

4.3 Measurement and expectation values

Slides: Video 4.3.1 Quantum-mechanical measurement

Text reference: Quantum Mechanics for Scientists and Engineers

Section 3.8





Measurement and expectation values



Quantum-mechanical measurement

Quantum mechanics for scientists and engineers

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Probabilities and expansion coefficients

Suppose we take some (normalized) quantum mechanical wave function $\Psi(\mathbf{r}, t)$

and expand it in some complete orthonormal set of spatial functions $\psi_n(\mathbf{r})$

At least if we allow the expansion coefficients c_n to vary in time we know we can always do this

$$\Psi(\mathbf{r}, t) = \sum_n c_n(t) \psi_n(\mathbf{r})$$

Probabilities and expansion coefficients

Then the fact that $\Psi(\mathbf{r}, t)$ is normalized means

we know the answer for the normalization integral

$$\int_{-\infty}^{\infty} |\Psi(\mathbf{r}, t)|^2 d^3\mathbf{r} = \int_{-\infty}^{\infty} \left[\sum_n c_n^*(t) \psi_n^*(\mathbf{r}) \right] \times \left[\sum_m c_m(t) \psi_m(\mathbf{r}) \right] d^3\mathbf{r} = 1$$

Because of the orthogonality of the basis functions

only terms with $n = m$ survive the integration

Because of the orthonormality of the basis functions

the result from any such term will simply be $|c_n(t)|^2$

Hence we have
$$\sum_n |c_n|^2 = 1$$

Measurement postulate



On measurement of a state
the system collapses into the n th
eigenstate of the quantity being
measured

with probability $P_n = |c_n|^2$

In the expansion of the state
in the eigenfunctions
of the quantity being measured
 c_n is the expansion coefficient
of the n th eigenfunction

Expectation value of the energy

Suppose do an experiment to measure the energy E of some quantum mechanical system

We could repeat the experiment many times
and get a statistical distribution of results

Given the probabilities P_n of getting a specific energy eigenstate, with energy E_n

we would get an average answer

$$\langle E \rangle = \sum_n E_n P_n = \sum_n E_n |c_n|^2$$

where we call this average $\langle E \rangle$ the “**expectation value**”

Energy expectation value example

For example, for the coherent state discussed above with parameter N , we have

$$\langle E \rangle = \sum_{n=0}^{\infty} E_n \frac{N^n \exp(-N)}{n!} = \hbar\omega \left[\sum_{n=0}^{\infty} n \frac{N^n \exp(-N)}{n!} \right] + \frac{1}{2} \hbar\omega = \left(N + \frac{1}{2} \right) \hbar\omega$$

