$

Where $|\alpha|^2$ is the mean number of photons in the cavity mode and uncertainty spread about the mean is equal to the square root of the mean number of photons.

ii)- Photon statistics: photon distribution function:

The probability of finding n-photons in the field $|\alpha\rangle$ is

$$p(n) = |\langle n | \alpha \rangle|^2$$

where

$$\langle n | \alpha \rangle = e^{-\frac{|\alpha|^2}{2}} \sum_{m=0}^{\infty} \frac{(\alpha)^m}{\sqrt{m!}} \langle n | m \rangle$$

$$\begin{aligned} p(n) &= \frac{e^{-|\alpha|^2} |\alpha|^{2n}}{n!} \\ &= \frac{e^{-\langle n \rangle} \langle n \rangle^n}{n!} \end{aligned}$$

is a poisson distribution.

iii)- Coherent state is the minimum energy state: i.e

$$\Delta p \Delta q = \frac{\hbar}{2}$$

\hat{a} and \hat{a}^\dagger are not hermitian but their combinations are.

Let $m_j = 1$ at $t = 0$ $r = 0$

$$\begin{aligned} \hat{a} &= \frac{1}{\sqrt{2\hbar\omega}} (\omega \hat{q} + i\hat{p}) \\ \hat{a}^\dagger &= \frac{1}{\sqrt{2\hbar\omega}} (\omega \hat{q} - i\hat{p}) \end{aligned}$$

adding these two we get

$$\begin{aligned} \frac{1}{2} (\hat{a} + \hat{a}^\dagger) &= \sqrt{\frac{\omega}{2\hbar}} \hat{q} \\ \hat{a} + \hat{a}^\dagger &= \sqrt{\frac{2\omega}{\hbar}} \hat{q} \end{aligned}$$

and

$$\begin{aligned}\hat{a} - \hat{a}^\dagger &= i\sqrt{\frac{2}{\hbar\omega}}\hat{p} \\ \frac{1}{2i}(\hat{a} - \hat{a}^\dagger) &= \sqrt{\frac{1}{2\hbar\omega}}\hat{p}\end{aligned}$$

\hat{p} and \hat{q} are hermitian and represent observable quantities,

$$\begin{aligned}[\hat{q}, \hat{p}] &= i\hbar \\ \Delta\hat{q}\Delta\hat{p} &\geq \frac{\hbar}{2}\end{aligned}$$

for a coherent state we have to prove that

$$\Delta p \Delta q = \frac{\hbar}{2}$$

$$\begin{aligned}\langle \hat{p} \rangle &= \langle \alpha | \hat{p} | \alpha \rangle = \frac{\sqrt{2\hbar\omega}}{2i} (\langle a \rangle - \langle a^\dagger \rangle) \\ \langle a \rangle &= \langle \alpha | a | \alpha \rangle = \alpha \langle \alpha | \alpha \rangle\end{aligned}$$

$$\begin{aligned}\langle p \rangle &= \frac{\sqrt{2\hbar\omega}}{2i} (\alpha - \alpha^*) \\ \langle p^2 \rangle &= \left(\frac{\sqrt{2\hbar\omega}}{2i} \right)^2 \langle (a - a^\dagger)(a - a^\dagger) \rangle \\ &= -\frac{2\hbar\omega}{4} \langle \alpha | a^2 + a^{\dagger 2} - aa^\dagger - a^\dagger a | \alpha \rangle \\ &= -\frac{2\hbar\omega}{4} (\alpha^2 + \alpha^{*2} - 2\alpha^* \alpha - 1)\end{aligned}$$

$$\begin{aligned}\Delta p^2 &= \langle p^2 \rangle - \langle p \rangle^2 \\ &= -\frac{\hbar\omega}{2} (\alpha^2 + \alpha^{*2} - 2\alpha^* \alpha - 1) + \frac{\hbar\omega}{2} (\alpha^2 + \alpha^{*2} - 2\alpha^* \alpha) \\ &= \frac{\hbar\omega}{2}\end{aligned}$$

Similarly

$$\begin{aligned}\Delta q^2 &= \frac{\hbar}{2\omega} \\ \Delta q \Delta p &= \frac{\hbar}{2}\end{aligned}$$

iv)- Coherent states are not orthogonal, but are normalized. i.e.

$$\langle \alpha | \alpha \rangle = e^{-|\alpha|^2} \sum_n \frac{(\alpha^* \alpha)^n}{n!} = 1$$

For two different complex numbers α and β .

$$\begin{aligned} | \alpha \rangle &= e^{-\frac{|\alpha|^2}{2}} \sum_n \frac{\alpha^n}{\sqrt{n!}} | n \rangle \\ | \beta \rangle &= e^{-\frac{|\beta|^2}{2}} \sum_m \frac{\beta^m}{\sqrt{m!}} | m \rangle \end{aligned}$$

$$\begin{aligned} \langle \alpha | \beta \rangle &= e^{-\frac{|\alpha|^2}{2}} e^{-\frac{|\beta|^2}{2}} e^{\alpha \times \beta} \neq 0 \\ \Rightarrow |\langle \alpha | \beta \rangle|^2 &= e^{-|\alpha - \beta|^2}. \end{aligned}$$

The $| \alpha \rangle$ form an over complete set of states and lack of orthogonality is a consequence of this. i, j, k are orthogonal and independent of each other. If we divide space in 5 directions they would not be orthogonal and independent of each other. Therefore over complete. i.e. there are many more coherent states $| \alpha \rangle$ than there are states $| n \rangle$.

