1 Introduction

This document should explain some essential technical details for symbolic mode programming in REDUCE for those who have some experience in REDUCE algebraic programming and need to write a program in symbolic mode. For a general introduction to REDUCE please consult the REDUCE User’s Manual or one of the available books, e.g. “Algebraic computing with REDUCE” by Malcolm MacCallum and Francis Wright (Oxford Press).

This text cannot be complete, as the set of facilities available in REDUCE is so rich that it would take years to describe all and months to read and understand such text. So a good policy for entering the business of symbolic mode programming is to study the source files - the liberal REDUCE distribution policy simplifies this - and to understand those parts which contribute to the field which one would like to use. This text tries to collect in one place some of the wide spread basic information in order to facilitate your walk through the REDUCE mountains.

When should you write a program in symbolic mode? Symbolic programs are not a priori better than algebraic programs - the essential thing is the mathematics which they implement. A common prejudice is that symbolic programs are more “efficient”. This is true only if you can save in symbolic mode substantial algebraic evaluation steps. As long as most of the computing time is needed for a few big calculations steps (integrals, equation solving, polynomial gcd etc.) you will gain nothing when calling the same procedures from symbolic mode. However, if you need structures which are
not present in algebraic mode or if you can design an evaluation avoiding lots of conversions, the step to symbolic mode programming is justified.

As it is very difficult to design non trivial but short examples for symbolic mathematical programming no attempt has been made in that direction. The examples in this text all are silly - please look at the sources of REDUCE for meaningful material. The following pieces of the sources are good points for first reading as they provide a substantial functionality within a few pages of code:

1. module \texttt{polrep} in package \texttt{poly}: routines \texttt{addf}, \texttt{addd} and \texttt{addm}, the heart of standard form and standard quotient arithmetic,

2. module \texttt{det} of package \texttt{matrix}: internal arithmetic with symbolic entities (standard quotients) and clever local data structure,

3. module \texttt{rational} in package \texttt{poly}: implementation of a typical REDUCE domain,

4. module \texttt{maxmin} in package \texttt{alg}: a typical simplification routine for an operator,

5. module \texttt{algbool} in package \texttt{alg}: demonstrates how to supply “small” pieces of symbolic code for algebraic use.

For symbolic mode programming you will urgently need the \textit{Standard LISP Report} which describes the basic LISP functions; these will be available under all REDUCE implementations and they guarantee an optimum of portability. However, in the course of the years REDUCE has imported some additional functions on the LISP level – they have been implemented on top of Standard LISP and live in the module \texttt{support} of the package \texttt{rlisp.red}. In order to prevent the reinvention of the wheel these functions are described in the appendix as an extension to the \textit{Standard LISP Report}.

The description is based on the recent version of REDUCE. Some of the described features are not available in earlier versions.

\section{Very short course on RLISP}

\subsection{What is RLISP}

As you will know REDUCE is based on the programming language LISP, or to be more exact, on the LISP dialect \texttt{Standard LISP}. Fortunately you need not learn the syntax of this language with its large number of brackets
- the REDUCE language is used for algebraic programming as well as for symbolic programs and in that mode it allows you to use all of the rich LISP data structures. It is LISP semantics in high level REDUCE syntax. In the following it is expected that you are familiar with the REDUCE programming language as far as it is used for algebraic mode programs. You should know how to write a (recursive) procedure with parameters and local variables, how to build loops, while blocks, if-then-else clauses and begin-end resp << -- >> blocks. If not, please study the REDUCE manual first or have a look at the source files of the REDUCE kernel - this is the best method for learning how to program in RLISP.

2.2 Modes, function classes

The symbols symbolic (or equivalent lisp) and algebraic play a double role: as a statement they switch the evaluation mode globally to symbolic or algebraic mode respectively, as a mode prefix they tell the REDUCE evaluator that one isolated evaluation should be done in the named mode. The scope of this prefix use can be rather narrow (e.g. one single expression) or cover a larger piece of code up to a complete procedure. Typically procedures in symbolic modules should be tagged “symbolic” explicitly although this might be redundant information in a totally symbolic context. If a procedure needs an evaluation mode different from the actual global one the mode prefix must be set.

In symbolic mode there are two additional procedure types available, macro and smacro. The discussion of macros is beyond the scope of this document - their use requires extensive knowledge of LISP. On the other hand smacros are frequently use in REDUCE and you will see lots of them in the sources. An smacro (an abbreviation for “substitution macro”) is an ordinary procedure tagged with the symbol “smacro” and usually with a rather small body. At source read time (or better: at REDUCE translator time) any call for an smacro will be replaced literally by the body of the smacro procedure which saves execution time for the call protocol at run time. So the purpose of smacros is twofold: encapsulation of frequently used pieces of code and an increased efficiency (avoiding function calls). Example:

smacro procedure my_diff(x,y); x-y;

symbolic procedure hugo(a,b); my_diff(a,b);

Here the formal function call my_diff in the procedure literally is replaced by the body of the smacro where x and y are replaced by a and b. Obviously
this translation can be done only if the smacro declaration is entered in a REDUCE session before the first reference. And later changes of an smacro don’t affect previously translated references.

Sometimes in older REDUCE sources you will find the symbol expr in front of a procedure declaration. This is more or less useless as the absence of macro or smacro symbols indicates that a procedure is of type expr - the default Standard LISP procedure type.

2.3 Evaluation model, symbols, and variables

The main difference between algebraic and symbolic mode lies in the evaluation model:

• In algebraic mode a symbol stands for itself as unknown as long as no value is assigned; after an assignment it plays the role of a representative for that value just like a variable in a standard programming language:

```
1: x;
X
2: x:=y+1$
3: x;
Y + 1
```

In symbolic mode there is a clear barrier between the role of a symbol as variable of the programming language RLISP, a named item which represents some variable value, and the role to be part of an algebraic expression. If you mean the symbol x you must write ’x; without the quote tag x is considered as variable x and it will be asked for its value which is NOT ’x initially. Uninitialized variables cause bugs.

• Consequently all variables must be declared.

• In algebraic mode u := (x + 1)^2; means it compute a formula by expanding (x+1)*(x+1); if a value had been assigned to x, substitute the value for x. In symbolic mode an algebraic expression is interpreted as statement to compute a numeric value, just like in C or Pascal. So u := (x + 1)^2; in symbolic mode means: “take the value of variable x which is expected to be number, add 1, square and assign the resulting number to the variable u”.

\footnote{In this text the caret symbol ^ is used for exponentiation instead of the older FORTRAN like operator **.}
• If you want to refer to an algebraic expression as a data object, you must code it as an algebraic form (see below) and mark it as constant by a quote character. The only constants which don’t need a quote are numbers and strings. Example:

\[ u := \text{'(expt (plus x 1) 2)}; \]

assigns the (algebraic) expression \((x + 1)^2\) to the variable \(u\).

• algebraic mode implicitly supports standard arithmetic and algebraic evaluation for mathematical expressions of arbitrary complexity and for numbers from various domains. In symbolic mode, arithmetic with infix operators \(+\, -\, \ast\, \wedge\, /\) is supported only for the basic LISP numbers (mainly integers). All arithmetic for formulas, even for domain elements such as rounded numbers, complex numbers etc., has to be performed by calling explicitly the functions which can do that job or by calling explicitly the REDUCE evaluator \text{reval resp aeval} (see below).

So symbolic mode programs are much more similar to programs in conventional programming languages such as Pascal or C. All algebraic functionality of REDUCE is available only as an explicit subroutine interface. In contrast to algebraic mode nothing is done implicitly for you.

2.4 Variable types

RLISP supports in symbolic mode the following classes of variables:

• Local variables in a procedure are those
  – declared in the parameter list,
  – declared in a scalar or integer statement in a begin-end block,
  – bound in the right-hand side of a where statement
  – implicitly introduced in a for statement.

These are valid only inside their local context. scalar variables are initialized to nil, integer to 0 unless they have a special initialization value (property \text{initvalue})*.

\begin{verbatim}
  symbolic procedure my_adder(x,y);
  begin integer r;
  \end{verbatim}
\r:=x+y;
\text{return } r:
\text{end;}

In this routine \(x, y\) and \(r\) are local variables.

In algebraic mode the \text{where} statement is used to activate one more rule locally. In symbolic mode only rules of the form \text{<name>=<value>} are allowed on the right-hand side of a \text{where} statement.

- \textbf{Global} variables have to be declared in a statement \text{global'}(v_1 v_2 \cdots); These can be accessed from everywhere once they have been declared. Important: declare them before you use them for the first time in a session. They are initially set to \text{nil}.

\text{global } '(my\_support!* my\_quotient!*);

It is a common practice to use a trailing asterisk in names for global and fluid variables such that they easily can be distinguished from locals. Names of global variables may not be used as local variables.

