# SCOPE 1.5 A Source-Code Optimization PackagE for REDUCE 3.7

User's Manual

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#### Abstract

The facilities, offered by SCOPE 1.5, a Source-Code Optimization PackagE for REDUCE 3.6, are presented. We discuss the user aspects of the package. The algorithmic backgrounds are shortly summarized. Examples of straightforward and more advanced usage are shown, both in algebraic and symbolic mode. Possibilities for a combined use of GENTRAN and SCOPE

are presented as well.

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# 1 Introduction

An important application of computer algebra systems is the generation of code for numerical purposes via automatic or semi-automatic program generation or synthesis. GENTRAN [11, 9, 10, 12], a flexible general-purpose package, was especially developed to assist in such a task, when using MAC-SYMA or REDUCE. Attendant to automatic program generation is the problem of automatic source-code optimization. This is a crucial aspect because code generated from symbolic computations often tends to be made up of lengthy arithmetic expressions. Such lengthy codes are grouped together in blocks of straight-line code in a program for numerical purposes. The main objective of SCOPE, our source-code optimization package, has been minimization of the number of (elementary) arithmetic operations in such blocks. This can be accomplished by replacing repeatedly occuring subexpressions, called common subexpressions or cse's for short, by placeholders. We further assume that new statements of the form "placeholder := cse" are inserted correctly in the code. This form of optimization is often helpful in reducing redundancy in (sets of) expressions. A recent application, code generation for an incompressible Navier-Stokes problem [15], showed reduction from 45.000 lines of FORTRAN code to 13.000 lines.

Optimizing compilers ought to deal effectively and efficiently with the average, hand coded program. The enormous, arithmetic intensive expressions, easily producable by a computer algebra system, fall outside the range of the FORTRAN programs, once analyzed and discussed by Knuth [22]. He suggested that optimization of the arithmetic in such a program is slightly overdone. The usual compiler optimization strategy is based on easy detection of redundancy, without assuming the validity of (some) algebric laws (see for instance [16]) Our optimization strategy however, requires the validity of some elementary algebraic laws. We employ heuristic techniques to reduce the arithmetic complexity of the given representation of a set  $E_{in}$ of input statements, thus producing a set  $E_{out}$  of output assignment statements.  $E_{in}$  and  $E_{out}$  define blocks of code, which would compute the same exact values for the same exact inputs, thus implicitly proving the correctness of the underlying software. Obviously the use of  $E_{out}$  ought to save a considerable amount of execution time in comparison with the application of  $E_{in}$ . Johnson et al [21] suggest that such transformations do not destabilize the computations. However this is only apparent after a careful error analysis of both  $E_{in}$  and  $E_{out}$ . In view of the size of both  $E_{in}$  and  $E_{out}$  such an

#### 1 INTRODUCTION

analysis has to be automatized as well. Work in this direction is in progress [20, 24, 25].

Although the use of SCOPE can consideribly reduce the arithmetic complexity of a given piece of code, we have to be aware of possible numerical side effects. In addition we have to realize that a mapping is performed from one source language to another source language, seemingly without taking into account the platform the resulting numerical code has to be executed on. Seemingly, because we implemented some facilities for regulating minimal expression length and for producing vector code.

This manual is organized as follows. We begin with some preliminary remarks in section 2. The history and the present status, the optimization strategy and the interplay between GENTRAN and SCOPE are shortly summarized. The basic algebraic mode facilities are presented in section 3. Special SCOPE features are discussed in section 4. Besides facilities for Horner-rules and an extended version of the REDUCE function structr, we introduce some tools for extending SCOPE with user defined specialties. File management follows in section 5. In section 6 declaration handling and related issues are discussed, before illustrating our strategy concerning data dependencies and generation of vector code in section 7. In section 8 is shown how a combined used of GENTRAN and SCOPE can be profitable for program-generation. The use of SCOPE in symbolic mode is presented in section 9. A SCOPE syntax summary is given in section 10. For completeness we present guidelines for installing the package in the last section.

#### Requests

- Comment and complaints about possible bugs can be send to the author using the e-mail address inflwh@cs.utwente.nl. A compact piece with REDUCE code, illustrating the bug, is prefered.
- When SCOPE 1.5 is used in prepairing results, presented in some publication, reference to its use is highly appreciated. A copy of the published document as well.

# 2 Preliminaries

For completeness we present a historical survey, a birds-eye view of the overall optimization strategy and the interplay between GENTRAN and SCOPE.

#### 2.1 History and Present Status

The first version of the package was designed to optimize the description of REDUCE-statements, generated by NETFORM [27, 26]. This version was tailored to a restrictive class of problems, mainly occurring in electrical network theory, thus implying that the right-hand sides (rhs's) in the input were limited to elements of  $\mathbf{Z}_{2}[V]$ , where V is a set of identifiers. The second version [30] allowed rhs's from  $\mathbf{Z}[V]$ . For both versions the validity of the commutative and the associative law was assumed. A third version evolved from the latter package by allowing to apply the distributive law, i.e. by replacing (sub)expressions like a.b + a.c by a.(b + c) whenever possible. But the range of possible applications of this version was really enlarged by redefining V as a set of kernels, implying that almost any proper REDUCE expression could function as a rhs. The mathematical capabilities of this version are shortly summarized in [33], in the context of code generation and optimization for finite-element analysis. This version was further extended [32] with a declaration-module, in accordance with the strategy outlined in [1], chapter 6. It is used in combination with GEN-TRAN, for the construction of Jacobians and Hessians [28, 2] and also in experiments with strategies for code vectorization [14]. In the meantime the Jacobian-Hessian production package, at present called GENJAC, is further extended with possibilities for global optimization and with some form of loop-differentiation. So in stead of optimizing separate blocks of arithmetic we are now able to optimize complete programs, albeit of a rather specific syntactical structure [2]. The present 1.5 version of SCOPE, is an intermediate between the distributed first version and the future, second version. Version 2 is currently in development and will contain, besides the already existing common sub expression (cse) searches, facilities for structure and pattern recognition. The 1.5 version permits -using the present REDUCE terminology-rounded coefficients, based on the domain features, described in [4], discovery and adequate treatement of a variety of data dependencies, and quotient-optimization, besides a collection of other improvements and

refinements for using the facilities in the algebraic mode. Furthermore, an increased flexibility in the interplay between GENTRAN and SCOPE is accomplished. It is used for experiments concerning automatic differentiation [13], symbolic-numeric approaches to stability analysis [8, 7] and for code generation for numerically solving the Navier-Stokes equations for incompressible flows [15]. An interesting example of its use elsewhere can be found in [6].

## 2.2 Acknowledgements

Many discussions with Victor V. Goldman, Jaap Smit and Paul S. Wang have contributed to the present status of SCOPE. I express my gratitude to the many students, who have also contributed to SCOPE, either by assisting in designing and implementing new facilities, or by applying the package in automated program generation projects in and outside university, thus showing shortcomings and suggesting improvements and extensions. I mention explicitly Frits Berger, Johan de Boer, John Boers, Pim Borst, Barbara Gates, Marcel van Heerwaarden, Pim van den Heuvel, Ben Hulshof, Emiel Jongerius, Steffen Posthuma, Anco Smit, Bernard van Veelen and Jan Verheul.

#### 2.3 The Optimization Strategy in a Birds-eye View

In [29, 30] we described the overall optimization strategy used for SCOPE as a composite function  $\mathbb{R}^{-1} \circ \mathbb{T} \circ \mathbb{R}$ . The function  $\mathbb{R}$  defines how to store the input  $\mathbb{E}_0$  in an expression database  $\mathbb{D}_0$ . The inverse function  $\mathbb{R}^{-1}$  defines the output production. The function  $\mathbb{T}$  defines the optimization process itself. It essentially consists of a heuristic remodeling of the (extendable and modifiable) expression database in combination with storing information required for a fast retrieval and correct insertion of the detected cse's in the output. This is accomplished by an iteratively applied search, resulting in a stepwise reduction of the arithmetic complexity of the input set, using an extended version of Breuer's grow factor algorithm [5, 29, 30]. It is applied until no further profit is gained, i.e. until the reduction in arithmetic complexity stops. Before producing output, a finishing touch can be performed to further reduce the arithmetic complexity with some locally applied techniques. Hence  $\mathbb{T}$  is also a composite function. The overall process can be summarized as follows:

$$\begin{array}{rcl} \mathbf{R} & : & \mathbf{E}_{\mathrm{in}} = \mathbf{E}_{0} & \rightarrow & (\mathbf{D}_{0}, \mathrm{profit}_{0}) \\ \mathbf{T}_{\beta} & : & (\mathbf{D}_{\mathrm{i}}, \mathrm{profit}_{\mathrm{i}}) & \rightarrow & (\mathbf{D}_{\mathrm{i+1}}, \mathrm{profit}_{\mathrm{i+1}}) \ , \ \mathbf{i} \ = \ 0, \dots, \ \lambda - 1. \\ \mathbf{F} & : & (\mathbf{D}_{\lambda}, \mathrm{profit}_{\lambda}) & \rightarrow & D_{\lambda} \\ \mathbf{R}^{-1} & : & D_{\lambda} & \rightarrow & \mathbf{E}_{\lambda} \ = \ \mathbf{E}_{\mathrm{out}} \end{array}$$

 $D_0$  is created as a result of an R-application performed on input  $E_0$ . The termination condition depends on some profit criterion related to the arithmetic complexity of the latest version of the input,  $D_i$ . Hence we assume profit<sub>i</sub> = true for  $i = 0, \dots, \lambda - 1$  and profit<sub> $\lambda$ </sub> = false. The function T is defined by  $T = F \circ T^{\lambda}_{\beta}$ , where  $T_{\beta}$  defines one iteration step, i.e. one application of the extended version of Breuer's algorithm, and where F defines a finishing touch, resulting in the final version  $D_{\lambda}$  of  $D_0$ , used to produce the output  $E_{\lambda}$ . It is stated in [30] that the computing time for  $T^{\lambda}_{\beta}$  is O(n.m), where n is the size of  $E_{in}$  and m the number of cse's found during this process. Practical experience showed that the finishing touch can take about 10 % of the actual cpu-time and that its real profit is limited. Therefore its use is made optional.

