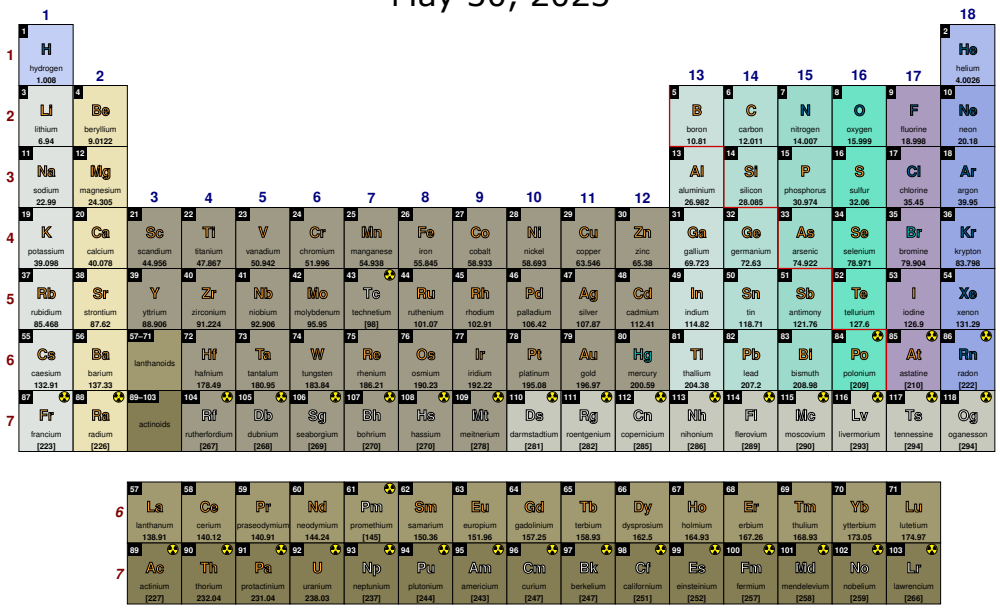
Manual for pgf-PeriodicTable 2.0.1

Hugo Gomes
hugo.parelho@gmail.com

May 30, 2023



\pgfPT[show title=false,show legend=false]

Abstract

The purpose of this package is to provide the Periodic Table of Elements in a simple way. It relies on pgf/TikZ to offer a full or partial periodic table with a variety of options and displaying the desired data. The data available, from all the actual 118 elements, is: atomic number, element name, chemical symbol, relative atomic mass, standard relative atomic mass, radioactivity, atomic radius (empirycal), covalente radius, ionic radius, first ionization energy, electronegativity (Pauling), electroaffinity, oxidation states, melting point (in Kelvin and Celsius degrees), boiling point (in Kelvin and Celsius degrees), electron distribuition, electronic configuration (increasing n and increasing $n+\ell$), density, specific heat capacity, thermal condutivity, lattice structure, lattice constants (a, b, c and c/a ratio), discovery year, discovery country and visible range spectral lines. It is possible to get the Periodic Table in six languages: English, French, German, Portuguese (from Portugal and from Brazil), Spanish and Italian.

Contents

Installation and usage 1

The data 2

The commands 3

\pgfPT . . . 3
\pgfPTstyle[options list] . . . 4
\pgfPTresetstyle . . . 5
\pgfPTbuildcell(nrows,ncolumns)[entries] . . . 6
\pgfPTresetcell . . . 6
\pgfPTbuildcellstyle{name}(nrows,ncolumns)[entries] . . . 6
\pgfPTpreviewcell . . . 6
\pgfPTpreviewcellstyle{name} . . . 7
\pgfPTnewColorScheme{name}{color list} . . . 7
\pgfPTnewZlist{name} . . . 9
\pgfPTsetLanguage{language flag} . . . 10

Options for \pgfPT: creating a «Periodic Table» 11

Periodic Table options: keys, styles and *pseudo styles* . . . 11
General layout . . . 11
Z list . . . 11
cell width . . . 12
cell height . . . 12
cell size . . . 12
cell line width . . . 13
cell line color . . . 13
cell style . . . 13
cell . . . 14
font . . . 14
back color scheme . . . 15
back color . . . 16
csSolid . . . 17
csSoft . . . 18
csJmol . . . 18
csCPK . . . 19
csRasmol . . . 19
csRasmolNew . . . 20
csWikipedia . . . 20
csMNM . . . 21
csPS . . . 21
csRadio . . . 22
csBlocks . . . 22
background . . . 23
IUPAC . . . 23
show label LaAc . . . 25
label LaAc font . . . 26
languages . . . 26
other languages font . . . 27
other languages color . . . 28
other lang . . . 28
show MNM line . . . 28
MNM line color . . . 29
MNM line width . . . 30
MNM . . . 30

Title and Legend 31
show title 31
title font 32
title color 32
title 32
show legend 33
legend acronyms 33
legend box 34
legend back color 35
legend radio color 35
legend CS color 36
legend Z color 37
show legend pins 38
legend pins 38
show extra legend 39
extra legend 39
legend 40
Periods and Groups 41
show period numbers 41
show group numbers 41
period label color 42
group label color 42
label font 43
per 43
gr 43
per+gr 44
Blocks and Families 44
show blocks 44
blocks font 46
s block color 46
s block font color 46
s block line width 46
p block color 46
p block font color 46
p block line width 46
d block color 46
d block font color 46
d block line width 46
f block color 46
f block font color 46
f block line width 47
blocks font color 47
blocks line width 47
blocks 48
show families 49
families font 50
r family color 50
r family font color 50
r family line width 51
tm family color 51
tm family font color 51
tm family line width 51
itm family color 51
itm family font color 51
itm family line width 51

➳ families font color . . . 51
➳ families line width . . . 51
➳ families . . . 52
➠ Periodic variations . . . 54
⇝ show periodic variations . . . 54
⇝ varR color . . . 55
⇝ varR font . . . 55
⇝ varR font color . . . 56
⇝ varEi color . . . 56
⇝ varEi font . . . 56
⇝ varEi font color . . . 56
⇝ vareaff color . . . 56
⇝ vareaff font . . . 56
⇝ vareaff font color . . . 56
➳ var font . . . 56
➳ var color . . . 57
➳ varR . . . 57
➳ varEi . . . 58
➳ vareaff . . . 59
➠ Dark mode . . . 60
➳ dark mode . . . 60
➠ *Exercise layout* . . . 60
⇝ only cells . . . 60
⇝ only cells plus Z . . . 61
⇝ only cells with periods and group numbers . . . 62
⇝ only cells with periods and group numbers plus Z . . . 63
⇝ Z exercise list . . . 63
⇝ exercise list in capitals . . . 64
⇝ exercise list color . . . 64
⇝ exercise list font . . . 64
➳ cells+Z . . . 65
➳ cells+p+g . . . 65
➳ cells+p+g+Z . . . 66
➳ exnocaps . . . 66
➳ exColor . . . 66
➳ exFont . . . 67
➳ ex . . . 67
✠ Cell contents options: keys, styles and *pseudo styles* . . . 67
➠ The atomic number . . . 67
⇝ Z backcolor . . . 67
⇝ Z color . . . 68
⇝ Z font . . . 68
⇝ Z use box width . . . 68
⇝ Z align . . . 68
⇝ Z padding . . . 69
➳ Z box . . . 69
➳ Z . . . 69
➠ The chemical symbol . . . 70
⇝ CS solid . . . 70
⇝ CS liquid . . . 70
⇝ CS gas . . . 71
⇝ CS synt . . . 71
➳ CS all . . . 71
⇝ CS font . . . 72
⇝ CS render mode . . . 72

⇝ CS outline color . . . 73
⇝ CS outline width . . . 73
➻ CS . . . 74
➠ The name . . . 74
⇝ name color . . . 74
⇝ name font . . . 74
⇝ name align . . . 75
⇝ capitalize element names . . . 75
➻ name . . . 75
➻ Name . . . 76
➻ NAME . . . 76
➠ The atomic weight . . . 76
⇝ Ar color . . . 76
⇝ Ar font . . . 77
⇝ Ar label . . . 77
⇝ Ar precision . . . 77
➻ Ar . . . 78
➠ The density . . . 79
⇝ d color . . . 79
⇝ d font . . . 79
⇝ d unit . . . 80
⇝ d precision . . . 81
➻ d . . . 83
➠ The lattice structure . . . 83
⇝ ls . . . 83
⇝ ls color . . . 84
⇝ ls font . . . 84
⇝ ls align . . . 85
⇝ ls unit . . . 85
⇝ ls precision . . . 85
➻ lat . . . 87
➠ The discovery year . . . 87
⇝ DiscY color . . . 87
⇝ DiscY font . . . 87
⇝ DiscY BC scale . . . 88
➠ *The electron distribution* . . . 88
⇝ eDist color . . . 88
⇝ eDist font . . . 89
⇝ eDist sep . . . 89
➠ *The other contents* . . . 90
⇝ <content name> color . . . 90
⇝ <content name> font . . . 91
➻ cell font . . . 91
➻ cell color . . . 91
⇝ E precision . . . 92
⇝ T precision . . . 93
⇝ Cp precision . . . 95
⇝ kT precision . . . 96

Designing cells with \pgfPTbuildcell **99**
✠ The cell contents . . . 100
✠ Built-in cell styles . . . 102

Designing color schemes **105**
✠ Designing a color scheme with \pgfPTnewColorScheme . . . 105
✠ Designing a color scheme with pgfPTcolorSchemes.html . . . 105

Libraries **109**
Color Schemes Library 109
\pgfPTGroupColors 109
\pgfPTPeriodColors 114
\pgfPTCScombine 117
\pgfPTCSwrite 120

A few more examples **122**

Index **129**

Installation and usage

pgf-PeriodicTable is placed under the terms of the LaTeX Project Public License, version 1.3 or later (http://www.latex-project.org/lppl.txt). pgf-PeriodicTable loads and only requires the TikZ package.

You need to put the package files (pgf-PeriodicTable.sty & *friends*) in a location where PDFLaTeX, LuaLaTeX or XeLaTeX can find them. According to the TDS conventions this may be a subdirectory named *tex/latex/pgf-PeriodicTable/* or *tex/latex/misc/* in your (site specific) installation tree (insert your appropriate directory delimiter instead of /, if needed).

If you are using PDFLaTeX, LuaLaTeX or XeLaTeX you can just simply include the style file without any option via the `\usepackage` command, `\usepackage{pgf-PeriodicTable}`

It can also be loaded with *one option* to select the desired language:
`\usepackage[language flag]{pgf-PeriodicTable}`

The language flags available are:

- ✓ **en** for English (default),
- ✓ **fr** for French,
- ✓ **de** for German,
- ✓ **pt** for Portuguese (Portugal),
- ✓ **br** for Portuguese (Brazil),
- ✓ **es** for Spanish and
- ✓ **it** for Italian.

The data

The data available in pgf-PeriodicTable was mainly compiled with selected and filtered data from Wikipedia, taken from November 2021 to July 2022.

acronym	description	unit	remarks (compiled from @date)
Ar	Relative Atomic Mass		(Wikidata @09/jan/2022)
Arstar	Standard Relative Atomic Mass		STANDARD ATOMIC WEIGHTS 2021, Commission on Isotopic Abundances and Atomic Weights, © **CIAAW, 2007–2022** (**https://ciaaw.org/impressum.htm**)
radio	Radioactivity		(gperiodic-3.0.3, Dec 26 2018)
R	Atomic Radius	pm	Calculated (Wikidata @04/jul/2022)
Rcov	Covalente Radius	pm	Single bond, Wikidata @04/jul/2022)
Rion	Ionic Radius	pm	(Wikidata @04/jul/2022)
Ei	First Ionization Energy	$kJ \cdot mol^{-1}$	(Wikidata @04/jul/2022)
eneg	Electronegativity (Pauling)		(Wikidata @04/jul/2022)
eaff	Electroaffinity	$kJ \cdot mol^{-1}$	(Wikidata @04/jul/2022)
O	Oxidation States		(Wikidata @09/jan/2022)
Tmelt	Melting Point	K	at standard pressure (Wikidata @21/dez/2021)
TmeltC	Melting Point	°C	at standard pressure (Wikidata @21/dez/2021)
Tboil	Boiling Point	K	at standard pressure (Wikidata @21/dez/2021)
TboilC	Boiling Point	°C	at standard pressure (Wikidata @21/dez/2021)
eDist	Electron Distribution		(Wikidata @01/nov/2021)
eConfign	Electronic Configuration (increasing n)		(Wikidata @01/nov/2021)
eConfign	Electronic Configuration (increasing $n + \ell$)		(Wikidata @01/nov/2021)
d	Density	$g \cdot dm^{-3}$ for gases $g \cdot cm^{-3}$ all other physical states	physical state at 25°C, 1 atm (Wikidata @01/nov/2021)
Cp	Specific heat capacity	$J \cdot mol^{-1} \cdot K^{-1}$	at 25°C and 100 kPa (Wikidata @20/nov/2021)
kT	Thermal Conductivity	$W \cdot m^{-1} \cdot K^{-1}$	at 25°C (Wikidata @21/nov/2021)
ls	Lattice Structure		(Wikidata @20/dez/2021 and University of Bielefeld)
lsa	Lattice constant: a	pm	(University of Bielefeld @21/dez/2021)
lsb	Lattice constant: b	pm	(University of Bielefeld @21/dez/2021)
lsc	Lattice constant: c	pm	(University of Bielefeld @21/dez/2021)
lsca	Lattice c/a ratio		Calculated from available data and rounded to two digits
DiscY	Discover Year		(Wikidata @22/dez/2021)
DiscC	Discover Country		(Wikidata @22/dez/2021)
spectra	Visible range spectral lines		Elements spectrum made with \pgfspectra. See the pgf-spectra manual for more details

The utilization of the *acronyms* will be explained in Designing cells with \pgfPTbuildcell.

The commands

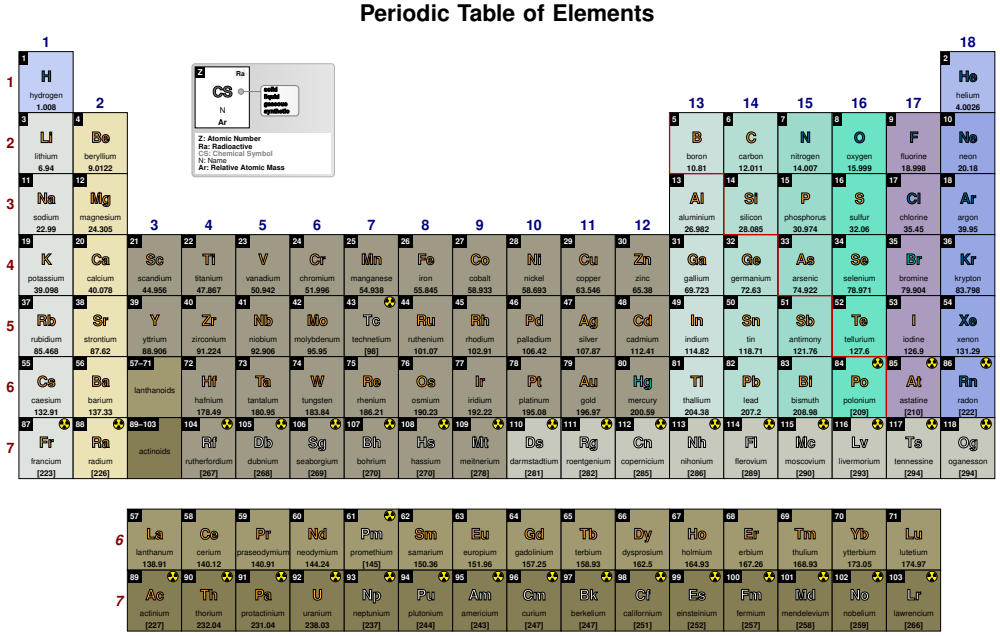
The commands to achieve the Periodic Table of Elements are:

- \pgfPT or \pgfPT[options list] – draws a full or partial graphical Periodic Table controlled by the optional keys.
- \pgfPTstyle[options list] – sets the global style for the Periodic Table.
- \pgfPTresetstyle – resets the style for the Periodic Table with the default values.
- \pgfPTbuildcell(nrows,ncolumns)[entries] – builds the contents of each cell in the Periodic Table.
- \pgfPTresetcell – resets the cell to its default layout.
- \pgfPTbuildcellstyle{name}(nrows,ncolumns)[entries] – builds the contents of each cell in the Periodic Table and stores it in a named style.
- \pgfPTpreviewcell or \pgfPTpreviewcell[scale factor] – preview the last unnamed built cell with an optional scale factor. If no cells have yet been built, the default cell is shown.
- \pgfPTpreviewcellstyle{name} or \pgfPTpreviewcellstyle[scale factor]{name} – preview the named builded cell with an optional scale factor.
- \pgfPTnewcolorscheme[trailing color]{name}{color list} – makes a color scheme to fill the cells along the Periodic Table.
- \pgfPTnewZlist{name} – create a user defined atomic numbers (Z) named list.
- \pgfPTsetLanguage{language flag} – globally change the default language.

► Utilization of \pgfPT

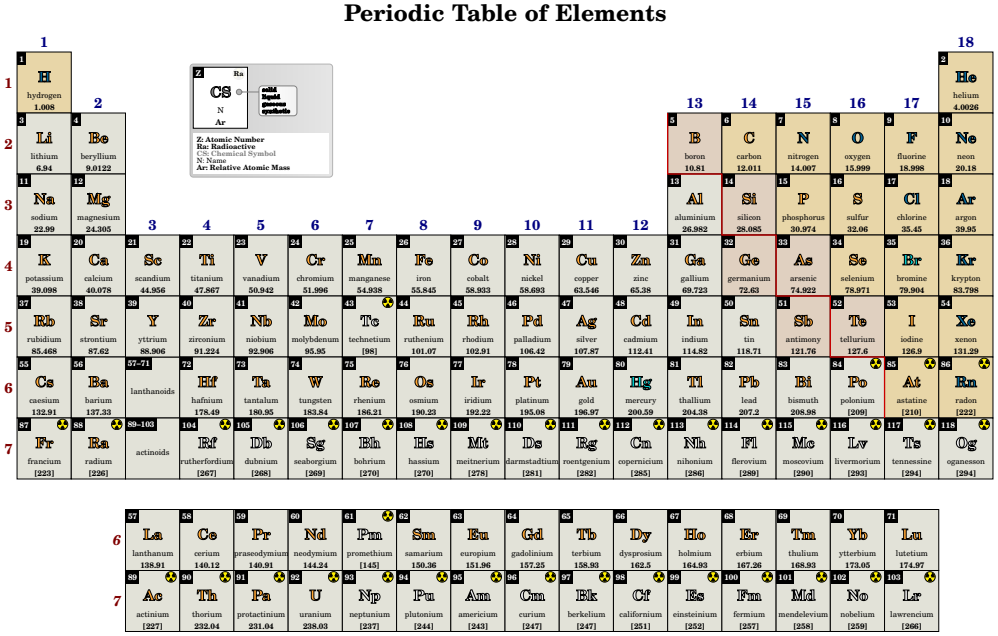
Use this command to draw the Periodic Table of Elements in the language selected at package inclusion (\usepackage[language flag]{pgf-PeriodicTable}):

```
\pgfPT
```



This command can also be used with options – as described in section Options for \pgfPT: creating a «Periodic Table» – to modify, for instance, the font of the Periodic Table or the colors of the cells:

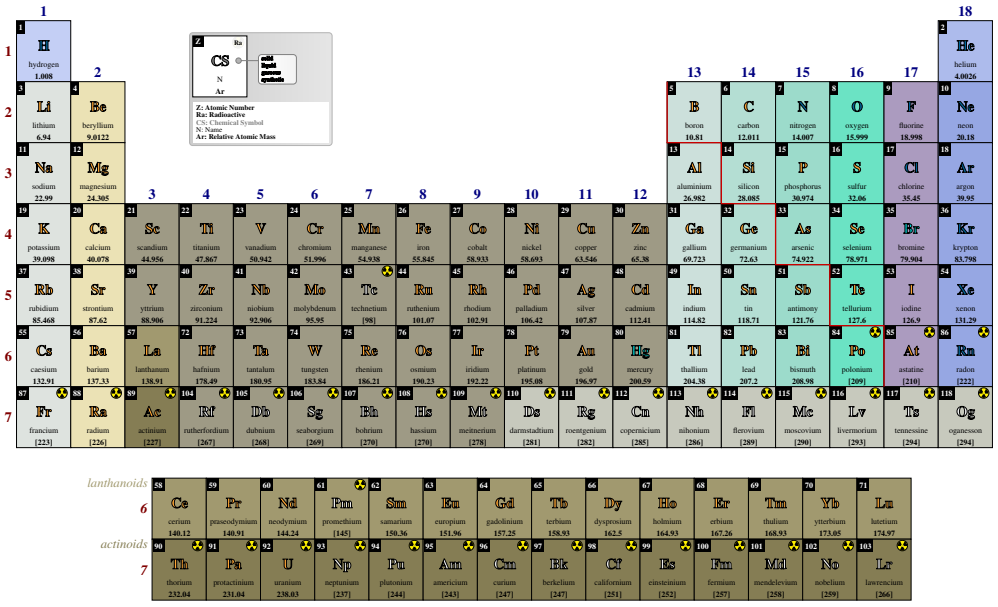
```
\pgfPT[font=pnc,back color scheme=MNM]
```



► Utilization of \pgfPTstyle[options list]

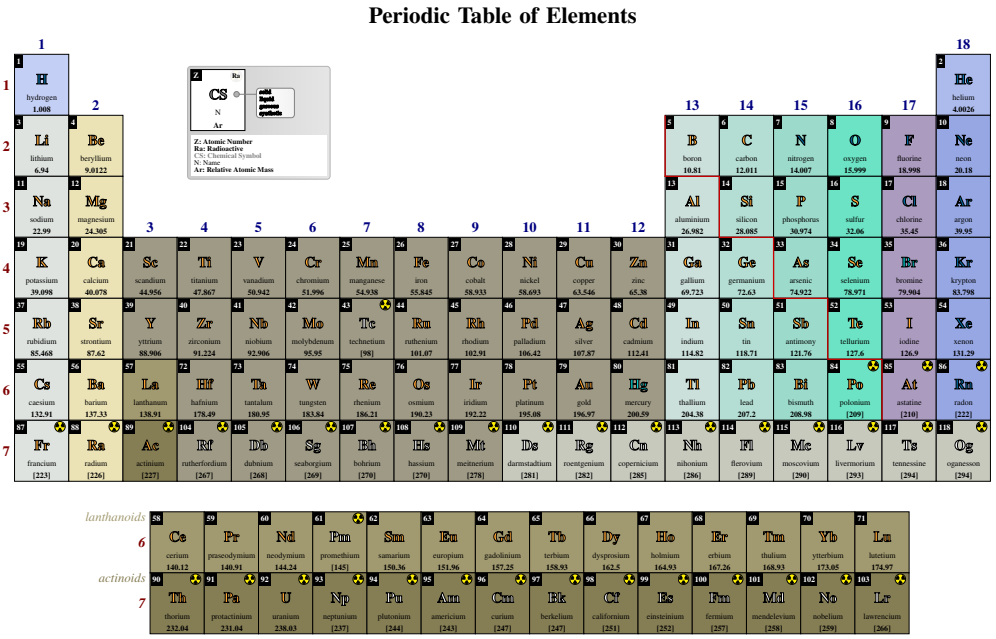
This command globally sets a style for the Periodic Table:

```
\pgfPTstyle[font=ptm,IUPAC=false,show title=false]
\pgfPT
```



It is possible to locally override the *global style* defined:

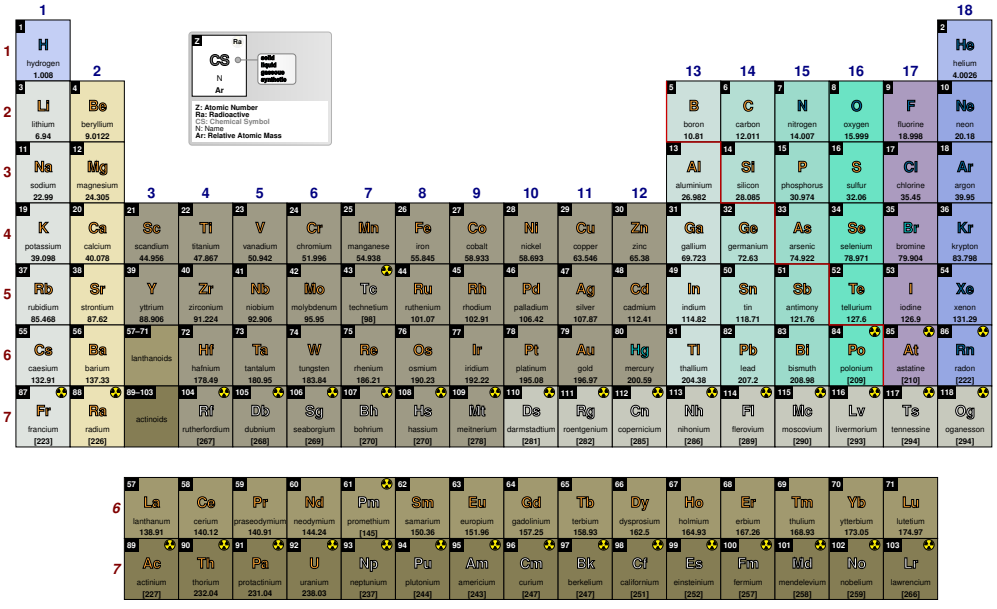
```
\pgfPT[show title]
```



► Utilization of \pgfPTresetstyle

This command resets the style used in the Periodic Table to default values:

```
\pgfPTresetstyle
\pgfPT
```



► Utilization of \pgfPTbuildcell(nrows,ncolumns)[entries]

With \pgfPTbuildcell it is possible to customize the *elementar* cell of the Periodic Table. Each cell is built on the given *number of rows* and *number of columns*. After that, each *entry* is constructed according to the structure row;column;what or initial row-final row;initial column-final column;what.

✓ The first *syntax* – row;column;what – puts «what» in the «row» row and in the «column» column with the height of one row and the width of one column:
– for example, 1;1;Z puts the atomic number Z in row 1 and column 1, witch actually corresponds to a box anchored to the top left corner of the cell and that *goes* below and to the right of that corner.

✓ The second *syntax* – initial row-final row;initial column-final column;what – puts «what» from «initial row» to «final row» with the height of $\text{final row} - \text{initial row} + 1$ and from «initial column» to «final column» with the width of $\text{final column} - \text{initial column} + 1$. It is important to keep in mind that when using this syntax the *row* and *column* could have any value between **1** and **number of rows** and **number of columns**, respectively.
– for example, 1;1-2.1;Z puts the atomic number Z in row 1 with the height of one row and from column 1 to *column* 2.1, with the width of $2.1 \times \text{column}$. Note that in this example the two *syntaxes* are mixed up.

The **default cell** of the Periodic Table is constructed with the command:

```
\pgfPTbuildcell(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-2.5;CS),(4;1-3;name),(5;1-3;Ar)]
```

► Utilization of \pgfPTresetcell

The \pgfPTresetcell resets the cell to its default layout.

► Utilization of \pgfPTbuildcellstyle{name}(nrows,ncol...)[entr...]

The \pgfPTbuildcellstyle command works like \pgfPTbuildcell, but stores the cell style under the name provided. It is only used when called via the cell style passed as an option to \pgfPT. Otherwise it remains unavailable, unlike the \pgfPTbuildcell command which immediately affects the cells of the Periodic Table.

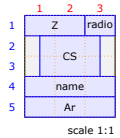
► Utilization of \pgfPTpreviewcell

The main purpose of this command is to show the built cell for *debugging*.
With \pgfPTpreviewcell you can preview the last unnamed built cell with an optional scale factor. If no cells have yet been built, the default cell is shown.

```
\pgfPTpreviewcell
```

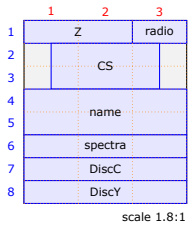
Using the last cell built
The build command:

```
\pgfPTbuildcell(5,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;Ar)]
```



```
\pgfPTbuildcell(8,3)% 8 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4-5;1-3;name), (6;1-3;spectra),(7;1-3;DiscC),(8;1-3;DiscY)]
                                        \pgfPTpreviewcell[1.8]
```

Using the last cell built
The build command:
\pgfPTbuildcell(8,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4-5;1-3;name),(6;1-3;spectra),(7;1-3;DiscC),(8;1-3;DiscY)]



► Utilization of \pgfPTpreviewcellstyle{name}

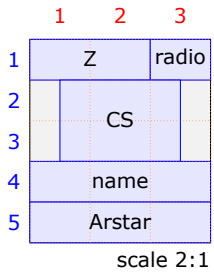
This previews a *named* cell, again with the optional scale factor.

```
\pgfPTpreviewcellstyle{myname}
```

User style **myname** *doesn't exist!*

```
\pgfPTbuilcellstyle{myname}(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;Ar*)]
                                        \pgfPTpreviewcellstyle[2]{myname}
```

User style **myname**
The build command:
\pgfPTbuildcell(5,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;Ar*)]



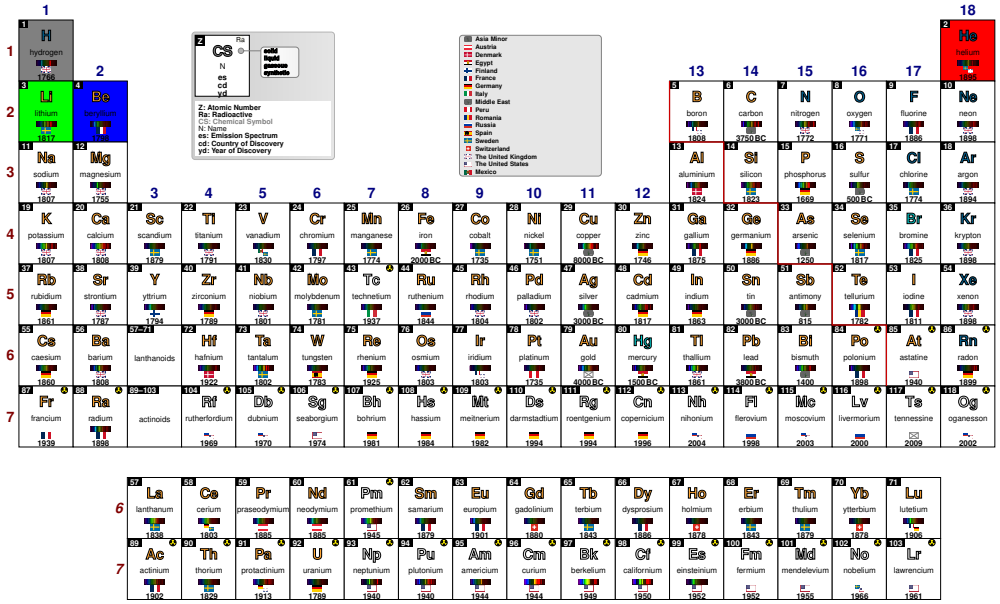
► Utilization of \pgfPTnewColorScheme{name}{color list}

Use this command to create a *color scheme* for cells in the Periodic Table. It has two mandatory arguments – name and color list – and an optional argument – trailing color.

The name is used to identify the *color scheme*. The color list is a comma-separated list of red, green and blue values written as r/g/b, defined in ascending order of Z and starting at Z=1. The optional argument trailing color is appended to the end of the list and is used for all cells starting from this point on. It also has the form r/g/b and its default value is 1/1/1 (white).

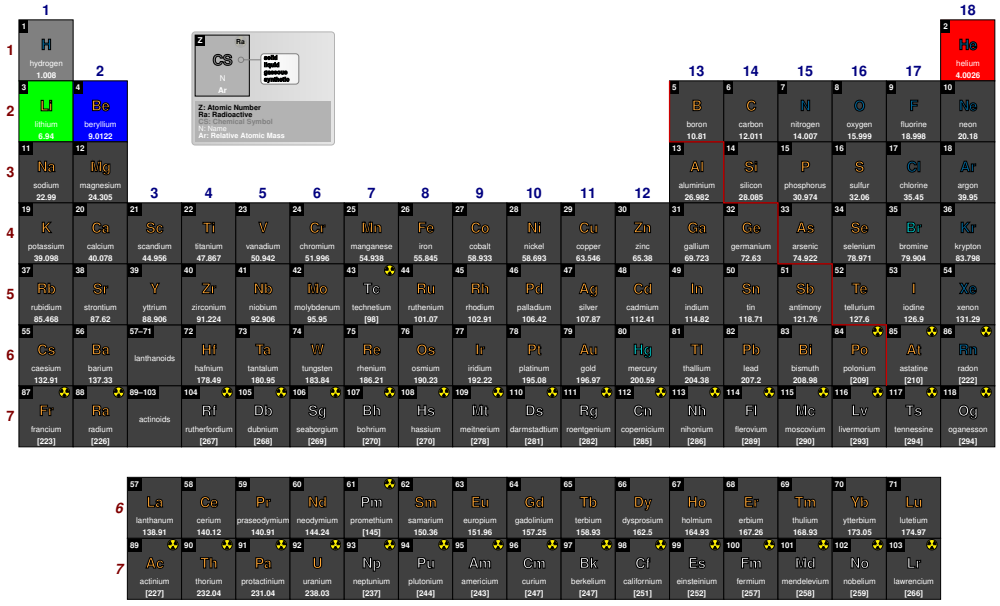
```
\pgfPTnewColorScheme{myname}{.5/.5/.5,1/0/0,0/1/0,0/0/1}
\pgfPT[back color scheme=myname]
```

Periodic Table of Elements



```
\pgfPTnewColorScheme[.25/.25/.25]{myname}{.5/.5/.5,1/0/0,0/1/0,0/0/1}
\pgfPTresetcell
\pgfPT[back color scheme=myname,name color=white, Ar color=white,legend back color=black!30]
```

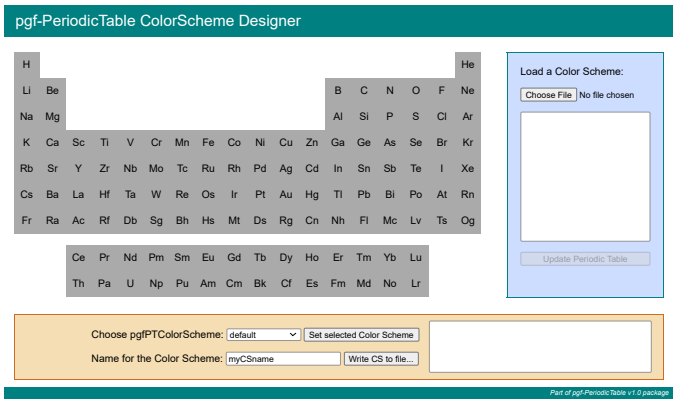
Periodic Table of Elements



There are a few *color schemes* predefined:

- ✓ **default**, the default built-in color scheme, which is loaded if no value is passed to the back color scheme key.
- ✓ **Soft**, a soft color pattern for cells, differentiating metals, non metals, semimetals, lanthanides and actinides.
- ✓ **Jmol**, a color scheme based upon Jmol: an open-source Java viewer for chemical structures in 3D.
- ✓ **CPK**, a color scheme that is based upon the colors of the popular plastic spacefilling models which were developed by Corey, Pauling and later improved by Kultun.
- ✓ **Rasmol** and **RasmolNew**, two color schemes based upon the computer program RasMol.
- ✓ **Wikipedia**, a color scheme built on the Periodic Table of Elements available at Wikipedia.
- ✓ **MNM**, a color pattern which distinguishes between **M**etals, semimetals and **N**on **M**etals.
- ✓ **PS**, a color scheme depicting the **P**hysical **S**tate at room temperature.
- ✓ **Radio**, a two color color scheme showing the radioactivity of the elements.
- ✓ **Blocks**, a four colored color scheme showing the *s*, *p*, *d* and *f* blocks of the Periodic Table.

Writing a color scheme can be painstaking work, so a *script* is provided for that:



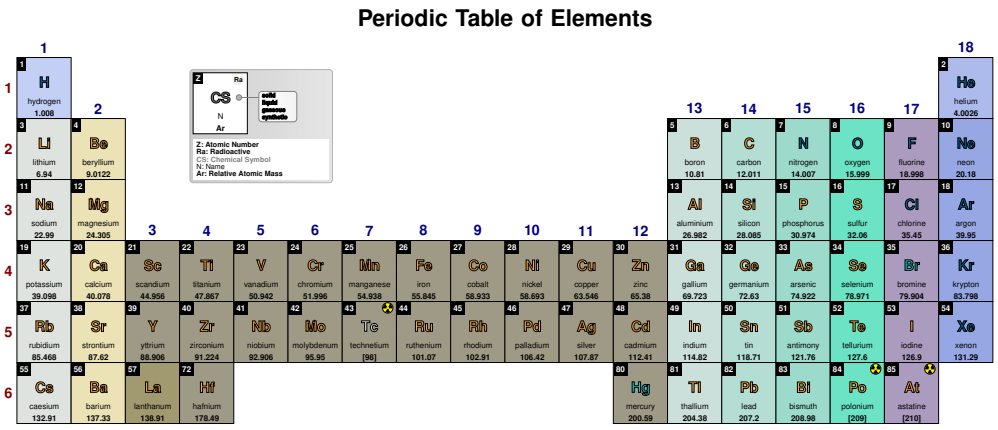
pgfPTcolorSchemes.html

► Utilization of \pgfPTnewZlist{name}

This command makes a user defined atomic numbers' list with the provided name. The list can be anything that the \foreach loop, defined in the TikZ package, can understand.

For more information on how to use \foreach loop refer to the section *Repeating Things: The Foreach Statement* in the pgfmanual.

```
\pgfPTnewZlist{myZlist}{1,...,57,72,80,81,...,85}
\pgfPT[Z list=myZlist,IUPAC=false]
```

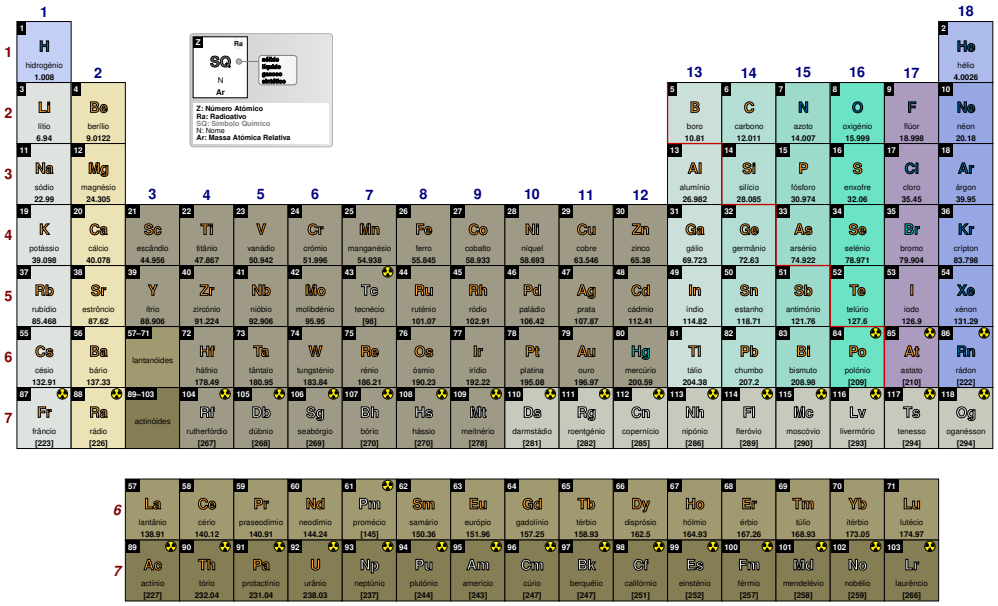


► Utilization of \pgfPTsetLanguage{language flag}

This command globally changes the default language of the Periodic Table.

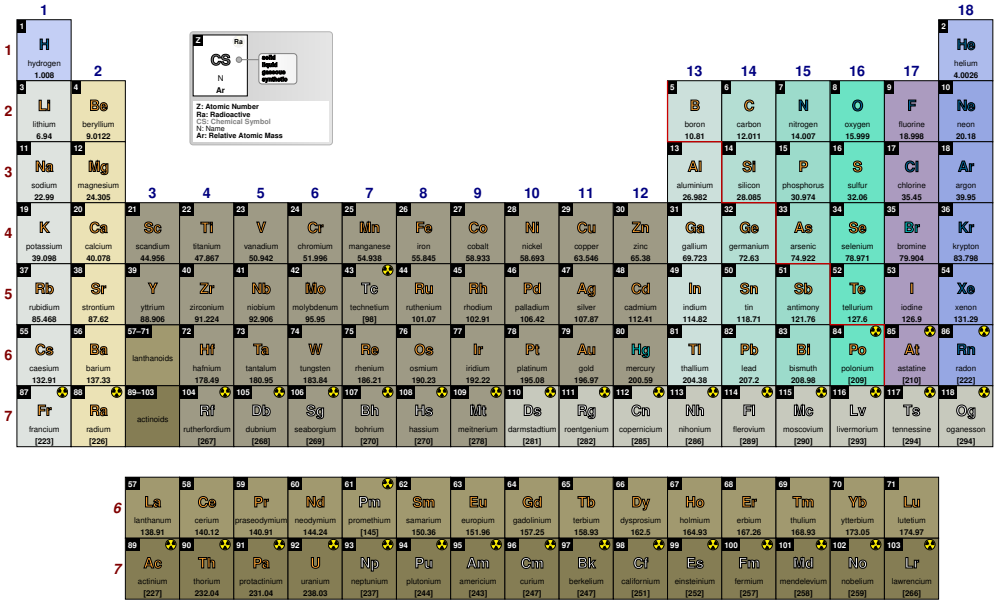
```
\pgfPTsetLanguage{pt}
\pgfPT
```

Tabela Periódica dos Elementos




```
\pgfPTsetLanguage{en}
\pgfPT
```

Periodic Table of Elements



Options for \pgfPT: creating a «Periodic Table»

For the commands \pgfPT and \pgfPTstyle there are a set of options available to draw the Periodic Table or any portion of the Periodic Table, as described below.

The list of options is a comma separated list of any of the following elements:

- a 'key' or a 'key=value' pair,
- a 'style' or a 'style=value' pair,
- a *pseudo style* with a proper syntax: 'style={key 1=value 1, key 2=value 2, … , key n=value n}', where none of the '*keys*' are mandatory.

The options *can be divided* in two subsets, one that affects the *appearance* of the *entire* Periodic Table, the other that concerns the *contents* of each cell of the Periodic Table.

✠ Periodic Table options: keys, styles and *pseudo styles*

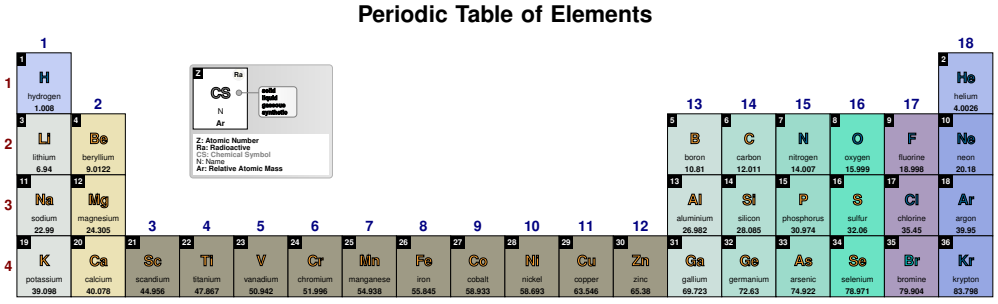
The following options and styles are used to *control* the Periodic Table *as a whole* in various aspects, such as the cell width or cell height, which elements are displayed (Z list), whether the title or legend are shown – show title or show legend – among others.

➠ General layout

Z list default: *all*

Set's the list of the elements to display in the Periodic Table. It could be a name or a comma separated list of atomic numbers, which in turn supports *the dots notation* as explained in the section *Repeating Things: The Foreach Statement* in the pgfmanual.

```
\pgfPT[Z list={1,...,36}]
```



The possible name is one of the following:

✓ **built-in**:

- ▷ 'all' is equivalent to Z list={1,...,118}, *i.e.*, all known elements.
- ▷ 's', 'p', 'd' or 'f', for the elements in the corresponding blocks.
- ▷ 'sp', 'spd', for the elements resulting from merging the corresponding blocks.
- ▷ 'lanthanoids' or simply 'La', for lanthanoids †.
- ▷ 'actinoids' or 'Ac', for actinoids †.
- ▷ 'G1*', 'G1', … , 'G18', which are used, respectively, for the elements of *group 1 without hydrogen*, *group 1*, … , *group 18*.
- ▷ 'P1', … , 'P7', 'P6*', 'P7*', which are used, respectively, for the elements of the 1^{st} *period*, … , 7^{th} *period*, 6^{th} *period and lanthanoids* †, 7^{th} *period and actinoids* †.

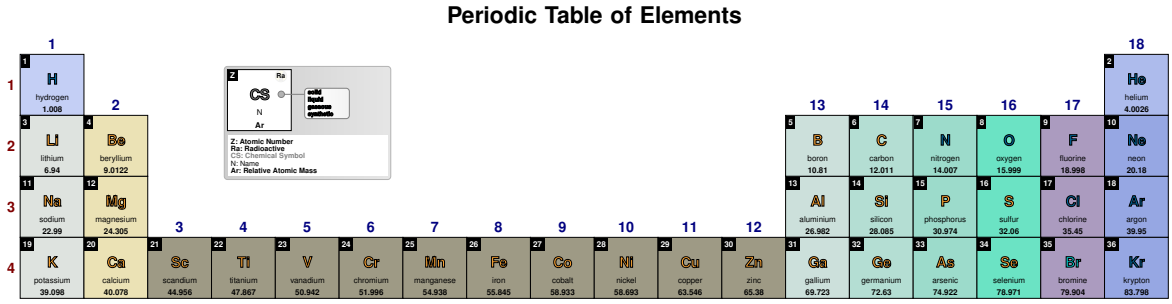
† *Depending on the value of the IUPAC key, the Lanthanum or Actinium are or are not included.*

✓ any **user defined** name via \pgfPTnewZlist{name}{list}

cell width default: *34pt*

Sets the width of each base cell of the Periodic Table.

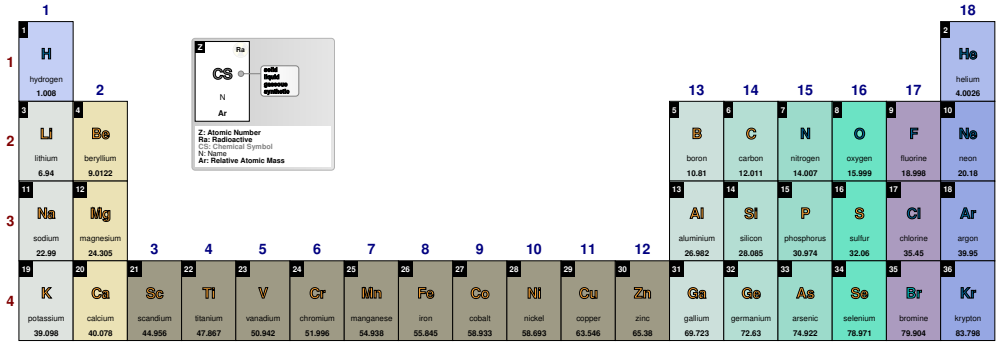
```
\pgfPT[Z list={1,...,36},cell width=40pt]
```



cell height default: *38.25pt*

Sets the height of each base cell of the Periodic Table.

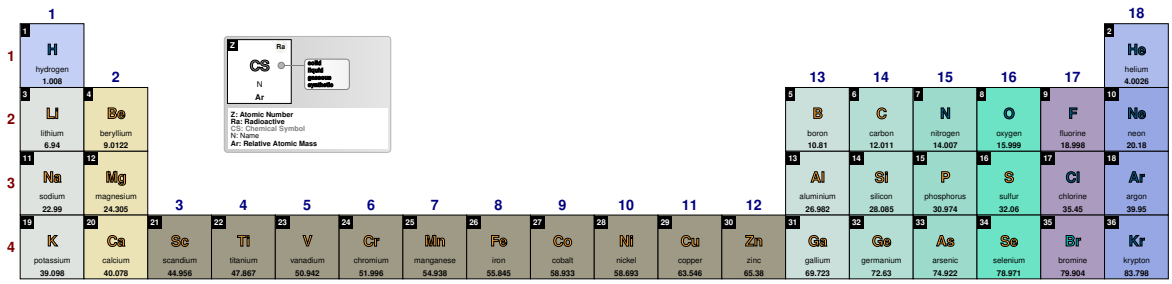
```
\pgfPT[Z list={1,...,36},cell height=50pt]
```



cell size default: *38.25pt*

Style to set both the width and the height of each base cell of the Periodic Table.

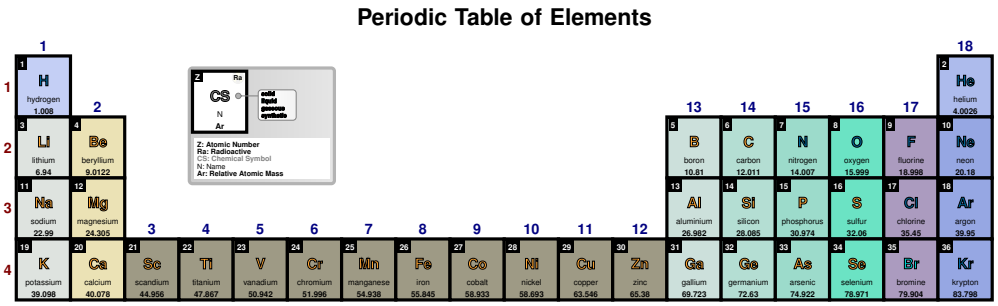
```
\pgfPT[Z list={1,...,36},cell size=40pt]
```



cell line width default: *0.4pt*

Sets the width of the line surrounding the base cell of the Periodic Table.

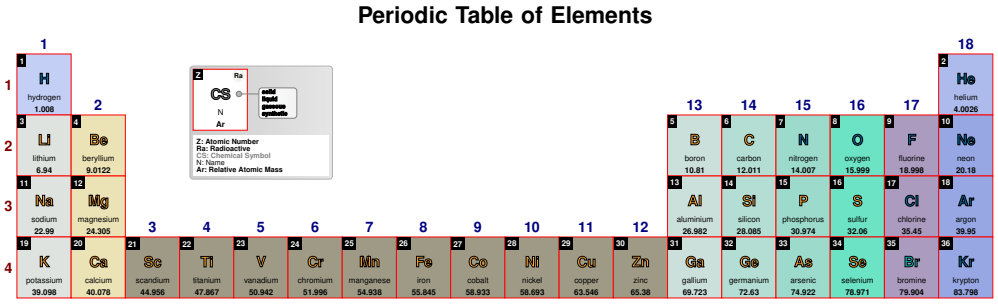
```
\pgfPT[Z list={1,...,36},cell line width=2pt]
```



cell line color default: *black*

Sets the color of the line surrounding the base cell of the Periodic Table.

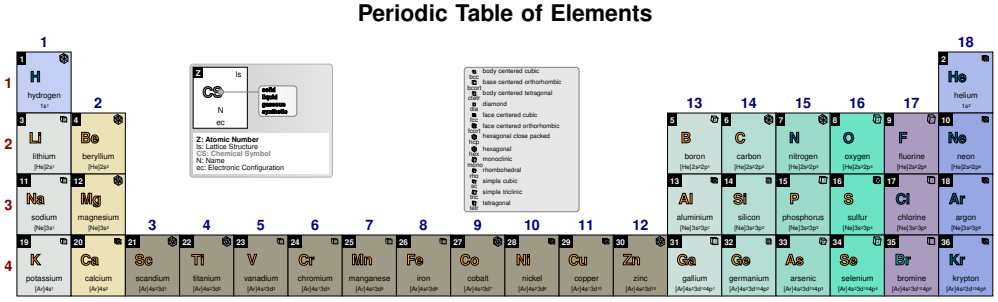
```
\pgfPT[Z list={1,...,36},cell line color=red]
```



cell style default: *{}*

Loads a named cell style, built via \pgfPTbuildcellstyle, to use as a layout for each cell of the Periodic Table.

```
\pgfPTbuilcellstyle{myname}(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;ls),(2-3;1.5-2.5;CS),(4;1-3;name),(5;1-3;eConfignl)]
\pgfPT[Z list={1,...,36},cell style=myname]
```

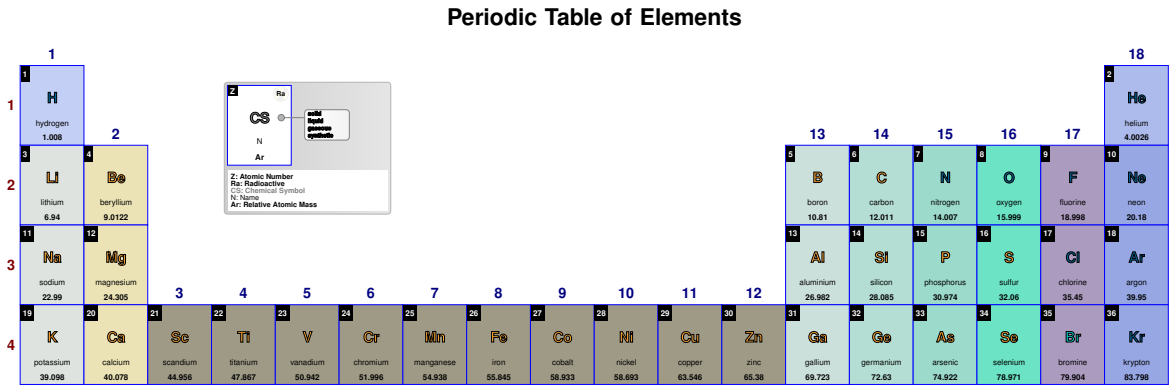


cell default: *{w=34pt,h=38.25pt,lw=.4pt,lc=black}*

Pseudo style to set the cell **w**idth, the cell **h**eight, the cell **s**ize, the cell **l**ine **w**idth, the cell **l**ine **c**olor and/or the cell **style**. None of the *keys* – w, h, s, lw, lc and style – are mandatory.

USAGE: cell={w=<length>,h=<length>,s=<length>,lw=<length>,lc=<color>,style=<name>}

```
\pgfPT[Z list={1,...,36},cell={w=40pt,h=50pt,lw=.6pt,lc=blue}]
```

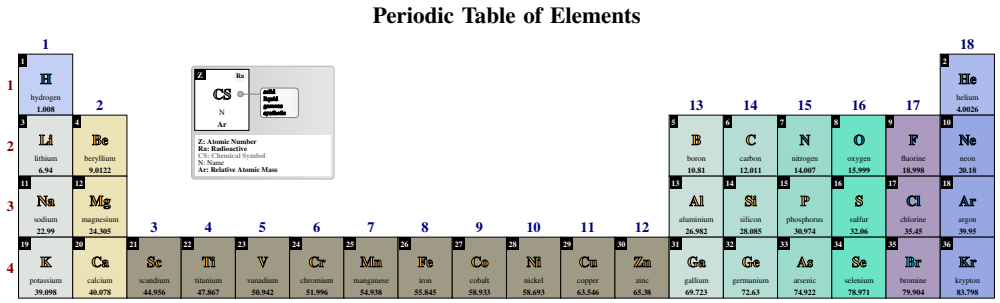


font default: *phv*

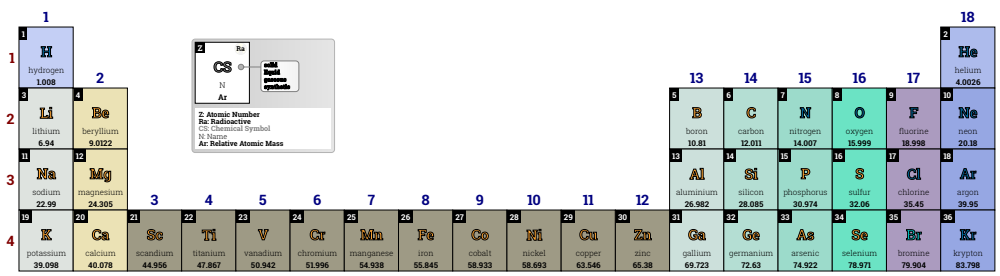
Sets the font family, via the proper *LaTeX font name*, to use in the Periodic Table. The *default* font is *phv*, *i.e.*, the Helvetica font. The value of the font key can be any *LaTeX font name* known to the local LaTeX installation.

See *LaTeX font names* below for further details.

```
\pgfPT[Z list={1,...,36},font=ptm]
```




```
\pgfPT[Z list={1,...,36},font=RobotoSlab-TLF]
```



LaTeX font names:

✓ The LaTeX font names commonly available in LaTeX distributions are:

- **Serif fonts**
  - ▷ cmr – Computer Modern Roman
  - ▷ lmr – Latin Modern Roman
  - ▷ pbk – Bookman
  - ▷ bch – Charter
  - ▷ pnc – New Century Schoolbook
  - ▷ ppl – Palatino
  - ▷ ptm – Times
- **Sans Serif fonts**
  - ▷ cmss – Computer Modern Sans Serif
  - ▷ lmss – Latin Modern Sans Serif
  - ▷ pag – Avant Garde
  - ▷ phv – Helvetica

✓ There are other fonts available to LaTeX that require installation of the corresponding packages:

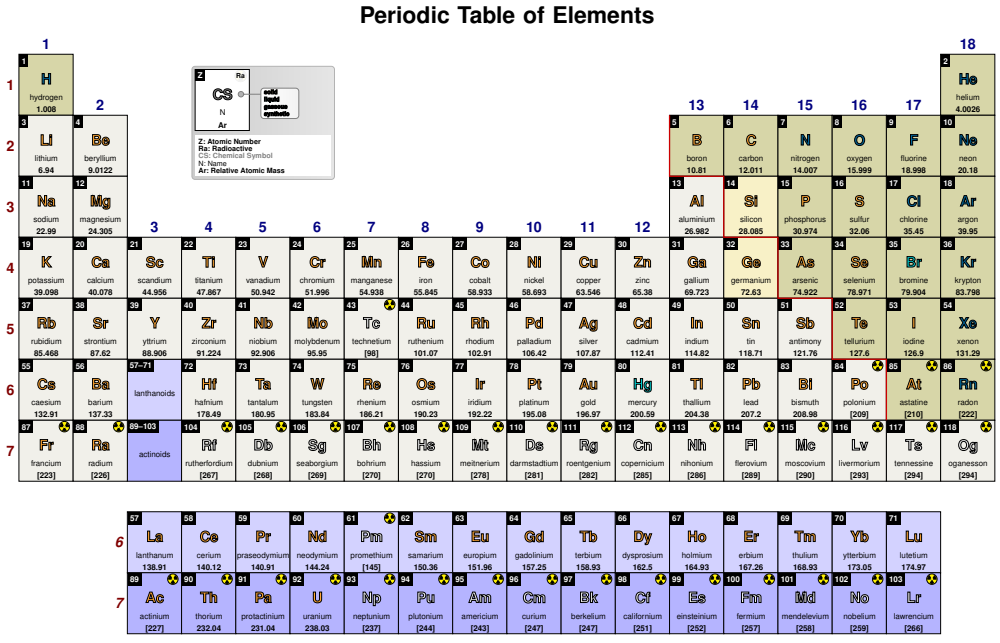
- ▷ the roboto package provides the following fonts:
  - Roboto-TLF – Roboto tabular lining
  - Roboto-LF – Roboto proportional lining
  - Roboto-OsF – Roboto proportional oldstyle
  - Roboto-TOsF – Roboto tabular oldstyle
  - RobotoSlab-TLF – RobotoSlab proportional lining
  - RobotoSlab-OsF – RobotoSlab proportional oldstyle
  - RobotoSlab-TOsF – RobotoSlab tabular oldstyle
  - RobotoMono-TLF – RobotoMono proportional lining
- ▷ the frcursive package provides the *frc - French Cursive* font.
- ▷ the miama package provides the *frm – Miama Nueva* font.
- ▷ ...

