

Tableau périodique

 : importation de la librairie tkinter à unifier + codes à améliorer

[tableau_periodique.py](#)

```
#!/usr/bin/env python
# -*- coding: utf-8 -*-
# Programme sur le tableau périodique
# MJ, Ba2 chimie 2010-2011

from tkinter import *
from element_liste import * #sert à importer la liste présente dans
l'autre fichier

#création de la commande générale du bouton
def elem(x):
    element=Tk()
    element.title("Proprietes")
    listbox=Listbox(element,height=10,width=40,fg="#070942")
    listbox.pack() #sert à ajuster la fenêtre
    listbox.insert(END)
    for item in tableau[x]:
        listbox.insert(END,item)
    quitter = Button(element,text='Quitter',command=element.destroy)
    quitter.pack(side=BOTTOM)
    mainloop()

fen1=Tk()
fen1.title("Tableau periodique")

#création de la légende (site www.jchr.be/python/tkinter.htm)
import tkinter
systemenu=tkinter.Menu(fen1)
def legende():
    legend1= Tk()
    legend1.title("Légende")
    frame= Frame(legend1, height=200, width= 300) #frame ouvre une
nouvelle fenêtre dans laquelle on peut insérer plusieurs widgets (ici
des messages)
    autre = Message(legend1, text="Autres", width=300, font="Arial 15",
bg="white") #font = taille et police
    autre.pack()
    metauxtransition= Message(legend1, text="Metaux de transition",
width=300, font="Arial 15", bg="#FFCCCC")
    metauxtransition.pack()
    gazrare = Message(legend1, text="Gaz rares", width=300, font="Arial
15", bg="#00CCFF")
    gazrare.pack()
    nonmetaux= Message(legend1, text="Non metaux", width=300,
```

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font="Arial 15", bg="#00FF66")
    nonmetaux.pack()
    metauxalcalins= Message(legend1, text="Metaux alcalins", width=300,
font="Arial 15", bg="#FF9900")
    metauxalcalins.pack()
    metauxalcalinoterreux= Message(legend1, text="Metaux alcalino
terreux", width=300, font="Arial 15", bg="#FFFF33")
    metauxalcalinoterreux.pack()
    lanthanides= Message(legend1, text="Lanthanides", width=300,
font="Arial 15", bg="#CCCC66")
    lanthanides.pack()
    actinides= Message(legend1, text="Actinides", width=300,
font="Arial 15", bg="#FF3300")
    actinides.pack()
    metaux= Message(legend1, text="Metaux", width=300, font="Arial 15",
bg="#9999FF")
    metaux.pack()

    frame.pack()
    mainloop()

Menu=tkinter.Menu(sysdemenu)
sysdemenu.add_cascade(label="Menu", menu=Menu)
Menu.add_command(label="Legende", command = legende)
Menu.add_command(label="Quitter", command=fen1.quit)

#création des boutons du tableau
bou_list=[]
for item in tableau[1:]:
    bou=Button(fen1, text=item[1], command=lambda
x=item[2]:elem(x), height=2, width=3, bg=item[6])
    bou.grid(row=item[4], column=item[5]) #grid est utilisé pour classer
les boutons par rangée et colonne
    bou_list.append(bou)

fen1.config(menu=sysdemenu)

#création de la listbox au sommet
(source:http://www.tkdocs.com/tutorial/grid.html)
centre=Listbox(fen1)
centre.grid(row=0, column=3, rowspan= 1, columnspan=8)
centre.insert(END)
for item in ["Nom:", "Symbole:", "Numero atomique:", "Masse
molaire:", "Ligne:", "Colonne:", "Couleur:"]:
    centre.insert(END, item)
mainloop()

fen1.destroy()
```

```
fen1.mainloop()
```

Nécessite ce fichier de données :

[element_liste.py](#)

