# CS 536 <br> Introduction to Programming Languages and Compilers <br> <br> Charles N. Fischer 

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## Lecture 7

## Java CUP

Java CUP is a parser- generation tool, similar to Yacc.
CUP builds a Java parser for LALR(1) grammars from production rules and associated Java code fragments.
When a particular production is recognized, its associated code fragment is executed (typically to build an AST).
CUP generates a Java source file parser.java. It contains a class parser, with a method
Symbol parse()
The symbol returned by the parser is associated with the grammar's start symbol and contains the AST for the whole source program.

The file sym.java is also built for use with a JLex- built scanner (so that both scanner and parser use the same token codes).
If an unrecovered syntax error occurs, Exception() is thrown by the parser.
CUP and Yacc accept exactly the same class of grammars-all LL(1) grammars, plus many useful nonLL(1) grammars.
CUP is called as
java java_cup. Main < file.cup

## Java CUP Specifications

Java CUP specifications are of the form:

- Package and import specifications
- User code additions
- Terminal and non- terminal declarations
- A context- free grammar, augmented with Java code fragments


## Package and Import Specifications

 You define a package name as:package name ;
You add imports to be used as: import java_cup.runtime.*;

## User Code Additions

You may define Java code to be included within the generated parser:
action code \{: /*java code */ :\} This code is placed within the generated action class (which holds user- specified production actions).
parser code \{: /*java code */ :\} This code is placed within the generated parser class.
init with\{: /*java code */ :\} This code is used to initialize the generated parser.
scan with\{: /*java code */ : \} This code is used to tell the generated parser how to get tokens from the scanner.

## Terminal and Non-terminal Declarations

You define terminal symbols you will use as:
terminal classname name ${ }_{1}$, name ${ }_{2}, \ldots$ classname is a class used by the scanner for tokens (CSXToken, CSXIdentifierToken, etc.)

You define non- terminal symbols you will use as: non terminal classname name ${ }_{1}$, name ${ }_{2}, \ldots$ classname is the class for the AST node associated with the non- terminal (stmtNode, exprNode, etc.)

## Production Rules

Production rules are of the form name ::= name ${ }_{1}$ name ${ }_{2}$... action ; or name ::= name ${ }_{1}$ action $_{1}$ name $_{2}$... | name $_{3}$ name $_{4} \ldots$ action $_{2}$ | ... ;
Names are the names of terminals or non- terminals, as declared earlier.
Actions are Java code fragments, of the form
\{: /*java code */ :\}
The Java object assocated with a symbol (a token or AST node) may be named by adding a :id suffix to a terminal or non- terminal in a rule.

Result names the left- hand side non- terminal.
The Java classes of the symbols are defined in the terminal and non- terminal declaration sections.
For example,
prog ::= LBRACE:l stmts:s RBRACE \{: RESULT =
new csxLiteNode(s,
l.linenum,l.colnum); :\}

This corresponds to the production prog $\rightarrow$ \{ stmts \}
The left brace is named 1 ; the stmts non- terminal is called $\mathbf{s}$.
In the action code, a new CSXLiteNode is created and assigned to prog. It is constructed from the AST node associated with s. Its line and column
numbers are those given to the left brace, 1 (by the scanner).
To tell CUP what non- terminal to use as the start symbol (prog in our example), we use the directive:
start with prog;

## Example

# Let's look at the CUP specification for CSX- lite. Recall its CFG is 

$$
\begin{aligned}
& \text { program } \rightarrow \text { \{stmts \} } \\
& \text { stmts } \rightarrow \text { stmt stmts } \\
& \mid \lambda \\
& \text { stmt } \rightarrow \text { id }=\text { expr ; } \\
& \text { | if ( expr ) stmt } \\
& \text { expr } \rightarrow \text { expr }+ \text { id } \\
& \text { | expr - id } \\
& \text { | id }
\end{aligned}
$$