- \textbf{Fluid} variables are similar to global variables as they can be accessed from everywhere. But in contrast to globals a \textbf{fluid} variable can occur as a local variable in a procedure. It then has temporarily the value assigned in the procedure (we say “it is rebound”); this value will be accessible globally by nested procedures as long as the rebinding procedure is not yet terminated. After termination the previous value is re-assigned. Fluid variables are declared in a statement \text{fluid'}(v_1 v_2 \cdots);. Like global variables they must be declared before use and they are initially set to \text{nil}.

\text{fluid } '(my\_var!*);
my\_var!*:= 'x;
\text{procedure compute(ex,my\_var!*);}
\text{do\_some\_work(ex);} 

In this case the variable \text{my\_var*} has been initialized to the symbol \(x\). If then a call \text{compute}'(a,z) is performed, the variable \text{my\_var*} will be set to \(z\) during the execution of the body of \text{compute}. 
• A **switch** can be declared using the **switch** statement `switch s_1, s_2, \ldots;` where \( s_1, s_2, \ldots \) are symbols. REDUCE automatically connects a fluid variable with each switch by prefixing an asterisk to the symbol name. This variable represents the *value* of the switch, which is either *nil* or *t* and is set by the statements **on** and **off**. E.g. the switch *nat* has the associated global variable \(*\text{nat}.*\)

• **share** variables are also global variables which are handled specially by the REDUCE evaluator such that their symbolic mode values are identical with their role as symbols in algebraic mode.

### 2.5 Symbols, numbers and strings

A **symbol** or **id**[^2] in LISP is a unit which has a name, in the standard case beginning with a letter and followed by letters and digits. Special characters in a name or a leading digit have to be flagged by a preceding exclamation mark. Symbols are the same entities as in algebraic mode. Symbols are unique in the system. You may view a symbol as a data structure with four slots, the **name** with is always a string, the **value cell** to store assigned values, the **function cell**[^3] to point to an assigned program structure and the **property cell**[^4] as anchor of the property list. Initially all cells except the name cell are marked empty.

In contrast to algebraic mode in LISP, only integers and floating point numbers (which don’t exist in that form in algebraic mode) are considered as numbers. All other numeric quantities from algebraic mode such as rationals, rounded numbers, Gaussian integers are composite data structures encoded in the REDUCE domain structure (see below).

Just as in algebraic mode, a sequence of characters enclosed in string quotes is a string. Functions for manipulating symbols, numbers and strings are[^5]

[^2]: “id” is an abbreviation for “identifier”.
[^3]: There are LISP systems which support only one cell for value and function – in such systems a symbol can represent only a variable or a function exclusively.
[^4]: Read the trailing \( p \) in names like “idp”, “numberp” etc. as “predicate”, e.g. “numberp” meaning “number predicate”.

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The functions gensym and compress make symbols which are not yet “internal”: they are unique, but there may be other symbols with the same name. It makes debugging especially hard if you cannot access the symbol by its name of course. The function intern can be used to convert such a symbol into an internal one - then every reference to the symbol name guarantees access to the desired object.

2.6 Boolean values

There are two special symbols named nil and t. nil is used at the same time for the empty list and the Boolean value “false”. Any value other than nil is considered as “true” in conditional statements. In order to facilitate programming, the symbol nil has the preassigned constant value nil such that a quote here is superfluous. And for the same reason the symbol t evaluates to itself - this symbol can be used if the Boolean value “true” is needed explicitly, e.g. as a return value. But please keep in mind that any non-nil value can play this role. In contrast to algebraic mode the number zero does not represent the Boolean value “false”.

2.6.1 Pairs, lists

The central structure of LISP is the pair. The external representation of a pair of two objects is (obj1, obj2). Here the infix dot is used as print symbol as well as infix constructor. It is often useful to use a double box for representing a pair graphically, e.g. for (10, 20):

```
    ,---------
   |   10    |
   |,--------|
```

Functions:
(infix) construct pair \((o_1, o_2)\)

\(\text{cons}(o_1, o_2)\) same as \(01.02\)

\(\text{pairp}(q)\) true if \(q\) is a pair

\(\text{car}(q)\) extract the first object from a pair

\(\text{cdr}(q)\) extract the second object from a pair

The pair is a very general construct - since its objects themselves can be pairs, arbitrary trees and data structures can be built. E.g. \(((a.b).(c.d))\)

```
\---------
| *     | * |
/--------\
/            \
/\            \
| a | b | c | d |
\--------' \--------'
```

On the other hand, structures with many members lead to a very complicated representation if printed using the above notation. Therefore the **list** concept has been introduced which allows one to use a special class of pair trees in a simplified form:

- in a pair with second element *nil* the dot and the nil are omitted: \((A.nil) = (A)\).

- for a pair with another pair as second element one bracket pair and the dot are omitted: \((B.(A)) = (B A), (C.(B A)) = (C B A)\).

So the **list** is a linear sequence of pairs, where the **cars** contain the “data” items, while the **cdrs** contain the reference to the successors. The last successor is *nil*, which signals the end of the linear list. For the graphical representation of linear lists horizontally aligned double boxes are useful, e.g. for \((C B A)\)

```
\---------
| *     | *+->| b | *+->| a |nil|
\--------' \--------'
```

or with omitting the final *nil*

```
\---------
| *     | *+->| b | *+->| a | / |
\--------' \--------'
```
The list notation is much simpler than the dot notation - unfortunately the dots cannot be avoided completely because pairs with id’s or numbers in the cdr part occur and they play an important role.

Lists can be nested; an example is the internal representation of algebraic forms (see below) which uses linear lists; e.g. \((x + 1)^2\) would be represented by lists as

\[
\text{((expt (plus x 1) 2)}
\]

Here the top level list has three elements \(\text{expt}\), \((\text{plus } x 1)\) and 2 where the second element itself is a linear list of length 3.

\[
\begin{array}{c}
\text{expt} \\
\text{plus} \\
\text{empty list}
\end{array}
\]

In this graph the cars contain the symbols or numbers or they point to substructures (downwards) while the cdrs point to the successors (to the right) until the end is reached.

As mentioned already there can be mixtures of both forms if one cdr is not nil: \((A(B.C)D.E)\) - here the second element of the list is a pair \((A.B)\), the third element is \(D\) and there is no fourth element - the list ends with a non nil symbol.

\[
\begin{array}{c}
A \\
B \\
empty list
\end{array}
\]

The empty list () is identical to the symbol nil. Note that the REDUCE algebraic forms are LISP lists where all lists are terminated by a nil.

Functions on lists:
| {o1,o2,..on} | construct list (o1 o2 ⋯ on) |
| list(o1,o2,..on) | same as {o1,o2,..on} |
| o1.l1 | (infix) put o1 in front of list l1 |
| pairp(q) | true if q is a pair (or a list) |
| atom(q) | true if q is NOT a pair/list |
| null(q) | true if q = nil (=empty list) |
| car(q) | extract the first object from a list |
| cdr(q) | extract the part of the list behind 1st element |
| nth(q,n) | extract the n-th element from list q |
| length(q) | count the number of elements of list q |
| reverse(q) | reverse a list q |
| append(a,b) | append b to the end of a copy of a |
| member(a,l) | test if a is equal to one member of l |
| memq(a,l) | test if a is identical one member of l |
| delete(a,l) | delete a from list l |

Remarks:

- All of the above functions are non destructive: an input list remains as it has been before - the value of each access function is a reference to the corresponding part of the structure.

- The access functions can operate only if the desired elements are really there; e.g. if nth with 4 is applied to a 3 element list produces an error, and car or cdr applied to any symbol or number (including nil) will fail.

- Nested calls of car and cdr can be abbreviated by combining the a’s and d’s between c and r up to four inner a/d letters. E.g. cadr(u) = car(cdr(u)) - this is the direct access to the second element, caddr returns the third element and cadddr the fourth element.

- Although the functions first, second, third, rest known from algebraic mode operate in some REDUCE implementations in symbolic mode as synonyms of car, cadr, caddr, cdr, they should not be used here as this is the case in all REDUCE versions.

- The functions member and memq not only test the membership: if a is member of the list l, member (or memq) returns the (first) pair which contains a as its car. E.g. memq('b,'(a b c)) returns (b c). This can be used either if you want to replace this item in the list, or if you want to process the part after the search item. The function memq
looks for a member of the list which is eq to the object, while member uses equal as test (see discussion of equality below). If applicable, memq is substantially faster than member.

- **Delete** produces a copy of the top level pair sequence leaving out the first occurrence of the search item: the original list is not changed, and a possible second occurrence of the search item in the list is not removed. delete('a,' (0 a b a)) will result in (0 b a) where the part (b a) is the original tail of the input list while the pairs in front of the original first a are new.

Examples: let u be a variable with value (A(B.C)NIL(D)). Then

\[
\begin{align*}
\text{pairp}(u) &= t \\
\text{length}(u) &= 4 \\
\text{null}(u) &= \text{nil} \\
\text{car}(u) &= A \\
\text{cadr}(u) &= (B.C) \\
\text{caadr}(u) &= B \\
\text{cdadr}(u) &= C \\
\text{cdr}(u) &= ((B.C) \text{NIL} (D)) \\
\text{length cdr}(u) &= 3 \\
\text{cddr}(u) &= (\text{nil} (D)) \\
\text{null cddr}(u) &= \text{nil} \\
\text{caddr}(u) &= \text{nil} \\
\text{null caddr}(u) &= t \\
\text{cdddr}(u) &= ((D)) \\
\text{cddddr}(u) &= \text{nil} \\
\text{null cddddr}(u) &= t
\end{align*}
\]

All data which are not pairs (or lists) are classified as atoms: symbols, numbers, strings, vectors(see below) and, just as in real world, they are not very atomic – there are tools to crack them.