Completeness relation:

For number states

$$\sum_n | n \rangle \langle n | = 1$$

Similarly the set of all coherent states $| \alpha \rangle$ is a complete set and satisfy the completeness relation.

$$\frac{1}{\pi} \int d^2 \alpha | \alpha \rangle \langle \alpha | = 1$$

Let

$$\begin{aligned} \alpha &= r e^{i\theta} \\ d^2 \alpha &= r dr d\theta \end{aligned}$$

$$\begin{aligned} \int d^2 \alpha | \alpha \rangle \langle \alpha | &= \int e^{-\frac{|\alpha|^2}{2}} \sum_m \frac{\alpha^m}{\sqrt{m!}} | m \rangle e^{-\frac{|\alpha|^2}{2}} \sum_n \frac{\alpha^{*n}}{\sqrt{n!}} \langle n | d^2 \alpha \\ &= \int_0^\infty \int_0^{2\pi} e^{-|\alpha|^2} \sum_{n,m} \frac{|\alpha|^{n+m+1}}{\sqrt{n!m!}} e^{i(m-n)\theta} | m \rangle \langle n | d|\alpha| d\theta \end{aligned}$$

as

$$\int_0^{2\pi} d\theta e^{i(m-n)\theta} = 2\pi \delta_{nm}$$

therefore

$$\begin{aligned} \int d^2\alpha \quad | \quad \alpha \rangle \langle \alpha | &= 2\pi \int_0^\infty \sum_n \frac{1}{n!} |\alpha|^{2n+1} e^{-|\alpha|^2} |n\rangle \langle n| d(|\alpha|) \\ &= \pi \sum_n \frac{1}{n!} \int_0^\infty 2|\alpha| d(|\alpha|) |\alpha|^{2n} e^{-|\alpha|^2} |n\rangle \langle n| \end{aligned}$$

putting

$$\begin{aligned} x &= |\alpha|^2 \\ dx &= 2|\alpha| d|\alpha| \end{aligned}$$

$$\int d^2\alpha \quad | \alpha \rangle \langle \alpha | = \pi \sum_n \frac{1}{n!} \int_0^\infty dx x^n e^{-x} |n\rangle \langle n|$$

$$\begin{aligned} \int_0^\infty dx x^n e^{-x} &= n! \\ \int d^2\alpha \quad | \quad \alpha \rangle \langle \alpha | &= \pi \sum_n \frac{1}{n!} n! |n\rangle \langle n| \\ &= \pi \sum_n |n\rangle \langle n| \\ \frac{1}{\pi} \int d^2\alpha \quad | \quad \alpha \rangle \langle \alpha | &= 1 \end{aligned}$$

The completeness property is essential for the utility of a set of states. $|\alpha\rangle$ is a complete but not orthogonal. As a result any coherent state can be expanded in terms of other states.

$$\begin{aligned} | \quad \alpha \rangle &= \frac{1}{\pi} \int d^2\alpha' | \alpha' \rangle \langle \alpha' | \alpha \rangle \\ &= \frac{1}{\pi} \int d^2\alpha' | \alpha' \rangle \exp \left[-\frac{1}{2} |\alpha|^2 + \alpha' \alpha^* - \frac{1}{2} |\alpha'|^2 \right] \end{aligned}$$

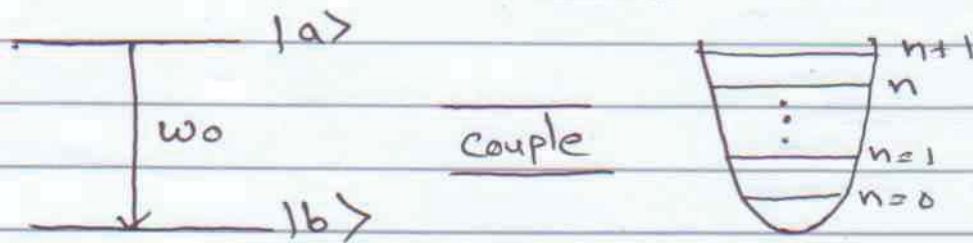
This shows that the coherent states are overcomplete.

ATOM-FIELD INTERACTION Quantum Theory.

⇒ ATOM → Quantum Mechanically

⇒ Field → Quantum Mechanically.

Interaction between a single-mode radiation field and a two-level atom inside a cavity



Atomic state-vector

$$|\psi_{\text{atom}}\rangle = c_a |a\rangle + c_b |b\rangle$$

Field state-vector

$$|\psi_{\text{field}}\rangle = \sum_n c_n |n\rangle$$

Atom-field state-vector

$$|\psi_{\text{a-f}}\rangle = \sum_n [c_{a,n} |a\rangle |n\rangle + c_{b,n} |b\rangle |n\rangle]$$

$c_{a,n}$ → probability amplitude

If at time $t=0$

$$|\psi(0)_{a-f}\rangle = |a\rangle |n\rangle$$

Then at later time t .

$$|\psi(t)_{a-f}\rangle = C_{a,n} |a\rangle |n\rangle + C_{b,n+1} |b\rangle |n+1\rangle$$

\Rightarrow a finite probability that atom has made a transition to the lower level and emitted a photon.