For more information about fonts visit the TUG Font Catalogue

back color scheme default: *default*

Sets a named back color scheme for the Periodic Table.

```
\pgfPT[back color scheme=Soft]
```



The possible name is one of the following:

✓ **built-in**:

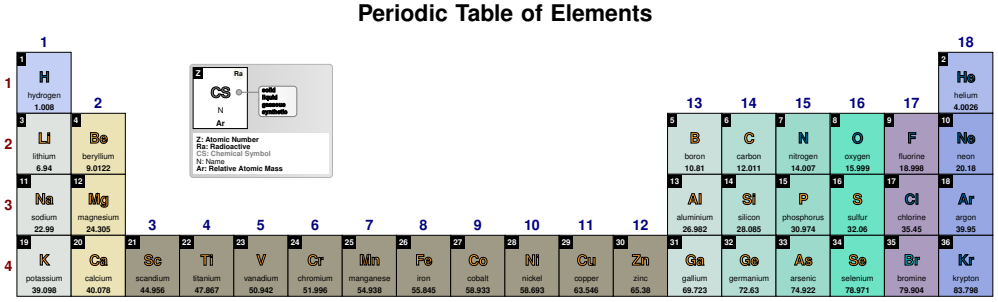
- ▷ 'Soft', a soft color scheme that distinguishes metal, non metals, silicon and germanium, lanthanoids and actinoids.
- ▷ 'Jmol', is the color scheme used in the computer software Jmol: an open-source Java viewer for chemical structures in 3D.
- ▷ 'CPK', is the color scheme of the popular color convention for distinguishing atoms of different chemical elements in molecular models. The scheme is named after the CPK molecular models designed by chemists Robert Corey and Linus Pauling, and improved by Walter Koltun.
- ▷ 'Rasmol', is the color scheme used in the computer software RasMol, a program for molecular graphics visualization originally developed by Roger Sayle.
- ▷ 'RasmolNew', is a color scheme used in RasMol with revision of CPK colors made by C. Chigbo (RasMol 2.7.3).
- ▷ 'Wikipedia', is the color scheme based on the Wikipedia Periodic Table of Elements.
- ▷ 'MNM', is designed to show **M**etals and **N**on **M**etals in two different colors, showing also the semimetals in a third color.
- ▷ 'PS', is designed to show the **P**hysical **S**tate of the elements at normal temperature and pressure (NTP) in different colors.
- ▷ 'Radio', is designed to show the **R**adioactive elements in one color and the non radioactive elements in another color.
- ▷ 'Blocks', for showing the elements in each block of the Periodic Table with the same color.
- ▷ 'solid', to show the background of each cell of the Periodic Table with the same color specified by the key 'back color'.

✓ any **user defined** name via \pgfPTnewColorScheme{name}{color list}

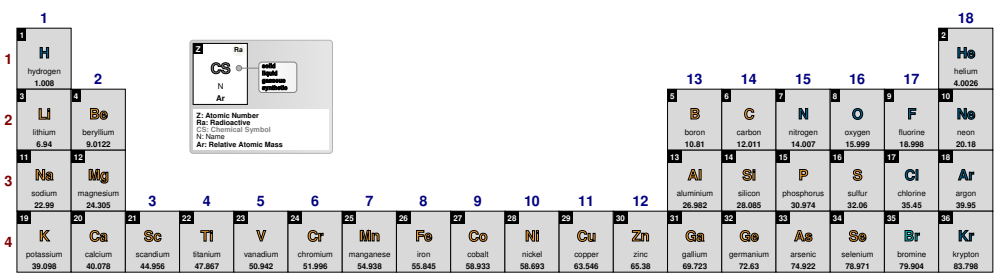
back color default: *white*

Sets the background of each cell of the Periodic Table. It only takes effect if the back color scheme key is set to solid

```
\pgfPT[Z list={1,...,36},back color=black!15]
```




```
\pgfPT[Z list={1,...,36},back color scheme=solid,back color=black!15]
```

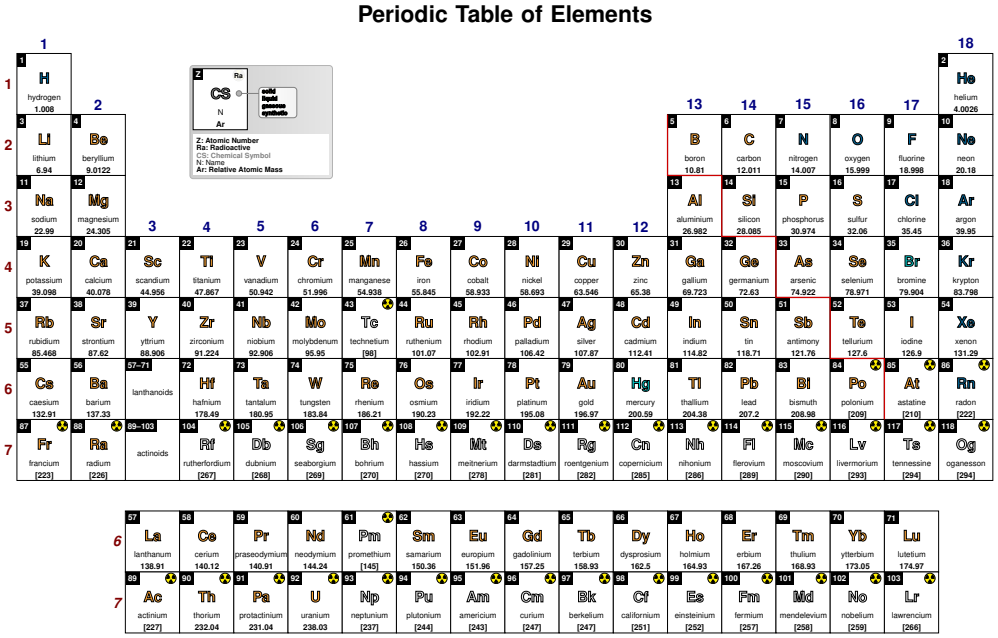


It is possible to set the back color scheme key with the built-in names using the following styles:

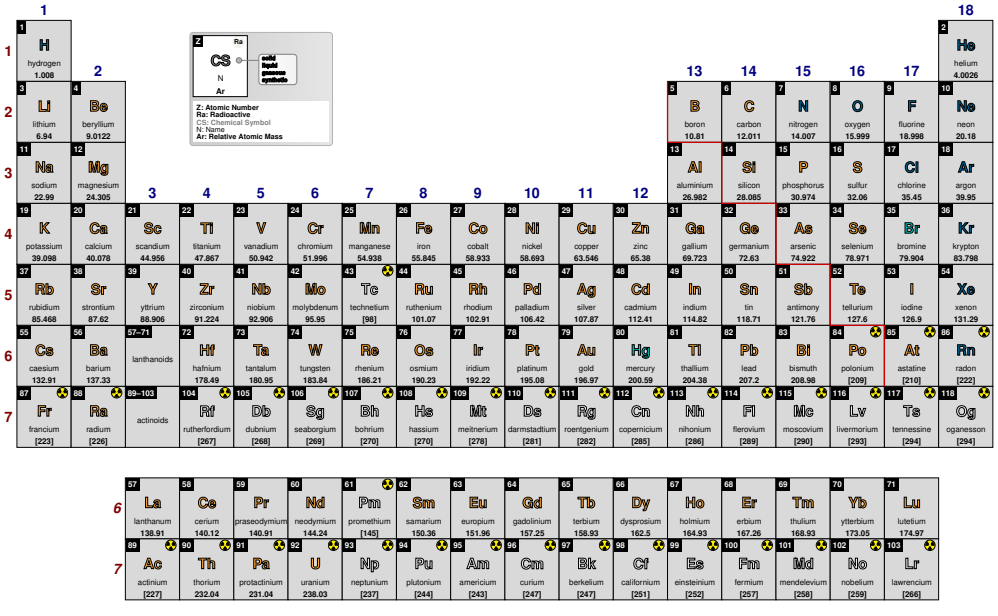
csSolid default: *white*

A style equivalent to back color scheme=solid,back color=#1

```
\pgfPT[csSolid]
```



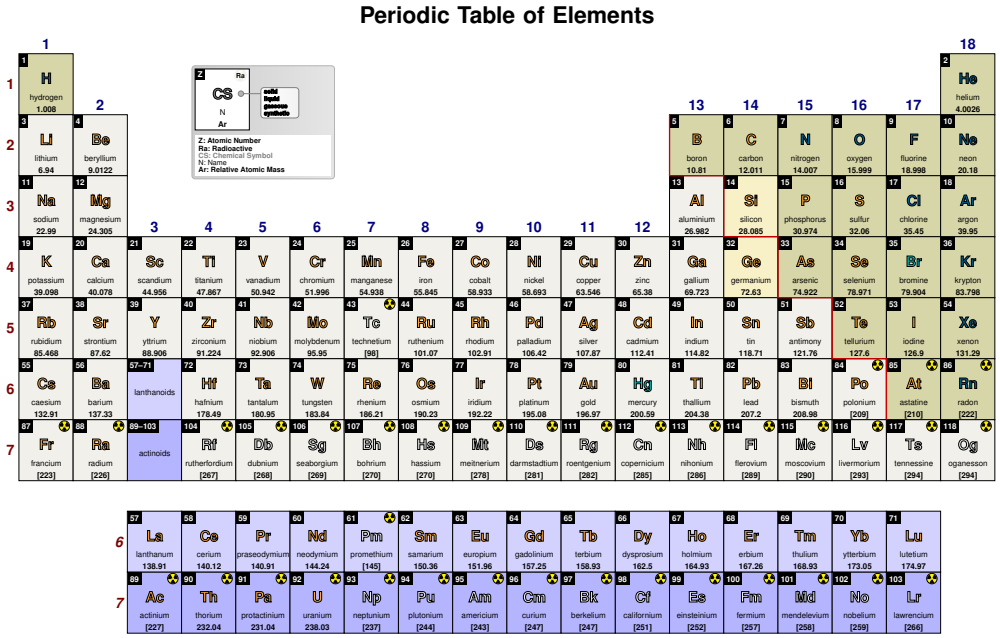

```
\pgfPT[csSolid=black!15]
```



csSoft *no value*

A style equivalent to back color scheme=Soft

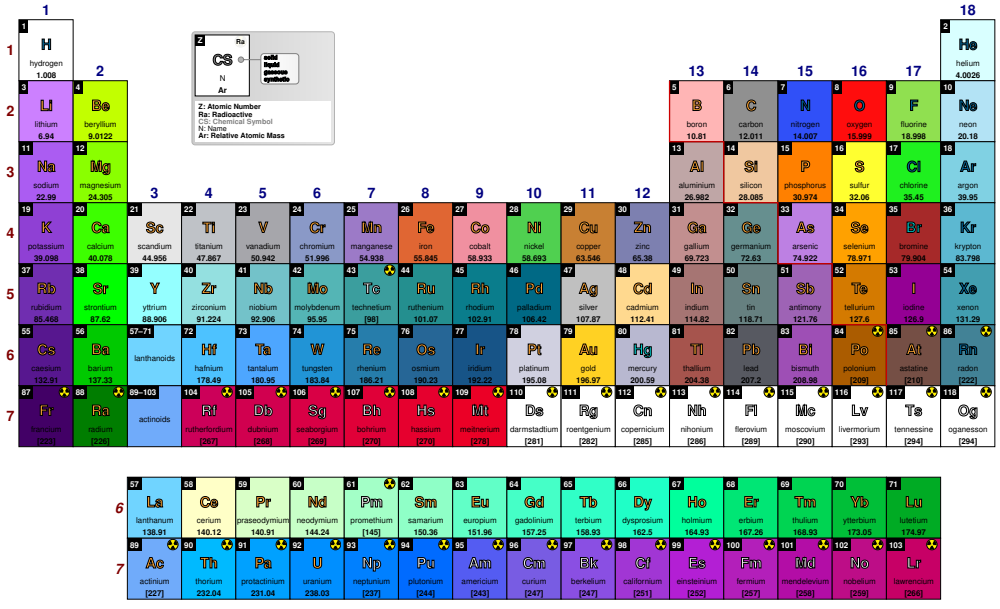
\pgfPT[csSoft]



csJmol *no value*

A style equivalent to back color scheme=Jmol

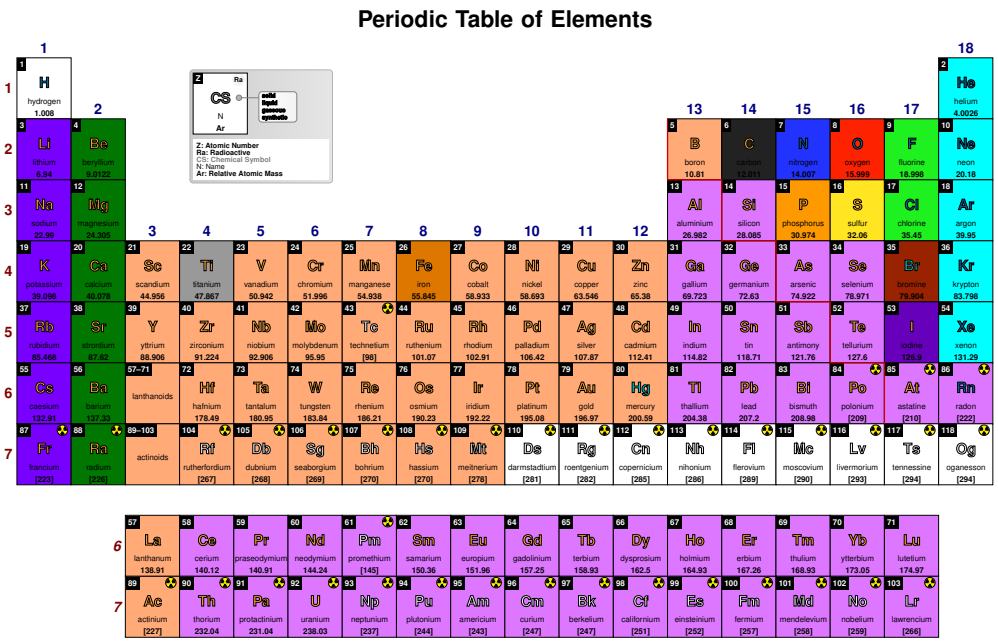
\pgfPT[csJmol]



csCPK *no value*

A style equivalent to back color scheme=CPK

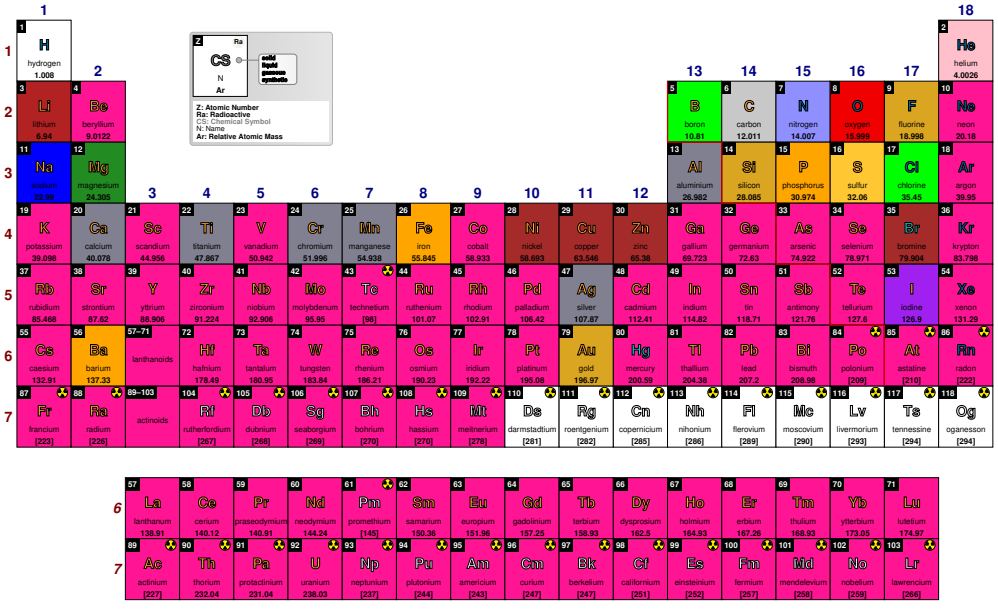
\pgfPT[csCPK]



csRasmol *no value*

A style equivalent to back color scheme=Rasmol

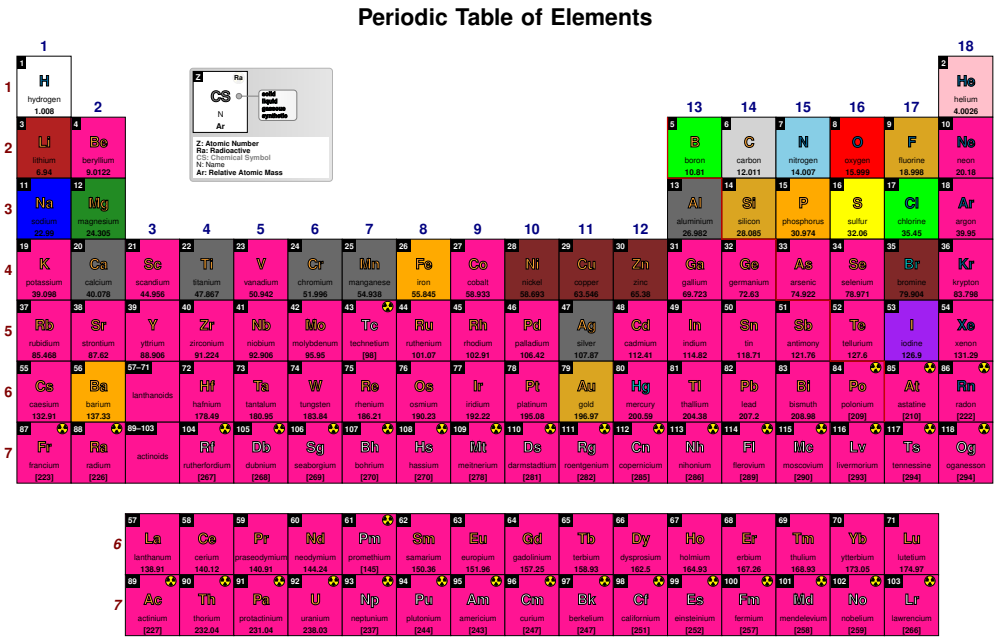
\pgfPT[csRasmol]



csRasmolNew *no value*

A style equivalent to back color scheme=RasmolNew

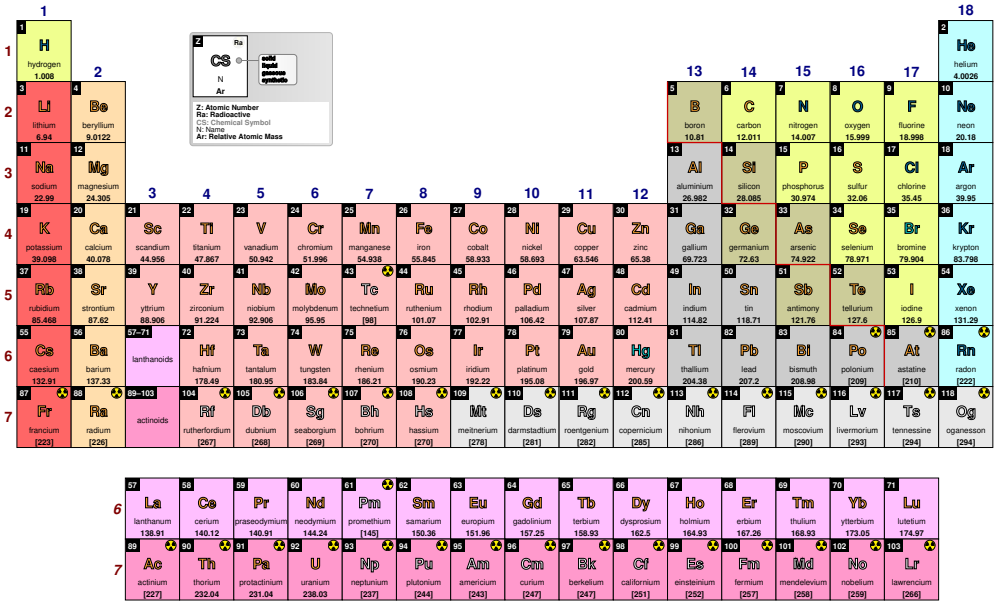
\pgfPT[csRasmolNew]



csWikipedia *no value*

A style equivalent to back color scheme=Wikipedia

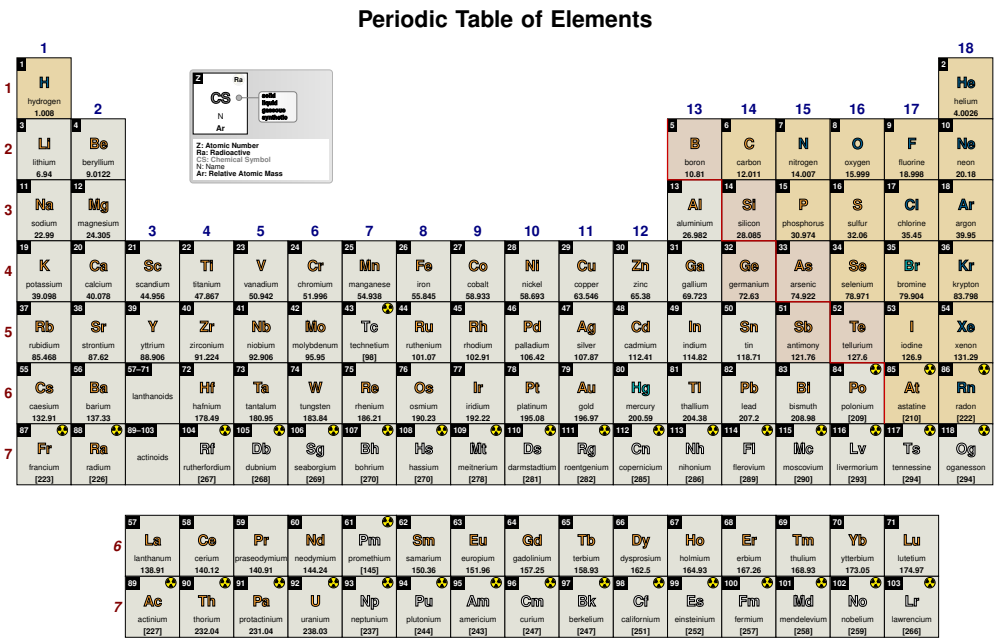
\pgfPT[csWikipedia]



csMNM *no value*

A style equivalent to back color scheme=MNM

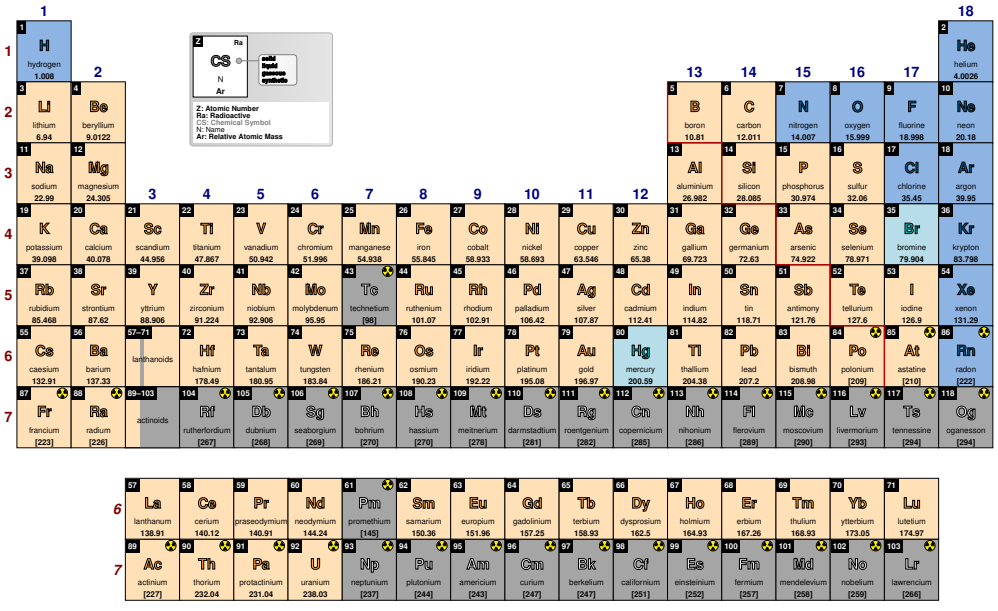
```
\pgfPT[csMNM]
```



csPS *no value*

A style equivalent to back color scheme=PS

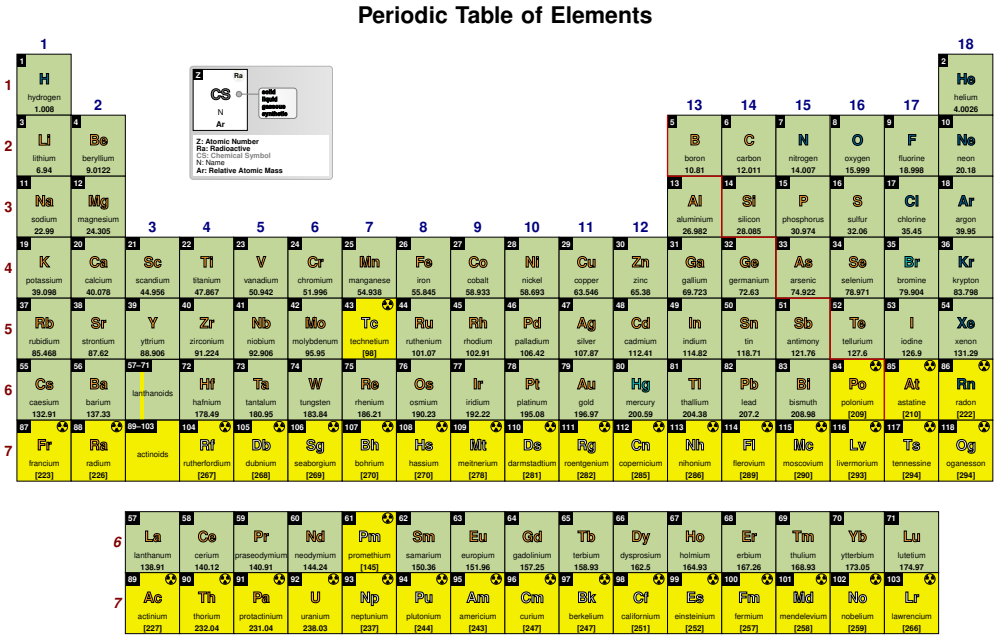
```
\pgfPT[csPS]
```



csRadio *no value*

A style equivalent to back color scheme=Radio

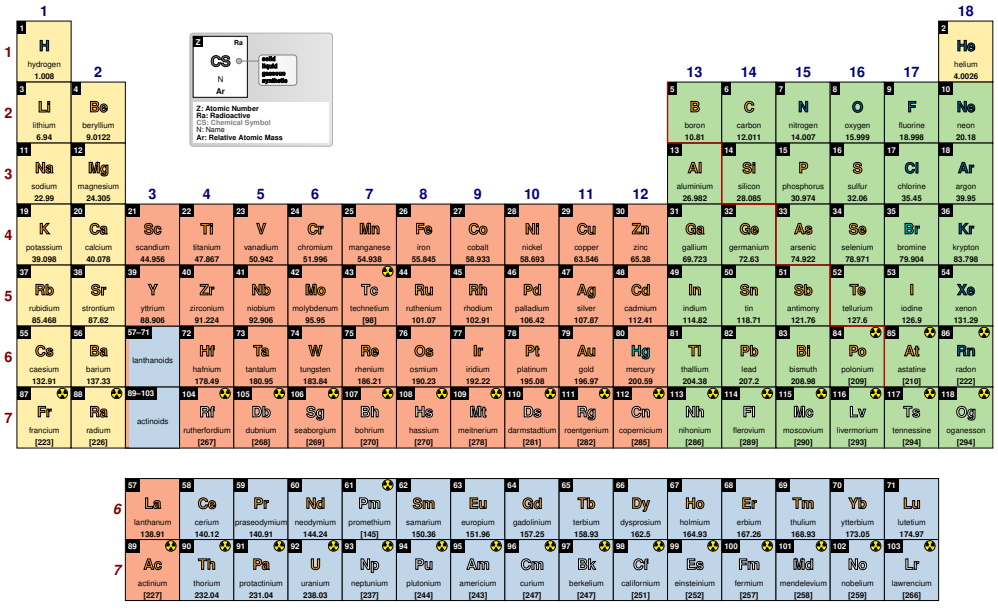
```
\pgfPT[csRadio]
```



csBlocks *no value*

A style equivalent to back color scheme=Blocks

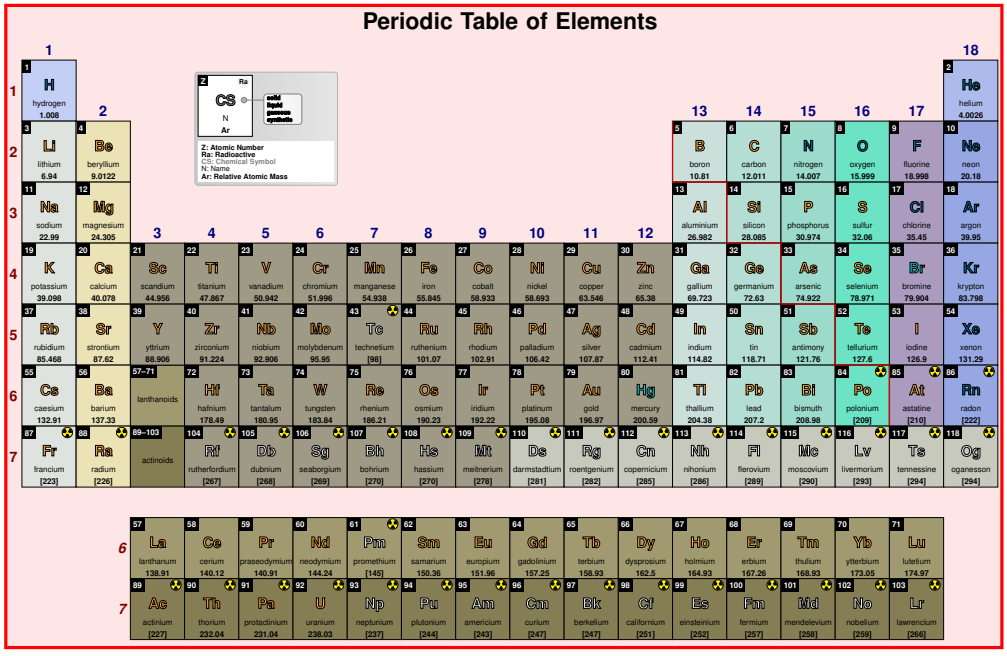
```
\pgfPT[csBlocks]
```



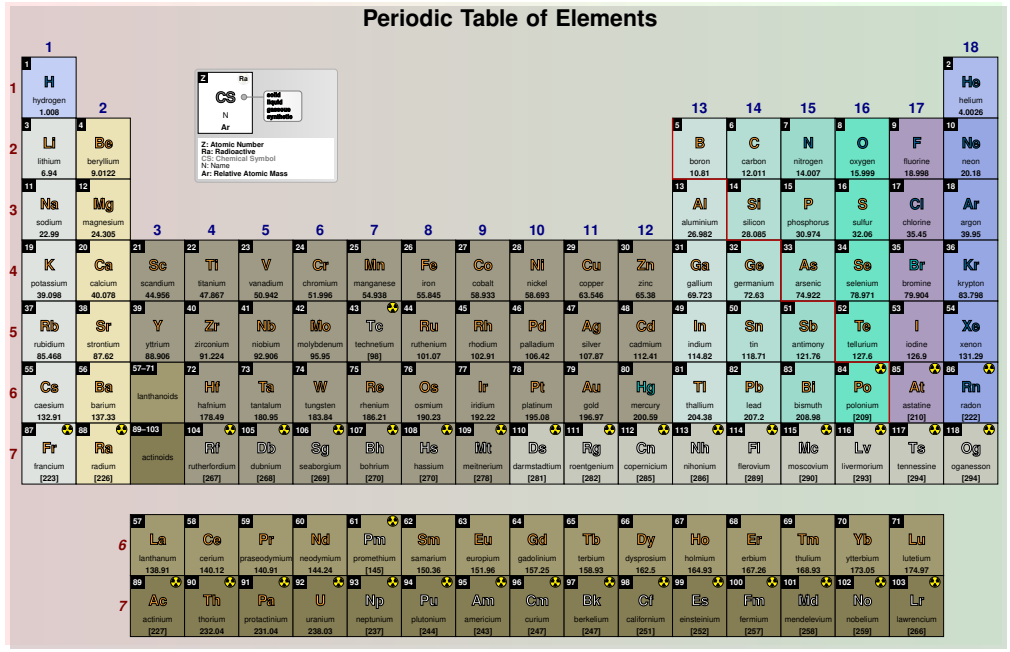
background default: {}

A style to set the background of the Periodic Table, built with any of the TikZ keys that can be applied to a path construction.

```
\pgfPT[background={draw=red,line width=2pt,fill=red!10}]
```



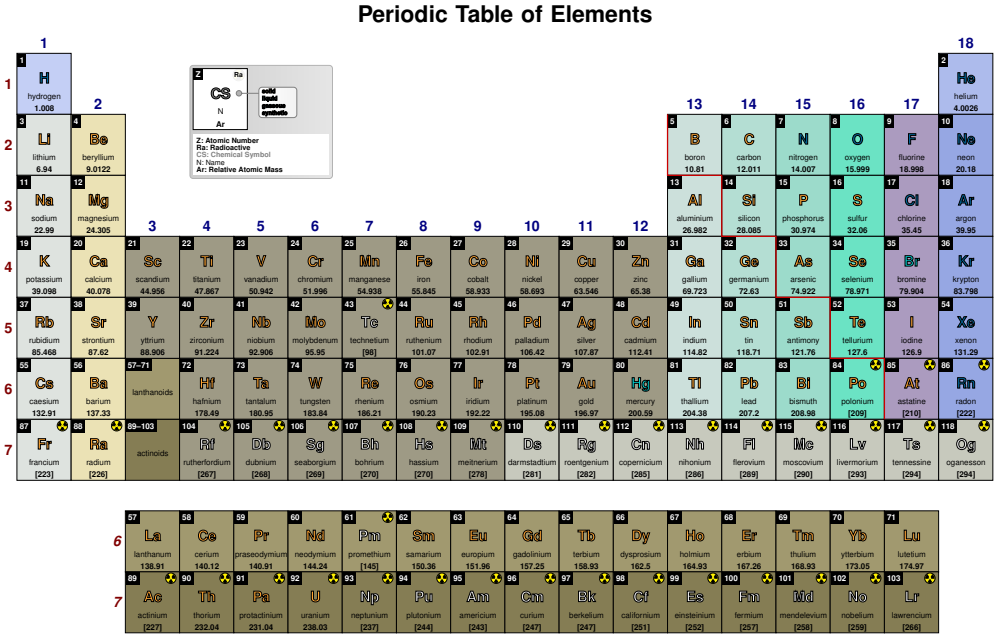

```
\usetikzlibrary{shadows}
\pgfPT[background={left color=red!10,right color=green!10,postaction={drop shadow={left color=red!10,right color=green!10}}}]
```



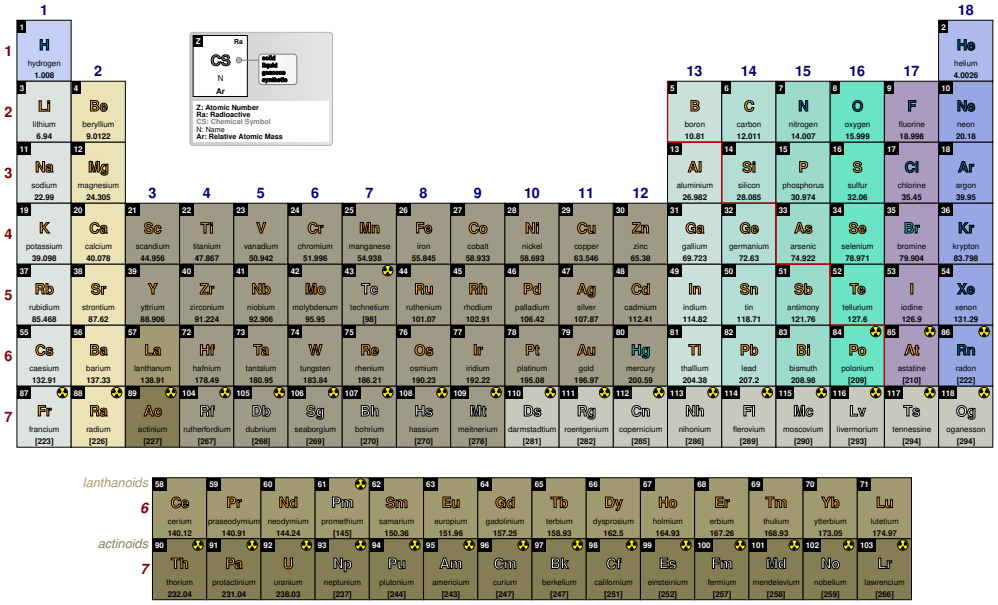
IUPAC default: *true*

When set to true draws the periodic table with *lanthanum* and *actinium* appended to block f and the labels *lanthanoids* and *actinoids* are placed at group 3, substituting *lanthanum* and *actinium*. When IUPAC is set to false, *lanthanum* and *actinium* are shown in group 3 and the labels *lanthanoids* and *actinoids* are place near the *f block* (if the key show label LaAc is set to true).

\pgfPT



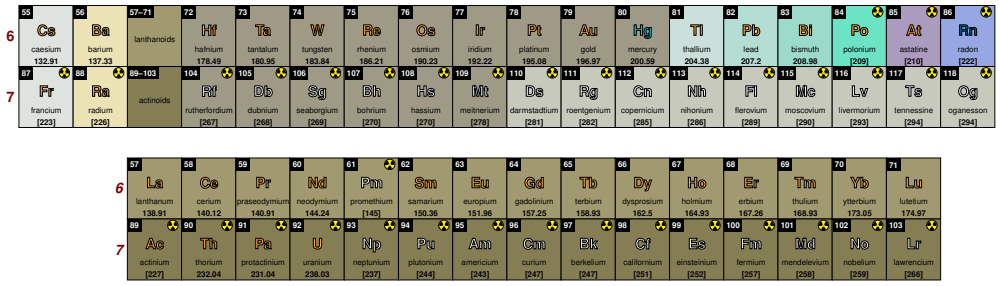
\pgfPT[IUPAC=false]



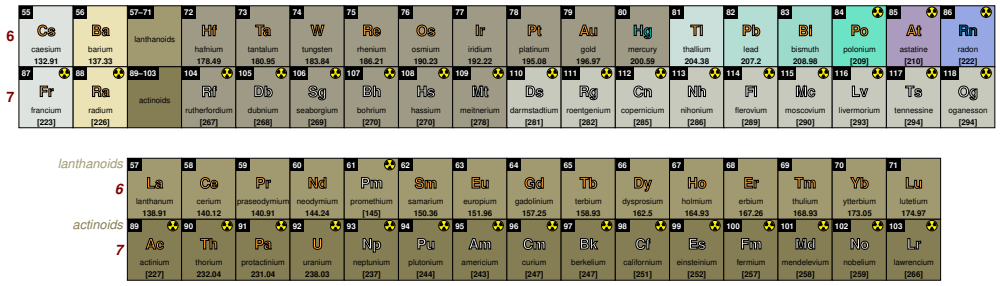
show label LaAc default: {}

Determines when the labels 'lanthanoids' and 'actinoids' are shown (true) or not shown (false) near the f block. When the IUPAC key is set to true, the default behavior is to show the labels and when the IUPAC key is set to false, the default behavior is to hide the labels. This *default behavior* **can be overridden by this key** setting it to true, to show the labels, or to false to hide them, independently of the value of the IUPAC key.

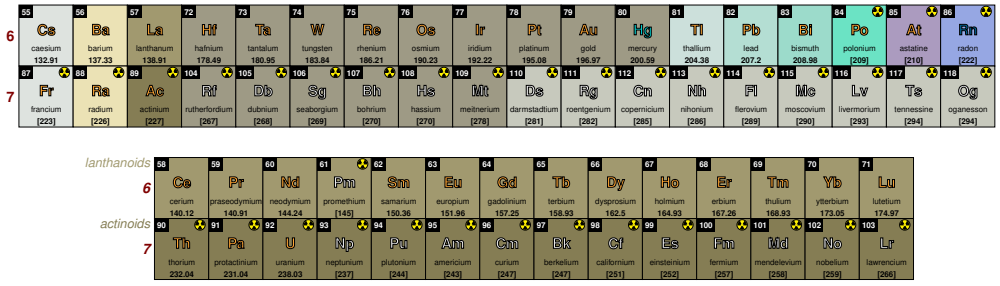
```
\pgfPTnewZlist{myZlist}{55,...,118}
\pgfPTstyle[show title=false,show legend=false,show group numbers=false]
\pgfPT[Z list=myZlist]
```



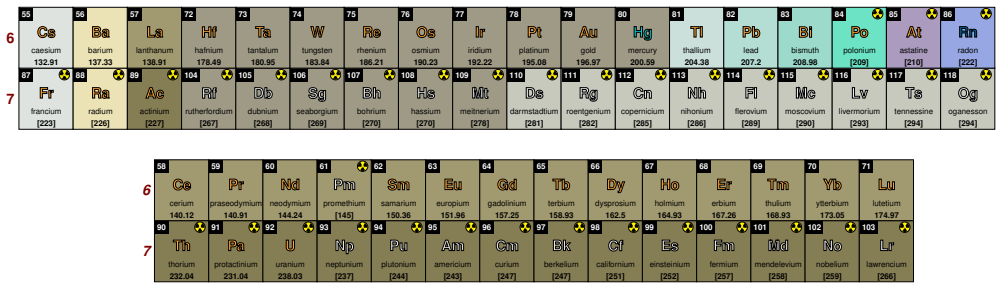
```
\pgfPT[Z list=myZlist,show label LaAc=true]
```



```
\pgfPT[Z list=myZlist,IUPAC=false]
```



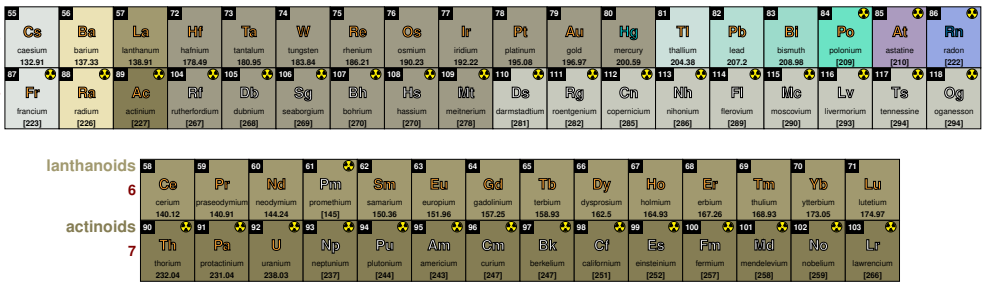
```
\pgfPT[Z list=myZlist,IUPAC=false,show label LaAc=false]
```



label LaAc font default: *\footnotesize\itshape*

Sets the font for the labels 'lanthanoids' and 'actinoids'.

```
\pgfPT[label LaAc font=\bfseries,Z list=myZlist,IUPAC=false]
```

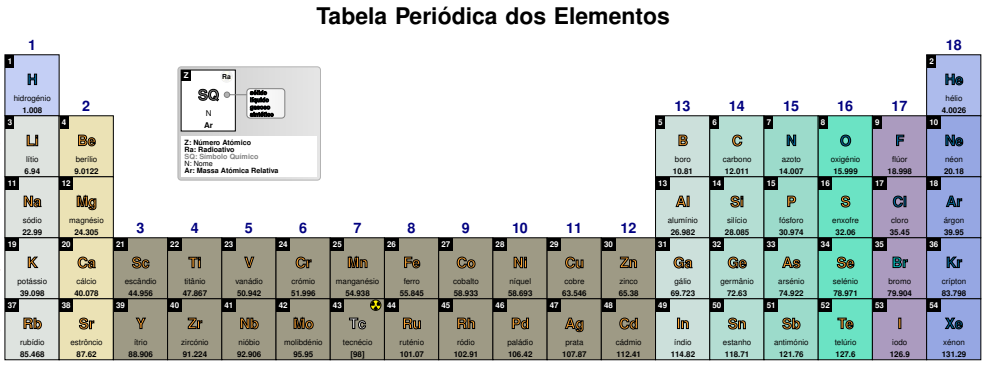


```
\pgfPTresetstyle
```

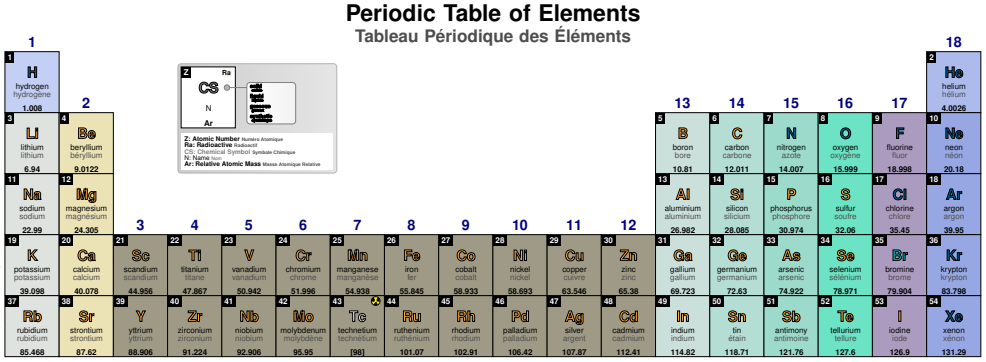
languages default: *{}*

Sets a language list to use in the Periodic Table. It is a comma separated list of language flags: 'pt', 'en', 'fr', 'de', 'it', 'es' or 'br'. *This key overrides the default language, that is, the language loaded at package inclusion*.

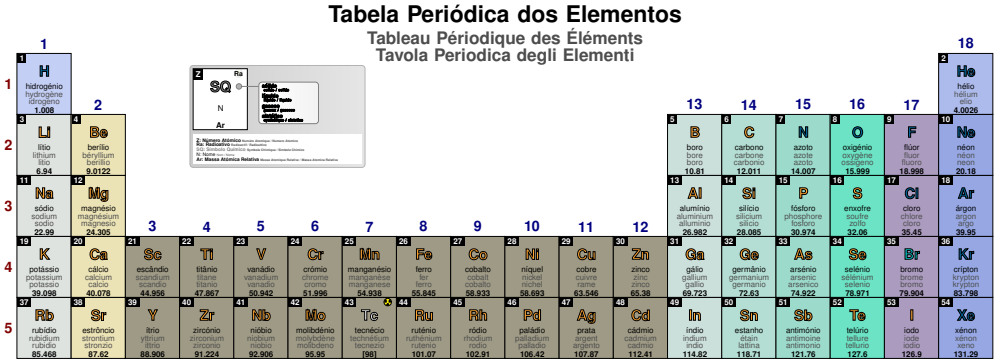
```
\pgfPT[Z list={1,...,54},languages=pt]
```




```
\pgfPT[Z list={1,...,54},cell style=pgfPT2lang,languages={en,fr}]
```

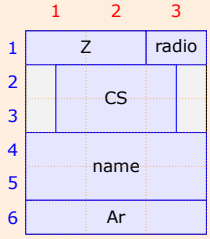


\pgfPT[Z list={1,...,54},cell style=pgfPT3lang,languages={pt,fr,it}]

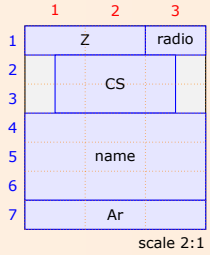


When using a set of languages, space to accommodate the names in each cell must be provided by building a suitable cell - typically one cell row per language. The cell styles used in the two examples above are built-in and serve this purpose.

✓ *Built-in style* **pgfPT2lang**
The build command:
\pgfPTbuildcell(6,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4-5;1-3;name),(6;1-3;Ar)]



✓ *Built-in style* **pgfPT3lang**
The build command:
\pgfPTbuildcell(7,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4-6;1-3;name),(7;1-3;Ar)]

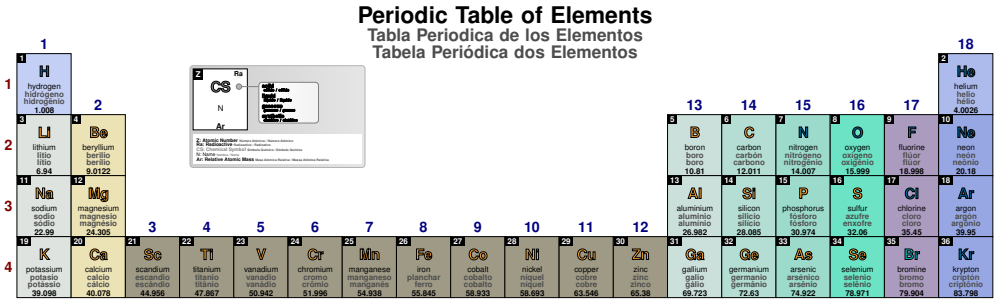


Also, the space for the title should be taken into account – if using more then three languages, the legend must be *turned off*, otherwise the title overlaps the legend.

other languages font default: *\tiny*

Sets the font used in *other languages*, *i.e.*, the languages started at the second entry of the list provide to the languages key.

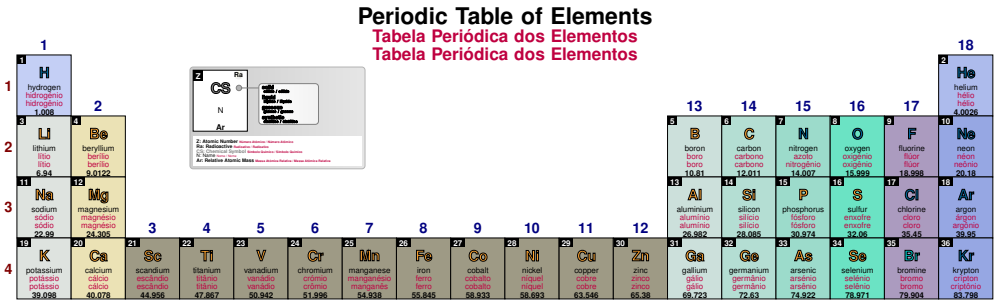
```
\pgfPT[Z list={1,...,36},cell style=pgfPT3lang,languages={en,es,br}, other languages font=\tiny\bfseries]
```



other languages color default: *black!70*

Sets the color of the font used in *other languages*.

```
\pgfPT[Z list={1,...,36},cell style=pgfPT3lang,languages={en,pt,br}, other languages color=purple]
```

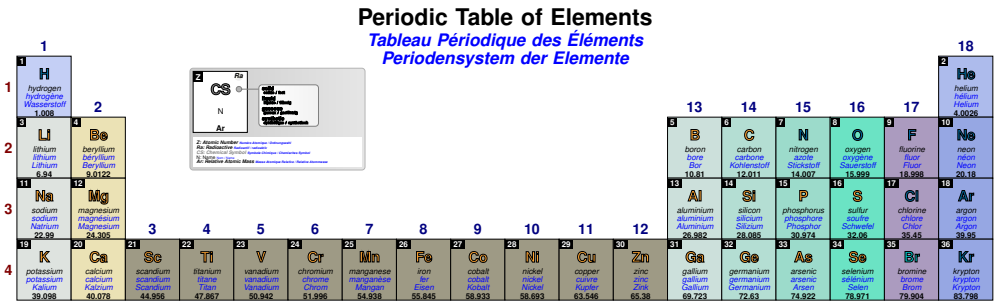


other lang default: *{f=\tiny,c=black!70}*

Pseudo style to set the keys: other languages **f**ont and/or other languages **c**olor. None of the *keys* – f and c – are mandatory.

USAGE: other lang={f=<font commands>,c=<color>}

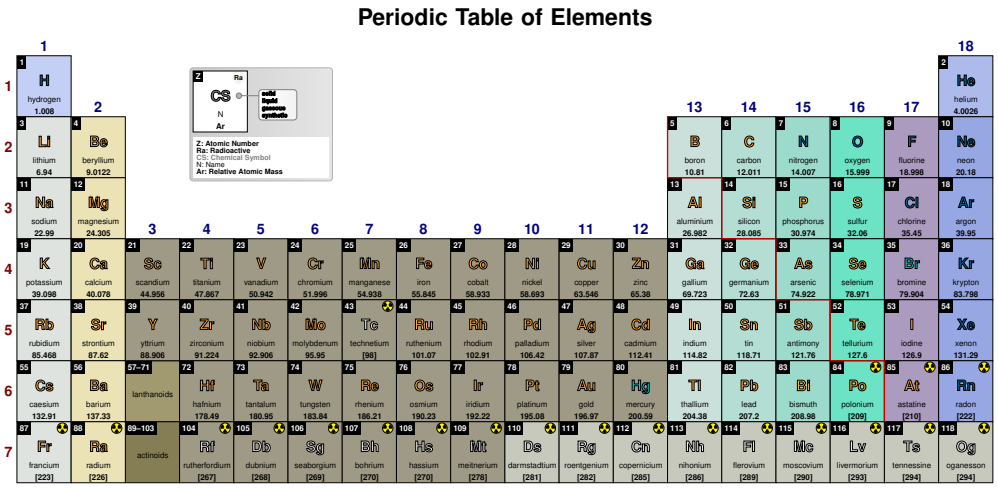
```
\pgfPT[Z list={1,...,36},cell style=pgfPT3lang,languages={en,fr,de}, other lang={f=\tiny\itshape,c=blue}]
```



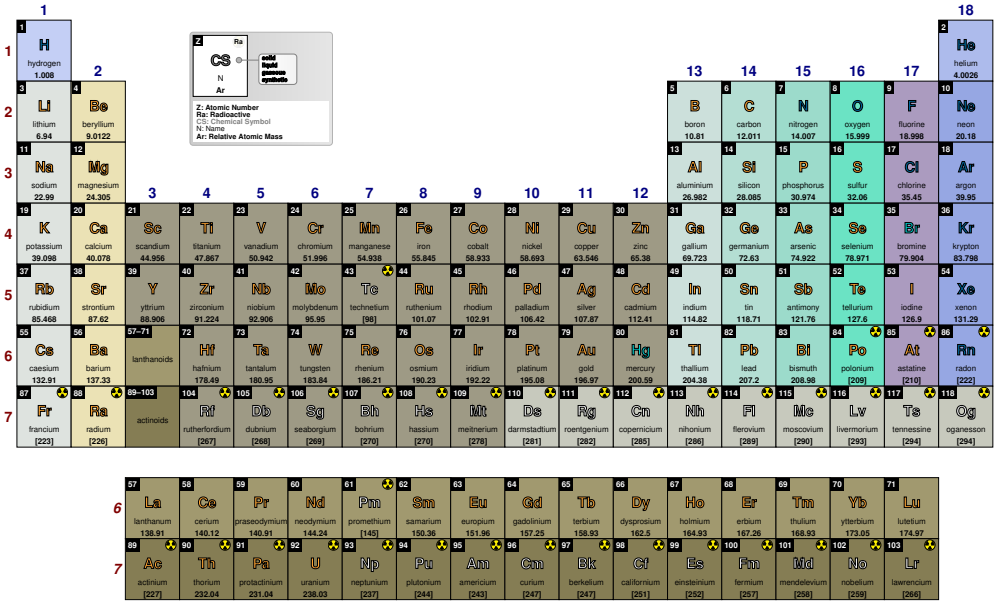
show MNM line default: *true*

If set to true a line separating metals from non metals is shown in the Periodic Table. The line starts at the upper left corner of the cell of boron (2nd period, group 13) and ends at the lower right corner of polonium (6th period, group 16). If set to false no line is drawn.