### 2.7 Programming with lists

There are several methods available to construct a list:

\[
\begin{align*}
u &:= \text{nil}; \\
u &:= 1 \cdot u; \\
u &:= 2 \cdot u;
\end{align*}
\]
Starting with an empty list, elements are pushed on its front one after the other, here giving (2 1). Please note the blank between the number and the dot: this is necessary as otherwise the dot would have been parsed as part of the number representing a floating point value.

\[
u := \{9,10\};
\]
\[
u := 1 . u;
\]
\[
u := 2 . u;
\]

The same, here not starting with the empty list giving (2 1 9 10). The \texttt{for statement} has special forms to handle lists:

- \texttt{for each x in l \ldots} performs an action elementwise for the elements in list \( l \); the elements are accessible by the automatically declared local variable \( x \) one after the other.

- the iterator actions \texttt{collect} and \texttt{join} allow you to construct lists:
  
  - \texttt{collect} generates a list where each iteration step produces exactly one element,
  
  - \texttt{join} expects that each iteration step produces a complete list (or \texttt{nil}) and joins them to one long list.

Examples: let e.g. \( l \) be a list of numbers (1 − 2 3 − 4 5 − 6):

\[
m := \text{for each } x \text{ in } l \text{ sum } x;
\]

\( m \) will be the sum of the elements,

\[
s := \text{for each } x \text{ in } l \text{ collect } x^2;
\]

\( s \) is computed as list with the numbers from \( x \) squared,

\[
p := \text{for each } x \text{ in } l \text{ join}
\]
\[
   \text{if } x > 0 \text{ then } \{x\} \text{ else nil};
\]

in this loop the positive numbers produce one element lists, while the negative numbers are mapped to nil; joined together the result is a list with only the positive elements,

\[
r := \text{for } i := 1:10 \text{ collect } -i;
\]
here the result is the list with numbers \((-1 - 2 - 3 \cdots - 10)\). The lists produced by \textit{collect} and \textit{join} are in “correct” order. \textbf{Warning:} In symbolic mode \textit{join} links the lists \textbf{in place} for maximum speed\(^5\) The last \textit{CDR} pointer of the first list is modified to point to the head of the second list, and so on. This is no problem if the joined objects are freshly built structures, or if there are no other references to them. If the in-place operation is dangerous for your application, you must copy the structure before, e.g. by a call to \textit{append} with \textit{NIL} as second argument:

\[
r:=\text{for } i:=1:10 \text{ join append}(\text{my extract}(u,r),\text{nil});
\]

In this example, the top level chain of the results of \textit{my extract} is copied such that the original structure is not modified when joining.

Another very common style for list operations is to use iterative or recursive programs. The following programs each sum the elements of the list:

\[
\text{symbolic procedure sum}_1(l);
\begin{align*}
&\text{begin } \text{integer } n; \\
&\quad \text{while } l \text{ do } <n:=n+\text{car } l; \ l:=\text{cdr } l>>; \\
&\quad \text{return } n; \\
&\text{end;}
\end{align*}
\]

This program picks in each step of the the while loop one element from the list and sets the variable \(l\) to the successor position; the loop terminates as soon as the successor \textit{nil} is found .

\[
\text{symbolic procedure sum}_2(l);
\begin{align*}
&\text{if null } l \text{ then } 0 \text{ else car } l + \text{sum}_2(\text{cdr } l);
\end{align*}
\]

This program uses recursion for the same task. The program will go down the list until \textit{nil} is found, then set the sum to 0 and when coming back from the recursion add the elements to the sum so far.

\[
\text{symbolic procedure sum}_3(l); \text{sum}_31(l,0); \ % \text{initializing} \\
\text{symbolic procedure sum}_31(l,n); \\
\text{if null } l \text{ then } n \text{ else } \text{sum}_31(\text{cdr } l,\text{car } l+n);
\]

The third example is a bit tricky: it initializes the sum to 0, and when stepping down in the recursion the sum is updated and passed to the next recursion level as a parameter.

\(^5\)In algebraic mode \textit{join} operates different: the operatend are copied before linking them to one result list

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It is also very common to produce lists in recursive style, e.g. inverting the signs in a list of numbers could be written as

```
symbolic procedure invertsgn(l);
    if null l then nil else -car l . invertsgn(cdr l);
```

and with most LISP compilers this program would as efficient as the corresponding loop

```
for each x in l collect -x;
```

but produce less code. Note that as with `for each - collect`, the elements in the resulting list will be in the original order. The recursive programming style is typical for LISP and it is used for hundreds of REDUCE routines.

### 2.8 In-place operations

All list construction described so far except `join` is of a copying nature: every construction by the dot or the curly brackets always creates fresh pairs and an old list is untouched. However, there are operations to modify elements or the structure of an existing list in place. These should be used only with greatest care: if there is another reference to the list, its structure will be modified too, a possibly unwanted side effect. The functions are:

<table>
<thead>
<tr>
<th>Function</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>car l := u</code></td>
<td>replace first element of <code>l</code> by <code>u</code></td>
</tr>
<tr>
<td><code>rplaca(l,u)</code></td>
<td>old form for <code>car l := u</code></td>
</tr>
<tr>
<td><code>cdr l := u</code></td>
<td>replace cdr part of <code>l</code> by <code>u</code></td>
</tr>
<tr>
<td><code>rplacd(l,u)</code></td>
<td>old form for <code>cdr l := u</code></td>
</tr>
<tr>
<td><code>nth(l,j) := u</code></td>
<td>replace the <code>j</code>-th element of <code>l</code> by <code>u</code></td>
</tr>
<tr>
<td><code>nconc(a,b)</code></td>
<td>insert <code>b</code> as cdr in the last pair of <code>a</code></td>
</tr>
<tr>
<td><code>reversip(l)</code></td>
<td>invert the sequence of <code>l</code> using the same pairs again</td>
</tr>
</tbody>
</table>

### 2.9 Equal tests: Equality vs. Identity

There are two classes of equality in LISP, `eq` and `equal`, in REDUCE written as the infix operators `eq`, and equal sign `=` respectively. As symbols are unique, both are equivalent for them: if symbols are `equal`, they are `eq`:

```
u :=' a; v :=' a; u eq v ⇒ t
```

here the variables `u` and `v` represent the same symbol and consequently they are `eq`.

---

6Read “reversip” as “reverse-in-place” and “nconc” as “concatenate”
For lists, things are different. Here we must distinguish between references to pairs which contain the same contents ("equal") or references to the same identical pair instance("eq"). A pair is eq to itself only, but it is equal to a different pair if their cars andcdrs are equal - this is a recursive definition. Example:

\[
\begin{align*}
u &:= \{1, 2\}; v := \{1, 2\}; w := u; \\
u &= v \Rightarrow t, u = w \Rightarrow t \\
u eq v &\Rightarrow nil, u eq w \Rightarrow t
\end{align*}
\]

Here u, v and w have as values pairs with same cars and equal cdrs, but only u and w refer to the identical pair because for v fresh pairs have been built. In some points REDUCE relies on identity of some structures, e.g. for composite kernels of standard forms.

2.10 Properties, Flags

Another important concept of LISP is the ability to assign properties and flags to each symbol. A property is a symbol (a “name”) together with a value. Both, property name and property value are completely free. Properties are set by the function put and read by the function get.

\[
\begin{align*}
\text{put('hugo,'evaluator,'compute_hugo);} \\
\ldots \\
\text{get('hugo,'evaluator)} &\Rightarrow \text{compute_hugo} \\
\text{get('otto,'evaluator)} &\Rightarrow \text{nil}
\end{align*}
\]

A property-value pair remains assigned for a symbol until the value is replaced by a different value or until the property is removed by remprop.

A flag is similar to a property, but it represents only the values “yes” or “no” (= flag is present or absent). The functions for setting and testing flags are flag and flagp, remflag for removal.

As most other LISP programs, REDUCE makes extensive use of properties. E.g. the algebraic properties of an operator are encoded in that way. So the symbol plus has among others the REDUCE properties prtch = + (the print character), \textit{infix} = 26\footnote{This number may change if additional operators of lower precedence are introduced} (the precedence among the infix operators), \textit{smpfn} = simpplus (the name of the routine responsible for its simplification) and the flags symmetric, realvalued, nary and some more. These allow a data driven programming model which has been used in the LISP world for decades already: if the simplifier operates on an expression, it asks the leading operator for its simplification function; if there is one (and there will be one for all operators), this function is invoked to do the job.
There are various fields where this type of programming is used inside REDUCE, e.g. the complete domain arithmetic has been implemented in that way: any non-integer domain element is encoded as a list with a leading symbol (the domain mode) which contains all knowledge about arithmetic, conversion, printing etc. for this domain as properties and flags.

