

# Normal Order

## Normal order

From Wikipedia, the free encyclopedia

*For the evaluation strategy, see [Normal order evaluation](#).*

When a [quantum mechanical Hamiltonian](#) is written in **normal order**, it means that all creation operators are placed to the left of all annihilation operators in the expression. The process of putting a Hamiltonian into normal order is called **normal ordering**. The terms **antinormal order** and **antinormal ordering** are analogously defined, where the annihilation operators are placed to the left of the creation operators.

When quantizing a [classical](#) Hamiltonian there is some freedom when choosing the operator order, and these choices lead to differences in the [ground state energy](#).

<http://books.nap.edu/html/biomems/gwick.html>



Gian-Carlo Wick made many fundamental contributions to nuclear and particle physics from the 1930s, when he was a close associate of Enrico Fermi in Rome, to the 1970s, when he worked with Tsung Dao Lee at Columbia. His landmark paper, entitled "Evaluation of the collision matrix" (1950), shows how to conduct explicit practical calculations starting from the formal relations of relativistic quantum field theory through expression of the chronological product of quantum fields in terms of a sum of normal products.

## Wick's Theorem (again, see Jacek's notes)

### Wick's Theorem for expectation values of operators

Consider an operator  $\hat{A}$  and a wave function  $|\Psi\rangle$

There always exists a decomposition:

$$\hat{A} = \hat{A}_0 + \hat{A}_+ + \hat{A}_-$$

where

$1^\circ \quad \hat{A}_0 = \text{const.}$
$2^\circ \quad \hat{A}_-  \Psi\rangle = 0,$
$3^\circ \quad \langle \Psi   \hat{A}_+ = 0.$

To demonstrate it, we introduce the projection operator:

$$\hat{P} \equiv |\Psi\rangle\langle\Psi|$$

In terms of this operator, the above decomposition can be written as:

$$\begin{aligned}\hat{A}_0 &= \langle\Psi|\hat{A}|\Psi\rangle \\ \hat{A}_- &= \left(\hat{A} - \langle\Psi|\hat{A}|\Psi\rangle\right) (1 - \hat{P}) \\ \hat{A}_+ &= (1 - \hat{P})\hat{A}\hat{P}\end{aligned}$$

Let us now calculate an expectation value of a product of two operators:

$$\begin{aligned}\langle \Psi | \hat{A} \hat{B} | \Psi \rangle &= \langle \Psi | (\hat{A}_0 + \hat{A}_+ + \hat{A}_-) (\hat{B}_0 + \hat{B}_+ + \hat{B}_-) | \Psi \rangle \\ &= \hat{A}_0 \hat{B}_0 + \langle \Psi | \hat{A}_- \hat{B}_+ | \Psi \rangle\end{aligned}$$

or

$$\langle \Psi | \hat{A} \hat{B} | \Psi \rangle = \langle \Psi | \hat{A} | \Psi \rangle \langle \Psi | \hat{B} | \Psi \rangle + \langle \Psi | \hat{A}_- \hat{B}_+ | \Psi \rangle$$

An average of a product

A product of averages

A “trouble” term

$$\begin{aligned}\langle \Psi | \hat{A}_- \hat{B}_+ | \Psi \rangle &= \langle \Psi | \hat{A}_- \hat{B} | \Psi \rangle = \langle \Psi | \hat{A} \hat{B}_+ | \Psi \rangle \\ &= \langle \Psi | \{ \hat{A}_-, \hat{B}_+ \} | \Psi \rangle = \langle \Psi | \{ \hat{A}_-, \hat{B} \} | \Psi \rangle = \langle \Psi | \{ \hat{A}, \hat{B}_+ \} | \Psi \rangle \\ &= \langle \Psi | [ \hat{A}_-, \hat{B}_+ ] | \Psi \rangle = \langle \Psi | [ \hat{A}_-, \hat{B} ] | \Psi \rangle = \langle \Psi | [ \hat{A}, \hat{B}_+ ] | \Psi \rangle.\end{aligned}$$