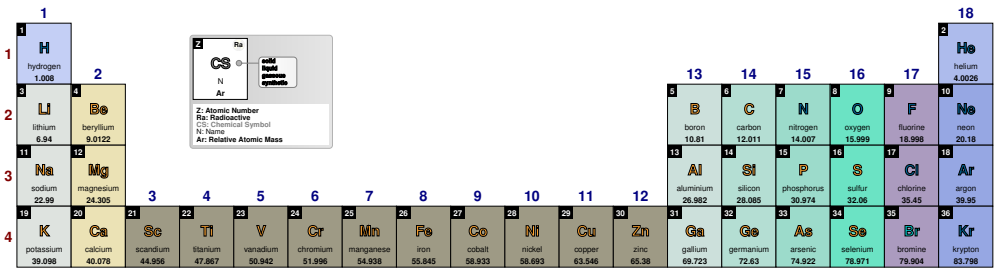
\pgfPT[Z list=spd]



\pgfPT[show MNM line=false]



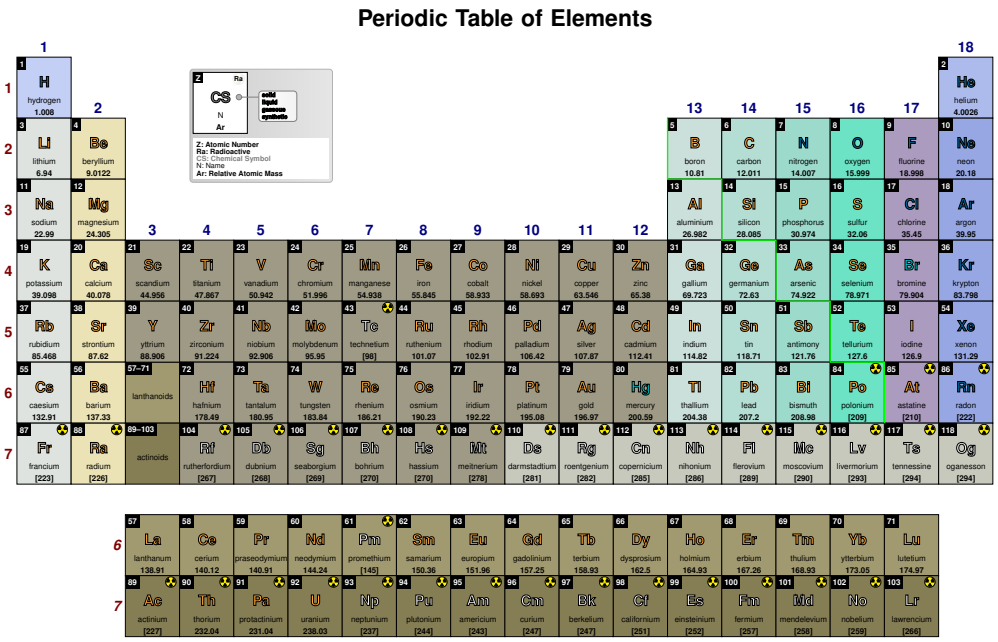
\pgfPT[Z list={1,...,36}]



MNM line color default: *red!80!black*

Sets the color of the *MNM line*.

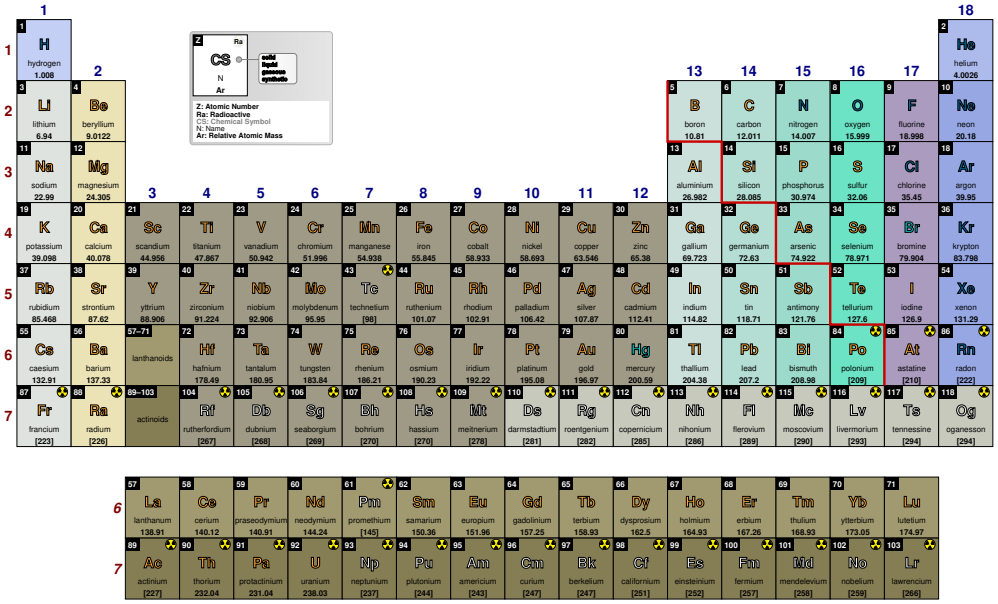
```
\pgfPT[MNM line color=green]
```



MNM line width default: *.8pt*

Sets the width of the *MNM line*.

```
\pgfPT[MNM line width=1.5pt]
```

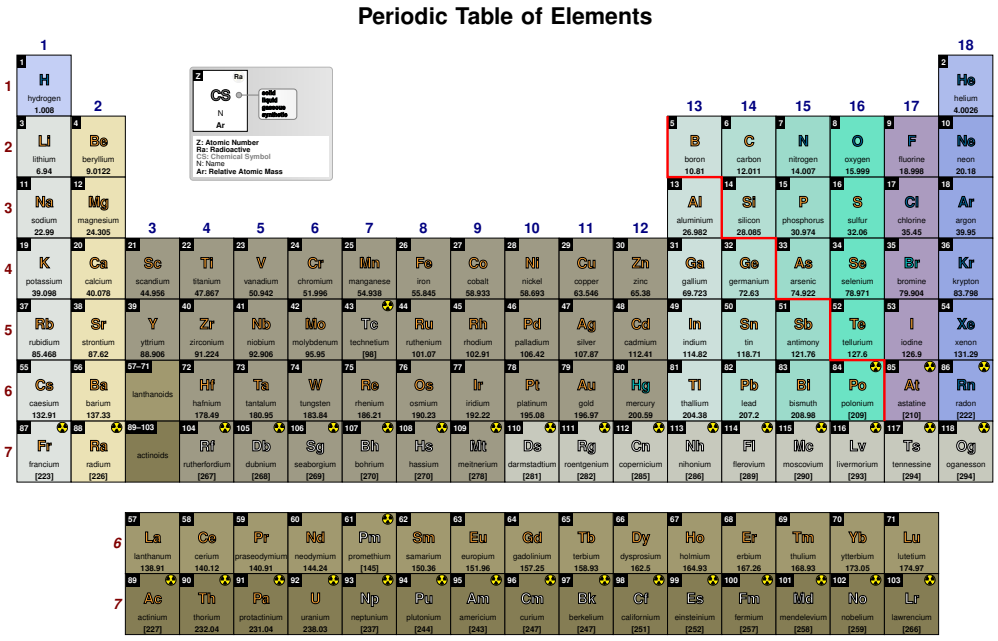


MNM default: *{c=red!80!black,w=.8pt}*

Pseudo style to set the *MNM line* **c**olor and/or **w**idth. None of the *keys* – c and w – are mandatory. The key show MNM line is set to true.

USAGE: MNM={c=<color>,w=<length>}

```
\pgfPT[MNM={w=1.5pt,c=red}]
```

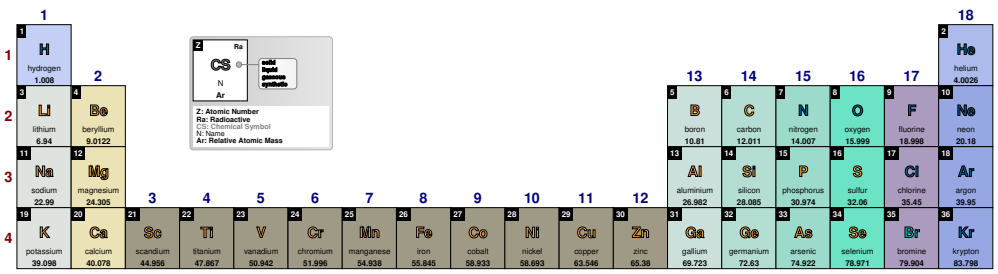


➠ Title and Legend

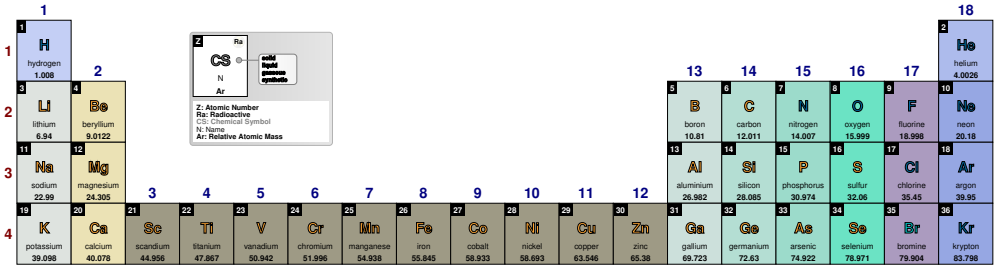
show title default: *true*

When set to true the title is shown, otherwise the title (Periodic Table of elements) is not shown.

```
\pgfPT[Z list={1,...,36}]
```



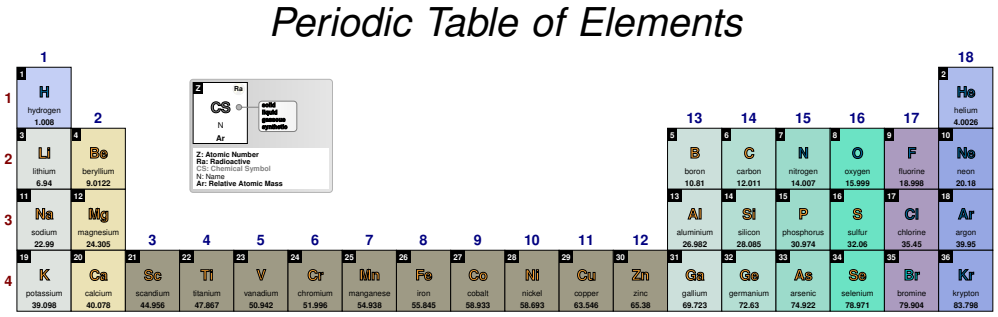
```
\pgfPT[Z list={1,...,36},show title=false]
```



title font default: *\Large\bfseries*

Sets the font used in the title.

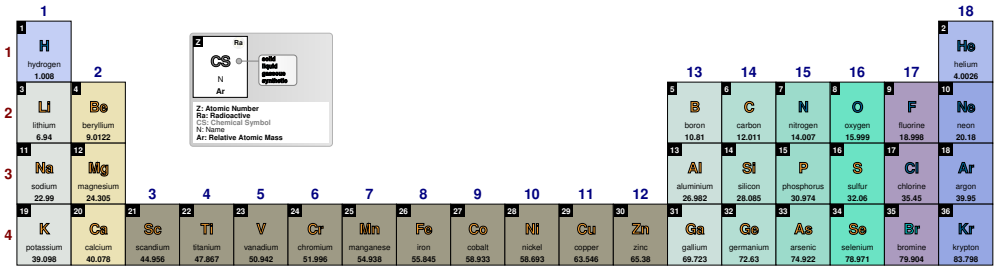
```
\pgfPT[Z list={1,...,36},title font=\Huge\itshape]
```



title color default: *black*

Sets the title color.

```
\pgfPT[Z list={1,...,36},title color=green!50!black]
```

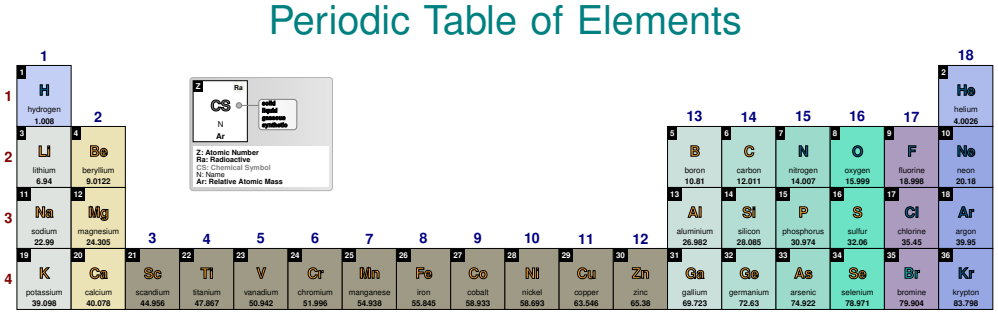


title default: *{f=\Large\bfseries,c=black}*

Pseudo style to set the keys: title **f**ont and/or title **c**olor. None of the *keys* – f and c – are mandatory. The key show title is set to true.

USAGE: title={f=<font commands>,c=<color>}

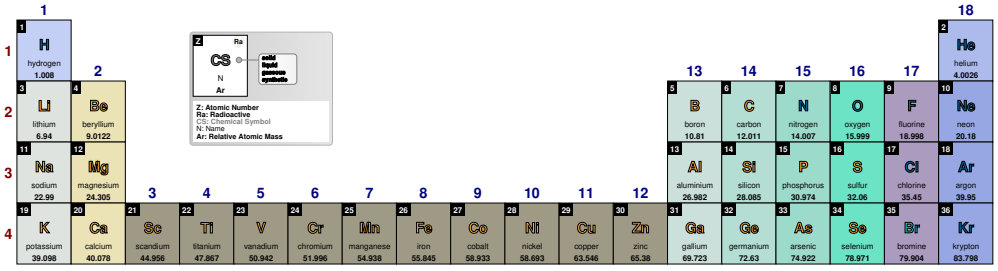
```
\pgfPT[Z list={1,...,36},title={f=\Huge,c=teal}]
```



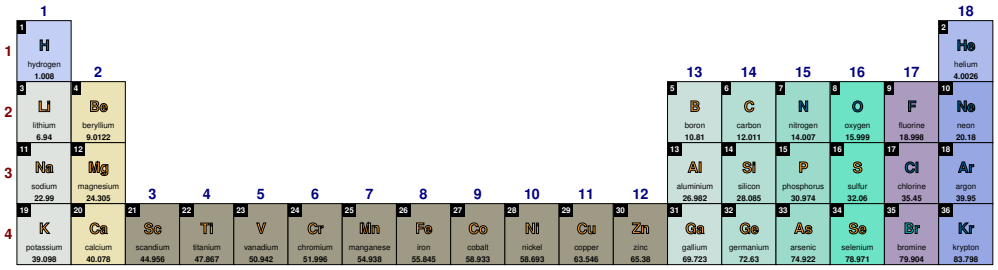
show legend default: *true*

When set to true the legend is shown, otherwise it is not shown.

```
\pgfPT[Z list={1,...,36}]
```



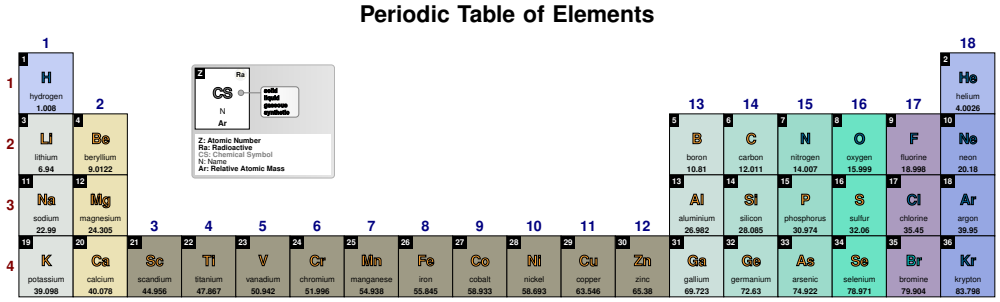

```
\pgfPT[Z list={1,...,36},show legend=false]
```



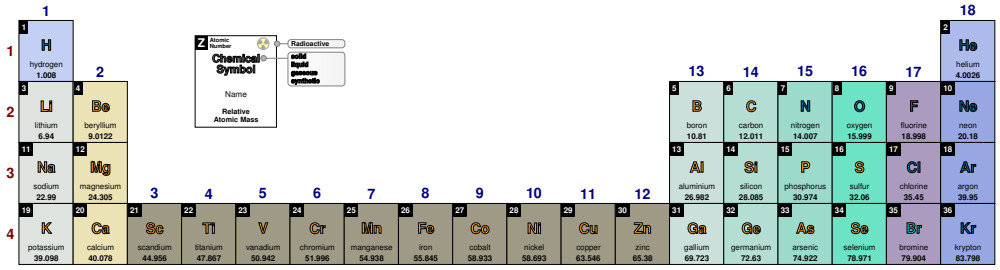
legend acronyms default: *true*

When set to true, the legend consists of a cell using acronyms for its contents and the corresponding descriptions below that cell. When set to false, only the cell is displayed with the descriptions in place of the acronyms. In the latter case, the description font size is automatically adjusted to the available box, which can *spoil the appearance of the whole caption*, depending on the described content.

```
\pgfPT[Z list={1,...,36}]
```



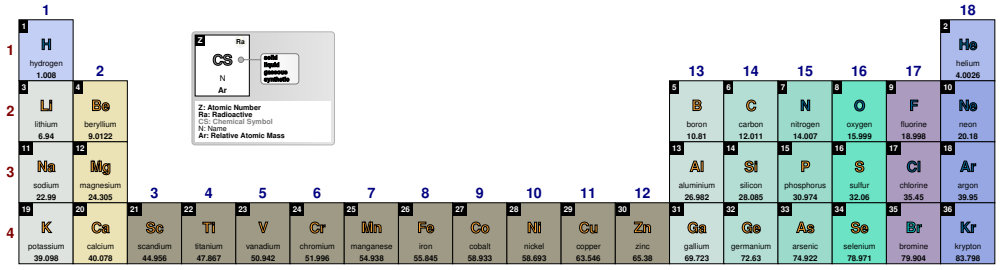

```
\pgfPT[Z list={1,...,36},legend acronyms=false]
```



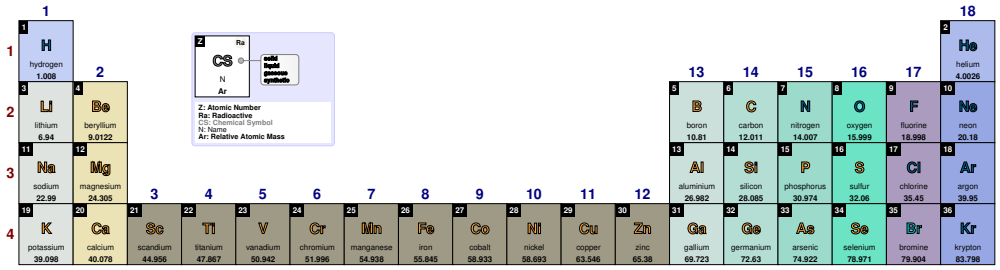
legend box default: *left color=black!20,right color=black!10,draw=black!30*

Style to define the appearance of the box around the legend, legend pins and acronym descriptions, built with any of the TikZ keys that can be applied to a path construction. *It only works when the key legend acronyms is set to true.*

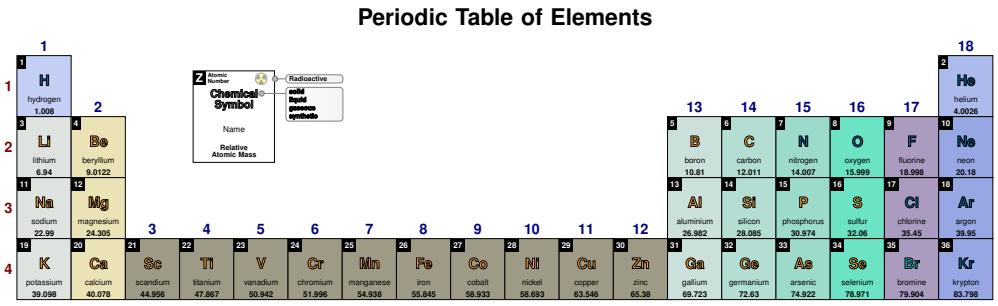
```
\pgfPT[Z list={1,...,36}]
```



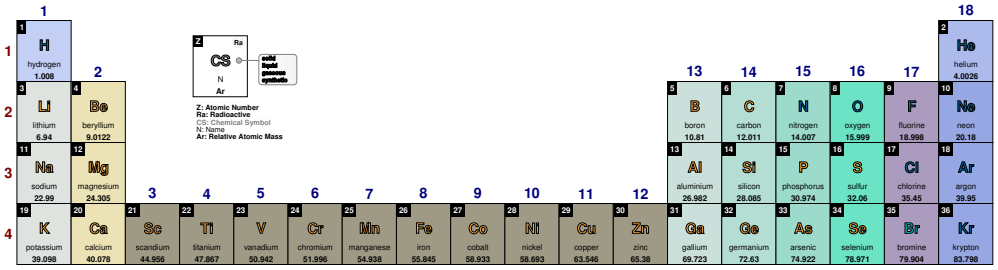

```
\pgfPT[Z list={1,...,36},legend box={draw=blue!20,fill=blue!10}]
```



```
\pgfPT[Z list={1,...,36},legend box={draw=blue!20,fill=blue!10,legend acronyms=false}]
```



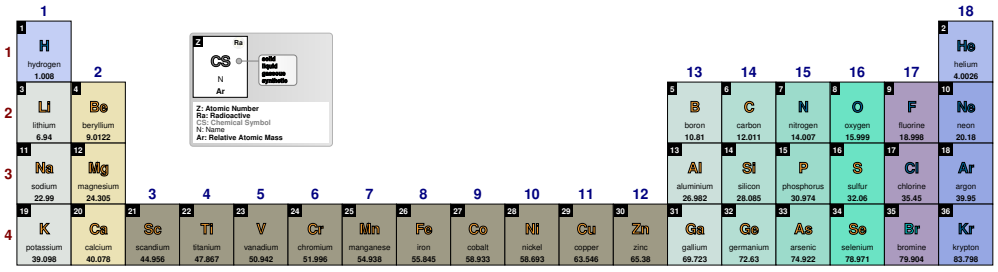

```
\pgfPT[Z list={1,...,36},legend box={}]
```



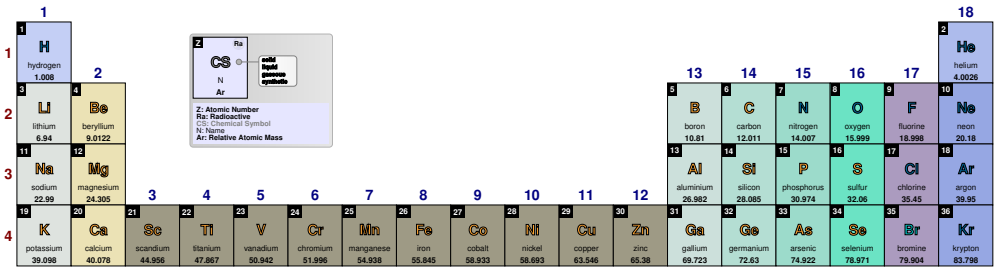
legend back color default: *white*

Sets the legend background color.

```
\pgfPT[Z list={1,...,36}]
```



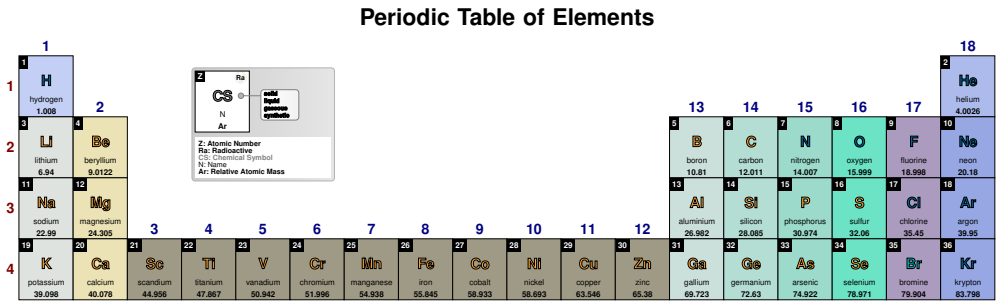

```
\pgfPT[Z list={1,...,36},legend back color=blue!10]
```



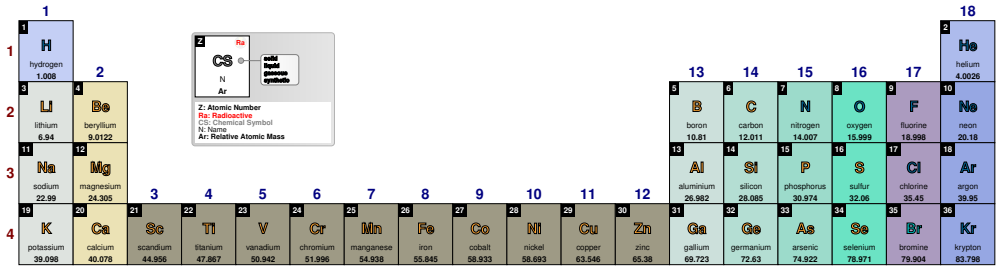
legend radio color default: *black*

Sets the color of the radioactivity acronym and corresponding description.

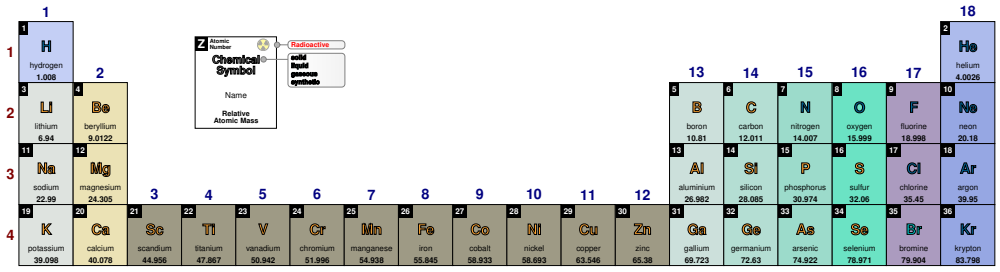
```
\pgfPT[Z list={1,...,36}]
```




```
\pgfPT[Z list={1,...,36},legend radio color=red]
```



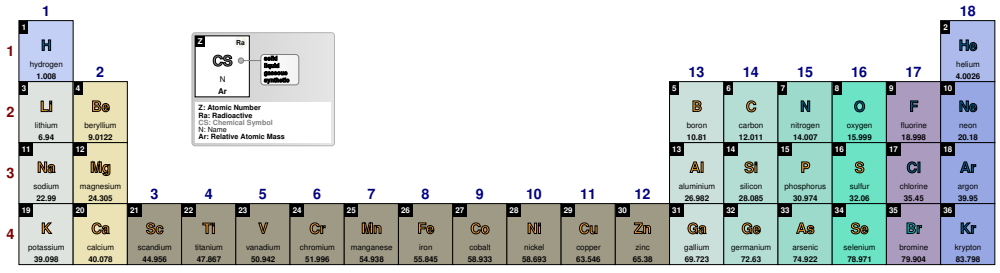

```
\pgfPT[Z list={1,...,36},legend radio color=red,legend acronyms=false]
```



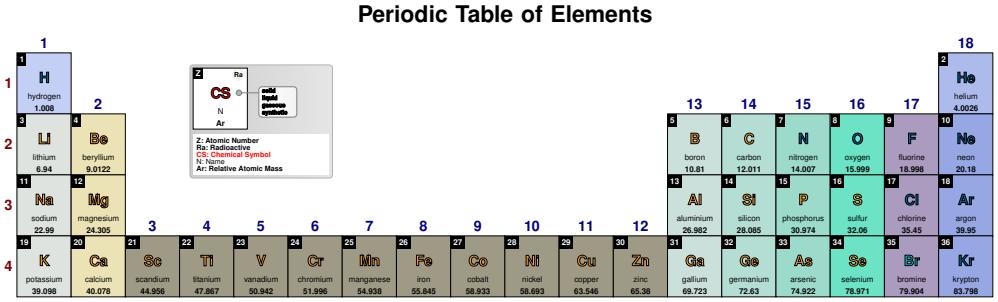
legend CS color default: *black!50*

Sets the color of the Chemical Symbol acronym and corresponding description.

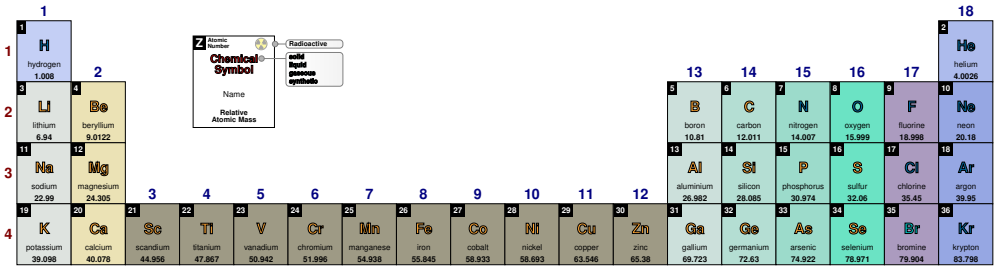
```
\pgfPT[Z list={1,...,36}]
```



\pgfPT[Z list={1,...,36},legend CS color=red]



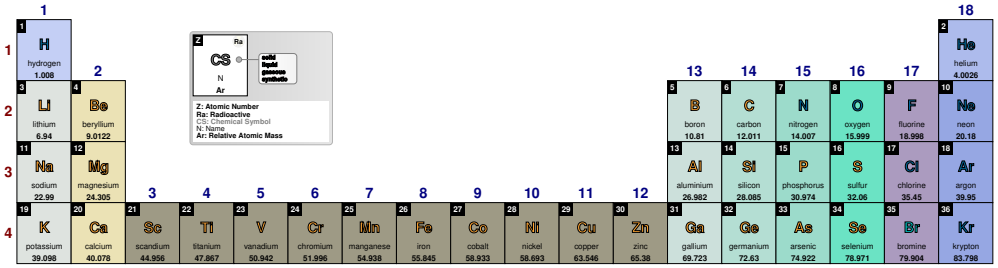
\pgfPT[Z list={1,...,36},legend CS color=red,legend acronyms=false]



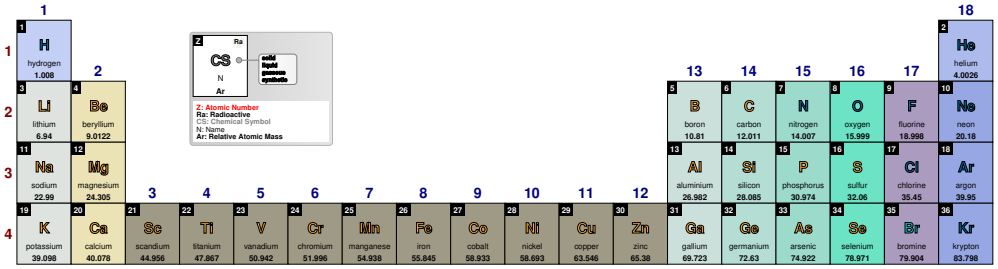
legend Z color default: {}

Sets the color of the atomic number description (only applies when the key legend acronyms is set to true.)

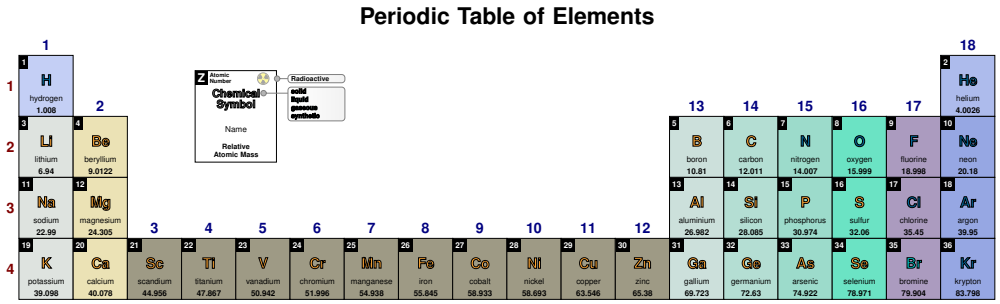
\pgfPT[Z list={1,...,36}]



\pgfPT[Z list={1,...,36},legend Z color=red]



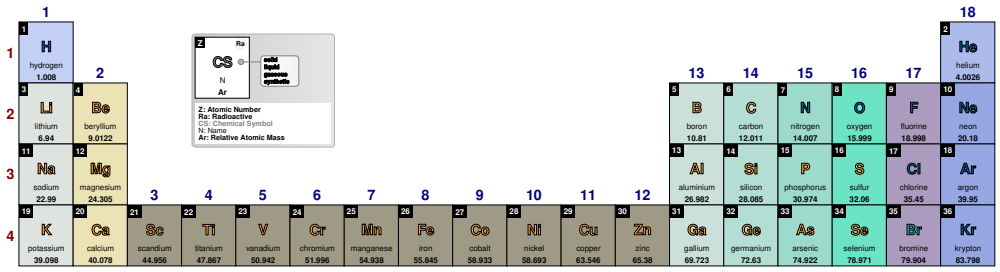
\pgfPT[Z list={1,...,36},legend Z color=red,legend acronyms=false]



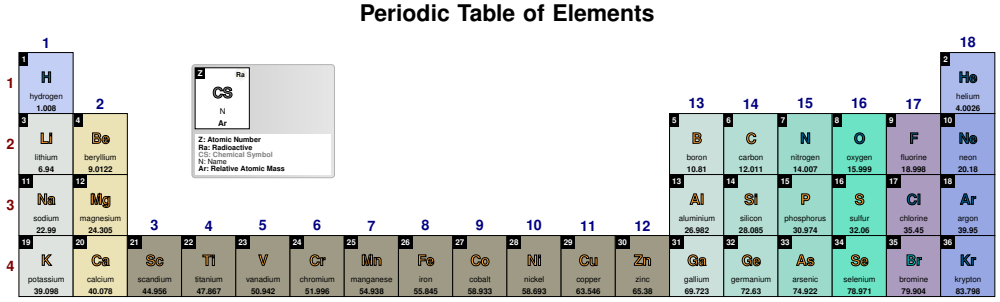
show legend pins default: *true*

When set to true the legend pins are shown, otherwise they are not shown.

```
\pgfPT[Z list={1,...,36}]
```



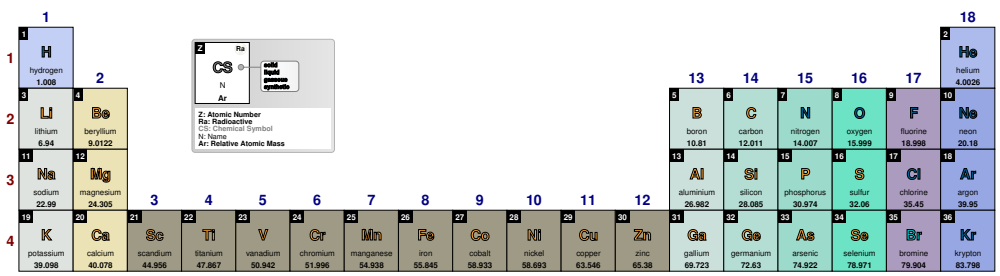

```
\pgfPT[Z list={1,...,36},show legend pins=false]
```



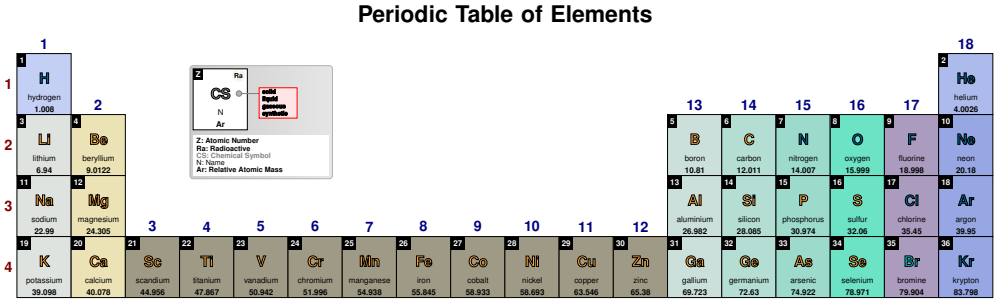
legend pins default: *{line width=.05pt,rounded corners=2pt,right color=black!5, left color=white,draw=black!50}*

Style to define the appearance of the legend pins, built with any of the TikZ keys that can be applied to a path construction.

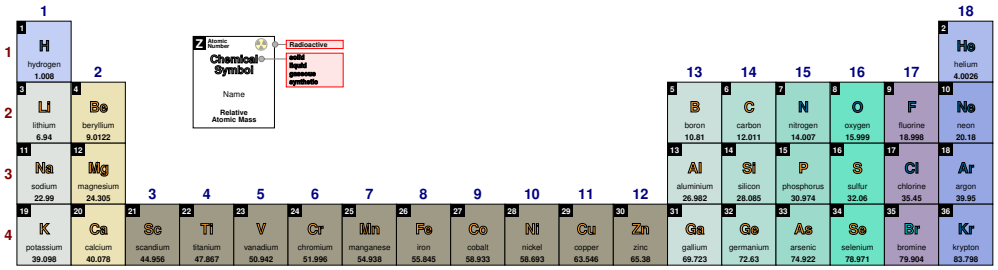
```
\pgfPT[Z list={1,...,36}]
```



```
\pgfPT[Z list={1,...,36},legend pins={draw=red,fill=red!10}]
```



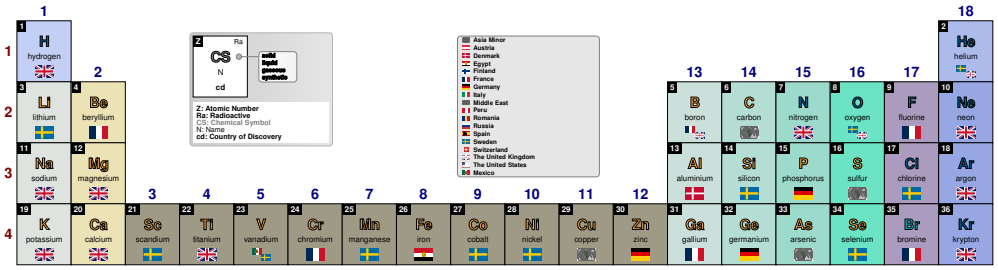

```
\pgfPT[Z list={1,...,36},legend pins={draw=red,fill=red!10},legend acronyms=false]
```



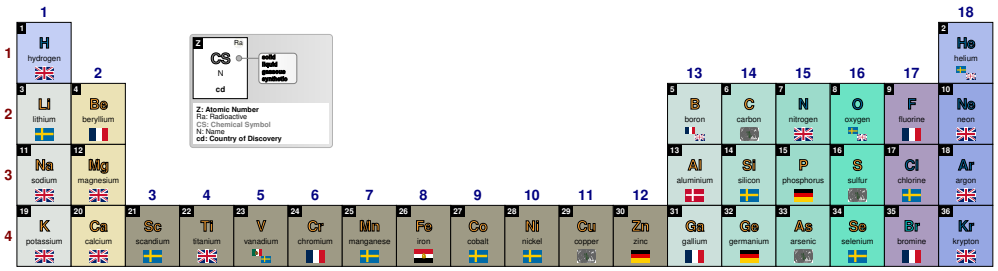
show extra legend default: *true*

When set to true the extra legend is shown, otherwise it is not shown.

```
\pgfPTbuilcellstyle{myname}(6,3)% 6 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5.25-6.75;1-3;DiscC)]
\pgfPT[Z list={1,...,36},cell style=myname]
```



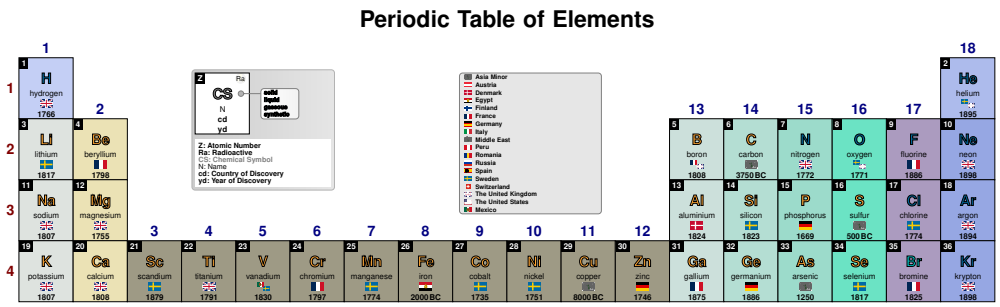

```
\pgfPT[Z list={1,...,36},cell style=myname,show extra legend=false]
```



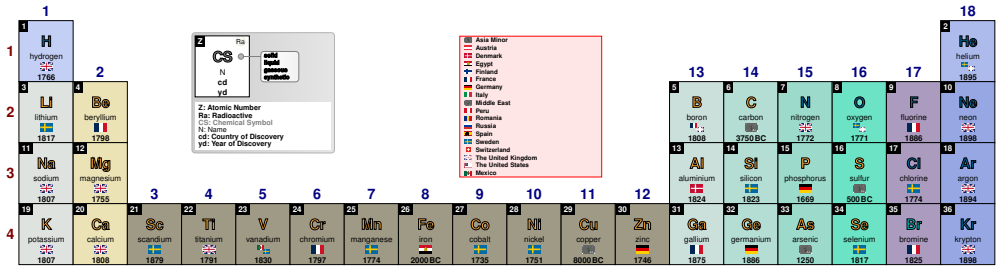
extra legend default: *{draw=black!50,fill=black!10,line width=.05pt, rounded corners=2pt}*

Style to define the appearance of the extra legend, built with any of the TikZ keys that can be applied to a path construction.

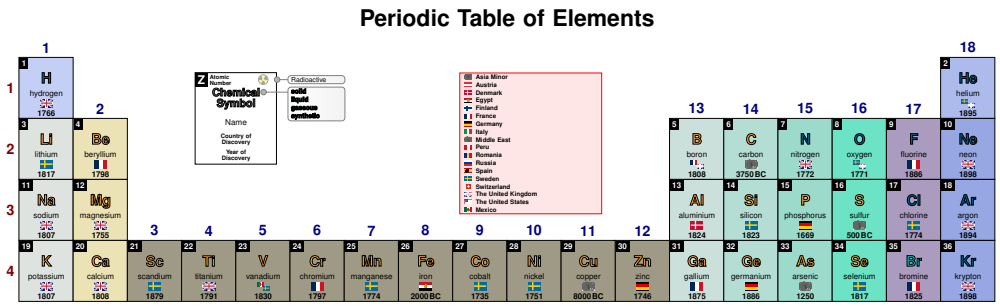
```
\pgfPT[Z list={1,...,36},cell style=pgfPTdisc]
```




```
\pgfPT[Z list={1,...,36},cell style=pgfPTdisc,extra legend={draw=red,fill=red!10}]
```




```
\pgfPT[Z list={1,...,36},cell style=pgfPTdisc,legend acronyms=false,extra
legend={draw=red,fill=red!10}]
```

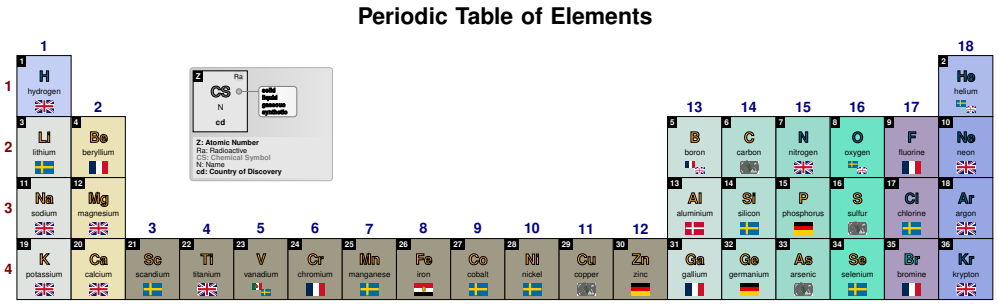


legend default: *{bc=white,pins=true,extra=true,acro=true}*

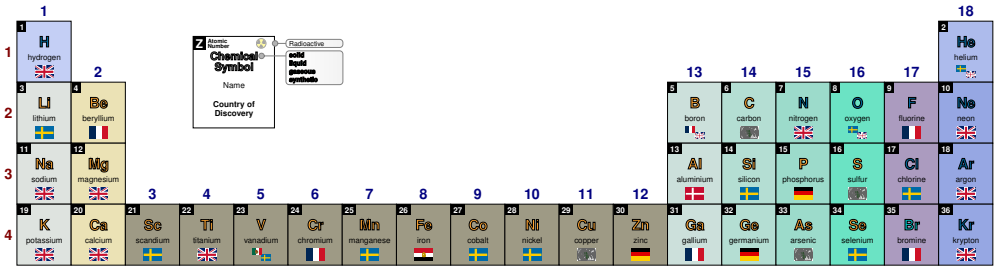
Pseudo style to set the keys: legend **b**ack **c**olor, show legend **pins**, show **extra** legend, legend **acro**nyms, legend **radio** color, legend **CS** color, legend **Z** color, legend **pins** (**style**), **extra** legend (**style**) and/or legend **box** (style). None of the *keys* – bc, pins, extra, acro, radio, CS, Z, pins style, extra style and box – are mandatory. The key show legend is set to true.

USAGE:	legend={bc=<color>,pins=<true|false>,extra=<true|false>,acro=<true|false>, radio=<color>,CS=<color>,Z=<color>,pins style=<tikz path keys>, extra style=<tikz path keys>,box=<tikz path keys>}

```
\pgfPT[Z list={1,...,36},cell style=myname,legend={bc=black!10,extra=false}]
```




```
\pgfPT[Z list={1,...,36},cell style=myname,legend={acro=false,extra=false}]
```

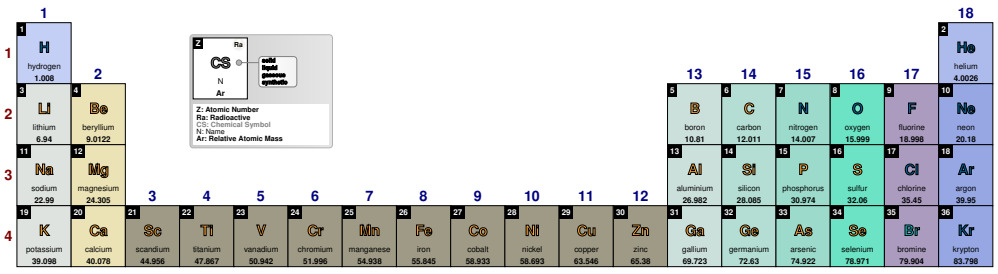


➠ Periods and Groups

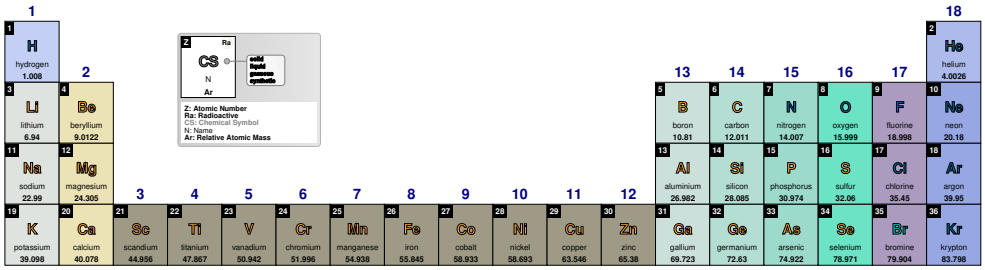
show period numbers default: *true*

When set to true the period numbers are shown, otherwise they are not shown.

```
\pgfPT[Z list={1,...,36}]
```



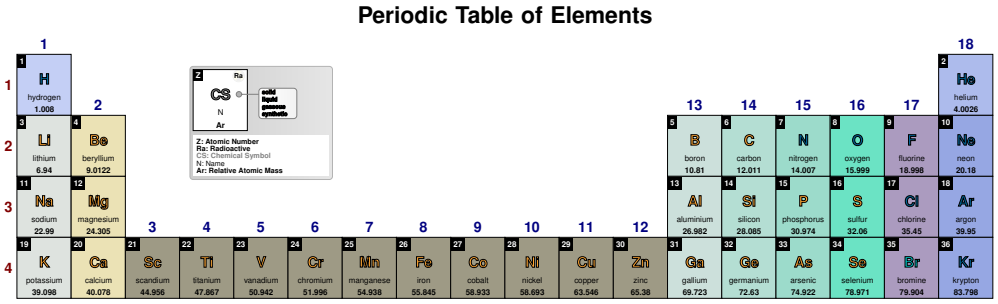

```
\pgfPT[Z list={1,...,36},show period numbers=false]
```



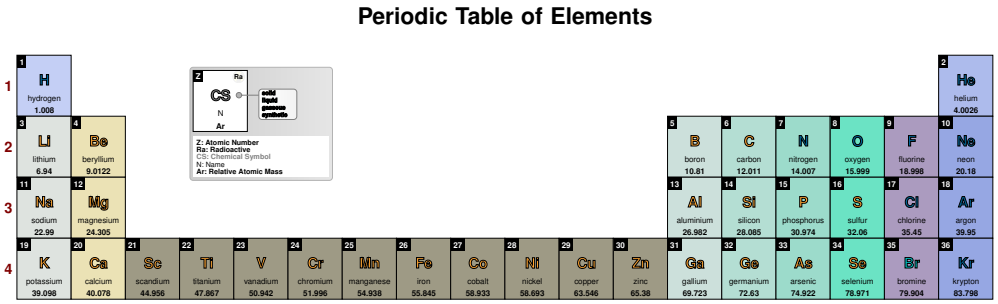
show group numbers default: *true*

When set to true the group numbers are shown, otherwise they are not shown.

```
\pgfPT[Z list={1,...,36}]
```



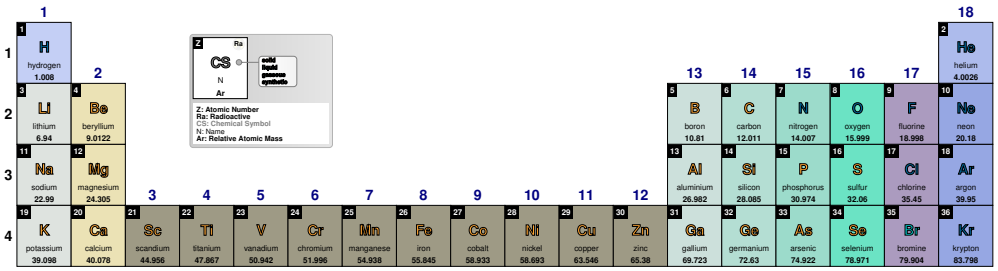

```
\pgfPT[Z list={1,...,36},show group numbers=false]
```



period label color default: *red!50!black*

Sets the period label color.

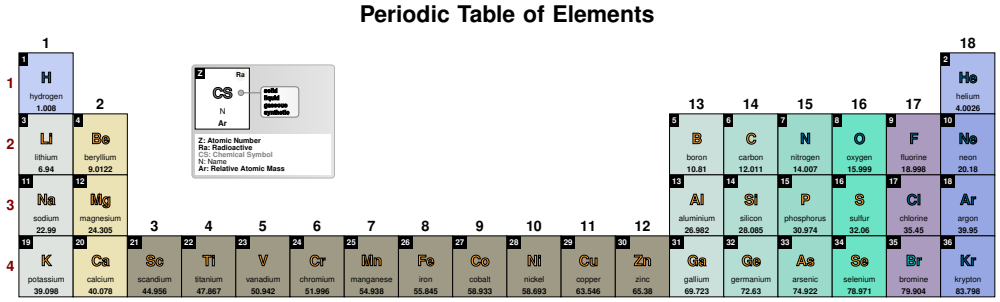
```
\pgfPT[Z list={1,...,36},period label color=black]
```



group label color default: *blue!50!black*

Sets the group label color.

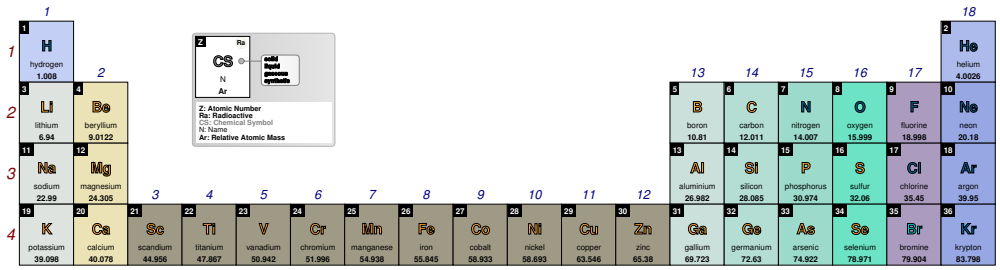
```
\pgfPT[Z list={1,...,36},group label color=black]
```



label font default: *\small\bfseries*

Sets the label font.

```
\pgfPT[Z list={1,...,36},label font=\itshape]
```

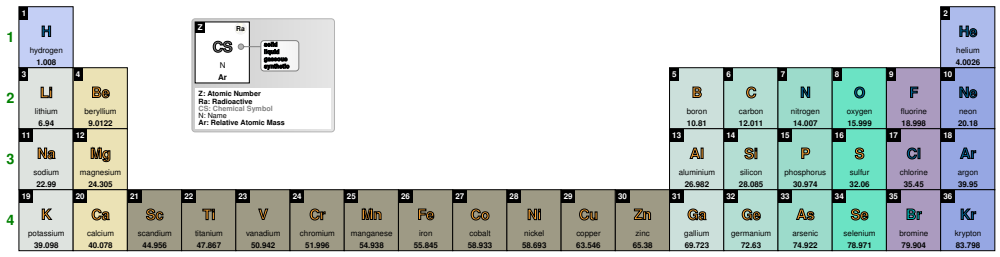


per default: *{gr=true,c=red!50!black,f=\small\bfseries}*

Pseudo style to set the keys: show **gr**oup numbers, period label **c**olor and/or label **f**ont. None of the *keys* – gr, c and f – are mandatory. The key show period numbers is set to true.

USAGE: per={gr=<true|false>,c=<color>,f=<font commands>}

```
\pgfPT[Z list={1,...,36},per={gr=false,c=green!50!black}]
```

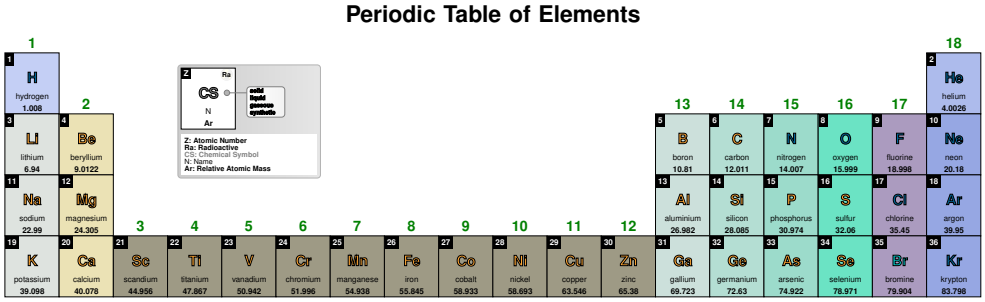


gr default: *{per=true,c=blue!50!black,f=\small\bfseries}*

Pseudo style to set the keys: show **per**iod numbers, group label **c**olor and/or label **f**ont. None of the *keys* – per, c and f – are mandatory. The key show group numbers is set to true.

USAGE: gr={per=<true|false>,c=<color>,f=<font commands>}

```
\pgfPT[Z list={1,...,36},gr={per=false,c=green!50!black}]
```

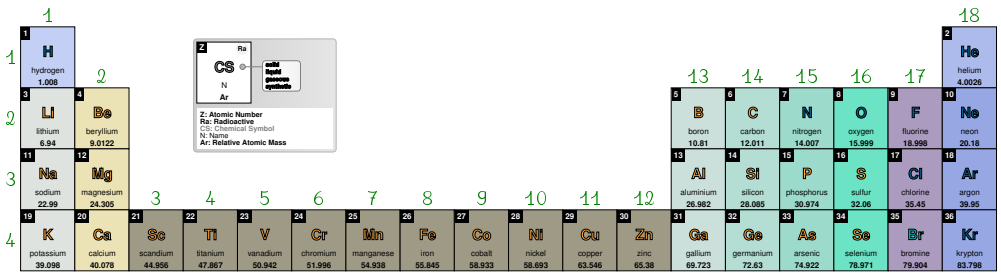


per+gr default: *{pc=red!50!black,gc=blue!50!black,f=\small\bfseries}*

Pseudo style: use **c** to set both keys group label color and period label color with the same color; use **pc** to set period label color, **gc** to set group label color and/or **f** to set label **f**ont. None of the *keys* – c, pc, gc and f – are mandatory. The keys show period numbers and show group numbers are set to true.