Total Hamiltonian of the system

$$H = H_A + H_F + H_I$$

H_A - Energy of free-atom

$$= \sum_i E_i |i\rangle \langle i|$$

H_F - Energy of free-field

$$= \sum_k \hbar \omega_k \left(a_k^\dagger a_k + \frac{1}{2} \right)$$

H_I - Interaction energy under dipole-approximation

$$H_I = -e \vec{r} \cdot \vec{E}$$

Define - atom transition operators

$$\sigma_{ij} = |i\rangle\langle j| \quad ij - \text{atomic level}$$

\Rightarrow

$$H_A = \sum_i E_i \sigma_{ii}$$

and

$$\begin{aligned} e\vec{r} &= \sum_i \sum_j |i\rangle\langle i| e\vec{r} |j\rangle\langle j| \\ &= \sum_{i,j} p_{ij} \sigma_{ij} \end{aligned}$$

with

$$p_{ij} = e\langle i|\vec{r}|j\rangle - \text{electric-dipole transition matrix element.}$$

Electric field operator

$$E = \sum_k \hat{e}_k \mathcal{E}_k (\hat{a}_k + \hat{a}_k^\dagger)$$

with

$$\mathcal{E}_k = (\hbar\omega_k / 2\epsilon_0 V)$$

\hat{e}_k — represents polarization

⇒ Complete Hamiltonian

$$H = \sum_k \hbar \omega_k \left(\hat{a}_k^\dagger \hat{a}_k + \frac{1}{2} \right) + \sum_i E_i \sigma_{ii} + \hbar \sum_{ij} \sum_k g_k^{ij} \sigma_{ij} (\hat{a}_k + \hat{a}_k^\dagger)$$

Here

$$g_k^{ij} = \frac{-\mathbf{p}_{ij} \cdot \hat{\mathbf{e}}_k \mathcal{E}_k}{\hbar}$$

- coupling constant.
- dimensions of frequency
- similar to Rabi-frequency

For a two-level atom and single-mode field

$$k=1$$

$$\left. \begin{array}{l} i = a, b \\ j = a, b \end{array} \right\} \text{ for two-level atom.}$$

$$\Rightarrow H = \hbar \omega \left(a^\dagger a + \frac{1}{2} \right) + E_a \sigma_{aa} + E_b \sigma_{bb} + \hbar (g^{aa} \sigma_{aa} + g^{ab} \sigma_{ab} + g^{ba} \sigma_{ba} + g^{bb} \sigma_{bb}) (a + a^\dagger)$$

$$g^{aa} = g^{bb} = 0$$

$$\because p_{aa} = \langle a | e r | a \rangle = 0$$

dipole-transitions

and

$$g^{ab} = g^{ba} = g \quad g \text{ is real.}$$

As $\sigma_{aa} + \sigma_{bb} = |a\rangle\langle a| + |b\rangle\langle b| = 1$

Hamiltonian reduces to

$$H = E_a \sigma_{aa} + E_b \sigma_{bb} + \hbar\omega \left(a^\dagger a + \frac{1}{2}\right) + \hbar g (\sigma_{ab} + \sigma_{ba}) (a + a^\dagger)$$

First term

$$E_a \sigma_{aa} + E_b \sigma_{bb} = \frac{1}{2} (E_a - E_b) (\sigma_{aa} - \sigma_{bb}) + \frac{1}{2} (E_a + E_b) (\sigma_{aa} + \sigma_{bb})$$

As $E_a - E_b = \hbar\omega_0$

$$\Rightarrow E_a \sigma_{aa} + E_b \sigma_{bb} = \frac{1}{2} \hbar\omega_0 (\sigma_{aa} - \sigma_{bb}) + \frac{1}{2} (E_a + E_b)$$

The total Hamiltonian

$$H = \hbar\omega \left(a^\dagger a + \frac{1}{2}\right) + \frac{1}{2} \hbar\omega_0 (\sigma_{aa} - \sigma_{bb}) + \frac{1}{2} (E_a + E_b) + \hbar g (\sigma_{ab} + \sigma_{ba}) (a + a^\dagger)$$

$$\sigma_{aa} - \sigma_{bb} = \sigma_z$$

$$\sigma_{ab} = |a\rangle\langle b| = \sigma_+$$

$$\sigma_{ba} = |b\rangle\langle a| = \sigma_-$$

Ignoring

$$\frac{1}{2}(E_a + E_b) \approx \frac{1}{2}\hbar\omega \quad \text{constant energy terms}$$

$$H = \hbar\omega a^\dagger a + \frac{1}{2}\hbar\omega\sigma_z + \hbar g(\sigma_+ + \sigma_-)(a + a^\dagger)$$

