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编程语言: python3.9

实验内容

• 阅读PLY使用手册

PLY (Python Lex-Yacc) — ply 4.0 documentation

Python Lex Yacc中文手册

PLY包含两个独立的模块: lex.py和yacc.py, 都定义在ply包下。lex.py模块用来将输入字符通过一系列的正则表达式分解成标记序列, yacc.py通过一些上下文无关的文法来识别编程语言语法。yacc.py使用LR解析法,并使用LALR(1)算法(默认)或者SLR算法生成分析表。

lex.py是用来将输入字符串标记化 yacc.py用来对语言进行语法分析。

熟悉yacc_example中分析四则运算的程序 标记列表

```
tokens = (
    'NUMBER',
    'PLUS',
    'MINUS',
    'TIMES',
    'DIVIDE',
    'LPAREN',
    'RPAREN',
)
```

标记的规则的形式有两种

```
t_PLUS = r'\+'

def t_NUMBER(t):
    r'\d+'
    t.value = int(t.value)
    return t
```

在yacc中每个语法规则被定义成一个Python的方法,方法的文档字符串描述了相应的上下文无关文法,方法的语句实现了对应规则的语义行为。每个方法接受一个单独的 p 参数, p 是一个包含有当前匹配语法的符号的序列, p[i] 与语法符号的对应关系如下:

• 编写程序, 计算化学分子式中元素的数目, 并完成以下测试:

```
atom\_count(He) == 1
atom\_count(H2) == 2
atom\_count(H2SO4) == 7
atom\_count(CH3COOH) == 8
atom\_count(NaCl) == 2
atom\_count(C60H60) == 120
```

实验步骤

使用lex进行序列标记

ply使用"t_"开头的变量来表示规则。如果变量是一个字符串,那么它被解释为一个正则表达式,匹配值是标记的值。 如果变量是函数,则其文档字符串包含模式,并使用匹配的标记调用该函数。该函数可以自由地修改序列或返回一个新的序列来代替它的位置。 如果没有返回任何内容,则忽略匹配。 通常该函数只更改"value"属性,它最初是匹配的文本。 在下面的 t_COUNT 将值转换为 int。

```
import ply.lex as lex
tokens = (
   "SYMBOL",
    "COUNT"
)
t_SYMBOL = (
   r"C[laroudsemf]?|Os?|N[eaibdpos]?|S[icernbmg]?|P[drmtboau]?|"
    r"H[eofgas]?|A[lrsgutcm]|B[eraik]?|Dy|E[urs]|F[erm]?|G[aed]|"
    r"I[nr]?|Kr?|L[iaur]|M[gnodt]|R[buhenaf]|T[icebmalh]|"
    r"U|V|W|Xe|Yb?|Z[nr]")
def t_COUNT(t):
   r"\d+"
    t.value = int(t.value)
    return t
def t_error(t):
    raise TypeError("Unknown text '%s'" % (t.value,))
```

对CH3COOH进行测试

```
lex.lex()

lex.input("CH3COOH")

for tok in iter(lex.token, None):
    print(repr(tok.type), repr(tok.value))
```

```
'SYMBOL' 'C'
'SYMBOL' 'H'
'COUNT' 3
'SYMBOL' 'C'
'SYMBOL' 'O'
'SYMBOL' 'O'
'SYMBOL' 'H'
```

使用yacc进行语法分析

PLY 的解析器适用于lex解析出的序列标记。 它使用 BNF 语法来描述这些标记是如何组装的。 解析器可以处理一些歧义。 在我们的化学分子式解析中,语法在阅读化学符号之后是不明确的。

解析算法的名称有 LALR(1)、SLR、LL 和 LR。

化学解析式有以下4种语法

```
species\_list: species\_listspecies species\_list: species species: SYMBOL species: SYMBOL\ COUNT
```

```
class Atom(object):
   def __init__(self, symbol, count):
       self.symbol = symbol
        self.count = count
   def __repr__(self):
        return "Atom(%r, %r)" % (self.symbol, self.count)
# When parsing starts, try to make a "chemical_equation" because it's
# the name on left-hand side of the first p_* function definition.
def p_species_list(p):
    "chemical_equation : chemical_equation species"
   p[0] = p[1] + [p[2]]
def p_species(p):
    "chemical_equation : species"
    p[0] = [p[1]]
def p_single_species(p):
    species : SYMBOL
   species : SYMBOL COUNT
   if len(p) == 2:
        p[0] = Atom(p[1], 1)
    elif len(p) == 3:
        p[0] = Atom(p[1], p[2])
def p_error(p):
```

```
print("Syntax error at '%s'" % p.value)
```

对H2SO4进行测试

```
yacc.yacc()
print(yacc.parse("H2SO4"))
```

结果如下

```
[Atom('H', 2), Atom('S', 1), Atom('0', 4)]
```

定义原子计数函数

```
def atom_count(s):
    count = 0
    for atom in yacc.parse(s):
        count += atom.count
    return count
```

实验结果

测试函数

为实验要求的几个化学分子式元素的数目进行测试

```
def test():
    assert atom_count("He") == 1
    assert atom_count("H2") == 2
    assert atom_count("H2SO4") == 7
    assert atom_count("CH3COOH") == 8
    assert atom_count("NaC1") == 2
    assert atom_count("C6OH6O") == 120
```

```
if __name__ == "__main__":
    test()
    print("All tests passed.")
```