The wish to optimize source code, defining arithmetic, usually leads an attempt to minimize the arithmetic complexity. This can be accomplished by replacing cse's by placeholders, assuming a new assignment statement "placeholder := cse" is correctly inserted in the code. So most of the cse-searches are done in right hand sides of arithmetic assignment statements.

The search strategy depends on the permissible structure of the arithmetic expressions. We assume these expressions to be multivariate polynomials or rational functions in a finite set of kernels, and presented in some normal form. Let us further assume that scalar placeholders are substituted for the non-scalar kernels, such that back-substitution remains possible, using an adequate information storage mechanism. Then we are left with the interesting question how to define a minimal set of constituents of multivariate polynomials in some normal form norm. Let us take as an example of such a polynomial or rational function  $p = 3a + 2b + 3b^2c(3a + 2b)(c + d)^2$ . We easily recognize linear forms, i.e. 3a + 2b (twice) and c + d, possibly raised to some power  $((c + d)^2)$ , power products, such as  $b^2c$ , or monomial parts of products, i.e.  $3b^2c$ . Hence with some imagination, one realizes that every polynomial can be decomposed in a set of linear forms and a set of power products. When assuming the validity of the commutative and the associative law, one can also realize that we can associate a coefficient matrix

with the linear forms and an exponent matrix with the power products. The rows can correspondent with (sub)expressions and the columns with scalar identifiers. The entries are either coefficients or exponents. It is therefore conceivable to make a parser, mapping a set of REDUCE expressions in a database, consisting of two incidence matrices and a function table, such that the original expressions can be retrieved. Taking a group of assignment statements or a list of equations, where in both cases the lhs's function as right hand side recognizers, does not confuse this idea. This rather informal indication merely serves as a suggestion how R and its inverse operation are designed.

So we suggest that we can consider any set of rhs's as being built with linear forms and power products only. An additional remark is worth being made: Non-scalar kernels will in general have non-commutable arguments. These arguments can in turn be arbitrary REDUCE-expressions, which also have to be incorporated in the database. Hence the function table is created recursively.

What is a cse and how do we locate its occurrences? A (sub)expression is common when it occurs repeatedly in the input. The occurrences are, as part of the input, distributed over the matrices, and shown as equivalent (sub)patterns. In fact, we repeatedly search for completely dense (sub)matrices of rank 1. The expression 2a + 3c is a cse of  $e_1 = 2a + 4b + 3c$ and  $e_2 = 4a + 6c + 5d$ , representable by (2,4,3,0) and (4,0,6,5), respectively. We indeed have to assume commutativity, so as to be able to produce new patterns (2,0,3,0,0), (0,4,0,0,1) and (0,0,0,5,2), representing s = 2a + 3c,  $e_1 = 4b + s$  and  $e_2 = 5d + 2s$ , respectively, and thus saving one addition and one multiplication. Such an additive cse can be a factor in a (sub)product, which in turn can extend its monomial part, when replacing the cse by a new symbol. Therefore an essential part of an optimization step is regrouping of information. This migration of information between the matrices is performed if the Breuer-searches are temporarily completed. After this regrouping the distributive law is applied, possibly also leading to a further regrouping. If at least one of these actions leads to a rearrangement of information the function  $T_{\beta}$  is again applied. In view of the iterative character of the optimization process we always accept minimal profits.

A similar search is performed to detect multiplicative cse's, for instance occuring in  $e_1 = a^2 b^4 c^3$  and  $e_2 = a^4 c^6 d^5$ . However, given a power product  $\prod_{i=1}^m x_i^{a_i}$ , any product  $\prod_{i=1}^m x_i^{b_i}$ , such that some  $b_i \leq a_i$ , for i = 1(1)m, can

function as a cse. We therefore extend the search for multiplicative cse's by employing this property, and as indicated in [30].

The finishing touch F is made to perform one-row and/or one-column searches. Once the extended Breuer-searches do not lead to further reduction in the arithmetic complexity we try -applying it- to improve what is left. The coefficients in (sub)sums can have, possibly locally, a gcd, which can be factored out. One-column operations serve to discover and to replace properly constant multiples of identifiers. As part of the output-process we subject all exponentiations left - at most one for each identifier - to an addition chain algorithm.

#### 2.4 The Interplay between GENTRAN and SCOPE 1.5

The current version of SCOPE is written in RLISP. Like GENTRAN, it can be used as an extension of REDUCE. When SCOPE is loaded GENTRAN is also activated.

If we start a REDUCE session, we create an initial algebraic mode programming environment. All switches get their initial value, such as ON EXP,PERIOD and OFF FORT. Certain REDUCE commands serve to modify or to enrich the current environment. Others are used to perform calculations, producing formulae. Such a calculation follows a standard pattern, although parts of this repertoire can be influenced by the user, for instance by changing the value of certain switches. Usually execution is a three-step process. First the infix text is parsed into a prefix form. Then the internal algebra is applied on this form, leading to a so-called standard quotient. This quotient is stored on the property list of the identifier functioning as assigned variable for this value. The last step defines the inverse route from internal existence to external presentation in infix form. Occurrences of identifiers are recursively replaced by their standard quotient representation when the internal algebra is applied. Hence the REDUCE simplification strategy follows the imperative programming paradigm.

When loading SCOPE, and thus GENTRAN, the environment is enriched with features for program generation and program optimization. Evaluation of GENTRAN and SCOPE commands differs from the standard REDUCE approach to evaluation. Both packages employ their own storage mechanism. The output is normally produced as a side-effect of the command evaluation. The output is directed to some medium, a file or a screen. Command evaluation is similar in GENTRAN and in SCOPE.

The code generation process of GENTRAN can be viewed as the application of a composite function to an argument, which is almost equivalent with a piece of REDUCE code. Almost, because some GENTRAN specific facilities can be used. We can distinguish between the preprocessing phase, the translation phase and the postprocessing phase. During preprocessing relevant parts of the input are evaluated prior to translation into prefix form. Such a locally performed evaluation can be accomplished through recognition of certain "evaluation markers", i.e. modifications of the traditional assignment symbol :=, such as ::=, :=: and ::=:. The := operator "orders" GEN-TRAN to translate the statement literally. Addition of an extra colon to the left hand side orders subscript expression evaluation before translation. An extra colon to the right hand side leads to right hand side evaluation, but without application of the storage mechanism of REDUCE. Hence evaluations remain anonymous and are only incorporated in the translatable "text". Another aspect of preprocessing is initialization of the symbol table, using information provided by a DECLARE statement. GENTRAN also allows to further rewrite (sets of) arithmetic assignment statements, using the switches GENTRANOPT and GENTRANSEG, introduced for code optimization (using SCOPE) and segmentation, respectively. It possibly leads to storage of additional information in the symbol table. During the translation phase the final internal form of the code is produced, in combination with formatting specifications and instructions to produce declarations. Postprocessing finally does produce formatted code strings. So essentially, each GENTRAN command has its own separate translation process, implying that the symbol table, required for the production of declarations, is fresh at the beginning of a GENTRAN command evaluation.

As stated before, a SCOPE command evaluation is also a composite operation. The role of the assignment operators in both GENTRAN and SCOPE is similar. In SCOPE, the locally performed evaluation provides information to be entered in the database  $D_0$ . If the declaration feature is activated, the symbol table generation and maintenance mechanism is borrowed from GENTRAN. For output production, we can make a choice from GENTRAN's target language repertoire. When declarations are required, we simply obey the GENTRAN regime as well.  $D_{\lambda}$  is used to update the symbol table. All cse-names, generated during the optimization process, are typed in accordance with the strategy for dynamic typing, which is discussed in [1], chapter 6. We assume all relevant identifiers of  $E_{in}$  to be adequately typed, using SCOPE's DECLARE facility, an equivalent of GEN-TRAN's DECLARE statement. The production of  $D_{\lambda}$  is completely decoupled from the normal REDUCE simplification strategy, because we employ our own expression database.

In principle, the status of REDUCE before and after a GENTRAN or SCOPE command execution is unaltered. In principle, because some minor modifications, although user controlable, may be necessary. The special assignment symbols -also usable in SCOPE- were only introduced as a syntactical instrument to allow internal algebraic actions, decoupled from the standard REDUCE expression processing.

This short excursion into the different evaluation strategies is added to assist in understanding the functioning of the different SCOPE commands and facilities, to be introduced in the next sections.

# 3 The Basic SCOPE 1.5 Facilities in the Algebraic Mode

REDUCE allows, roughly speaking, two modes of operation in algebraic mode: ON EXP or OFF EXP. The first is the default setting, leading to expanded forms. The latter gives unexpanded forms, as discussed by Hearn in some detail [18, 19]. It is obvious that the OFF EXP setting is in general preferable over the ON EXP setting when attempting to optimize the description of a set of assignment statements.

The result of an application of SCOPE can be influenced by the use of certain REDUCE- or SCOPE-switches. The influence of EXP is obvious: unexpanded input is more compact than expanded. ON ACINFO serves to produce tables with the numbers of arithmetic operations, occuring in  $E_0$  and  $E_{\lambda}$ , respectively. ON INPUTC serves to echo the input, processed by SCOPE. The actual form of the input can be the consequence of locally performed evaluations, before the actual parsing into the database takes place. ON PRIMAT can be used to visualize both  $D_0$  and  $D_{\lambda}$ . ON PRIALL finally, can be used instead of ON ACINFO, INPUTC, PRIMAT. These SCOPE-switches are initially all turned OFF. SCOPE has a facility to visualize the status of all SCOPE-switches and some relevant REDUCE-switches. The current status of all relevant switches can be presented with the command

#### SCOPE\_SWITCHES\$

## Example 1

The start of a REDUCE session shows the initial state of REDUCE, directly after loading the SCOPE package. The set of relevant switches is made visible. Besides the REDUCE switches EVALLHSEQP, EXP, FORT, NAT, PERIOD,  $\widehat{ROUNDBF}$  and ROUNDED six additional SCOPE switches, i.e. AGAIN, FTCH, INTERN, PREFIX, SIDREL and VECTORC, and the GENTRAN switches DOUBLE and GENTRANOPT are thus presented. They all wil be discussed in more detail below.

