

# physics2 manual for the legacy physics users

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## Abstract

This short document describes `physics2` package for those who are used to the `physics` package. This document is only a simple reference manual for:

- Frequent users of the legacy `physics` package;
- Those who have to maintain a document written with `physics`;
- Users who failed to use `unicode-math` with `physics`.

It seems no reason for any other user to read *this* document instead of the [package documentation](#) of `physics2`, because this document cannot describe the package in detail.

In this document, the modules of `physics2` will be introduced in the same order as the `physics` documentation.

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\*<https://www.github.com/AlphaZTX/physics2>

# 1 Before you start

## 1.1 Legacy problems with **physics** package

The **physics** package provides `\qty` command for automatic-sizing braces. The `\qty` command would cause conflict with the **siunitx** package, which provides a unified method to typeset numbers and units correctly.

Besides, after you loaded **physics**, when you type `\homework` you will get Maxwell equations and Schrödinger equation. The `\homework` command is “declared” in `physics.sty` but it was not described in the documentation. That is, if you have defined `\homework` before loading **physics** package, **physics** would overwrite the definition “silently”.

The vector-notation part of **physics** uses **amsmath**’s (more exactly, `amsbsy.sty`’s) `\boldsymbol` command to generate bold vectors. Commands for cross/dot product are defined with `\boldsymbol`. `\boldsymbol` uses `\mathversion`, a  $\TeX$  2<sub>ε</sub> kernel command that works well with traditional TFM-based fonts but fails when using **unicode-math**.

In the definition of `\imat`, `\xmat`, `\dmat` and `\admat` commands from **physics**, there is a `\newtoks` command which allocates a token list register and two `\newcount` commands allocating two count registers. Every time you write a command like `\imat` in your document, then one token list register and two count registers will be wasted. What’s even worse is that, if you wrote really too many matrix commands from **physics** (for example, 32767 `\imats` in Lua $\TeX$ ), there’d be no room for a new `\count`.

**physics** integrated all the functions in one file (`physics.sty`), that is, you cannot load one of the total seven parts of functions; you have to load the seven parts altogether, even included the extra `\homework` command we mentioned in the first paragraph.

Moreover, the code of `physics.sty` “abuses” the g-type arguments of **xparse** package. Therefore the syntax of **physics** package looks kind of weird. See [here](#) for more.

## 1.2 Loading **physics2**

The **physics2** package includes different modules, among which every module focuses on one single function.

Write the following line in the preamble to load **physics2**:

```
\usepackage{physics2}
```

But this is not enough. `physics2` contains different modules, among which, only the `common` module would be loaded automatically by the package. If you want to load other modules of `physics2`, write this after loading `physics2` package:

```
\usephysicsmodule{<module list>}
```

For example, “`\usephysicsmodule{ab,doubleprod}`” loads the `ab` module and the `doubleprod` module.

You can also load a module with options:

```
\usephysicsmodule[<option list>]{<module>}
```

For example, “`\usephysicsmodule[legacy]{ab}`” loads `ab` with the option “`legacy`”.



Attention, if you used any font package in your document, remember that `physics2` requires to be loaded *after* font packages.

## 2 List of commands

### 2.1 Automatic bracing

As mentioned in §1.1, the `\qty` command from `physics` would cause conflicts with `siunitx`. The command for automatic braces in `physics2` is `\ab`, a shorthand for automatic braces.

The `\ab` command requires the `ab` module, so don’t forget to write `\usephysicsmodule{ab}` in the preamble after you loaded `physics2`. Always remember, *do not put an `\ab` separately in the end of a math formula*. Take some examples:

[2.1.1]      `\[ \ab ( \frac{1}{2} ) \quad`  
               `\ab [ \frac{1}{2} ] \quad`  
               `\ab{ \frac{1}{2} } \]`

$$\left(\frac{1}{2}\right) \quad \left[\frac{1}{2}\right] \quad \left\{\frac{1}{2}\right\}$$

`\ab` can modify a delimiter-braced subformula. But the delimiters should not be out of the range described by the following chart:

<code>(, )</code>	
<code>[, ]</code>	
<code>\{, \}</code>	or <code>\lbrace, \rbrace</code>
<code>&lt;, &gt;</code>	or <code>\langle, \rangle</code>
<code> ,  </code>	or <code>\vert, \vert</code>
<code>\ , \ </code>	or <code>\Vert, \Vert</code>

For example,  $\backslash ab\{foo\}$  and  $\backslash ab(foo)$  are illegal, but  $\backslash ab\{foo\}$  and  $\backslash ab(foo)$  are okay;  $\backslash ab([])$  is okay but  $\backslash ab(())$  is illegal.