Here

$$\sigma_- = \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix} \quad \text{lowering operator}$$

$$\sigma_+ = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} \quad \text{raising operator}$$

$$\sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

$$\Rightarrow H = \hbar\omega a^\dagger a + \frac{1}{2}\hbar\omega\sigma_z + \hbar g(\sigma_+ a^\dagger + \sigma_+ a + \sigma_- a^\dagger + \sigma_- a)$$

$\left. \begin{array}{l} \sigma_+ a^\dagger \\ \sigma_- a \end{array} \right\}$ energy non-conserving terms
Dropping under RWA

$$\Rightarrow H = \hbar\omega a^\dagger a + \frac{1}{2}\hbar\omega\sigma_z + \hbar g(\sigma_+ a + a^\dagger \sigma_-)$$

$$H_0 = \hbar\omega a^\dagger a + \frac{1}{2}\hbar\omega\sigma_z$$

$$H_I = \hbar g(\sigma_+ a + a^\dagger \sigma_-)$$

For multimode field

$$H = \sum_k \hbar \omega_k a_k^\dagger a_k + \frac{1}{2} \hbar \omega_0 \sigma_z + \hbar \sum_k g_k (\sigma_+ a_k + a_k^\dagger \sigma_-)$$

Hamiltonian

$$H = \hbar \omega a^\dagger a + \frac{1}{2} \hbar \omega_0 \sigma_z + \hbar g (\sigma_+ a + a^\dagger \sigma_-)$$

Describes the atom-field interaction under dipole and rotating-wave approximation.

This Hamiltonian is exactly solvable is called Jaynes-Cummings Model (JCM)

INTERACTION PICTURE

operator in Interaction Picture

$$\hat{O}_I = e^{iH_0 t/\hbar} \hat{O}(0) e^{-iH_0 t/\hbar}$$

Hamiltonian in Interaction picture

$$V = e^{iH_0 t/\hbar} H_I e^{-iH_0 t/\hbar}$$

with

$$H_0 = \frac{1}{2} \hbar \omega_0 \sigma_z + \hbar \omega a^\dagger a$$

$$\Rightarrow V = \hbar g \left\{ e^{i(\omega a^\dagger a + \frac{\omega_0 \sigma_z}{2})t} (\sigma_+ \hat{a} + a^\dagger \sigma_-) e^{-i(\omega a^\dagger a + \frac{\omega_0 \sigma_z}{2})t} \right\}$$

Atomic and field operators commute

$$\Rightarrow V = \hbar g \left\{ \left(e^{i\frac{\omega_0 \sigma_z}{2}t} \sigma_+ e^{-i\frac{\omega_0 \sigma_z}{2}t} \right) \left(e^{i\omega a^\dagger a t} \hat{a} e^{-i\omega a^\dagger a t} \right) + \left(e^{i\omega a^\dagger a t} a^\dagger e^{-i\omega a^\dagger a t} \right) \left(e^{i\frac{\omega_0 \sigma_z}{2}t} \sigma_- e^{-i\frac{\omega_0 \sigma_z}{2}t} \right) \right\}$$

Using

$$e^{\alpha A} B e^{-\alpha A} = B + \alpha [A, B] + \frac{\alpha^2}{2!} [A, [A, B]] + \dots$$

First consider

$$e^{i\omega a^\dagger t} \hat{a} e^{-i\omega a^\dagger t} = ?$$

here $\alpha = i\omega t$; $A = \hat{a}^\dagger \hat{a}$ and $B = \hat{a}$

$$e^{i\omega a^\dagger t} \hat{a} e^{-i\omega a^\dagger t} = \hat{a} + i\omega t [\hat{a}^\dagger \hat{a}, \hat{a}] + \frac{(i\omega t)^2}{2!} [\hat{a}^\dagger \hat{a}, [\hat{a}^\dagger \hat{a}, \hat{a}]] + \dots$$

$$\text{As } [\hat{a}, \hat{a}^\dagger] = 1 \text{ \& } [\hat{a}^\dagger, \hat{a}] = -1$$

$$\begin{aligned} [\hat{a}^\dagger \hat{a}, \hat{a}] &= \hat{a}^\dagger \hat{a} \hat{a} - \hat{a} \hat{a}^\dagger \hat{a} = [\hat{a}^\dagger \hat{a} - \hat{a} \hat{a}^\dagger] \hat{a} \\ &= [\hat{a}^\dagger, \hat{a}] \hat{a} = -\hat{a} \end{aligned}$$

And

$$[\hat{a}^\dagger \hat{a}, [\hat{a}^\dagger \hat{a}, \hat{a}]] = [\hat{a}^\dagger \hat{a}, -\hat{a}] = \hat{a}$$

Putting in above eqn we get.

$$\begin{aligned} e^{i\omega a^\dagger t} \hat{a} e^{-i\omega a^\dagger t} &= \hat{a} - i\omega t \hat{a} + \frac{(i\omega t)^2}{2!} \hat{a} - \frac{(i\omega t)^3}{3!} \hat{a} + \dots \\ &= \hat{a} \left[1 - i\omega t + \frac{(i\omega t)^2}{2!} - \frac{(i\omega t)^3}{3!} + \dots \right] \\ &= \hat{a} e^{-i\omega t} = \hat{a}_I \quad \text{— destruction in I.P} \end{aligned}$$