REDUCE 3.7, 15-Apr-1999 ...

- 1: load\_package nscope\$
- 2: SCOPE\_SWITCHES\$

ON : exp ftch nat period

OFF : acinfo again double evallhseqp fort gentranopt inputc intern prefix priall primat roundbf rounded sidrel vectorc

```
3: % etc. ...
```

Output is by default given in REDUCE syntax, but FORTRAN syntax is possible in the usual way, e.g. ON FORT and OFF PERIOD, for instance. The use of other target languages from the GENTRAN repertoire is discussed in section 6.

## 3.1 The OPTIMIZE command: Straightforward use

A SCOPE application is easily performed and based on the use of the following syntax:

$<$ SCOPE_application $>$	::=	$OPTIMIZE < object\_seq> [INAME < cse\_prefix>]$
$<$ object_seq $>$	::=	<object>[,<object_seq>]</object_seq></object>
< object >	::=	$\langle stat \rangle \mid \langle alglist \rangle \mid \langle alglist\_production \rangle$
$\langle stat \rangle$	::=	<name> <assignment operator=""> <expression></expression></assignment></name>
<assignment operator=""></assignment>	::=	:=   ::=   ::=:   :=:
<alglist></alglist>	::=	$\{< eq\_seq>\}$
<eq_seq></eq_seq>	::=	$<$ name $> = <$ expression $>$ [, $<$ eq_seq $>$ ]
$< alglist\_production >$	::=	<name>   <function_application></function_application></name>
<name></name>	::=	$\langle id \rangle   \langle id \rangle (\langle a\_subscript\_seq \rangle)$
$<$ a_subscript_seq>	::=	<a_subscript>[,<a_subscript_seq>]</a_subscript_seq></a_subscript>
<a_subscript></a_subscript>	::=	<integer>   <integer infix_expression=""></integer></integer>
$< cse\_prefix >$	::=	<id></id>

A SCOPE action can be applied on one assignment statement. The assigned variable is either a scalar identifier, like z in example 2, or a name with subscripts, such as a(1,1) in example 3. In stead of one statement a sequence of such statements, separated by comma's, is possible. An alternative is provided by the use of an algebraic mode list, consisting of REDUCE equations. An assigned variable, identifying such a list, is allowed as well. Examples are presented in section 3.2. The function\_application is introduced in section 4. Such an application ought to produce an algebra.

The expressions, i.e. rhs's in assignments or equations are legal REDUCE expressions or ought to evaluate to such expressions. Statements inside expressions are allowed, but only useful if these expressions are evaluated, before being optimized. Only integer or rounded coefficients are supported by SCOPE. So we either suppose the default integer setting or allow the switch ROUNDED to be turned ON.

The optional use of the INAME extension in an OPTIMIZE command is introduced to allow the user to influence the generation of cse-names. The cse\_prefix is an identifier, used to generate cse-names, by extending it with an integer part. If the cse\_prefix consists of letters only, the initially selected integer part is 0. All following integer parts are obtained by incrementing the previous integer part by 1. If the user-supplied cse\_prefix ends with an integer its value functions as initial integer part. The gensym-function is applied when the INAME-extension is omitted. The three alternatives are illustrated in example 2.

As stated before minimal profits are accepted during all stages of the optimization process: many small steps may lead to impressive final results. But it can also lead to unwanted details. Therefore, it can be desirable to rerun an optimization request with a restriction on the minimal size of the rhs's. The command

#### SETLENGTH <integer>\$

can be used to produce rhs's with a minimal arithmetic complexity, dictated by the value of its integer argument. Statements, used to rename function applications, are not affected by the SETLENGTH command. The default setting is restored with the command

#### **RESETLENGTH**\$

We now illustrate the use of the OPTIMIZE command through a number of small examples, being parts of REDUCE sessions. We show in example 2 the effect of the different visualization switches, the use of SETLENGTH and RESETLENGTH and of the three INAME alternatives. In example 3 the effect of some of the GENTRAN and SCOPE input processing features is presented. Some finishing touch activities are illustrated in the examples 4 and 5. The approach towards rational exponents is presented in example 6, while some form of quotient optimization is illustrated in example 7. Finally, we present the differences in ON/OFF EXP processing in example 8.

#### Example 2

The multivariate polynomial z is a sum of 6 terms. These terms, monomials, are constant multiples of power products. A picture of  $D_0$  is shown after the input echo. The sums-matrix consists of only one row, identifiable by its Fa(the)r z, the lhs. Its exponent is given in the EC (Exponent or Coefficient) field. The 6 monomials are stored in the products-matrix. The coefficients are stored in the EC-fields and the predecessor row index, 0, is given in the Far-field. Before the  $D_{\lambda}$  picture is given the effect of the optimization process, the output and the operator counts are shown. The optimized form of z is obtained by applying the distributive law. The output also shows applications of an addition chain algorithm ([23] 441-466) as part of  $\mathbb{R}^{-1}$ , although its use in example 4 is more apparent.

Observe that the output illustrates the heuristic character of the optimization process: In this particular case the rhs can be written as a polynomial in g4, thus saving one extra multiplication.

The SETLENGTH command is illustrated too. See also example 12. Application of a Horner-rule may be profitable as well. See, for instance example 16.

```
ON PRIALL$
z:=a<sup>2</sup>*b<sup>2</sup>+10*a<sup>2</sup>*m<sup>6</sup>+a<sup>2</sup>*m<sup>2</sup>+2*a*b*m<sup>4</sup>+2*b<sup>2</sup>*m<sup>6</sup>+b<sup>2</sup>*m<sup>2</sup>;
        2 2
                      2 6
                                 2 2
                                                    4
                                                             2 6
                                                                        2 2
z := a *b + 10*a *m + a *m + 2*a*b*m + 2*b *m + b *m
OPTIMIZE z:=:z$
        2 2
                      2 6
                                                             2
                                                                        2 2
                                 2 2
                                                    4
                                                                6
z := a *b + 10*a *m + a *m + 2*a*b*m + 2*b *m + b *m
Sumscheme :
    || EC|Far
  0|| 1| z
 _____
```

Productscheme :

```
| 0 1 2| EC|Far
_____
 1|
       2 2 1 0
 2 6 2 10 0
 3 2 2 1 0
 4 4 1 1 2 0
 5 6 2 | 2 0
 6 2 2 1 0
_____
0 : m
1 : b
2 : a
Number of operations in the input is:
Number of (+/-) operations
                        : 5
Number of unary - operations : 0
Number of * operations : 10
Number of integer ^ operations : 11
Number of / operations : 0
Number of function applications : 0
g1 := b*a
g5 := m*m
g2 := g5*b*b
g3 := g5*a*a
g4 := g5*g5
z := g2 + g3 + g1*(2*g4 + g1) + g4*(2*g2 + 10*g3)
Number of operations after optimization is:
Number of (+/-) operations
                        : 5
Number of unary - operations : 0
Number of * operations : 12
Number of integer ^ operations : 0
Number of / operations : 0
Number of function applications : 0
Sumscheme :
  | 0 3 4 5| EC|Far
```

\_\_\_\_\_

(	) (			1	1	1	z
1	51			2	10	1	14
1	7	2	1			1	16
0	:	g4					
3	:	g1					
4	:	g2					
5	:	gЗ					

Productscheme :

	8	9	10	11	17	18	19	20	EC	Far		
7  8  9  10  11  14  16	1 1 2	1			1	2	1 2	 2  	1  1  1  1			
8 : 9 : 10 : 11 : 17 : 18 : 19 : 20 :	g4 g3 g2 g1 m b											
OFF PRIALL\$ SETLENGTH 2\$												
OPTIMIZE z:=:z INAME s\$												
2 2 s1 := b *m 2 2 s2 := a *m 4 4 z := (a*b + 2*m)*a*b + 2*(s1 + 5*s2)*m + s1 + s2												
RESETLENGTH\$												
OPTIMIZE z:=:z INAME s1\$												
s1 := b*a s5 := m*m s2 := s5*b*b s3 := s5*a*a s4 := s5*s5 z := s2 + s3 + s1*(2*s4 + s1) + s4*(2*s2 + 10*s3)												

#### Example 3

The input echo below shows the literal copy of the first assignment. This is in accordance with role the GENTRAN assignment operator := ought to play. The second assignment, this time using the operator ::=:, leads to rhs evaluation (expansion) and lhs subscript-value substitution. Application of the distributive law is refected by the rhs of a(1,1) in the presented result.