## The corresponding CUP specification is:

/***
This Is A Java CUP Specification For CSX-lite, a Small Subset of The CSX Language, Used In Cs536 ***/
/* Preliminaries to set up and use the scanner. */
import java_cup.runtime.*;
parser code \{: public void syntax_error (Symbol cur_token) \{
report_error(
"CSX syntax error at line " + String.valueOf(( CSXToken)
cur_token.value).linenum), null); $\}$
: \};
init with \{: : \};
scan with \{:
return Scanner.next_token();
: \};
/* Terminals (tokens returned by the scanner). */
terminal CSXIdentifierToken IDENTIFIER; terminal CSXToken SEMI, LPAREN, RPAREN, ASG, LBRACE, RBRACE;
terminal CSXToken PLUS, MINUS, rw_IF;
/* Non terminals */
non terminal csxLiteNode prog;
non terminal stmtsNode stmts;
non terminal stmtNode stmt;
non terminal exprNode exp;
non terminal nameNode ident;
start with prog;
prog: := LBRACE:l stmts:s RBRACE
\{: RESULT=
new csxLiteNode(s,
l.linenum,l.colnum); : \}
;
stmts: : $=$ stmt:s1 stmts:s2
\{: RESULT= new stmtsNode(s1,s2, s1.linenum, s1.colnum);
: \}

```
|
    {: RESULT= stmtsNode.NULL; : }
;
stmt::= ident:id ASG exp:e SEMI
    {: RESULT=
                        new asgNode(id,e,
                        id.linenum,id.colnum);
    :}
| rw_IF:i LPAREN exp:e RPAREN stmt:s
    {: RESULT=new ifThenNode(e,s,
                        stmtNode.NULL,
                            i.linenum,i.colnum); :}
;
exp::=
    exp:leftval PLUS:Op ident:rightval
    {: RESULT=new binaryOpNode(leftval,
    sym.PLUS, rightval,
            op.linenum,op.colnum); :}
    | exp:leftval MINUS:op ident:rightval
    {: RESULT=new binaryOpNode(leftval,
                                    sym.MINUS,rightval,
                        op.linenum,op.colnum); :}
    | ident:i
    {: RESULT = i; :}
;
```

ident::= IDENTIFIER:i
\{: RESULT = new nameNode ( new identNode(i.identifierText, i.linenum,i.colnum), exprNode.NULL, i.linenum,i.colnum); : \} ;
Let's parse

$$
\{\mathbf{a}=\mathbf{b} ;\}
$$

First, a is parsed using
ident::= IDENTIFIER:i
\{: RESULT = new nameNode (
new identNode(i.identifierText,
i.linenum,i.colnum),
exprNode.NULL

$$
\text { i.linenum,i.colnum); :\} }
$$

We build


Next, b is parsed using
ident::= IDENTIFIER:i
\{: RESULT = new nameNode ( new identNode(i.identifierText, i.linenum,i.colnum), exprNode.NULL, i.linenum,i.colnum); :\}

We build


Then b's subtree is recognized as an exp:
| ident:i
\{: RESULT = i; :\}
Now the assignment statement is recognized:
stmt::= ident:id ASG exp:e SEMI
\{: RESULT=
new asgNode(id,e,
id.linenum,id.colnum);
: \}
We build


The stmts $\rightarrow \lambda$ production is matched (indicating that there are no more statements in the program).
CUP matches
stmts: : =
\{: RESULT= stmtsNode.NULL; : \}
and we build

## nullStmtsNode

Next,
stmts $\rightarrow$ stmt stmts is matched using
stmts: : = stmt:s1 stmts:s2
\{: RESULT=
new stmtsNode(s1,s2, s1.linenum, s1.colnum);
: \}

This builds


As the last step of the parse, the parser matches
program $\rightarrow$ \{ stmts \} using the CUP rule
prog: := LBRACE:l stmts:s RBRACE
\{: RESULT=
new csxLiteNode(s,
l.linenum,l.colnum); : \}
;

