

# Code documentation to the `physics2` package

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# File I

## The bare **physics2**

### 1 The **physics2** package

```
1 (*package)
2 \NeedsTeXFormat{LaTeX2e}[2020/10/01]
3 \ProvidesPackage{physics2}
4 [2024/01/10 v1.0.1 Tools for typesetting math for physics.]
```

#### 1.1 Common variables

---

```
\phy@temp.. \phy@temp@register type}{a or b}
```

Some  $\text{\LaTeX} 2\epsilon$  variables starting with “`\phy@temp`”. These variables can be shared by any module of **physics2**.

```
5 \newcount \phy@tempcnta
6 \newdimen \phy@tempdima
7 \newdimen \phy@tempdimb
8 \newskip \phy@tempskipa
9 \newmuskip \phy@tempmuskipa
10 \newbox \phy@tempboxa
11 \newif \ifphy@tempswa
12 \newtoks \phy@toksa
```

#### 1.2 Package requirements and module-loading methods

**physics2** requires **keyval** (part of the graphics bundle) to process options of modules.

```
13 \RequirePackage{keyval}
14 \def\phy@true{true}
15 \def\phy@false{false}
```

---

```
\phy@define@key \phy@define@key {\langle module \rangle} {\langle key \rangle} [{\langle default value \rangle}] {\langle code \rangle}
\phy@setkeys \phy@setkeys {\langle module \rangle} {\langle key-val list \rangle}
\phy@processkeyopt \phy@processkeyopt {\langle module \rangle}
```

The position of `\phy@processkeyopt` in a **physics2** module is just the same as the position of `\ProcessOptions` in a regular  $\text{\LaTeX}$  package.

```
16 \long\def\phy@define@key#1{\define@key{phy-#1}}
17 \long\def\phy@setkeys#1{\setkeys{phy-#1}}
18 \def\phy@processkeyopt#1{\let\reserved@a\empty%
19   \edef\reserved@a{\optionlist{\currname.\@currext}}%
20   \edef\reserved@a{\noexpand\phy@setkeys{#1}\f\reserved@a}%
21   \reserved@a% the next line thanks to `geometry'
22   \AtEndOfPackage{\let\unprocessedoptions\relax}}
```

We use almost the same way to load **physics2** modules as  $\text{\LaTeX} 2\epsilon$  kernel does. We use a lot of kernel commands in  $\text{\LaTeX} 2\epsilon$ .

---

```
\usephysicsmodule \usephysicsmodule [<key-val options>] {<module>} [<key-val options>]
\phy@requiremodule \phy@requiremodule [<key-val options>] {<module>} [<key-val options>]

\usephysicsmodule is a user command, and \phy@requiremodule is a developer com-
mand.

23 \def\usephysicsmodule{\phy@FWoptions\@pkgextension}
24 \let\phy@requiremodule\usephysicsmodule
25 \onlypreamble\usephysicsmodule
26 \def\phy@FWoptions#1{\@ifnextchar[%]
27   {\phy@FWoptions#1}{\phy@FWoptions#1[]}}
28 \onlypreamble\phy@FWoptions
29 \def\phy@FWoptions#1[#2]{\@ifnextchar[%]
30   {\phy@FWoptions@ns#1[##2]}{\phy@FWoptions@ns#1[##2]#3[]}}
31 \onlypreamble\phy@FWoptions
32 \def\phy@FWoptions@ns#1[#2]{%
33   \def\reserved@b##1{%
34     \ifx\@nnil##1\relax\else
35       \ifx\@nnil##1\@nnil\else
36         \noexpand\@onefilewithoptions{phy-##1}[\{\unexpanded{##2}\}][{##4}]%
37         \noexpand\@pkgextension
38       \fi
39       \expandafter\reserved@b
40     \fi}%
41   \edef\reserved@a{\zap@space#3 \@empty}%
42   \edef\reserved@a{\expandafter\reserved@b\reserved@a,\@nnil,}%
43   \reserved@a}
44 \onlypreamble\phy@FWoptions@ns
```

---

### 1.3 The (used to be) **common** module

The code below used to be the automatically-loaded **common** module, but now we load it together with **physics2**'s code. This change may bring better performance in Windows system.

Check if **unicode-math** loaded and (re)define the vert symbols. The `\relax`'s at the ends of `\vert` and `\Vert`'s definitions must not be removed. They are for `\ifx` to compare. **unicode-math** sets these symbols `\fam1`, `\symoperators` is equal to 1 in L<sup>A</sup>T<sub>E</sub>X 2 <sub>$\varepsilon$</sub>  kernel. Moreover, we make `\mid` as a delimiter but it may not work.

```
45 \AtBeginDocument{\ifcsname symrm\endcsname
46   \protected\def\|{\Udelimiter 0 \symoperators "2016 }%
47   \protected\def\vert{\Udelimiter 0 \symoperators "007C\relax}%
48   \protected\def\Vert{\Udelimiter 0 \symoperators "2016\relax}%
49   \protected\def\mid{\Udelimiter 3 \symoperators "007C }%
50 \fi}
51 \protected\def\Vert{\delim{"026B30D}\relax}
52 \protected\def\mid{\delim{"326A30C }
```

---

```
\delopen \delopen <left delimiter>
\delclose \delclose <right delimiter>
```

---

Actually in T<sub>E</sub>X, `\left` and `\right` will enclose the subformula as “inner”, but `\delopen` and `\delclose` will make the subformula an empty open node and a non-empty close node.

```

53 \DeclareRobustCommand\delopen{\mathopen{}\mathclose\bgroup\left}
54 \DeclareRobustCommand\delclose{\aftergroup\egroup\right}
55 % Extension to 2e kernel's or amsmath's bigggg commands.

\bBigg@ is a command from amsmath. The code below should update with amsmath together.

56 \ifdefined\bBigg@
57   \DeclareRobustCommand\biggg{\bBigg@{3}}
58   \DeclareRobustCommand\Biggg{\bBigg@{3.5}}
59 \else
60   \DeclareRobustCommand\biggg[1]{\leavevmode
61     {\hbox{$\left#1\vbox{to20.5\p@\right.\n@space$}}}}
62   \DeclareRobustCommand\Biggg[1]{\leavevmode
63     {\hbox{$\left#1\vbox{to23.5\p@\right.\n@space$}}}}
64   \AtBeginDocument{\ifdefined\bBigg@
65     \DeclareRobustCommand\biggg{\bBigg@{3}}%
66     \DeclareRobustCommand\Biggg{\bBigg@{3.5}}%
67   \fi}
68 \fi
69 \DeclareRobustCommand\bigggl{\mathopen\biggg}
70 \DeclareRobustCommand\bigggm{\mathrel\biggg}
71 \DeclareRobustCommand\bigggr{\mathclose\biggg}
72 \DeclareRobustCommand\Bigggl{\mathopen\Biggg}
73 \DeclareRobustCommand\Bigggm{\mathrel\Biggg}
74 \DeclareRobustCommand\Bigggr{\mathclose\Biggg}