#### Example 4

The effect is shown of a finishing touch application, in combination with FORTRAN output. The value of S0 is rewritten during output preparation, and the earlier mentioned addition chain algorithm is applied. When turning OFF the switch FTCH the latter activity is skipped.

```
ON FORT$ OFF PERIOD$
```

OPTIMIZE z:=(6\*a+18\*b+9\*c+3\*d+6\*f+18\*g+6\*h+5\*j+5\*k+3)^13 INAME s;

S0=5\*(J+K)+3\*(3\*C+D+1+6\*(B+G)+2\*(A+F+H)) S3=S0\*S0\*S0 S2=S3\*S3 Z=S0\*S2\*S2 OFF FTCH\$

```
OPTIMIZE z:=(6*a+18*b+9*c+3*d+6*f+18*g+6*h+5*j+5*k+3)^13 INAME s;
```

```
Z=(5*(J+K)+3*(3*C+D+1+6*(B+G)+2*(A+F+H)))**13
```

#### Example 5

Recovery of repeatedly occurring integer multiples of identifiers, as part of the finishing touch, is illustrated. This facility is part of the finishing touch and will seemingly be neglected when using SETLENGTH 2\$ instruction in stead of OFF FTCH.

```
OPTIMIZE x:=3*a*p, y:=3*a*q, z:=6*a*r+2*b*p,
         u:=6*a*d+2*b*q, v:=9*a*c+4*b*d, w:=4*b INAME s;
s2 := 3*a
x := s2*p
y := s2*q
s0 := 2*b
s3 := 6*a
z := s0*p + s3*r
u := s0*q + s3*d
w := 4*b
v := w*d + 9*c*a
OFF FTCH$
OPTIMIZE x:=3*a*p, y:=3*a*q, z:=6*a*r+2*b*p,
         u:=6*a*d+2*b*q, v:=9*a*c+4*b*d, w:=4*b INAME t;
x := 3*p*a
y := 3*q*a
z := 2*b*p + 6*r*a
u := 2*b*q + 6*d*a
v := 4*d*b + 9*c*a
w := 4*b
ON FTCH$
SETLENGTH 2$
OPTIMIZE x:=3*a*p, y:=3*a*q, z:=6*a*r+2*b*p,
         u:=6*a*d+2*b*q, v:=9*a*c+4*b*d, w:=4*b INAME t;
```

```
x := 3*p*a
y := 3*q*a
z := 2*b*p + 6*r*a
u := 2*b*q + 6*d*a
v := 4*d*b + 9*c*a
w := 4*b
```

### Example 6

This example serves to show how SCOPE deals with rational exponents. All rational exponents of an identifier are collected. The least common multiple lcm of the denominators of these rationals is computed and the variable is replaced by a possibly newly selected variable name, denoting the variable raised to the power 1/lcm. This facility is only efficient for what we believe to be problems occurring in computational practice. That is easily verified by extending the sum we are elaborating here with some extra terms.

```
ON INPUTC, FORT$
```

OPTIMIZE z:=:FOR j:=2:6 SUM q^(1/j) INAME s;

```
1/6 1/5 1/4 1/3
z := q + q + q + q + q + sqrt(q)
S0=Q**(1.0/60.0)
S8=S0*S0
S7=S8*S0
S5=S8*S7
S3=S5*S5
S2=S8*S3
S1=S7*S2
S4=S5*s1
Z=S4+S1+S2+S3+S4*S3
```

# Example 7

The special attention, given to rational exponents, is not extended to rational coefficients. The script in this example shows four different approaches for dealing with such coefficients using the expressions assigned to **f** and **g**. We start with a literal parsing of the two assignments, leading to a form of D<sub>0</sub>, which is based on the present REDUCE strategy for dealing with fixed float numbers in the default integer coefficient domain setting. The four rational numbers  $\frac{31}{5}$ ,  $\frac{31}{10}$ ,  $\frac{93}{5}$  and  $\frac{93}{10}$  are just like  $b^{\frac{1}{5}}$ ,  $\sqrt{\sin(\cdots)}$  and  $\sin(\cdots)^{\frac{5}{3}}$ considered as kernels.

The second approach illustrates the effect of simplification in an OFF ROUNDED mode prior to parsing. The input expressions are remodeled into rational expressions, the usual internal standard quotient form.

After turning ON the switch ROUNDED we repeat the previous commands. Again some differences in evaluation can be observed. Literally taken input, the third approach, shows rational exponent optimizations prior to the production of rounded exponents in the output. The last approach, simplification before parsing, leads to a float representation for the rational exponents. SCOPE's exponent optimization features are designed for integer and rational exponents only. Floating point exponentiation is therefore assumed to be a function application.

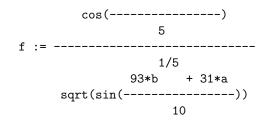
Further illustrations of operations on quotients are shown in example 22.

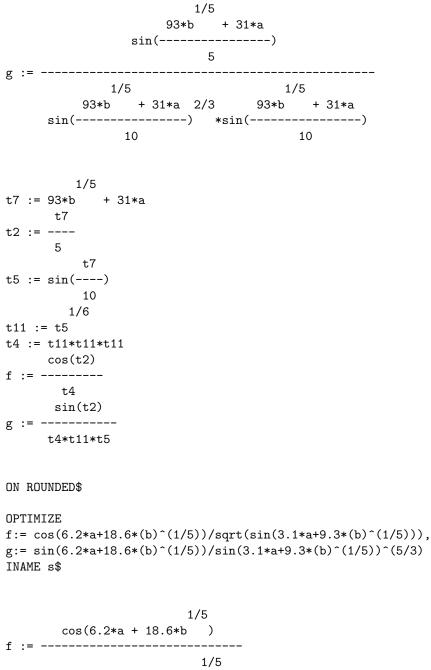
ON INPUTC\$

```
OPTIMIZE
f:= cos(6.2*a+18.6*(b)^(1/5))/sqrt(sin(3.1*a+9.3*(b)^(1/5))),
g:= sin(6.2*a+18.6*(b)^(1/5))/sin(3.1*a+9.3*(b)^(1/5))^(5/3)
INAME s$
```

```
31 93 1/5
      cos(----*a + ----*b )
         5
                5
f := -----
           31 93 1/5
    sqrt(sin(----*a + ----*b ))
           10 10
         31 93 1/5
     sin(----*a + ----*b )
        5
              5
g := -----
        31 93 1/5 5/3
    sin(----*a + ----*b )
       10 10
     1/5
s15 := b
        93 31
s12 := s15*---- + a*----
        5 5
93 31
s6 := sin(----*s15 + ----*a)
               10
       10
      1/6
s14 := s6
s5 := s14*s14*s14
   cos(s12)
f := -----
      s5
   sin(s12)
g := -----
    s5*s14*s6
OPTIMIZE
f:=: cos(6.2*a+18.6*(b)^(1/5))/sqrt(sin(3.1*a+9.3*(b)^(1/5))),
g:=: sin(6.2*a+18.6*(b)^(1/5))/sin(3.1*a+9.3*(b)^(1/5))^(5/3)
INAME t$
```

1/5 93\*b + 31\*a





sqrt(sin(3.1\*a + 9.3\*b ))

```
0.2
s5 := 9.3*b + 3.1*a
s8 := 2*s5
s4 := sin(s5)
    0.16666666666
s10 := s4
s3 := s10*s10*s10
   cos(s8)
f := -----
     s3
    sin(s8)
g := -----
    s3*s10*s4
OPTIMIZE
f:=: cos(6.2*a+18.6*(b)^(1/5))/sqrt(sin(3.1*a+9.3*(b)^(1/5))),
g:=: sin(6.2*a+18.6*(b)^(1/5))/sin(3.1*a+9.3*(b)^(1/5))^(5/3)
INAME t$
           0.2
    cos(18.6*b + 6.2*a)
f := -----
     0.2 0.5
    sin(9.3*b + 3.1*a)
                0.2
        sin(18.6*b + 6.2*a)
g := -----
        0.2 1.66666666666
    sin(9.3*b + 3.1*a)
       0.2
t6 := 9.3*b + 3.1*a
t9 := 2*t6
t5 := sin(t6)
   cos(t9)
f := -----
     0.5
     t5
       sin(t9)
g := -----
```

1.66666666667 t5

#### Example 8

The effect of ON EXP or OFF EXP on the result of a SCOPE-application is illustrated by optimizing the representation of the determinant of a symmetric (3,3) matrix m. Besides differences in computing time we also observe that the arithmetic complexity of the optimized version of the expanded representation of the determinant is about the same as the not optimized form of the unexpanded representation.

```
MATRIX m(3,3)$
```

```
m(1,1):=18*cos(q3)*cos(q2)*m30*p^2-sin(q3)^2*j30y+sin(q3)^2*j30z-
        9*sin(q3)^2*m30*p^2+j1oy+j30y+m10*p^2+18*m30*p^2$
m(2,1):=
m(1,2):=9*cos(q3)*cos(q2)*m30*p^2-sin(q3)^2*j30y+sin(q3)^2*j30z-
        9*sin(q3)^2*m30*p^2+j30y+9*m30*p^2$
m(3,1):=
m(1,3):=-9*sin(q3)*sin(q2)*m30*p^2$
m(2,2):=-sin(q3)<sup>2</sup>*j30y+sin(q3)<sup>2</sup>*j30z-9*sin(q3)<sup>2</sup>*m30*p<sup>2</sup>+j30y+
        9*m30*p^2$
m(3,2):=
m(2,3):=0$
m(3,3):=9*m30*p^2+j30x$
ON ACINFO, FORT$ OFF PERIOD$
OPTIMIZE detm:=:det(m) INAME s;
Number of operations in the input is:
Number of (+/-) operations
                            : 36
Number of unary - operations : 0
Number of * operations : 148
Number of integer ^ operations : 84
Number of / operations
                                : 0
Number of function applications : 32
      S2=SIN(REAL(Q2))
      S30=S2*S2
```

```
S3=SIN(REAL(Q3))
     S29=S3*S3
     S31=P*P
     S8=S31*M30
     S32=S8*S8
      S4=S32*J30Y
     S28=S32*S8
     S9=S29*M10
     S10=S30*S29*S29
     S44=COS(REAL(Q3))*COS(REAL(Q2))
     S11=S44*S44
     S20=S31*S8
     S23=S31*J30X
     S22=S29*J30X
     S24=S8*J10Y
     S19=M10*J30Y
     S43=81*S32*J30X
     S35=-S43-(81*S32*J10Y)
     S36=-(729*S29*S28)-(81*S29*S4)
     S37=J30Z-J30Y
     S39=9*S37
     S40=9*J30X
     S41=81*S32*J30Z
     S42=81*S4
     DETM=S42+S36-S35+729*S28+S37*(S22*J10Y+9*S29*S24+S23*S9)+S10*(S42-
     . S41)+S20*S8*81*(M10-S9)+S20*S9*(S39-S40)+S22*S8*(S39-(9*J10Y))+
     . S20*(9*S19+S40*M10)+S24*(S40+9*J30Y)+J30Y*J30X*(J10Y+9*S8)+S28*
     . 729*(S10-S11)+S29*(S41+S35)+S36*S30+S23*S19-(S43*S11)
Number of operations after optimization is:
Number of (+/-) operations : 30
Number of unary - operations : 0
                              : 59
Number of * operations
Number of integer ^ operations : 0
Number of / operations
                       : 0
Number of function applications : 4
OFF EXP$
OPTIMIZE detm:=:det(m) INAME t;
Number of operations in the input is:
```

```
Number of (+/-) operations
                                : 23
Number of unary - operations
                                : 1
Number of * operations
                                : 38
Number of integer ^ operations : 21
Number of / operations
                               : 0
Number of function applications : 10
      T1=SIN(REAL(Q3))
      T9=T1*T1
      T8=P*P
      T5=T8*M30
      T16=9*T5
     T10=-T16-(9*T5*COS(REAL(Q3))*COS(REAL(Q2)))
      T13=(T16+J30Y-J30Z)*T9
      T15=T13-J30Y
      T0=T15+T10
      T14=T13-T16-J30Y
      T17=T5*SIN(REAL(Q2))
      DETM=81*T17*T17*T14*T9-((T16+J30X)*(T0*T0-(T14*(T15+2*T10-J10Y-(T8
     . *M10)))))
```