Similarly,

$$e^{i\omega_0 t} \hat{a} e^{-i\omega_0 t} = \hat{a} e^{i\omega t} = \hat{a}_I^+$$

$$e^{\frac{i\omega_0 \sigma_z}{2}} \sigma_+ e^{-\frac{i\omega_0 \sigma_z}{2}} = \sigma_+ e^{i\omega t} = \sigma_{+I}$$

$$\text{And } e^{\frac{i\omega_0 \sigma_z}{2}} \sigma_- e^{-\frac{i\omega_0 \sigma_z}{2}} = \sigma_- e^{-i\omega t} = \sigma_{-I}$$

Putting in V we get.

$$\begin{aligned} V &= \hbar g \left[\sigma_+ \hat{a} e^{i(\omega_0 - \omega)t} + \hat{a}^\dagger \sigma_- e^{-i(\omega_0 - \omega)t} \right] \\ &= \hbar g (\sigma_+ \hat{a} e^{i\Delta t} + \hat{a}^\dagger \sigma_- e^{-i\Delta t}) \end{aligned}$$

where $\Delta = \omega_0 - \omega$ — detuning.

The non-conservative terms

$$\left. \begin{aligned} \sigma_+ \hat{a}^\dagger &\sim e^{i(\omega_0 + \omega)t} \\ \sigma_- \hat{a} &\sim e^{-i(\omega_0 + \omega)t} \end{aligned} \right\} \begin{array}{l} \text{Rapidly oscillating} \\ \text{neglected in R.W.A.} \end{array}$$

\Rightarrow In I. P

$$\hat{a}_I^+ = \hat{a}^\dagger e^{i\omega t} ; a_I = a e^{-i\omega t}$$

and

$$\sigma_{+I} = \sigma_+ e^{i\omega t} ; \sigma_{-I} = \sigma_- e^{-i\omega t}.$$

At exact resonance

$$\Delta = \omega_0 - \omega = 0 \Rightarrow \omega = \omega_0$$

$$\Rightarrow V = \hbar g (a \sigma_+ + a^\dagger \sigma_-)$$

Is the interaction part of Hamiltonian in Interaction picture in RWA approximation and at exact resonance.

Equation of motion in Interaction picture is written as,

$$i \hbar \dot{|\psi_I\rangle} = -\frac{i}{\hbar} V |\psi\rangle$$

The state-vector in I.P

$$|\psi(t)\rangle_I = \sum_n [C_{an}(t) |a, n\rangle + C_{bn}(t) |b, n\rangle]$$

C_{an} & C_{bn} — slowly varying probability amplitudes.

The interaction energy can only cause transitions b/w $|a, n\rangle$ & $|b, n+1\rangle$

\Rightarrow

$$|\Psi(t)\rangle = C_a |a, n\rangle + C_{b, n+1} |b, n+1\rangle$$

Putting in eqn. of motion and multiplying the result by $\langle a, n|$

$$\begin{aligned}\dot{C}_a &= -ig \langle a, n | (a\sigma_+ + a^\dagger\sigma_-) | a, n \rangle C_a \\ &\quad -ig \langle a, n | a\sigma_+ + a^\dagger\sigma_- | b, n+1 \rangle C_{b, n+1}\end{aligned}$$

using

$$a|n\rangle = \sqrt{n}|n-1\rangle \quad ; \quad a^\dagger|n\rangle = \sqrt{n+1}|n+1\rangle$$

$$\sigma_+|b\rangle = |a\rangle \quad ; \quad \sigma_-|a\rangle = |b\rangle$$

$$\text{and } \langle a, n | a, n \rangle = 1$$

$$\Rightarrow \dot{C}_a(t) = -ig\sqrt{n+1} C_{b, n+1}(t)$$

$$\dot{C}_{b, n+1}(t) = -ig\sqrt{n+1} C_a(t)$$

Coupled differential equations.

For atom initially in level $|a\rangle$

$$C_{an}(0) = C_n(0) \quad \text{where } C_a(0) = 1$$

$$C_{bn}(0) = 0 \quad \text{" } C_b(0) = 0$$

$$\Rightarrow C_{an}(t) = C_{an}(0) \cos(g\sqrt{n+1}t) - C_n(0) \cos g\sqrt{n+1}t$$

and

$$C_{b,n+1}(t) = -i C_{an}(0) \sin(g\sqrt{n+1}t) = -i C_n(0) \sin(g\sqrt{n+1}t)$$

These are the conditions for emission

For stimulated absorption

$$|\Psi_I(0)\rangle = |b\rangle |n+1\rangle$$

Solutions are

$$C_{a,n}(t) = -i \sin(g\sqrt{n+1}t) C_{b,n+1}(0)$$

$$C_{b,n+1}(t) = \cos(g\sqrt{n+1}t) C$$

The factor $g\sqrt{n+1}$ is the Rabi
flopping frequency.

Population Inversion.