If your REDUCE system is based on PSL, you can inspect a property list by calling the function prop which has one argument (the symbol to be inspected) and which returns the properties and flags as list. E.g.

```reduce
a := mat((1,2),(3,4));
prop 'a;
prop 'mat;
prop 'matrix;
```

### 2.11 Alists

Another frequently used LISP structure is the association list (alist), which is an associative memory: a list where each element is a pair with a search item and a value item. The function assoc picks the pair for a given search item out of the alist. For example constructing a multiplier with memory:

```reduce
c> fluid '(product_database*);
c> symbolic procedure multiply(u,v);
c> begin scalar w,r;
c> w:=u.v;
c> r:=assoc(w,product_database*);
c> if r then return cdr r; % found
+c> r:=u*v;
c> product_database*:=(w.r).product_database*;
c> return r;
c> end;
```

here the pair u.v is used as search item. If an entry is found in the database, its value part is returned. Otherwise the product is computed and the new entry is added to the database. After calling this function with arguments 2 and 3, and then with 2 and 6 the structure would be

```
+-----------------------.
|   |   |   |   |   |   |
| * | * | +-----| * | / | |
| - | ----- | - | ----- |
```

17
Here each element in the top level list points to one generalized name-value pair, where the name itself is a structure (here a pair of two numbers). REDUCE uses association lists for many different purposes; e.g. if an expression is simplified, the expression and its simplified variant are stored in an association list such that for a second access to the same expression its simplified form is immediately accessible.

2.12 Vectors

Lists enable access of items of a linear collection by position number using the function \texttt{nth}; however, this access is inefficient if used frequently because for every access the sequence of pairs has to be traversed until the desired object is reached. An alternative structure for direct access by element number in symbolic mode is the \texttt{vector}. The first element in a vector has the index 0.

\begin{verbatim}
v:=mkvect(n)  create a vector of length n
putv(v,m,u)   put value u into mth slot of vector v
getv(v,m)     extract element m from vector v
upbv(v)       returns upper bound for v
\end{verbatim}

Note that \texttt{putv} is a “destructive” operation.

3 Algebraic data formats and evaluation

3.1 Prefix forms

The basic format for representing mathematical formulae is the algebraic form: an expression is internally encoded as a list with prefix operators:

- the first element of a list is an operator,
- the following elements of a list are the arguments which can be symbols (unknowns), numbers or algebraic forms in prefix notation.
Examples:
\[ x + 1 \Rightarrow (PLUS X 1) \]
\[ x + y + 1 \Rightarrow (PLUS X Y 1) \]
\[ x + y * z + 1 \Rightarrow (PLUS X (TIMES Y Z) 1) \]
\[ x^2(y + 1) \Rightarrow (EXPT X (PLUS Y 1)) \]

Algebraic forms are used for many purposes, for data input, for transferring data between various system components and for output. To get a feel as to how algebraic forms look like, you can make REDUCE display them for you, e.g. by the following sequence:

\[
\begin{align*}
  u &:= (x+y)^3/(\log z-sqrt 2); \\
  \text{symbolic reval 'u;} 
\end{align*}
\]

The first statement assigns an algebraic expression to a variable as usual in algebraic mode, and the second statement forces REDUCE to print the algebraic form corresponding to this expression in symbolic mode.

3.2 *SQ

REDUCE uses internally a different representation for algebraic expressions, the standard quotient(SQ, see below). As conversion between algebraic forms and standard quotients is expensive, REDUCE tries to keep the data as long in SQ form as possible. For this purpose there is a special variant of the algebraic form (*SQ u ind): here the internal operator *SQ indicates that the parameter u is not an algebraic form but a standard quotient. If such expression is evaluated, REDUCE detects immediately that it will find the standard quotient already as the second element of the list and no more conversion is necessary.

However, if since creation of this form some values in the system have been changed, the form u might be no longer the “correct” standard quotient; e.g. one of its unknowns might have a value now. For this purpose the third element ind in the list has been introduced: As long as the form u is the correct one, it will be T. As soon as a significant value in the system changes which might affect the validity of u, REDUCE switches all ind elements in the *SQ forms swimming around to nil. So the evaluator decides from the second parameter whether to use u unchanged or to re-evaluate.

By the way, the global replacement of the second parameter in the *SQ forms is a good example of an in-place list operation with a wanted side effect: REDUCE maintains a unique list *sqvar* = (t) and all *SQ forms are built with this list as last pair. If the environment changes it is sufficient
to replace this single \( t \) in \( *sqvar* \) by \( nil \) and all \( *SQ \) forms built with the actual \( *sqvar*-tail \) inherit this automatically switching from \( t \) to \( nil \); REDUCE then creates a fresh list \( *sqvar* = (t) \) to be used for the next generation of \( *SQ \) forms.

### 3.3 Composite forms

#### 3.3.1 Lists, equations

For algebraic mode, lists are represented by the prefix operator \( LIST \). From the symbolic mode standpoint it may appear a bit curious that one has to insert the symbol \( LIST \) into a list in order to make it a correct algebraic object, but this is necessary in the prefix algebraic context. E.g. try in algebraic mode

```lisp
u:={a,b,{c,d},e};
lisp reval 'u;
```

you will see the result \( (LIST A B (LIST C D) E) \).

Correspondingly equations are represented by the prefix operator \( EQUAL \) followed by the right-hand and the left-hand sides as parameters; e.g. \( x = \sqrt{2} \) is represented as \( (EQUAL X (SQRT 2)); \)

The result of \( \text{solve} \ (x^2=2,x); \) is an instructive mixture of all these forms:

```lisp
(LIST (EQUAL X (!*SQ (((((EXPT 2 (QUOTIENT 1 2)). 1). -1)). 1) T))
     (EQUAL X (!*SQ (((((EXPT 2 (QUOTIENT 1 2)). 1). 1)). 1) T)))
```

this form is a list with two equations and the equations have \( *SQ \) forms as right-hand sides.

#### 3.3.2 Matrices

A matrix too is represented internally by a prefix form with operator \( MAT \) followed by the rows; each row is a list containing the elements as algebraic forms. The elements can be accessed easily using the access function \( nth \). Example: the matrix \( \text{mat}((1,2),(3,4)) \) is encoded as \( (MAT (1 2) (3 4)) \). The form \( nth(nth(m,3),2) \) gives access to the element \( (2,2) \). Be careful if you use this form for an assignment: \( nth(nth(m,3),2) := 17; \) is an in-place operation. You can create a matrix easily by a symbolic program, e.g. a \( n*n \) unit matrix
symbolic procedure my_unit_matrix(n);

'mat . for i:=1:n collect
  for j:=1:n collect if i=j then 1 else 0;

3.4 Algebraic evaluator: reval, aeval, aeval*

The main entries to the algebraic evaluator are the symbolic mode functions reval and aeval. Both have as input an algebraic expression and return the fully evaluated result. Reval always returns an algebraic form in full prefix operator form, while aeval tries to keep the result in *SQ-form whenever possible. Almost all algebraic functionality of REDUCE can be invoked by calling these procedures, e.g.

reval '(solve (plus (expt x 2) 3) x);

If you turn on the switch defn in algebraic mode you will see that REDUCE translates most statements into calls for aeval.

Reval and aeval are also adequate tools for accessing the values of algebraic quantities and for evaluation of unevaluated procedure parameters; they do the complete job for you.

However, there is one exception: the left-hand side of an equation in general is not evaluated by reval and aeval. E.g.

l := a=b;
a := 17; b := 4;
symbolic reval 'l;

produces the output (EQUAL A 4). If you write a program for equations you either should evaluate the left-hand sides yourself by calling reval or rebind the internal switch *evallhseqp locally(!) by t - then REDUCE will evaluate both sides of an equation.

REDUCE maintains an internal buffer alglist* where all evaluation results are stored such that a repeated request for the same evaluation can be satisfied by a simple and fast table lookup. As a side effect, aeval resets alglist* empty before starting a computation. There is a second function

---

8 Read “evallhseqp” as “evaluate left-hand side of equations predicate”.
9 The use of alglist* is controlled by an internal switch *uncached: if this variable has a non-nil value, the caching is suppressed.
10 Of course, there are many more places where alglist* needs to be reset; e.g. any assignment causes alglist* to be reset because the assigned value might make one of the stored values invalid.
aeval*: this function performs the same operation, but in contrast to aeval it does not reset alglist* and can profit from previous results. The REDUCE uses both, aeval and aeval* depending of the specific environment of a translated statement.

4 Standard Forms, Standard Quotients

4.1 Introduction

One of the most important tasks performed by a computer algebra system is the reduction of an algebraic expression to a canonical form. A reduction process is canonical if it guarantees that two expressions are transformed to the same form if and only if they are algebraically identical. Unfortunately canonical forms are not known for all algebraic objects, e.g. for surds. The canonical form of REDUCE is the standard quotient (“SQ”) which is a quotient of two standard forms (“SF”). An SF is a polynomial in internal representation and an SQ is a quotient of two polynomials, a rational function. At the same time SQ and SF are the forms used internally for most arithmetic.