```

\phy@mathvphantom \phy@mathvphantom {*math mode material*}

This command is just like \vphantom in L<sup>A</sup>T<sub>E</sub>X 2 <sub>$\varepsilon$</sub>  kernel but only works in math mode.

```

75 \def\phy@mathvphantom#1{\setbox\phy@tempboxa=\hbox{}%
76   \mathchoice
77     {\setbox\@tempboxa\hbox{$\displaystyle#1$}%
78      \ht\phy@tempboxa=\ht\@tempboxa
79      \dp\phy@tempboxa=\dp\@tempboxa
80      \box\phy@tempboxa}
81     {\setbox\@tempboxa\hbox{$\textstyle#1$}%
82      \ht\phy@tempboxa=\ht\@tempboxa
83      \dp\phy@tempboxa=\dp\@tempboxa
84      \box\phy@tempboxa}
85     {\setbox\@tempboxa\hbox{$\scriptstyle#1$}%
86      \ht\phy@tempboxa=\ht\@tempboxa
87      \dp\phy@tempboxa=\dp\@tempboxa
88      \box\phy@tempboxa}
89     {\setbox\@tempboxa\hbox{$\scriptscriptstyle#1$}%
90      \ht\phy@tempboxa=\ht\@tempboxa
91      \dp\phy@tempboxa=\dp\@tempboxa
92      \box\phy@tempboxa}%
93 }

```

## 1.4 The (used to be) **explsetup** module

Some common variables and functions for experimental L<sup>A</sup>T<sub>E</sub>X3 syntax.

```
94 <@@=phy>
```

```

95 \ExplSyntaxOn
96 \int_new:N \l__phy_tmpa_int
97 \int_new:N \l__phy_tmpb_int
98 \tl_new:N \l__phy_tmpa_tl
99 \tl_new:N \l__phy_tmpb_tl

The function that can gobble one token.

100 \cs_new:Npn \__phy_gobble_i:n #1 { }
101 \ExplSyntaxOff
102 <@=>
103 </package>

```

## File II

# Modules written in L<sup>A</sup>T<sub>E</sub>X 2 <sub>$\epsilon$</sub> syntax

### 1 The ab module

(\*gibberish)

This module is important but the code is hard to read. One of the motivations I manage **physics2** with **DocStrip** is that, when I tried to write a new module based on **ab** after 5 months when I maintained **physics2** the last time, I found that I could not understand the code I wrote at all! Therefore, it's significant to comment out the alien code in **ab**.

```

</gibberish>

1 <*ab>
2 \ProvidesFile{phy-ab.sty}
3 [2023/10/24 `ab' (autobraces) module of physics2]

```

If you don't know when to use `\phy@define@key`, `\phy@setkeys` and `\phy@processkeyopt` in a module, see ahead. In **ab**, the `tightbraces` option can control if the automatically-sized braces are tight or not. Do you remember `\delopen` and `\delclose`?

```

4 \phy@define@key{ab}{tightbraces}[true]{\def\@phy@abtight{\#1}}
5 \phy@setkeys{ab}{tightbraces=true}
6 \phy@processkeyopt{ab}

```

---

```

\phy@abopen \phy@abopen <left delimiter>
\phy@abclose \phy@abclose <right delimiter>

```

They are defined either `\delopen`, `\delclose` or `\left`, `\right`. If a module requires **ab**, these two commands are likely to be used.

```

7 \ifx\@phy@abtight\phy@true
8   \let\phy@abopen\delopen
9   \let\phy@abclose\delclose
10 \else
11   \let\phy@abopen\left
12   \let\phy@abclose\right
13 \fi

```

## 1.1 The implementation of \ab

This is the alienest part of `ab`. It's better to draw something rather than write boring comments. First let's take a look at `\ab`'s syntax. After `\ab` should be a pair of delimiters; take `()` as an example. Between `\ab` and `("")` can be a biggg command or star, or even nothing. `\ab` is defined as follows:

```
\ab ← begindef
    \phy@d@lx {mb} {ab}
enddef
```

where `ab` is the branch name of `\ab()`, and `mb` is the branch name of `\ab\big()` and `\ab*`. Then let's see the syntax of `\phy@d@lx`.

```
\phy@d@lx {\biggg or star branch name} {\automatic branch name} {#3}
```

Here exists an `#3`. `#3` is one token immediately following `\ab`, which can be `{ a biggg command or a star }` or a `"("`, under our assumption.

`\phy@d@lx` is defined as follows:

```
\phy@d@lx ← begindef (#1: biggg or star branch name, <mb>; #2:
    automatic branch name, <ab>; #3, the token after
    \ab)
    if #3 == biggg or #3 == star (↔ csname
        {phy@del\string#3} is defined) then
            let <next cs> = csname {phy@d@lx<mb>}
    else
        let <next cs> = csname {phy@d@lx<ab>}
    endif
    <next cs> #3
enddef
```

The condition should be true when `#3` is `\big` or `*`, and it should be false when `#3` is `"("`. Accordingly, in math mode,

```
\ab \big ( → \phy@d@lxmb \big (
\ab      ( → \phy@d@lxab      (
```

Then we meet two new commands — `\phy@d@lxmb` and `\phy@d@lxab`. Syntax is as follows.

```
\phy@d@lxmb <biggg or *> <left delimiter> <subformula> <right delimiter>
\phy@d@lxab           <left delimiter> <subformula> <right delimiter>
```

Notice that `ab` and `mb` in the above commands are names of `\ab`'s two branches — they are like namespaces. `\phy@d@lxmb` and `\phy@d@lxab` are defined by the following two lines:

```
\phy@d@l@genxm{mb}
\phy@d@l@genxa{ab}
```

\phy@d@l@genxm and \phy@d@l@genxa are defined as follows:

```

\phy@d@l@genxm ← begindef (#1: biggg or star branch name, ⟨mb⟩)
    \phy@d@lx⟨mb⟩ ← begindef (##1: biggg or star;
        ##2: left delimiter)
            \begingroup
                if ##1 == star then
                    ⟨temp⟩ ← \relax
                else
                    ⟨temp⟩ ← ##1
                endif
                csname {phy@⟨mb⟩@\string##2}
                    ⟨temp⟩ ##2
                % requires an \endgroup after the right delimiter
            
```

**enddef**

**enddef**

```

\phy@d@l@genxa ← begindef (#1: automatic branch name, ⟨ab⟩)
    \phy@d@lx⟨ab⟩ ← begindef (##1: left delimiter)
        csname {phy@⟨ab⟩@\string##1}
            ##1
    
```

**enddef**

**enddef**

So we can get

```

\ab \big ( → \begingroup csname {phy@mb@() \big (
\ab * ( → \begingroup csname {phy@mb@() \relax (
\ab ( → csname {phy@ab@() (

```

The csnames above (\phy@mb@() and \phy@ab@()) are generated with \phy@AB@gen.

```
\phy@AB@gen {⟨branch name⟩} ⟨left delimiter⟩ {⟨arg spec⟩} {⟨definition⟩}
```

If ⟨branch name⟩ is mb, {⟨arg spec⟩} should be mr(), where m is for biggg or star; If ⟨branch name⟩ is ab, {⟨arg spec⟩} should be r().