Attention, if you want to delimit a subformula with “{” and “}”, you can only write  $\backslash \{$ ,  $\backslash \}$  or  $\backslash lbrace$ ,  $\backslash rbrace$  around it. { and } are not supported in **ab** module.

Between  $\backslash ab$  and the first delimiter can be a “biggg” command, that is, from  $\backslash big$  to  $\backslash Bigg$ . Actually, you can also write  $\backslash biggg$  and  $\backslash Biggg$  because **physics2** defines these after you load it. For example,

[2.1.2] 
$$\begin{aligned} &\backslash [ \backslash ab\backslash Big \backslash | \backslash frac{1}{2} \backslash | \backslash quad \\ &\backslash ab\backslash Bigg < \backslash frac{1}{2} > \backslash quad \\ &\backslash ab\backslash Biggg | \backslash frac{1}{2} | \backslash \end{aligned}$$

$$\left\| \frac{1}{2} \right\| \quad \left\langle \frac{1}{2} \right\rangle \quad \left| \frac{1}{2} \right|$$

Between  $\backslash ab$  and the first delimiter can also be a star (\*), which means “use the default size of delimiters”. But in this situation, you needn’t use the  $\backslash ab$  command at all.

The **physics** package provides commands like  $\backslash pqty$ ,  $\backslash bqty$ . In the **ab** module of **physics2**, these commands have changed to  $\backslash pab$ ,  $\backslash bab$ , etc. The following example shows all the  $\backslash Xab$  commands in **ab** module:

[2.1.3] 
$$\begin{aligned} &\backslash def\backslash 0\{\backslash frac{1}{2}\} \\ &\backslash [ \backslash pab\{\backslash 0\} \backslash quad \backslash bab\{\backslash 0\} \\ &\quad \quad \quad \backslash quad \backslash Bab\{\backslash 0\} \backslash \\ &\backslash [ \backslash aab\{\backslash 0\} \backslash quad \backslash vab\{\backslash 0\} \\ &\quad \quad \quad \backslash quad \backslash Vab\{\backslash 0\} \backslash \end{aligned}$$

$$\begin{aligned} &\left(\frac{1}{2}\right) \quad \left[\frac{1}{2}\right] \quad \left\{\frac{1}{2}\right\} \\ &\left\langle \frac{1}{2} \right\rangle \quad \left|\frac{1}{2}\right| \quad \left\|\frac{1}{2}\right\| \end{aligned}$$

$\backslash Xab$  can take an optional star and an optional [*biggg*] argument. For example,

[2.1.4] 
$$\begin{aligned} &\backslash def\backslash 0\{\backslash frac{1}{2}\} \\ &\backslash [ \backslash pab[Big]\{\backslash 0\} \backslash quad \backslash bab*\{\backslash 0\} \backslash \end{aligned}$$

$$\left(\frac{1}{2}\right) \quad \left[\frac{1}{2}\right]$$

**physics** also provides the following commands:

$\backslash abs$     $\backslash norm$     $\backslash eval$     $\backslash order$     $\backslash comm$     $\backslash acomm$     $\backslash pb$



These commands are not originally supported by **physics2**, but the first four commands can be used through the **ab.legacy** module of **physics2**:

$\backslash usephysicsmodule\{ab.legacy\}$

For example,

[2.1.5] 
$$\begin{aligned} &\backslash def\backslash 0\{\backslash frac{1}{2}\} \\ &\backslash [ \backslash abs\{\backslash 0\} \backslash quad \backslash abs[big]\{\backslash 0\} \\ &\quad \quad \quad \backslash quad \backslash abs*\{\backslash 0\} \backslash \end{aligned}$$

$$\left|\frac{1}{2}\right| \quad \left|\frac{1}{2}\right| \quad \left|\frac{1}{2}\right|$$