In this section the structures of SF and SQ are described, followed by a list of operations on these. For programming and debugging in symbolic mode it is often necessary to decipher an expression of this type. However, the knowledge about this part of the internal REDUCE structure should never be used to access SF of SQ directly by LISP primitives. Instead use only “official” primitives supplied by REDUCE.

4.2 Coefficients

The coefficients in SF and SQ are the “numbers” from various algebraic contexts. They are called domain elements and uniformly represented either as integer numbers or pairs where the car part contains an identifier describing the domain and its arithmetic by its properties; the cdr part describes the value.

1. Integers: integer numbers ∈ Z represent integer coefficients directly,

2. Gaussian integers: complex numbers with integer real and imaginary part (:gi: re . im) with re, im ∈ Z,

3. Rational numbers: quotients of integers (:rn: nu . de) with nu, de ∈ Z,
4. Modular numbers: numbers modulo the current prime $(\mod: . j)$ where $0 \leq j < \text{currentmodulus}$.\[11\]

5. Rounded numbers: floating point numbers extended in precision and magnitude $(\rd: . fp)$ with LISP floating point number $fp$ or $(\rd: \text{mant} . \text{exp})$ with integers $\text{mant}$ and $\text{exp}$ representing an extended floating point value.

6. Rounded complex numbers: numbers with rounded real and imaginary parts $(\gr: \text{fre} . \text{fim})$ with LISP floating point numbers $\text{fre}$ and $\text{fim}$ or $(\gr: (\text{rmant} . \text{rexp}) . (\text{imant} . \text{iexp}))$ with extended real and imaginary parts.

The neutral elements $\text{zero}$ and $\text{one}$ for addition and multiplication respectively are identical for all domains: $\text{zero}$ is represented as $\text{nil}$ and $\text{one}$ by the integer number 1.

The list of domains is open ended: if a new coefficient domain is required for an application field, this can be created and linked to the system at any time. The necessary structure is described in the source file for the module $dmodeop$.

The tag of the currently active domain is stored in the value of the variable $\text{dmode*}$. $\text{nil}$ represents the default domain (integers). This variable should not be modified directly - instead use the function $\text{setdmode}$:

$$\text{setdmode}(\text{DOM:domain}, \text{MODE:bool});$$

where $\text{domain}$ is a domain name (e.g. 'rational) and the Boolean second parameter tells whether to activate ($t$) or to deactivate ($\text{nil}$) the domain. Note that $\text{setdmode}$ does not set on the corresponding switches. You must do that yourself. E.g. for a proper activation and deactivation of complex rounded call

```lisp
!*rounded := t;
setdmode('rounded,t);
!*complex := t;
setdmode('complex,t);
.... % do your computation
setdmode('complex,nil);
```

[11] The upper half of the modular numbers are printed as negative numbers if the switch $\text{balanced_mod}$ is on; however, that does not affect the internal representation which uses always integers $\geq 0$
/*complex := nil;
setdmode('rounded,nil);
!*rounded := nil;

In order to convert the domain tag into the domain name use the property dname, e.g.

get(dmode!*,'dname);

will result in the name of the current domain.

4.3 Generic arithmetic with domain elements

There is a complete set of arithmetic functions for performing arithmetic operations with domain elements:

<table>
<thead>
<tr>
<th>Function</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>!:zerop(a)</td>
<td>test $a = 0$</td>
</tr>
<tr>
<td>!:plus(a,b)</td>
<td>$a + b$</td>
</tr>
<tr>
<td>!:difference(a,b)</td>
<td>$a - b$</td>
</tr>
<tr>
<td>!:minus(a)</td>
<td>$-a$</td>
</tr>
<tr>
<td>!:minusp(a)</td>
<td>test $a &lt; 0$</td>
</tr>
<tr>
<td>!:times(a,b)</td>
<td>$a \cdot b$</td>
</tr>
<tr>
<td>!:quotient(a,b)</td>
<td>$a/b$</td>
</tr>
<tr>
<td>!:recip(a)</td>
<td>$1/b$</td>
</tr>
<tr>
<td>!:expt(a,n)</td>
<td>$a^n$</td>
</tr>
<tr>
<td>!:gcd(a,b)</td>
<td>greatest common divisor of $(a, b)$</td>
</tr>
</tbody>
</table>

4.4 Kernels, kernel order

Kernels are the “unknowns” in a polynomial. A kernel can be an

- identifier: e.g. $x, \alpha$, represented as LISP symbols $X, ALPHA$,
- operator form: e.g. $\sin(\alpha),\sqrt{x + 1},\text{bessel}(j,x)$, represented as
  algebraic expression $\text{(SINALPHA)},(\text{EXPT (PLUS X 1) (QUOTIENT 1 2))},$
  $\text{(BESSEL J X)}$.
- A standard form, e.g. $\text{(((A . 1) . 1) . 1)}$ This option is used only if the
default mode of full expanions is turned off (see below: non-expanded standard form).

It is essential for the operation of REDUCE that kernels are unique. This is
no problem for id’s as these are unique by LISP definition. For composite
kernels REDUCE ensures that these are made unique before they enter a **standard form**. Never construct a composite kernel by hand. Instead use the function *a2k for converting an algebraic expression into a kernel or use simp (see below). On the other hand, the uniqueness of kernels is of big advantage for comparisons: the equality can be tested with eq instead of = and faster memq and atsoc can be used instead of member and assoc.

The set of all kernels is totally ordered by the kernel order. REDUCE has by default some ordering which is roughly based on lexicographic comparison. The ordering can be changed internally by calling setkorder (corresponding to the algebraic statement korder) with a list of kernels as argument. These kernels precede all other kernels in the given sequence. Setkorder returns as result the previous order and it is a good policy to restore the old order after you have changed it:

```reduce
begin scalar oldorder, ....;
    oldorder:=setkorder '(x y z);
    ....
    setkorder oldorder;
    ....
end;
```

The current kernel order list is the value of the variable kord*. The value nil signals that the default ordering is active.

To check the system order of two algebraic expressions use the function ordp with two arguments, which returns true if the first argument is lower in the ordering sequence as the second one. This check does not consider the actual kord*. For kernels there is an extended order check function ordop which does consider the actual kord*. However, you need to call these functions only rarely as the algebraic engine of REDUCE maintains a proper order in any context.

### 4.5 Standard form structure

A **standard form** is a polynomial in internal recursive representation where the kernel order defines the recursive structure. E.g. with kernel order (x y) the polynomial $x^2y + x^2 + 2xy + y^2 + x + 3$ is represented as a polynomial in $x$ with coefficients which are polynomials in $y$ and integer coefficients: $x^2 * (y * 1 + 1) + x * (y * 2 + 1) + (y^2 * 1 + 3)$; for better correspondence with the internal representation here the integer coefficients are in the trailing position and the trivial coefficients 1 are included. A standard form is

- **<domain element>**
• \(<\text{mvar}>.**<\text{ldeg}>.*<\text{lc}>.+<\text{red}>,\) internally represented as 
\(((\text{mvar}\.\text{ldeg}\.\text{lc}).\text{red})\)

with the components

• \text{mvar}: the leading kernel,

• \text{ldeg}: an integer representing the power of \text{mvar} in the leading term,

• \text{lc}: the leading coefficient is itself a standard form, not containing \text{mvar} any more,

• \text{red}: the reductum too is a standard form, containing \text{mvar} at most in powers below \text{ldeg}.

Note that any standard form ends with a domain element which is \text{nil} if there is no constant term. E.g. the above polynomial will be represented internally by

\(((X\.2)((Y\.1).1)\.1)((X\.1)((Y\.1)\.2))((Y\.2)\.1)\.3)\)

with the components

• \text{mvar}: X

• \text{ldeg}: 2

• \text{lc}: ((Y\.1).1)\.1 = (y + 1)

• \text{red}: ((((X\.1)((Y\.1).2))((Y\.2).1))\.3) = xy^2 + 3.

4.6 Operations on standard forms

Arithmetic with standard forms is performed by the functions\footnote{Be careful not to intermix the names “minusf” and “negf”. And take into account that \text{minusf} will return a positive result only if the value is definitely known to be less that 0.}
null(f)  test $f = 0$
minusf(f)  test $f < 0$ (formally)
domainp(f)  test if $f$ is a domain element (e.g. number)
addf(f1,f2)  $f_1 + f_2$
negf(f)  $-f$
addf(f1,negf f2)  $f_1-f_2$
multf(f1,f2)  $f_1*f_2$
mksp**(f1,n)  compute $f_1^{**n}$
quotf(f1,f2)  $f_1/f_2$ if divisible, else nil
quotfx(f1,f2)  $f_1/f_2$ divisibility assumend, no check
qremf(f1,f2)  pair (quotient.remainder)

As standard forms are a superset of domain elements these functions also can be used for arithmetic with domain elements or for combining standard forms with integer numbers. They can even be used to compute with integers as long as these are regarded as “degenerate” polynomials.