Probability of finding atom in state $|a\rangle$

- taking trace over field variables

$$\begin{aligned} P(a) &= \text{Tr}_f |C_{an}|^2 = \text{Tr}_f \sum_n |n\rangle \langle n| |C_{an}|^2 \\ &= \sum_n \langle n|n\rangle |C_{an}|^2 = \sum_n |C_{an}|^2 \end{aligned}$$

Probability of finding the atom in state $|b\rangle$

$$P(b) = \sum_n |C_{bn}|^2$$

\Rightarrow Population Inversion

$$W = P(a) - P(b) = \sum_n [|C_{an}|^2 - |C_{bn}|^2]$$

Probability of finding n -photons in the field at time t is

$$\begin{aligned} P(n, t) &= \text{Tr}_\alpha |C_{\alpha, n}(t)|^2 = \sum_{\alpha=a, b} |C_{\alpha, n}(t)|^2 \\ &= |C_{an}(t)|^2 + |C_{bn}(t)|^2 \end{aligned}$$

$$\Rightarrow P(n,t) = P_{nn}(t)$$

$$\Rightarrow P_{nn}(0) = |C_{an}(0)|^2 + |C_{bn}(0)|^2$$

For initial condition

$$C_{an}(0) = C_n(0) \text{ \& } C_{b,n+1}(0) = 0$$

$$\Rightarrow P_{nn}(0) = |C_{an}(0)|^2 = |C_n(0)|^2 \because C_n(0) = 1$$

It gives the probability that there are n -photons in the field at time $t=0$

$$\begin{aligned} \Rightarrow P(n,t) &= |C_{an}(t)|^2 \cos^2(g\sqrt{n+1}t) + |C_{a,n-1}(t)|^2 \sin^2 g\sqrt{n}t \\ &= P_{nn}(0) \cos^2(g\sqrt{n+1}t) + P_{n-1,n-1}(0) \sin^2 g\sqrt{n}t \end{aligned}$$

Using $P_{nn}(0) = |C_{an}(0)|^2$ Population inversion can be written as.

$$\begin{aligned} W &= P(a) - P(b) = \sum_n [|C_{an}|^2 - |C_{bn}|^2] \\ &= \sum_n [P_{nn}(0) \cos^2 g\sqrt{n+1}t - P_{n-1,n-1}(0) \sin^2 g\sqrt{n}t] \end{aligned}$$

We need $P_{nn}(0) = ?$ for W

In semiclassical theory probability amplitudes are

$$C_a(t) = \cos\left(\frac{\Omega_R t}{2}\right)$$

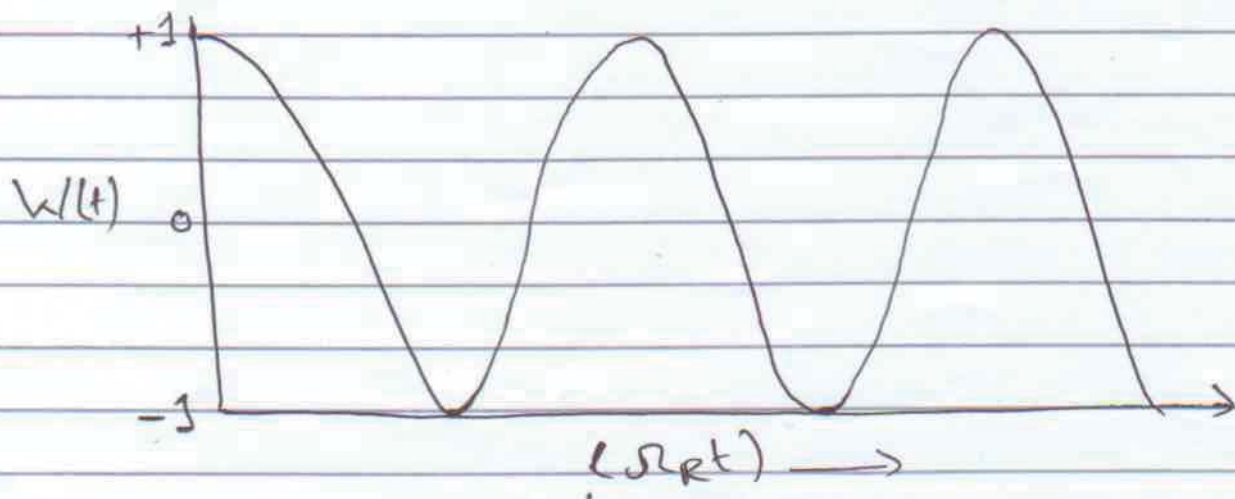
$$C_b(t) = i \sin\left(\frac{\Omega_R t}{2}\right)$$

$$\Omega_R = \frac{p \cdot E}{\hbar} = \text{Rabi-freq}$$

\Rightarrow

Population inversion in semiclassical theory

$$\begin{aligned} W(t) &= P_a(t) - P_b(t) = \cos^2\left(\frac{\Omega_R t}{2}\right) - \sin^2\left(\frac{\Omega_R t}{2}\right) \\ &= \cos(\Omega_R t) \end{aligned}$$



Population inversion oscillates b/w -1 & +1 at
at freq. Ω_R . Atom undergoes a Rabi flopping
b/w the upper and lower level under the action of
field.