USAGE: per+gr={c=<color>,pc=<color>,gc=<color>,f=<font commands>}

```
\pgfPT[Z list={1,...,36},per+gr={c=green!50!black,
f=\fontfamily{frc}\selectfont\normalsize\bfseries}]
```

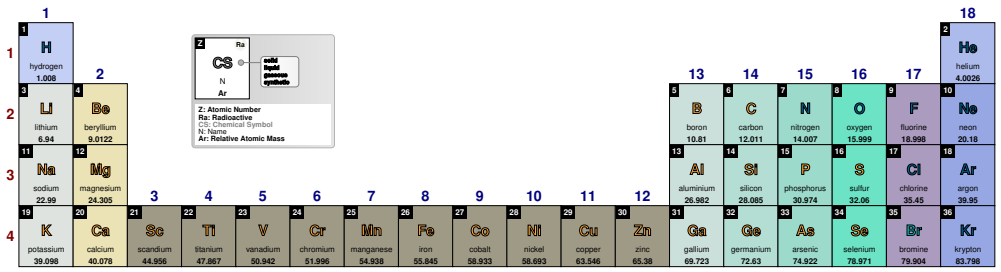


➠ Blocks and Families

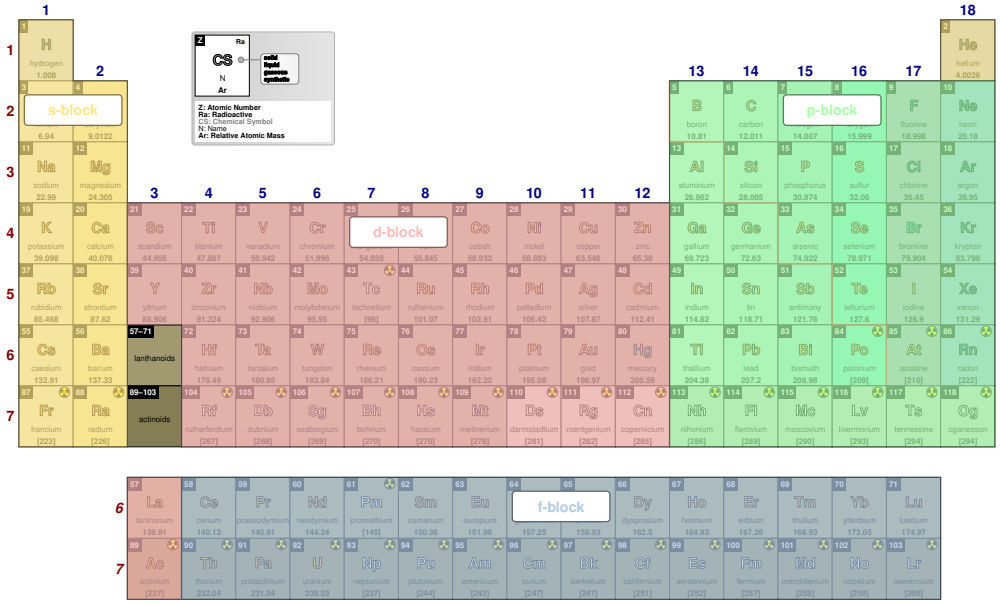
show blocks default: *false*

When set to true the blocks s, p, d and f are drawn overlaying the Periodic Table and their labels are shown. *Note that blocks are only shown when the Z list contains, at least, all elements of blocks s, p and d.*

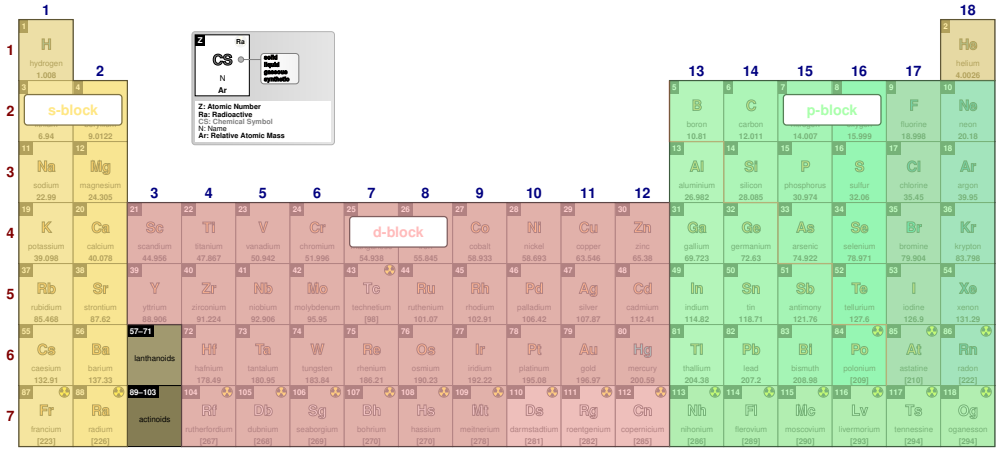
```
\pgfPT[Z list={1,...,36},show blocks=true,show title=false]
```



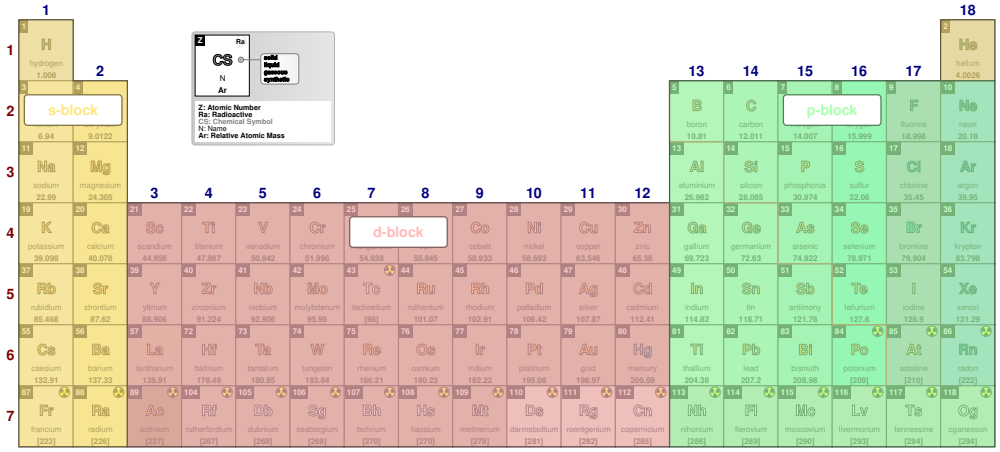
\pgfPT[show blocks,show title=false]



\pgfPT[Z list=spd,show blocks,show title=false]



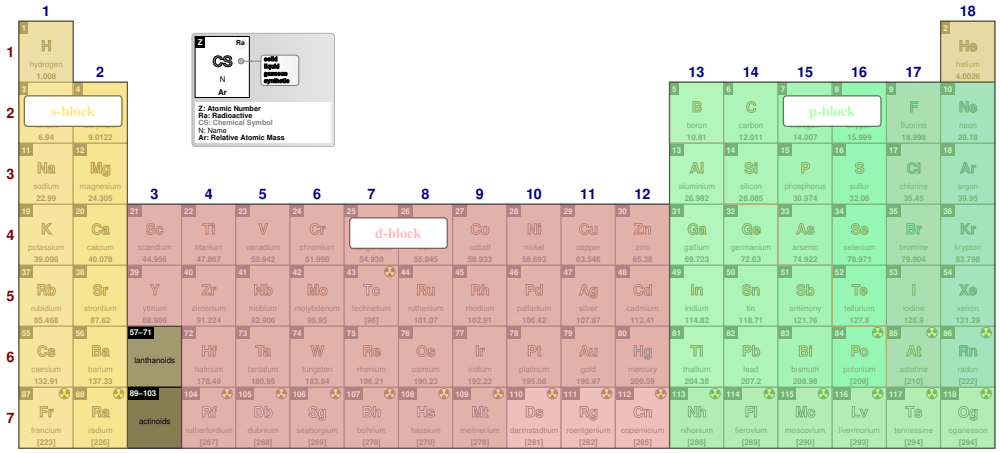
\pgfPT[Z list=spd,show blocks,show title=false,IUPAC=false]



blocks font default: *\small\bfseries*

Sets the font used in the block labels.

```
\pgfPT[Z list=spd,show blocks,show title=false,blocks
font=\small\bfseries\fontfamily{ptm}\selectfont]
```



s block color default: RGB: 255,231,132

Sets the block s color.

s block font color default: *{}*

Sets the s block label font color. If no color is provided, the s block color will be used as the font color.

s block line width default: *0.8pt*

Sets the width of the line surrounding the s block.

p block color default: RGB: 170,255,172

Sets the block p color.

p block font color default: *{}*

Sets the p block label font color. If no color is provided, the p block color will be used as the font color.

p block line width default: *0.8pt*

Sets the width of the line surrounding the p block.

d block color default: RGB: 255,187,187

Sets the block d color.

d block font color default: *{}*

Sets the d block label font color. If no color is provided, the d block color will be used as the font color.

d block line width default: *0.8pt*

Sets the width of the line surrounding the d block.

f block color default: RGB: 177,203,228

Sets the block f color.

f block font color default: *{}*

Sets the f block label font color. If no color is provided, the f block color will be used as the font color.

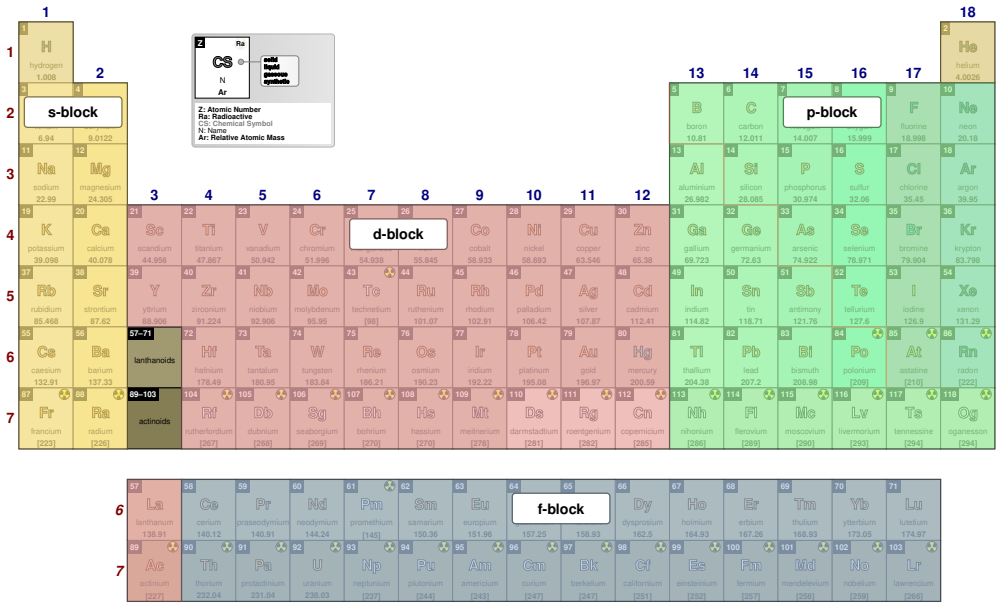
f block line width default: *0.8pt*

Sets the width of the line surrounding the f block.

blocks font color default: *black*

Style to set a common color for the labels of s, p, d and f blocks. The key show blocks is set to true.

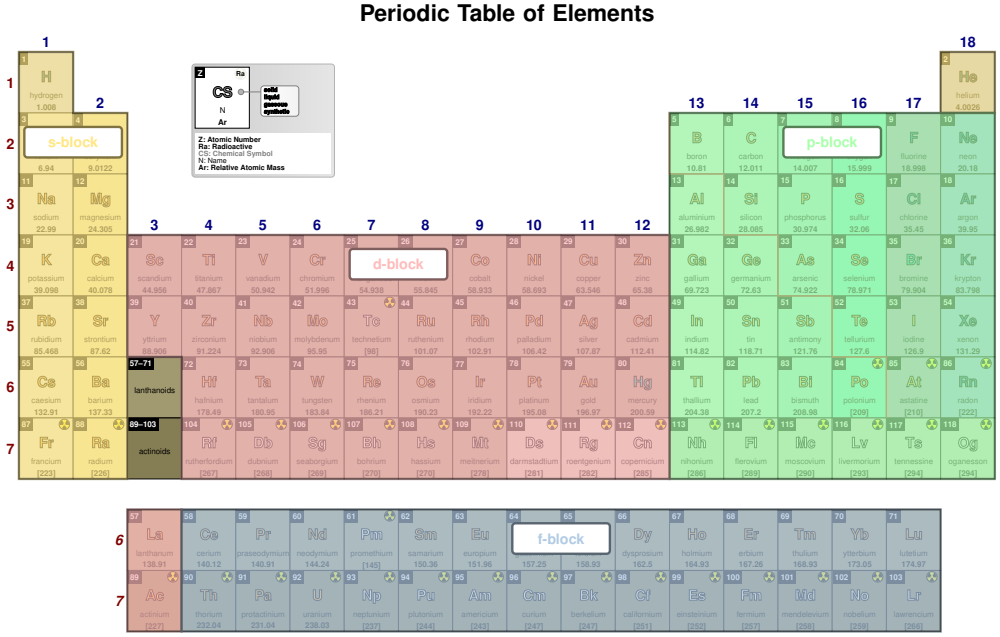
```
\pgfPT[blocks font color,show title=false]
```



blocks line width default: *0.8pt*

Style to set a common width of the lines surrounding the s, p, d and f blocks. The key show blocks is set to true.

```
\pgfPT[blocks line width=1.5pt]
```



blocks default: *{sc=blocos,pc=blocop,dc=blocod,fc=blocof,lw=.8pt,font=\small\bfseries}*

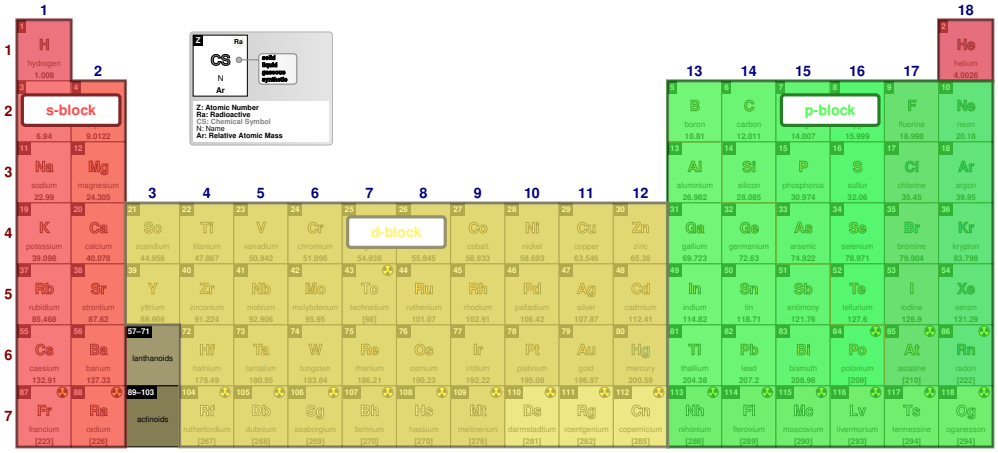
Pseudo style to set the keys: block **s** color, block **p** color, block **d** color, block **f** color, the common **l**ine **w**idths of the blocks, the **s** block **l**ine **w**idth, the **p** block **l**ine **w**idth, the **d** block **l**ine **w**idth, the **f** block **l**ine **w**idth, blocks **f**ont, **s** block **f**ont **c**olor, **p** block **f**ont **c**olor, **d** block **f**ont **c**olor and/or **f** block **f**ont **c**olor. None of the *keys* – sc, pc, dc, fc, lw, slw, plw, dlw, flw, sfc, pfc, dfc and ffc – are mandatory. The key show blocks is set to true.

NOTE:
The colors provided to the color keys of the blocks – sc, pc, dc and fc – could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2` or `color1!value`, as explained in the xcolor package documentation.

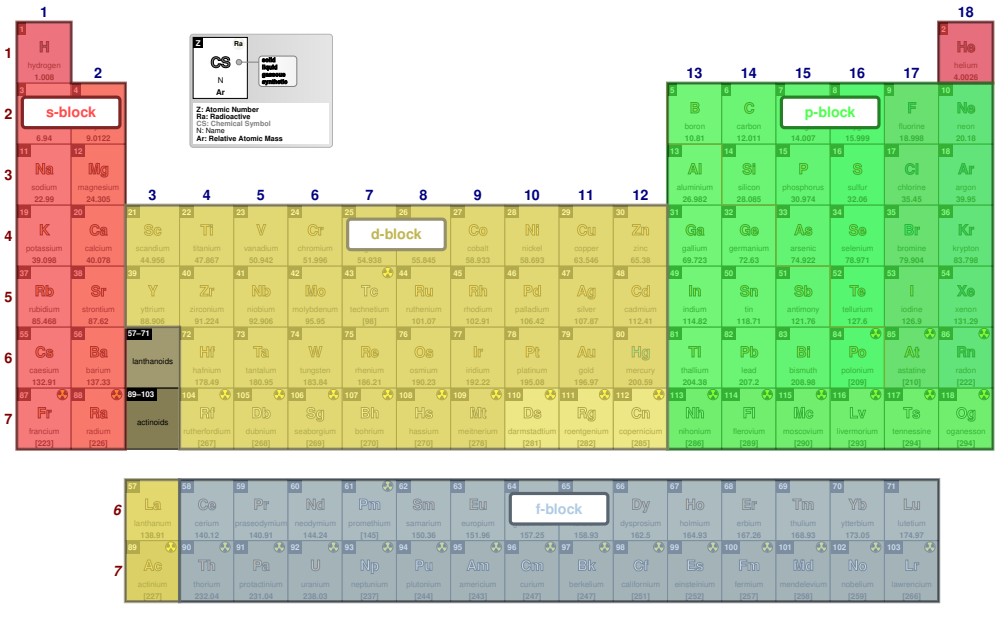
USAGE:

```
blocks={sc=<color>,pc=<color>,dc=<color>,fc=<color>,lw=<lenght>,
        slw=<lenght>,plw=<lenght>,dlw=<lenght>,flw=<lenght>,f=<font commands>,
        sfc=<color>,pfc=<color>,dfc=<color>,ffc=<color>}
```

```
\pgfPT[blocks={sc=red!70!white,pc=green!70!white,dc=yellow!70!white,lw=2pt},
show title=false,Z list=spd]
```



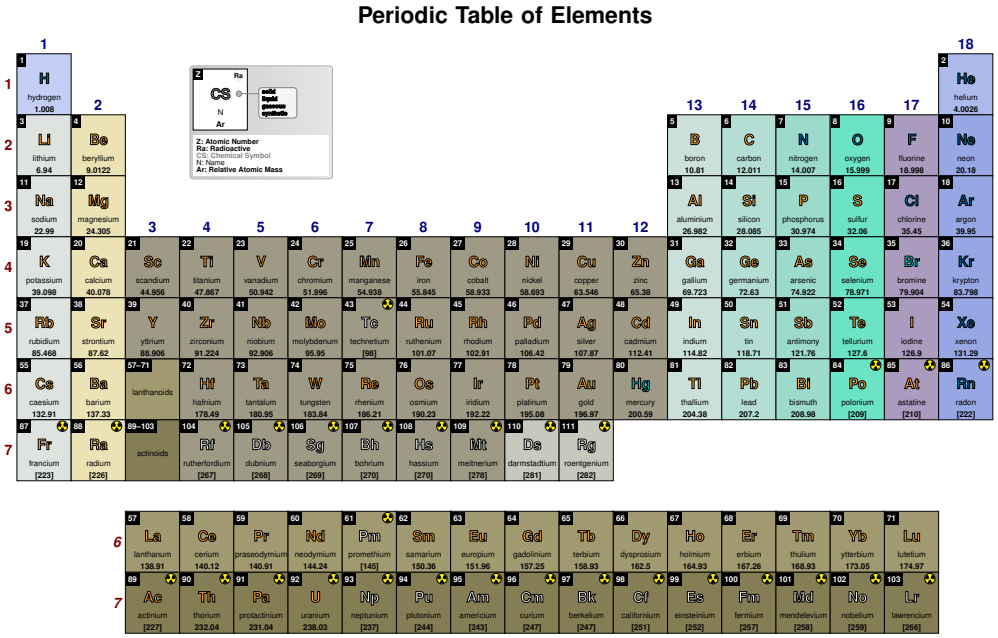

```
\pgfPT[blocks={sc=red!70!white,pc=green!70!white,dc=yellow!70!white,
dfc=yellow!70!black,lw=2pt},show title=false]
```



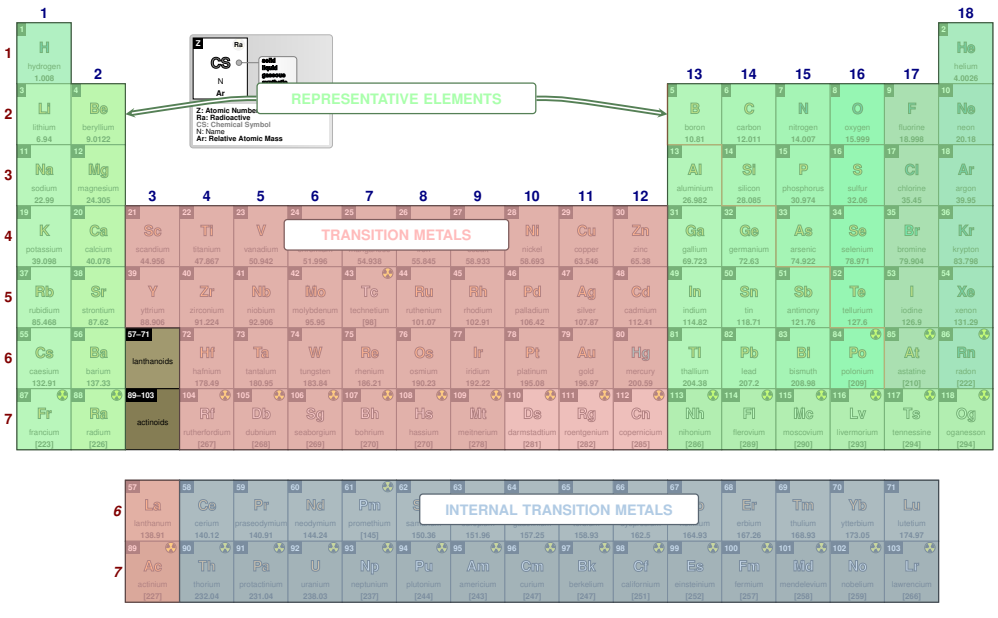
show families default: *false*

When set to true the main families – representative elements, transition metals and internal transition metals – are drawn overlaying the Periodic Table and their labels are shown. *Note that families are only shown when the Z list contains, at least, all elements of blocks s, p and d.*

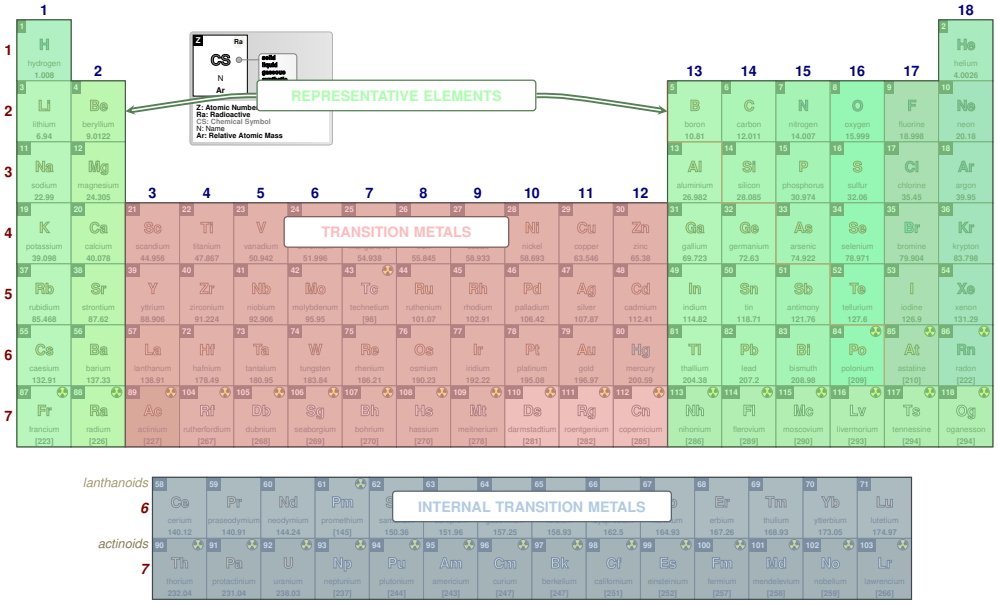
```
\pgfPT[Z list={1,...,111},show families]
```




```
\pgfPT[show families]
```



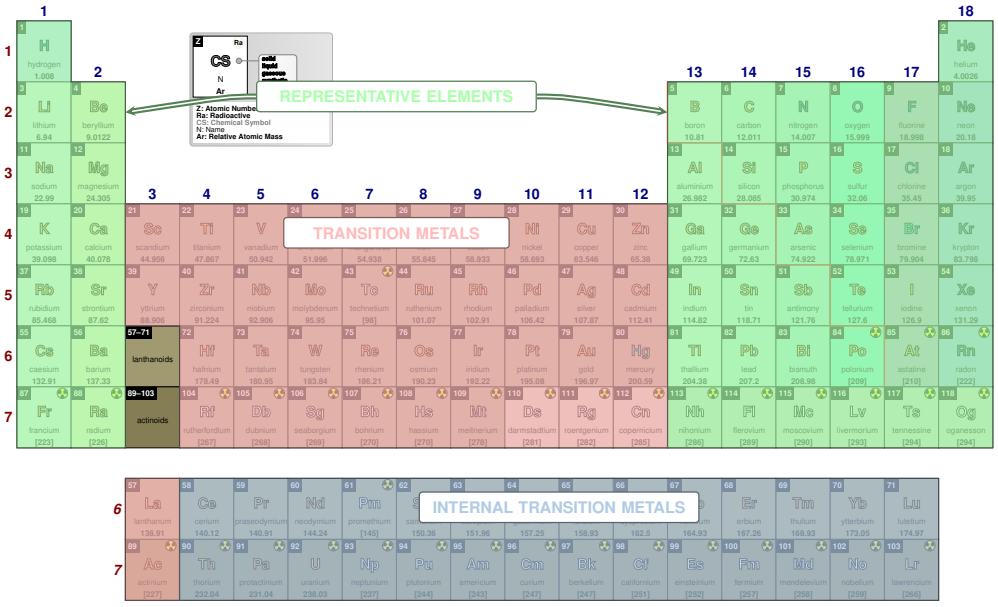
\pgfPT[show families,show title=false,IUPAC=false]



families font default: \small\bfseries

Sets the font used in the family labels.

\pgfPT[show families,show title=false,families font=\normalsize]



r family color default: RGB: 170,255,172

Sets the representative elements *block* color.

r family font color default: {}

Sets the representative elements *block* label font color. If no color is provided, the r family color will be used as the font color.

r family line width default: *0.8pt*

Sets the width of the line surrounding the representative elements *block*.

tm family color default: *RGB: 255,187,187*

Sets the transition metals *block* color.

tm family font color default: *{}*

Sets the transition metals *block* label font color. If no color is provided, the tm family color will be used as the font color.

tm family line width default: *0.8pt*

Sets the width of the line surrounding the transition metals *block*.

itm family color default: *RGB: 177,203,228*

Sets the internal transition metals *block* color.

itm family font color default: *{}*

Sets the internal transition metals *block* label font color. If no color is provided, the itm family color will be used as the font color.

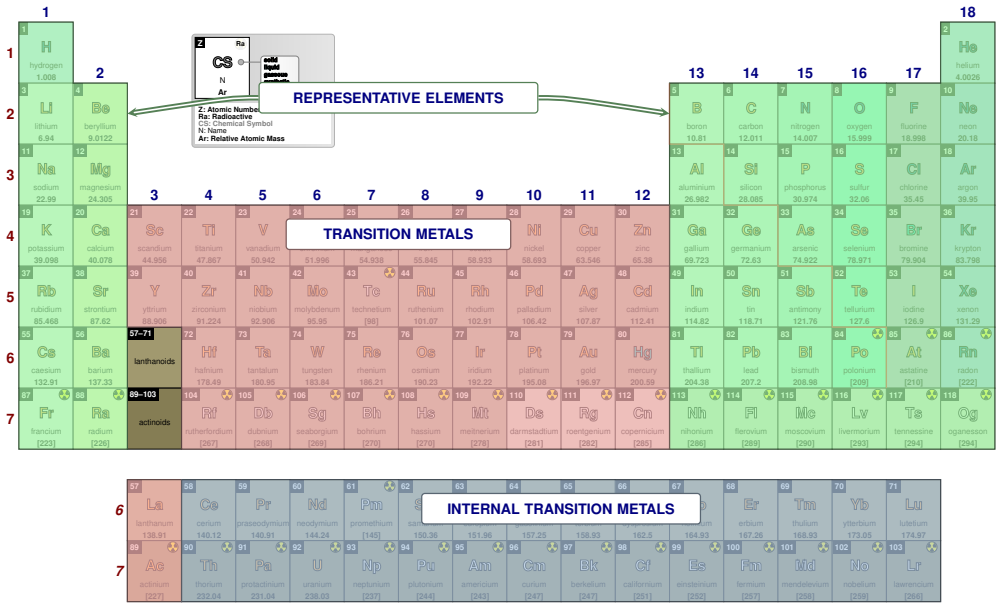
itm family line width default: *0.8pt*

Sets the width of the line surrounding the internal transition metals *block*.

families font color default: *black*

Style to set a common color for the labels of representative elements, transition metals and internal transition metals *blocks*. The key show blocks is set to true.

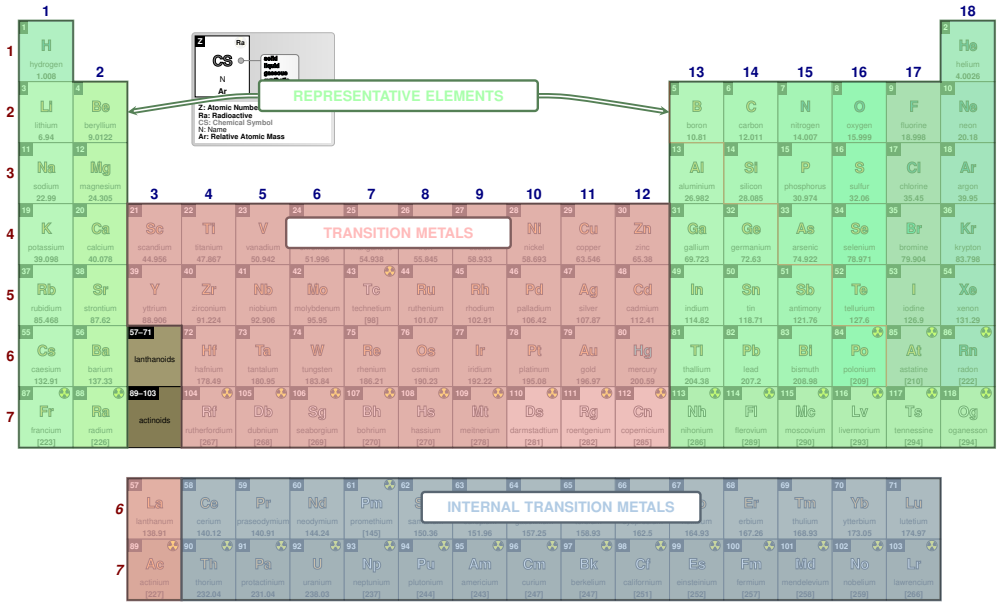
```
\pgfPT[show title=false,families font color=blue!50!black]
```



families line width default: *0.8pt*

Style to set a common width of the lines surrounding the representative elements, transition metals and internal transition metals *blocks*. The key show families is set to true.

```
\pgfPT[show title=false,show families,families line width=1.5pt]
```



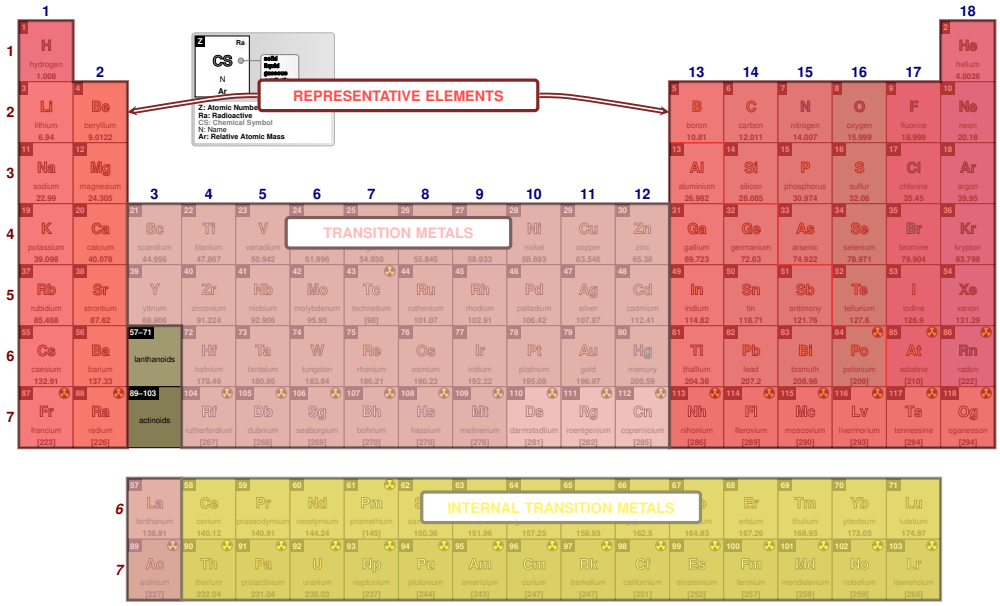
families default: *{rc=blocor,tc=blocot,ic=blocoi,lw=.8pt,f=\small\bfseries}*

Pseudo style to set the keys: **r** family **c**olor, **t**m family **c**olor, **i**tm family **c**olor, the common **l**ine **w**idths of the families, the **r** family **l**ine **w**idth, the **t**m family **l**ine **w**idth, the **i**tm family **l**ine **w**idth, the families **f**ont, **r** family **f**ont **c**olor, **t**m family **f**ont **c**olor and/or **i**tm family **f**ont **c**olor. None of the *keys* – rc, tc, ic, lw, rlw, tlw, ilw, f,rfc, tfc and ifc – are mandatory. The key show families is set to true.

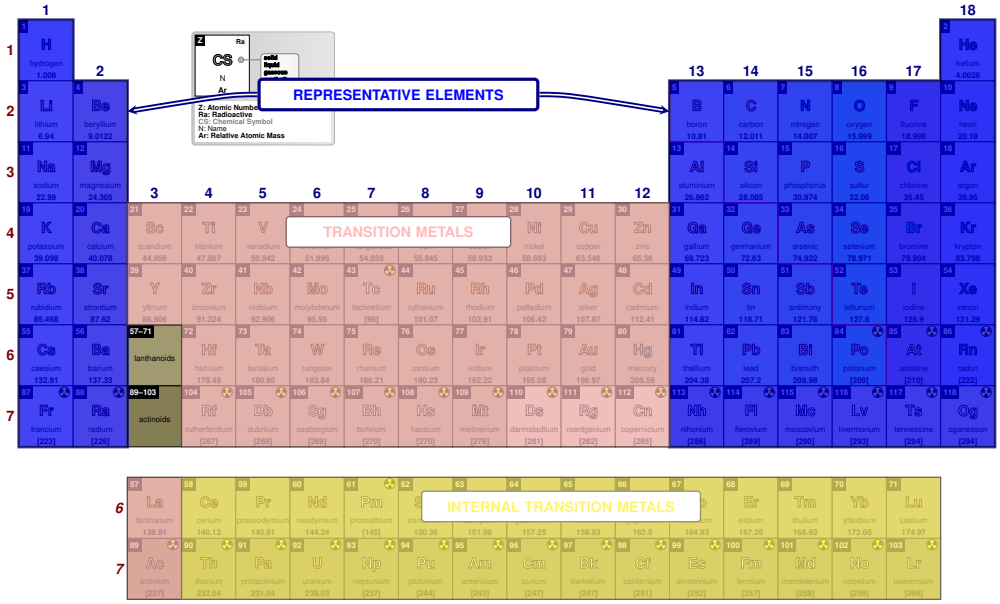
NOTE:
The colors provided to the color keys of the families could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2`, as explained in the xcolor package documentation.

USAGE:
families={rc=<color>,tc=<color>,ic=<color>,lw=<lenght>,rlw=<lenght>,
tlw=<lenght>,ilw=<lenght>,f=<font commands>,rfc=<color>,
tfc=<color>,ifc=<color>}

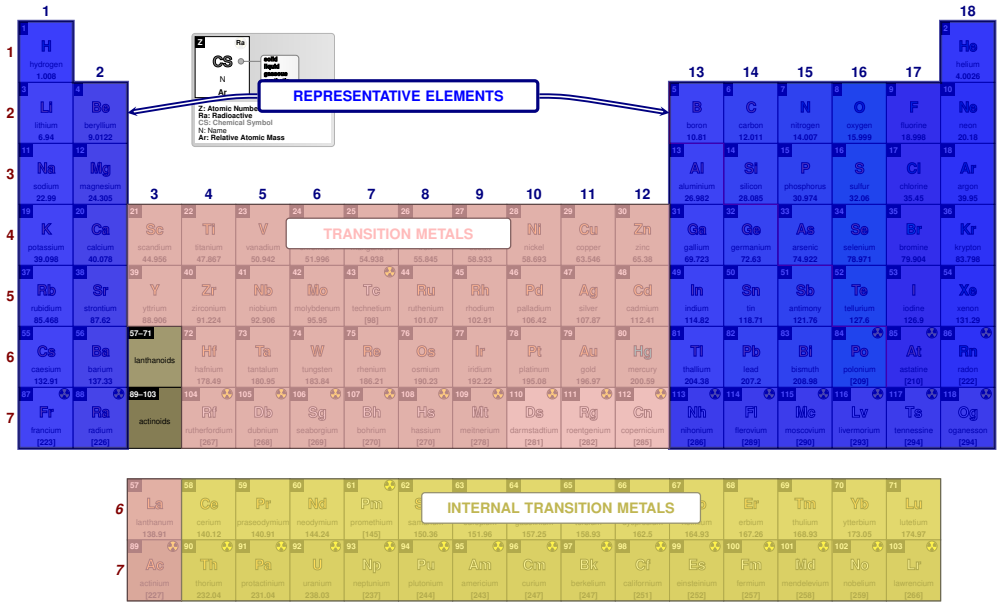
```
\pgfPT[families={rc=red!70!white,ic=yellow!70!white,lw=2pt},show title=false]
```



\pgfPT[families={rc=blue,ic=yellow!70!white,rlw=2pt},show title=false]



\pgfPT[families={rc=blue,ic=yellow!70!white,rlw=2pt,ifc=yellow!70!black},show title=false]



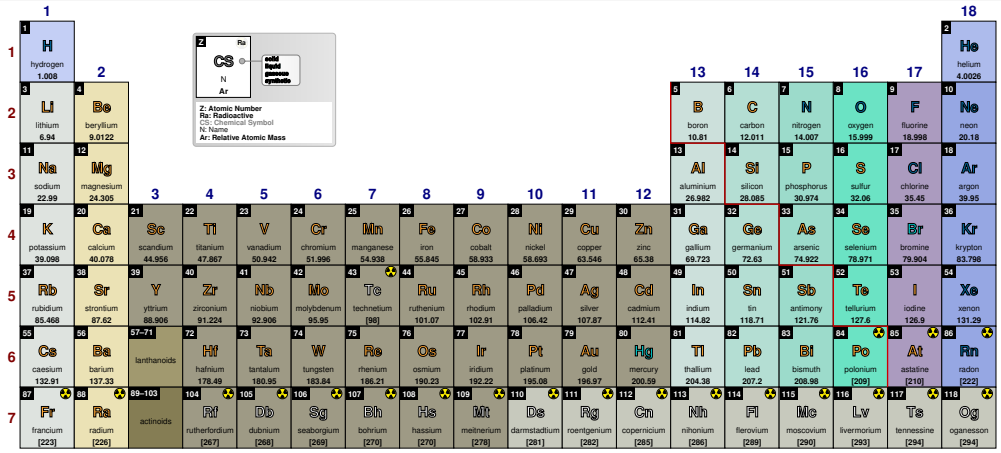
Periodic variations

show periodic variations default: *false*

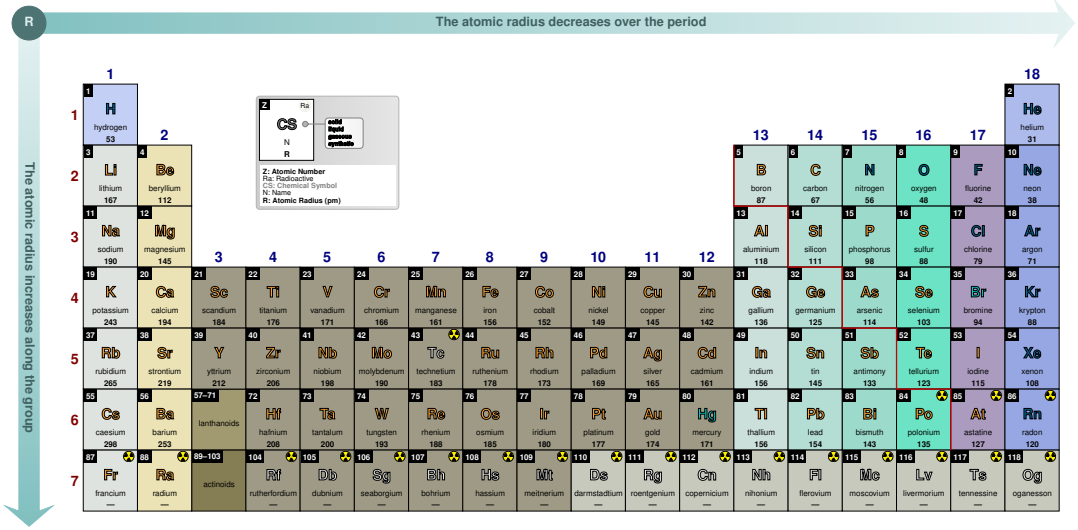
When set to true the periodic variations – for atomic radius, ionization energy and/or electron affinity – are shown with two *arrows*. One horizontal arrow is placed at the top of the Periodic Table for the variation over the period and the other vertically to the left of the Periodic Table for the variation over the group.

NOTE:
The variations are only shown when the *base cell* of the Periodic Table contains the atomic radius, the ionization energy and/or the electron affinity. If none of them is present setting this key (show periodic variations) has no effect.

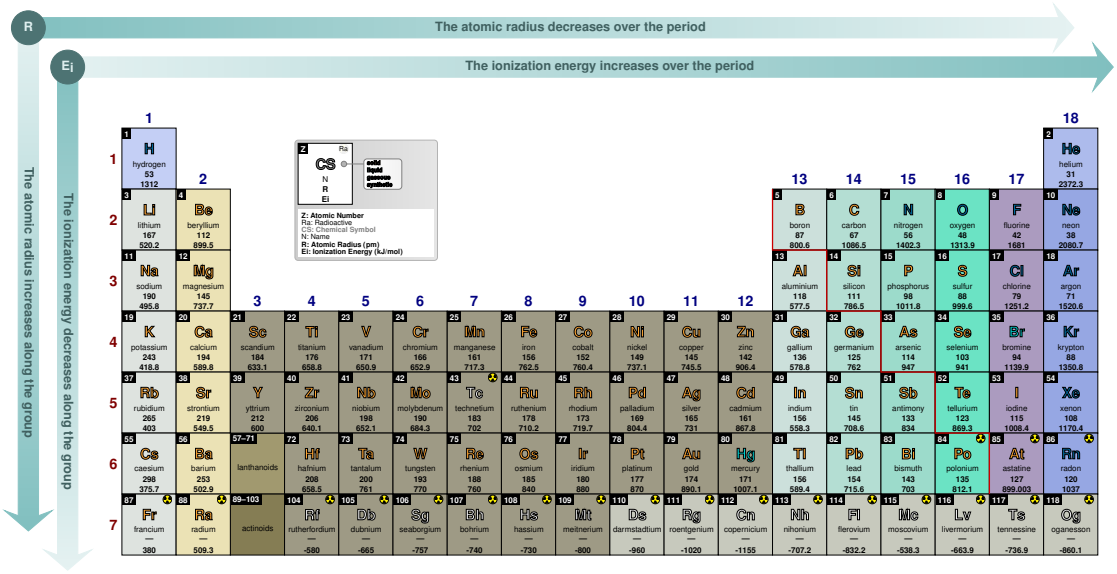
```
\pgfPTstyle[Z list=spd,show title=false]
\pgfPT[show periodic variations]
```




```
\pgfPT[show periodic variations,cell style=pgfPTR]
```



\pgfPT[show periodic variations,cell style=pgfPTREi]

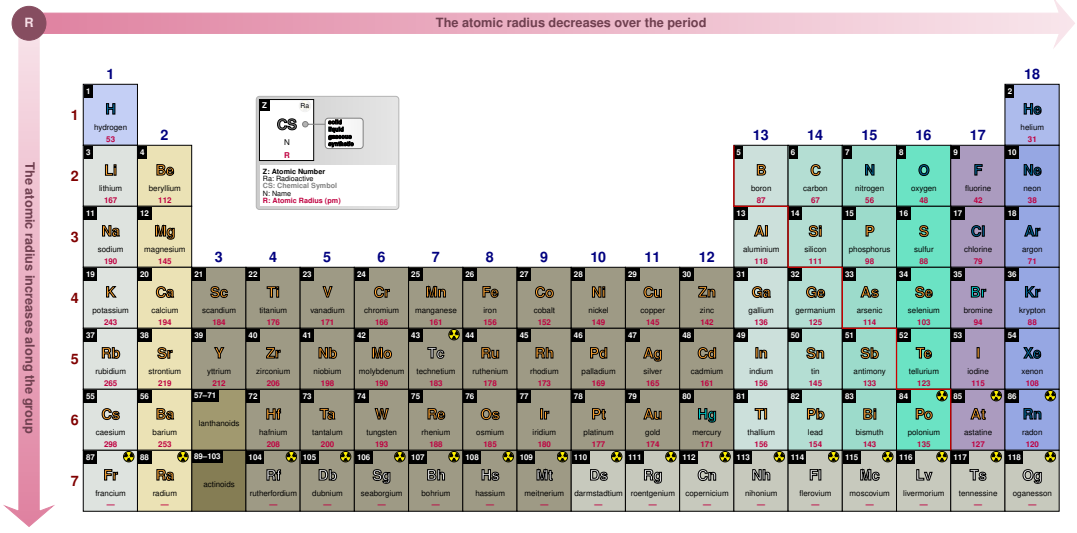


varR color default: RGB: 128,191,191

Sets the color used in the filling of the *arrows* for the atomic radius variations.

NOTE:
The color provided to varR color could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2`, as explained in the xcolor package documentation.

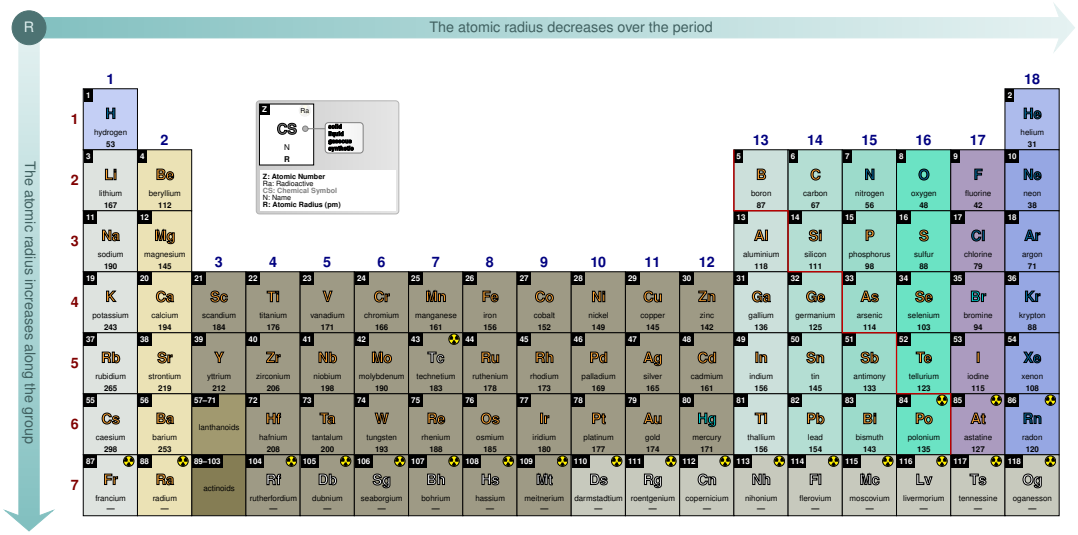
\pgfPT[show periodic variations,cell style=pgfPTR,varR color=teal,R color=purple]



varR font default: *\footnotesize\bfseries*

Sets the font for the text displayed inside the arrow, describing the variation of the atomic radius.

\pgfPT[show periodic variations,cell style=pgfPTR,varR font=\small\itshape]



varR font color default: *(value of varR color)!50!black*

Sets the color of the text showing the atomic radius variations displayed inside the corresponding arrows.
See the note in varR color.

varEi color default: *RGB: 128,191,191*

Sets the color used in the filling of the *arrows* for the ionization energy variations.
See the note in varR color.

varEi font default: *\footnotesize\bfseries*

Sets the font for the text displayed inside the arrow, describing the variation of the ionization energy.

varEi font color default: *(value of varEi color)!50!black*

Sets the color of the text showing the ionization energy variations displayed inside the corresponding arrows.
See the note in varR color.

vareaff color default: *RGB: 128,191,191*

Sets the color used in the filling of the *arrows* for the electron affinity variations.
See the note in varR color.

vareaff font default: *\footnotesize\bfseries*

Sets the font for the text displayed inside the arrow, describing the variation of the electron affinity.

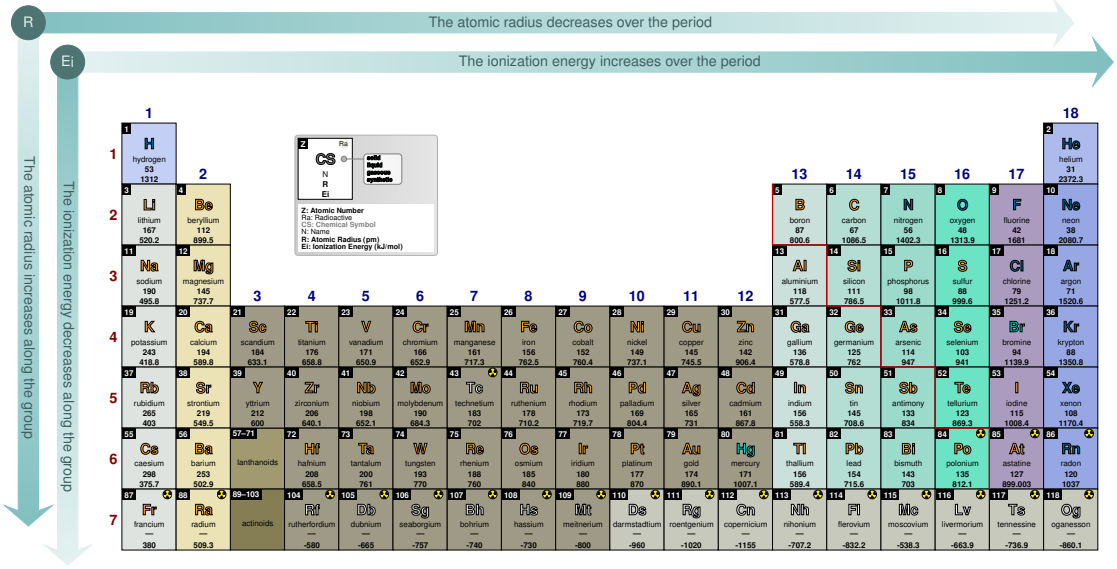
vareaff font color default: *(value of vareaff color)!50!black*

Sets the color of the text showing the electron affinity variations displayed inside the corresponding arrows.
See the note in varR color.

var font default: *\footnotesize\bfseries*

Style to set a common font for the variations along the Periodic Table.
Setting var font=<font commands> is equivalent to setting {varR font=<font commands>, varEi font=<font commands>,vareaff font=<font commands>}.

\pgfPT[show periodic variations,cell style=pgfPTREi,var font=\small\itshape]

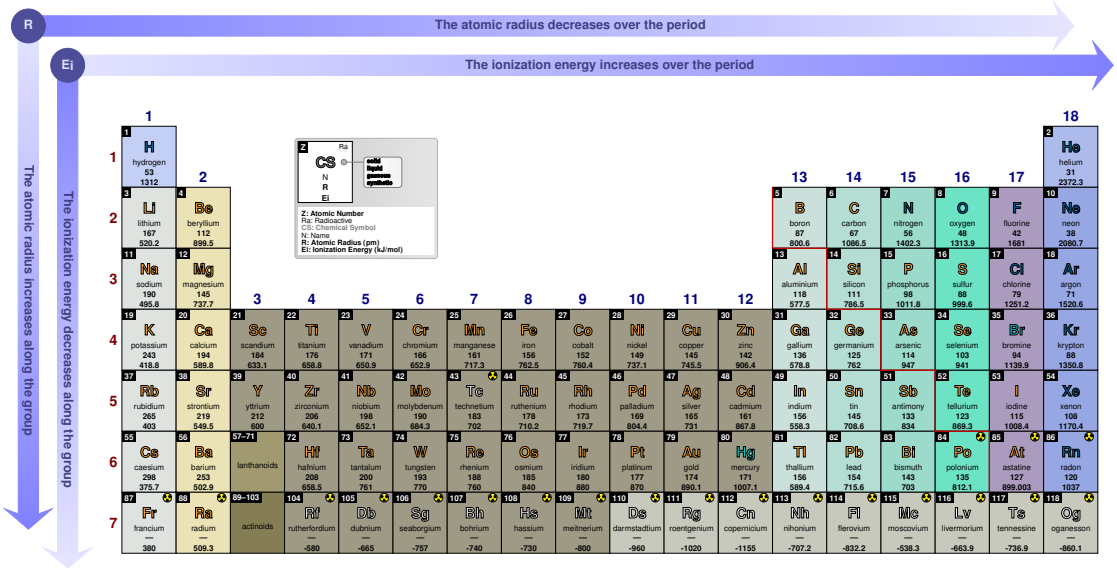


var color default: RGB: 128,191,191

Style to set a common color for the variations along the Periodic Table.
Setting var color=<color> is equivalent to setting {varR color=<color>,varEi color=<color>, vareaff color=<color>}.

NOTE:
The color provided to var color could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2`, as explained in the xcolor package documentation.
Keep in mind that setting the variations colors also changes the default text colors for them.

\pgfPT[show periodic variations,cell style=pgfPTREi,var color=blue!50!white]



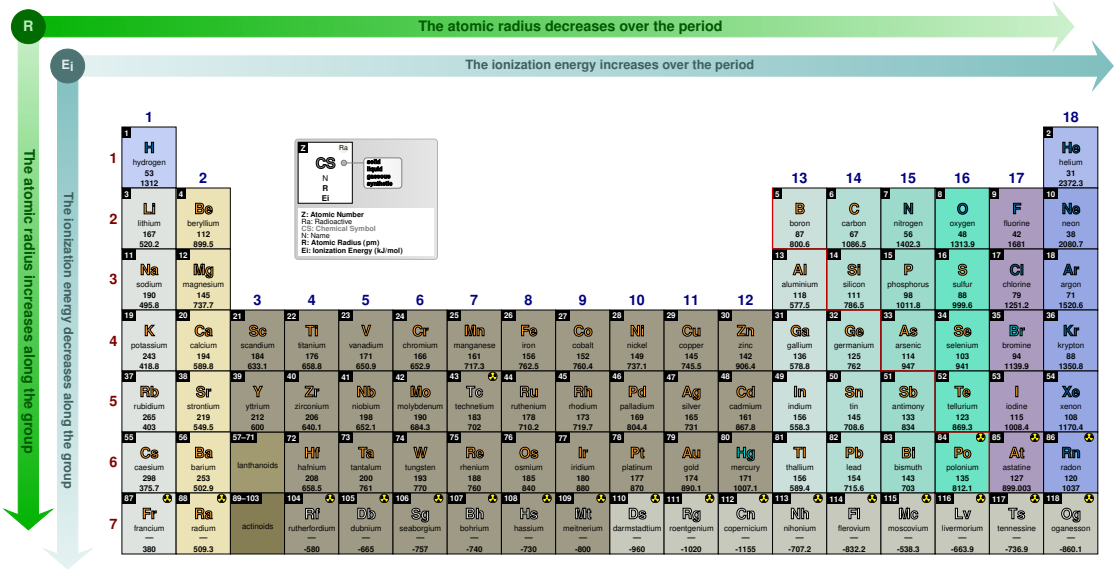
varR default: *{c=colorvariations,f=\footnotesize\bfseries}*

Pseudo style to set the keys: varR **c**olor, varR **f**ont and/or varR **f**ont **c**olor. None of the *keys* – c, f and fc – are mandatory.

NOTE:
The color provided to varR color could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2`, as explained in the xcolor package documentation.

USAGE: varR={c=<color>,f=<font commands>,fc=<color>}

```
\pgfPT[show periodic variations,cell style=pgfPTREi,
varR={c=green!70!black,f=\small\bfseries}]
```



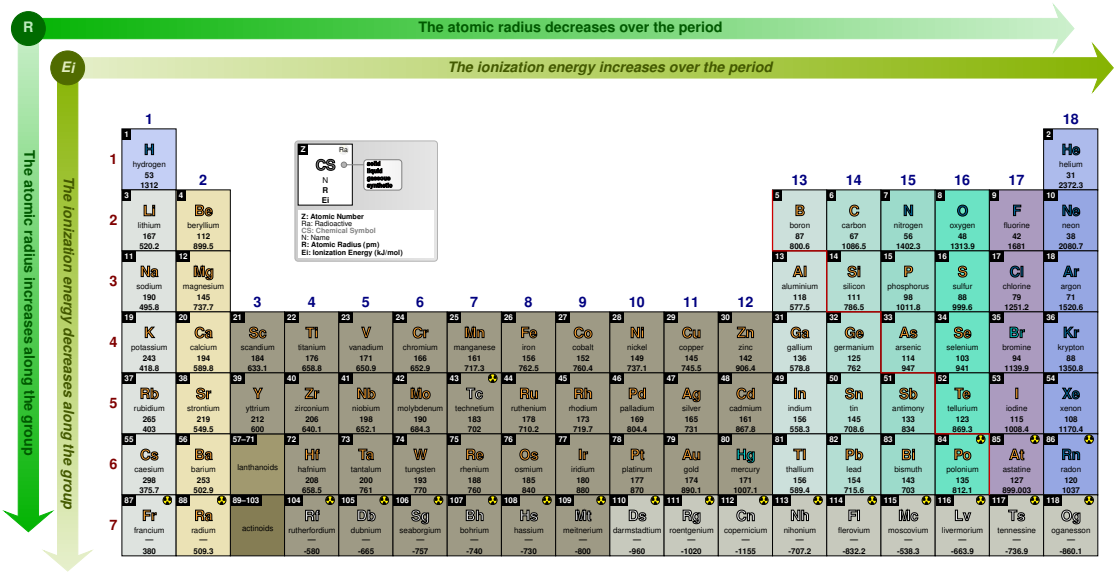
varEi default: *{c=colorvariations,f=\footnotesize\bfseries}*

Pseudo style to set the keys: varEi **c**olor, varEi **f**ont and/or varEi **f**ont **c**olor. None of the *keys* – c, f and fc – are mandatory.

NOTE:
The color provided to varEi color could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2`, as explained in the xcolor package documentation.