结果

结果如下

```
All tests passed.

Process finished with exit code 0
```

其他

自动生成的语法分析表

```
# parsetab.py
# This file is automatically generated. Do not edit.
```

```
# pylint: disable=W,C,R
_tabversion = '3.10'
_lr_method = 'LALR'
_lr_signature = 'COUNT SYMBOL'n chemical_equation :\n chemical_equation :
species_list\n species_list : species_list speciesspecies_list : species\n
_lr_action_items = {'$end':([0,1,2,3,4,5,6,],[-1,0,-2,-4,-5,-3,-6,]),'SYMBOL':
([0,2,3,4,5,6,],[4,4,-4,-5,-3,-6,]),'COUNT':([4,],[6,]),
_1r_action = {}
for _k, _v in _lr_action_items.items():
   for _x,_y in zip(_v[0],_v[1]):
     if not _x in _lr_action: _lr_action[_x] = {}
     _lr_action[_x][_k] = _y
del _lr_action_items
_lr_qoto_items = {'chemical_equation':([0,],[1,]),'species_list':([0,],
[2,]), 'species':([0,2,],[3,5,]),}
ln=0
for _k, _v in _lr_goto_items.items():
   for _x, _y in zip(_v[0], _v[1]):
      if not _x in _lr_goto: _lr_goto[_x] = {}
      _lr_goto[_x][_k] = _y
del _lr_goto_items
_lr_productions = [
  ("S' -> chemical_equation", "S'", 1, None, None, None),
  ('chemical_equation ->
<empty>','chemical_equation',0,'p_chemical_equation','main.py',48),
  ('chemical_equation ->
species_list','chemical_equation',1,'p_chemical_equation','main.py',49),
  ('species_list -> species_list
species','species_list',2,'p_species_list','main.py',59),
  ('species_list -> species','species_list',1,'p_species','main.py',64),
  ('species -> SYMBOL', 'species',1,'p_single_species', 'main.py',70),
  ('species -> SYMBOL COUNT', 'species', 2, 'p_single_species', 'main.py', 71),
]
```

yacc.py在生成分析表时会创建出一个调试文件

文件中出现的不同状态,代表了有效输入标记的所有可能的组合,这是依据文法规则得到的。当得到输入标记时,分析器将构造一个栈,并找到匹配的规则。每个状态跟踪了当前输入进行到语法规则中的哪个位置,在每个规则中, 2.2 表示当前分析到规则的哪个位置,而且,对于在当前状态下,输入的每个有效标记导致的动作也被罗列出来。

```
Created by PLY version 3.11 (http://www.dabeaz.com/ply)

Grammar

Rule 0 S' -> chemical_equation
Rule 1 chemical_equation -> <empty>
Rule 2 chemical_equation -> species_list
Rule 3 species_list -> species_list species
Rule 4 species_list -> species
```

```
Rule 5 species -> SYMBOL
Rule 6
         species -> SYMBOL COUNT
Terminals, with rules where they appear
COUNT
                    : 6
SYMBOL
                    : 5 6
error
Nonterminals, with rules where they appear
chemical_equation : 0
species
                     : 3 4
species_list
                   : 2 3
Parsing method: LALR
state 0
    (0) S' -> . chemical_equation
    (1) chemical_equation -> .
    (2) chemical_equation -> . species_list
    (3) species_list -> . species_list species
    (4) species_list -> . species
    (5) species -> . SYMBOL
    (6) species -> . SYMBOL COUNT
    $end
                   reduce using rule 1 (chemical_equation -> .)
    SYMBOL
                   shift and go to state 4
   chemical_equation
                                   shift and go to state 1
    species_list
                                  shift and go to state 2
   species
                                  shift and go to state 3
state 1
    (0) S' -> chemical_equation .
state 2
    (2) chemical_equation -> species_list .
    (3) species_list -> species_list . species
    (5) species -> . SYMBOL
    (6) species -> . SYMBOL COUNT
    $end
                   reduce using rule 2 (chemical_equation -> species_list .)
    SYMBOL
                   shift and go to state 4
   species
                                   shift and go to state 5
state 3
    (4) species_list -> species .
    SYMBOL
                    reduce using rule 4 (species_list -> species .)
                    reduce using rule 4 (species_list -> species .)
    $end
```

```
state 4
    (5) species -> SYMBOL .
   (6) species -> SYMBOL . COUNT
                   reduce using rule 5 (species -> SYMBOL .)
   SYMBOL
                  reduce using rule 5 (species -> SYMBOL .)
   $end
                   shift and go to state 6
   COUNT
state 5
   (3) species_list -> species_list species .
   SYMBOL
                   reduce using rule 3 (species_list -> species_list species .)
   $end
                   reduce using rule 3 (species_list -> species_list species .)
state 6
   (6) species -> SYMBOL COUNT .
                   reduce using rule 6 (species -> SYMBOL COUNT .)
   SYMBOL
                   reduce using rule 6 (species -> SYMBOL COUNT .)
    $end
```