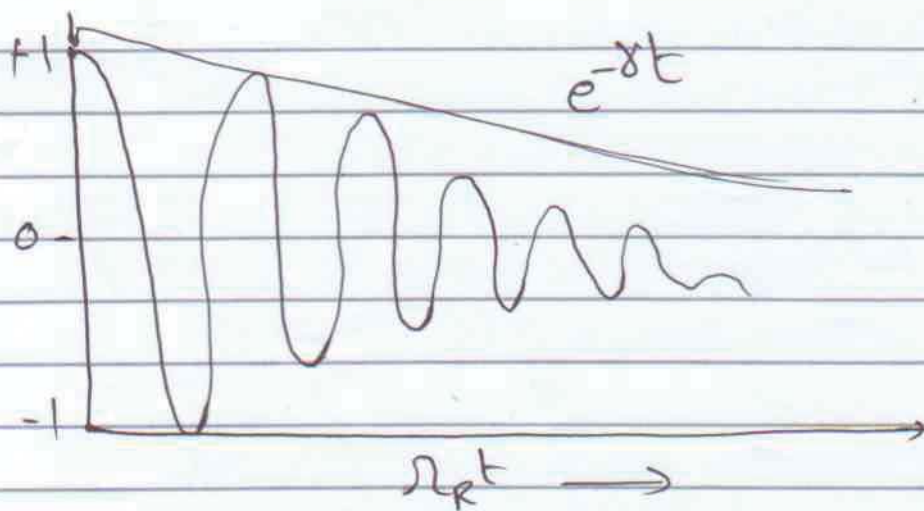
If atomic is included the probability amplitudes are written as,

$$C_a(t) = e^{-\gamma/2 t} \cos\left(\frac{\Omega_R t}{2}\right)$$

$$C_b(t) = i e^{-\gamma/2 t} \sin\left(\frac{\Omega_R t}{2}\right)$$

⇒ Population inversion at time t is

$$W(t) = e^{-\gamma t} \cos(\Omega_R t)$$



Rabi-oscillations are damped due to atomic decay.

In quantum theory of atom-field interaction

For atom initially in excited state we have

$$W(t) = \sum_n \left[P_{nn}(0) \cos^2 g \sqrt{n+1} t - P_{n-1, n-1}(0) \sin^2 g \sqrt{n+1} t \right]$$

↓ shifting by one

$P_{nn}(0) = |c_n(0)|^2$ - probability that there are n -photons at time $t=0$

Field can be

- i, Vacuum
- ii, Fock state
- iii, Coherent state.

i, For a vacuum state

$$\rho = |0\rangle\langle 0|$$

$$\Rightarrow P_{nn}(0) = \langle n|0\rangle\langle 0|n\rangle = \delta_{n0} = 1$$

for $n=0$ only

As

$$\begin{aligned} W(t) &= \sum_{n=0}^{\infty} \left[P_{nn}(0) \cos^2 g \sqrt{n+1} t - P_{nn}(0) \sin^2 g \sqrt{n+1} t \right] \\ &= \sum_{n=0}^{\infty} P_{nn}(0) \left[\cos^2 g \sqrt{n+1} t \right] \end{aligned}$$

$$\Rightarrow W(t) = \sum_{n=0}^{\infty} \delta_{n0} \cos(2gt\sqrt{n+1})$$

$$= \cos 2gt\sqrt{1} \quad \text{for } n=0$$

\Rightarrow The Rabi-oscillations take place even when there is no field.

\Rightarrow The square root of 1 i.e., $\sqrt{1}$ corresponds to spontaneous emission.

This is drastically different from the predictions of semi-classical theory.

\Rightarrow In Semi-classical theory the probability of finding the atom in state $|b\rangle$ is

$$P_b(t) = \sin^2(\Omega_R t) \quad \text{where } \Omega_R = \frac{\mathbf{p} \cdot \mathbf{E}}{\hbar}$$

If there is no field.

$$P_b(t) = \sin^2(\Omega_R t) = 0$$

\Rightarrow No transition in the absence of field.

In quantum theory, the probability of finding atom in state $|b\rangle$ is

$$P_b(t) = \sum_{n=0}^{\infty} |C_{bn}(t)|^2$$
$$= \sum_{n=0}^{\infty} S_{nn}(0) \sin^2(g\sqrt{n+1}t)$$

For ~~vacuum~~ vacuum = no-field

$$S_{nn}(0) = \delta_{n0} \quad \forall n \geq 0$$

$$\Rightarrow P_b(t) = \sum_{n=0}^{\infty} \delta_{n0} \sin^2(g\sqrt{n+1}t) = \sin^2 gt$$

vacuum Rabi-freq.

→ In semi-classical theory - atom in excited state cannot make a transition to lower-level in the absence of field.

In quantum treatment transition from $|a\rangle \rightarrow |b\rangle$ in vacuum becomes possible due to spontaneous emission.