**Note:** The “(” in the example above must not be replaced by a subformula braced by a pair of {}.

---

```
\phy@AB@gen \phy@AB@gen {⟨branch name⟩} ⟨left delimiter⟩ {⟨arg spec⟩} {⟨definition⟩}
```

```

14 \def\phy@AB@gen#1#2{\expandafter\DeclareDocumentCommand\csname phy@#1@\string#2\endcsname}
15 \phy@AB@gen{ab}({r()}{\phy@abopen(#1\phy@abclose)}
16 \phy@AB@gen{ab}[{r[]}]{\phy@abopen[#1\phy@abclose]}
17 \phy@AB@gen{ab}\{{r\{\}}{\phy@abopen\{#1\phy@abclose\}}
18 \phy@AB@gen{ab}|{r||}{\phy@abopen|#1\phy@abclose|}
```

```

19 \phy@AB@gen{ab}\{|{r\\|}{\phy@abopen\\#1\phy@abclose\\|}
20 \phy@AB@gen{ab}<{r>}{\phy@abopen\\#1\phy@abclose>}
21 \phy@AB@gen{ab}\lbrace{r\lbrace\rbrace}{\phy@abopen\lbrace\\#1\phy@abclose\rbrace}
22 \phy@AB@gen{ab}\vert{r\vert\vert}{\phy@abopen\vert\\#1\phy@abclose\vert}
23 \phy@AB@gen{ab}\Vert{r\Vert\Vert}{\phy@abopen\Vert\\#1\phy@abclose\Vert}
24 \phy@AB@gen{ab}\langle{r\langle}{\phy@abopen\langle\\#1\phy@abclose\rangle}

\endgroup's in the end of the following definitions are corresponding to \begingroup's
in the definition of \phy@d@l@genxm.

25 \phy@AB@gen{mb}(\{mr()\}){\mathopen{#1}\mathclose{#1}\endgroup}
26 \phy@AB@gen{mb}[\{mr[]\}]{\mathopen{#1}\mathclose{#1}\endgroup}
27 \phy@AB@gen{mb}\{{\{mr\\}\}}{\mathopen{#1}\lbrace\\#2\mathclose{#1}\rbrace\endgroup}
28 \phy@AB@gen{mb}|{\{mr\\|\}}{\mathopen{#1}\vert\\#2\mathclose{#1}\vert\endgroup}
29 \phy@AB@gen{mb}\{|{\{mr\\|\\|}\}}{\mathopen{#1}\Vert\\#2\mathclose{#1}\Vert\endgroup}
30 \phy@AB@gen{mb}<{\{mr>\}}{\mathopen{#1}\langle\\#2\mathclose{#1}\rangle\endgroup}
31 \phy@AB@gen{mb}\lbrace{\{mr\lbrace\rbrace}{\mathopen{#1}\lbrace\\#2\mathclose{#1}\rbrace\endgroup}
32 \phy@AB@gen{mb}\vert{\{mr\vert\vert}{\mathopen{#1}\vert\\#2\mathclose{#1}\vert\endgroup}
33 \phy@AB@gen{mb}\Vert{\{mr\Vert\Vert}{\mathopen{#1}\Vert\\#2\mathclose{#1}\Vert\endgroup}
34 \phy@AB@gen{mb}\langle{\{mr\langle}{\mathopen{#1}\langle\\#2\mathclose{#1}\rangle\endgroup}

```

\phy@del\string. The syntax seems not important. These following lines seems only for \ifcsname to judge if the commands are defined.

```

35 \def\phy@del#1#2#3{\phy@abopen#1#3\phy@abclose#2}
36 \expandafter\def\csname phy@del\string*\endcsname#1#2#3{\mathopen{#1}\mathclose{#2}}
37 \expandafter\def\csname phy@del\string\big\endcsname#1#2#3{\bigl#1\big#3\bigr#2}
38 \expandafter\def\csname phy@del\string\Big\endcsname#1#2#3{\Bigl#1\Big#3\Bigr#2}
39 \expandafter\def\csname phy@del\string\bigg\endcsname#1#2#3{\biggl#1\big#3\biggr#2}
40 \expandafter\def\csname phy@del\string\Bigg\endcsname#1#2#3{\Biggl#1\Big#3\Biggr#2}
41 \expandafter\def\csname phy@del\string\biggg\endcsname#1#2#3{\bigggl#1\big#3\bigggr#2}
42 \expandafter\def\csname phy@del\string\Biggg\endcsname#1#2#3{\Bigggl#1\Big#3\Bigggr#2}

```

\phy@d@lx \phy@d@lx {\{biggg or star branch name\}} {\{automatic branch name\}} {\#3}

```

43 \def\phy@d@lx#1#2#3{%
44   \ifcsname phy@del\string#3\endcsname
45     \def\reserved@a{#1}% #3 is star or \biggg
46   \else
47     \def\reserved@a{#2}% #3 is delimiter
48   \fi
49   \csname phy@d@lx\reserved@a\endcsname#3}

```

\phy@d@l@genxm \phy@d@l@genxm {\{biggg or star branch name\}}
\phy@d@l@genxa \phy@d@l@genxa {\{automatic branch name\}}

```

50 \def\phy@d@l@genxm#1{%
51   \expandafter\def\csname phy@d@lx#1\endcsname##1##2{%
52     \begingroup % \endgroup is at the end of #4 of \phy@AB@gen
53     \ifx##1*\let\phy@tempa=\relax\else\let\phy@tempa=##1\fi
54     \csname phy@#1@string##2\endcsname\phy@tempa##2}
55 \def\phy@d@l@genxa#1{%
56   \expandafter\def\csname phy@d@lx#1\endcsname##1{%
57     \csname phy@#1@string##1\endcsname##1}}

```

---

```
\phy@d@lxmb \phy@d@lxmb {biggg or *} <left delimiter> <subformula> <right delimiter>
\phy@d@lxab \phy@d@lxab <left delimiter> <subformula> <right delimiter>

58 \phy@d@l@genxm{mb}
59 \phy@d@l@genxa{ab}
```

---

\ab The users' command \ab.

```
60 \DeclareRobustCommand\ab{\phy@d@lx{mb}{ab}}
```

## 1.2 \pab-like commands

This is so simple. No need to comment a lot.