Number of operations after optimization is:

```
Number of (+/-) operations: 13Number of unary - operations: 0Number of * operations: 18Number of integer ^ operations: 0Number of / operations: 0Number of function applications: 4
```

We can also use this example to show that correctness of results is easily verified. When storing the result of a SCOPE application in a file, it is of course possible to read the result in again. Then we apply a normal RE-DUCE evaluation strategy. This implies that all references to cse-names are automatically replaced by their values. We show the "correctness" of SCOPE by storing the optimized version of the expanded form of the determinant of M, called detm1 in file out.1 and the result of a SCOPE-application on the unexpanded form, detm2, in file out.2, followed by reading in both files and by subtracting detm2 from detm1, resulting in the value 0. This is of course an ad hoc correctness-proof for one specific example. It is in fact another way of testing the code of the package. We show it as a direct continuation of the previous determinant calculations.

This example also serves to show that the OPTIMIZE command can be extended with the OUT option. The keyword OUT has to be followed by a file-name. This file is properly closed and left in a readable form, assuming printing is produced in a OFF NAT fashion. SCOPE's file management features are discussed in more detail in section 5.

```
OFF ACINFO, FORT, NAT$ ON EXP$
```

```
OPTIMIZE detm1:=:det(M) OUT "out.1" INAME s;
OFF EXP$
OPTIMIZE detm2:=:det(M) OUT "out.2" INAME t;
ON NAT$
IN "out.1","out.2"$
detm1-detm2;
0
```

So far we presented via some examples straightforward algebraic mode use. The output is produced as a side-effect. However, optimization results can easily be made operational in algebraic mode. The parameterless function

#### ARESULTS

delivers the result of the directly preceding OPTIMIZE command in the form of a list of equations, corresponding with the sequence of assignment statements, shown either in REDUCE syntax or in the syntax of one of GEN-TRAN's target languages. But, we need to operate carefully. Application of a variety of assignment operators can easily bring in identifiers, representing earlier produced algebraic values. They will be substituted automatically, when referenced in rhs's in the list, produced with an ARESULTS application. Therefore, we implemented a protection mechanism. Before delivering output produced by ARESULTS, we make the system temporarily deaf for such references. The possibly game-spoiling algebraic values are stored at a seemingly anonymous place. All identifiers, subjected to this special treatement, can be made visible with the command

#### RESTORABLES;

Their original status can be restored, either globally with the command

#### **RESTOREALL\$**

or selectively with the instruction

ARESTORE <subsequence>\$

This subsequence is built with names, selected from the list of RESTORABLES, and separated by comma's. Information restoration is only possible before the next OPTIMIZE command.

#### Example 9

The use of these commands is now illustrated. A further explanation is given in the form of comment in the script.

```
u:=a*x+2*b$ v:=sin(u)$ w:=cos(u)$ f:=v^2*w;
                                  2
f := cos(a*x + 2*b)*sin(a*x + 2*b)
OFF EXP$
OPTIMIZE f:=:f,g:=:f^2+f INAME s;
s3 := x*a + 2*b
s2 := sin(s3)
f := s2*s2*cos(s3)
g := f*(f + 1)
alst:=ARESULTS;
alst := {s3=a*x + 2*b,
         s2=sin(s3),
                     2
         f=cos(s3)*s2,
         g=(f + 1)*f
% ---
\% SCOPE is made deaf for the standard reference mechanism for algebraic
```

% variables. However the rhs's in the list alst are simplified before % being shown. It explains the differences between the layout in the % alst items and the results, presented by the OPTIMIZE-command itself. % ---

```
RESTORABLES;
{f}
f;
f
ARESTORE f$
f;
                             2
\cos(a*x + 2*b)*\sin(a*x + 2*b)
% ----
\% f is re-associated with its original value. This can lead to a modified
% presentation of some of the rhs's of alst.
% ---
alst;
{s3=a*x + 2*b,
s2=sin(s3),
             2
 f=cos(s3)*s2,
                                2
                                                                      2
 g=(cos(a*x + 2*b)*sin(a*x + 2*b) + 1)*cos(a*x + 2*b)*sin(a*x + 2*b) }
OPTIMIZE f:=:f,g:=:f^2+f INAME s;
s3 := x*a + 2*b
s2 := sin(s3)
f := s2*s2*cos(s3)
g := f*(f + 1)
alst2:=ARESULTS$
OPTIMIZE f:=:f,g:=:f^2+f INAME s;
g := f*(f + 1)
```

```
% ----
% The algebraic value, which was associated with f, is permanently
% lost. It ought to be restored before a new OPTIMIZE command is given.
% Therefore f:=:f produced an identity, which is redundant in terms of
% code production. More details about removal of redundant code are
% given in section 7, when discussing data dependencies and related topics.
% ----
RESTOREALL$
f;
f
```

## 3.2 The function ALGOPT: Straightforward use

The function ALGOPT accepts up to three arguments. It can be used in stead of the OPTIMIZE command. It returns the optimization result, like ARESULTS, in the form of a list of equations. Since the ARESULTS mechanism is applied as well, the pre-ALGOPT-application situation can be restored with RESTOREALL or partly and selective with ARESTORE, using information, providable by an application of the function RESTORABLES.

The first argument of ALGOPT, like the other two optional, is the equivalent of the alglist or alglist\_production in the earlier introduced syntax of a SCOPE\_application. The second argument can be used to inform SCOPE that input from file(s) have to be processed. We survey SCOPE's file management features in section 5. So we omit a further discussion now. The last argument correspondents with the cse\_prefix of the INAME option of the OPTIMIZE command. The extension of the SCOPE\_application syntax, needed to include possible ALGOPT activities, is:

```
<SCOPE_application> ::= ··· |
ALGOPT(<a_object_list>[,<string_id_list>][,<cse_prefix>]) |
ALGOPT([<a_object_list>,]<string_id_list>[,<cse_prefix>]))
<a_object_list> ::= <a_object> | {<a_object>[,<a_object_seq>]}
<a_object_seq> ::= <a_object>[,<a_object_seq>]
<a_object> ::= <id> | <alglist> | <alglist_production> | {<a_object>}
```

We require at least one actual parameter, here the a\_object\_list. Its syntactical structure allows to apply a GENTRAN-like repertoire in an algebraic mode setting. The a\_object's can either be an alglist identifier, an alglist producing function application, or an alglist itself. An alglist identifier can be either a scalar or a matrix or array entry. The alglist producing functions will be discussed in section 4. An algebraic that the structure of an algebraic mode list; its elements are either a\_object's or equations of the form lhs = rhs. Such equations correspondent with the "take literal" GENTRAN operator := facility in the setting of an OPTIMIZE command (see also section 4 for a further discussion). The alternatives, i.e. uses of ::=, :=: or ::=:, are also covered by the a\_object syntax. The examples, given in this subsection, show that simplification of an algebraic list of equations leads to right hand side simplification, corresponding with the effect of the colon-addedto-the-right-extension of the assignment operator. However, as illustrated by example 13, some care has to be taken when operating in OFF EXP mode. Turning ON the switch EVALLHSEXP, can lead to lhs evaluations, corresponding with the extra-colon-to-the-left strategy. But we have to be aware of the instanteneous evaluation mechanism for matrix and array entries, when referenced.

We present some examples of possible use of the ALGOPT function. In example 10 a straightforward application is given. In example 11 follows an illustration of a possible strategy concerning optimizing sets of array- and/or matrix-entries. Then, in example 12, possible SCOPE assistance in problem analysis is shown. Finally in example 13 some differences in simplification and their influence on optimization are discussed. We also introduce and explain the role of the SCOPE switch SIDREL.

#### Example 10

A number of possible alglist elements is presented in the script. The first three elements of the actual parameter define values, obtained via the usual algebraic mode list evaluation mechanism. The last two will be processed literally. So, the actual parameter for ALGOPT is composed of the scalar alist, a list consisting of the matrix element m(1,1), the array element ar(2,2), nested even deeper, and two equations. Before an ALGOPT argument is optimized it is flattened by the SCOPE parser into a list of equations, the algebraic mode equivalent of the sequence of assignments in the OPTIMIZE context. Evaluation of an ALGOPT application leads to an algebraic mode list of equations, with optimized rhs's. The cse\_prefix was seemingly su-

perfluous, because all its references disappeared by back-substitution before output-processing started. See also example 13.