## The final AST reurned by the parser is



## Errors in Context-Free Grammars

Context- free grammars can contain errors, just as programs do. Some errors are easy to detect and fix; others are more subtle. In context- free grammars we start with the start symbol, and apply productions until a terminal string is produced.
Some context- free grammars may contain useless non- terminals.
Non- terminals that are unreachable (from the start symbol) or that derive no terminal string are considered useless.
Useless non- terminals (and productions that involve them) can be safely removed from a grammar without changing the
language defined by the grammar.
A grammar containing useless
non-terminals is said to be nonreduced.
After useless non- terminals are removed, the grammar is reduced.
Consider
$\mathbf{S} \rightarrow \mathbf{A}$
| $\mathbf{x}$
$B \rightarrow b$
$\mathrm{A} \rightarrow \mathrm{a}$ A
$C \rightarrow d$

Which non- terminals are
unreachable? Which derive no terminal string?

## Finding Useless Non- terminals

To find non- terminals that can derive one or more terminal strings, we'll use a marking algorithm.
We iteratively mark terminals that can derive a string of terminals, until no more non- terminals can be marked. Unmarked nonterminals are useless.
(1) Mark all terminal symbols
(2) Repeat

If all symbols on the righthand side of a production are marked Then mark the lefthand side Until no more non-terminals can be marked

We can use a similar marking algorithm to determine which non- terminals can be reached from the start symbol:
(1) Mark the Start Symbol
(2) Repeat

If the lefthand side of a production is marked
Then mark all non-terminals in the righthand side Until no more non- terminals can be marked

## $\lambda$ Derivations

When parsing, we'll sometimes
need to know which nonterminals can derive $\lambda$. ( $\lambda$ is
"invisible" and hence tricky to parse).
We can use the following marking algorithm to decide which nonterminals derive $\lambda$
(1) For each production $A \rightarrow \lambda$ mark A
(2) Repeat

If the entire righthand side of a production is marked
Then mark the lefthand side Until no more non- terminals can be marked

## As an example consider $S \rightarrow$ AB C $A \rightarrow a B$ <br> $$
\rightarrow C \quad D
$$ <br> $$
D \rightarrow \mathbf{d}
$$ <br> $$
\mid \lambda
$$ <br> $$
C \rightarrow c
$$ <br> $$
\mid \lambda
$$

Recall that compilers prefer an unambiguous grammar because a unique parse tree structure can be guaranteed for all inputs.
Hence a unique translation,
guided by the parse tree structure, will be obtained.
We would like an algorithm that checks if a grammar is ambiguous.
Unfortunately, it is undecidable whether a given CFG is ambiguous, so such an algorithm is impossible to create.
Fortunately for certain grammar classes, including those for which we can generate parsers, we can prove included grammars are unambiguous.

Potentially, the most serious flaw that a grammar might have is that it generates the "wrong language."
This is a subtle point as a grammar serves as the definition of a language.
For established languages (like C or Java) there is usually a suite of programs created to test and validate new compilers. An incorrect grammar will almost certainly lead to incorrect compilations of test programs, which can be automatically recognized.
For new languages, initial implementors must thoroughly test the parser to verify that inputs are scanned and parsed as expected.

## Parsers and Recognizers

Given a sequence of tokens, we can ask:
"Is this input syntactically valid?"
(Is it generable from the grammar?).
A program that answers this question is a recognizer.
Alternatively, we can ask:
"Is this input valid and, if it is, what is its structure (parse tree)?"
A program that answers this more general question is termed a parser.
We plan to use language structure to drive compilers, so we will be especially interested in parsers.

## Two general approaches to parsing exist.

The first approach is top-down.
A parser is top- down if it "discovers" the parse tree corresponding to a token sequence by starting at the top of the tree (the start symbol), and then expanding the tree (via predictions) in a depth- first manner.
Top- down parsing techniques are predictive in nature because they always predict the production that is to be matched before matching actually begins.