---

```
\phy@d@l@geny \phy@d@l@geny <command> <left delimiter> <right delimiter>
```

This command used to define commands like \pab.

```
61 \def\phy@d@l@geny#1#2#3{%
62   \DeclareDocumentCommand#1{som}{%
63     \IfBooleanTF{##1}{%
64       {##2##3}%
65       {\IfValueTF{##2}{%
66         {\csname##2\endcsname##3\csname##2r\endcsname##3}%
67         {\phy@abopen##2##3\phy@abcclose##3}%
68       }%
69     }%
70   }%
71   \phy@d@l@geny\pab()
72   \phy@d@l@geny\bab[]
73   \phy@d@l@geny\Bab\lbrace\rbrace
74   \phy@d@l@geny\vab\vert\vert
75   \phy@d@l@geny\aab\langle\rangle
76   \phy@d@l@geny\Vab\Vert\Vert
77   \langle\ab\rangle}
```

## 2 The ab.braket module

```
1 <*ab.braket>
2 \ProvidesFile{phy-ab.braket.sty}
3 [2023/10/24 `ab.braket' module of physics2]
```

This module requires \phy@abopen and \phy@abcclose from **ab**. This module may have conflict with **braket**.

```
4 \phy@requiremodule{ab}
5 \ifdefined\phy@bra@@
6 \PackageWarning{physics2}{You cannot load `ab.braket' and `braket'
7 modules together.\MessageBreak Only `ab.braket' module works now.}
8 \fi
```

---

\bra \bra < <subformula> |

```

9 \phy@AB@gen{br.m}<{mr<|}{\mathopen#1\langle\mathclose#1\vert\endgroup}
10 \phy@AB@gen{br.a}<{r<|}{\phy@abopen\langle\mathopen#1\phy@abclose\vert}
11 \phy@d@l@genxm{br.m}
12 \phy@d@l@genxa{br.a}
13 \DeclareRobustCommand\bra{\phy@d@lx{br.m}{br.a}}


\underline{\ket} \ket | subformula >

14 \phy@AB@gen{kt.m}|{mr|>}{\mathopen#1\vert#2\mathclose#1\rangle\endgroup}
15 \phy@AB@gen{kt.a}|{r|>}{\phy@abopen\vert#1\phy@abclose\rangle}
16 \phy@d@l@genxm{kt.m}
17 \phy@d@l@genxa{kt.a}
18 \DeclareRobustCommand\ket{\phy@d@lx{kt.m}{kt.a}}


\underline{\braket} \braket < subformula 1 | subformula 2 [| subformula 3 ...] >

19 \begingroup
20 \catcode`\|=active
21 \gdef\phy@@mb@bk#1#2{\begingroup
22   \mathcode`|=8000\def|\egroup#1\vert\bgroup}%
23   \def\<{\mathrel{<}}\def\>{\mathrel{>}}%
24   \mathopen#1\langle\mathopen\bgroup#2\egroup\mathclose#1\rangle\endgroup}
25 \gdef\phy@@ab@bk#1{\begingroup
26   \mathcode`|=8000\def|\egroup\phy@abb@bkv\bgroup}%
27   \def\<{\mathrel{<}}\def\>{\mathrel{>}}%
28   \phy@abopen\langle\mathopen\bgroup#1\egroup\phy@abclose\rangle\endgroup}
29 \endgroup
30 \def\phy@abb@bkv{\middle\vert}
31 \phy@AB@gen{bk.m}<{mr|>}{\phy@@mb@bk#1#2}\endgroup}
32 \phy@AB@gen{bk.a}<{r|>}{\phy@ab@bk{#1}}
33 \phy@d@l@genxm{bk.m}
34 \phy@d@l@genxa{bk.a}
35 \DeclareRobustCommand\braket{\phy@d@lx{bk.m}{bk.a}}


\underline{\ketbra} \braket | subformula 1 > subformula 2 < subformula 3 |

36 \begingroup
37 \catcode`\<=active
38 \catcode`\>=active
39 \gdef\phy@@mb@kb#1#2{\begingroup
40   \mathcode`<=8000 \mathcode`>="8000%
41   \def\<{\mathopen\langle\mathopen\bgroup} \def\>{\mathclose\rangle\mathclose\bgroup}%
42   \def\<{\phy@abb@l}\def\>{\phy@abb@r}%
43   \mathopen#1\vert#2\mathclose#1\vert\endgroup}
44 \endgroup
45 \gdef\phy@@ab@kb#1#2<#3\phy@@end{\begingroup
46   \def\<{\phy@abb@l}\def\>{\phy@abb@r}%
47   \phy@abopen\vert\mathopen{\phy@mathvphantom{#3}}\#1\phy@abclose\rangle\mathclose#2%
48   \phy@abopen\langle\mathopen\mathclose{\phy@mathvphantom{#1}}\phy@abclose\vert
49 \endgroup}
50 \AtBeginDocument{\ifcsname symbf\endcsname
51   \def\phy@abb@l{\Umathchar 3 \symoperators "003C }%
```

```

52 \def\phy@abb@r{\Umathchar 3 \symoperators "003E }%
53 \fi}
54 \def\phy@abb@l{\mathchar"313C }
55 \def\phy@abb@r{\mathchar"313E }
56 \phy@AB@gen{kb.m}|{mr||}{\phy@mb@kb#1{#2}\endgroup}%
57 \phy@AB@gen{kb.a}|{r||}{\phy@ab@kb#1\phy@end}%
58 \phy@d@l@genxm{kb.m}%
59 \phy@d@l@genxa{kb.a}%
60 \DeclareRobustCommand\ketbra{\phy@d@lx{kb.m}{kb.a}}%
61 \langle ab.braket\rangle

```

### 3 The **braket** module

```

1 <*>braket>
2 \ProvidesFile{phy-braket.sty}
3 [2023/10/24 `braket' module of physics2]

```

This module requires `\phy@abopen` and `\phy@abcclose` from **ab**. This module may have conflict with **ab.braket**.

```

4 \phy@requiremodule{ab}
5 \ifdef{\phy@abb@bkv}
6 \PackageWarning{physics2}{You cannot load `ab.braket' and `braket'
7 modules together.\MessageBreak Only `braket' module works now.}%
8 \fi

```

---

```

\bra \bra * [<biggg>] {<subformula>}

9 \DeclareDocumentCommand\bra{ s o m }{%
10 \IfBooleanTF{#1}{%
11   {\mathopen\langle#3\mathclose\rvert}%
12 }{\IfValueTF{#2}{%
13   {\csname#21\endcsname\langle#3\csname#2r\endcsname\rvert}%
14   {\phy@abopen\langle#3\phy@abcclose\rvert}}%
15 }%
16 }