USAGE: varEi={c=<color>,f=<font commands>,fc=<color>}

```
\pgfPT[show periodic variations,cell style=pgfPTREi,
varR={c=green!70!black,f=\small\bfseries},
varEi={c=lime!70!black,f=\small\bfseries}]
```



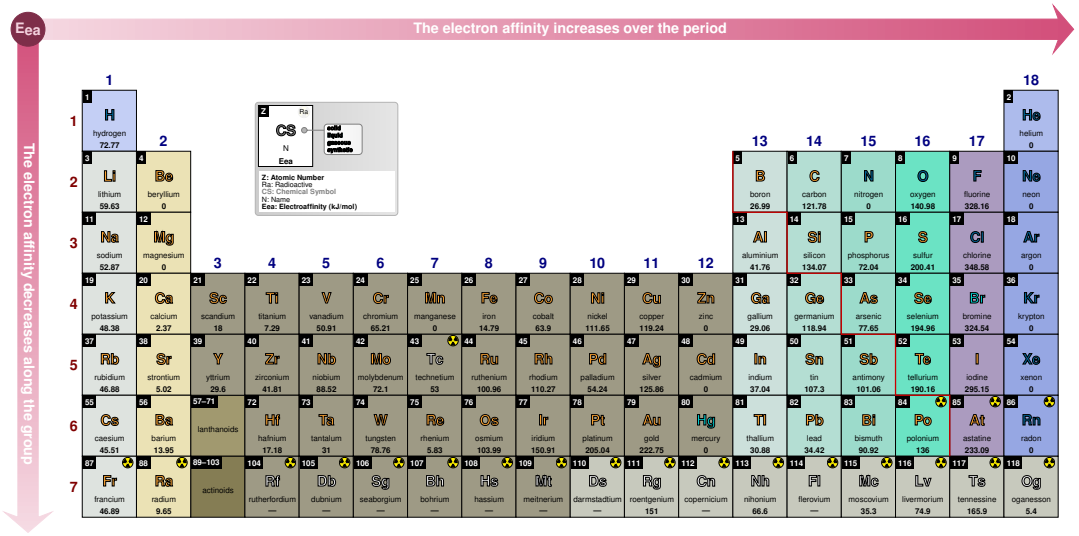
vareaff default: *{c=colorvariations,f=\footnotesize\bfseries}*

Pseudo style to set the keys: vareaff **c**olor, vareaff **f**ont and/or vareaff **f**ont **c**olor. None of the *keys* – c, f and fc – are mandatory.

NOTE:
The color provided to vareaff color could be any defined color via the command `\definecolor` or by *mixing* colors, using, for instance, the syntax `color1!value!color2`, as explained in the xcolor package documentation.

USAGE: vareaff={c=<color>,f=<font commands>,fc=<color>}

```
\pgfPT[show periodic variations,cell style=pgfPTeaff,
vareaff={c=purple!70!white,f=\small\bfseries,fc=white}]
```




```
\pgfPTresetstyle
```

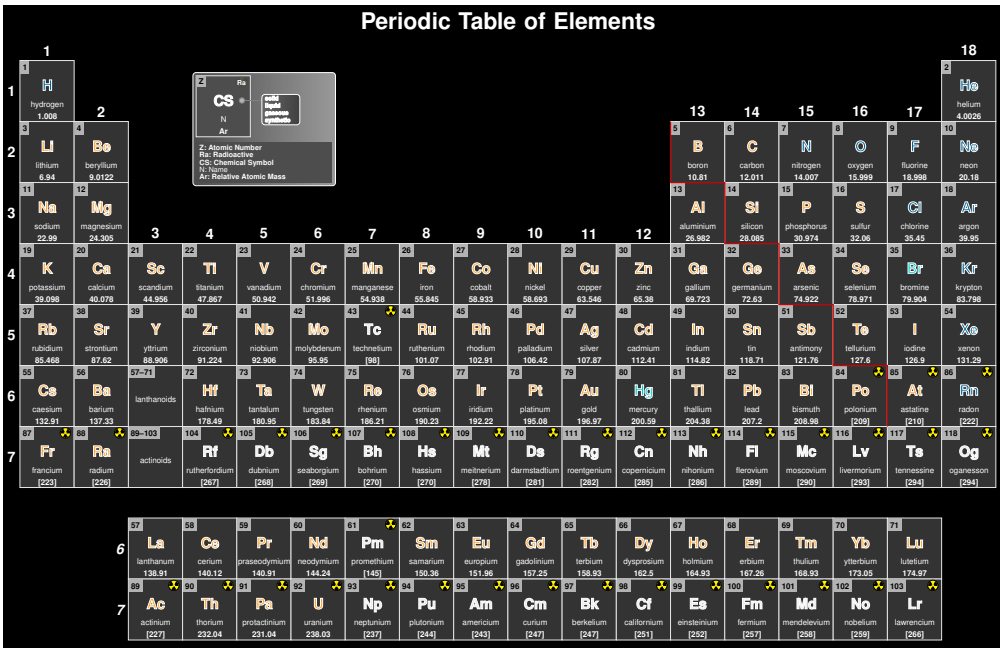
Dark mode

dark mode default: *no value*

Style to change the overall appearance of the Periodic Table to a dark mode suitable for on-screen viewing.

***This style sets the following keys with the values**:*
back color scheme=solid, back color=black!80, cell line color=black!10, CS outline color=white, cell color=white, Z backcolor=black!30, Z color=black, background={fill=black}, varR font color=black!20, varEi font color=black!20, vareaff font color=black!20, per+gr={c=white}, title color=white, other languages color=black!40, legend={bc=black!70,radio=white,CS=white,Z=white,pins style={draw=white,right color=black!75, left color=black!60,line width=.05pt,rounded corners=2pt},extra style={draw=white,fill=black!70,line width=.05pt, rounded corners=2pt},box={left color=black!70,right color=black!40,draw=white}}

```
\pgfPT[dark mode]
```



Exercise layout

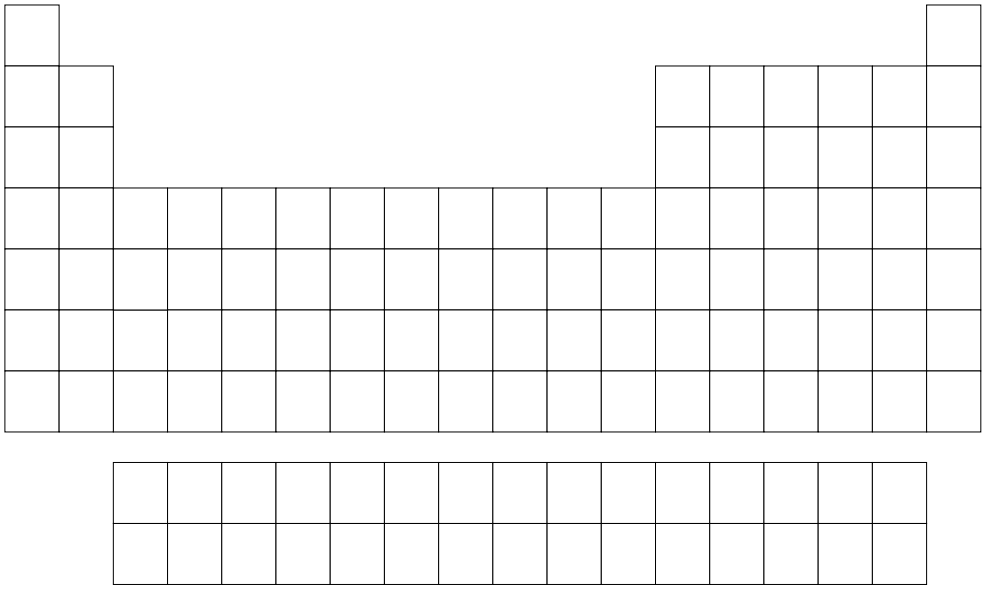
The keys described in this section enable the *exercise layout* of the Periodic Table, *i.e.*, in this mode the *structure* of the Periodic Table is drawn, but there are only a few contents available in the cells.

only cells default: *false*

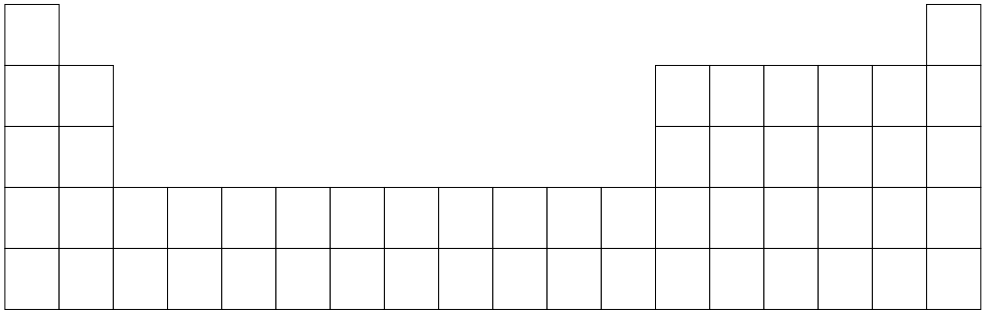
When set to true the Periodic Table is drawn with only the cells without any contents.

NOTE:
The following keys are also set: back color scheme=solid, show title=false, show period numbers=false, show group numbers=false, show legend=false, show MNM line=false

\pgfPT[only cells]



\pgfPT[Z list={1,..,54},only cells]

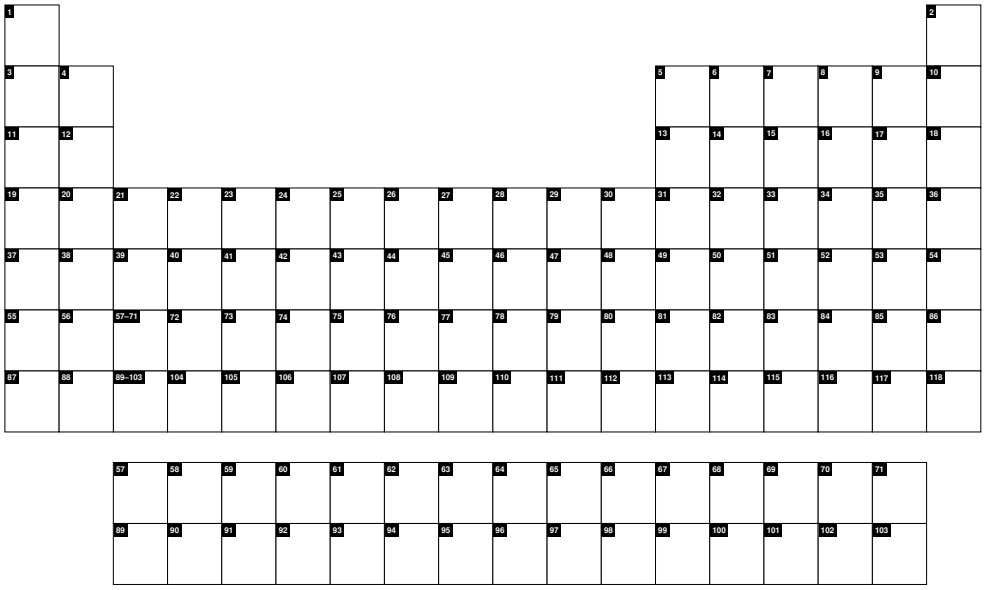


only cells plus Z default: *false*

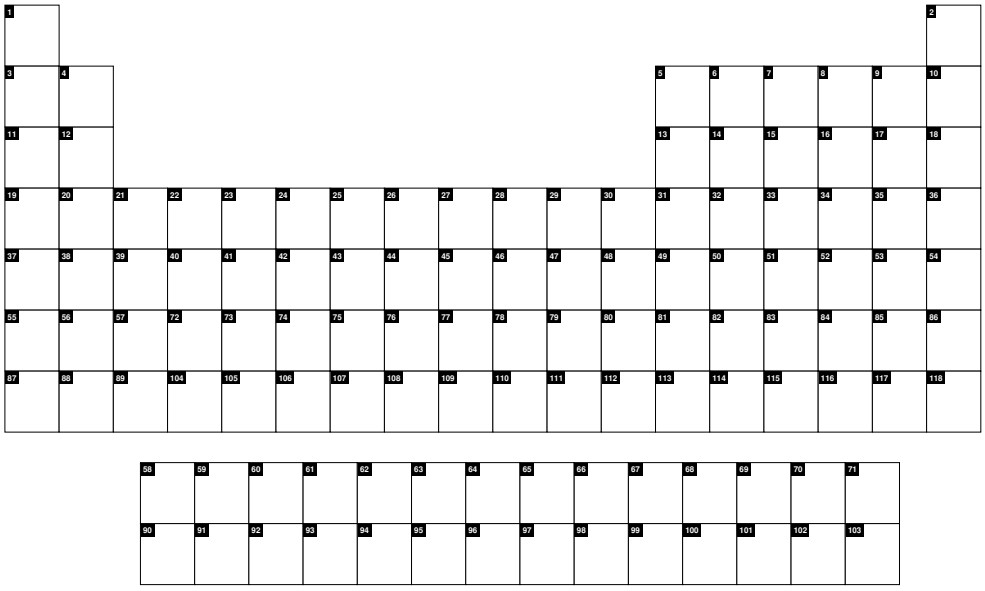
When set to true the Periodic Table is drawn with only the cells without any contents, except the atomic number (Z).

NOTE:
The following keys are also set: back color scheme=solid, show title=false, show period numbers=false, show group numbers=false, show legend=false, show MNM line=false

\pgfPT[only cells plus Z]



\pgfPT[only cells plus Z,IUPAC=false]

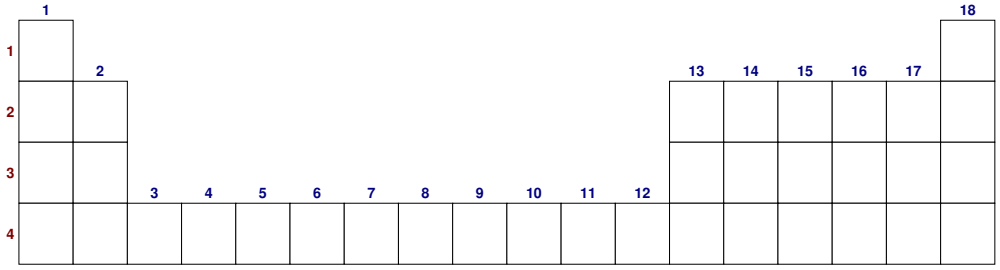


only cells with periods and group numbers default: *false*

When set to true the Periodic Table is drawn with only the cells without any contents. The period and group numbers are shown.

NOTE:
The following keys are also set: back color scheme=solid, show title=false, show legend=false, show MNM line=false

\pgfPT[Z list={1,...,36},only cells with periods and group numbers]

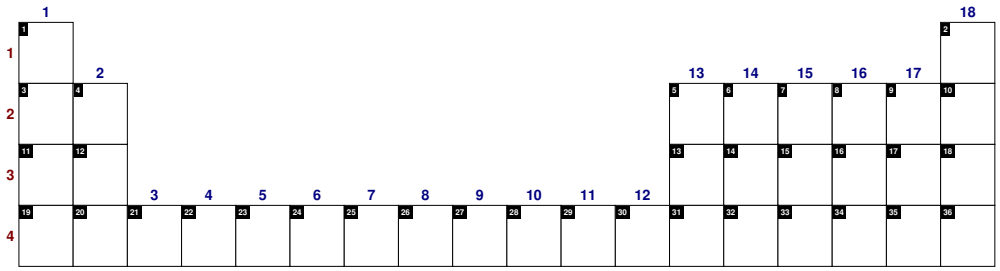


only cells with periods and group numbers plus Z default: *false*

When set to true the Periodic Table is drawn with only the cells without any contents, except the atomic number (Z). The period and group numbers are shown.

NOTE:
The following keys are also set: back color scheme=solid, show title=false, show legend=false, show MNM line=false

\pgfPT[Z list={1,...,36},only cells with periods and group numbers plus Z]



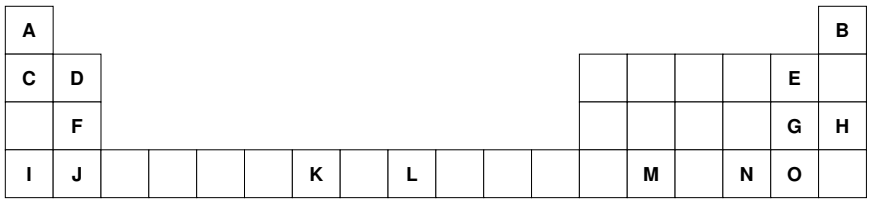
Z exercise list default: *{}*

Sets the list of atomic numbers to display as letters instead of their chemicals symbols.

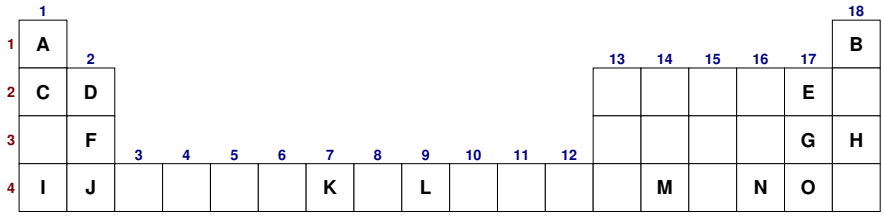
NOTES:

- ✓ When values are provided to the Z exercise list and none of the above *exercise layout* is set, the *exercise layout* only cells is used.
- ✓ The line dots – ... – notation is not available in the *Z exercise list*, mainly to avoid *errors* on the desired list. For example {1,...,4,8,...,16} is expanded by the `\foreach` statement of TikZ to {1,2,3,4,8,15} instead of {1,2,3,4,8,9,10,11,12,13,14,15,16}. For achieving that purpose it must be typed {1,...,4,8,9,...,16}. Since the goal of *Z exercise list* is typing only a list of specific elements, it will often be easier to type element by element.

\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87}, cell size=3em,Z list={1,...,36}]



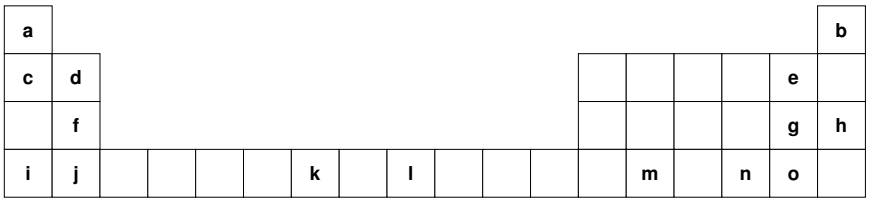
```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36},only cells with periods and group numbers]
```



exercise list in capitals default: *true*

When set to true the *letters* are typed in capitals, otherwise they are typed as lowercase letters.

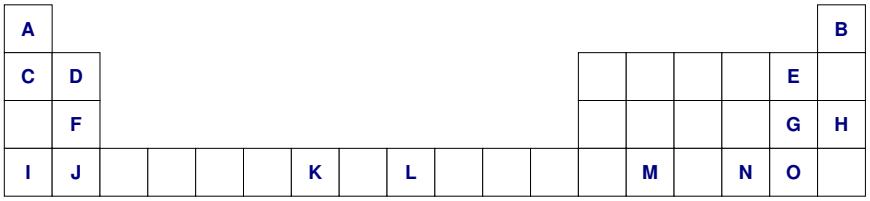
```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36},exercise list in capitals=false]
```



exercise list color default: *black*

Sets the color of the displayed *letters* in the *exercise layout*.

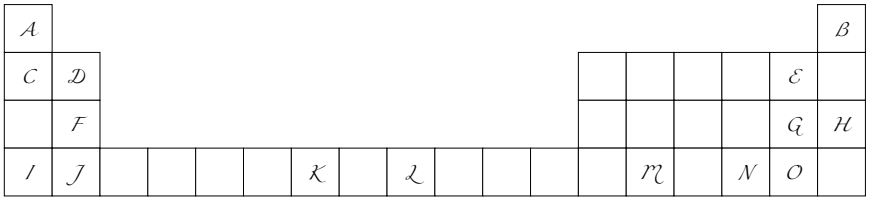
```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36}, exercise list color=blue!50!black]
```



exercise list font default: *\bfseries\large*

Sets the font of the displayed *letters* in the *exercise layout*.

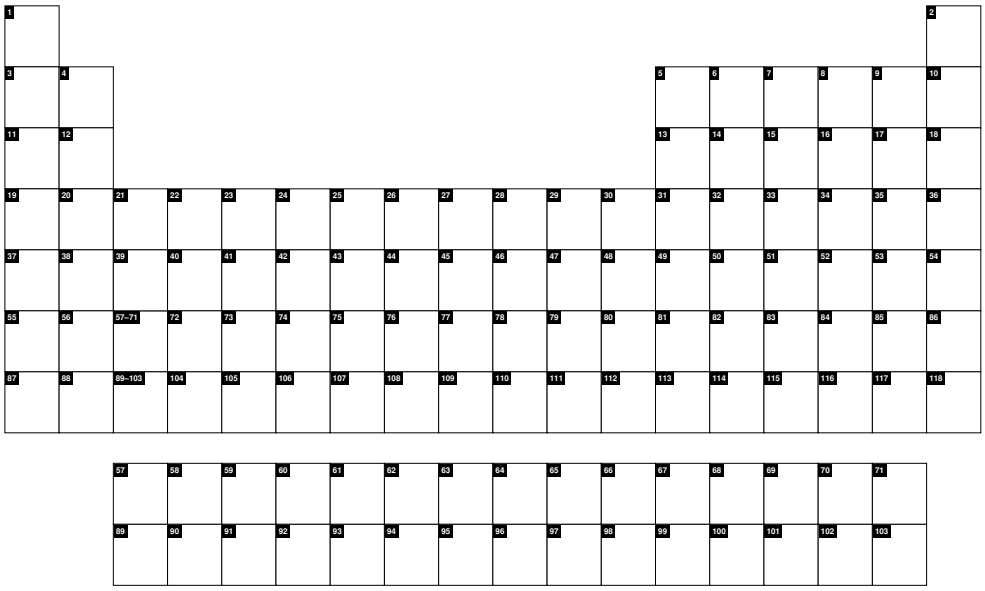
```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36}, exercise list font=\fontfamily{fmm}\selectfont]
```



cells+Z *no value*

Style to set the key only cells plus Z to true.

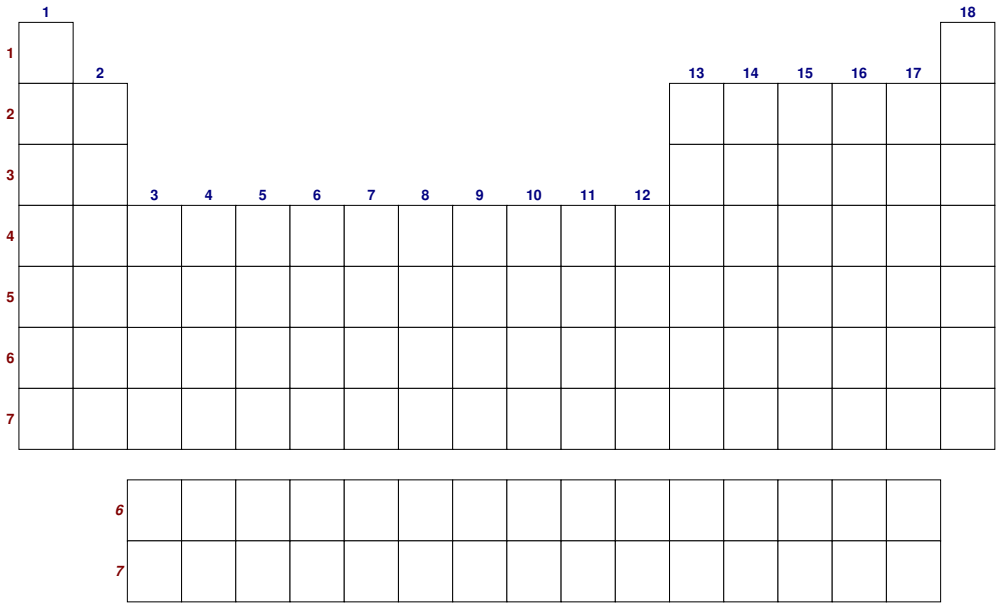
```
\pgfPT[cells+Z]
```



cells+p+g *no value*

Style to set the key only cells with periods and group numbers to true.

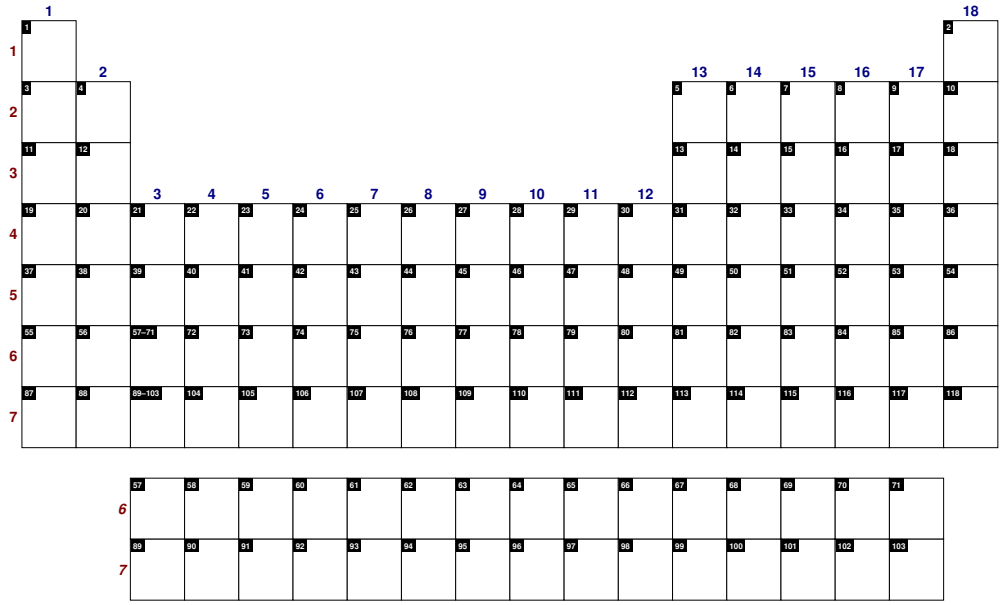
```
\pgfPT[cells+p+g]
```



cells+p+g+Z *no value*

Style to set the key only cells with periods and group numbers plus Z to true.

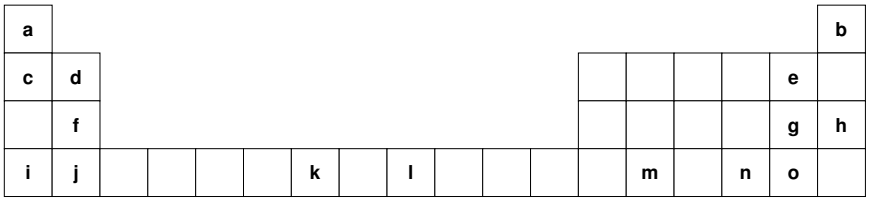
```
\pgfPT[cells+p+g+Z]
```



exnocaps *no value*

Style to set the key exercise list in capitals to false.

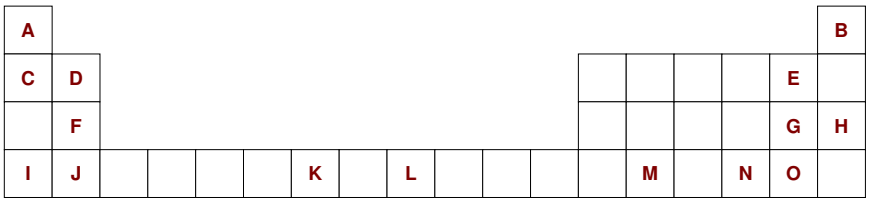
```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36},exnocaps]
```



exColor default: *black*

Style to set the key exercise list color.

```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36},exColor=red!50!black]
```



exFont default: *\bfseries\large*

Style to set the key exercise list font.

```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36},exFont=\Large]
```

A																	B
C	D															E	
	F															G	H
I	J					K		L					M		N	O	

ex default: *{caps=true,c=black,f=\bfseries\large}*

Pseudo style to set the keys: exercise list in **cap**ital**s**, exercise list **c**olor and/or exercise list **f**ont. None of the *keys* – caps, c and f – are mandatory.

USAGE: ex={caps=<true|false>,c=<color>,f=<font commands>}

```
\pgfPT[Z exercise list={1,2,3,4,9,12,17,18,19,20,25,27,32,34,35,49,54,74,86,87},
cell size=3em,Z list={1,...,36},ex={c=blue,f=\Large\bfseries}]
```

A																	B
C	D															E	
	F															G	H
I	J					K		L					M		N	O	

✠ Cell contents options: keys, styles and *pseudo styles*

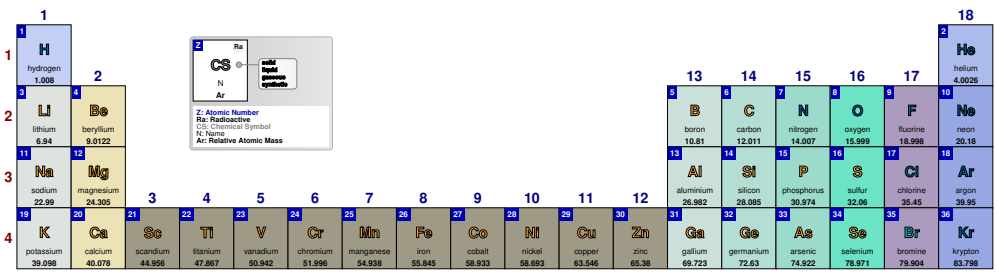
The following options and styles are used for customizing the contents available in each individual cell of the Periodic Table, like the *fonts* or the *colors* used in the shown contents.

➠ The atomic number

Z backcolor default: *black*

Sets the background color of the box where the atomic number is displayed.

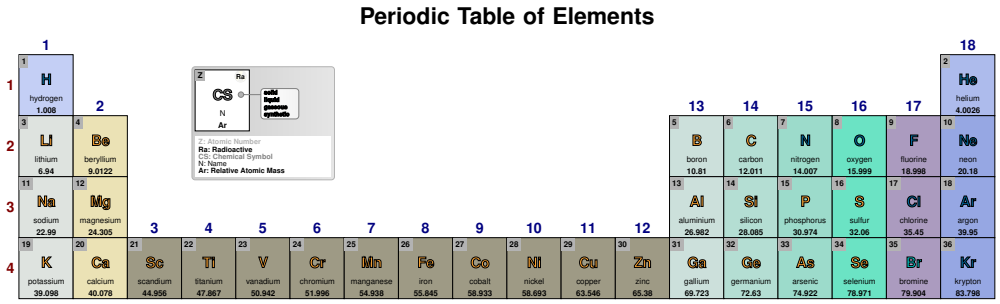
```
\pgfPT[Z list={1,...,36},Z backcolor=blue!70!black]
```



Z color default: *white*

Sets the color of the atomic number.

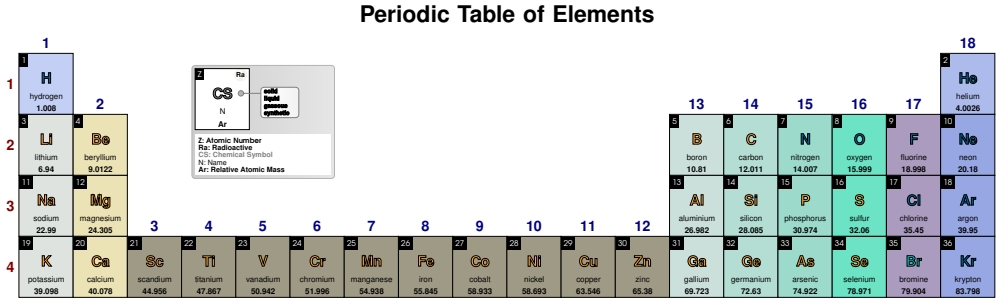
```
\pgfPT[Z list={1,...,36},Z backcolor=black!30,Z color=black]
```



Z font default: *\tiny\bfseries*

Sets the font of the atomic number.

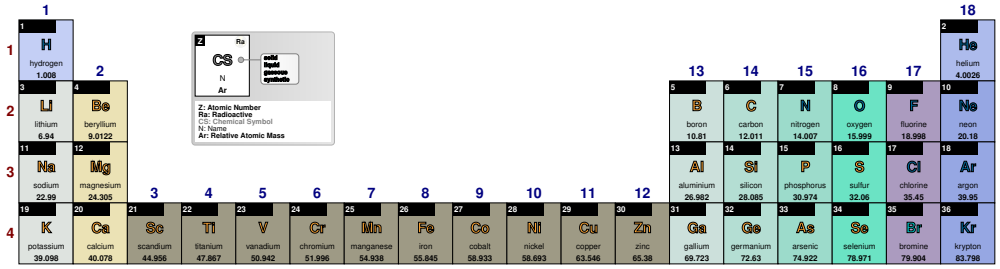
```
\pgfPT[Z list={1,...,36},Z font=\fontfamily{pag}\selectfont\tiny]
```



Z use box width default: *false*

If true, the width specified in the constructed cell is used, otherwise, the *natural* width of the box containing Z value is used.

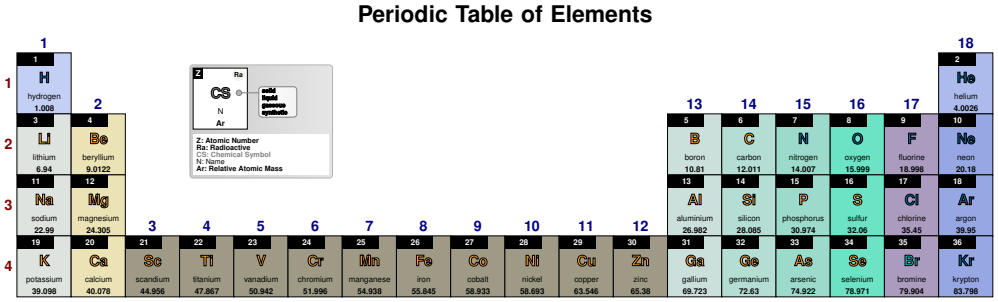
```
\pgfPT[Z list={1,...,36},Z use box width]
```



Z align default: *left*

Sets the alignment of the atomic number value to *left*, *center* or *right* with respect to its containing box. It only takes effect when Z use box width is true.

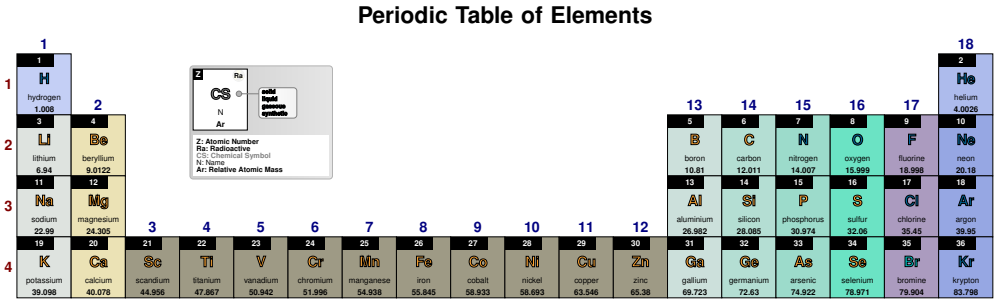
\pgfPT[Z list={1,...,36},Z use box width,Z align=center]



Z padding default: *0.25ex*

Sets the padding between the atomic number value and the box that contains it. It only takes effect when Z use box width is true.

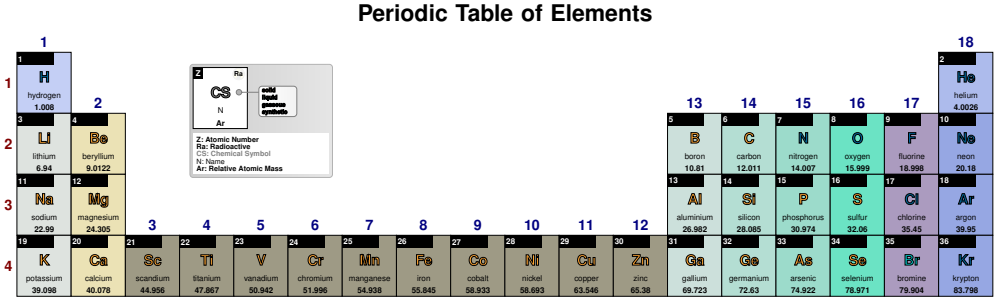
\pgfPT[Z list={1,...,36},Z use box width,Z align=right, Z padding=1em]



Z box *no value*

Style equivalent to Z use box width=true.

\pgfPT[Z list={1,...,36},Z box]

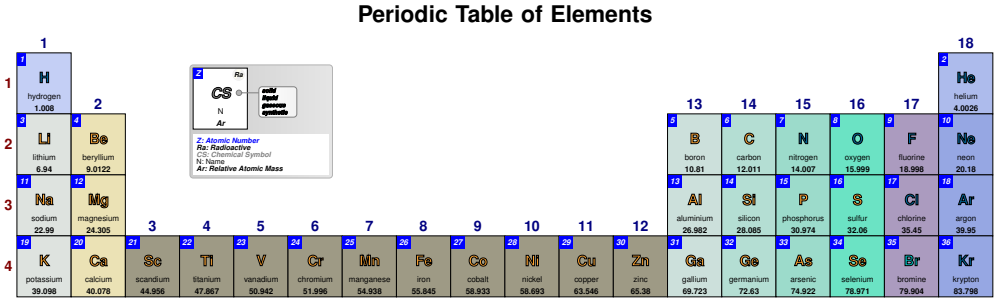


Z default: *{bc=black,c=white,f=\tiny\bfseries,boxwd=false,align=left,pad=.25ex}*

Pseudo style to set the keys: Z **b**ack**c**olor, Z **c**olor, Z **f**ont, Z use **box wi**dth, Z **align** and/or Z **pad**ding. None of the *keys* – bc, c, f, boxwd, align and pad – are mandatory.

USAGE: Z={bc=<color>,c=<color>,f=<font commands>,boxwd=<true|false>, align=<left|center|right>,pad=<length>}

```
\pgfPT[Z list={1,...,36},Z={bc=blue,f=\tiny\bfseries\itshape}]
```

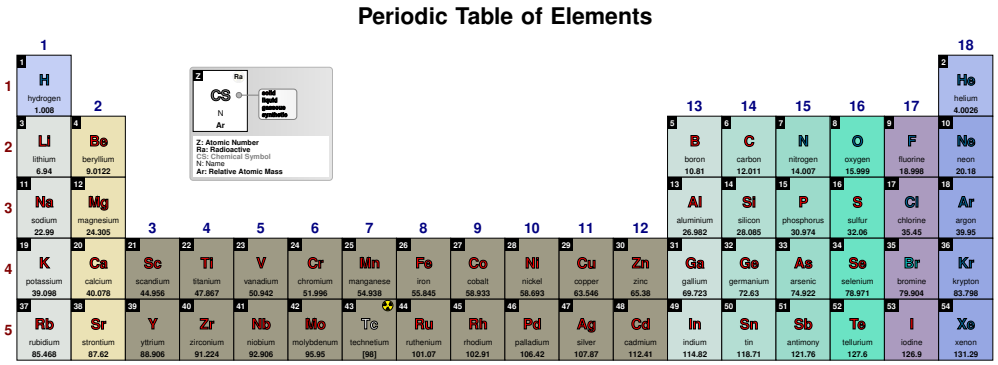


➠ The chemical symbol

CS solid default: RGB: 255,166,51

Sets the color of the chemical symbol for elements that are in the solid state at normal temperature and pressure (NTP).

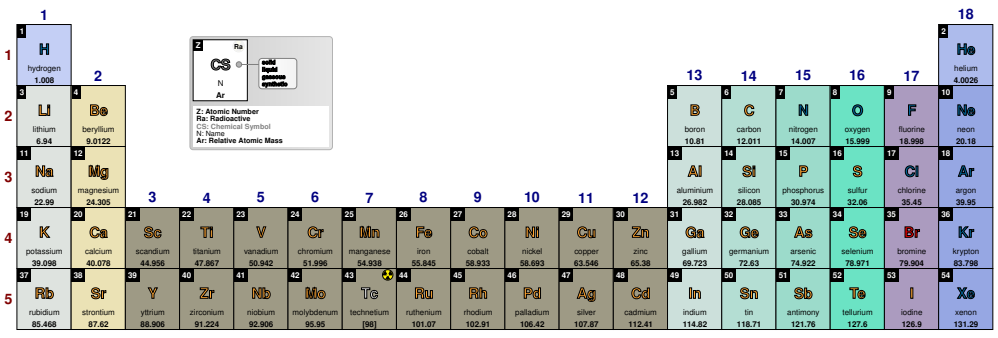
```
\pgfPT[Z list={1,...,54},CS solid=red]
```



CS liquid default: RGB: 0,204,204

Sets the color of the chemical symbol for elements that are in a liquid state at normal temperature and pressure (NTP).

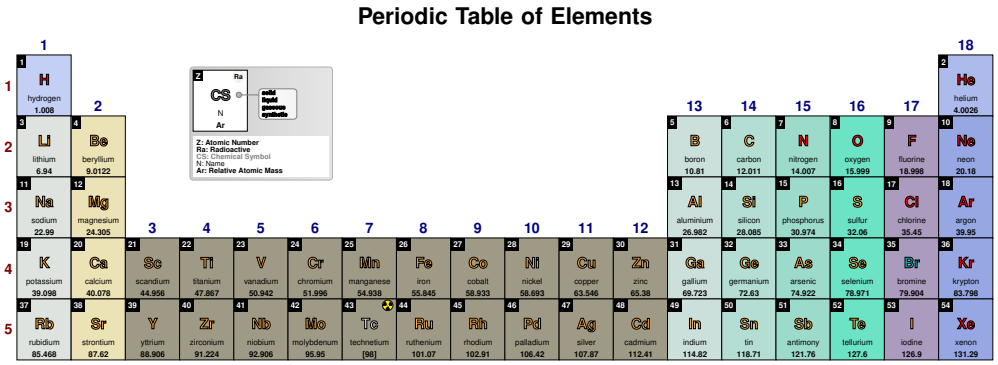
```
\pgfPT[Z list={1,...,54},CS liquid=red]
```



CS gas default: *RGB: 0,102,153*

Sets the color of the chemical symbol for elements that are in a gaseous state at normal temperature and pressure (NTP).

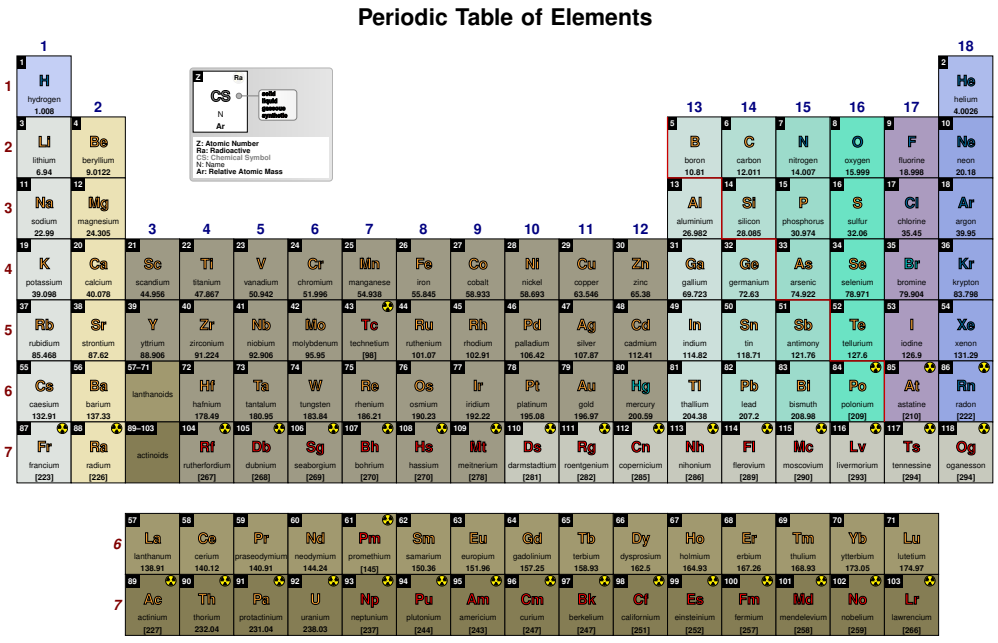
```
\pgfPT[Z list={1,...,54},CS gas=red]
```



CS synt default: *RGB: 236,236,236*

Sets the color of the chemical symbol for elements that are synthetic.

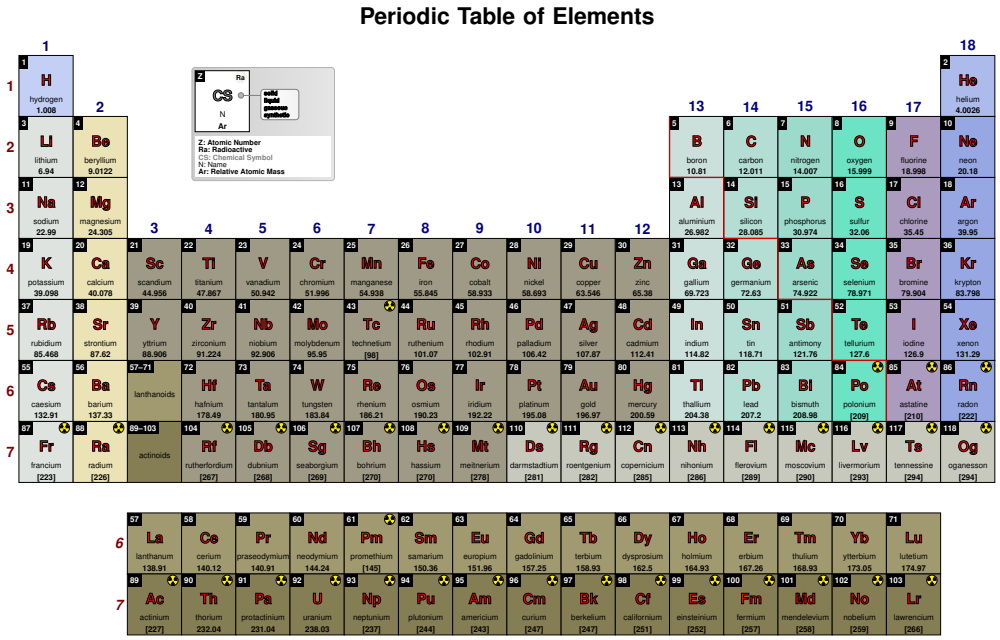
```
\pgfPT[CS synt=red]
```



CS all default: *black*

Style to set a common color to the chemical symbols, equivalent to CS solid=<color>, CS liquid=<color>,CS gas=<color>,CS synt=<color>.

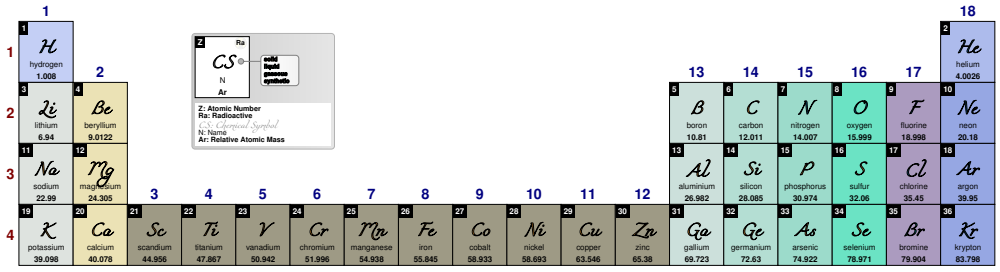
```
\pgfPT[CS all=red]
```



CS font default: *\small\bfseries*

Sets the font for the chemical symbol.

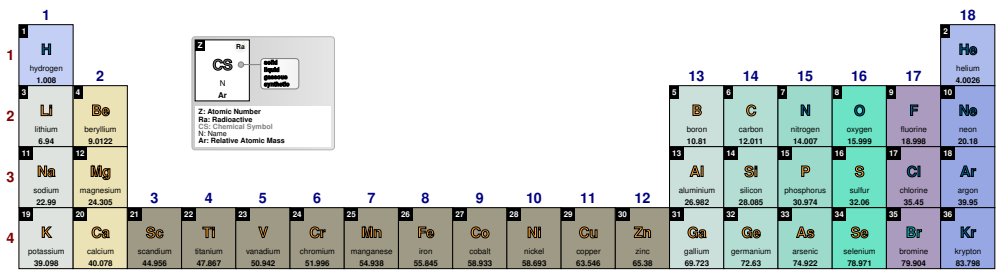
```
\pgfPT[Z list={1,...,36},CS font=\small\fontfamily{fmm}\selectfont]
```



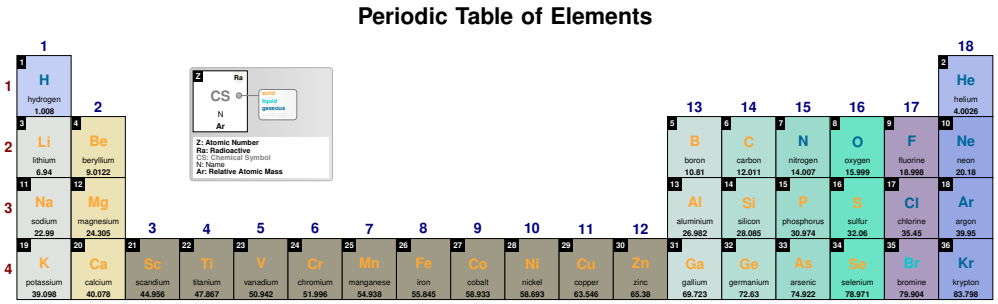
CS render mode default: *fill and outline*

Sets the chemical symbol render mode. Available modes are fill, outline or fill and outline.

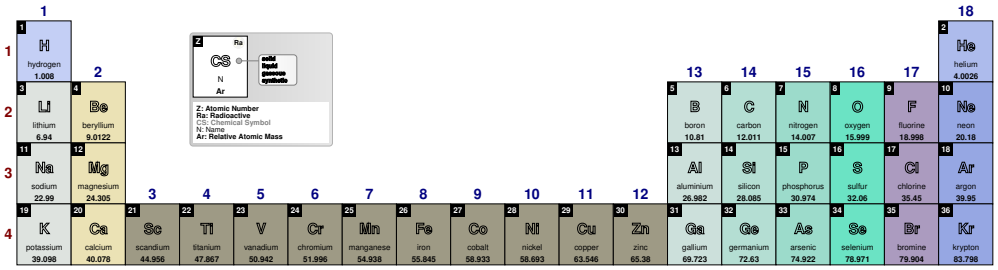
```
\pgfPT[Z list={1,...,36}]
```



\pgfPT[Z list={1,...,36},CS render mode=fill]



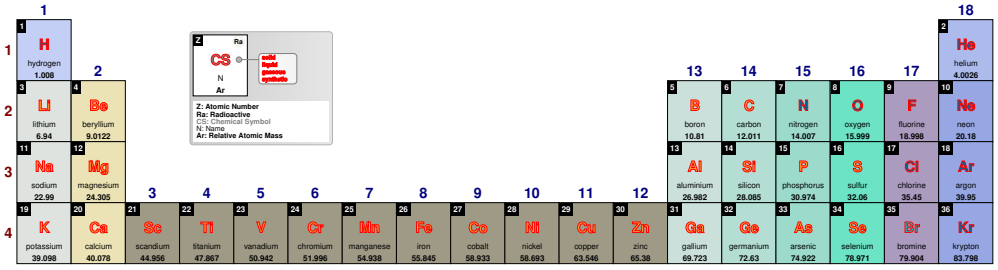
\pgfPT[Z list={1,...,36},CS render mode=outline]



CS outline color default: *black*

Sets the outline color for the chemical symbol.

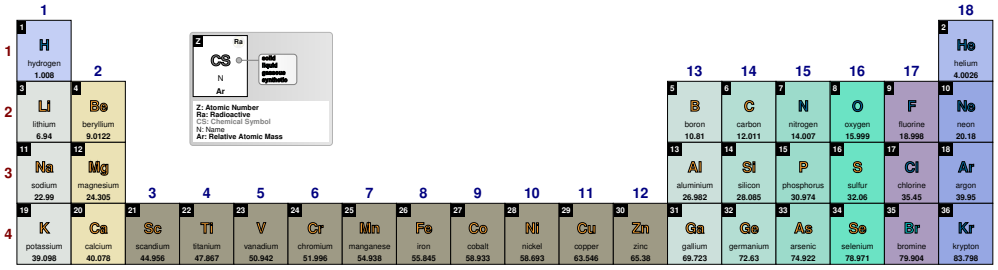
\pgfPT[Z list={1,...,36},CS outline color=red]



CS outline width default: *0.05*

Sets the outline width of the chemical symbol. It is any positive numerical value **without dimensions** (1.0 is roughly 1.0pt).

\pgfPT[Z list={1,...,36},CS outline width=.2]

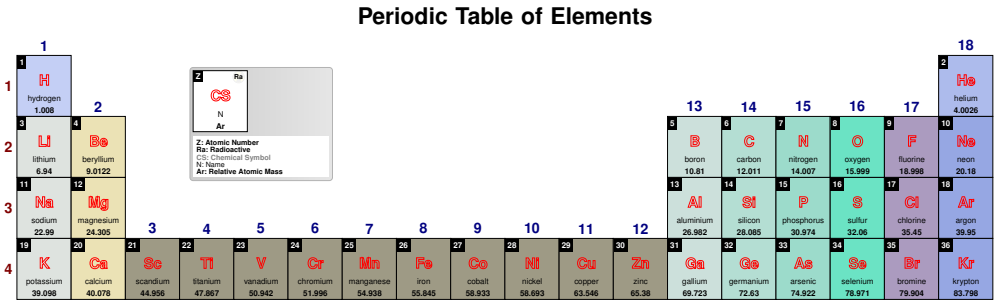


CS default: *{r=fill and outline,c=black,w=.05,f=\small\bfseries,s=solido,l=liquido, g=gasoso,sy=sintetico}*

Pseudo style to set the keys: CS **r**ender mode, CS **o**ut**l**ine **c**olor, CS **o**ut**l**ine **w**idth, CS font, CS **s**olid, CS **l**iquid, CS **g**as and/or CS **sy**nt and/or the style CS **all**. None of the *keys* – r, olc, olw , f, s, l, g, sy and all – are mandatory.

USAGE: CS={r=<fill|outline|fill and outline>,olc=<color>,olw=<positive numerical value> f=<font commands>,s=<color>,l=<color>,g=<color>,sy=<color>,all=<color>}

```
\pgfPT[Z list={1,...,36},CS={r=outline,olc=red,olw=.4},show legend pins=false]
```

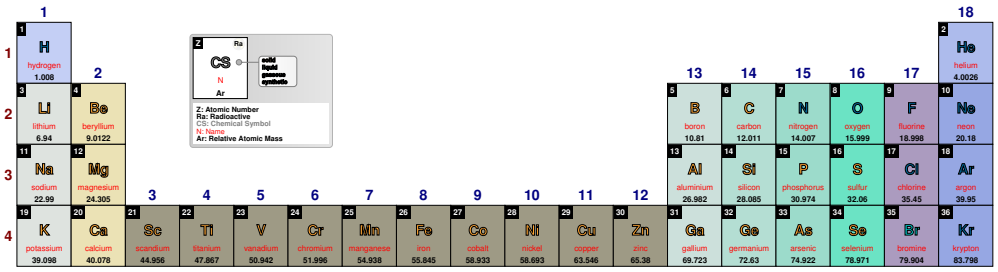


➠ The name

name color default: *black*

Sets the color of the element name.

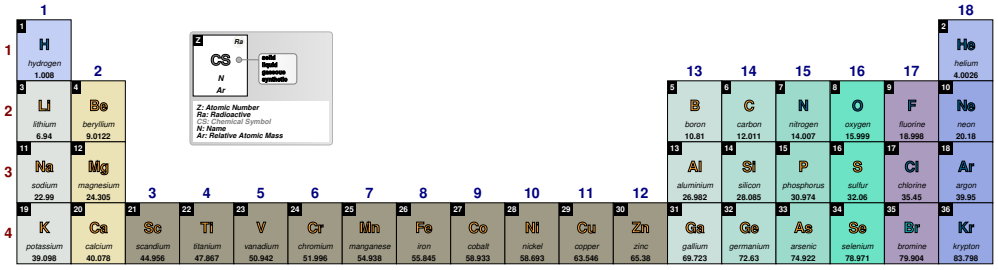
```
\pgfPTstyle[show title=false]
\pgfPT[Z list={1,...,36},name color=red]
```



name font default: *\fontseries\selectfont\tiny*

Sets the font of the element name.