```
tableau=[
    ["Nom", "symbole", "Numero atomique", "Masse atomique",
    "ligne", "colonne", "couleur"],
    ["Hydrogene", "H", 1, 1.00794, 1, 0, "white"],
    ["Helium", "He", 2, 4.0026, 1, 17, "#00CCFF"],
    ["Lithium", "Li", 3, 6.9412, 2, 0, "#FF9900"],
    ["Beryllium", "Be", 4, 9.012182, 2, 1, "#FFFF33"],
    ["Bore", "B", 5, 10.811, 2, 12, "#00FF66"],
    ["Carbone", "C", 6, 12.0107, 2, 13, "#00FF66"],
    ["Azote", "N", 7, 14.00674, 2, 14, "#00FF66"],
    ["Oxygene", "O", 8, 15.9994, 2, 15, "#00FF66"],
    ["Fluor", "F", 9, 18.9984, 2, 16, "#00FF66"],
    ["Neon", "Ne", 10, 20.1797, 2, 17, "#00CCFF"],
    ["Sodium", "Na", 11, 22.9897, 3, 0, "#FF9900"],
    ["Magnesium", "Mg", 12, 24.3050, 3, 1, "#FFFF33"],
    ["Aluminium", "Al", 13, 26.9815, 3, 12, "#9999FF"],
    ["Silicium", "Si", 14, 28.0855, 3, 13, "#00FF66"],
    ["Phosphore", "P", 15, 30.973761, 3, 14, "#00FF66"],
    ["Soufre", "S", 16, 32.066, 3, 15, "#00FF66"],
    ["Chlore", "Cl", 17, 35.453, 3, 16, "#00FF66"],
    ["Argon", "Ar", 18, 39.948, 3, 17, "#00CCFF"],
    ["Potassium", "K", 19, 39.0963, 4, 0, "#FF9900"],
    ["Calcium", "Ca", 20, 40.078, 4, 1, "#FFFF33"],
    ["Scandium", "Sc", 21, 44.9559, 4, 2, "#FFCCCC"],
    ["Titane", "Ti", 22, 47.867, 4, 3, "#FFCCCC"],
    ["Vanadium", "V", 23, 50.9415, 4, 4, "#FFCCCC"],
    ["Chrome", "Cr", 24, 51.9961, 4, 5, "#FFCCCC"],
    ["Manganese", "Mn", 25, 54.9380, 4, 6, "#FFCCCC"],
    ["Fer", "Fe", 26, 55.8457, 4, 7, "#FFCCCC"],
    ["Cobalt", "Co", 27, 58.9332, 4, 8, "#FFCCCC"],
    ["Nickel", "Ni", 28, 58.6934, 4, 9, "#FFCCCC"],
    ["Cuivre", "Cu", 29, 63.546, 4, 10, "#FFCCCC"],
    ["Zinc", "Zn", 30, 65.409, 4, 11, "#FFCCCC"],
    ["Gallium", "Ga", 31, 69.723, 4, 12, "#9999FF"],
    ["Germanium", "Ge", 32, 72.64, 4, 13, "#9999FF"],
    ["Arsenic", "As", 33, 74.9216, 4, 14, "#00FF66"],
    ["Selenium", "Se", 34, 78.96, 4, 15, "#00FF66"],
    ["Brome", "Br", 35, 79.904, 4, 16, "#00FF66"],
    ["Krypton", "Kr", 36, 83.798, 4, 17, "#00CCFF"],
    ["Rubidium", "Rb", 37, 85.4678, 5, 0, "#FF9900"],
    ["Strontium", "Sr", 38, 87.62, 5, 1, "#FFFF33"],
    ["Yttrium", "Y", 39, 88.9059, 5, 2, "#FFCCCC"],
    ["Zirconium", "Zr", 40, 91.224, 5, 3, "#FFCCCC"],
    ["Niobium", "Nb", 41, 92.9063, 5, 4, "#FFCCCC"],
```

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[ "Molybdene", "Mo", 42, 95.94, 5, 5, "#FFCCCC" ],
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[ "Ruthenium", "Ru", 44, 101.07, 5, 7, "#FFCCCC" ],
[ "Rhodium", "Rh", 45, 102.9055, 5, 8, "#FFCCCC" ],
[ "Palladium", "Pd", 46, 106.42, 5, 9, "#FFCCCC" ],
[ "Argent", "Ag", 47, 107.8682, 5, 10, "#FFCCCC" ],
[ "Cadmium", "Cd", 48, 112.411, 5, 11, "#FFCCCC" ],
[ "Indium", "In", 49, 114.818, 5, 12, "#9999FF" ],
[ "Etain", "Sn", 50, 118.710, 5, 13, "#9999FF" ],