```

---

```

\ket \ket * [<biggg>] {<subformula>}

17 \DeclareDocumentCommand\ket{ s o m }{%
18 \IfBooleanTF{#1}{%
19   {\mathopen\lvert#3\mathclose\rangle}%
20 }{\IfValueTF{#2}{%
21   {\csname#21\endcsname\lvert#3\csname#2r\endcsname\rangle}%
22   {\phy@abopen\lvert#3\phy@abcclose\rangle}}%
23 }%
24 }

```

---

```

\braket \braket * [<biggg>, n \in {1,2,3}] {<subformula 1>} ... {<subformula n>}

25 \DeclareDocumentCommand\braket{ s O{} }{%
26 \IfBooleanTF{#1}{%
27 }{%
28   \gdef\@phy@bk@argnum{ii}%
}

```

```

29   \phy@bk@doopt{#2}%
30   \gdef\@phy@bk@l{\mathopen}%
31   \gdef\@phy@bk@m{\mathord}%
32   \gdef\@phy@bk@r{\mathclose}%
33 }%
34 {%
35   \gdef\@phy@bk@argnum{ii}%
36   \gdef\@phy@bk@l{\phy@abopen}%
37   \gdef\@phy@bk@m{\middle}%
38   \gdef\@phy@bk@r{\phy@abcclose}%
39   \phy@bk@doopt{#2}%
40 }%
41 \csname phy@bk@in@ \@phy@bk@argnum \endcsname%
42 }



---


\nphy@bk@in@i \phy@bk@in@<n.roman> {\<subformula 1>} ... {\<subformula n>}
\nphy@bk@in@ii <n.roman> is n in roman lowercase, where n  $\in \{1, 2, 3\}.
\nphy@bk@in@iii

43 \def\phy@bk@in@i#1{%
44   \csname\@phy@bk@l\endcsname\langle{#1}%
45   \csname\@phy@bk@r\endcsname\rangle}%
46 \def\phy@bk@in@ii#1#2{%
47   \csname\@phy@bk@l\endcsname\langle{#1}%
48   \csname\@phy@bk@m\endcsname\vert{#2}%
49   \csname\@phy@bk@r\endcsname\rangle}%
50 \def\phy@bk@in@iii#1#2#3{%
51   \csname\@phy@bk@l\endcsname\langle{#1}%
52   \csname\@phy@bk@m\endcsname\vert{#2}%
53   \csname\@phy@bk@m\endcsname\vert{#3}%
54   \csname\@phy@bk@r\endcsname\rangle}



---


\nphy@bk@doopt \phy@bk@doopt {\<clist>}
\nphy@bk@do@pt Parse the optional argument of \braket. This will add 3 entries to hash.
55 \def\@phy@bk@do@pt#1,{\ifx#1\relax\@empty\else
56   \edef\reserved@a{\zap@space#1 \@empty}%
57   \ifx\reserved@a\@empty\else
58     \ifcsname phy@del\expandafter\string\csname\reserved@a\endcsname\endcsname
59       \xdef\@phy@bk@l{\reserved@a 1}%
60       \xdef\@phy@bk@m{\reserved@a} but not m (m stands for \mathrel)
61       \xdef\@phy@bk@r{\reserved@a r}%
62   \else
63     \ifnum\reserved@a>3%
64       \PackageError{physics2}{\string\braket\space can only take 3
65         mandatory arguments at most}{Check if you had written a number
66         more than 3 in the [optional] argument.}%
67     \fi
68     \xdef\@phy@bk@argnum{\romannumeral\reserved@a}%
69   \fi
70 \fi
71 \expandafter\@phy@bk@do@pt\fi}
72 \def\phy@bk@doopt#1{\@phy@bk@do@pt#1,\relax,}$ 
```

---

```

\ketbra \ketbra * [<biggg>] {<subformula 1>} [<between 1 and 2>] {<subformula 2>}

73 \DeclareDocumentCommand\ketbra{ s o m O{} m }{%
74   \IfBooleanTF{#1}{%
75     \mathopen\vert#3\mathclose\rangle\mathopen\langle#5\mathclose\vert\%%
76     \IfValueTF{#2}{%
77       \csname#21\endcsname\vert#3\csname#2r\endcsname\rangle\mathopen\langle#4\%
78       \csname#21\endcsname\langle#5\csname#2r\endcsname\vert\%%
79       \begingroup
80         \phy@abopen\vert\mathopen{\phy@mathvphantom{#5}}\#3\phy@abclose\rangle\mathopen\langle#4\%
81         \phy@abopen\langle#5\mathclose{\phy@mathvphantom{#3}}\phy@abclose\vert\%
82       \endgroup\%
83     }%
84   }%
85 }
```

## 4 The **doubleprod** module

```

1  {*doubleprod}
2  \ProvidesFile{phy-doubleprod.sty}
3  [2023/10/24 `doubleprod' (vertically stacked binary operators) module of physics2]
```

Boolean options.

```

4  \phy@define@key{doubleprod}{crosssymbol}{\def\@phy@dbl@c{#1}}
5  \phy@define@key{doubleprod}{dotsymbol}{\def\@phy@dbl@d{#1}}
6  \phy@define@key{doubleprod}{crossscale}{\def\@phy@dbl@sc{#1}}
7  \phy@define@key{doubleprod}{dotscale}{\def\@phy@dbl@sd{#1}}
8  \phy@define@key{doubleprod}{crossopenup}{\def\@phy@dbl@oc{#1}}
9  \phy@define@key{doubleprod}{dotopenup}{\def\@phy@dbl@od{#1}}
10 \phy@setkeys{doubleprod}{crosssymbol=\times,dotsymbol=\ldotp,
11   crossscale=0.8,dotscale=1,crossopenup=.02,dotopenup=.2}
12 \phy@processkeyopt{doubleprod}
13 \def\phy@dbl@gen#1#2#3#4{%
14   \ DeclareRobustCommand#1{\mathbin{\vcenter{\baselineskip\z@skip\%
15     \lineskip#4\phy@dblcurr@size\%
16     \setbox\@tempboxa=\hbox{\fontsize{#2}\phy@dblcurr@size}\z@\#3\%%
17     \copy\@tempboxa\box\@tempboxa}}}}
18 \def\phy@dblcurr@size{\dimexpr\f@size pt\relax}
19 \phy@dbl@gen\doublecross\@phy@dbl@sc\@phy@dbl@c\@phy@dbl@oc
20 \phy@dbl@gen\doubledot\@phy@dbl@sd\@phy@dbl@d\@phy@dbl@od
21 }/doubleprod}
```

## File III

# Modules written in L<sup>A</sup>T<sub>E</sub>X3 syntax

We use `phy` as the namespace for `physics2` modules.

```
1  <@@=phy>
```

## 1 The **diagmat** module

```
1  {*diagmat}
```

```

2 \ProvidesExplFile{phy-diagmat.sty}{2024/01/10}{}
3   {'diagmat' module of physics2}
4 \RequirePackage { amsmath }
5 \phy@define@key { diagmat } { empty } [ 0 ] { \tl_gset:Nn \l__phy_mat_empty_tl { #1 } }