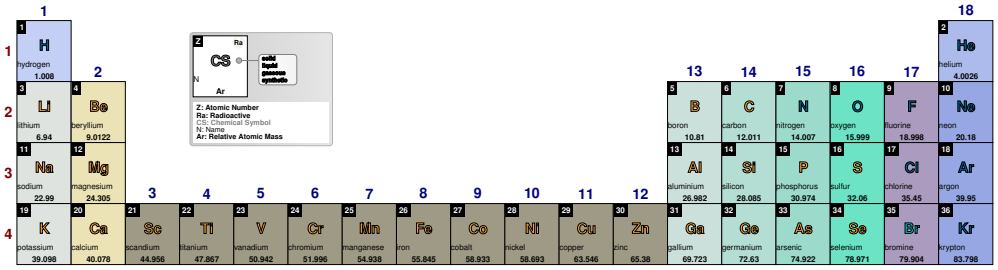
```
\pgfPT[Z list={1,...,36},name font=\itshape\tiny]
```



name align default: *center*

Sets the alignment of the element name to *left*, *center* or *right* with respect to its containing box. *(new in v1.0.1)*

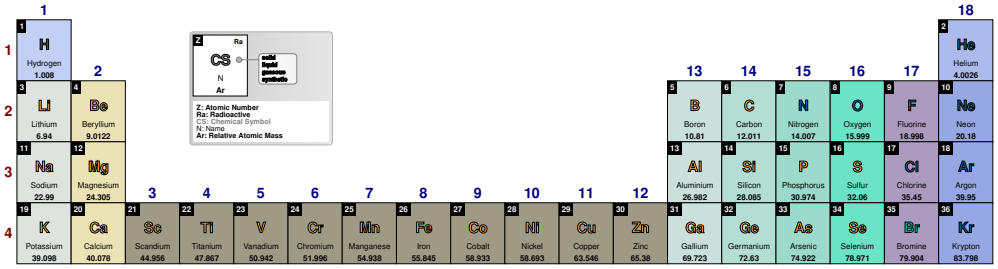
\pgfPT[Z list={1,...,36},name align=left]



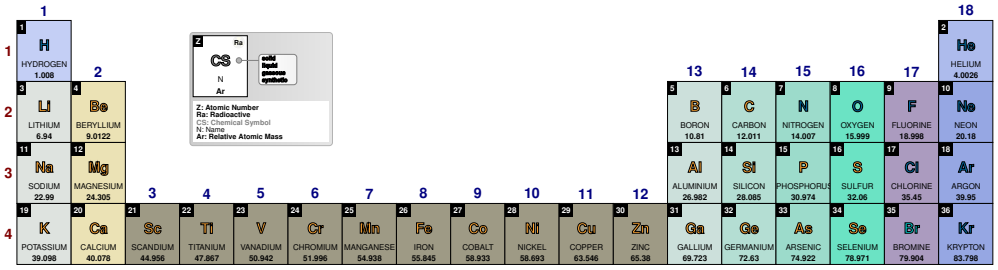
capitalize element names default: *false*

If set to true the first letter of the name of the elements is a capital letter (except in German where names start with a capital letter since version 2.0.1). If set to TRUE the whole name of the elements is in capital letters *(changed in v1.0.1)*

\pgfPT[Z list={1,...,36},capitalize element names=true]



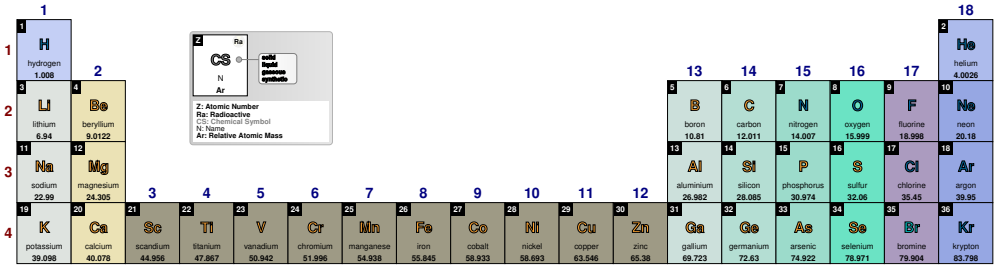
\pgfPT[Z list={1,...,36},capitalize element names=TRUE]



name *no value*

A style equivalent to capitalize element names=false

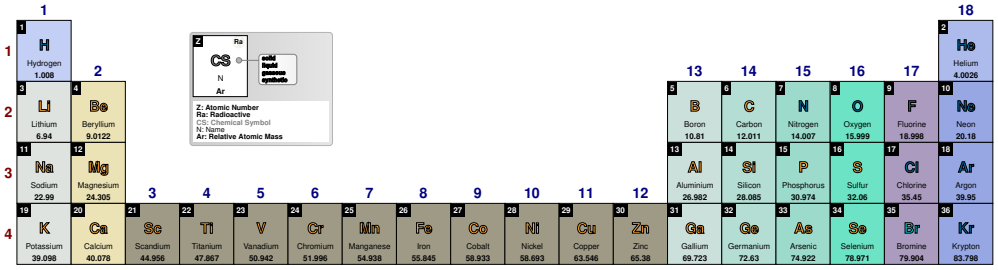
\pgfPT[Z list={1,...,36},name]



Name *no value*

A style equivalent to capitalize element names=true

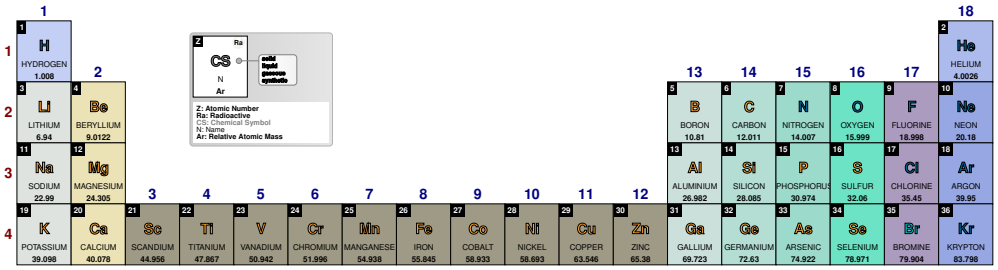
\pgfPT[Z list={1,...,36},Name]



NAME *no value*

A style equivalent to capitalize element names=TRUE *(new in v1.0.1)*

\pgfPT[Z list={1,...,36},NAME]

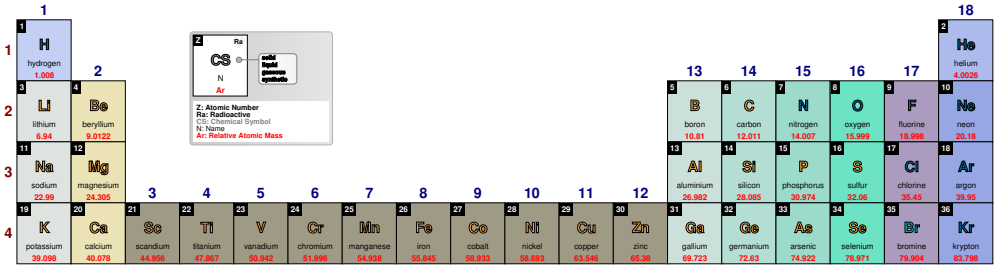


➠ The atomic weight

Ar color default: *black*

Sets the relative atomic mass color.

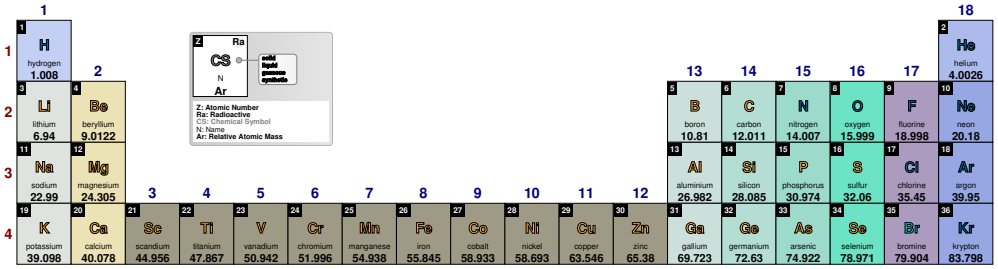
\pgfPT[Z list={1,...,36},Ar color=red]



Ar font default: *\tiny\bfseries*

Sets the relative atomic mass font.

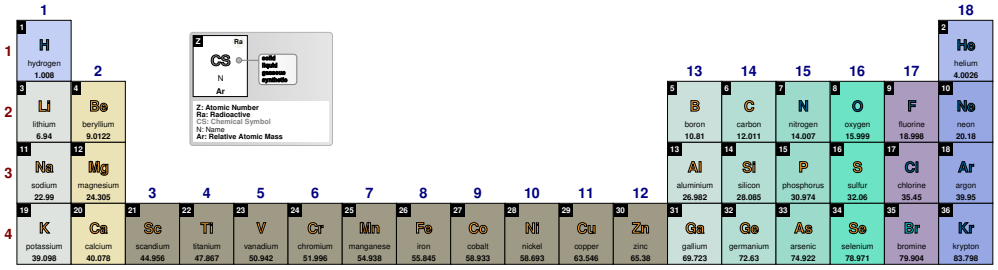
```
\pgfPT[Z list={1,...,36},Ar font=\scriptsize\bfseries]
```



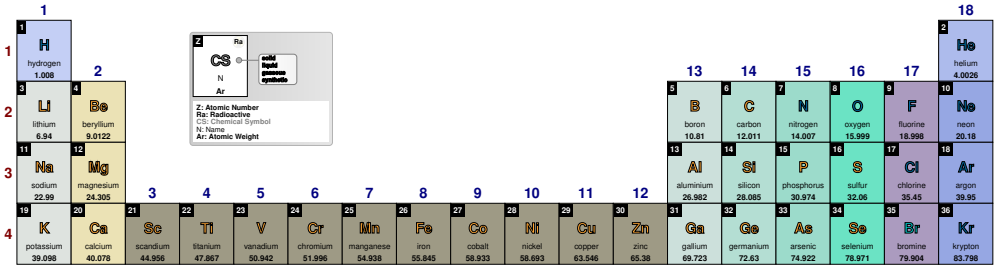
Ar label default: *m*

Sets the label to be used within the relative atomic mass description. When set to 'm' the term mass is used and when set to 'w' the term weight is used, resulting in *Relative Atomic Mass* and *Atomic Weight* labels respectively.

```
\pgfPT[Z list={1,...,36}]
```



```
\pgfPT[Z list={1,...,36},Ar label=w]
```

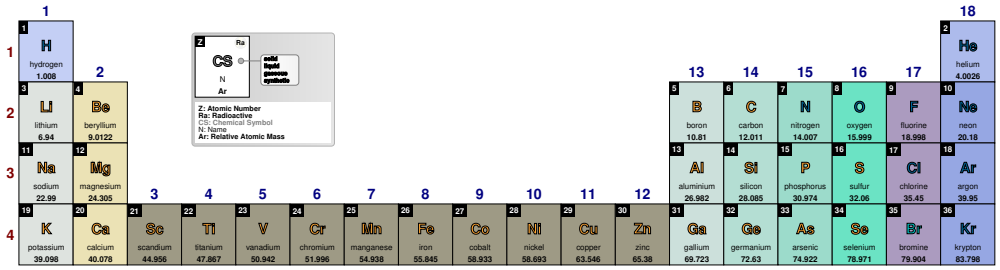


Ar precision default: *-1*

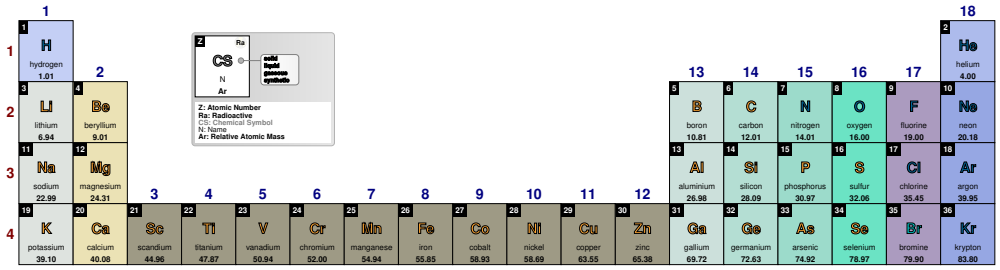
Sets the relative atomic mass precision, *i.e.*, the decimal places displayed in the relative atomic mass value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over the relative atomic mass data values witch actually have a maximum of 4 decimal places. So giving this key a value of -1 (the value of relative atomic mass as-is) or 4 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 3, i.e., -1, 0, 1, 2 or 3. Any other integer provided will be processed as -1.

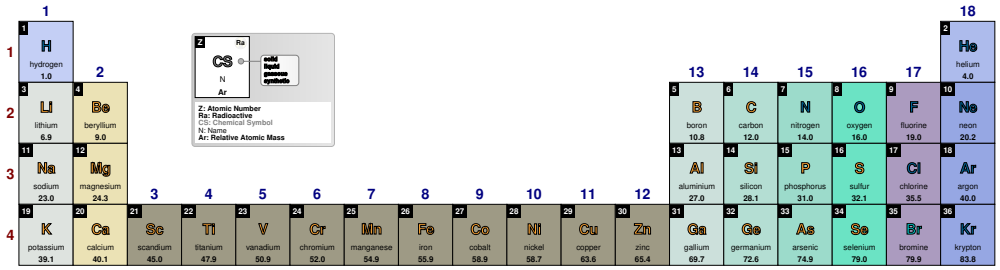
\pgfPT[Z list={1,...,36}]



\pgfPT[Z list={1,...,36},Ar precision=2]



\pgfPT[Z list={1,...,36},Ar precision=1]

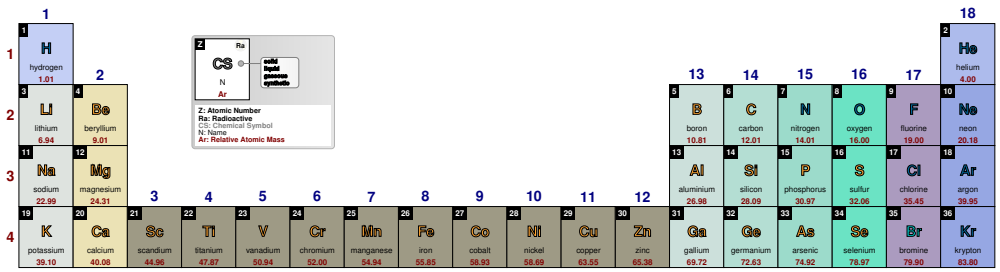


Ar default: *{c=black,f=\tiny\bfseries,l=m,p=-1}*

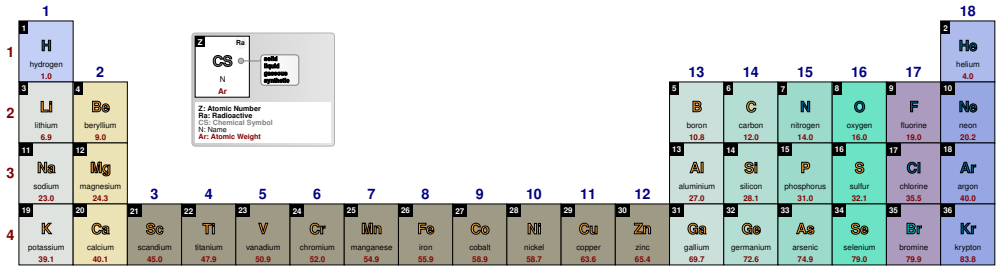
Pseudo style to set the keys: Ar **c**olor, Ar **f**ont, Ar **l**abel and/or Ar **p**recision. None of the *keys* – c, f, l and p – are mandatory.

USAGE: Ar={c=<color>,f=<font commands>,l=<m|w>p=<integer value>}

\pgfPT[Z list={1,...,36},Ar={c=red!50!black,p=2}]



```
\pgfPT[Z list={1,...,36},Ar={c=red!50!black,p=1,l=w}]
```

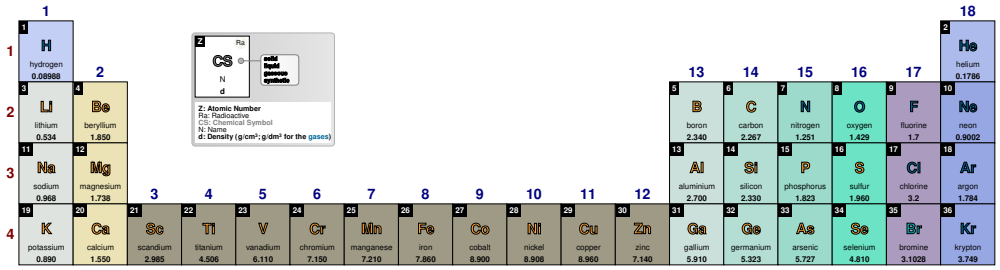


➠ The density

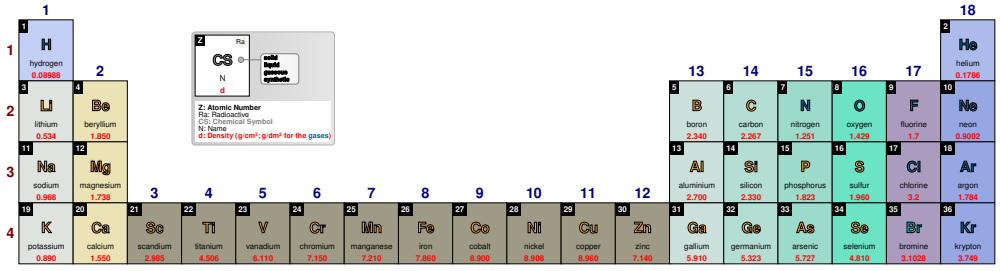
d color default: *black*

Sets the density value text color.

```
\pgfPTbuilcellstyle{myd}(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;d)]
\pgfPT[Z list={1,...,36},cell style=myd,show title=false]
```



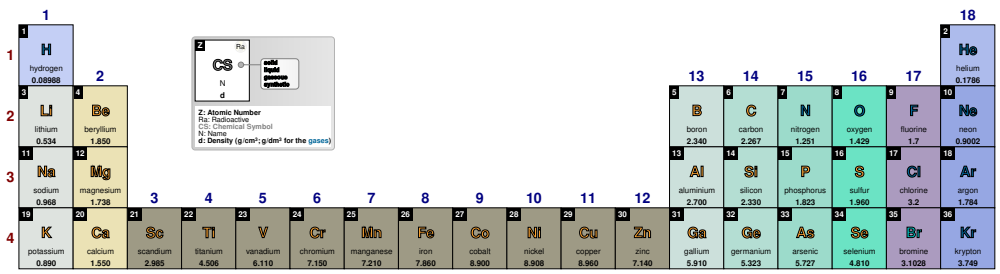
```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false,d color=red]
```



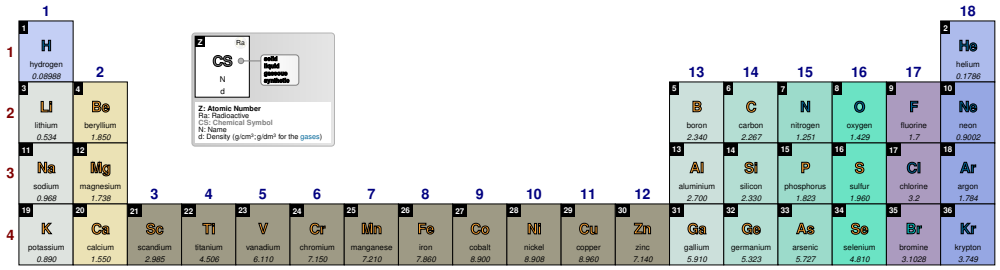
d font default: *\tiny\bfseries*

Sets the density value text font.

```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false]
```



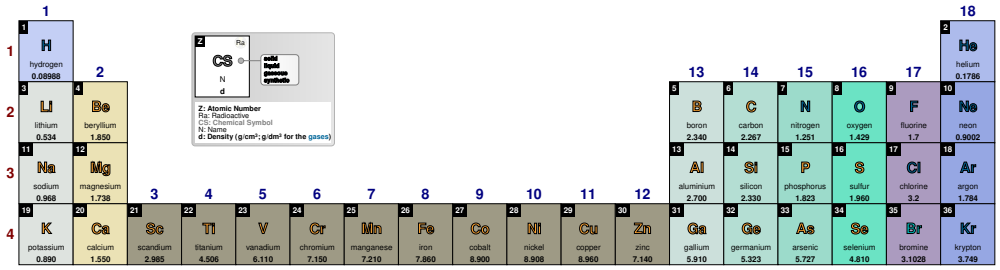
```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false,d font=\tiny\itshape]
```



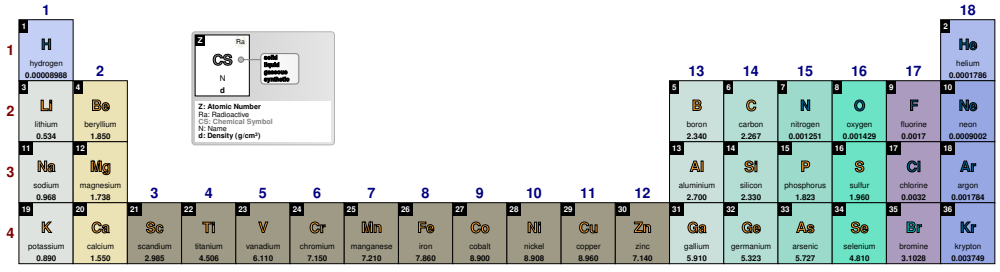
d unit default: *both*

Sets the unit for the density of the elements. The two possible values to this key are g/dm3 (g/dm^3), g/cm3 (g/cm^3) and both (g/dm^3 for elements in the gaseous state and g/cm^3 for all other elements). *(new in v1.0.1)*

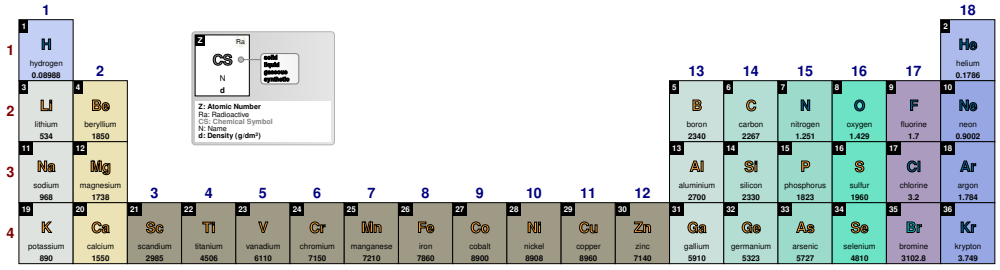
```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false]
```



```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false,d unit=g/cm3]
```



```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false,d unit=g/dm3]
```

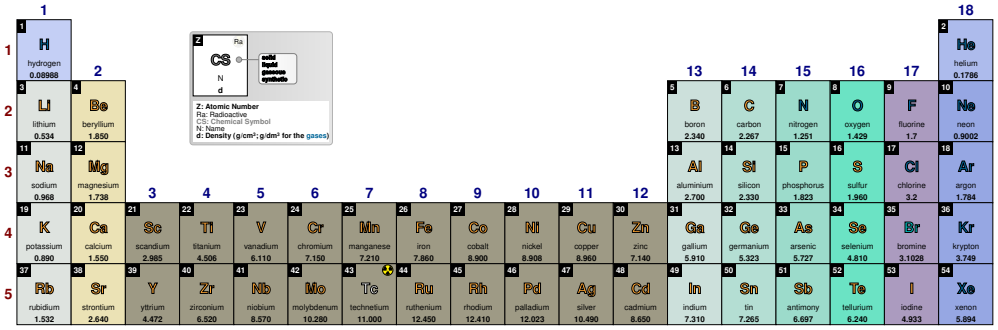


d precision default: *-1*

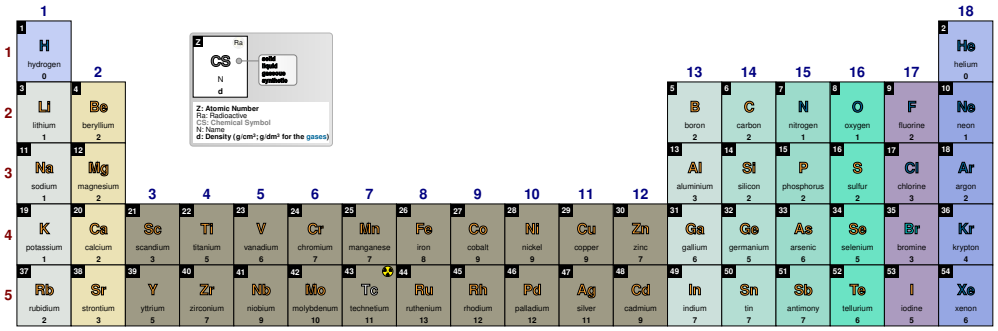
Sets the density precision, *i.e.*, the decimal places displayed in their value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over density values witch actually have a maximum 5 or 8 decimal places, when the values are in g/dm^3 or in g/cm^3, respectively.. So giving this key a value of -1 (the value of the melting or boiling point as-is) or 5 or 8 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 4 (g/dm^3) or 7 (g/cm^3). Any other integer provided will be processed as -1.

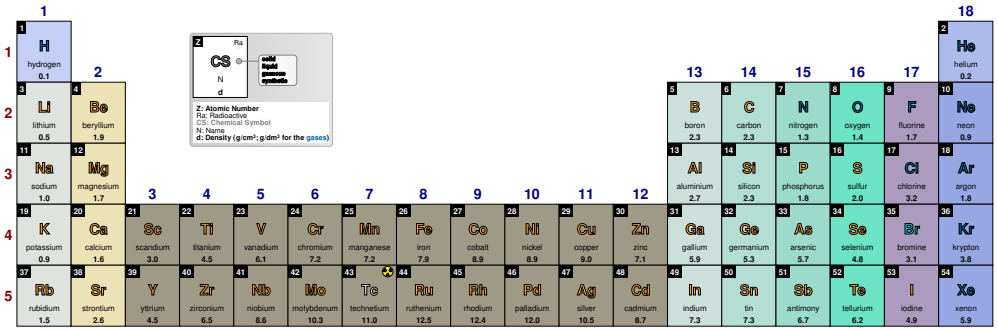
```
\pgfPTstyle[Z list={1,...,54},cell style=myd,show title=false]
\pgfPT
```



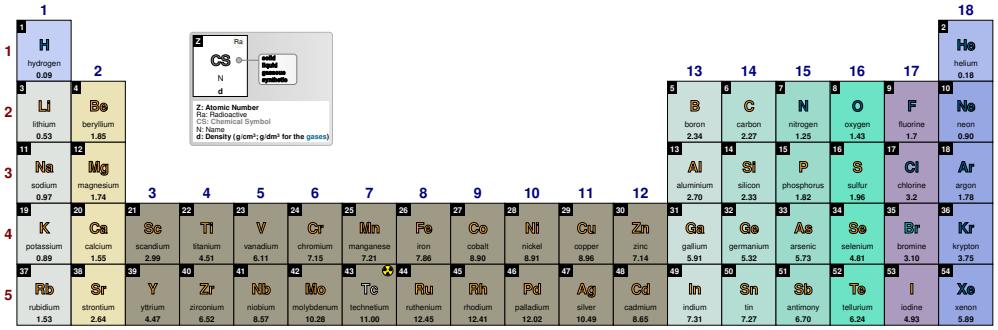
```
\pgfPT[d precision=0]
```



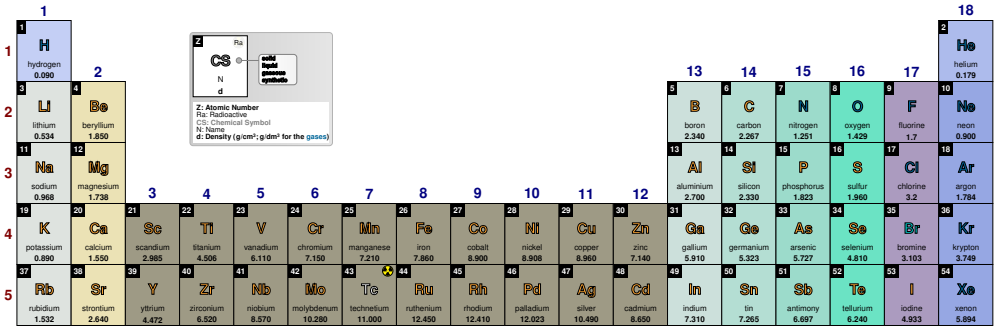
```
\pgfPT[d precision=1]
```



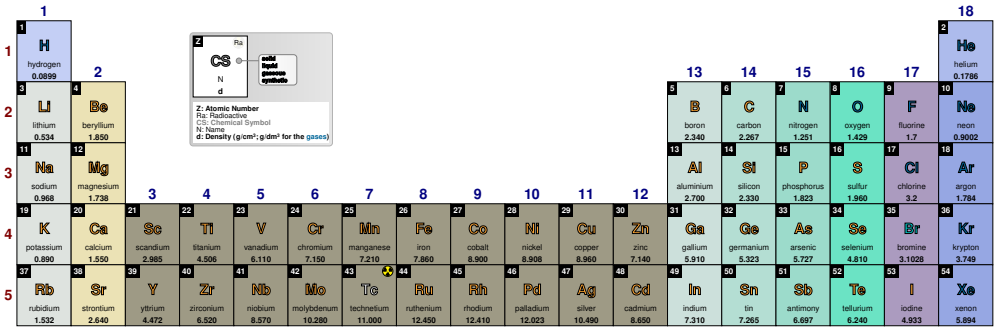
\pgfPT[d precision=2]



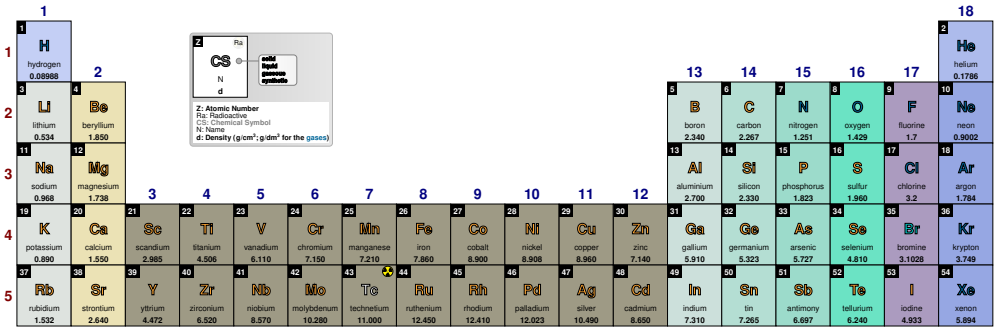
\pgfPT[d precision=3]



\pgfPT[d precision=4]



\pgfPT[d precision=5]

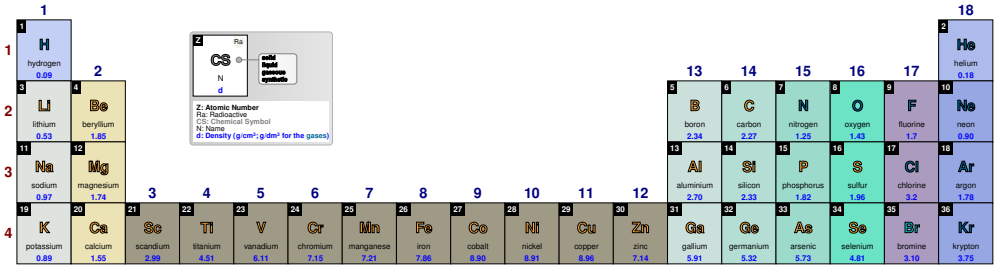


d default: *{c=black,f=\tiny\bfseries,p=-1,u=both}*

Pseudo style to set the keys: d **c**olor, d **f**ont, d **p**recision and/or d **u**nit. None of the *keys* – c, f, p and u – are mandatory. *(new in v1.0.1)*

USAGE: d={c=<color>,f=<font commands>,p=<integer value>,u=<pm|A>}

```
\pgfPT[Z list={1,...,36},cell style=myd,show title=false,d={c=blue,p=2}]
```

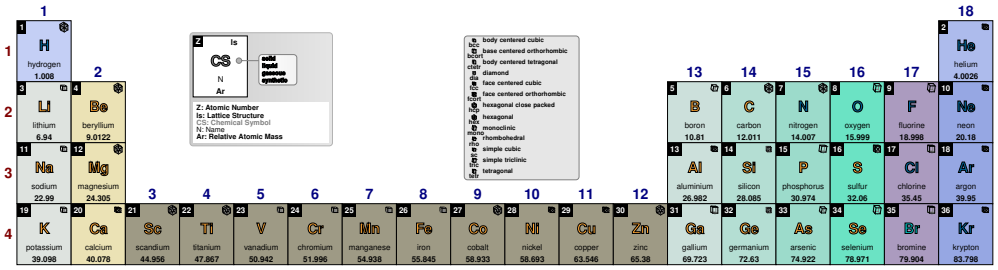


➠ The lattice structure

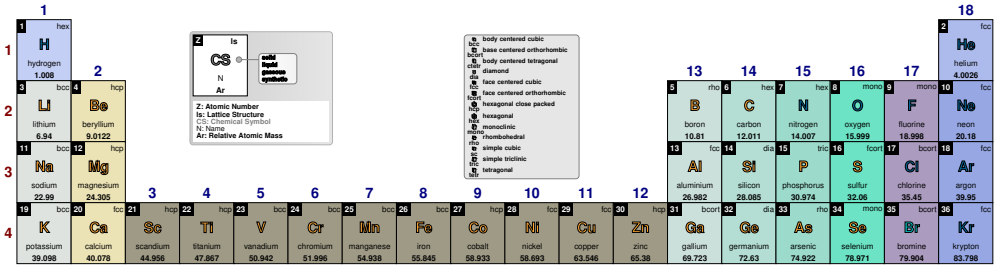
ls default: *fig*

Sets what is displayed for the lattice structure: a figure (fig) or text (txt) or both (fig+txt or txt+fig).

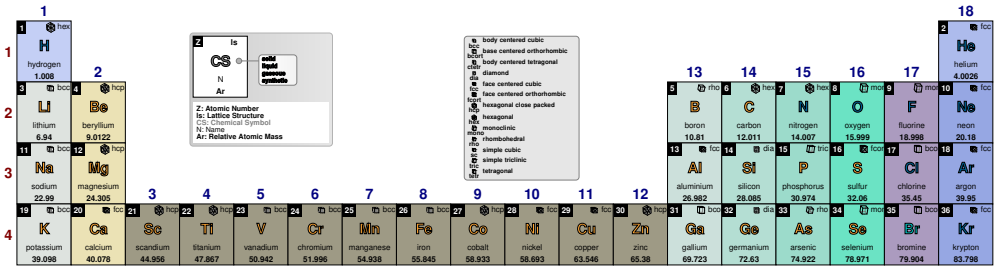
```
\pgfPT[Z list={1,...,36},cell style=pgfPTls]
```



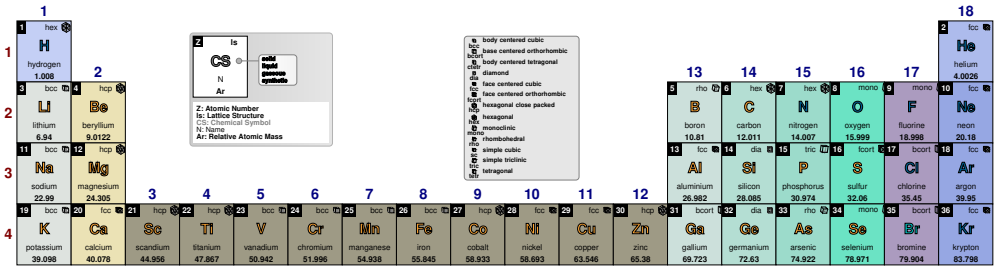
```
\pgfPT[Z list={1,...,36},cell style=pgfPTls,ls=txt]
```



```
\pgfPT[Z list={1,...,36},cell style=pgfPTls,ls=fig+txt]
```



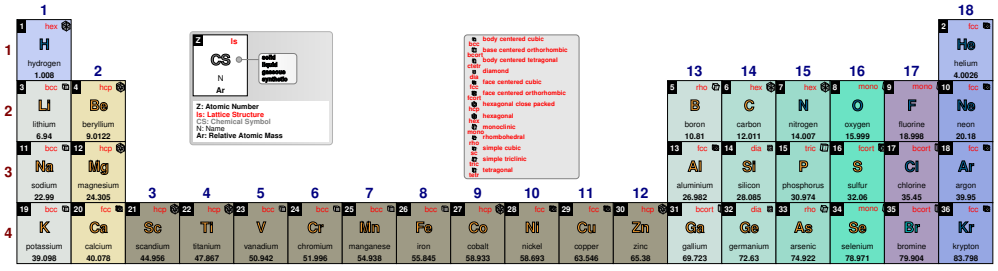
```
\pgfPT[Z list={1,...,36},cell style=pgfPTls,ls=txt+fig]
```



ls color default: *black*

Sets the lattice structure text color.

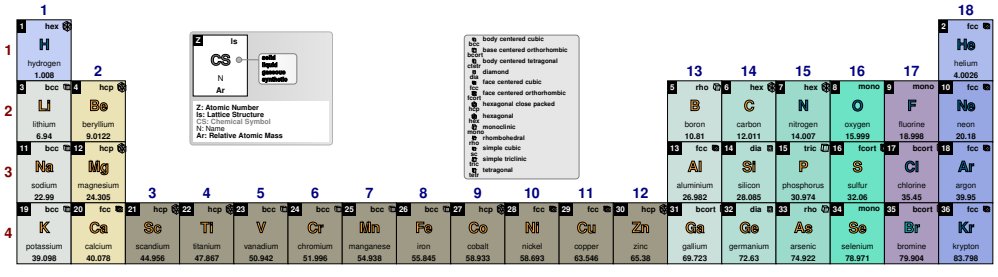
```
\pgfPT[Z list={1,...,36},cell style=pgfPTls,ls=txt+fig,ls color=red]
```



ls font default: *\tiny*

Sets the lattice structure text font.

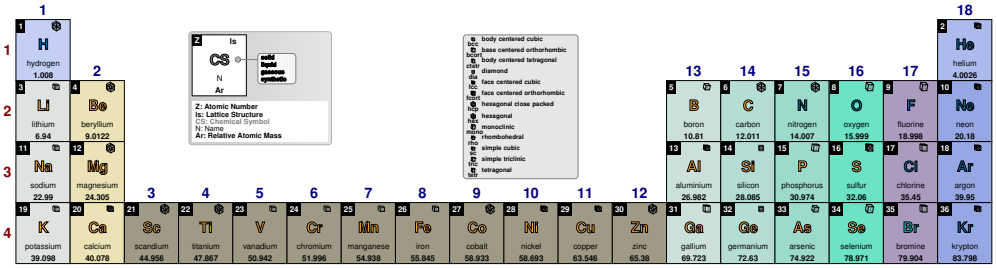
```
\pgfPT[Z list={1,...,36},cell style=pgfPTls,ls=txt+fig,ls font=\tiny\bfseries]
```



ls align default: *right*

Sets the alignment of the lattice structure to left or center or right.

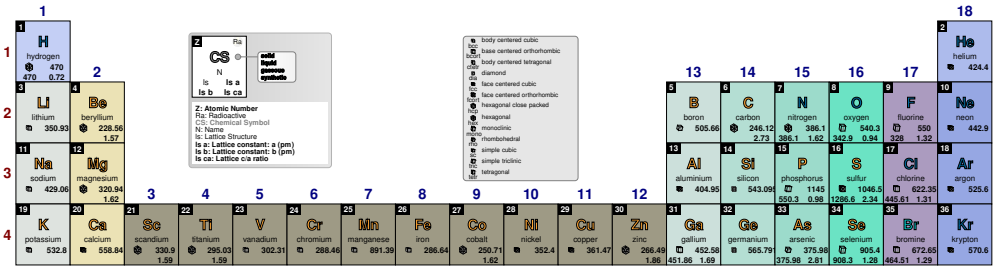
```
\pgfPT[Z list={1,...,36},cell style=pgfPTls,ls align=center]
```



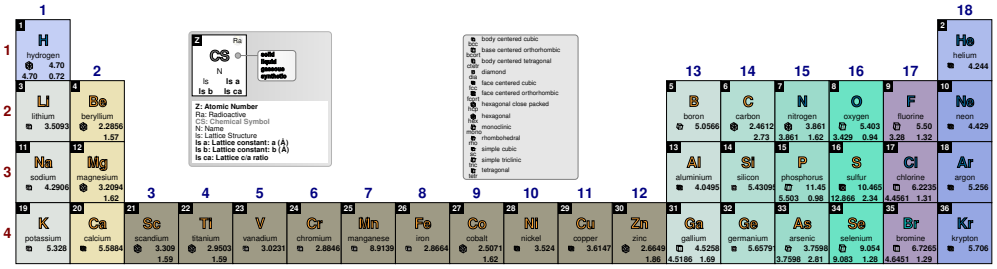
ls unit default: *pm*

Sets the unit for the lattice structure constants: a, b and c. The two possible values to this key are pm (picometers) and A (Å – angstroms). *(new in v1.0.1)*

```
\pgfPTbuilcellstyle{myls}(6,3)% 6 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-2.5;ls),(5;2.5-3;lsa),
(6;1-2.5;lsb),(6;2.5-3;lsca)]
\pgfPTstyle[Z list={1,...,36},cell style=myls,show title=false,ls align=center]
\pgfPT
```



```
\pgfPT[ls unit=A]
```



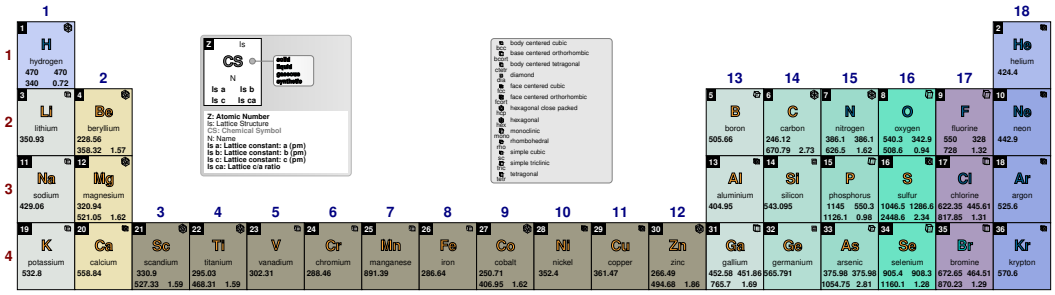
ls precision default: *-1*

Sets the lattice structure constants - a, b, and c - precision, as also the lattice c/a ratio, *i.e.*, the decimal places displayed in their value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over the constants data or c/a ratio values witch actually have a maximum of 2 or 4 decimal places, when the values are in picometers or in angstroms, respectively. So giving this key a value of -1 (the value of the constants or c/a ratio as-is) or 2 or 4 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 1 (pm) or 3 (Å). Any other integer provided will be processed as -1.

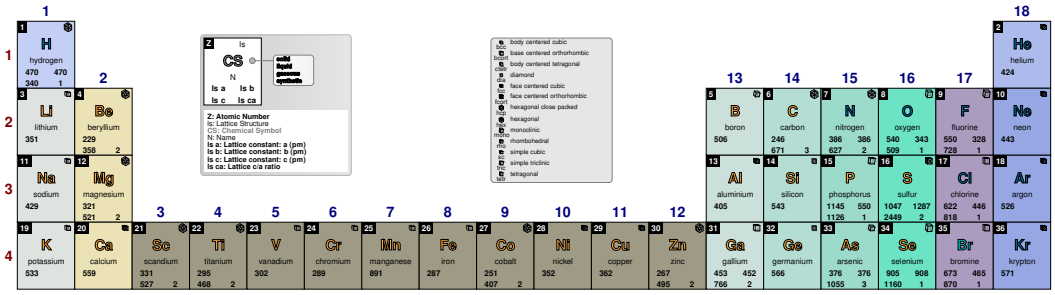
```
\pgfPTbuilcellstyle{myls}(6,3)% 6 rows by 3 columns
[(1;1-2;Z),(1;2-3;ls),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-2.5;lsa), (5;2.5-3;lsb),(6;1-2.5;lsc),(6;2.5-3;lsca)]
\pgfPTstyle[Z list={1,...,36},cell={w=36pt,h=42pt,style=myls}]
\pgfPT
```

Periodic Table of Elements



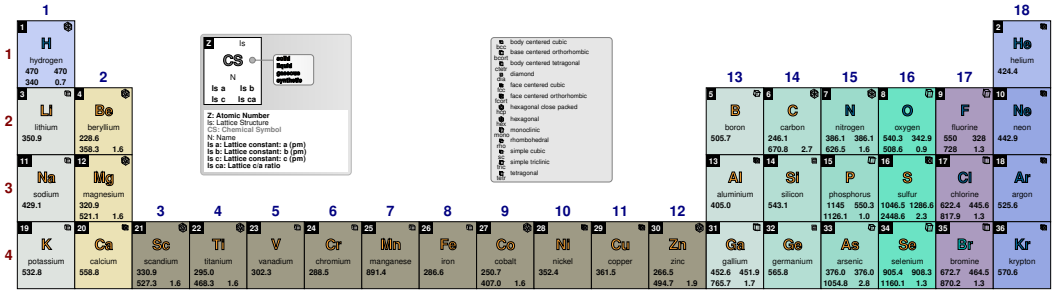
```
\pgfPT[ls precision=0]
```

Periodic Table of Elements



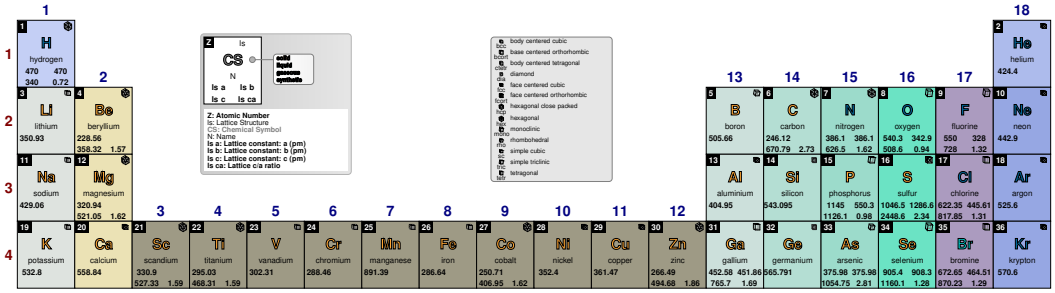
```
\pgfPT[ls precision=1]
```

Periodic Table of Elements

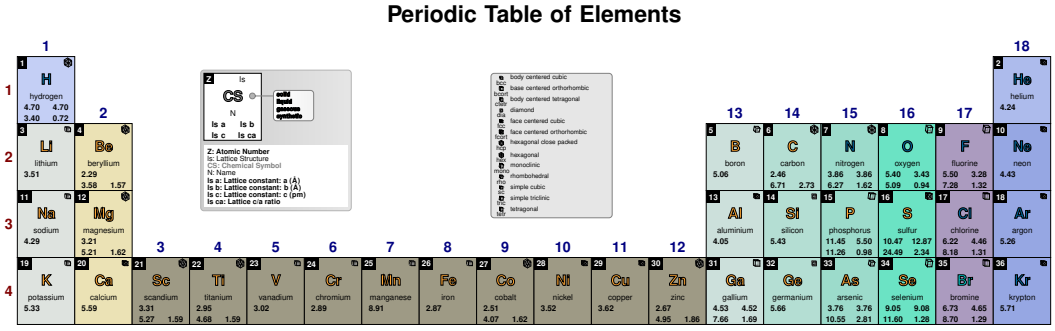


```
\pgfPT[ls precision=2]
```

Periodic Table of Elements



```
\pgfPT[ls precision=2,ls unit=A]
```

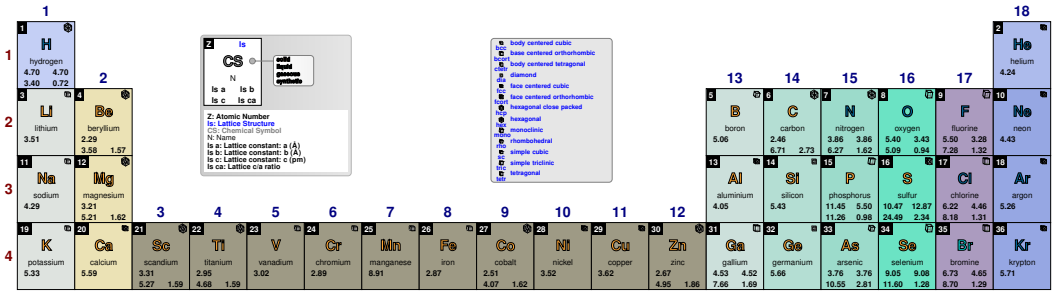


lat default: *{ls=fig,c=black,f=\tiny,align=right,p=-1,u=pm}*

Pseudo style to set the keys: **ls**, ls **c**olor, ls **f**ont, ls **align**, ls **p**recision and/or ls **u**nit. None of the *keys* – ls, c, f, align, p and u – are mandatory. *(new in v1.0.1)*

USAGE: lat={ls=<fig|txt|fig+txt|txt+fig>,c=<color>,f=<font commands>, align=<left|center|right>,p=<integer value>,u=<pm|A>}

```
\pgfPT[Z list={1,..,36},lat={c=blue,f=\tiny\bfseries,u=A,p=2}]
```

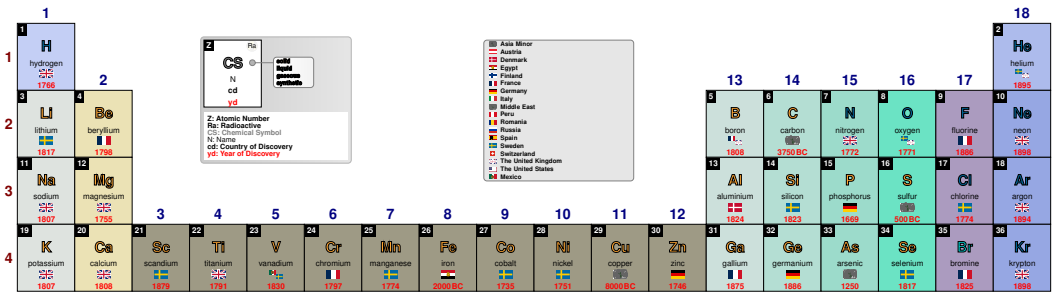


The discovery year

DiscY color default: *black*

Sets the discovery year color.

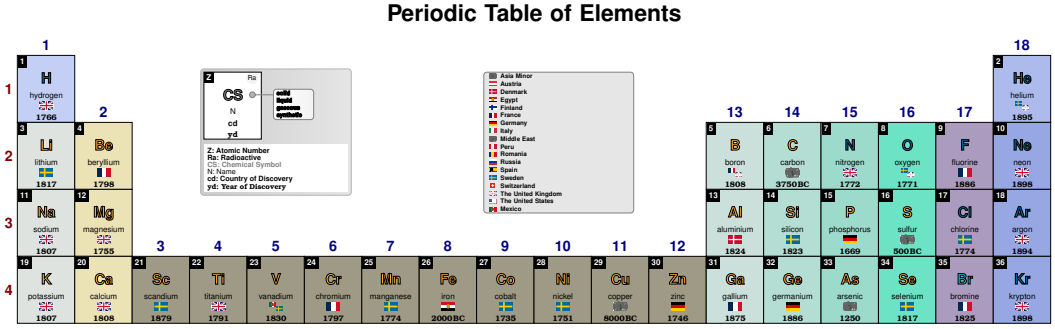
```
\pgfPT[Z list={1,..,36},cell style=pgfPTdisc,DiscY color=red]
```



DiscY font default: *\tiny\bfseries*

Sets the discovery year font.

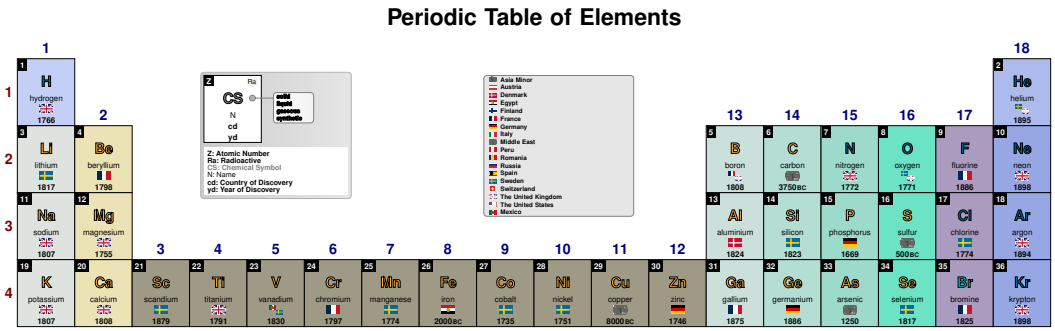
```
\pgfPT[Z list={1,...,36},cell style=pgfPTdisc,DiscY
font=\fontfamily{pbk}\selectfont\tiny\bfseries]
```



DiscY BC scale default: *1*

Sets the font factor scaling for the Before Christ (BC) acronym in the discovery year.

```
\pgfPT[Z list={1,...,36},cell style=pgfPTdisc,DiscY BC scale=.8]
```



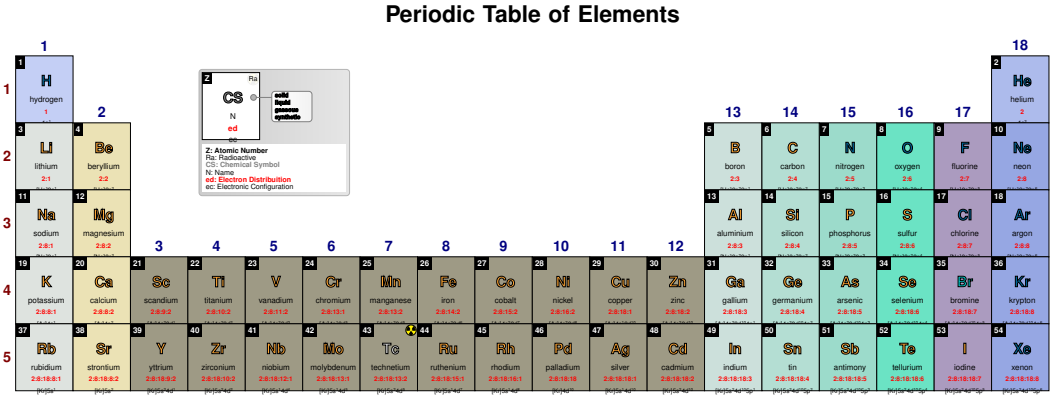
➠ *The electron distribution*

```
\pgfPTbuilcellstyle{electron}(6,3)% 6 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name), (5;1-3;eDist),(6;1-
3;eConfignl)]
```

eDist color default: *black*

Sets the electron distribution color.

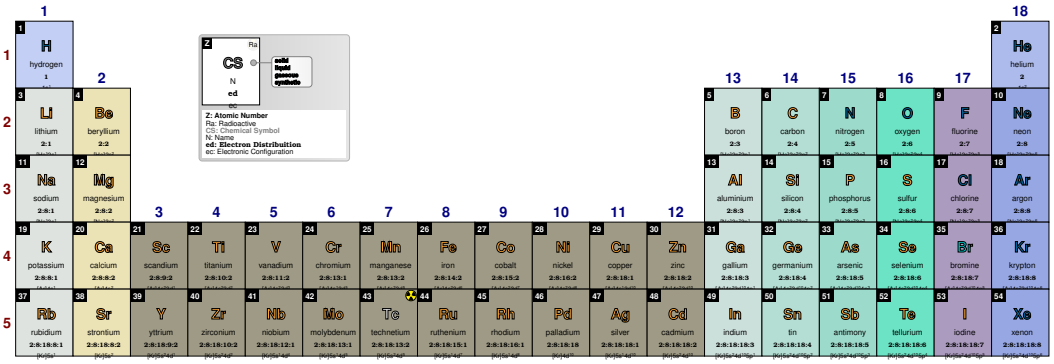
```
\pgfPT[Z list={1,...,54},cell style=electron,eDist color=red]
```



eDist font default: *\tiny\bfseries*

Sets the electron distribution font.

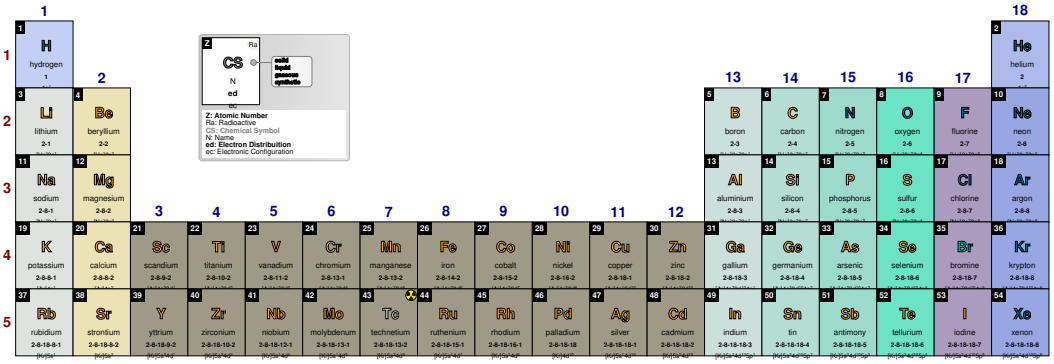
```
\pgfPT[Z list={1,...,54},cell style=electron,eDist
font=\fontfamily{pbk}\selectfont\tiny\bfseries]
```



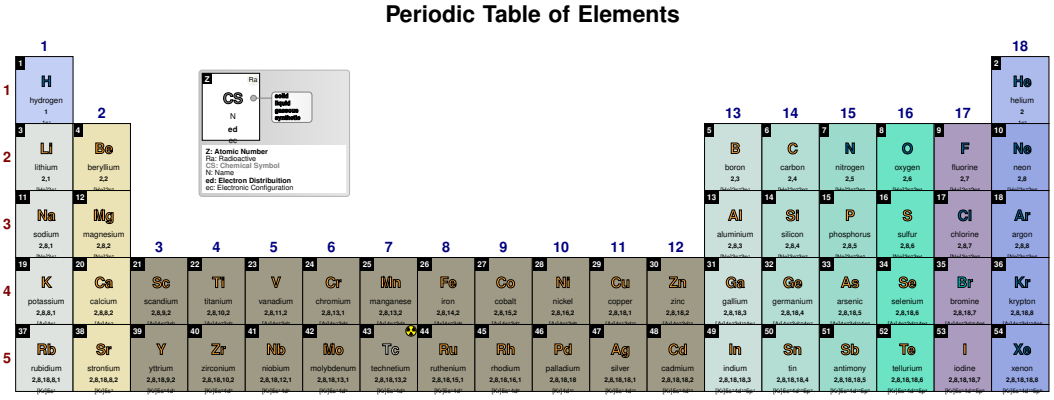
eDist sep default: *:*

Sets the separator character between energy levels in electron distribution. *If the separator character is a comma it must be provided between curly braces – {,}.*

```
\pgfPT[Z list={1,...,54},cell style=electron,eDist sep=-]
```



```
\pgfPT[Z list={1,...,54},cell style=electron,eDist sep={,}]
```



➠ *The other contents*

For all the *other contents* available for the cells of the periodic table, two keys can be set: <content name> color and <content name> font.

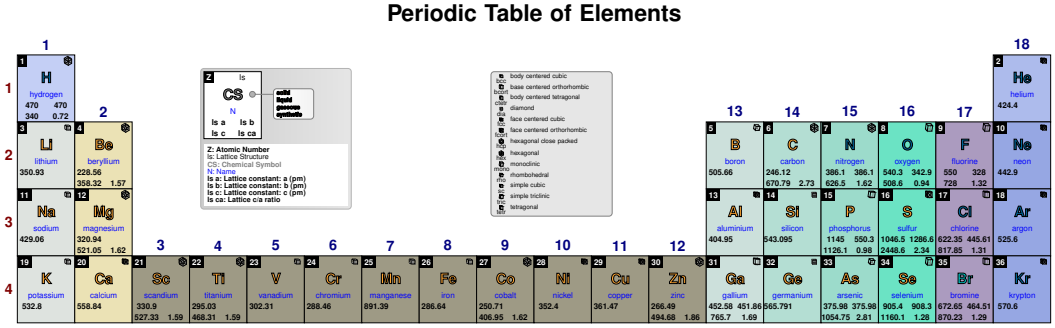
The <content name>'s list:

- ✓ **R**: atomic radius
- ✓ **Rcov**: covalente radius
- ✓ **Rion**: ionic radius
- ✓ **Ei**: first ionization energy
- ✓ **eneg**: electronegativity (Pauling)
- ✓ **eaff**: electroaffnity
- ✓ **O**: oxidation states
- ✓ **Tmelt**: melting point (Kelvin)
- ✓ **TmeltC**: melting point (Celsius degrees)
- ✓ **Tboil**: boiling point (Kelvin)
- ✓ **TboilC**: boiling point (Celsius degrees)
- ✓ **eConfign**: electronic configuration (increasing n)
- ✓ **eConfignl**: electronic configuration (increasing n+ℓ)
- ✓ **Cp**: specific heat capacity
- ✓ **kT**: thermal conductivity
- ✓ **lsa**: lattice constant – a
- ✓ **lsb**: lattice constant – b
- ✓ **lsc**: lattice constant – c
- ✓ **lsca**: lattice c/a ratio
- ✓ **DiscC**: discover country
- ✓ **spectra**: visible range spectral lines

<content name> color default: *black*

Sets the <content name> color.