[ "Antimoine", "Sb", 51, 121.760, 5, 14, "#9999FF" ],
[ "Tellure", "Te", 52, 127.60, 5, 15, "#00FF66" ],
[ "Iode", "I", 53, 126.9045, 5, 16, "#00FF66" ],
[ "Xenon", "Xe", 54, 131.293, 5, 17, "#00CCFF" ],
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[ "Neodyme", "Nd", 60, 144.24, 8, 6, "#CCCC66" ],
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[ "Samarium", "Sm", 62, 150.36, 8, 8, "#CCCC66" ],
[ "Europium", "Eu", 63, 151.964, 8, 9, "#CCCC66" ],
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[ "Terbium", "Tb", 65, 158.9253, 8, 11, "#CCCC66" ],
[ "Dysprosium", "Dy", 66, 162.500, 8, 12, "#CCCC66" ],
[ "Holmium", "Ho", 67, 164.9303, 8, 13, "#CCCC66" ],
[ "Erbium", "Er", 68, 167.259, 8, 14, "#CCCC66" ],
[ "Thulium", "Tm", 69, 168.9342, 8, 15, "#CCCC66" ],
[ "Ytterbium", "Yb", 70, 173.04, 8, 16, "#CCCC66" ],
[ "Lutecium", "Lu", 71, 174.967, 8, 17, "#CCCC66" ],
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[ "Tantale", "Ta", 73, 180.9479, 6, 4, "#FFCCCC" ],
[ "Tungstene", "W", 74, 183.84, 6, 5, "#FFCCCC" ],
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[ "Osmium", "Os", 76, 190.23, 6, 7, "#FFCCCC" ],
[ "Iridium", "Ir", 77, 192.217, 6, 8, "#FFCCCC" ],
[ "Platine", "Pt", 78, 195.078, 6, 9, "#FFCCCC" ],
[ "Or", "Au", 79, 166.9665, 6, 10, "#FFCCCC" ],
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[ "Protactinium", "Pa", 91, 231.0359, 9, 5, "#FF3300" ],
[ "Uranium", "U", 92, 238.02891, 9, 6, "#FF3300" ],
[ "Neptunium", "Np", 93, 237, 9, 7, "#FF3300" ],
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[ "Berkelium", "Bk", 97, 247, 9, 11, "#FF3300" ],
[ "Californium", "Cf", 98, 251, 9, 12, "#FF3300" ],
[ "Einsteinium", "Es", 99, 252, 9, 13, "#FF3300" ],
[ "Fermium", "Fm", 100, 257, 9, 14, "#FF3300" ],
[ "Mendelevium", "Md", 101, 258, 9, 15, "#FF3300" ],
[ "Nobelium", "No", 102, 259, 9, 16, "#FF3300" ],
[ "Lawrencium", "Lr", 103, 262, 9, 17, "#FF3300" ],
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[ "Seaborgium", "Sg", 106, 266, 7, 5, "#FFCCCC" ],
[ "Bohrium", "Bh", 107, 264, 7, 6, "#FFCCCC" ],
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[ "Meitneium", "Mt", 109, 268, 7, 8, "#FFCCCC" ],
[ "Darmstadtium", "Ds", 110, 271, 7, 9, "#FFCCCC" ],
[ "Roentgenium", "Rg", 111, 272, 7, 10, "#FFCCCC" ],
[ "Copernicium", "Cn", 112, 285, 7, 11, "#FFCCCC" ],
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[ "Ununquadium", "Uuq", 114, 289, 7, 13, "#9999FF" ],
[ "Ununpentium", "Uup", 115, 288, 7, 14, "#9999FF" ],
[ "Ununhexium", "Uuh", 116, 292, 7, 15, "#9999FF" ],
[ "Ununseptium", "Uus", 117, "-", 7, 16, "white" ],
[ "Ununoctium", "Uuo", 118, "-", 7, 17, "white" ]
]
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Last update: **2021/04/19 21:00**