```

This module requires some new variables.

```

6 \clist_new:N \l__phy_mat_diag_clist
7 \int_new:N \l__phy_mat_dim_int
8 \tl_new:N \l__phy_mat_line_tl
9 \tl_new:N \l__phy_diagmat_tl
10 \tl_new:N \l__phy_mat_empty_tl
11 \tl_gset:Nn \l__phy_mat_empty_tl { 0 }
12 \phy@processkeyopt { diagmat }
13 \keys_define:nn { phy/diagmat }
14 {
15   empty .tl_set:N = \l__phy_mat_empty_tl ,
16 }

```

---

\diagmat \langle delimiter type\rangle diagmat [\langle key-val list\rangle] {\langle diagonal\rangle}

---

```

17 \DeclareDocumentCommand \diagmat { O{} m }
18   { \__phy_diagmat_type:nnn { } { #1 } { #2 } }
19 \DeclareDocumentCommand \pdiagmat { O{} m }
20   { \__phy_diagmat_type:nnp { p } { #1 } { #2 } }
21 \DeclareDocumentCommand \bdiagmat { O{} m }
22   { \__phy_diagmat_type:nmb { b } { #1 } { #2 } }
23 \DeclareDocumentCommand \Bdiagmat { O{} m }
24   { \__phy_diagmat_type:nmm { B } { #1 } { #2 } }
25 \DeclareDocumentCommand \vdiagmat { O{} m }
26   { \__phy_diagmat_type:nnn { v } { #1 } { #2 } }
27 \DeclareDocumentCommand \Vdiagmat { O{} m }
28   { \__phy_diagmat_type:nnn { V } { #1 } { #2 } }

```

---

\\_\_phy\_diagmat\_type:nnn \\_\_phy\_diagmat\_type:nnn {\langle delimiter type\rangle} {\langle key-val list\rangle} {\langle diagonal\rangle}

---

```

29 \cs_new:Npn \__phy_diagmat_type:nnn #1#2#3
30 {
31   \group_begin:
32   \clist_set:Nn \l__phy_mat_diag_clist { #3 }
33   \int_set:Nn \l__phy_mat_dim_int { \clist_count:N \l__phy_mat_diag_clist }
34   \int_compare:nNnT { \l__phy_mat_dim_int } > { \value { MaxMatrixCols } }
35     { \setcounter { MaxMatrixCols } { \l__phy_mat_dim_int } }
36   \keys_set:nn { phy/diagmat } { #2 }
37   \tl_gclear:N \l__phy_diagmat_tl
38   \int_step_inline:nnn { 1 } { \l__phy_mat_dim_int }
39   {
40     \int_step_inline:nnn { 1 } { \l__phy_mat_dim_int }
41     {
42       \int_compare:nNnTF { ##1 } = { #####1 }
43       {
44         \clist_gpop:NN \l__phy_mat_diag_clist \l__phy_tmpa_tl
45         \tl_if_empty:NTF \l__phy_tmpa_tl
46           { \tl_gput_right:Nn \l__phy_mat_line_tl { \l__phy_mat_empty_tl } }

```

Maybe it's better to use `\expandafter\scantokens\expandafter{\l_@@_tmpa_t1}` in the next line.

```

47          { \tl_gput_right:No \l_@_phy_mat_line_tl { \l_@_phy_tmpa_t1 } }
48      }
49      { \tl_gput_right:Nn \l_@_phy_mat_line_tl { \l_@_phy_mat_empty_tl } }
50      \int_compare:nNnTF { #####1 } = { \l_@_phy_mat_dim_int }
51      {
52          \tl_gput_right:Nn \l_@_phy_mat_line_tl { \\ }
53      }
54      {
55          \tl_gput_right:Nn \l_@_phy_mat_line_tl { & }
56      }
57  }
58  \tl_gput_right:No \l_@_phy_diagmat_tl { \l_@_phy_mat_line_tl }
59  \tl_gclear:N \l_@_phy_mat_line_tl
60  }
61  \begin { #1 matrix }
62  \tl_use:N \l_@_phy_diagmat_tl
63  \end { #1 matrix }
64  \group_end:
65 }
66 
```

## 2 The **xmat** module

```

1  {*xmat}
2  \ProvidesExplFile{phy-xmat.sty}{2023/10/24}{}
3  {`xmat' module of physics2}
4  \RequirePackage { amsmath }
5  \phy@define@key { xmat } { showtop }
6  { \int_gset:Nn \l_@_phy_xmat_showtop_int { #1 } }
7  \phy@define@key { xmat } { showleft }
8  { \int_gset:Nn \l_@_phy_xmat_showleft_int { #1 } }

```

This module requires some new variables.

```

9  \bool_new:N \l_@_phy_xmat_extra_vdots_bool
10 \bool_new:N \l_@_phy_xmat_extra_cdots_bool
11 \int_new:N \l_@_phy_xmat_showtop_int
12 \int_new:N \l_@_phy_xmat_showleft_int
13 \tl_new:N \l_@_phy_xmat_tl
14 \int_gset:Nn \l_@_phy_xmat_showtop_int { \value { MaxMatrixCols } - 2 }
15 \int_gset:Nn \l_@_phy_xmat_showleft_int { \value { MaxMatrixCols } - 2 }
16 \cs_new:Npn \__phy_xmat_entry_format:nnn #1#2#3
17  {
18      #1 \c_math_subscript_token { #2 #3 }
19  }
20 \phy@processkeyopt { xmat }
21 \DeclareDocumentCommand \xmat { O{} m m m }
22 { \__phy_xmat_type:nnnnn { } { #1 } { #2 } { #3 } { #4 } }
23 \DeclareDocumentCommand \pxmat { O{} m m m }
24 { \__phy_xmat_type:nnnnn { p } { #1 } { #2 } { #3 } { #4 } }
25 \DeclareDocumentCommand \bxmat { O{} m m m }
26 { \__phy_xmat_type:nnnnn { b } { #1 } { #2 } { #3 } { #4 } }
27 \DeclareDocumentCommand \Bxmat { O{} m m m }
28 { \__phy_xmat_type:nnnnn { B } { #1 } { #2 } { #3 } { #4 } }

```

```

29 \DeclareDocumentCommand \vxmat { O{} m m m }
30 { \_phy_xmat_type:nnnnn { v } { #1 } { #2 } { #3 } { #4 } }
31 \DeclareDocumentCommand \Vxmat { O{} m m m }
32 { \_phy_xmat_type:nnnnn { V } { #1 } { #2 } { #3 } { #4 } }
33 \keys_define:nn { phy/xmat }
34 {
35   format .cs_set:Np = \_phy_xmat_entry_format:nnn #1#2#3 ,
36   showtop .int_set:N = \l_phy_xmat_showtop_int ,
37   showleft.int_set:N = \l_phy_xmat_showleft_int ,
38 }