Consider

$$
\begin{array}{l|l}
\mathrm{E} \rightarrow \mathrm{E}+\mathbf{T} & \mathbf{T} \\
\mathbf{T} \rightarrow \mathbf{T} \text { id } & \text { id }
\end{array}
$$

To parse id+id in a top- down manner, a parser build a parse tree in the following steps:

$$
\begin{aligned}
& \mathbf{E} \Rightarrow \mathrm{F} \Rightarrow \mathrm{E} \quad \Rightarrow \\
& \mathbf{E}+\mathbf{T} \\
& \begin{array}{cc}
\mathbf{E}+\mathrm{T} \\
\mathrm{~T} & \mathrm{E}
\end{array} \\
& \text { /1 } \backslash \\
& \text { ハ } \\
& \begin{array}{l}
\mathrm{E}+\mathrm{T} \\
\underset{T}{2}
\end{array} \\
& \text { id }
\end{aligned}
$$

A wide variety of parsing techniques take a different approach.
They belong to the class of bottom- up parsers.
As the name suggests, bottom- up parsers discover the structure of a parse tree by beginning at its bottom (at the leaves of the tree which are terminal symbols) and determining the productions used to generate the leaves.
Then the productions used to generate the immediate parents of the leaves are discovered.
The parser continues until it reaches the production used to expand the start symbol.
At this point the entire parse tree has been determined.

A bottom- up parse of $\mathrm{id}_{1}+\mathrm{id}_{2}$ would follow the following steps:


## A Simple Top-Down Parser

 We'll build a rudimentary topdown parser that simply tries each possible expansion of a nonterminal, in order of production definition.If an expansion leads to a token sequence that doesn't match the current token being parsed, we backup and try the next possible production choice.
We stop when all the input tokens are correctly matched or when all possible production choices have been tried.

## Example

Given the productions
$\mathbf{S} \rightarrow \mathbf{a}$
I (S)
we try a, then (a), then ((a)), etc.
Let's next try an additional alternative:

$$
S \rightarrow a
$$

$$
\begin{array}{ll}
\text { ( } \mathrm{S} \text { ) } \\
\text { | } & \text { ( }
\end{array}
$$

Let's try to parse a, then (a], then ((a]], etc.
We'll count the number of productions we try for each input.

- For input = a

We try $\mathbf{S} \rightarrow$ a which works.
Matches needed = 1

- For input = ( a ]

We try $\mathbf{S} \rightarrow$ a which fails.
We next try $\mathbf{S} \rightarrow$ ( $\mathbf{S}$ ).
We expand the inner S three
different ways; all fail.
Finally, we try S $\rightarrow$ ( J .
The inner $S$ expands to a, which works.
Total matches tried $=$
$1+(1+3)+(1+1)=7$.

- For input = (( a ]]

We try $\mathbf{S} \rightarrow$ a which fails.
We next try $\mathbf{S} \rightarrow(\mathbf{S})$.
We match the inner S to (a] using 7 steps, then fail to match the last ].
Finally, we try $S \rightarrow$ ( S ].
We match the inner S to (a] using 7
steps, then match the last ].
Total matches tried $=$

$$
1+(1+7)+(1+7)=17
$$

- For input = ((( a ] ]]

We try $\mathbf{S} \rightarrow$ a which fails.
We next try $\mathbf{S} \rightarrow$ ( S ).
We match the inner $S$ to ((a]] using
17 steps, then fail to match the last ].
Finally, we try S $\rightarrow$ ( S ]. We match the inner $S$ to ((a]] using 17 steps, then match the last ]. Total matches tried $=$

$$
1+(1+17)+(1+17)=37
$$

Adding one extra ( ... ] pair doubles the number of matches we need to do the parse.
In fact to parse ( ${ }^{\text {i }}{ }^{\text {j }}$ itakes $5^{*} 2^{i}-3$ matches. This is exponential growth!