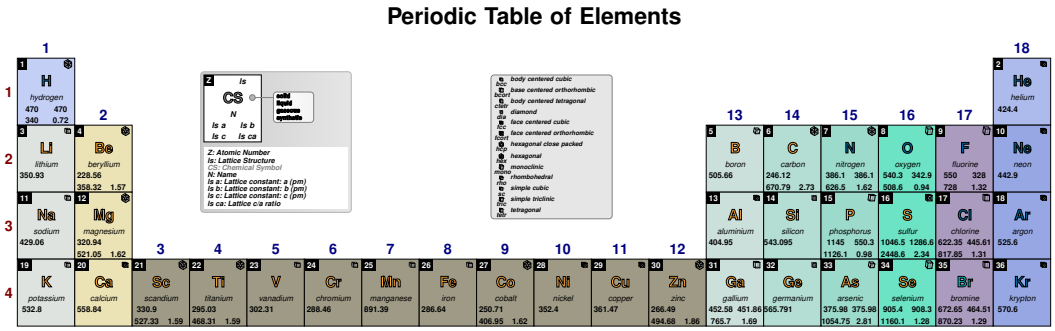
```
\pgfPT[Z list={1,...,36},name color=blue]
```



<content name> font default: *\tiny\bfseries*

Sets the <content name> font.

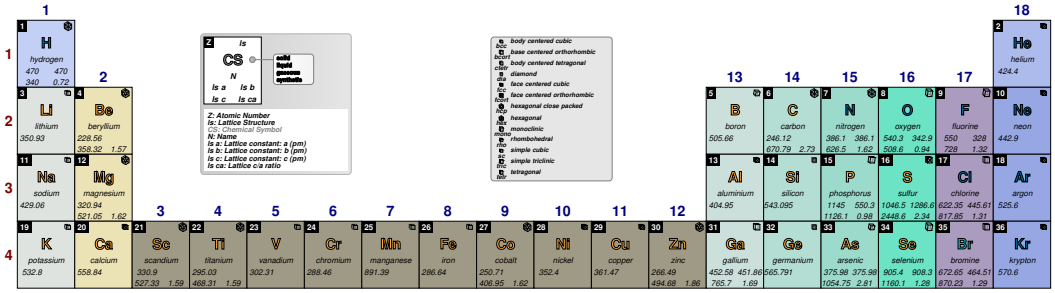
```
\pgfPT[Z list={1,...,36},name font=\tiny\itshape]
```



cell font default: *\bfseries\tiny*

Style to set the font for all cell contents, except for the Z and Chemical Symbol fonts.

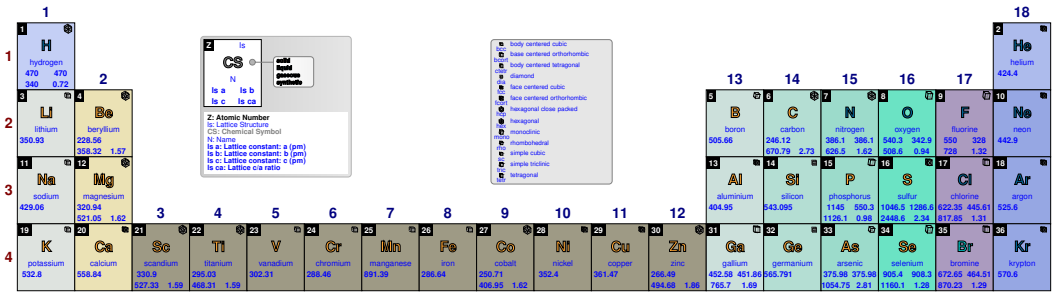
```
\pgfPT[Z list={1,...,36},cell font=\tiny\itshape]
```



cell color default: *black*

Style to set the color for all cell contents, except for the Z and Chemical Symbol colors.

```
\pgfPT[Z list={1,...,36},cell color=blue]
```



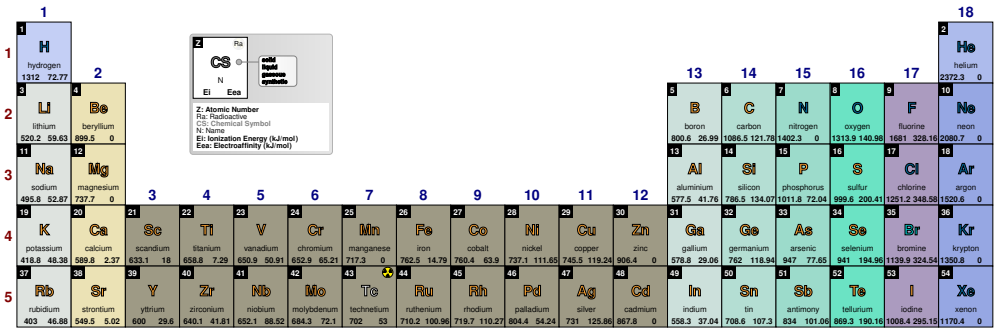
The precision of the *other contents*, which have numerical values, can also be set by a key. *Atomic radius, covalent radius, and ionic radius all have integer values, so precision does not apply to them*.

E precision default: *-1*

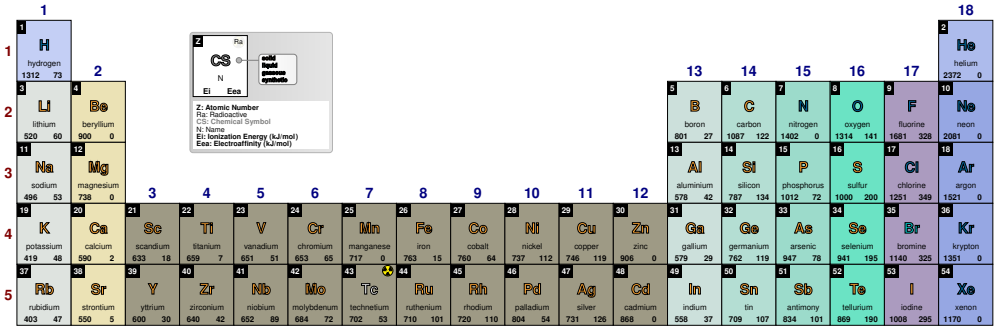
Sets the first ionization energy and the electroaffnity precision, *i.e.*, the decimal places displayed in their value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over energy values witch actually have a maximum of 3 decimal places. So giving this key a value of -1 (the value of the energy as-is) or 3 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 2, i.e., -1, 0, 1 or 2. Any other integer provided will be processed as -1.

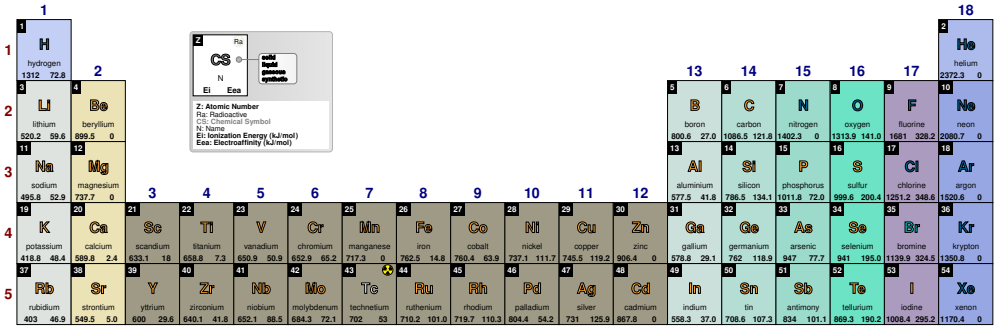
```
\pgfPTbuilcellstyle{myE}(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-2.5;Ei), (5;2.5-3;eaff)]
\pgfPTstyle[Z list={1,...,54},cell style=myE,show title=false]
\pgfPT
```



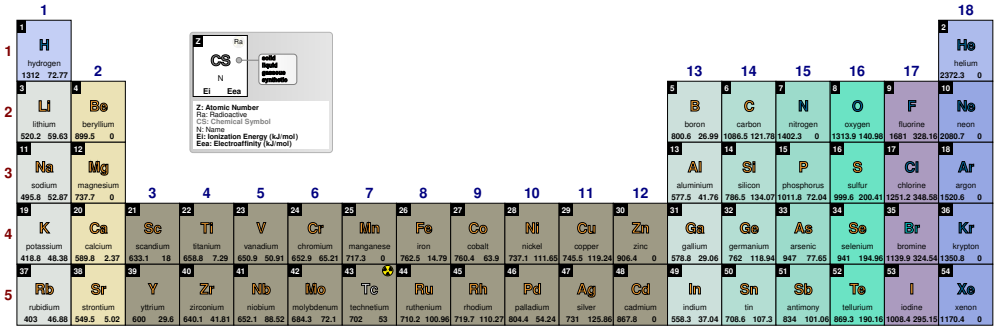
```
\pgfPT[E precision=0]
```



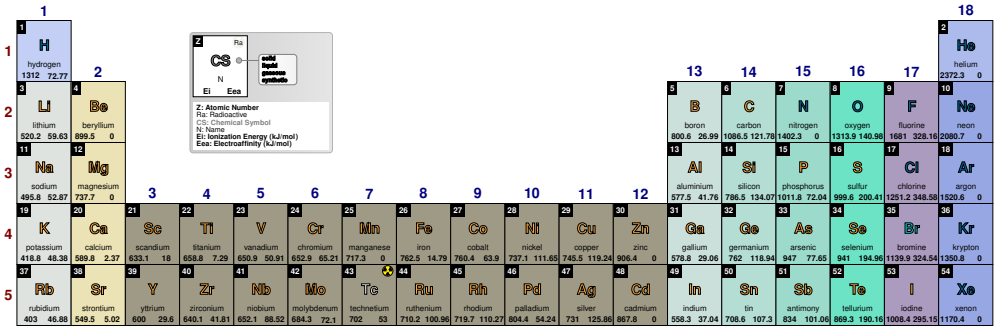
```
\pgfPT[E precision=1]
```



\pgfPT[E precision=2]



\pgfPT[E precision=3]

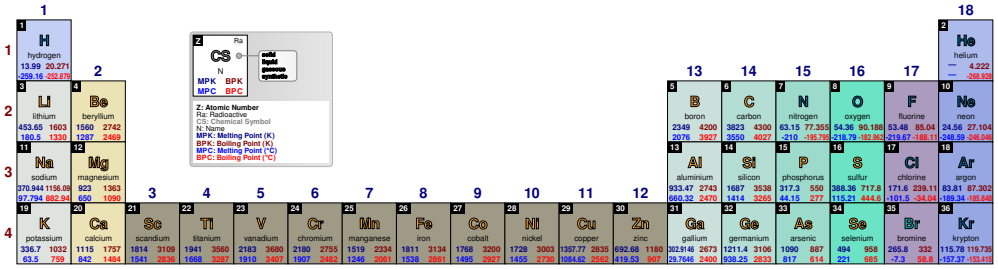


T precision default: *-1*

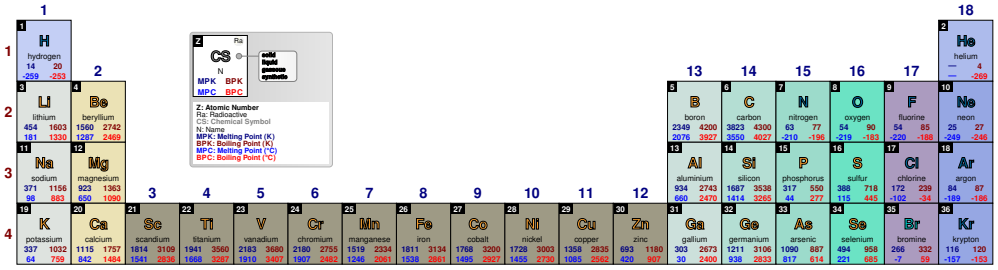
Sets the melting point an boiling point precision, *i.e.*, the decimal places displayed in their value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over melting or boiling point values witch actually have a maximum, respectively, of 4 or 2 decimal places. So giving this key a value of -1 (the value of the melting or boiling point as-is) or, respectively, 4 or 2 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 3 or 2. Any other integer provided will be processed as -1.

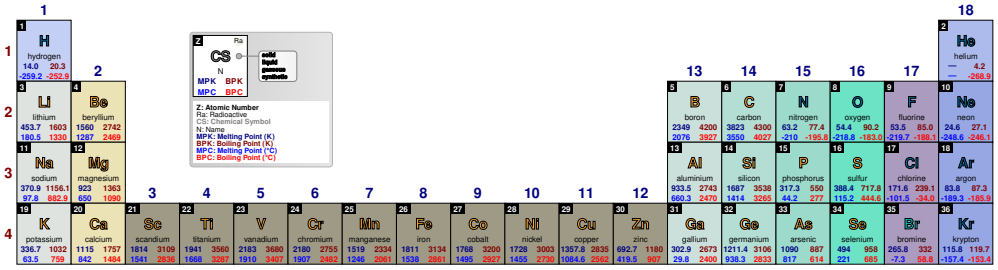
```
\pgfPTbuilcellstyle{myT}(6,3)% 6 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-2.5;Tmelt), (5;2.5-3;Tboil),(6;1-2.5;TmeltC),(6;2.5-3;TboilC)]
\pgfPTstyle[Z list={1,...,36},cell style=myT,Tmelt color=blue!50!black,TmeltC color=blue,Tboil color=red!50!black,TboilC color=red,show title=false]
\pgfPT
```



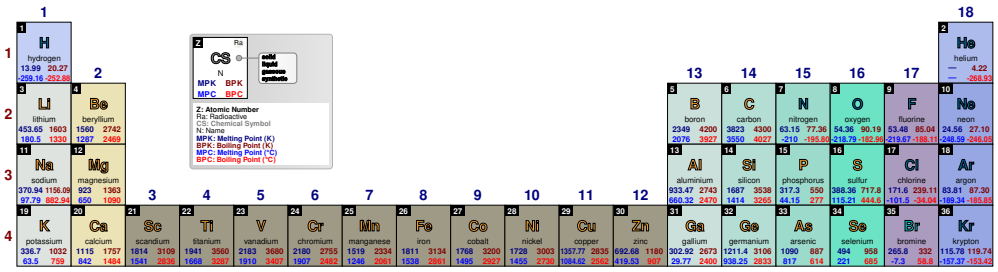
\pgfPT[T precision=0]



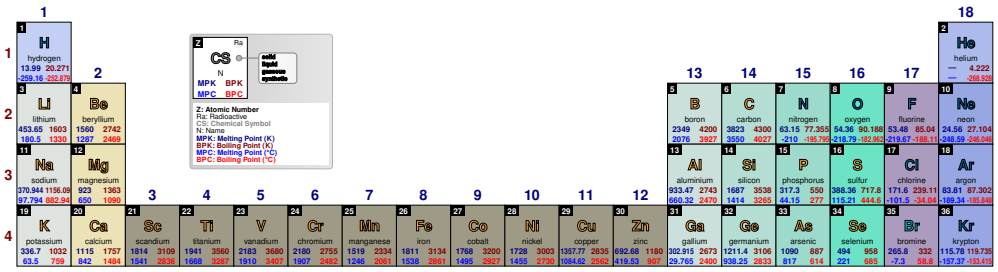
\pgfPT[T precision=1]



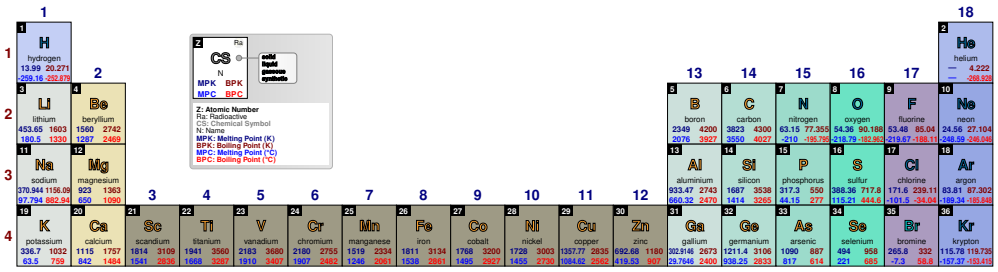
\pgfPT[T precision=2]



\pgfPT[T precision=3]



\pgfPT[T precision=4]

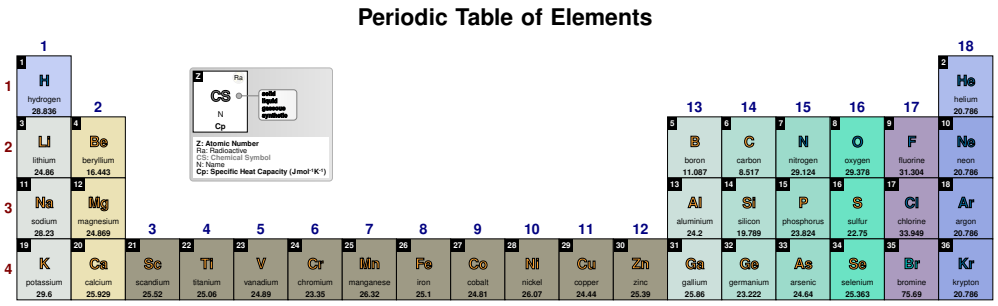


Cp precision default: *-1*

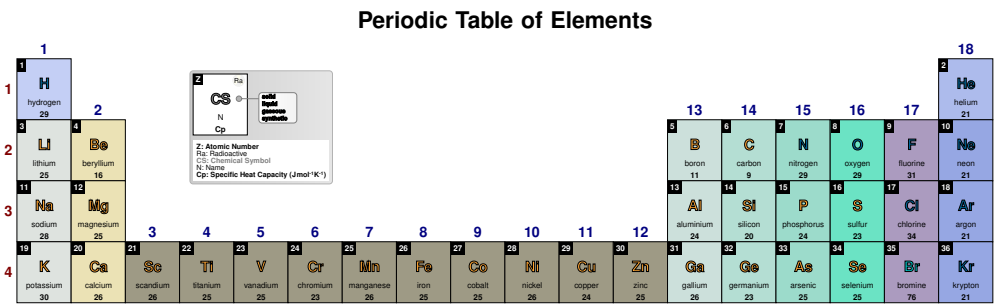
Sets the specific heat capacity precision, *i.e.*, the decimal places displayed in their value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over density values witch actually have a maximum 3 decimal places. So giving this key a value of -1 (the value of the melting or boiling point as-is) or 3 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 2. Any other integer provided will be processed as -1.

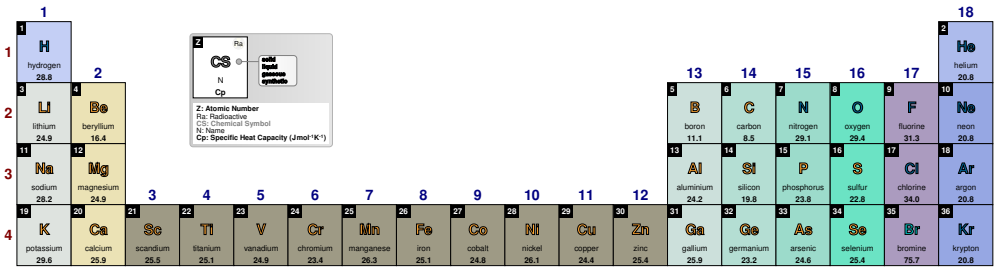
```
\pgfPTbuilcellstyle{myCp}(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;Cp)]
\pgfPTstyle[Z list={1,...,36},cell style=myCp]
                                                \pgfPT
```



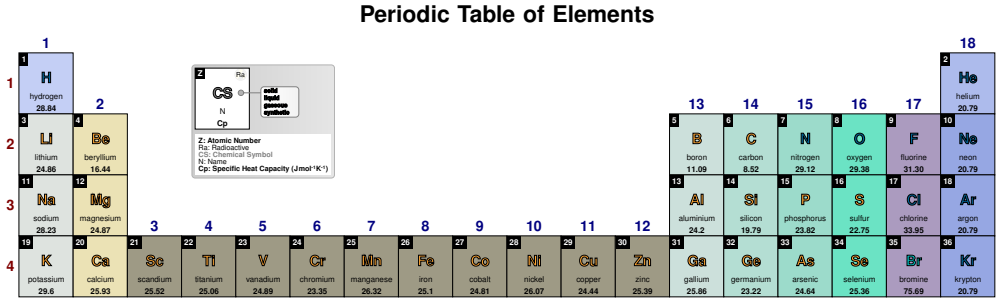

```
\pgfPT[Cp precision=0]
```



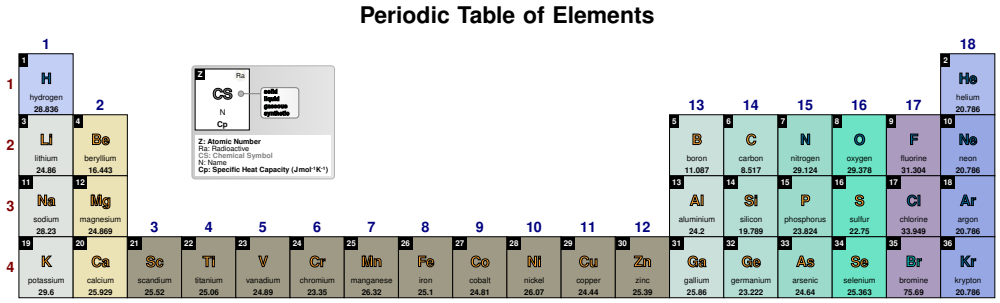

```
\pgfPT[Cp precision=1]
```



```
\pgfPT[Cp precision=2]
```




```
\pgfPT[Cp precision=3]
```

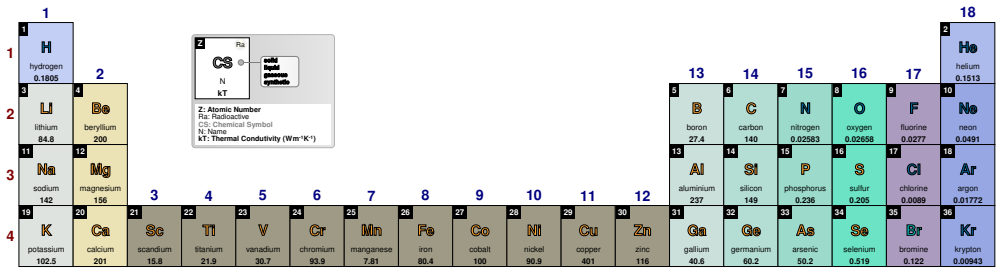


kT precision default: *-1*

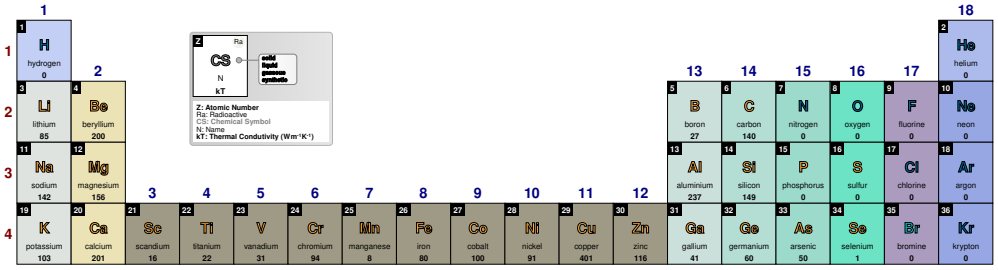
Sets the thermal condutivity precision, *i.e.*, the decimal places displayed in their value, performing the respective rounding, without zero padding the value.

NOTE:
Rounding is performed over density values witch actually have a maximum 5 decimal places. So giving this key a value of -1 (the value of the melting or boiling point as-is) or 5 has the same effect.
Therefore the values provided to this key should be any integer between -1 and 4. Any other integer provided will be processed as -1.

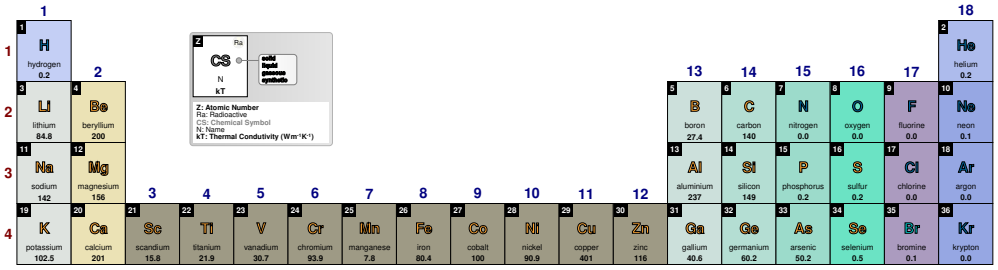
```
\pgfPTbuilcellstyle{mykT}(5,3)% 5 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;kT)]
\pgfPTstyle[Z list={1,...,36},cell style=mykT,show title=false]
                                                    \pgfPT
```



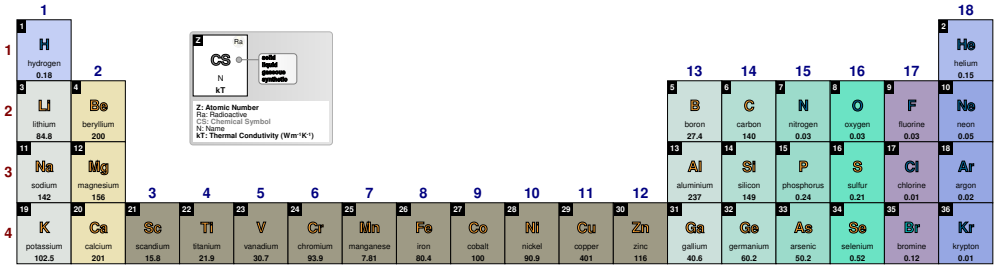
\pgfPT[kT precision=0]



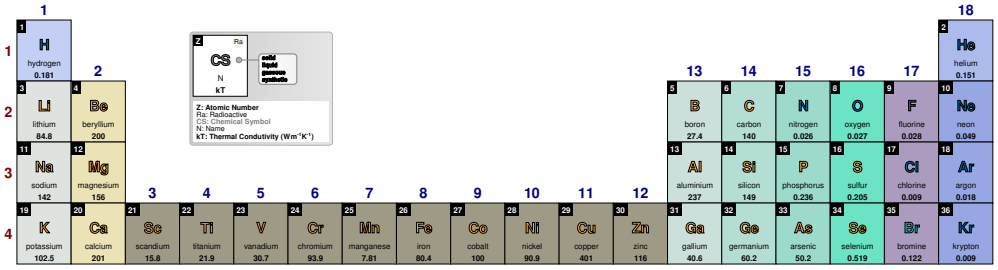
\pgfPT[kT precision=1]



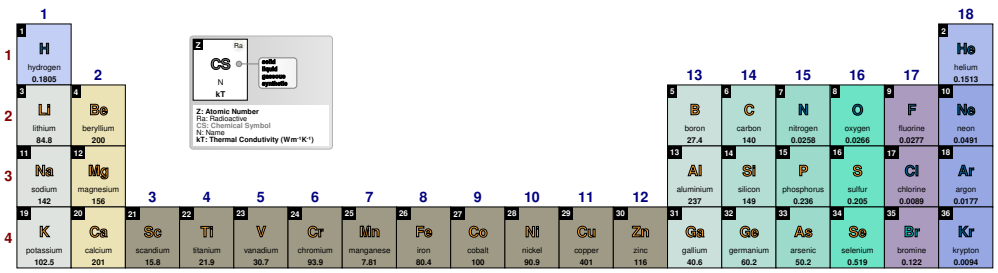
\pgfPT[kT precision=2]



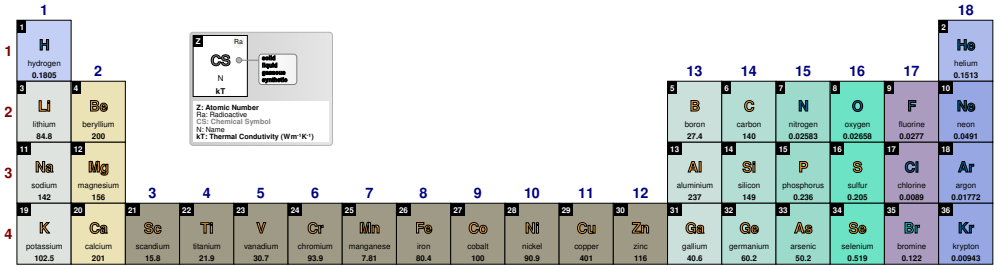
\pgfPT[kT precision=3]



\pgfPT[kT precision=4]



\pgfPT[kT precision=5]

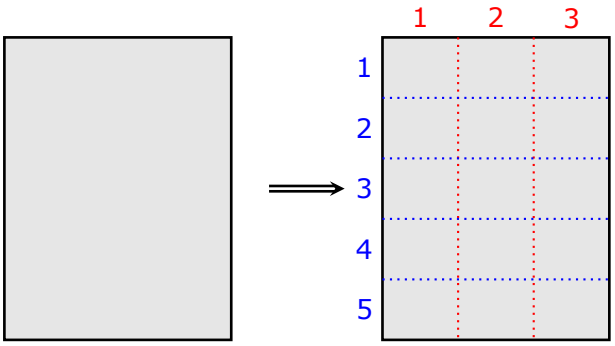


\pgfPTresetstyle

Designing cells with \pgfPTbuildcell

To start designing the *base cell* of the Periodic Table it is necessary to keep in mind that each cell will be split into **n** rows and **k** columns.

As a running example, 5 rows and 3 columns will be used:

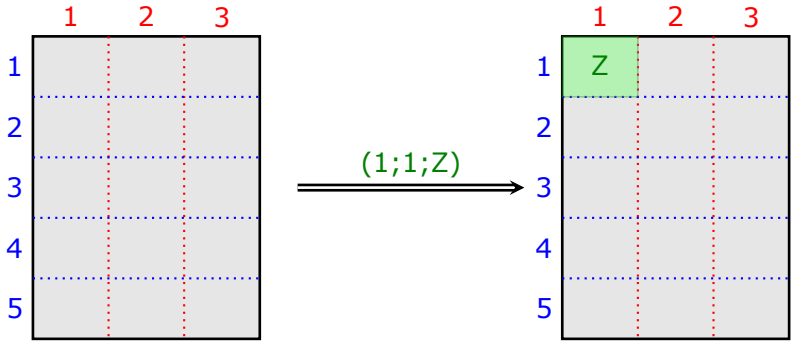


The next task is to assign contents to the cell by typing *trios* with the structure

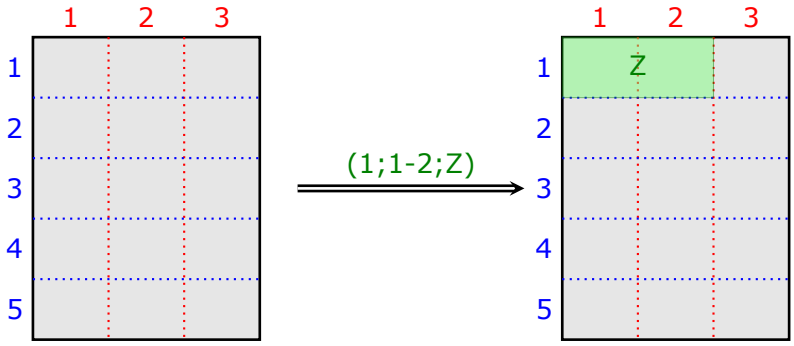
- **(row;column;content)**
- or **(start row-end row;start column-end column;content)**
- or a combination of both.

The available **contents** are: Z, name, CS, Ar, Ar*, radio, R, Rcov, Rion, Ei, eneg, eaff, O, Tmelt, TmeltC, Tboil, TboilC, eDist, eConfign, eConfignl, d, Cp, kT, ls, lsa, lsb, lsc, lsca, DiscY, DiscC and spectra.

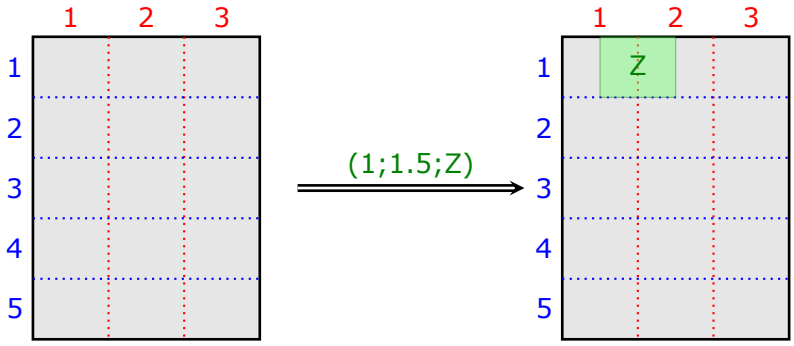
Assigning, for instance, (1;1;Z) will show the atomic number in the first row and in the first column,



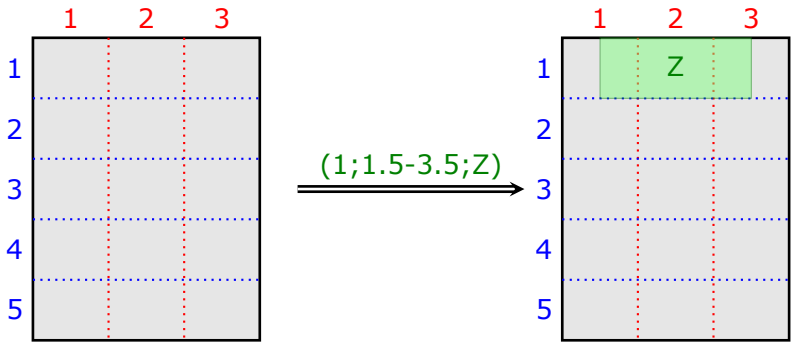
while the assignment (1;1-2;Z) will show the atomic number in the first row and filling the first and second columns,



It is also possible to start at a *fraction* of a line or column. If it is intended to start a line at the middle of the first column the value used should be 1.5, which means that the start value is at the half (*0.5*) of the first column (*1*), observing that *1.5* is *0.5* plus *1*:



As in the second example above it is possible to end up in a specified *fraction* of a line or column:



The row, column *syntax*

Both lines and columns share the same syntax, where n is any integer between 1 and the number of rows and f is the fractional part of any number between 0 and 1:

(1) If only the row number n is provided the *content* is placed at the row n .

(2) If the row number n is provided followed by a dot and a number f, the *content* is placed at the fraction f of the row n.

(3) If the start row n_s and the end row n_e are provided separated by a dash, *i.e.*, n_s-n_e, the *content* is placed filling all the rows from n_s to n_e.
The dot notation described in **(2)** *can be used both on* n_s *and* n_e.

(4) All of the items above apply to columns in the same way.

✠ The cell contents

- ✓ **Z** – the atomic number of the elements.
- ✓ **name** – the name of the elements.
- ✓ **CS** – the chemical symbol of the elements.
- ✓ **Ar** – the relative atomic mass (atomic weight) of the elements.
- ✓ **Ar*** –the standard relative atomic mass (standard atomic weight) of the elements.

- ✓ **radio** – radioactivity of the elements. If the element is radioactive the figure ☢ is placed in the cell, otherwise nothing is shown.
- ✓ **R** – the atomic radius of the elements. The atomic radius shown is the calculated radius and is expressed in picometers.
- ✓ **Rcov** – the covalent radius of the elements. The covalent radius shown is for single bonds and is expressed in picometers.
- ✓ **Rion** – the ionic radius of the elements. The radius shown is the effective ionic radius in picometers.
- ✓ **Ei** – the first ionization energy of the elements, measured in $\mathrm{kJ \cdot mol^{-1}}$. All data from rutherfordium onwards is predicted.
- ✓ **eneg** – the Pauling electronegativity of the elements.
- ✓ **eaff** – the electroaffinity (electron affinity) of the elements, measured in $\mathrm{kJ \cdot mol^{-1}}$. Estimated negative values have been replaced by zero, since the negative ions formed in these cases are always unstable (they may have lifetimes of the order of microseconds to milliseconds, and invariably autodetach after some time).
- ✓ **O** – the common oxidation states of the elements.
- ✓ **Tmelt** – the melting point, in Kelvin, of the elements.
- ✓ **TmeltC** – the melting point, in degrees Celsius, of the elements.
- ✓ **Tboil** – the boiling point, in Kelvin, of the elements.
- ✓ **TboilC** – the boiling point, in degrees Celsius, of the elements.
- ✓ **eDist** – the electron distribution of the elements.
- ✓ **eConfign** – the electronic configuration, in increasing n (principal quantum number), of the element, corresponding to the *spectroscopic* order of orbital energies, that is, the reverse of the order in which electrons are removed from a given atom to form positive ions.

  ***Note**: the short version of the electronic configuration is used, i.e., [previous noble gas]remaining electrons. For example, for scandium it is:* [Ar]$3d^1 4s^2$
- ✓ **eConfignl** – the electronic configuration, in increasing sum of n and ℓ (azimuthal quantum number), of the element, following the order based on the Madelung rule.

  ***Note**: the short version of the electronic configuration is used, i.e., [previous noble gas]remaining electrons. For example, for scandium it is:* [Ar]$4s^2 3d^1$
- ✓ **d** – the density of the elements, in the corresponding physical state, at 25°C and 1 atm.
- ✓ **Cp** – the specific heat capacity of the elements in $\mathrm{J \cdot mol^{-1} \cdot K^{-1}}$ at 25°C and 100 kPa.
- ✓ **kT** – the thermal conductivity of the elements in $\mathrm{J \cdot m^{-1} \cdot K^{-1}}$ at 25°C.
- ✓ **ls** – the lattice structure of the elements at 1 bar and mostly at 25°C.
- ✓ **lsa** – the lattice constant a of the elements in picometers at 1 bar and mostly at 25°C.
- ✓ **lsb** – the lattice constant b of the eligible elements in picometers at 1 bar and mostly at 25°C.
- ✓ **lsc** – the lattice constant c of the eligible elements in picometers at 1 bar and mostly at 25°C.
- ✓ **lsca** – the lattice c/a ratio of the eligible elements at 1 bar and mostly at 25°C.
- ✓ **DiscY** – the discovery year of the elements.
- ✓ **DiscC** – the discovery country or in, a few cases, region (Middle East or Asia Minor) of the elements.
- ✓ **spectra** – the emission spectrum of the elements. The spectrum is only shown if available. The spectra are pre-built using the package pgf-spectra via the commands:

```
\pgfspectraStyle[back=visible40,line width=1pt,width=180pt,height=45pt,%
                 relative intensity,relative intensity threshold=.375,%
                 brightness=.5,charge=all,Imin=.125,gamma=1]
\foreach \SQ in {H,He,(...),Bi,Po,Rn,Fr,(...),Es}% Z=1,2,(...),83,84,86,87,(...),99
         {%
         \pgfspectra[element=\SQ]%
         }%
```

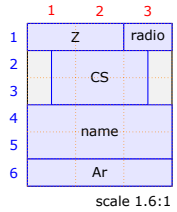
✠ Built-in cell styles

There is a set of *built-in* cell styles that could be used for the described purposes:

✓ **pgfPT2lang** – a cell layout to use with the name in two languages.

Built-in style **pgfPT2lang**
The build command:

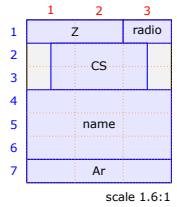
```
\pgfPTbuildcell(6,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4-5;1-3;name),(6;1-3;Ar)]
```



✓ **pgfPT3lang** – a cell layout to use with the name in three languages.

Built-in style **pgfPT3lang**
The build command:

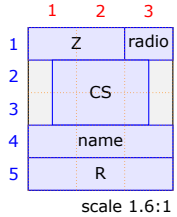
```
\pgfPTbuildcell(7,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4-6;1-3;name),(7;1-3;Ar)]
```



✓ **pgfPTR** – a cell layout to display the atomic radius and its periodic variations (if of course the show periodic variations key is set to true).

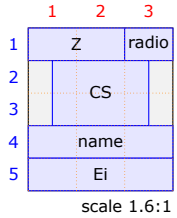
Built-in style **pgfPTR**
The build command:

```
\pgfPTbuildcell(5,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;R)]
```



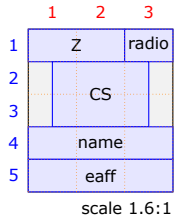
✓ **pgfPTEi** – a cell layout to display the first ionization energy and its periodic variations (if of course the show periodic variations key is set to true).

Built-in style **pgfPTEi**
The build command:
```
\pgfPTbuildcell(5,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;Ei)]
```



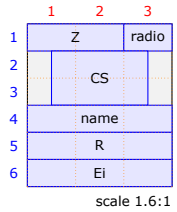
✓ **pgfPTeaff** – a cell layout to display the electron affinity and its periodic variations (if of course the show periodic variations key is set to true).

Built-in style **pgfPTeaff**
The build command:
```
\pgfPTbuildcell(5,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;eaff)]
```



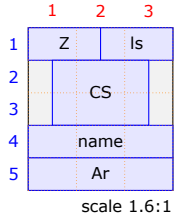
✓ **pgfPTREi** – a cell layout to display the atomic radius and first ionization energy and their periodic variations (if of course the show periodic variations key is set to true).

Built-in style **pgfPTREi**
The build command:
```
\pgfPTbuildcell(6,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;R),(6;1-3;Ei)]
```



✓ **pgfPTls** – a cell layout to display the lattice system.

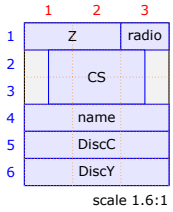
Built-in style **pgfPTls**
The build command:
```
\pgfPTbuildcell(5,3)%
[(1;1-2.5;Z),(1;2.5-3;ls),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;Ar)]
```



✓ **pgfPTdisc** – a cell layout to display the discovery country and discovery year.

Built-in style **pgfPTdisc**
The build command:

```
\pgfPTbuildcell(6,3)%
[(1;1-2;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4;1-3;name),(5;1-3;DiscC),(6;1-3;DiscY)]
```



Designing color schemes

There are three ways to make a new color scheme:

- with the command \pgfPTnewColorScheme
- using the *script* in the file pgfPTcolorSchemes.html
- with the commands provided by the colorschemes library (see the libraries section).

✠ Designing a color scheme with \pgfPTnewColorScheme

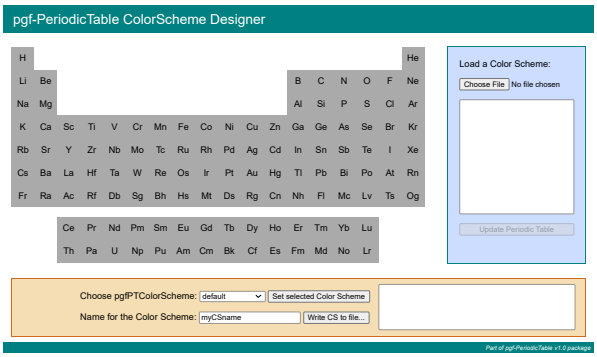
This command provides a way to set the cell background color of each of the 118 elements of the Periodic Table. *If the intention is to set the background color for all of them, it is highly recommended to use the file pgfPTcolorSchemes.html*, unless the trailing color begin at a small atomic number.

Despite that, this command can always be used taking into account:

1. It has the form \pgfPTnewColorScheme[trailing color]{name}{color list} where:
   - the first argument (enclosed by square brackets) is optional. If provided, the specified trailing color will be used, otherwise the default color (white) will be used as trailing color.
   - the second and third arguments are mandatory and specify, respectively, the color scheme name and the color list.
2. The name is any name made up of letters (only the characters a,…,z and A,…,Z).
3. The color list is a comma-separated list where each entry has the format r/g/b, representing the red, blue and green values, between 0 and 1, of the color: the first entry of the list will be the background color used in the cell of the element with atomic number 1, the second entry, the background color of the cell of the element with atomic number 2, and so on.

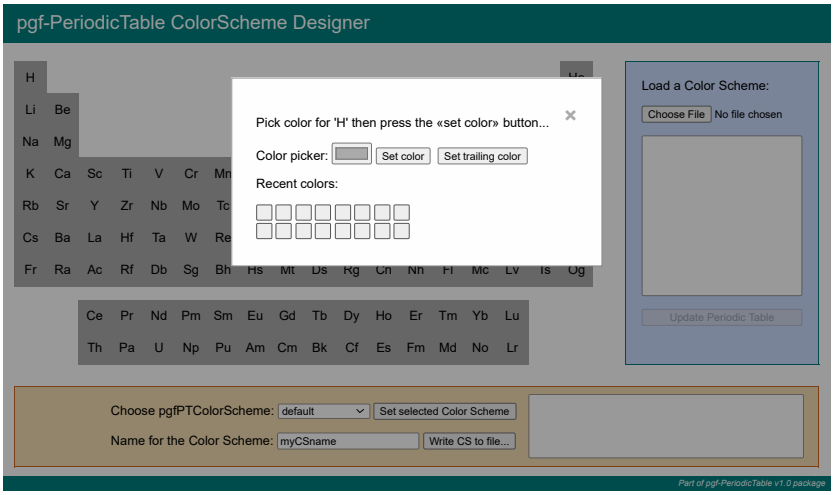
   If the color list has ten entries, these entries will set the background colors of the elements with atomic numbers from 1 to 10. For the following atomic numbers, greater than or equal to 11, the trailing color will be used in the color background.

✠ Designing a color scheme with pgfPTcolorSchemes.html

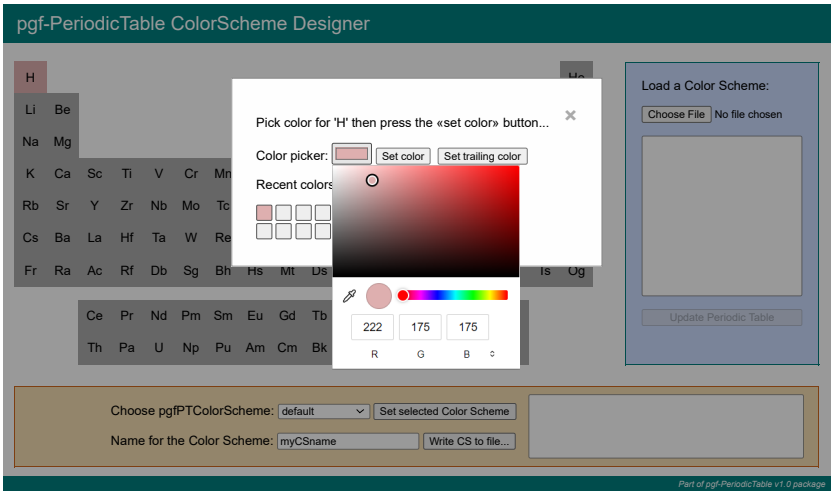


The pgfPTcolorSchemes.html *designer* is an *html* file with a little *javascript* code to perform the task of building a color scheme to use with the back color scheme key associated with the \pgfPT command.

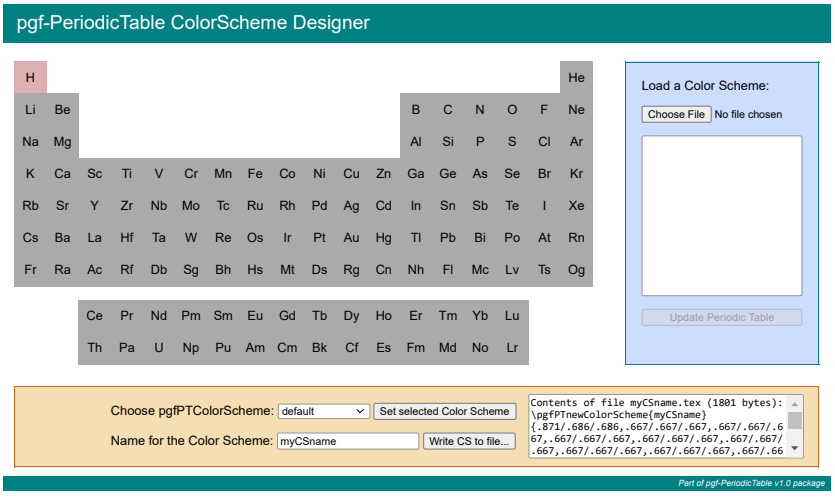
The Periodic Table of the Elements is displayed on the page and clicking on an element opens a color dialog:



Clicking on the Color picker: button opens a color dialog, where there is the possibility to choose the desired color or manually enter one color using one of the three models available (RGB, HSL or HEX):



After changing the desired colors it is possible to save the color scheme in a file by clicking on Write CS to file...:

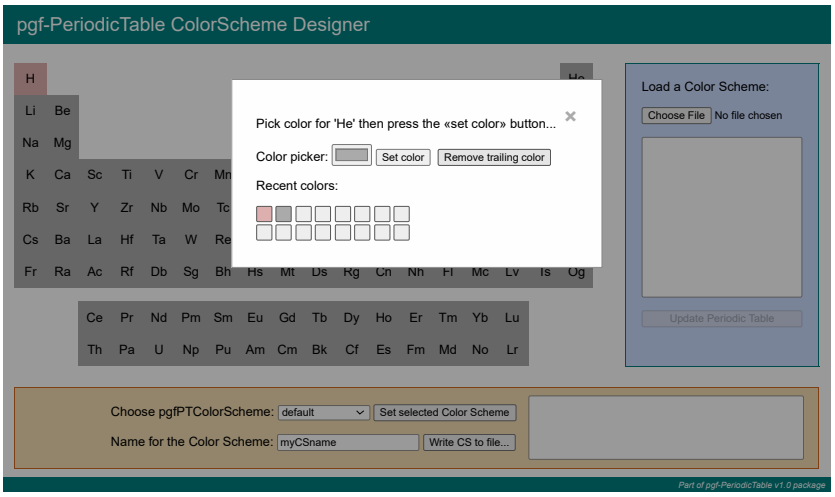


To use a color scheme saved in a file there are two possible ways:

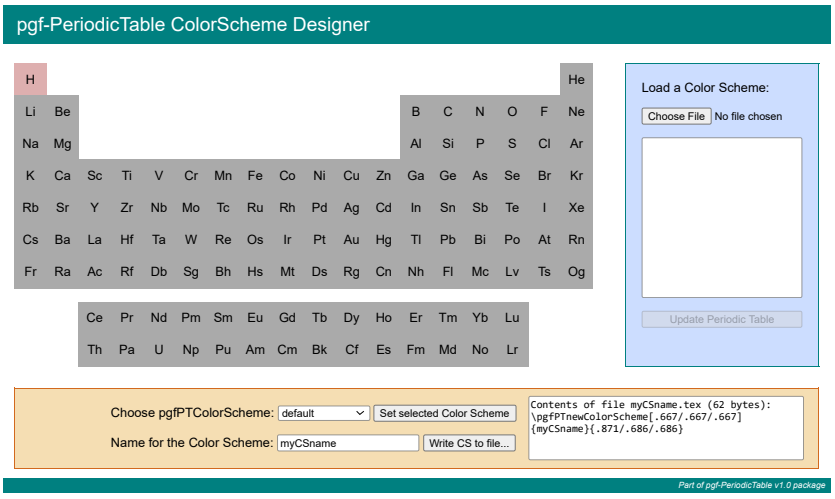
- loading the file in the working document via the \input LaTeX command, for instance, \input{myCSname.tex}.
- or by opening the file and copying and pasting its contents into the working document.

In either case, the operation can be performed at any location in the document, but before the named color scheme is used.

Note that in the previous example there is only one color that has been defined (for hydrogen). In that case, it is useful to set the trailing color in helium by clicking in Set trailing color (which automatically changes to Remove trailing color). After that only the hydrogen and helium are clickable, all the other elements are locked to click:

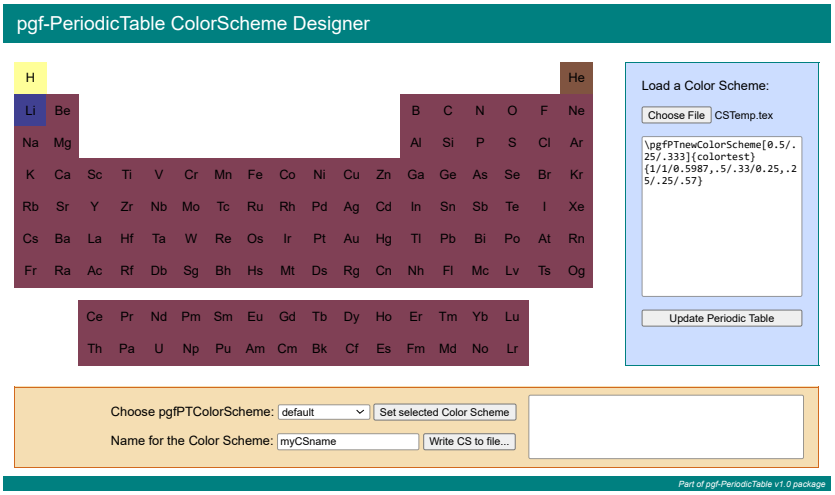


Then the saved color scheme will have the optional trailing color and the size will be smaller as only the color codes of the changed elements are stored:

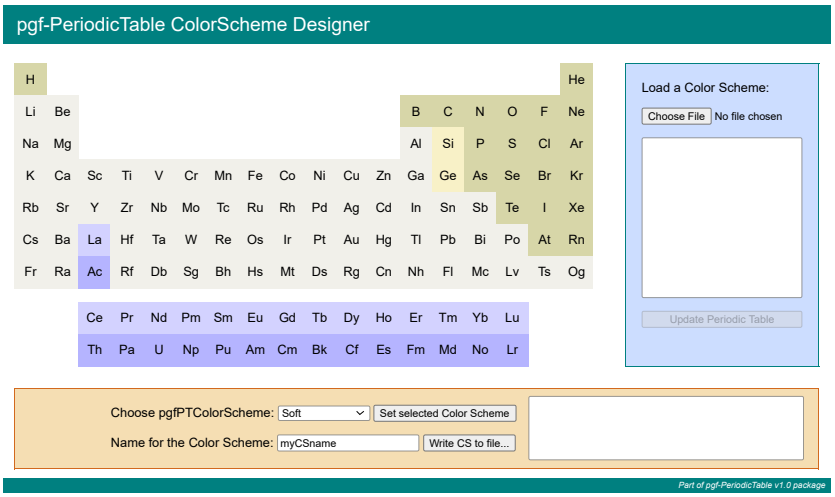


To remove the trailing color click on the last enabled element (in the above case helium) and then click on Remove trailing color. After that, all elements can be clicked again.

It is also possible to load a color scheme saved to a file by clicking on Choose File and then clicking on Update Periodic Table for the color scheme to take effect:



Finally its possible to load a built-in color scheme by choosing a named *pgfPTColorScheme* in the corresponding combo box and then clicking on Set selected Color Scheme:



All the operations described are always available.

Libraries

In this part the library packages are documented. They provide additional commands to extend the capabilities provided by this package out of the box. The libraries are not loaded by default since many users will not need them.

▶ Color Schemes Library

pgf-PeriodicTable Library colorschemes

USAGE: \usepgfPTlibrary{colorschemes}

This library extends the features provided by the command \pgfPTnewColorScheme. It defines a set of commands that automatically generate a new color scheme.

- \pgfPTGroupColors{name of the new color scheme}{list of colors,options}
- \pgfPTPeriodColors{name of the new color scheme}{list of colors,options}
- \pgfPTCScombine[proportion,mode]{name of the first color scheme,name of the second color scheme,name of the new color scheme}
- \pgfPTCSwrite[filename]{list of color schemes names}

Color arguments for this library's commands can use both the base package syntax – namedColor or namedColorA!##!namedColorB<!##><!named...> – or any color model supported by the `xcolor` package[a] using the *special syntax* *[model:values], *e.g.*, *[rgb:.5;.2;.3] or *[cmyk:.5;.2;.3;.3] or *[HTML:5FA287]. **The values for the individual color components of a color specified this way must be separated by semicolons instead of commas**, except for the HTML, Gray and wave color models as explained in the `xcolor` package.

[a] See *Table 3: Supported color models* on page 10 of the documentation of xcolor v2.14 2022/06/12

\pgfPTGroupColors[default group color]{name of the new color scheme}{list of colors,options}

This command **creates a Color Scheme** with the name name of the new color scheme. **Group colors** can be configured in three different ways:

- **setting the colors one by one**, using the *key=value* mechanism in the list of colors. For example:

```
\pgfPTGroupColors{name of the new color scheme}%
{G1=red,G2=red!50,G3=orange,<...>,G18=blue,options}
```

This will set the specified color for each group. If no color is specified for a group, default group color will be used.

NOTE: default group color is initially set to white.

- **defining a gradient** using the keys left color=<color>, middle color=<color> and right color=<color> as the list of colors. Note that all the keys are optional, but at least one of them is required. This produces a gradient starting from group 1, with *left color*, to group 18, with *right color*. If the *middle color* key is used then the gradient starts at group 1 with *left color*, goes to the middle position of the groups (between groups 9 and 10) with *middle color* and ends at group 18 with *right color*. For example:

```
\pgfPTGroupColors{name of the new color scheme}%
{left color=red,right color=blue,options}
```

defines a gradient from red (group 1) to blue (group 18).

- **defining a custom gradient** as the list of colors by using the *key=value* mechanism inside the gradient key. For example:

```
\pgfPTGroupColors{name of the new color scheme}%
{gradient={G1=red,G4=red!50,G18=blue},options}
```

defines a gradient from red (group 1) to red!50 (group 4) and to blue (group 18).