```

---

\\_phy\_if\_digits\_only\_p:n ★ \\_phy\_if\_digits\_only:nTF {{*token list*}} {{*true code*}} {{*false code*}}

---

\\_phy\_if\_digits\_only:nTF \*

Use L<sup>A</sup>T<sub>E</sub>X3 regular expression to tell if *<token list>* (the numbers of rows or columns) contain digits only.

```

39 \prg_new_conditional:Npnn \_phy_if_digits_only:n #1 { TF }
40 {
41   \regex_match:nnTF { \A [[:digit:]]* \Z } { #1 }
42   { \prg_return_true: } { \prg_return_false: }
43 }

```

---

\\_phy\_xmat\_type:nnnn \\_phy\_xmat\_type:nnnnn {{*delimiter type*}} {{*key-val list*}} {{*common entry*}} {{*row number*}} {{*column number*}}

```

44 \cs_new:Npn \_phy_xmat_type:nnnnn #1#2#3#4#5
45 {
46   \group_begin:
47   \tl_gclear:N \l_phy_xmat_tl
48   \keys_set:nn { phy/xmat } { #2 } %
49   \_phy_if_digits_only:nTF { #4 }
50   {
51     \int_compare:nNnTF { #4 } < { \l_phy_xmat_showtop_int + 1 }
52     {
53       \int_set:Nn \l_phy_xmat_showtop_int { #4 }
54       \bool_set_false:N \l_phy_xmat_extra_vdots_bool
55     }
56     {
57       \bool_set_true:N \l_phy_xmat_extra_vdots_bool
58     }
59   }
60   {
61     \bool_set_true:N \l_phy_xmat_extra_vdots_bool
62   }
63   \_phy_if_digits_only:nTF { #5 }
64   {
65     \int_compare:nNnTF { #5 } < { \l_phy_xmat_showleft_int + 1 }
66     {
67       \int_set:Nn \l_phy_xmat_showleft_int { #5 }
68       \bool_set_false:N \l_phy_xmat_extra_cdots_bool
69     }
70     {
71       \bool_set_true:N \l_phy_xmat_extra_cdots_bool

```

```

72         }
73     }
74     {
75         \bool_set_true:N \l__phy_xmat_extra_cdots_bool
76     }
77 \int_step_inline:nn { \l__phy_xmat_showtop_int }
78     {
79         \tl_put_right:Nn \l__phy_xmat_tl
80             { \_phy_xmat_entry_format:nnn { #3 } { ##1 } { 1 } }
81         \int_step_inline:nnn { 2 } { \l__phy_xmat_showleft_int }
82             {
83                 \tl_put_right:Nn \l__phy_xmat_tl
84                     { & \_phy_xmat_entry_format:nnn { #3 } { ##1 } { #####1 } }
85             }
86         \bool_if:NT \l__phy_xmat_extra_cdots_bool
87             {
88                 \tl_put_right:Nn \l__phy_xmat_tl
89                     { & \cdots & \_phy_xmat_entry_format:nnn { #3 } { ##1 } { #5 } }
90             }
91         \tl_put_right:Nn \l__phy_xmat_tl { \\ }
92     }
93 \bool_if:NT \l__phy_xmat_extra_vdots_bool
94     {
95         \tl_put_right:Nn \l__phy_xmat_tl { \vdots }
96         \prg_replicate:nn { \l__phy_xmat_showleft_int - 1 }
97             {
98                 \tl_put_right:Nn \l__phy_xmat_tl { & \vdots }
99             }
100        % Add \ddots if vdots_bool and cdots_bool be true simultaneously.
101        \bool_if:NT \l__phy_xmat_extra_cdots_bool
102            {
103                \tl_put_right:Nn \l__phy_xmat_tl { & \ddots & \vdots }
104            } % else relax
105        \tl_put_right:Nn \l__phy_xmat_tl { \\ }
106        % The last row.
107        \tl_put_right:Nn \l__phy_xmat_tl
108            { \_phy_xmat_entry_format:nnn { #3 } { #4 } { 1 } }
109        \int_step_inline:nnn { 2 } { \l__phy_xmat_showleft_int }
110            {
111                \tl_put_right:Nn \l__phy_xmat_tl
112                    { & \_phy_xmat_entry_format:nnn { #3 } { #4 } { ##1 } }
113            }
114        \bool_if:NT \l__phy_xmat_extra_cdots_bool
115            {
116                \tl_put_right:Nn \l__phy_xmat_tl
117                    { & \cdots & \_phy_xmat_entry_format:nnn { #3 } { #4 } { #5 } }
118            }
119        } % else relax
120        \begin{#1 matrix}
121            \tl_use:N \l__phy_xmat_tl
122        \end{#1 matrix}
123        \group_end:
124    }
125 
```

This part ends here.

126 ⟨@@=⟩

## File IV

# Legacy modules written in L<sup>A</sup>T<sub>E</sub>X 2<sub>ε</sub> syntax

## 1 The ab.legacy module

```
1  {*ab.legacy}
2  \ProvidesFile{phy-ab.legacy.sty}
3    [2023/10/24 `ab.legacy' module of physics2]

Requires ab's tight option.
4  \phy@requiremodule{ab}
5  \phy@define@key{ab.legacy}{order}[\mathcal{O}]{\def\phy@ab@ordersym{\#1}}
6  \phy@setkeys{ab.legacy}{order}
7  \phy@processkeyopt{ab.legacy}
8  \phy@d@l@geny\abs\vert\vert
9  \phy@d@l@geny\norm\Vert\Vert
10 \DeclareDocumentCommand\order{som}{%
11   \phy@ab@ordersym
12   \IfBooleanTF{\#1}
13   {(\#3)}
14   {\IfValueTF{\#2}
15    {\csname#2\endcsname(\#3\csname#2r\endcsname)}
16    {\phy@abopen(\#3\phy@abclose)}%
17   }%
18 }
19 \phy@d@l@geny\eval.\vert
20 \phy@d@l@geny\peval(\vert
21 \phy@d@l@geny\beval[\vert
22 ⟨/ab.legacy⟩
```