For access of the components of a standard form there are functions

| mvar(f) | extract leading kernel |
| ldeg(f) | extract leading degree |
| lc(f)  | extract leading coefficient |
| red(f) | extract reductum $f$ |

These functions can be applied only if the form is not a domain element. The following example presents a typical function evaluating a standard form; it looks for the maximum power of a specified kernel $y$ in a standard form $f$:

```
symbolic procedure my_maxpow(f,y);
    if domainp f then 0 else
        if mvar f eq y then ldeg f else
            max(my_maxpow(lc f,d),my_maxpow(red f,d));
```

This function has a double recursion which is typically for functions traversing standard forms: if the top node is not trivial (domainp) or not a power of $y$ the procedure branches to the coefficient form and to the reductum form - both can contain references to the kernel $y$. Kernels are unique in the system and therefore the equality test is performed by eq.

For the initial construction of a standard form a safe and efficient way is to build an algebraic expression first and convert it using simp (see below) or to use the arithmetic functions above too. For elementary conversions use
Note that integers are already an SF and so need not be converted.

Although the are infix macros \( \times \), \( \div \), \( + \) to construct standard forms, they should be avoided unless you are completely sure what you are doing: they construct without any consistency checks, and wrong structures may lead to inconsistent results; such cases are difficult to debug.

### 4.7 Reordering

If you change the kernel order it is necessary to reorder standard forms if they are processed under the new order by calling

\[
\text{reorder}(f) \quad \text{reorders } f \text{ to the current order}
\]

Please ensure that you never use a standard form which does not correspond to the actual order - otherwise very strange results can occur. Freshly created standard forms will always have the current order.

### 4.8 Standard Quotients

A standard quotient is a pair of standard forms \((\text{numr} \div \text{denr})\), in LISP represented as \((\text{numr} \div \text{denr})\). If the denominator is trivial, the standard form 1 is used. The constructor, selector and conversion functions are

\[
\begin{array}{|c|c|}
\hline
\text{numr}(q) & \text{select the numerator part} \\
\text{denr}(q) & \text{select the denominator part} \\
f2q(f) & \text{convert a standard form } f \text{ to SQ} \\
simp(a) & \text{convert algebraic form } a \text{ to SQ} \\
\text{prepsq}(q) & \text{convert SQ to algebraic form} \\
\hline
\end{array}
\]

Arithmetic with standard quotients is performed by the functions

\[
\begin{array}{|c|c|}
\hline
\text{addsq}(q1,q2) & q1 + q2 \\
\text{negsq}(q) & -q \\
\text{subtrsq}(q1,q2) & q1 - q2 \\
\text{multsq}(q1,q2) & q1 * q2 \\
\text{quotsq}(q1,q2) & q1 / q2 \\
\hline
\end{array}
\]

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Note that there is no zero test for standard quotients; instead use the zero test for the numerator $\text{null}(\text{numr}(q))$.

When should you use standard quotients and when standard forms? If you are sure that no denominator other than one can occur you should use standard forms, otherwise standard quotients are the only possible choice. You can of course mix arithmetic with both, however that strategy is not recommended unless you keep track carefully; debugging is hard. Arithmetic functions have been coded for maximum efficiency: none of them test the validity of their arguments.

4.9 Substitutions

For substituting kernels in standard forms or standard quotients by other kernels or arbitrary algebraic expressions the following functions are available:

- $\text{subf}(f, a)$ substitute $a$ in SF $f$
- $\text{subsq}(q, a)$ substitute $a$ in SQ $q$

where $a$ is a list of pairs, each with a kernel to be replaced in the car and its replacement as algebraic form in the cdr. E.g.

\[
\text{uhu} := \text{simp}(\text{'(expt (plus x y) 10))};
\]
\[
\text{otto} := \text{subsq} (\text{uhu}, \text{'((x . 5) (y . (plus a b))));}
\]

Here $x$ is replaced by 5 while $y$ is replaced by $a + b$.

Note that

- both $\text{subf}$ and $\text{subsq}$ return a standard quotient (the substitution might have caused a non trivial denominator),

- $\text{subf}$ and $\text{subsq}$ also introduce substitutions which have been coded in rules - so it makes sense to call these routines with $\text{nil}$ as the second argument if you have introduced a new rule and some standard forms or standard quotients need to be re simplified under these.

4.10 Useful facilities for standard forms

4.10.1 Kernel list

The function $\text{kernels}$ extracts all kernels from a standard form and returns them as list.

[13]The second argument of $\text{subf}$ and $\text{subsq}$ is a typical association list structure.
kernels numr simp '(plus (expt x 3)(expt y 4));

will result in (x y).

4.10.2 Greatest common divisor

The greatest common divisor of two polynomials is computed by the function *gcd*. It takes as arguments two standard forms and returns a standard form or a 1 if there is no non trivial common divisor.

4.10.3 Factorization

The function fctrf is the interface for calling the factorizer. It takes a standard form as parameter and returns a list which contains

- as first element the content of the polynomial - that is the gcd of the coefficients or the “common numeric factor”,
- as second and following elements the factors in the form (SF.m) where SF is the factor as standard form and m is its multiplicity.

So calling fctrf for \(2x^2 - 2\) will result in

\[
(2 (((X . 1) . 1) . 1) . 1) (((X . 1) . 1) . -1) . 1))
\]

Here you see the common domain factor 2 and the two factors \(x + 1\) and \(x - 1\) as standard forms paired with their multiplicity 1.

4.11 Variant: non-expanded standard form

The standard form structure described so far is used by default for the fully expanded expressions with common denominator. By setting on factor, off exp, or off mcd the slots mvar and ldeg may be used in an exended style:

- ldeg can be a negative number to represent a reciprocal factor,

```lisp
(off mcd; (x*a+y*b)/(a*b);

-1   -1
a *y + b *x
```

```lisp
lisp ws;
(!*sq (((a . -1) ((y . 1) . 1)) ((b . -1) ((x . 1) . 1))) . 1) t)
```

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Here the numerator form of the standard quotient is a sum where each term is a product of one kernel with a positive and one kernel with a negative exponent.

- *mvar* may be a standard form to represent a factored polynomial,

\[
\text{on factor; } (x^2+y)^2*(y+1);
\]

\[
\begin{align*}
2 & \quad 2 \\
(x + y) & *(y + 1)
\end{align*}
\]

\[
\text{lisp ws; }
\]

\[
(!*sq (((((((x . 2) . 1) ((y . 1) . 1)) . 2) (((((y . 1) . 1) . 1) . 1) . 1) . 1))) . 1) \ t)
\]

Here the numerator is a standard form in factored form where the inner polynomials \((x + y)\) and \((y + 1)\) are used as *mvars*, the first one with the exponent 2 and the second one with exponent 1.

Special functions to work with composite standard forms:

<table>
<thead>
<tr>
<th>Function</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>sfp(m)</td>
<td>test if <em>m</em> is a standard form (T) or a kernes(NIL)</td>
</tr>
<tr>
<td>expnd(f)</td>
<td>expand a standard form <em>f</em></td>
</tr>
</tbody>
</table>

To distinguish between factored and expanded standard forms, use the predicated *sfp*; *sfp* applied to *mvar* of a standard form returns *T* if the main variable slot is occupied by a nested standard form. To expand a factored standard form you may use the function *expnd*; however, you need to turn the switch *exp* on during that call, e.g.

\[
u:=\text{expnd u where } !*\text{exp}=t;\]

Don’t build non–expanded standard forms yourself – otherwise you run the risk to produce objects with a bad structure (e.g. a wrong term order) resulting in wrong computational results. You better rely on the standard routines for manipulating these objects – as soon as *exp* is off and *factor* is on they produce the product forms anyway.

5 **Communication between algebraic and symbolic modes**

The main aim of programs written in symbolic mode is to implement the evaluation steps to be invoked from algebraic mode (top level REDUCE).
This section describes how to link symbolic routines to algebraic mode and how to exchange data between the modes.

5.1 Algebraic variables

A variable which has been declared share by the REDUCE share command is accessible from both modes in the same style and it represents the same algebraic value. A share variable is the easiest way to transfer data between the modes. The only restriction is that a symbolic program should assign only a legal algebraic form to it - otherwise a later access in algebraic mode can lead to problems. Example:

```
share hugo;
hugo :=(x+y)**2$
hugo;
symbolic;
hugo := reval {'sqrt,hugo};
algebraic;
hugo;
```

Variables which have not been declared share have different values in symbolic and algebraic mode. Nevertheless a symbolic mode program has access to the algebraic value of a symbol:

- **reval(x)**: if the value of \( x \) is a symbol or if the parameter of reval is a directly quoted symbol the function returns the algebraic value associated with this symbol, e.g. `reval('y),

- **setk(x,y)** sets the algebraic value of the expression \( x \) which then must be a symbol (or more general: a kernel) to \( y \), e.g. `setk('z, reval('plus u 17))

Of course a clever LISP programmer easily sees how REDUCE organizes the assignment and storage of values to algebraic variables in property lists, and he might be attempted to use this knowledge for “easier” access; please resist: otherwise your program might not run in a future version of REDUCE.

\[^{14}\] Read “setk” as “set kernel value”, the REDUCE equivalent to the Standard LISP assignment function “setq”.