With a more effective dynamic programming approach, in which results of intermediate parsing steps are cached, we can reduce the
number of matches needed to $n^{3}$ for an input with n tokens.
Is this acceptable?
No!
Typical source programs have at least 1000 tokens, and $1000^{3}=10^{9}$ is a lot of steps, even for a fast modern computer.
The solution?
-Smarter selection in the choice of productions we try.

## Reading Assignment Read Chapter 5 of Crafting a Compiler.

## Prediction

We want to avoid trying productions that can't possibly work.
For example, if the current token to be parsed is an identifier, it is useless to try a production that begins with an integer literal. Before we try a production, we'll consider the set of terminals it might initially produce. If the current token is in this set, we'll try the production.
If it isn't, there is no way the production being considered could be part of the parse, so we'll ignore it.
A predict function tells us the set of tokens that might be initially generated from any production.

For $A \rightarrow X_{1} \ldots X_{n}$, Predict $(A \rightarrow$
$\left.X_{1} \ldots X_{n}\right)=$ Set of all initial (first) tokens derivable from $A \rightarrow X_{1} \ldots X_{n}$
$=\left\{a \operatorname{in} V_{t} \mid A \rightarrow X_{1} \ldots X_{n} \Rightarrow^{*} a \ldots\right\}$
For example, given
Stmt $\rightarrow$ Label id = Expr ;
| Label if Expr then Stmt ; | Label read ( IdList ) ; | Label id (Args );
Label $\rightarrow$ intlit :

$$
\mid \lambda
$$

| Production | Predict Set |
| :--- | :--- |
| Stmt $\rightarrow$ Label id = Expr ; | \{id, intlit \} |
| Stmt $\rightarrow$ Label if Expr then Stmt ; | \{if, intlit\} |
| Stmt $\rightarrow$ Label read ( IdList ) ; | \{read, intlit \} |
| Stmt $\rightarrow$ Label id ( Args ) ; | \{id, intlit \} |

We now will match a production $p$ only if the next unmatched token is in p's predict set. We'll avoid trying productions that clearly won't work, so parsing will be faster.
But what is the predict set of a $\lambda$ - production?
It can't be what's generated by $\lambda$ (which is nothing!), so we'll define it as the tokens that can follow the use of a $\lambda$ - production.
That is, $\operatorname{Predict}(\mathrm{A} \rightarrow \lambda)=\operatorname{Follow}(\mathrm{A})$ where (by definition)
Follow $(A)=\left\{a\right.$ in $V_{t} \mid S \Rightarrow^{+}$...Aa... $\}$
In our example,
Follow(Label $\rightarrow \lambda$ ) $=\{$ id, if, read $\}$
(since these terminals can immediately follow uses of Label in the given productions).

Now let's parse id (intlit) ;
Our start symbol is Stmt and the initial token is id.
id can predict
Stmt $\rightarrow$ Label id = Expr;
id then predicts Label $\rightarrow \lambda$
The id is matched, but "(" doesn't match "=" so we backup and try a different production for Stmt. id also predicts
Stmt $\rightarrow$ Label id (Args );
Again, Label $\rightarrow \lambda$ is predicted and used, and the input tokens can match the rest of the remaining production.
We had only one misprediction, which is better than before.
Now we'll rewrite the productions a bit to make predictions easier.