Since an ALGOPT application always results in an algebraic mode list, one can not use this feature for production of code in one of GENTRAN's target languages. To facilitate the translation of the result of an ALGOPT application, we extended the syntax of the OPTIMIZE input repertoire, such that alglst\_production's are processable by OPTIMIZE as well, as illustrated in the script of this example and in example 13

```
OFF EXP$
ARRAY ar(2,2)$ MATRIX m(2,2)$
alst:={p1=a+b,p2=(a+b)^2}$
m(1,1):={q1=c+d,q2=(c+d)^2}$
ar(2,2):={r1=(a+b)*(c+d),r2=(a+b)^2*(c+d)^2}$
optlst:=ALGOPT({alst,{m(1,1)},{{ar(2,2)}},
                t1=(a+b)*(c+d)^2,t2=(c+d)*(a+b)^2},s);
optlst := {p1=a + b,
                2
           p2=p1 ,
           q1=c + d,
                2
           q2=q1 ,
           r1=p1*q1,
                2
           r2=r1 ,
           t1=q1*r1,
           t2=p1*r1}
OPTIMIZE optlst$
p1 := a + b
p2 := p1*p1
q1 := c + d
q2 := q1*q1
r1 := q1*p1
r2 := r1*r1
t1 := r1*q1
t2 := r1*p1
```

#### Example 11

In example 8 we introduced a symmetric (3,3)-matrix m. We present an alternative computation of its determinant. We start with building a list of equations, with rhs's, being the non-zero entries of m, relevant for the computation. The lhs's are produced with the mkid function. These newly generated names are assigned to the matrix-entries as well. Finally we add the definition of the computation of the determinant of m, in terms of the redefined entries, to this list. For the construction of the value of mlst we applied both the lhs and rhs evaluation mechanism. Observe also that, due to the redefinition process, the original values of the entries of the matrix m are lost. We can optimize mlst, using either an OPTIMIZE command or an ALGOPT application. The reduction in arithmetic is not yet impressive here, certainly comparing it with the non-expanded, optimized form in the earlier example 8. (See also example 13 for additional comment). However, working with larger and non-symmetric matrices will certainly improve results, when applying a comparable strategy.

Observe that the syntax of permissible ALGOPT a\_object's does not allow to use matrix or array names to compactly identify the complete set of their entries. The script in this example shows that such a facility is easily made. This possibility exists already for matrices in a GENTRAN setting (see also example 22 in section 8).

```
% ----
% We assume the matrix m to be known already.
% ----
mlst:={}$ l:=-1$
OFF EXP$ ON EVALLHSEQP$
FOR j:=1:3 DO FOR k:=j:3 DO
IF m(j,k) neq 0 THEN
<< s:=mkid(t,l:=l+1);
    mlst:=append(mlst,{s=m(j,k)});
    m(j,k):=m(k,j):=s
>>$
```

OFF EVALLHSEQP\$

```
m;
[t0 t1 t2]
Γ
          ]
[t1 t3 0]
[
         ]
[t2 0 t4]
mlst:=append(mlst,{detm=det(m)});
                                     2
                                               2
mlst := {t0= - (j30y - j30z + 9*m30*p )*sin(q3)
                                                           2
         + 18*cos(q2)*cos(q3)*m30*p + j10y + j30y + m10*p
                   2
          + 18*m30*p ,
                                     2
                                               2
         t1= - (j30y - j30z + 9*m30*p )*sin(q3)
                                                     2
                                   2
         + 9*cos(q2)*cos(q3)*m30*p + j30y + 9*m30*p ,
                                      2
        t2 = -9 * sin(q2) * sin(q3) * m30 * p,
                                     2
                                              2
                                                                 2
         t3= - (j30y - j30z + 9*m30*p )*sin(q3) + j30y + 9*m30*p ,
                          2
        t4=j30x + 9*m30*p ,
                         2
                                 2
        detm=(t0*t3 - t1 )*t4 - t2 *t3}
ON ACINFO, FORT$ OFF PERIOD$
OPTIMIZE mlst INAME s;
Number of operations in the input is:
```

```
Number of (+/-) operations: 19Number of unary - operations: 1Number of * operations: 33Number of integer ^ operations: 16Number of / operations: 0Number of function applications: 9
```

```
S0=SIN(REAL(Q3))

S7=P*P

S5=S7*M30

S4=S5*COS(REAL(Q3))*COS(REAL(Q2))

S13=9*S5

S11=(S13+J30Y-J30Z)*S0*S0

T0=J30Y+J10Y+18*(S4+S5)+S7*M10-S11

T3=S13+J30Y-S11

T1=T3+9*S4

T2=-(S13*SIN(REAL(Q2))*S0)

T4=S13+J30X

DETM=T4*(T3*T0-(T1*T1))-(T2*T2*T3)
```

Number of operations after optimization is:

```
Number of (+/-) operations: 13Number of unary - operations: 1Number of * operations: 17Number of integer ^ operations: 0Number of / operations: 0Number of function applications: 4
```

### Example 12

We now illustrate that information, produced by SCOPE, can possibly also play a role in computations in algebraic mode. Let  $A.\vec{x} = \vec{b}$  be given by

Γ	-1	2	-2	1	3	2 ]		$\begin{bmatrix} x1 \end{bmatrix}$		5	1
	-2	4	-4	2	-2	3		x2		1	
	1	1	1	1	2	4		x3		10	
	2	-2	-1	1	-1	-2	•	x4	=	-3	.
	3	1	-4	1	1	2		x5		4	
L	-1	-5	1	1	3	6		<i>x</i> 6		5	

This artificial system is constructed for illustrative purposes. Its solution is

simply  $x_i = 1, i = 1, ..., 6$ . But straightforward inspection shows that

$$A = \begin{bmatrix} A_1 & A_2 \\ A_3 & A_4 \end{bmatrix} \text{ where } A_1 = \begin{bmatrix} -1 & 2 & -2 & 1 \\ -2 & 4 & -4 & 2 \end{bmatrix} \text{ and } A_4 = \begin{bmatrix} 2 & 4 \\ -1 & -2 \\ 1 & 2 \\ 3 & 6 \end{bmatrix}.$$

We can use ALGOPT to "discover" and thus to employ this information. The system is introduced in the form of assignment statements  $e_i = (\sum_{j=1}^6 a_{ij} \cdot x_j) - b_i$ ,  $i = 1, \dots, 6$ . We use alst, identifying the set of equations (see command 9), as actual parameter for ALGOPT, leading to an algebraic list, identified by reslst (see command 10). We recognize g2 = g6 + x5 (= x5 + 2x6) and g1 = g3 + g5 + -2x3 (= -x1 + 2x2 - 2x3 + x4). Through command 12 we require cse's to have an arithmetic complexity of a least 4. We then find g1 directly, now called g8, because we continue applying the function gensym; the cse\_prefix was left out as actual parameter. The solve function is applied (command 14) to obtain rootset1, a list of values for x5 and x6, expressed in the parameter g8. After assigning g8 its value in algebraic mode and resetting the algebraic values of ei,  $i = 1, \dots, 6$  with RESTOREALL instructions (the commands 11 and 16), we can obtain the solution of the subsets, denoted by rootset1 and rootset2.

- 1: LOAD\_PACKAGE nscope\$
- 2: e1:=2\*x6+3\*x5+x4-2\*x3+2\*x2-x1-5\$
- 3: e2:=3\*x6-2\*x5+2\*x4-4\*x3+4\*x2-2\*x1-1\$
- 4: e3:=2\*x5+4\*x6+x1+x2+x3+x4-10\$
- 5: e4:=-x5-2\*x6+2\*x1-2\*x2-x3+x4+3\$
- 6: e5:=x5+2\*x6+3\*x1+x2-4\*x3+x4-4\$
- 7: e6:=3\*x5+6\*x6-x1-5\*x2+x3+x4-5\$
- 8: solve({e1,e2,e3,e4,e5,e6},{x1,x2,x3,x4,x5,x6});

 $\{x1=1, x2=1, x3=1, x4=1, x5=1, x6=1\}$ 

- 9: alst:={e1=e1,e2=e2,e3=e3,e4=e4,e5=e5,e6=e6}\$
- 10: reslst:=ALGOPT alst;

```
reslst := \{g3 = -x1 + x4,
           g5=2*x2,
           g1=g3 + g5 - 2*x3,
           g6=2*x6,
           e1=g1 + g6 + 3*x5 - 5,
           e2=2*g1 - 2*x5 + 3*x6 - 1,
           g2=g6 + x5,
           g4=x2 + x4,
           e3=2*g2 + g4 + x1 + x3 - 10,
           e4= - g2 - g5 + 2*x1 - x3 + x4 + 3,
           e5=g2 + g4 + 3*x1 - 4*x3 - 4,
           e6=3*g2 + g3 - 5*x2 + x3 - 5
11: RESTOREALL$
12: SETLENGTH 4$
13: reslst:=ALGOPT alst;
reslst := {g8= - x1 + 2*x2 - 2*x3 + x4,
           e1=g8 + 3*x5 + 2*x6 - 5,
           e^{2}=2*g^{3} - 2*x^{5} + 3*x^{6} - 1,
           e3=x1 + x2 + x3 + x4 + 2*x5 + 4*x6 - 10,
           e4=2*x1 - 2*x2 - x3 + x4 - x5 - 2*x6 + 3,
           e5=3*x1 + x2 - 4*x3 + x4 + x5 + 2*x6 - 4,
           e6= -x1 - 5*x2 + x3 + x4 + 3*x5 + 6*x6 - 5
```