Users of the legacy `physics` package should notice that the syntax of `\eval` has been changed. The `ab.legacy` module abandoned the `\eval(foo|`-like syntax. The new `\eval`’s syntax is just like other commands in this module. There are also two variants of `\eval` — `\peval` and `\beval`. For example,

[2.1.6]

```

\def\0{1+\frac{1}{2}x}
\[\eval{\0}_a^b \quad
\peval*{\0}_a^b \quad
\beval[big]{\0}_a^b \quad \]
```

$$1 + \frac{1}{2}x \Big|_a^b \quad \left(1 + \frac{1}{2}x\right)_a^b \quad \left[1 + \frac{1}{2}x\right]_a^b$$

The `\comm`, `\acomm` and `\pb` (Poisson bracket) are not supported. But you can write like `\ab[foo,baz]` or `\bab{foo,baz}` instead.

By the way, you can set the “order” symbol in `ab.legacy` through the `order` option like this:

```
\usephysicsmodule[order=0]{ab.legacy}
```

Then `\order(N)` yields  $O(N)$ .

## 2.2 Vector notation

Unfortunately, there is not a plan for `physics2` to support this part of `physics` completely, but the rest of this section will show some methods to maintain the document written with `physics`.

The `\vb(*)`, `\va(*)` and `\vu(*)` are not supported in any module of `physics2`. But these commands can be defined by copying the following lines below and pasting them in the preamble:

```

\makeatletter
\newcommand\vb{\@ifstar\boldsymbol\mathbf}
\newcommand\va[1]{\@ifstar{\vec{#1}}{\vec{\mathrm{#1}}}}
\newcommand\vu[1]{%
  \@ifstar{\hat{\boldsymbol{#1}}}{\hat{\mathbf{#1}}}}
\makeatother
```

The `\boldsymbol` command requires the `amsmath` or `bm` package. If you prefer to use `bm`, you can also use the `\bm` command. What’s more, if you tried the commands above, you might find that, the result of `\va` above is different from that of `physics`. This is because, if you choose to present a vector in bold, there’s almost no need to put a `\vec` (“”) sign above it.

However, the method above may not work well with `unicode-math` because there are so many OpenType math fonts without a bold version. When using

`unicode-math`, it's recommended to use `\symbf` and `\symbfitt` for a separate vector. For example, `\symbf{0}` yields **0**, and `\symbfitt{A}` yields **A**.

The `\vdot` and `\cross` commands are not supported in any module of `physics2`. Actually, there is no need to use a bold “ $\cdot$ ” or “ $\times$ ” for the products of two vectors. Using `\cdot` and `\times` is enough.

The commands related to “ $\nabla$ ” are supported through `nabla.legacy` module. These commands are `\grad`, `\div` and `\curl`. These commands should not be put in the end of a math formula either (just like `\ab`). Notice that the former `\div` command for a “ $\div$ ” (if there exists one) is redefined as `\divsymbol`. For example,

[2.2.1]      % \usephysicsmodule{nabla.legacy}  
              \[ \grad F          \quad  
                  \grad(\frac{G2})        \  
              \[ \div\Bigg[X]     \quad  
                  \curl\*\{\frac{Y2}\} \  
              \[ 2 \divsymbol 1        \]

$$\nabla F \quad \nabla \left( \frac{G}{2} \right) \\ \nabla \cdot \left[ X \right] \quad \nabla \times \left\{ \frac{Y}{2} \right\} \\ 2 \div 1$$

Actually, the nabla-related commands end with `\ab`. Thus, the subformula after these commands can be delimited with `()`, `[]` and `\{\}`.

The `nabla.legacy` requires the `fixdf` package at least version 2.0 (file date: 2023/01/31 or after 2023/01/31).

By the way, if you are used to writing `\bm` for a vector but interested in `unicode-math`, the `bm-um.legacy` module would be a passable alternative to `bm` package. Notice that the `\bm` command from the `bm-um.legacy` module can only take *one* letter (or *one* digit) as its argument.

## 2.3 Operators

There's no plan for `physics2` to support this part of `physics` completely. The syntax in this part of `physics` (such as `\tan[2](x)`) abuses `xparse`.