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5.2 Calling symbolic routines from algebraic mode

REDUCE offers various ways to use symbolic routines in algebraic mode such that they appear on the algebraic level as genuine parts of REDUCE. These mainly differ in their strategy of evaluating parameters (and results).

5.2.1 Common protocol

Some general protocol rules apply for REDUCE symbolic mode routines:

- A routine should check its parameters carefully; in the case of illegal parameters the REDUCE error handlers should be invoked (see below).

- If a routine cannot process its parameters for mathematical reasons it best returns the expression which had caused the call unchanged.

- Evaluators should operate silently; events such as messages needed for testing and program development should be carried out in dependency of a switch.

- A routine should return its result as a value rather than printing it; for an isolated call in interactive mode a printed result would be sufficient, but for following evaluation steps the availability of an algebraic value is essential.

5.2.2 Symbolic operator

The simplest way to link a symbolic procedure to algebraic mode is the symbolic operator declaration. In that case the routine can be called by its proper name in algebraic mode with REDUCE first evaluating its parameters to fully simplified algebraic forms. Upon return the result should also be an algebraic form (or NIL). Example:

```reduce
symbolic procedure my_plus(a,b);
begin scalar r;
  print a; print b;
  r := reval{'plus,a,b};
  print r;
  return r;
end;
symbolic operator my_plus;
```

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This routine receives two algebraic expressions and computes their sum by calling the algebraic evaluator. The calls the the LISP function print have been inserted here and in the following examples to show you the forms passed to the routine.

5.2.3 Polyfn

If the symbolic routine is specialized for computations with pure polynomials it can be declared polyfn. REDUCE will evaluate the arguments for such a routine to standard forms and expects that the result also will have that form. Example:

```reduce
symbolic procedure poly_plus(a,b);  
begin scalar r;  
    print a; print b;  
    r := addf(a,b);  
    print r;  
    return r;  
end;  
put('poly_plus,'polyfn,'poly_plus);
```

This routine also adds its arguments but it is specialized for polynomials. If called with a non-polynomial form an error message will be generated. In the put statement the first argument is the algebraic operator name and the third one is the name of the associated evaluator function. These may differ.

5.2.4 Psopfn

The most general type of function is that of a psopfn. REDUCE will not evaluate the parameters for such a routine, instead it passes the unevaluated list of parameters to the function. The routine has to evaluate its parameters itself (of course using the services of reval and friends). So a psopfn can handle variable numbers of parameters. The result must be an algebraic expression or nil. Example:

```reduce
symbolic procedure multi_plus0 u;  
begin scalar r;  
r:=0;  
for each x in u do  
<<x:=reval x; print x;
```
This routine can be called with an arbitrary number of parameters; it will evaluate them one after the other, add them by calling \texttt{reval} and return the sum as result. Note that the name used in algebraic mode is \textit{multi\_plus} which is different from the name of the symbolic routine doing the work. This is necessary because otherwise the argument count check of REDUCE would complain as soon as the routine is called with more than one argument.

In the next example a \texttt{psopfn} implements an operator with a fixed number of arguments; it expects two numbers as arguments and performs an extensive checking as any such routine should do; again the name of the routine and the algebraic mode operator are different:

\begin{verbatim}
symbolic procedure bin\_plus0 u;
  begin scalar p1,p2,r;
    if length u neq 2 then rederr "illegal number of arguments";
    p1:=reval car u; p2:=reval cadr u;
    if not numberp p1 then typerr(p1,"number");
    if not numberp p2 then typerr(p2,"number");
    r:=p1+p2;
    return r;
  end;

  put(\texttt{\textquoteleft bin\_plus\textquoteleft,\textquoteleft psopfn,\textquoteleft bin\_plus0});
\end{verbatim}

The functions \texttt{typerr} and \texttt{rederr} are explained in the section \textit{error management}.

### 5.2.5 Simpfn

When you declare a new algebraic operator and want to assign a special evaluation mode for it which cannot be written in algebraic mode (e.g. as a rule set) you can create a simplifier which has the task to convert each operator expression to a \textit{standard quotient}. A simplifier function is linked
to the operator by the property simpfn and has one argument which is the unevaluated parameter list of the operator expression. It will be called every time the operator appears in an expression. It must return a standard quotient. Example:

```plaintext
algebraic operator op_plus;

symbolic procedure simp_op_plus u;
begin scalar r;
   r:=simp 0;
   for each x in u do
      <<x:=simp x; print x;
      r:=addsq(r,x)
      >>;
   return print r;
end;

put('op_plus,'simpfn,'simp_op_plus);
```

In many cases the simpfn is the method of choice as it is best integrated into the REDUCE evaluation process: its results can immediately be used for subsequent calculations while algebraic form results need to be converted to standard quotients before combining them with other formulae.

Note that a simpfn which cannot simplify its argument must nevertheless return a standard quotient, then with the unevaluable form as kernel. E.g. a function for supporting the algebraic evaluation of the operator “<”:

```plaintext
algebraic operator <;

symbolic procedure simp_lessp u;
begin scalar a1,a2,d;
   a1:=simp car u; a2:=simp cadr u;
   d:=subtrsq(a1,a2);
   if domainp denr d and domainp numr d then
      return if !:minusp numr d then simp 1 else simp 0;
   return mksq({'lessp,prepsq a1,prepsq a2},1);
end;

put('lessp,'simpfn,'simp_lessp);
```

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Here all comparisons with numeric items are reduced to zero or one because of the equivalence of zero and “false” in algebraic mode, while in all other cases the non-evaluable expression is returned mapped by the function \texttt{mksq} which converts an algebraic expression into a standard quotient, ensuring the identity of composite kernels (see the section \texttt{standard forms})

5.3 Statements, forms

An additional way to link symbolic entries to algebraic mode is to invent a new syntactical unit. A REDUCE statement is introduced by assigning the property \texttt{stat = rlis} to the leading keyword of the new statement. The corresponding function will be called with one parameter which is a list of the (unevaluated) parameters of the statement. A statement normally is called for its side effect rather than for its value - so the result will be not interpreted as an algebraic quantity (\texttt{ws} not set, printed as symbolic list etc.). Example:

\begin{verbatim}
put('my_add,'stat,'rlis);

symbolic procedure my_add u;
begin scalar r;
  r:= 0;
  for each x in u do
    <<x:=reval x;
    r:=reval{'plus,x,r}
    >>;
  return r;
end;
\end{verbatim}

to be called by a statement

\begin{verbatim}
my_add x,x^2,x^3;
\end{verbatim}

A statement without parameters is introduced by the property - value pair \texttt{stat = endstat} because the statement “end” is the most prominent representative of this class.

For a more complicated syntax you can write a function which preprocesses a statement before evaluating it. Such preprocessors are linked to the leading keyword by the property \texttt{formfn}. However writing a formfn
is rather difficult as you will have to organize the reading of the rest of the statement yourself. This approach is not recommended as the formfn interface is rather delicate and might be subject to future changes.

5.4 Printing and messages

For printing in symbolic mode you can use the Standard LISP functions

<table>
<thead>
<tr>
<th>Function</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>print(u)</td>
<td>print u with LISP syntax and following linefeed</td>
</tr>
<tr>
<td>prin1(u)</td>
<td>print u with LISP syntax but without linefeed</td>
</tr>
<tr>
<td>prin2(u)</td>
<td>print u without LISP syntax and without linefeed</td>
</tr>
<tr>
<td>prin2t(u)</td>
<td>print u without LISP syntax but with linefeed</td>
</tr>
<tr>
<td>terpri()</td>
<td>print a single linefeed</td>
</tr>
</tbody>
</table>

*LISP syntax* here means that a string is surrounded by string quotes and special characters in a symbol are tagged by an exclamation mark.