14: rootset1:=solve({part(reslst,2,2),part(reslst,3,2)},{x5,x6});

g8 + 13 - 8\*g8 + 13 rootset1 := {{x5=------,x6=-------}} 13 13 15: g8:=part(reslst,1,2); g8 := - x1 + 2\*x2 - 2\*x3 + x4 16: RESTOREALL\$ 17: rootset2:=solve(sub(rootset1,{e3,e4,e5,e6}),{x1,x2,x3,x4}); rootset2 := {{x1=1,x2=1,x3=1,x4=1}} 18: rootset1:=sub(rootset2,rootset1); rootset1 := {{x5=1,x6=1}}

Example 13

The script in example 11 suggests that we can easily copy GENTRAN's assignment features by some listprocessing in algebraic mode. However, we have to operate carefully. In the script of the present example we introduce an expression denoted by f. Production of a number of its partial (higher) derivatives is a straightforward mechanism to assist in constructing a set of assignment statements, containing lots of cse's. Inspection of the values, in OFF EXP mode assigned to faa, tst1 and tst2, respectively, learns that the value of mlst in example 11 may be improvable.

u:=a\*x+2\*b\$ v:=sin(u)\$ w:=cos(u)\$ f:=v^2\*w\$

OFF EXP\$

faa:=df(f,a,2);

2 2 2 2 faa := (2\*cos(a\*x + 2\*b) - 7\*sin(a\*x + 2\*b))\*cos(a\*x + 2\*b)\*x tst1:={faa=df(f,a,2)};

3 2

2 2

tst1 := {faa=2\*cos(a\*x + 2\*b) \*x - 7\*cos(a\*x + 2\*b)\*sin(a\*x + 2\*b) \*x }

tst2:={faa=(faa:=df(f,a,2))};

2 2 2 2 tst2 := {faa=(2\*cos(a\*x + 2\*b) - 7\*sin(a\*x + 2\*b) )\*cos(a\*x + 2\*b)\*x }

We produce an optimized version of the value of tlst, using ALGOPT. Switching ON INPUTC and PRIMAT results in an input echo, indeed showing expanded rhs's and a vizualized picture of the Sumscheme of  $D_{\lambda}$ . We skipped the D<sub>0</sub>-picture and the rest of the  $D_{\lambda}$ -picture from the script. The value of reslst shows the patterns s14 = -7.f.x + 2.s9.x and fbb = -28.f +8.s9. The presented Sumscheme of  $D_{\lambda}$  suggests that fbb and s13 (see the Fa(the)r entries) seemingly have nothing in common. But s14 stands for 2.s6 - 7.s7, because, column 8 has to be identified with s7, etc. Since both S6 and S7 occur only once in  $D_{\lambda}$ , their value replaces them in the output. It is an illustration of the heuristic character of the optimization process. Optimization of the value of reslst shows that the repeated pattern is now recognized.

```
ON INPUTC, PRIMAT$
```

```
reslst:=ALGOPT(tlst,s);
```

2 f := cos(a\*x + 2\*b)\*sin(a\*x + 2\*b)

 $2 \qquad 3$ fa := 2\*cos(a\*x + 2\*b) \*sin(a\*x + 2\*b)\*x - sin(a\*x + 2\*b) \*x
fb := 4\*cos(a\*x + 2\*b) \*sin(a\*x + 2\*b) - 2\*sin(a\*x + 2\*b)
faa := 2\*cos(a\*x + 2\*b) \*x - 7\*cos(a\*x + 2\*b)\*sin(a\*x + 2\*b) \*x
fab := 4\*cos(a\*x + 2\*b) \*x - 14\*cos(a\*x + 2\*b)\*sin(a\*x + 2\*b) \*x
fab := 4\*cos(a\*x + 2\*b) \*x - 14\*cos(a\*x + 2\*b)\*sin(a\*x + 2\*b) \*x

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3 2 fbb := 8\*cos(a\*x + 2\*b) - 28\*cos(a\*x + 2\*b)\*sin(a\*x + 2\*b) Sumscheme : | 3 4 5 6 7 8 9 10 11 12 32| EC|Far ------\_\_\_\_\_ 1 2| 1| s3 3| -28 8 20| | 1| fbb -7 2 | 1| s14 -1 2 | 1| s15 37| 38| \_\_\_\_\_ 3 : s3 4 : s15 5 : s14 6 : s4 7 : s9 8 : s7 9 : s6 10 : s10 11 : s5 12 : s8 32 : b reslst := {s3=a\*x + 2\*b, s0=cos(s3), 3 s9=s0 , s2=sin(s3), s12=s0\*s2, f=s12\*s2, 3 s15=2\*s0\*s12 - s2, fa=s15\*x, fb=2\*s15,

```
s14= - 7*f*x + 2*s9*x,
faa=s14*x,
fab=2*s14,
fba=fab,
fbb= - 28*f + 8*s9}
OFF INPUTC,PRIMAT$
OPTIMIZE res1st$
```

```
s3 := 2*b + x*a
s0 := cos(s3)
s9 := s0*s0*s0
s2 := sin(s3)
s12 := s2*s0
f := s12*s2
s15 := 2*s12*s0 - s2*s2*s2
fa := s15*x
fb := 2*s15
g3 := 2*s9 - 7*f
s14 := g3*x
faa := s14*x
fab := 2*s14
fba := fab
fbb := 4*g3
```

Repeating this process, this time with an OPTIMIZE command to begin with, learns that the OFF EXP mode is now effective. But this time, and for similar reasons, the assignments fbb = 4.s9.s0 and s12 = s9.s0.x still have a subexpression in common. Now the Productscheme of  $D_{\lambda}$  helps understanding the phenomenon; again we skipped for shortness the rest of the information, provided by the ON PRIMAT status of SCOPE. Internally s12 denotes the product s4.s9, where s4 = x.s0. The cse s4 disappeared from the output. An ALGOPT application leads to the "discovery" of the cse g10 = s0.s9.

2

 $f:=v^2*w;$ 

f := cos(a\*x + 2\*b)\*sin(a\*x + 2\*b)

ON INPUTC, PRIMAT\$

```
OPTIMIZE f:=:f,fa:=:df(f,a),fb:=:df(f,b),faa:=:df(f,a,2),
        fab:=:df(f,a,b),fba:=:df(f,b,a),fbb:=:df(f,b,2) INAME s$
                                 2
f := cos(a*x + 2*b)*sin(a*x + 2*b)
                                        2
                      2
fa := (2*cos(a*x + 2*b) - sin(a*x + 2*b) )*sin(a*x + 2*b)*x
                        2
                                         2
fb := 2*(2*cos(a*x + 2*b) - sin(a*x + 2*b) )*sin(a*x + 2*b)
                       2
                                          2
                                                              2
faa := (2*cos(a*x + 2*b) - 7*sin(a*x + 2*b) )*cos(a*x + 2*b)*x
                                             2
                         2
fab := 2*(2*cos(a*x + 2*b) - 7*sin(a*x + 2*b) )*cos(a*x + 2*b)*x
                         2
                                             2
fba := 2*(2*cos(a*x + 2*b) - 7*sin(a*x + 2*b))*cos(a*x + 2*b)*x
                        2
                                             2
fbb := 4*(2*cos(a*x + 2*b) - 7*sin(a*x + 2*b) )*cos(a*x + 2*b)
s3 := x*a + 2*b
s0 := cos(s3)
s2 := sin(s3)
s6 := s2*s2
f := s6*s0
s14 := 2*s0*s0
s13 := (s14 - s6) * s2
fa := s13*x
fb := 2*s13
s9 := s14 - 7*s6
s12 := s9*s0*x
faa := s12*x
fab := 2*s12
fba := fab
fbb := 4*s9*s0
Productscheme :
   0 2 3 4 5 12 14 18 19 20 21 22 23 EC Far
```

0  5  9  13  17  21  25  29  30	1	1 1	1	1			1				1 1 1 1	 1 1	             1	1  1  2  1  1  1  4  1	f fa fb faa fab fba fbb s4 s5
31  33  37  38  39  40			1	1	1	1			1	2	2			1  1  1  1  2  2	s6 s8 s12 s13 s14 s15
0 :: 2 :: 3 :: 4 :: 5 :: 12 :: 14 :: 18 :: 19 :: 20 :: 21 :: 22 :: 23 :: ALGO	s13 s12 s9 s10 s8 s6 s5 s4 s2= s0= x	sin	(s3)	)											
{s3=a*x + 2*b,															
s0=cos(s3),															
s6=	sin( 2 s2 , 0*s6		,												

2 s14=2\*s0, s13=(s14 - s6)\*s2, fa=s13\*x, fb=2\*s13, s9=s14 - 7\*s6, g10=s0\*s9, s12=g10\*x, faa=s12\*x, fab=2\*s12, fba=fab, fbb=4\*g10}

This script is shown for different reasons. It illustrates the heuristic character of the optimization process. We optimize, but do not guarantee the optimal solution. It also shows how easily repeated SCOPE applications can be accomplished. Hence commands like "ALGOPT ARESULTS;", "ALGOPT ALGOPT  $\cdots$ ;" or "OPTIMIZE ALGOPT  $\cdots$ ;" are all possible. However, it is sometimes better to avoid such a combination when a RESTOREALL instruction has to follow the first application. A more detailed discussion about these possibilities is given in section 4, and especially in section 4.1. An additional reason was, to stipulate that SCOPE's actual parameters have to be built carefully.