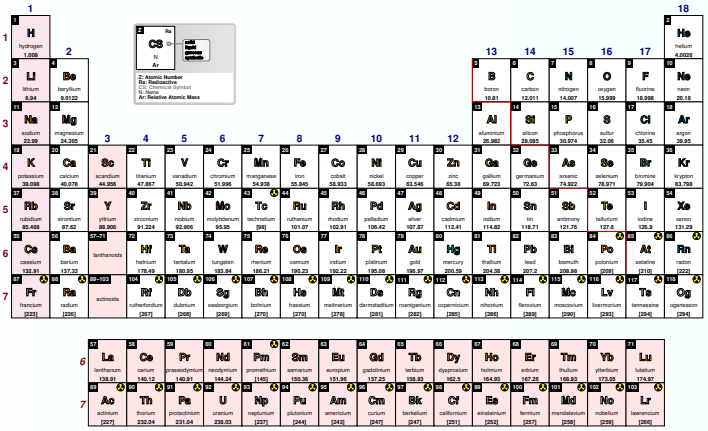
The options available to this command are:

- H=<color>, sets the color of the *hydrogen* cell. If not set, group 1's color will be used. If set, the color of the *hydrogen* cell won't be affected by period blending.
- La=<color>, sets the color of the *lanthanum* cell. If not set, group 3's color will be used.
- Lanta=<color>, sets the color of the *lanthanoids* cells. If not set, *lanthanum*'s color will be used.
- Ac=<color>, sets the color of the *actinium* cell. If not set, group 3's color will be used.
- Actin=<color>, sets the color of the *actinoids* cells. If not set, *actinium*'s color will be used.
- period blending={color=<color>, percentage=<positive or negative integer>, mode=<add|sub|linear>}, performs a *mode* blend over the periods up to the specified percentage with the provided color.

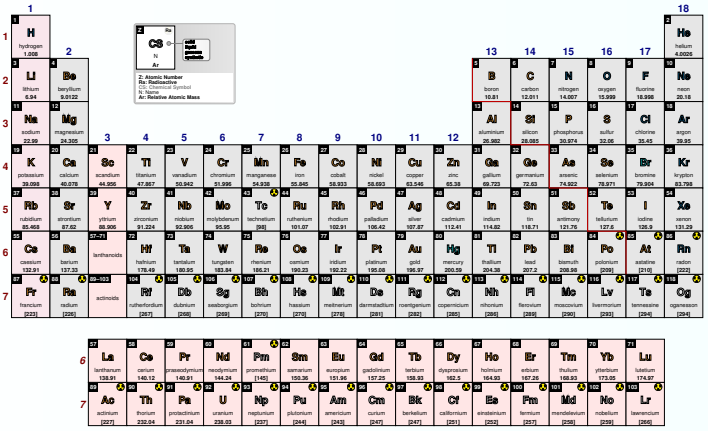
  NOTES:

  - percentage refers to how much of the color, in total, was mixed over the 7 periods. For example 60% adds 10% to each period: P1►0% ⇝ P2►10% ⇝ P3►20% ⇝ ... ⇝ P7►60%. If the percentage is positive, the mixing is done in descending order (from P1 to P7); if the percentage is negative, the mixing is done in ascending order (from P7 to P1).
  - The mode's values are add for *additive* blending, sub for *subtractive* blending and linear for *linear* blending (as in the `xcolor` package).
  - **If period blending is used without further options** all the default values are used, so period blending is equivalent to period blending={color=white,percentage=60,mode=linear}.
  - None of the keys color, percentage and mode are mandatory. If omitted the default value is used.

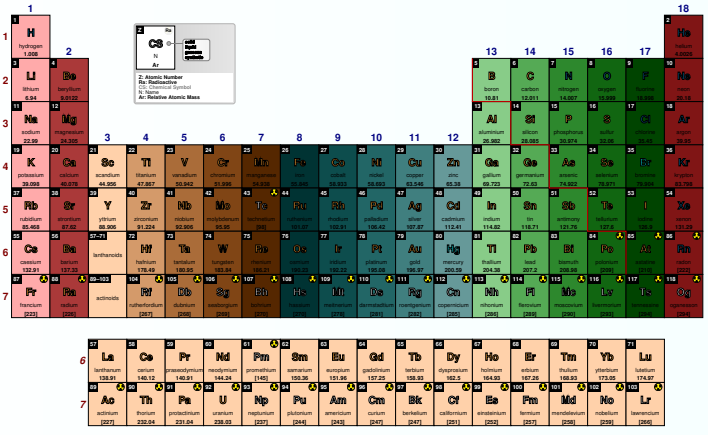
```
\pgfPTGroupColors{example}{G1=purple!10,G3=red!10}
\pgfPT[back color scheme=example,show title=false]
```



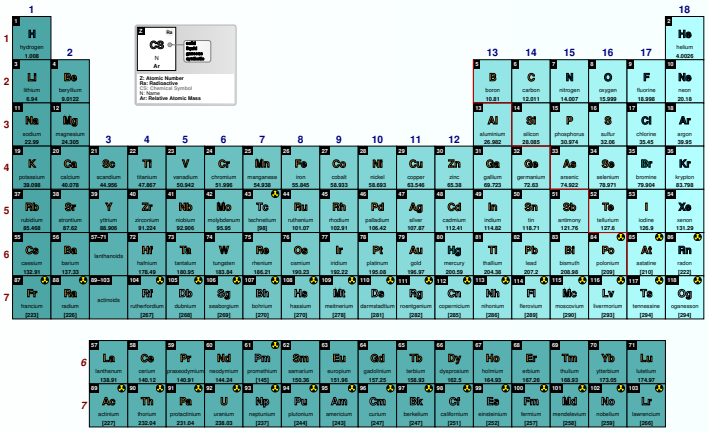
```
\pgfPTGroupColors[black!10]{example}{G1=purple!10,G3=red!10}
\pgfPT[back color scheme=example,show title=false]
```



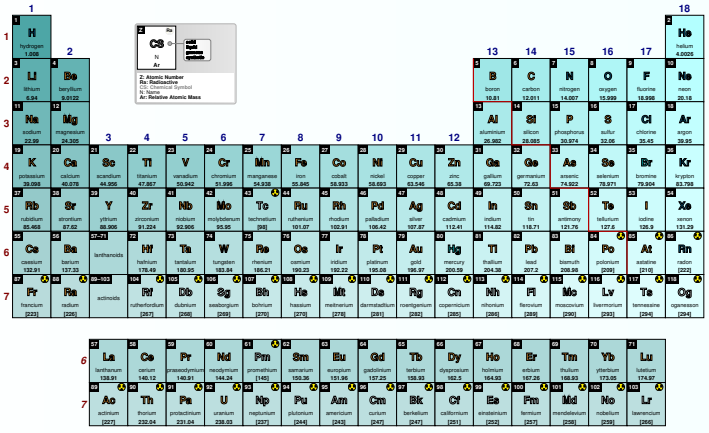
```
\pgfPTGroupColors{example}{G1=*[HTML:FFAAAA],G2=*[HTML:AA3939],
G3=*[HTML:FFD1AA],G4=*[HTML:D49A6A],G5=*[HTML:AA6C39],
G6=*[HTML:804515],G7=*[HTML:552700],G8=*[HTML:003333],
G9=*[HTML:0D4D4D],G10=*[HTML:226666],G11=*[HTML:407F7F],
G12=*[HTML:669999],G13=*[HTML:88CC88],G14=*[HTML:55AA55],
G15=*[HTML:2D882D],G16=*[HTML:116611],G17=*[HTML:004400],
G18=*[HTML:801515] }
\pgfPT[back color scheme=example,show title=false]
```



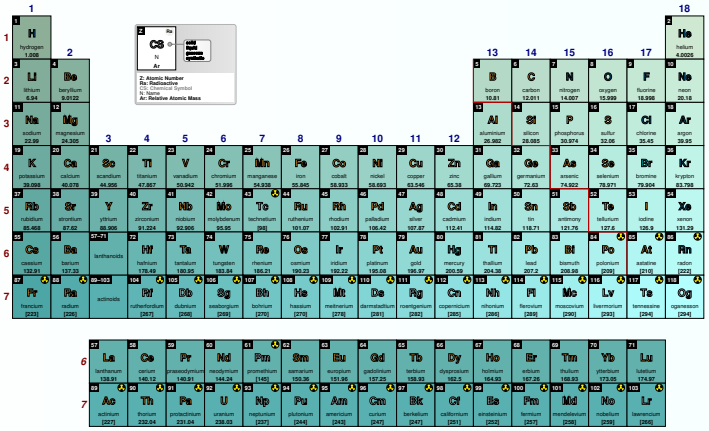
```
\pgfPTGroupColors{example}{left color=teal!70,right color=cyan!30}
\pgfPT[back color scheme=example,show title=false]
```



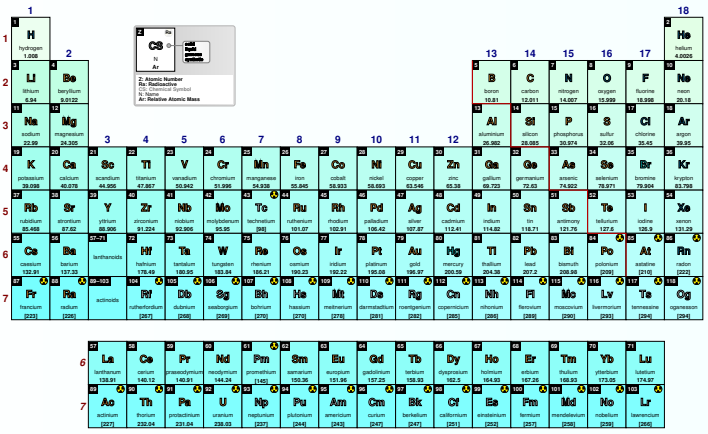
```
\pgfPTGroupColors{example}{left color=teal!70,right color=cyan!30,period blending}
\pgfPT[back color scheme=example,show title=false]
```



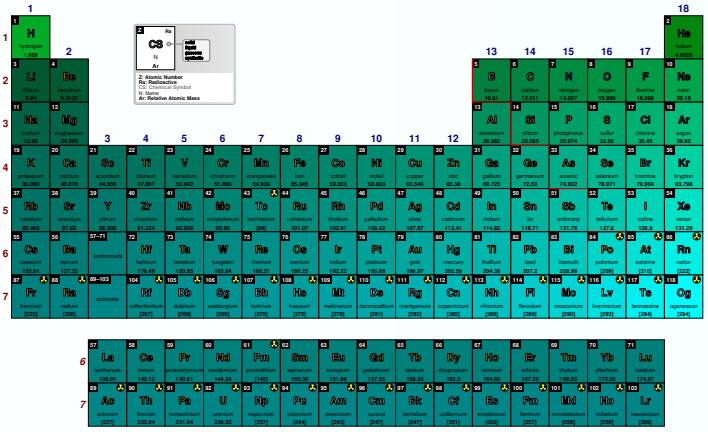
```
\pgfPTGroupColors{example}{left color=teal!70,right color=cyan!30,
period blending={color=orange!50,percentage=-40}}
\pgfPT[back color scheme=example,show title=false]
```



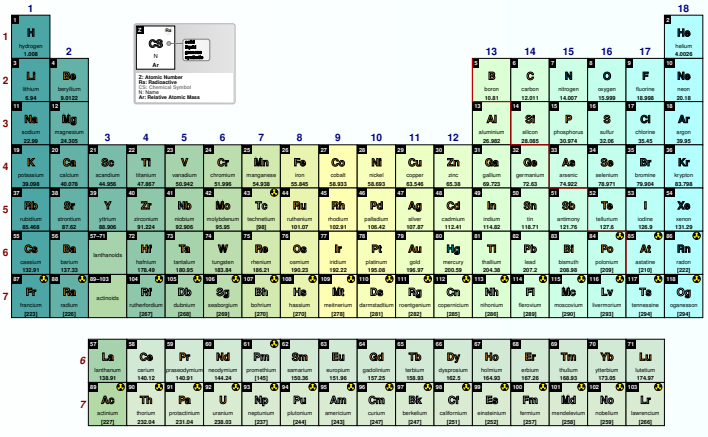
```
\pgfPTGroupColors{example}{left color=teal!70,right color=cyan!30,
period blending={color=orange!50,percentage=-40,mode=add},
H={*[cmyk:.071,0,.055,.035]}}
\pgfPT[back color scheme=example,show title=false]
```



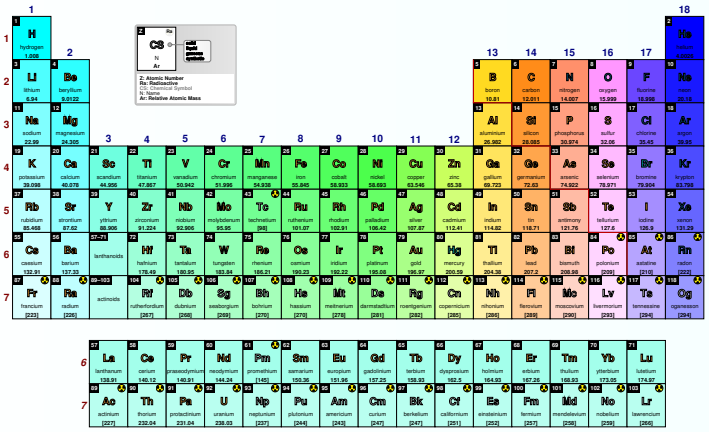
```
\pgfPTGroupColors{example}{left color=teal!70,right color=cyan!30,
period blending={color=orange!50,percentage=-40,mode=sub},
H=*[cmyk:.071;0;.055;.035]}
\pgfPT[back color scheme=example,show title=false]
```



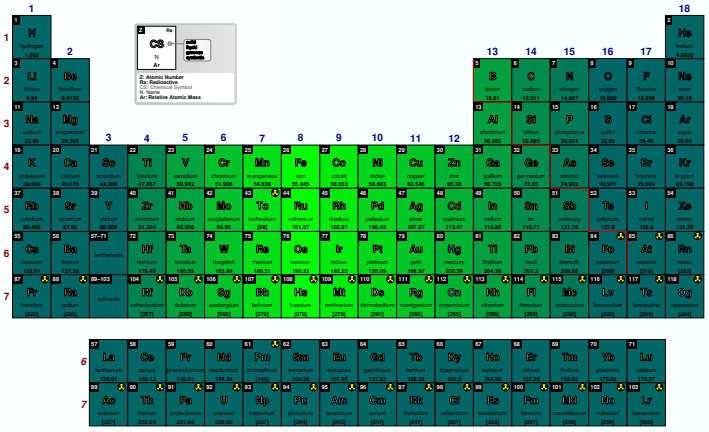
```
\pgfPTGroupColors{example}{left color=teal!70,middle color=yellow!30,right
color=cyan!30,La=teal!70!yellow!50,Ac=teal!60!yellow!50,
Lanta=teal!70!yellow!50!white!50,Actin=teal!60!yellow!50!white!50}
\pgfPT[back color scheme=example,show title=false]
```



```
\pgfPTGroupColors{example}{gradient={G1=teal!50!black,G2=teal,G10=green,
G14=orange,G18=blue},period blending={mode=add}}
\pgfPT[back color scheme=example,show title=false]
```



```
\pgfPTGroupColors{example}{gradient={G3=teal!80!black,G16=teal!80!black,
G8=green}}
\pgfPT[back color scheme=example,show title=false]
```



Note: the group numbers can be specified in any order and the gradient can start or end in any group. *In this example, the smallest group number is 3 and the greatest is 16, so the gradient is built from group 3 to group 16 and the colors from group 1 to 3 are equal to group 3's color, just like the colors from group 16 to 18 are equal to group 16's color.*

\pgfPTPeriodColors[default period color]{name of the new color scheme}{list of colors,options}

This command **creates a Color Scheme** with the name name of the new color scheme. **Period colors** can be configured in three different ways:

✓ **setting the colors one by one**, using the *key=value* mechanism in the list of colors. For example:

```
\pgfPTPeriodColors{name of the new color scheme}%
{P1=red,P2=red!50,<...>,P7=blue,options}
```

This will set the specified color for each period. If no color is specified for a period,

default period color will be used.

NOTE: default period color is initially set to white.

✓ **defining a gradient** using the keys top color=<color>, middle color=<color> and bottom color=<color> as the list of colors. Note that all the keys are optional, but at least one of them is required. This produces a gradient starting from period 1, with *top color*, to period 7, with *bottom color*. If the *middle color* key is used then the gradient starts at period 1 with *top color*, goes to the middle position of the periods (period 4) with *middle color* and ends at period 7 with *bottom color*. For example:

```
\pgfPTPeriodColors{name of the new color scheme}%
{top color=red,middle color=yellow,bottom color=blue,options}
```

defines a gradient from red (period 1) to yellow (period 4) and from yellow (period 4) to blue (period 7).

✓ **defining a custom gradient** as the list of colors by using the *key=value* mechanism inside the gradient key. For example:

```
\pgfPTPeriodColors{name of the new color scheme}%
{gradient={P1=red,P3=red!50,P7=blue},options}
```

defines a gradient from red (period 1) to red!50 (period 3) and to blue (period 7).

The options available to this command are:

✓ H=<color>, sets the color of the *hydrogen* cell. If not set, period 1's color will be used. If set, the color of the *hydrogen* cell won't be affected by group blending.

✓ La=<color>, sets the color of the *lanthanum* cell. If not set, period 6's color will be used.

✓ Lanta=<color>, sets the color of the *lanthanoids* cells. If not set, *lanthanum*'s color will be used.

✓ Ac=<color>, sets the color of the *actinium* cell. If not set, period 7's color will be used.

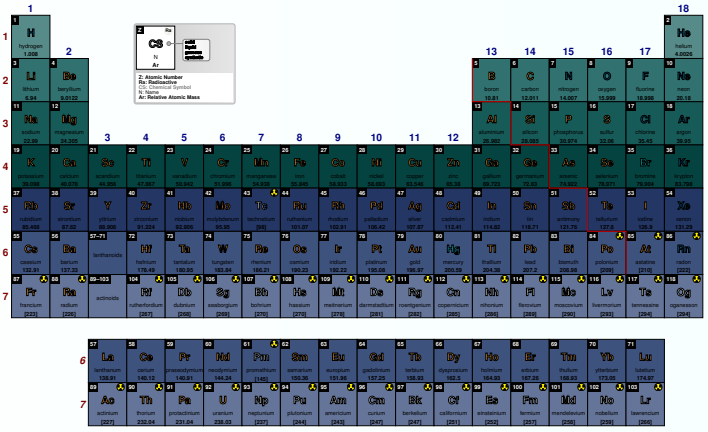
✓ Actin=<color>, sets the color of the *actinoids* cells. If not set, *actinium*'s color will be used.

✓ group blending={color=<color>, percentage=<positive or negative integer>, mode=<add|sub|linear>}, performs a *mode* blend over the groups up to the specified percentage with the provided color.

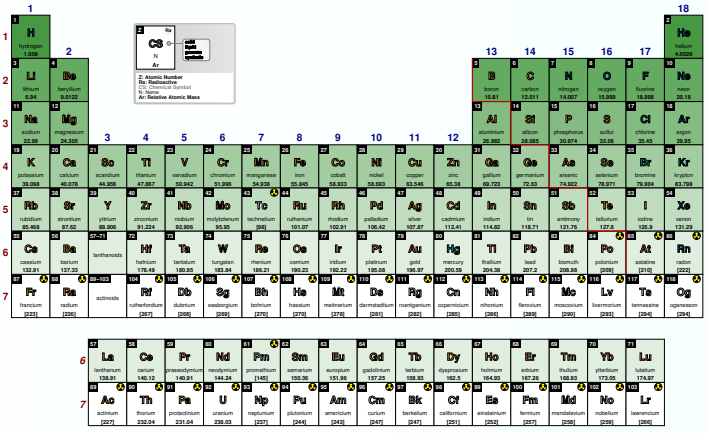
NOTES:

- ✓ percentage refers to how much of the color, in total, was mixed over the 18 groups. For example 68% adds 4% to each period: G1▶0% ⇝ G2▶4% ⇝ G3▶8% ⇝ … ⇝ G18▶68%. If the percentage is positive, the mixing is done from left to right (from G1 to G18); if the percentage is negative, the mixing is done from right to left (from G18 to G1).
- ✓ The mode's values are add for *additive* blending, sub for *subtractive* blending and linear for *linear* blending (as in the `xcolor` package).
- ✓ **If group blending is used without further options** all the default values are used, so group blending is equivalent to group blending={color=white,percentage=68,mode=linear}.
- ✓ None of the keys color, percentage and mode are mandatory. If omitted the default value is used.

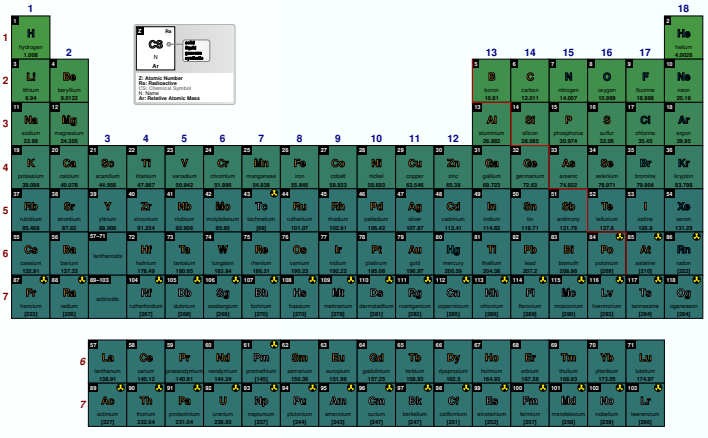
```
\pgfPTPeriodColors{example}{P1=*[RGB:86;139;137],P2=*[RGB:49;114;112],
P3=*[RGB:23;91;88],P4=*[RGB:5;67;64],P5=*[RGB:35;54;100],
P6=*[RGB:62;82;126],P7=*[RGB:101;117;153]}
\pgfPT[back color scheme=example,show title=false]
```



```
\pgfPTPeriodColors{example}{top color=*[Hsb:117;.57;.6]}
\pgfPT[back color scheme=example,show title=false]
```



```
\pgfPTPeriodColors{example}{gradient={P1=*[Hsb:117;.57;.6],
P5=*[Hsb:178;.57;.45]}}
\pgfPT[back color scheme=example,show title=false]
```



\pgfPTCScombine[prop1:prop2,mode]{name of color scheme one,name of color scheme two,name of the new color scheme}

This command **combines two named Color Schemes** and merges the result in a new Color Scheme with name of the new color scheme.
For example \pgfPTCScombine{myCSA,myCSB,myCSC} adds the color scheme myCSA to the color scheme myCSB and their sum will be available as the color scheme myCSC.

NOTE: if the Color Schemes have different sizes (*i.e.*, different number of colors), the last color from the color scheme that ends first will be used until the other color scheme also ends.

The optional parameters [prop1:prop2,mode] are for controlling how the two Color Schemes are combined:

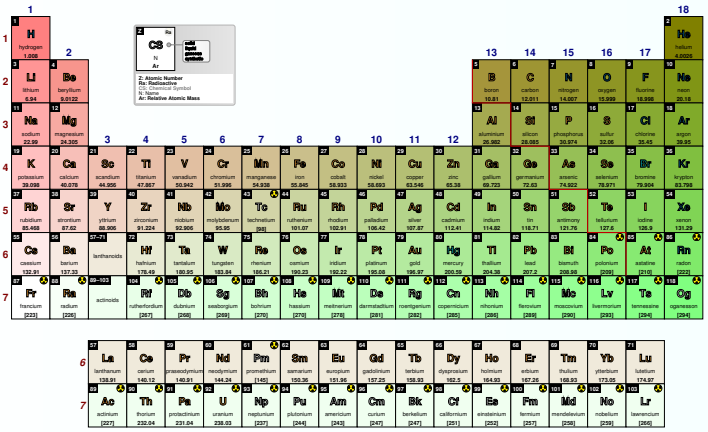
- ✓ The first parameter – prop1:prop2 – controls the proportions used to mix the color schemes: prop1 parts of name of color scheme one and prop2 parts of name of color scheme two. Both prop1 and prop2 must be integer values between 1 and 999.

  NOTE: default proportion is 1:1.
  For example, 1:4 *will mix each color in the ratio of 1 to 4, i.e., the nth-color from the first color scheme is used as 1/5 of the mixed color and the nth-color from the second color scheme is used as 4/5 of the mixed color*.

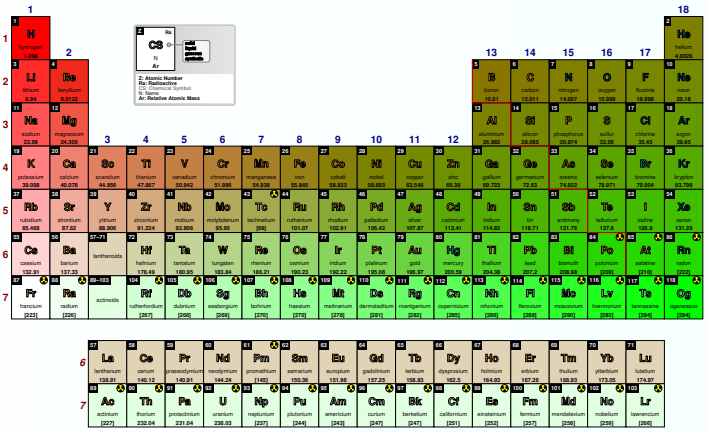
- ✓ The mode refers to how the colors are mixed: use add for *additive* mixing, sub for *subtractive* mixing and linear for *linear* mixing (as in the `xcolor` package).

  NOTE: default mode is linear.

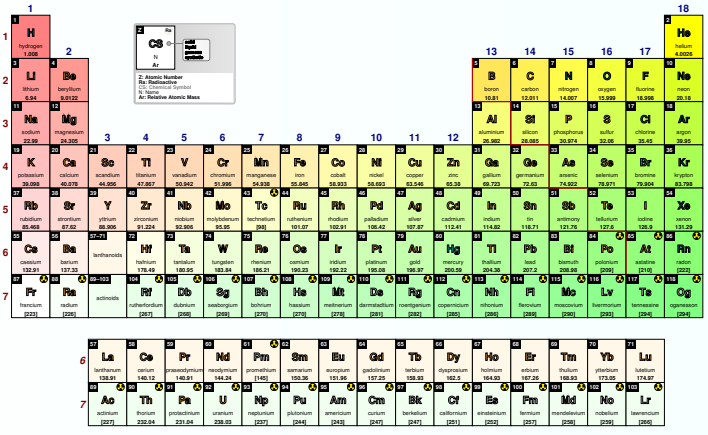
```
\pgfPTPeriodColors{period}{top color=red}
\pgfPTGroupColors{group}{right color=green}
\pgfPTCScombine{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```



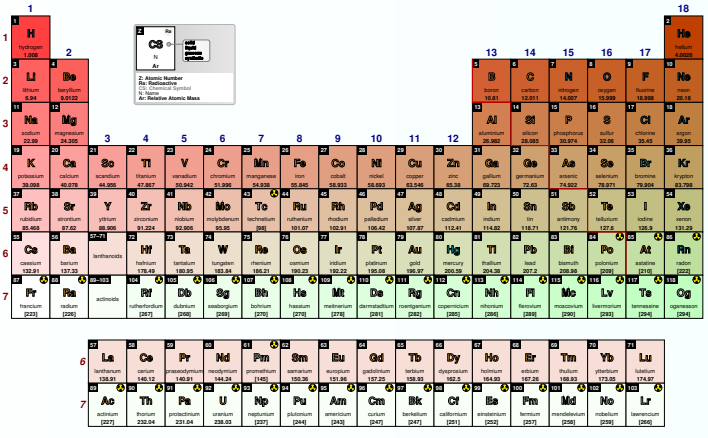
```
\pgfPTCScombine[sub]{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```



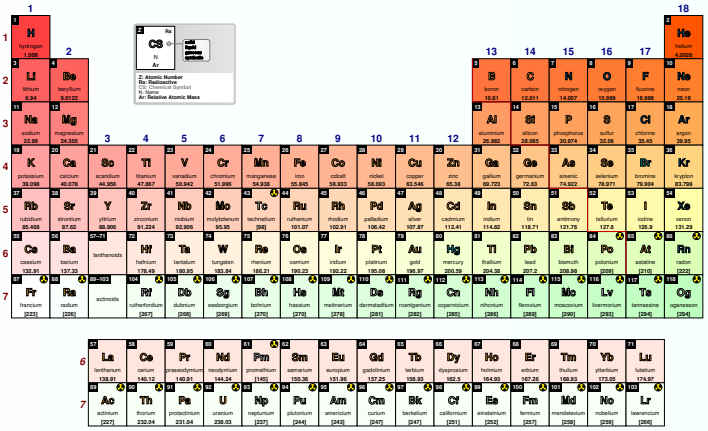
```
\pgfPTCScombine[add]{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```



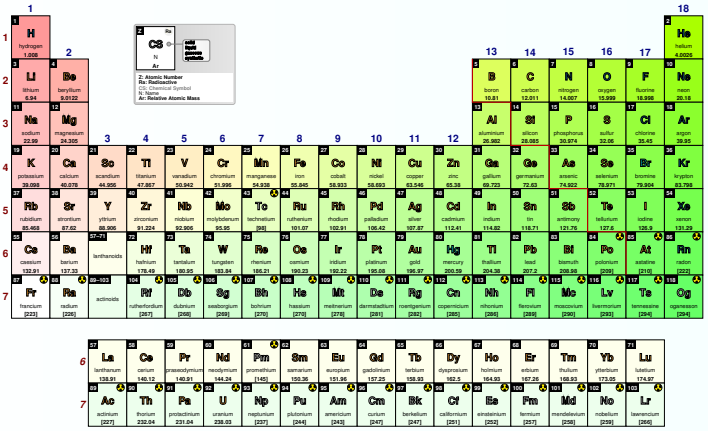
```
\pgfPTCScombine[3:1]{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```



```
\pgfPTCScombine[3:1,add]{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```

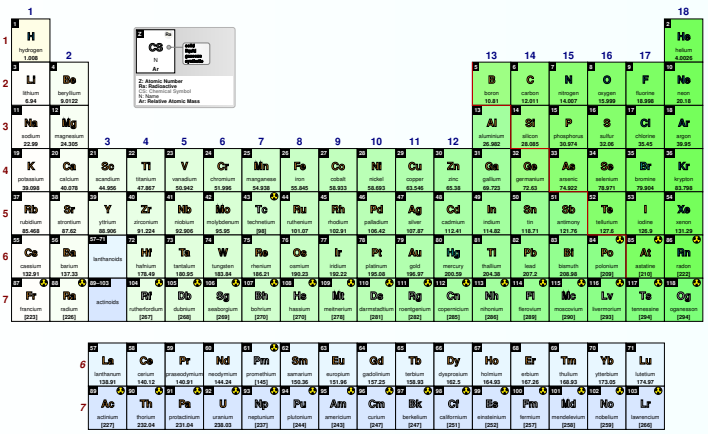


```
\pgfPTCScombine[add,2:3]{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```

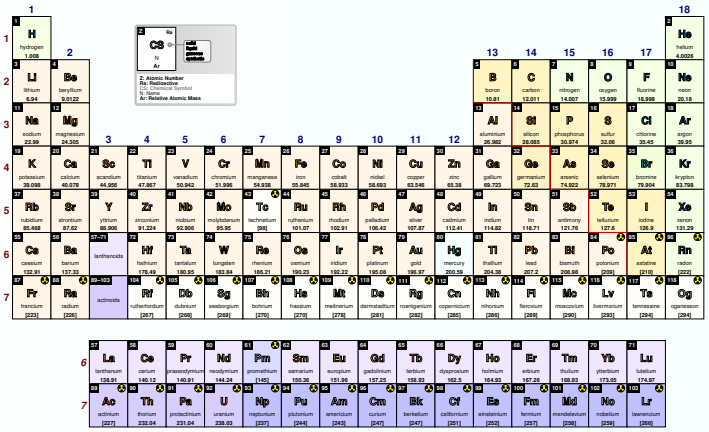


Built-in color schemes can also be mixed:

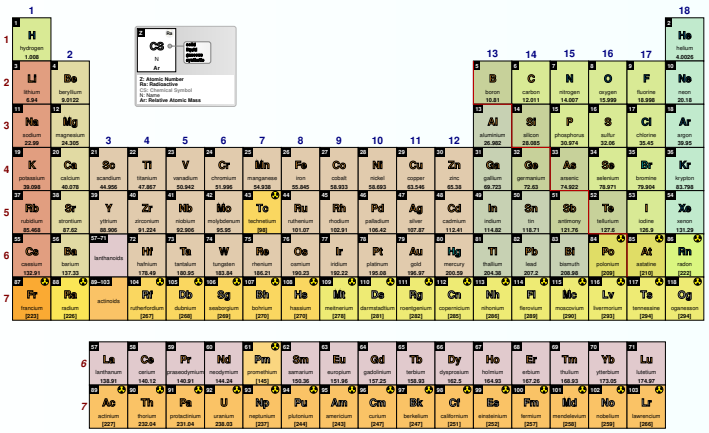
```
\pgfPTCScombine[add]{Soft,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```



```
\pgfPTCScombine[add,3:1]{Soft,PS,mix}
\pgfPT[back color scheme=mix,show title=false]
```



```
\pgfPTCScombine[add]{Radio,Wikipedia,mix}
\pgfPT[back color scheme=mix,show title=false]
```



```
\pgfPTCSwrite[filename]{list of color schemes names}
```

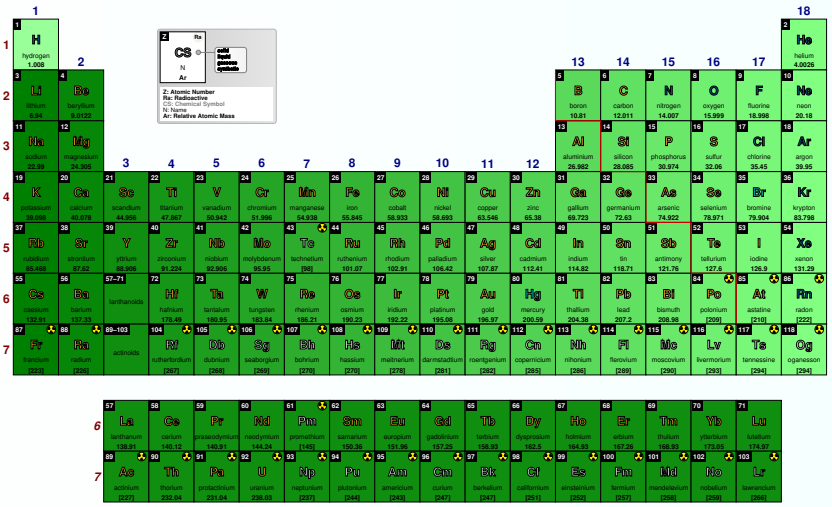
This command **writes the provided Color Schemes to a file** for later use without loading this library. It has a mandatory argument, the list of the color schemes names to be written and an optional argument, the filename. If no filename is provided the first name on the list of the color schemes names is used.
For example, `\pgfPTCSwrite[myGroupColors]{myGroupGradGreenToRed,myGroupGreens, myGroupGradYellowToRed}`, *will create (or overwrite), in the current working directory, a file with name* `myGroupColors.tex` *with the follwing contents*:

```
\pgfPTnewColorScheme{myGroupGradGreenToRed}{0/1/0,...
\pgfPTnewColorScheme{myGroupGreens}{0/1/.1,...
\pgfPTnewColorScheme{myGroupGradYellowToRed}{1/1/0,...
```

After that, it's possible to use `\input{myGroupColors.tex}`, anywhere in any document (in the same working directory). The named color schemes defined in the loaded file are now available for use as usual:

```
\pgfPTPeriodColors{myGroupGradGreenToRed}{gradient={G1=green!50!black,
G18=red!30!black},H=green!40!white}
\pgfPTPeriodColors{myGroupGreens}{gradient={G1=green!50!black,
G18=green!50!white},H=green!40!white}
\pgfPTPeriodColors{myGroupGradYellowToRed}{gradient={G1=yellow!50!white,
G18=red!30!black},H=yellow!40!white}
\pgfPTCSwrite[myGroupColors]{myGroupGradGreenToRed,myGroupGreens, myGroup-
GradYellowToRed}
```

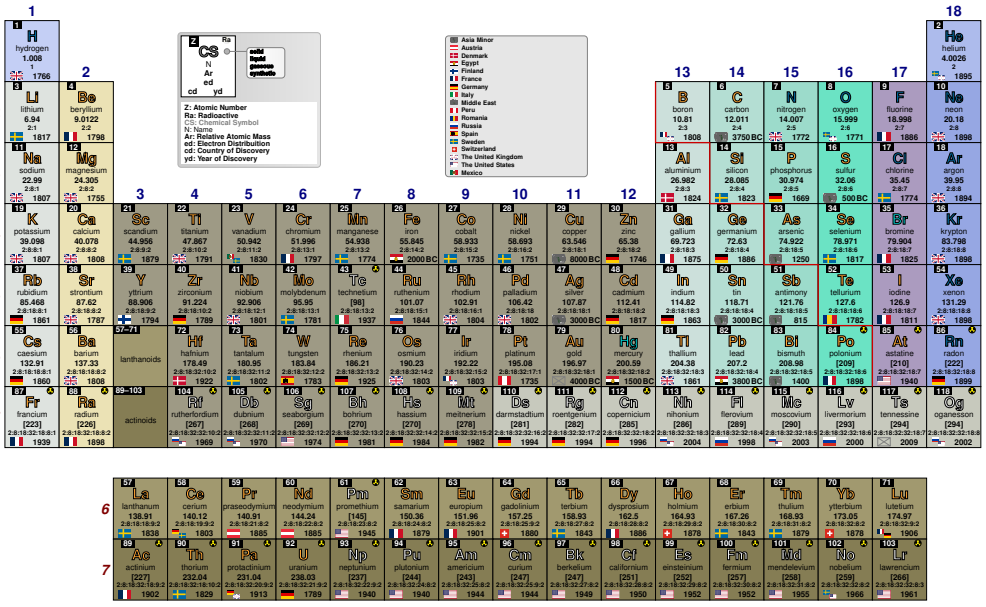
```
%\usepgfPTlibrary{colorschemes}
\input{myGroupColors.tex}%
\pgfPT[back color scheme=myGroupGreens,show title=false]
```



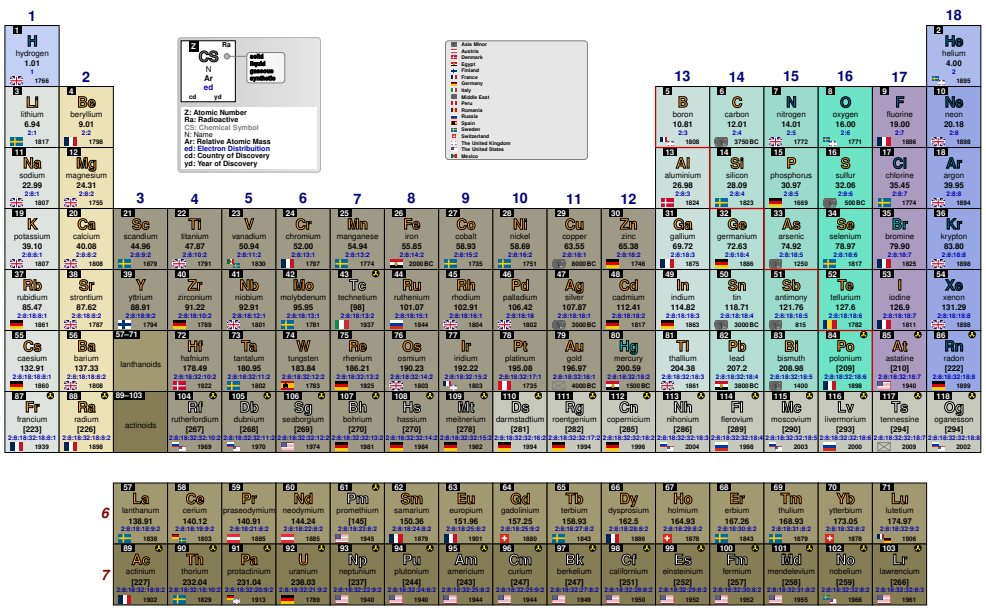
A few more examples

The following examples could be used for students or for any other purposes.

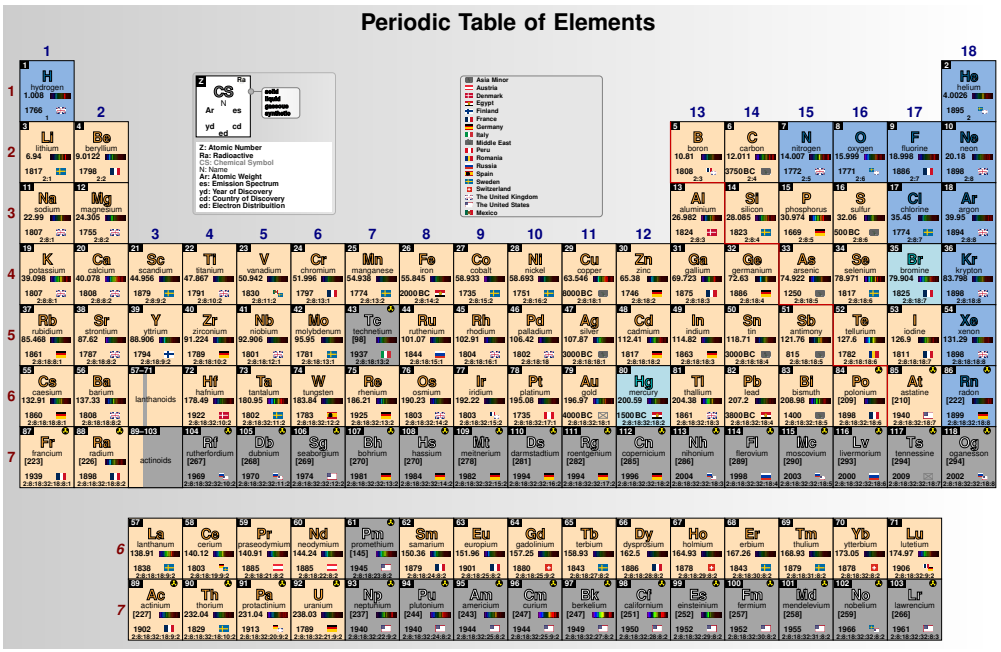
```
\pgfPTbuildcell(8,3)% 8 rows by 3 columns
[(1;1.4-2.8;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4.2;1-3;name), (5.4;1-3;Ar),(6.5;1-3;eDist),(7.55-8.95;1-2.25;DiscC),(7.55-8.95;2.25-3.8;DiscY)]
\pgfPT
```



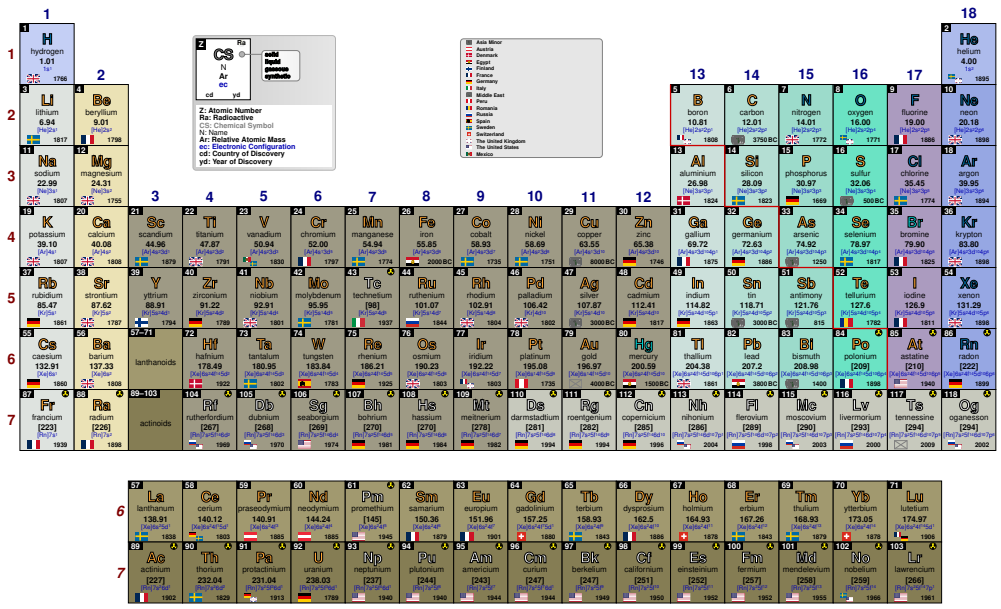

```
\pgfPT[eDist color=blue!70!black,Ar precision=2,DiscC font=\fontsize{4}{4}\selectfont,DiscY font=\fontsize{4}{4}\selectfont\bfseries]
```



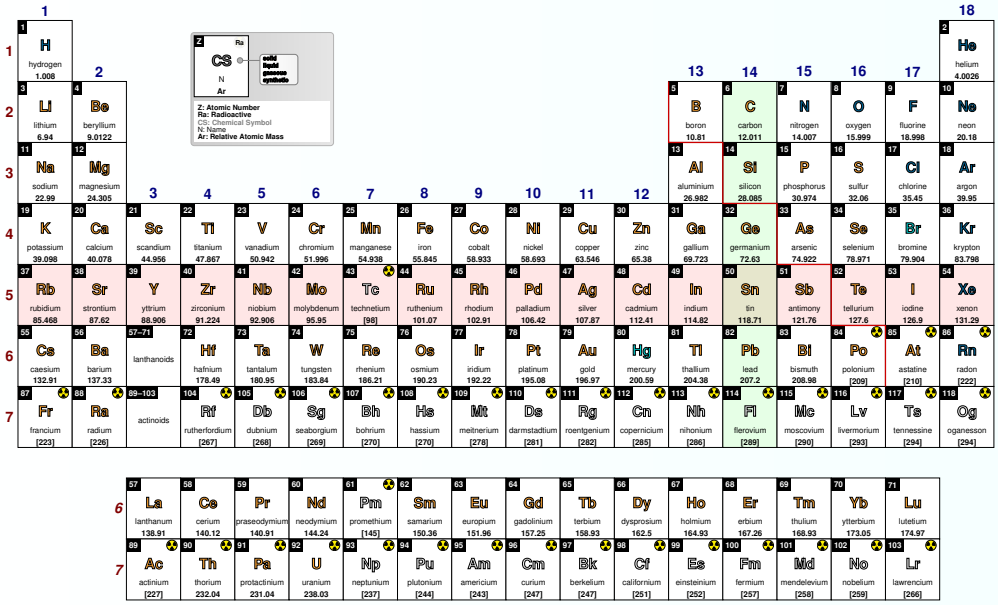
```
\pgfPTbuildcell(8,3)% 8 rows by 3 columns
[(1;1-2;Z),(1;3;radio),(2-3;1-3;CS),(4;1-3;name),(5;1-2.5;Ar),(5;2.5-3;spectra),
(7;1-2.5;DiscY),(7;2.5-3;DiscC),(8;1-3;eDist)]
        \pgfPT[csPS,Ar label=w,background={left color=black!20}]
```




```
\pgfPTbuildcell(8,3)% 8 rows by 3 columns
[(1;1-3;Z),(1;3;radio),(2-3;1.5-3.5;CS),(4.2;1-3;name),(5.4;1-3;Ar), (6.5;1-
3;eConfignl),(7.55-8.95;1-2.45;DiscC),(7.55-8.95;2.45-3;DiscY)]
        \pgfPT[eConfignl color=blue!70!black,Ar precision=2,DiscC
    font=\fontsize{4}{4}\selectfont,DiscY font=\fontsize{4}{4}\selectfont\bfseries]
```



```
\usepgfPTlibrary{colorschemes}
\pgfPTPeriodColors{period}{P5=red!20}
\pgfPTGroupColors{group}{G14=green!20}
\pgfPTCScombine{period,group,mix}
\pgfPT[back color scheme=mix,show title=false]
```



In the Periodic Table, a row is called a **period** and a column is called a **group**.

Representative elements: element families

For the ***representative elements*** (groups **1**, **2** and **13** to **18**) it is common to speak of families that reflect their common characteristics. So we have **the families**:

GROUP **1**: **Alkali metals**

► *lithium, sodium, potassium, rubidium, cesium and francium.*

The atoms of these elements **have** only **one valence electron**.

- ✓ They react violently with water to form hydroxides.
- ✓ They have a silver-gray color, with the exception of cesium, which has a golden hue.

GROUP **2**: **Alkaline earth metals**

► *beryllium, magnesium, calcium, strontium, barium and radium.*

The atoms of these elements **have two valence electrons**.

- ✓ Their oxides remain solid at high temperatures and form alkaline solutions.
- ✓ They react violently with water to form hydroxides.
- ✓ When they burn, they have reddish flames, excluding barium, which presents a greenish flame.

GROUP 1**3**: ***Boron*** **group**

► *boron, aluminium, gallium, indium, thallium and nihonium.*

The atoms of these elements **have three valence electrons**.

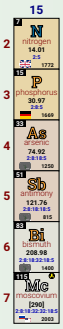
- ✓ Boron is a metalloid and the other are metals.
- ✓ Boron, aluminium, gallium, indium and thallium are often used as p-type silicon dopants.
- ✓ Aluminium is the third most abundant element in the Earth's crust (7.4%)

GROUP 1**4**: ***Carbon*** **group**

► *carbon, silicon, germanium, tin, lead and flerovium.*

The atoms of these elements **have four valence electrons**.

- ✓ Carbon is a non-metal, silicon and germanium are metalloids, and tin and lead are metals.
- ✓ Silicon and germanium are used in semiconductors.

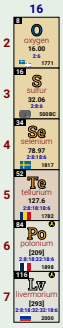


GROUP 1**5**: **Pnictogens**

► *nitrogen, phosphorus, arsenic, antimony, bismuth and moscovium*.

The atoms of these elements **have five valence electrons**.

- ✓ Nitrogen and phosphorus are non-metals, arsenic and antimony are metalloids and bismuth is a metal.
- ✓ Phosphorus, arsenic, antimony and bismuth are often used as n-type silicon dopants.
- ✓ Diatomic nitrogen is the main constituent of the Earth's atmosphere (78%).



GROUP 1**6**: **Chalcogens**

► *oxygen, sulfur, selenium, tellurium, polonium and livermorium*.

The atoms of these elements **have six valence electrons**.

- ✓ Oxygen, sulfur and selenium are non-metals, tellurium is a metalloid and polonium is a metal.
- ✓ Diatomic oxygen is the second constituent of the Earth's atmosphere (21%).

GROUP 1**7**: **Halogens**

► *fluorine, chlorine, bromine, iodine, astatine and tennessine*.

The atoms of these elements **have seven valence electrons**.

- ✓ They are extremely reactive elements, as they are very electronegative.
- ✓ Fluorine is able to *attack* inert substances, including the heavier noble gas atoms.

GROUP 1**8**: **Noble gases**

► *helium, neon, argon, krypton, xenon, radon and oganesson*.

The atoms of these elements have the valence shell fully filled, which corresponds to **eight valence electrons**, with the exception Helium, which has only one shell and, consequently, has **two valence electrons**.

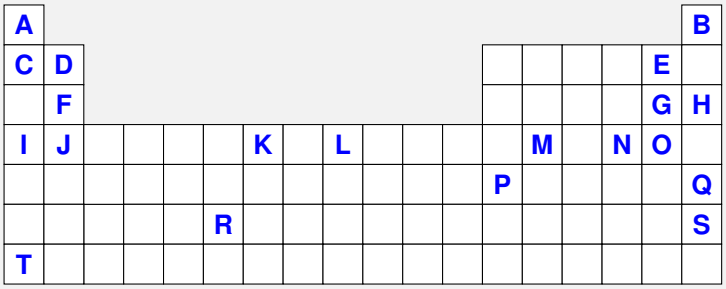
- ✓ They are extremely inert elements, that is, they do not react with other elements, as they are the most stable elements in Nature.

For the source of this example please see the file pgf-PeriodicTableManual_Examples.tex

EXERCISE:

In the following scheme of the Periodic Table, the positions of some chemical elements are represented by letters:

THE LETTERS DO NOT CORRESPOND TO THE CHEMICAL SYMBOLS OF THE ELEMENTS.



Using the letters shown:

1. identify group 2 elements of the Periodic Table.
2. identify the elements of the 2nd period of the Periodic Table.
3. identify group 17 elements of the Periodic Table.
4. identify the elements of block s.
5. identify the elements of block p.
6. identify the elements of block d.
7. identify the metallic elements.
8. identify the non-metallic elements.
9. identify the transition metals.
10. identify the alkaline earth metals.
11. identify the noble gases.
12. tell which element belongs, simultaneously, to the 4th period and to group 14.
13. identify the representative elements that tend to generate positive ions.
14. indicate an element that forms binegative ions.
15. indicate the halogen whose mononegative ion has the largest radius.
16. write the chemical formula of the compound formed by the elements **F** and **O**.
17. identify, justifying, the element with the largest atomic radius.
18. identify, justifying, the element with the lowest 1st ionization energy.

For the source of this example please see the file pgf-PeriodicTableManual_Examples.tex

EXERCISE:

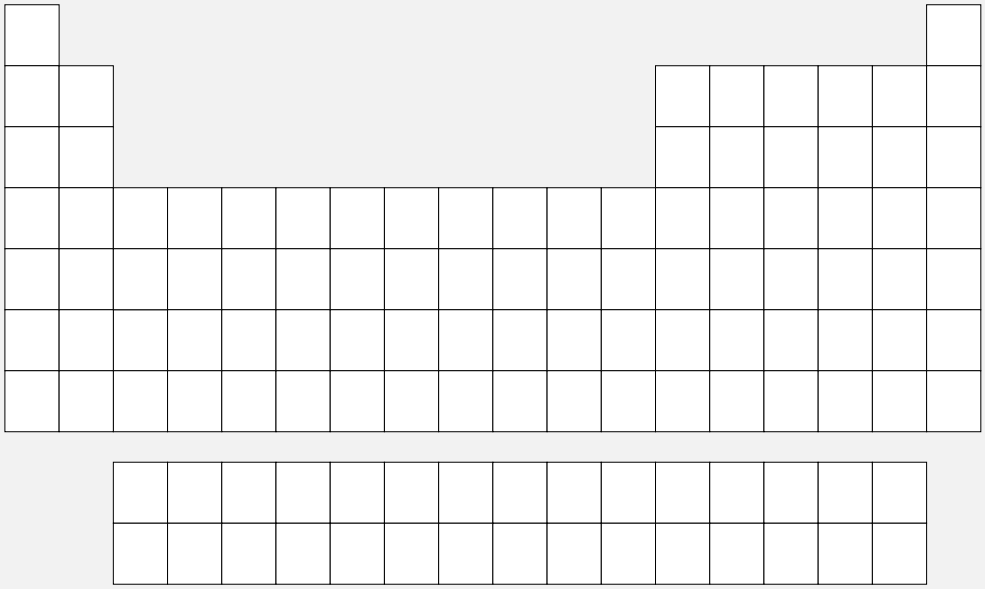
Using the following notation,

X for the elements in the gaseous state (NTP),

○ for the elements in the liquid state (NTP) and

△ for the synthetic elements,

fill in the following Periodic Table:



For the source of this example please see the file pgf-PeriodicTableManual_Examples.tex

Index

BUILT-IN

cell styles 102
color schemes 9

COMMANDS

\pgfPT 3
\pgfPTbuildcell 6
designing cells with 99
row, column syntax 100
\pgfPTbuildcellstyle 6
\pgfPTnewColorScheme 7
\pgfPTnewZlist 9
\pgfPTpreviewcell 6
\pgfPTpreviewcellstyle 7
\pgfPTresetcell 6
\pgfPTresetstyle 5
\pgfPTsetLanguage 10
\pgfPTstyle 4

LIBRARIES

Color Schemes Library 109
\pgfPTGroupColors 109
\pgfPTPeriodColors 114
\pgfPTCScombine 117
\pgfPTCSwrite 120

OPTIONS

<content name> color 90
<content name> font 91
Ar color 76
Ar font 77
Ar label 77
Ar precision 77
back color 16
back color scheme 15
blocks font 46
capitalize element names 75
cell height 12
cell line color 13
cell line width 13
cell style 13
cell width 12
Cp precision 95
CS font 72
CS gas 71
CS liquid 70
CS outline color 73
CS outline width 73
CS render mode 72
CS solid 70
CS synt 71
d block color 46
d block font color 46
d block line width 46
d color 79
d font 79
d precision 81
d unit 80
DiscY BC scale 88
DiscY color 87
DiscY font 87
E precision 92
eDist color 88
eDist font 89
eDist sep 89
exercise list color 64
exercise list font 64
exercise list in capitals 64
f block color 46
f block font color 46
f block line width 47
families font 50
font 14
group label color 42
itm family color 51
itm family font color 51
itm family line width 51
IUPAC 23
kT precision 96
label font 43
label LaAc font 26
languages 26
legend acronyms 33
legend back color 35
legend CS color 36
legend radio color 35
legend Z color 37
ls 83
ls align 85
ls color 84
ls font 84
ls precision 85
ls unit 85
MNM line color 29
MNM line width 30
name align 75
name color 74
name font 74
only cells 60
only cells plus Z 61
only cells with periods and group numbers plus Z 63
only cells with periods and group numbers 62
other languages color 28
other languages font 27
p block color 46
p block font color 46
p block line width 46
period label color 42
r family color 50
r family font color 50
r family line width 51
s block color 46
s block font color 46
s block line width 46

show blocks 44
show extra legend 39
show families 49
show group numbers 41
show label LaAc 25
show legend 33
show legend pins 38
show MNM line 28
show period numbers 41
show periodic variations 54
show title 31
T precision 93
title color 32
title font 32
tm family color 51
tm family font color 51
tm family line width 51
vareaff color 56
vareaff font 56
vareaff font color 56
varEi color 56
varEi font 56
varEi font color 56
varR color 55
varR font 55
varR font color 56
Z align 68
Z backcolor 67
Z color 68
Z exercise list 63
Z font 68
Z list 11
Z padding 69
Z use box width 68

STYLES

Ar 78
background 23
blocks 48
blocks font color 47
blocks line width 47
cell 14
cell color 91
cell font 91
cell size 12
cells+p+g 65
cells+p+g+Z 66
cells+Z 65
CS 74
CS all 71
csBlocks 22
csCPK 19
csJmol 18
csMNM 21
csPS 21
csRadio 22
csRasmol 19
csRasmolNew 20
csSoft 18
csSolid 17
csWikipedia 20
d 83
dark mode 60
ex 67
exColor 66
exFont 67
exnocaps 66
extra legend 39
families 52
families font color 51
families line width 51
gr 43
lat 87
legend 40
legend box 34
legend pins 38
MNM 30
NAME 76
Name 76
name 75
other lang 28
per 43
per+gr 44
title 32
var color 57
var font 56
vareaff 59
varEi 58
varR 57
Z 69
Z box 69