We remove the Label prefix from all the statement productions (now intlit won't predict all four productions).
We now have
Stmt $\rightarrow$ Label BasicStmt BasicStmt $\rightarrow$ id = Expr ;

$$
\begin{aligned}
& \text { if Expr then Stmt ; } \\
& \quad \text { read (IdList) ; } \\
& \text { id (Args ); }
\end{aligned}
$$

Label $\rightarrow$ intlit :
| $\lambda$
Now id predicts two different BasicStmt productions. If we rewrite these two productions into
BasicStmt $\rightarrow$ id StmtSuffix StmtSuffix $\rightarrow=$ Expr ;
| (Args );

# we no longer have any doubt over which production id predicts. We now have 

| Production | Predict Set |
| :--- | :--- |
| Stmt $\rightarrow$ Label BasicStmt | Not needed! |
| BasicStmt $\rightarrow$ id StmtSuffix | \{id \} |
| BasicStmt $\rightarrow$ if Expr then Stmt ; | \{if $\}$ |
| BasicStmt $\rightarrow$ read ( IdList ) ; | \{read \} |
| StmtSuffix $\rightarrow$ (Args ); | $\{( \}$ |
| StmtSuffix $\rightarrow$ Expr ; | $\{=\}$ |
| Label $\rightarrow$ intlit : | \{intlit $\}$ |
| Label $\rightarrow \lambda$ | \{if, id, read $\}$ |

> This grammar generates the same statements as our original grammar did, but now prediction never fails!

Whenever we must decide what production to use, the predict sets for productions with the same lefthand side are always disjoint.
Any input token will predict a unique production or no production at all (indicating a syntax error).
If we never mispredict a production, we never backup, so parsing will be fast and absolutely accurate!

## LL(1) Grammars

A context- free grammar whose Predict sets are always disjoint (for the same non-terminal) is said to be $L L(1)$.
LL(1) grammars are ideally suited for top- down parsing because it is always possible to correctly predict the expansion of any nonterminal. No backup is ever needed.
Formally, let
$\operatorname{First}\left(X_{1} \ldots X_{n}\right)=$
$\left\{\mathrm{a}\right.$ in $\left.\mathrm{V}_{\mathrm{t}} \mid \mathrm{A} \rightarrow \mathrm{X}_{1} \ldots \mathrm{X}_{\mathrm{n}} \Rightarrow^{*} \mathrm{a} \ldots\right\}$
Follow $(A)=\left\{a\right.$ in $V_{t} \mid S \Rightarrow^{+}$...Aa... $\}$
$\operatorname{Predict}\left(A \rightarrow X_{1} \ldots X_{n}\right)=$
If $X_{1} \ldots X_{n}{ }^{*} \lambda$
Then First $\left(X_{1} \ldots X_{n}\right)$ U Follow(A)
Else First $\left(X_{1} \ldots X_{n}\right)$
If some CFG, G, has the property that for all pairs of distinct productions with the same lefthand side,
$A \rightarrow X_{1} \ldots X_{n}$ and $A \rightarrow Y_{1} \ldots Y_{m}$ it is the case that $\operatorname{Predict}\left(A \rightarrow X_{1} \ldots X_{n}\right) \cap$ $\operatorname{Predict}\left(A \rightarrow Y_{1} \ldots Y_{m}\right)=\phi$ then G is $\mathrm{LL}(1)$.
$\mathrm{LL}(1)$ grammars are easy to parse in a top- down manner since predictions are always correct.

## Example

| Production | Predict Set |
| :--- | :--- |
| $\mathbf{S} \rightarrow \mathbf{A} \mathbf{a}$ | $\{b, \mathrm{~d}, \mathrm{a}$ |
| $\mathbf{A} \rightarrow \mathbf{B} \mathbf{D}$ | $\{b, \mathbf{d}, \mathrm{a}\}$ |
| $\mathbf{B} \rightarrow \mathbf{b}$ | $\{b\}$ |
| $\mathbf{B} \rightarrow \lambda$ | $\{d, a\}$ |
| $\mathbf{D} \rightarrow \mathbf{d}$ | $\{d\}$ |
| $\mathbf{D} \rightarrow \lambda$ | $\{a\}$ |

Since the predict sets of both $B$ productions and both D productions are disjoint, this grammar is LL(1).