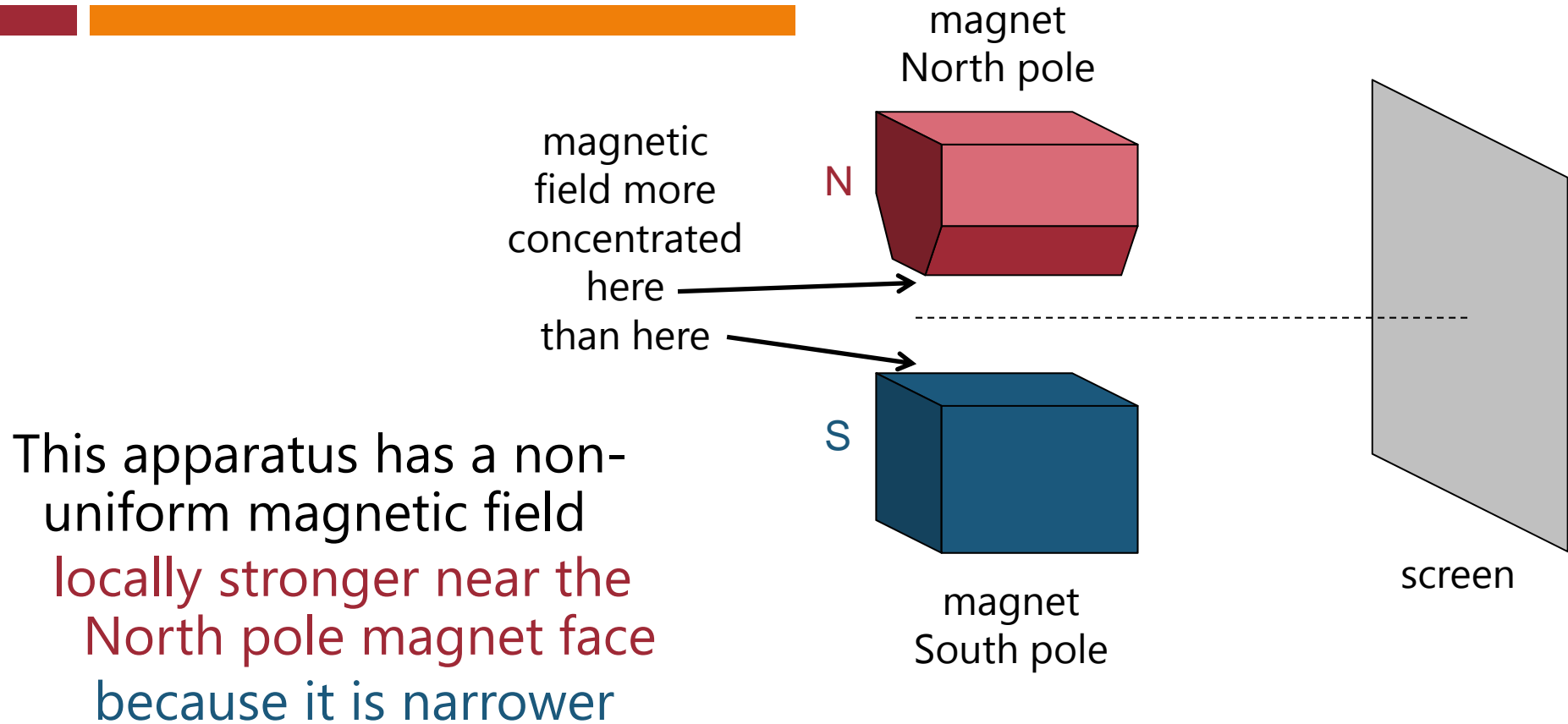
where we use the result that the average in a Poisson statistical distribution is just the parameter N