## 2 The nabla.legacy module

```
1  {*nabla.legacy}
2  \ProvidesFile{phy-nabla.legacy.sty}
3    [2023/10/24 `nabla.legacy' module of physics2]
4  \phy@requiremodule{ab}

Requires fixdif version 2.x.
5  \RequirePackage{fixdif}[2023/01/31]
6  \letdif\phy@nl@nabla{nabla}
7  \AtBeginDocument{\ifcsname div\endcsname\let\divsymbol\div\fi
8  \ DeclareRobustCommand\grad{\phy@nl@nabla\ab}%
9  \ DeclareRobustCommand\div{\phy@nl@nabla\cdot\ab}%
10 \ DeclareRobustCommand\curl{\phy@nl@nabla\times\ab}%
11 \ DeclareRobustCommand\laplacian{\phy@nl@nabla^2\ab}%
12 }
13 ⟨/nabla.legacy⟩
```

### 3 The **op.legacy** module

```
1 <*op.legacy>
2 \ProvidesFile{phy-op.legacy.sty}
3   [2023/10/24 `op.legacy' module of physics2]
4 \phy@define@key{op.lega}{ReIm}[true]{\def\phy@reserveda{\#1}}
5 \phy@define@key{op.lega}{PV}{\def\@phy@oplega@PV{\#1}}
6 \phy@define@key{op.lega}{pv}{\def\@phy@oplega@pv{\#1}}
7 \phy@setkeys{op.lega}{PV=\mathcal{P},pv={p.v.},ReIm=true}
8 \phy@processkeyopt{ab}
9 \DeclareRobustCommand\asin{\mathop{\operator@font asin}\nolimits}
10 \DeclareRobustCommand\acos{\mathop{\operator@font acos}\nolimits}
11 \DeclareRobustCommand\atan{\mathop{\operator@font atan}\nolimits}
12 \DeclareRobustCommand\acsc{\mathop{\operator@font acsc}\nolimits}
13 \DeclareRobustCommand\asec{\mathop{\operator@font asec}\nolimits}
14 \DeclareRobustCommand\acot{\mathop{\operator@font acot}\nolimits}
15 \DeclareRobustCommand\Tr{\mathop{\operator@font Tr}\nolimits}
16 \DeclareRobustCommand\tr{\mathop{\operator@font tr}\nolimits}
17 \DeclareRobustCommand\rank{\mathop{\operator@font rank}\nolimits}
18 \DeclareRobustCommand\erf{\mathop{\operator@font erf}\nolimits}
19 \DeclareRobustCommand\Res{\mathop{\operator@font Res}\nolimits}
20 \DeclareRobustCommand\res{\mathop{\operator@font res}\nolimits}
21 \DeclareRobustCommand\PV{\mathord{\@phy@oplega@PV}}
22 \DeclareRobustCommand\pv{\mathop{\operator@font @phy@oplega@pv}\nolimits}
```

Restore  $\text{Re}$  and  $\text{Im}$  in  $\text{\Resymbol}$  and  $\text{\Imsymbol}$ . The  $\text{\AtBeginDocument}$  hook is used for the compatibility of **unicode-math**.

```
23 \ifx\phy@reserveda\phy@true
24 \AtBeginDocument{%
25   \let\Resymbol\Re%
26   \let\Imsymbol\Im%
27   \DeclareRobustCommand\Re{\mathop{\operator@font Re}\nolimits}%
28   \DeclareRobustCommand\Im{\mathop{\operator@font Im}\nolimits}%
29 }
30 \fi
31 </op.legacy>
```

### 4 The **qtext.legacy** module

This module is written for the compatibility with the bad commands provided by **physics** only. The commands in this module should NEVER be used!

```
1 <*qtext.legacy>
2 \ProvidesFile{phy-qtext.legacy.sty}
3   [2023/10/24 `qtext.legacy' module of physics2.sty]
4 \RequirePackage{amstext}
5 \def\phy@qtext@#1#2{\text{\#2}\quad}
6 \DeclareRobustCommand\qqtext{\@ifstar{\phy@qtext@{}{\phy@qtext@\quad}}{\phy@qtext@\quad}}
7 \DeclareRobustCommand\qq{\qqtext}
8 \DeclareRobustCommand\qcomma{,\quad}
9 \DeclareRobustCommand\qc{\qcomma}
10 \DeclareRobustCommand\qcc{\@ifstar{\phy@qtext@{}{c.c}}{\phy@qtext@\quad{c.c}}}
11 \def\phy@qtext@lega@gen@#1{%
12   \expandafter\ DeclareRobustCommand\csname q#1\endcsname{%
13     \@ifstar{\phy@qtext@{}{\#1}}{\phy@qtext@\quad{\#1}}}}
```

```

14 \phy@qtext@lega@gen@if}
15 \phy@qtext@lega@gen{then}
16 \phy@qtext@lega@gen{else}
17 \phy@qtext@lega@gen{otherwise}
18 \phy@qtext@lega@gen{unless}
19 \phy@qtext@lega@gen{give}
20 \phy@qtext@lega@gen{using}
21 \phy@qtext@lega@gen{unless}
22 \phy@qtext@lega@gen{assume}
23 \phy@qtext@lega@gen{since}
24 \phy@qtext@lega@gen{let}
25 \phy@qtext@lega@gen{for}
26 \phy@qtext@lega@gen{all}
27 \phy@qtext@lega@gen{even}
28 \phy@qtext@lega@gen{odd}
29 \phy@qtext@lega@gen{integer}
30 \phy@qtext@lega@gen{and}
31 \phy@qtext@lega@gen{or}
32 \phy@qtext@lega@gen{as}
33 \phy@qtext@lega@gen{in}
34 
```

# File V

## Legacy modules written in L<sup>A</sup>T<sub>E</sub>X3 syntax

1 <@@=phy>

### 1 The **bm-um.legacy** module

```

1 <*bm-um.legacy>
2 \ProvidesExplFile{phy-bm-um.legacy.sty}{2023/10/24}{}%
3 {`bm-um.legacy' module of physics2}
4 \AtBeginDocument
5 {
6     \cs_if_exist:cF { symbf }
7     {
8         \PackageError { physics2 }
9         {
10             The ~ `bm-um.legacy' ~ module ~ requires ~
11             `unicode-math' ~ package
12         }
13         {
14             Have ~ you ~ used ~ `unicode-math' ~
15             in ~ the ~ preamble?
16         }
17     }
18 }
19 \DeclareDocumentCommand \bm { m }
20 {

```

```

21   \mode_if_math:TF
22   {
23     \tl_if_head_eq_catcode:nNTF { #1 } A
24     {
25       \symbfit { #1 }
26     }
27     {
28       \symbf { #1 }
29     }
30   }
31 {
32   \PackageError { physics2 }
33   {
34     The ~ \string\bm\space command ~ should ~ be ~
35     used ~ in ~ math ~ mode ~ only. \MessageBreak
36     This ~ is ~ an ~ error ~ from ~ `bm-um.legacy' ~ module
37   }
38   {
39     Check ~ if ~ any ~ ` \string\bm' ~ is ~ out ~
40     of ~ math ~ mode.
41   }
42 }
43 }
44 /bm-um.legacy

```

This part ends here.

45 <@@=