However if you want to print an algebraic expression in the same style as the REDUCE write statement you should use the function writepri. This function needs two arguments, the first one is the object to be printed which must be a string, a number or an algebraic expression tagged with a dynamic quote, e.g. using the function mkquote. The second argument is one of only, first, nil, and last and controls linefeeds. Example:

```lisp
<<u:='(expt x 2);
  writepri(" u= ",’first);
  writepri(mkquote u,nil);
  writepri(” printed in algebraic style”,’last);
>>;
```

For printing one single algebraic form you can also use the function mathprint:

```lisp
  u := ’(expt x 3);
  print u;
  mathprint u;
```

6 Error management

6.1 Issue an error message

When programming an application in symbolic mode, it may be necessary to report an exception to the user, e.g. if an algebraic prerequisite for applying
the algorithm is not fulfilled. There are several routines for signalling an error in symbolic mode:

- The easiest way to complain about a bad algebraic form is to call \texttt{typerr}. This function has two parameters \texttt{typerr(exp, str)}: the first one is an algebraic expression, the bad object, and the second one is a string which describes what the object should have been. When printed \textsc{reduce} will report \textit{exp} in mathematical notation and insert a string “illegal as”. The current program is terminated.

\begin{verbatim}
u := '(plus x 1);
    . . .
if not numberp u then typerr(u,"number");
\end{verbatim}

- A general error can be reported using the function \texttt{rederr}; it prints some asterisks and then its argument in pure LISP style (no mathematical format); the argument can be a list - in that case the surrounding brackets are omitted in output. The current program is terminated.

\begin{verbatim}
rederr "algorithm not applicable if denominator neq 1";
rederr {"never call me with ",u," again"};
\end{verbatim}

- Similar to \texttt{rederr} the routine \texttt{rerror} terminates the run with an error message; it has been designed for usage in a \textsc{reduce} package enabling the error handler to localize the error reason and to provide additional information to the end user (e.g. via a \texttt{help} system). It has 3 parameters: \texttt{rerror(pack, nbr, mess)} where \textit{pack} is the name of the package (usually a quoted identifier), \textit{nbr} is an integer error number local to the package, beginning with 1 and \textit{mess} is the error message just like \texttt{rederr}.

\begin{verbatim}
rerror('asys,13,"remainder not zero");
\end{verbatim}

Wherever possible \texttt{rerror} should be used instead of \texttt{rederr}.

### 6.2 Catching errors

On the other hand a symbolic program might want to react to an exception which occurred in a calculation branch. \textsc{reduce} supports this by encapsulation with the routines \texttt{errorset*} and \texttt{errorset}. You have to construct the action to be encapsulated as a subroutine call in prefix LISP syntax.
which is identical to algebraic form syntax: a list where the first element
is the name of the function (usually a quoted symbol) followed by the ar-
guments; these must be tagged by quote, best using the function mkquote.
The second parameter describes the error mode: if $t$, an eventual error dur-
ing the evaluation will be reported to the user, if $nil$ the error message will
be suppressed. errorset additionally has a third parameter which tells the
error manager whether to print a backtrace or not in an error case. The
result of these functions will be

- in the case of an error an error code (its form depends on the underlying
  LISP system),
- otherwise the result of the computed value as car of a list.

You should use the function errorp for testing whether the call failed or
not; in the positive case pick out the result by a car access. Example:

```reduce
begin scalar u,v;
. . .
q := errorset!*[{'quotsq,mkquote(u),mkquote(v)},nil];
if errorp q then rederr "division failed" else
  q:=car q;
. . .
```

Note that you cannot write the expression to be encapsulated in REDUCE
syntax here. And also `(quotsq u v)` would be incorrect because then $u$ and
$v$ were constants and not references to the local variables $u$ and $v$. Only the
`mkquote` expressions guarantee that at run time a list is generated where
the values of $u$ and $v$ are inserted as parameters for the call.

7 Compiling, modules, packages

7.1 Program modules

It is a good practice to design a complex program group in separate mod-
ules each residing in a separate source file. You can use the REDUCE state-
ment pair module - endmodule to reflect this structure. A module has a
name which often will be identical with its source file (without extension,
of course). This correspondence is not essential, but is very practical for
the daily work. As a side effect module switches automatically to symbolic
mode during input and endmodule switches back to the original mode. A
typical module in a file would look like
module alggeo1; % Primitives for algebraic geometry.

% Author: Hugo Nobody, Nirwanatown, Utopia.

....... the symbolic mode code

endmodule;
end;

If this module resides in a file alggeo1.red it can either be read into a
REDUCE session by

    in "alggeo1.red"

or it can be compiled to a fast loadable binary by

    faslout "alggeo1";
    in "alggeo1.red"
    faslend;

You then can load this module by load alggeo.

Although REDUCE currently does not yet rely exclusively on declarations
exports and imports, these should be used in order to be safe for the future.
exports should name all entry points which the module offers for external
access and imports is the list of modules or packages which the module
needs at run time or during compilation.

## 7.2 Packages

If your application consists of several modules you can design it as a package.
A package is a collection of modules which form a unit but allow you an
individual management, e.g. updating and recompilation. A package is defined
by a header module which should have the name of the package as the module
name. This module then contains a declaration for the package by the function create-package, which expects as first parameter a list of the modules
which belong to the package (this list must begin with the header package)
and a second dummy parameter to be set to nil (this parameter is meaningful only for modules in the REDUCE kernel). Additionally the header
module should contain all global declarations for the complete package, e.g.
fluid, global declarations and exports and imports statements. Also,
all macros, smacros used in more than one module in the package should be
defined in the header module.
As long as you develop your application and frequently recompile single modules it would be practical to replace the function `create-package` by a loop calling `load-package` (please note the underscore: `load-package` is a different function not usable here). You then can load the complete package by one call to `load-package` for the header module. Example:

```lisp
module alggeo; % Algebraic geometry support.

% Author: Hugo Nobody, Nirwanatown, Utopia.

% create_package('(alggeo alggeo1 alggeo2 alggeo3));
for each x in '(alggeo1 alggeo2 alggeo3) do load x;

fluid '(alggeo_vars!* ......);
exports ...
    ... and so on

endmodule;
end;
```

Later you can replace the load loop `load` by a correct `create-package` expression - you will then have to compile the object as one big binary with the header module as the first entry.

For compiling you use the function `faslout` which creates a fast loading file - see the corresponding section in the User’s manual.

8 Coding suggestions

In order to guarantee that your program is as portable and version independent as possible the following rules may be useful. Note that these are recommendations only - your program may do a good job for you even if you ignore all of them.

- Don’t use functions from the underlying LISP which are not explicitly part of REDUCE or `standard lisp` - they will probably be missing in other support systems.

- Don’t rely on character case. Some implementations support upper case output, others prefer lower case. As the tendency goes towards lower case, the best policy is to use lower case characters only, except in strings or comments.
• Use proper indentation. Take the REDUCE sources as example. If the indentation corresponds to the block structure your programs will be better readable.

• Don’t produce lines with more than 72 characters. There are still computers and networks around which have problem if the size of the good old punched card is exceeded.

• Write comments. The major entry point routines of your symbolic code should have one comment line following the procedure declaration; this comment should be one or more complete English sentences, beginning with an upper case character and ending with a dot.

• Try to avoid name conflicts. It is a common policy to name global and fluid variables with a trailing asterisk and to use a common name prefix for the procedure names.

• If you produce a large piece of software, organize it as a REDUCE package.

9 Appendix: Standard LISP extensions

The following typical LISP functions are defined in the support module of the rlisp package.

\['\text{atsoc}(u:\text{any},p:\text{alist})\]
Same as \text{assoc}, but using the operator \text{eq} instead of \text{equal} for the search. \text{atsoc} can be applied only if the identity of the object to be searched for is guaranteed, e.g. for symbols. It is substantially faster than \text{assoc}.

\['\text{copyd}(u:\text{id},v:\text{id})\]
Copies the procedure definition of \( v \) to the symbol \( u \). Afterwards the procedure can be called by using the name \( u \). Frequently used for embedding a function, e.g. for implementing a private trace:

\[
\begin{align*}
\text{if null getd 'myprog_old then copyd('myprog_old,'myprog);} \\
\text{symbolic procedure myprog(u);} \\
\text{<<print {"calling myprog with parameter ",u};} \\
\text{myprog_old(u)>>}
\end{align*}
\]

\['\text{eqcar}(u:\text{any},v:\text{any})\]

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Implements the frequently needed test \textit{pairp} $u$ and \textit{car} $u$ \textit{eq} $v$.

\textbf{listp($u$:any)}
Returns \texttt{t} if $u$ is a top level list ending with a \texttt{nil}.

\textbf{mkquote($u$:any)}
tags $u$ with a quote. This function is useful for cases where a LISP eval situation is given, e.g. for coding for an \texttt{errorset} or for \texttt{writepri}.

\textbf{reversip($u$:list)}
Reverses the elements of $l$ in place - faster than \texttt{reverse}, but desctroys the input list.

\textbf{smember($u$:any,$v$:any)}
Similar to \texttt{member} this routine tests whether $u$ occurs in the structure $v$, however in an arbitrarily deep nesting. While \texttt{member} tests only the top level list, \texttt{smember} descends down all branches of the tree.

\textbf{smemq($u$:any,$v$:any)}
Same as \texttt{smemr}, however using \texttt{eq} as test instead of \texttt{equal}.

The following routines handle lists as sets (every element at most once):

\textbf{union($u$:any,$v$:any)}
\textbf{intersection($u$:any,$v$:any)}
\textbf{setdiff($u$:any,$v$:any)}

The Standard LISP \texttt{apply} function requires that the parameters for the function to be applied are encoded as a list. For standard cases of 1, 2 or 3 parameters the following variants allow an easier encoding:

\textbf{apply1($u$:id,$u$:any)}
\textbf{apply2($u$:id,$u$:any,$v$:any)}
\textbf{apply3($u$:id,$u$:any,$v$:any,$w$:any)}
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