Note that N does not have to be an integer

This is an expectation value, not an eigenvalue

We can have states with any expectation value we want

Stern-Gerlach experiment



Stern-Gerlach experiment

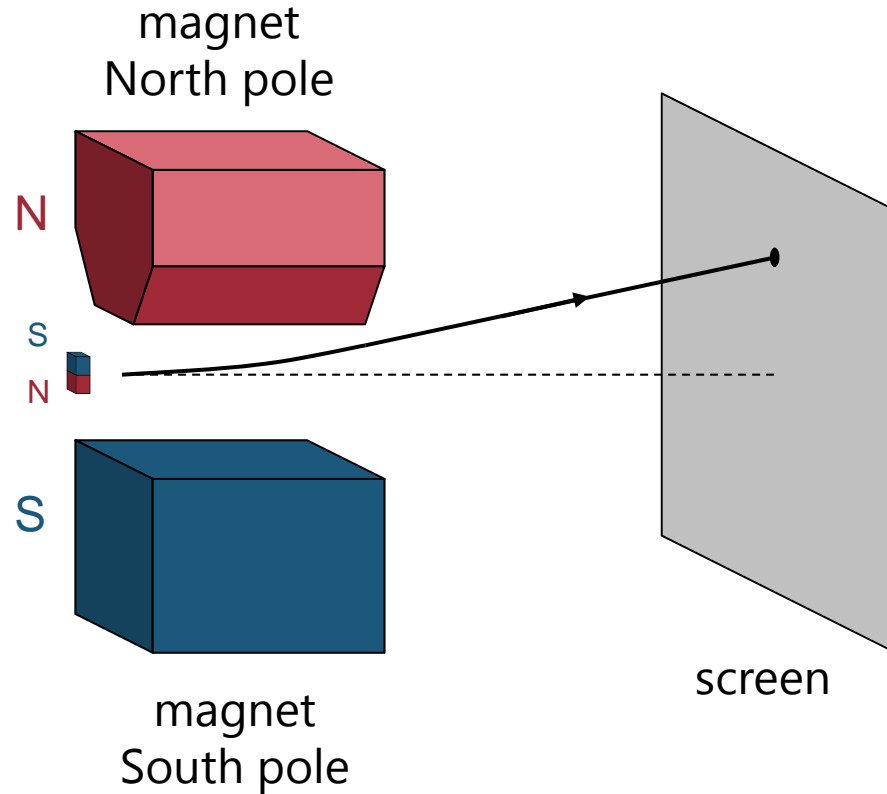
Imagine firing some small magnets

initially along the dashed line

Because the magnetic field is non-uniform

stronger near the North pole than near the South pole

a vertical magnet will be deflected up



Stern-Gerlach experiment

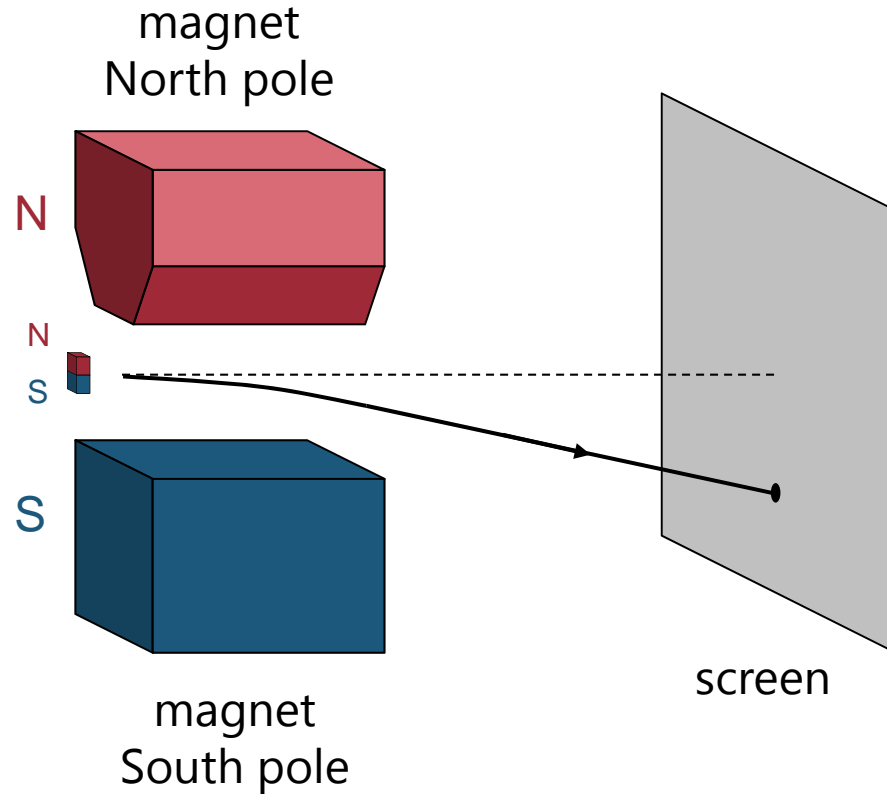
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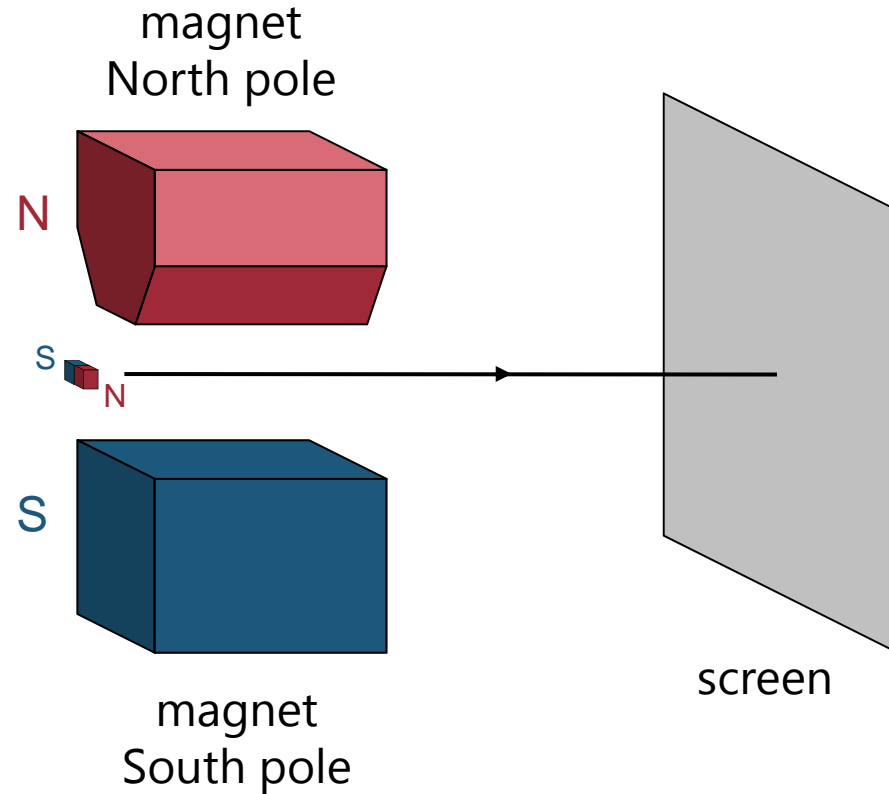
stronger near the North pole than near the South pole