Useful when dealing with fermions

Useful when dealing with bosons

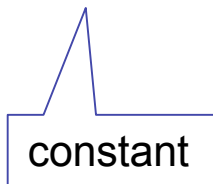
How to calculate the product of many operators?

$$\langle \Psi | \hat{A}_1 \hat{A}_2 \dots \hat{A}_n | \Psi \rangle$$

$$\begin{aligned} & \langle \Psi | \hat{A}_1 \hat{A}_2 \dots \hat{A}_n | \Psi \rangle \\ &= \langle \Psi | \hat{A}_1 | \Psi \rangle \langle \Psi | \hat{A}_2 \dots \hat{A}_n | \Psi \rangle + \langle \Psi | \hat{A}_1 - \hat{A}_2 \dots \hat{A}_n | \Psi \rangle \end{aligned}$$

Let us introduce **the contraction**:

$$\overline{\hat{A}\hat{B}} \equiv \hat{A}\hat{B} - c\hat{B}\hat{A}$$


  
constant

For  $c=-1$  ( $+1$ ), contraction becomes an anticommutator (commutator)

It is also convenient to introduce **the self-contraction**:

$$\overline{\hat{A}} \equiv \hat{A}_0 \quad \text{or} \quad \overline{\hat{A}} = \langle \Psi | \hat{A} | \Psi \rangle$$

Why is contraction useful?

$$\begin{aligned} \langle \Psi | \hat{A}_1 \hat{A}_2 \hat{A}_3 \dots \hat{A}_n | \Psi \rangle &= \overline{\hat{A}_1} \langle \Psi | \hat{A}_2 \hat{A}_3 \dots \hat{A}_n | \Psi \rangle + \langle \Psi | \overline{\hat{A}_1 \hat{A}_2} \hat{A}_3 \dots \hat{A}_n | \Psi \rangle \\ &+ c \langle \Psi | \hat{A}_2 \overline{\hat{A}_1 \hat{A}_3} \dots \hat{A}_n | \Psi \rangle + \dots + c^{n-2} \langle \Psi | \hat{A}_2 \hat{A}_3 \dots \overline{\hat{A}_1 \hat{A}_n} | \Psi \rangle, \end{aligned}$$

**Are we breaking the First Weinberg's Law of Progress in Theoretical Physics ?**

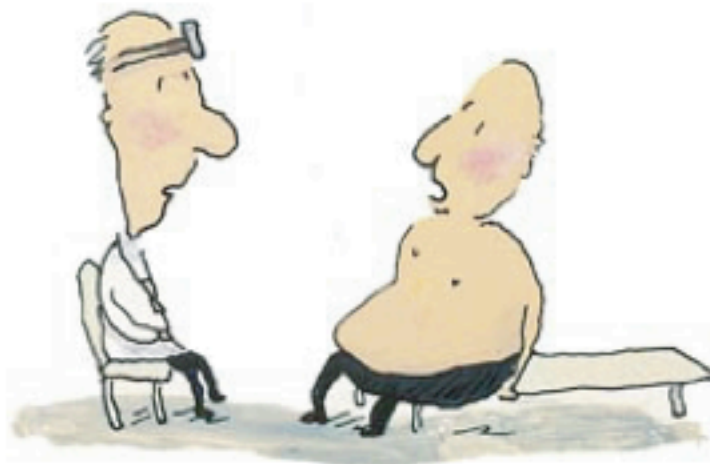
# Weinberg's Laws of Progress in Theoretical Physics

From: "Asymptotic Realms of Physics" (ed. by Guth, Huang, Jaffe, MIT Press, 1983)

**First Law:** "The conservation of Information" (*You will get nowhere by churning equations*)

**Second Law:** "Do not trust arguments based on the lowest order of perturbation theory"

**Third Law:** "You may use any degrees of freedom you like to describe a physical system, but if you use the wrong ones, you'll be sorry!"



**Patient:** Doctor, doctor, it hurts when I do this!

**Doctor:** Then don't do that.