It's suggested to write like this if you used the `ab` module:

$$\sin^2 \left( \frac{\alpha}{2} \right)$$

Rather than take the superscript as an optional argument of the command of log-like functions.

The `physics` package provides a bundle of commands for log-like functions that have not been defined in the  $\text{\LaTeX}_{\mathcal{E}}$  kernel. Those log-like functions can be used with the `op.legacy` module; this module does not support the syntax of `physics` either. For example:

[2.3.1] `% \usephysicsmodule{op.legacy}`  
`\[ \asin^2 x \quad \text{rank} \{ A \} \]`

$\text{asin}^2 x$      $\text{rank}\{A\}$

The `\Re` and `\Im` commands are redefined as operators “Re” and “Im”, while  $\Re$  and  $\Im$  are reserved as `\Resymbol` and `\Imsymbol`.  $\Re$  and  $\Im$  are ordinary symbols but `Re` and `Im` are operators.

## 2.4 Quick quad text

The `qtext.legacy` module provides the `\q{foo}` commands for `\quad`-wrapped texts. These commands have the same syntax as `physics`. For example,

[2.4.1] `% \usephysicsmodule{qtext.legacy}`  
`\[ A \qq {foo bar} B \]`  
`\[ A \qq*{foo bar} B \]`  
`\[ C \qcc D \qcc* E \]`  
`\[ F \qif G \qthen H \]`

$A$      $\text{foo bar}$      $B$

$A\text{foo bar}$      $B$

$C$      $c.c$      $Dc.c$      $E$

$F$      $\text{if } G$      $\text{then } H$

All the commands described in §2.4 of [physics documentation](#) are supported when using `qtext.legacy` module, but I don’t recommend using this module unless you are maintaining a document written with `physics`’s `\q{foo}` commands.

## 2.5 Derivatives

There is no plan for `physics2` to support this part of `physics`. If you want to typeset the differential operators on a better sense, you can try the `fixdif` package; if you want an easy way to type derivatives, you can try the `derivative` package. These two packages can be used together. For example,

[2.5.1] `% \usepackage{fixdif,derivative}`  
`\[ \pdv{f}{x,y,z} \, \text{d} x \]`  
`Math (\text{d} x) \text{ v.s. } \text{Text} (\text{d} x)`

$\frac{\partial^3 f}{\partial x \partial y \partial z} \text{d} x$

Math ( $\text{d} x$ ) v.s. Text ( $\text{d} x$ )

Here are the documentations of [fixdif](#) and [derivative](#).

[fixdif](#)’s commands behave better in superscripts and subscripts.

## 2.6 Dirac bra-ket notation

There are two solutions to Dirac bra-ket in `physics2` — `ab.braket` and `braket`. These two modules are *not* compatible and neither of them supports `physics`’s syntax completely. Click [here](#) to see the `ab.braket` module and [here](#) to see the `braket` module.

**The `ab.braket` module** This module provides four commands — `\bra`, `\ket`, `\braket` and `\ketbra`. After these commands can be a star (\*) or a “biggg” command. These commands share similar syntaxes like `\ab`’s syntax. But, *the bra-ket commands from `ab.braket` module are completely different from `\ab`*. Their internal structures are different.

The argument of `\bra` should be delimited with `<` and `|`, that is,

$$\backslash\bra < \langle subformula \rangle |$$

For example,

[2.6.1] `\[ \bra < \frac \phi 2 | \]`  
`\[ \bra* < \frac \phi 2 | \]`  
`\[ \bra\Big < \phi | \]`

$$\begin{array}{c} \left\langle \frac{\phi}{2} \right| \\ \left\langle \frac{\phi}{2} \right| \\ \left\langle \phi \right| \end{array}$$

The argument of `\ket` should be delimited with `|` and `>`, that is,

$$\backslash\ket | \langle subformula \rangle >$$

For example,

[2.6.2] `\[ \ket | \frac \psi 2 > \]`  
`\[ \ket* | \frac \psi 2 > \]`  
`\[ \ket\Big | \psi > \]`

$$\begin{array}{c} \left| \frac{\psi}{2} \right\rangle \\ \left| \frac{\psi}{2} \right\rangle \\ \left| \psi \right\rangle \end{array}$$



If you want to write “>” and “<” for relations in the argument of `\bra` and `\ket`, you can write `\mathrel{>}` and `\mathrel{<}` (although there is almost no such need).