a vertical magnet will be deflected up or down



Stern-Gerlach experiment

A horizontally-oriented magnet
will not be deflected



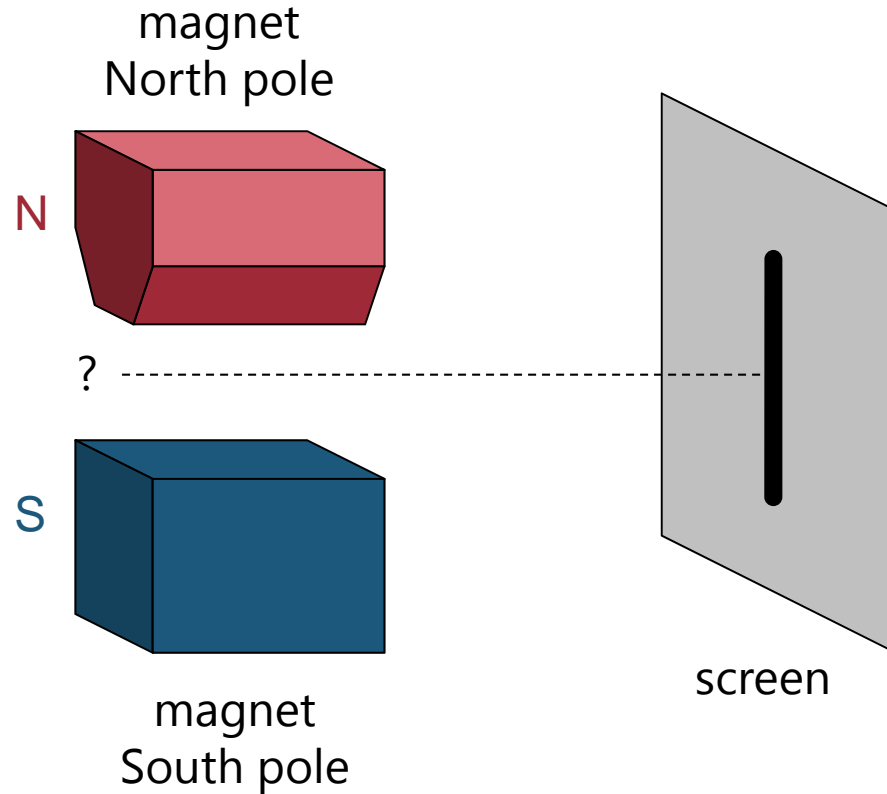
Stern-Gerlach experiment

A horizontally-oriented magnet
will not be deflected
and magnets of other
orientations

should be deflected by
intermediate amounts

After “firing” many randomly
oriented magnets

we should end up with a line
on the screen



Electrons and the Stern-Gerlach experiment

Electrons have a quantum mechanical property called spin

It gives them a “magnetic moment”
just like a small magnet

What will happen if we fire electrons
with no particular “orientation” of their spin
into the Stern-Gerlach apparatus?

We might expect the “line” on the screen

(Note: the actual experiment used silver atoms,
which behave the same as electrons in this case)

Stern-Gerlach experiment

With electrons

we get two dots!

"Explanation"

We are measuring the vertical component of the spin

There are two eigenstates of this component

up and down

so we have collapse to the eigenstates