For field initially in number state

$$\rho = |n_0\rangle\langle n_0|$$

$$\Rightarrow \rho_{nn|0} = \langle n|n_0\rangle\langle n_0|n\rangle = \delta_{nn_0}$$

$$\Rightarrow W(t) = \sum_{n=0}^{\infty} \delta_{nn_0} \cos(2g\sqrt{n+1}t)$$

$$= \cos 2gt\sqrt{n_0+1}$$

for $n_0 \gg 1$

$$W(t) \approx \cos(2gt\sqrt{n_0})$$

\Rightarrow This is like semi-classical result

for $n_0 \gg 1$

$$W(t) \approx \cos(2gt\sqrt{n_0})$$

here for $n_0 = 0$

$W(t) = 0$ like classical treatment.

For the field to be initially in the coherent state.

$$P_{nn}(0) = \frac{|\alpha|^{2n} e^{-|\alpha|^2}}{n!}$$

\Rightarrow

$$W(t) = \sum_{n=0}^{\infty} \left[\frac{e^{-|\alpha|^2} |\alpha|^{2n}}{n!} \cos^2 g\sqrt{n+1}t - \frac{e^{-|\alpha|^2} |\alpha|^{2n-1}}{(n-1)!} \sin^2 g\sqrt{n}t \right]$$

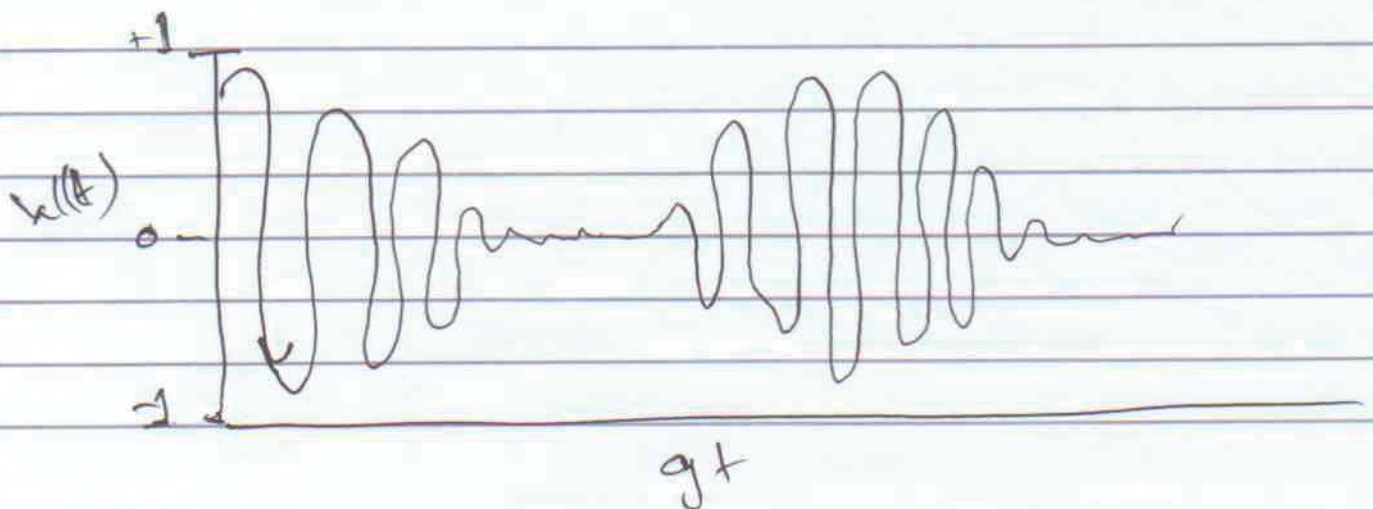
here $|\alpha|^2 = \bar{n}$

$$W(t) = \sum_{n=0}^{\infty} \left(\frac{e^{-\bar{n}} \bar{n}^n}{n!} \cos^2 g\sqrt{n+1}t - \frac{e^{-\bar{n}} \bar{n}^{n-1}}{(n-1)!} \sin^2 g\sqrt{n}t \right)$$

\hookrightarrow shifting by one

$$= \sum_{n=0}^{\infty} \frac{e^{-\langle n \rangle} \langle n \rangle^n}{n!} \left[\cos^2 g\sqrt{n+1}t - \sin^2 g\sqrt{n+1}t \right]$$

$$= \sum_{n=0}^{\infty} P_{nn}(0) \left[\cos^2 g\sqrt{n+1}t \right]$$



The phenomena of collapse and revival can be understood from

$$W(t) = \sum_{n=0}^{\infty} P_{nn}(0) \left[\cos 2g\sqrt{n+1}t \right]$$

\Rightarrow Each term in the summation represents Rabi-oscillations, for a definite value of n . At time $t=0$ all terms are correlated. As time increases the Rabi-oscillations associated with different excitations have different frequencies and therefore become un-correlated leading to a collapse of inversion. As time further increases the correlation is restored and revival occurs.

\Rightarrow Revival is pure Q.M phenomenon and occurs due to the discrete values of n .

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The number state behaves like semi-classical state because both have definite Intensity, needed to avoid the interference leading to a collapse. The random phase associated with number state (but not with the classical field) is not important for Rabi-flopping since the atom and field maintain a precise relative phase in the absence of decay processes.

While ~~coher~~ minimum uncertainty in ~~the~~ intensity in coherent state, causes the atom-field relative phase to diffuse away i.e any spread in field strength will dephase Rabi-oscillations.