The argument of `\braket` should be delimited with `<` and `>`, that is,

$$\backslash\braket < \langle subformula \rangle >$$

In the  $\langle subformula \rangle$  argument, every “|” will be regarded as an extensible vertical bar. For example,



[2.6.3] `\[ \braket< \phi > \]`  
`\[ \braket< \phi | \psi > \]`  
`\[ \braket< \phi | A | \psi > \]`

$$\langle \phi \rangle$$

$$\langle \phi | \psi \rangle$$

$$\langle \phi | A | \psi \rangle$$

[2.6.4] `\def\0{\frac\phi2}`  
`\[ \braket < \0 | \psi > \]`  
`\[ \braket* < \0 | \psi > \]`  
`\[ \braket\Bigg< \0 | \psi > \]`

$$\left\langle \frac{\phi}{2} \middle| \psi \right\rangle$$

$$\left\langle \frac{\phi}{2} \middle| \psi \right\rangle$$

$$\left\langle \frac{\phi}{2} \middle| \psi \right\rangle$$

The argument of `\ketbra` should be delimited with `|` and `|`. In the argument, `>` and `<` will be regarded as extensible `\rangle` and `\langle`. That is,

$$\ketbra{| \langle subformula_1 \rangle > \langle optional \rangle < \langle subformula_2 \rangle |}$$

For example,

[2.6.5] `\def\0{\frac\phi2}`  
`\[ \ketbra | \0 > \psi | \]`  
`\[ \ketbra* | \0 > \psi | \]`  
`\[ \ketbra\Bigg| \0 > \psi | \]`

$$\left| \frac{\phi}{2} \right\rangle \left\langle \psi \right|$$

$$\left| \frac{\phi}{2} \right\rangle \left\langle \psi \right|$$

$$\left| \frac{\phi}{2} \right\rangle \left\langle \psi \right|$$

[2.6.6] `\def\0{\frac\phi2}`  
`\[ \ketbra| \0 >_x^y < \psi | \]`

$$\left| \frac{\phi}{2} \right\rangle_x^y \left\langle \psi \right|$$



If you want to write “`>`” and “`<`” for relations in the argument of `\braket` and `\ketbra`, you can write `\>` and `\<` (although there is almost no such need). It is quite different from `\mathrel{>}` or `\mathrel{<}` because in these commands’ argument, `>` and `<` will be redefined.

**The `braket` module** This module contains four commands — `\bra`, `\ket`, `\braket` and `\ketbra`. After these commands can be a star (\*) or a square-bracket-delimited size option, the size option can take the following values:

big, Big, bigg, Bigg, biggg or Biggg.

Star stands for “do not size the bra-ket automatically”.

The argument(s) of these four commands are braced with { and }. \bra and \ket take one mandatory argument. For example,

[2.6.7] 

```
\def\0{\frac\phi2}
\[ \bra {\0} \quad \bra* {\0}
\quad \bra[Big] {\0} \]
\[ \ket {\0} \quad \ket* {\0}
\quad \ket[Big] {\0} \]
```

$$\begin{array}{ccc} \left\langle \frac{\phi}{2} \right| & \left\langle \frac{\phi}{2} \right| & \left\langle \frac{\phi}{2} \right| \\ \left| \frac{\phi}{2} \right\rangle & \left| \frac{\phi}{2} \right\rangle & \left| \frac{\phi}{2} \right\rangle \end{array}$$

The \braket command, in default, can take two arguments.

[2.6.8] 

```
\def\0{\frac\phi2}
\[ \braket {\0} {\psi} \quad
\braket* {\0} {\psi} \quad
\braket[big] {\0} {\psi} \]
```

$$\left\langle \frac{\phi}{2} \right| \psi \rangle \quad \langle \frac{\phi}{2} | \psi \rangle \quad \langle \frac{\phi}{2} | \psi \rangle$$

If you want \braket to take one or three arguments, you can write the number of arguments in the square bracket. If you need to specify the size of bra-ket simultaneously, you need to separate the number and the size with a comma:

[2.6.9] 

```
\def\0{\frac\phi2}
\[ \braket [1] {\0} \quad
\braket*[1] {\0} \]
\[ \braket [3] {\0}{A}{\psi} \quad
\braket[3,big] {\0}{A}{\psi}
\quad
\braket[Big,3] {\0}{A}{\psi} \]
```

$$\begin{array}{cc} \left\langle \frac{\phi}{2} \right\rangle & \left\langle \frac{\phi}{2} \right\rangle \\ \left\langle \frac{\phi}{2} \right| A | \psi \rangle & \\ \left\langle \frac{\phi}{2} \right| A | \psi \rangle & \left\langle \frac{\phi}{2} \right| A | \psi \rangle \end{array}$$

The \ketbra command takes two mandatory arguments. It can also take an optional argument between the two mandatory arguments. The optional argument will be placed between > and <:

[2.6.10] 

```
\def\0{\frac\phi2}
\[ \ketbra {\0} {\psi} \quad
\ketbra* {\0} {\psi} \quad
\ketbra [Bigg] {\0} {\psi} \quad
\ketbra {\0} [_x^y] {\psi} \]
```

$$\begin{array}{cc} \left| \frac{\phi}{2} \right\rangle \langle \psi | & \left| \frac{\phi}{2} \right\rangle \langle \psi | \\ \left| \frac{\phi}{2} \right\rangle \langle \psi | & \\ \left| \frac{\phi}{2} \right\rangle_x^y \langle \psi | & \end{array}$$

## 2.7 Matrix macros

Unfortunately, **physics2** do not support the `\mqty` command from **physics**. If you are used to this command, you can write like this:

```
\newcommand\mqty[1]{\begin{matrix}#1\end{matrix}}
\newcommand\pmqty[1]{\begin{pmatrix}#1\end{pmatrix}}
$\ab(\mqty{foo})$ or $\pmqty{foo}$
```

These are equal to **physics**'s `\mqty(foo)` (require **amsmath**).

**physics2**'s **diagmat** module provides `\diagmat` command for diagonal matrices. For example,

[2.7.1] `\[`  
`\diagmat { 1, 2, 3 }`  
`\]`

$$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 2 & 0 \\ 0 & 0 & 3 \end{pmatrix}$$

[2.7.2] `\[`  
`\pdiagmat [ empty = {} ]`  
`{ a, b, c, d }`  
`\]`

$$\begin{pmatrix} a & & & \\ & b & & \\ & & c & \\ & & & d \end{pmatrix}$$

`\pdiagmat`, `\bdiagmat`, `\Bdiagmat`, `\vdiagmat` and `\Vdiagmat` are also available.

**physics2**'s **xmat** module provides `\xmat` command for matrices with formatted entries. For example,

[2.7.3] `\[`  
`\xmat{a}{2}{3}`  
`\]`

$$\begin{matrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \end{matrix}$$

[2.7.4] `% \usephysicsmodule`  
`% [showleft=3,showtop=3]{xmat}`  
`\[`  
`\pxmat{X}{m}{n}`  
`\]`

$$\begin{pmatrix} X_{11} & X_{12} & X_{13} & \cdots & X_{1n} \\ X_{21} & X_{22} & X_{23} & \cdots & X_{2n} \\ X_{31} & X_{32} & X_{33} & \cdots & X_{3n} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ X_{m1} & X_{m2} & X_{m3} & \cdots & X_{mn} \end{pmatrix}$$

[2.7.5] `\[`  
`\xmat [showleft=2,showtop=2,`  
`format=\texttt{#1[#2][#3]}}`  
`{x}{m}{n}`  
`\]`

$$\begin{matrix} x[1][1] & x[1][2] & \cdots & x[1][n] \\ x[2][1] & x[2][2] & \cdots & x[2][n] \\ \vdots & \vdots & \ddots & \vdots \\ x[m][1] & x[m][2] & \cdots & x[m][n] \end{matrix}$$

`\pxmat`, `\bxmat`, `\Bxmat`, `\vxmat` and `\Vxmat` are also available.