This example is also used to illustrate the role, which the switch SIDREL can possibly play. When turned it ON the finishing touch F (see subsection 2.3) is omitted and all non-additive cse's are substituted back, thus producing a possibly still rewritten input set, which consists of toplevel input and additive cse's only. A simple straightforward backsubstitution mechanism is applied on the optimization result before it is presented to the user. Seemingly, it can lead to surprises as shown below by the differences between the presentations of s15, s14 and fbb when again optimizing the contents of

### tlst. This effect disappeares when using SETLENGTH.

The switch SIDREL was introduced in SCOPE quite long ago. By that time Hearn was wondering [18, 19] if (parts of) SCOPE output, presented in algebraic mode, can be used as input for a Gröbner-base algorithm application, thus attempting to assist in expression restructuring leading to improved expression representations.

```
ON SIDREL$
ALGOPT(tlst,s);
\{s3=a*x + 2*b,
                  2
 f=\cos(s3)*\sin(s3),
              2
                                  3
 s15=2*cos(s3) *sin(s3) - sin(s3) ,
 fa=s15*x,
 fb=2*s15,
                                        2
                                                        3
 s14= - 7*cos(a*x + 2*b)*sin(a*x + 2*b) *x + 2*cos(s3) *x,
 faa=s14*x,
 fab=2*s14,
 fba=2*s14,
                                         2
                                                       3
 fbb= - 28*cos(a*x + 2*b)*sin(a*x + 2*b) + 8*cos(s3) }
SETLENGTH 4$
ALGOPT(tlst,s);
                                 2
{f=cos(a*x + 2*b)*sin(a*x + 2*b) ,
                     2
                                                        3
 s15=2*\cos(a*x + 2*b) *\sin(a*x + 2*b) - \sin(a*x + 2*b),
```

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```
fa=s15*x,
fb=2*s15,
s14=2*\cos(a*x + 2*b)^{3}*x - 7*\cos(a*x + 2*b)*\sin(a*x + 2*b)^{2}*x,
faa=s14*x,
fab=2*s14,
fbb=2*s14,
fbb=8*cos(a*x + 2*b)^{3} - 28*cos(a*x + 2*b)*sin(a*x + 2*b)^{2}}
```

# 4 Special SCOPE 1.5 Features

Part of the input syntax for the function ALGOPT was left undiscussed in section 3.2. It was the permissable form for (parts of) the actual parameter, defining function applications, producing an alglist. The alglist is an algebraic mode list, consisting either of equations of the form lhs = rhs or of constructs evaluating into an alglist or referencing an alglist. In section 4.1 is explained which type of user defined functions lead to permissable function applications as (part of an) actual parameter for an OPTIMIZE command or an ALGOPT application. Tools are provided for building a SCOPE library. Already available facilities, designed along these lines, cover structure recognition, presented in section 4.2 and Horner-rule based expression rewriting, surveyed in section 4.3.

# 4.1 Towards a SCOPE 1.5 Library

Design and implementation of an algebraic or symbolic procedure, returning a list of equations in algebraic mode, is straightforward as long as the number of formal parameters is exactly known. Let us call such procedures functions of NORMAL-type. When formal parameters are not required, the so-called ENDSTAT-variant can be used. One simply associates an indicator stat, with value endstat, with the function name f-name, using lisp(put('f-name,'stat,'endstat))\$ as instruction. Such a function will be said to be of ENDSTAT-type.

The so-called PSOPFN-type function is similar to the FEXPR-type function in symbolic mode. It may have an arbitrary number of unevaluated parameters. Special attention is made possible by modifying the function reval1, used in both reval and aeval. The relevant section of the evaluator is:

The actual parameter u of reval1 is a function application in prefixform: (function-name arg1 ... argn). The function-name is replaced by the value of the indicator psopfn. The thus modified S-expression is evaluated. This mechanism leaves control over evaluation of (part of) the arguments, collected in cdr u, to the designer of the function hidden behind the psopfn value. We employed this simple mechanism to implement ALGOPT and some of the features, to be discussed in section 4.2 and section 4.3. A possible re-evaluation step, based on the use of the value of the indicator cleanupfn was not necessary; it is not yet allowed in the SCOPE context.

The different combinations, suggested in example 13, such as ALGOPT ALGOPT or ALGOPT ARESULTS, are merely examples of a general rule:

ENDSTAT-, NORMAL- and PSOPFN-type functions, delivering an alglist when applied, are all applicable as actual parameter or as element of an alglist, functioning as actual parameter in an OPTIMIZE command or an ALGOPT application.

So, in principle, special features, providing a form of preprocessing, can be designed and implemented as extension of the default optimization repertoire. Of course additional function types are conceivable. We illustrate the potential of this facility with a simple example. Further examples follow in section 4.2 and section 4.3.

#### Example 14

The procedures asquares and repeated\_squaring define the production of lists of equations. The lhs's function as name and the rhs's as the computational rules. Application of these functions shows how easy a user can provide new features, usable in a SCOPE context. The procedures asquares and repeated\_squaring are essentially different The first has one parameter, a list of equations, while the latter accepts an arbitrary number of such lists as actual parameters. The psopfn indicator value is repeated\_squaringeval, the name of the function, which is actually introduced. asquares is of NORMAL-type and applicable in both algebraic and symbolic mode.

**OPERATOR** square\$

sq\_rule:={square(~u) => u^2}\$
ALGEBRAIC PROCEDURE asquare u;
square(u) WHERE sq\_rule\$

```
SYMBOLIC PROCEDURE rsquare u;
reval asquare aeval u$
SYMBOLIC PROCEDURE asquares u;
 append(list('list),
        FOREACH el IN cdr u COLLECT list('equal,cadr el,rsquare caddr el))$
SYMBOLIC OPERATOR asquares$
SYMBOLIC PROCEDURE repeated_squaringeval u;
BEGIN SCALAR res; INTEGER j;
 j=0;
 FOREACH el IN u DO
  << j:=j+1; el:=asquares el;
    FOR k:=2:j DO el:=asquares el;
    res:= IF j=1 THEN el ELSE append(res,cdr el)
 >>;
 RETURN res
END$
LISP( put('repeated_squaring,'psopfn,'repeated_squaringeval))$
% ----
% Examples of the use of asquares and repeated_squaring.
% Although the psopfn-mechanism can be easily avoided,
% it is used for illustrative purposes here.
% ----
OFF EXP$
asquare sin(x);
      2
sin(x)
LISP(rsquare('(sin x)));
(expt (sin x) 2)
asq:=asquares {s1=a+b,s2=(a+b)^2,s3=(a+b)^3};
                  2
                            4
                                          6
asq := {s1=(a + b) ,s2=(a + b) ,s3=(a + b) }
```

```
repeated_squaring({s1=a+b,s2=(a+b)^2},{s3=(a+b)^3,s4=(a+b)^4}, {s5=(a+b)^5,s6=(a+b)^6});
```

```
2
{s1=(a + b) ,
         4
s2=(a + b) ,
         12
s3=(a + b) ,
         16
s4=(a + b) ,
        40
s5=(a + b) ,
          48
s6=(a + b) }
% ---
\% The "ALGOPT asquares ...;" application is similar to the "ALGOPT asq;"
% instruction.
% ---
ALGOPT(asquares {s1=a+b,s2=(a+b)^2,s3=(a+b)^3},t);
              2
                    2
{t2=a + b,s1=t2 ,s2=s1 ,s3=s1*s2}
ALGOPT asq;
```

```
2
                      2
{g6=a + b,s1=g6 ,s2=s1 ,s3=s1*s2}
% ---
% The OPTIMIZE variant is now applied on a repeated_squaring application.
% ----
OPTIMIZE repeated_squaring({s1=a+b,s2=(a+b)^2},{s3=(a+b)^3,s4=(a+b)^4},
                           {s5=(a+b)^5,s6=(a+b)^6}) INAME t;
t5 := a + b
s1 := t5*t5
s2 := s1*s1
t12 := s2*s2
s3 := s2*t12
s4 := s2*s3
s5 := t12*t12*t12*s4
s6 := t12*s5
```

### 4.2 Structure Recognition: GSTRUCTR and ALGSTRUCTR

The structr command in REDUCE 3.7 (see the manual, section 8.3.8) can be used to display the skeletal structure of its evaluated argument, a single expression. After setting ON SAVESTRUCTR a structr command will return a list, whose first element is a presentation for the expression and subsequent elements are the subexpression relations.

A special SCOPE feature provides an extended display facility, called GSTRUCTR. The syntax of this generalized command is:

$<$ REDUCE_command>	::=	
		$\texttt{GSTRUCTR} < \texttt{stat\_group} > [\texttt{NAME} < \texttt{cse\_prefix} > ]$
<stat_group></stat_group>	::=	$\ll$ <stat_list> <math>\gg</math></stat_list>
$<$ stat_list $>$	::=	$\langle gstat \rangle [; \langle stat\_list \rangle]$
$\langle gstat \rangle$	::=	$<\!\!\mathrm{name}\!\!> := <\!\!\mathrm{expression}\!\!> \mid <\!\!\mathrm{matrix\_id}\!\!>$

The stat\_group consists of one assignment statement or a group of such statements. Application of a GSTRUCTR command provides a display of the structure of the whole set of assignments. Such an assignment can be replaced by a matrix reference. That leads to the display of all the non-zero entries of the referenced matrix as well. The NAME part is optional. The cse-name mechanism is applied in the usual way.

The equivalent of a possible ON SAVESTRUCTR setting is provided in the form of a PSOPFN-type function, called ALGSTRUCTR. Its syntax is:

$<$ function_application>	::=	ALGSTRUCTR ( <arg_list> [, <cse_prefix> ])</cse_prefix></arg_list>
<arg_list></arg_list>	::=	$< arg\_list\_name >   {< arg\_seq >}$
<arg_seq></arg_seq>	::=	$< arg > [, < arg\_seq >]$
<arg></arg>	::=	<matrix_id>   <name>=<expression></expression></name></matrix_id>
$< arg\_list\_name >$	::=	<id></id>

The result is presented in the form of an algebraic mode list.

Earlier SCOPE-versions allowed to use a GSTRUCTR command as (part of an) actual parameter for an OPTIMIZE command. This facility is not longer supported. In stead, an ALGSTRUCTR application can now be used as (part of an) actual parameter in both an OPTIMIZE command or an ALGOPT application. We now illustrate these features in:

#### Example 15

The script hardly requires explanation. However, observe that v1, v3, v4, v6 and v7 occur only once in the result of the GSTRUCTR application. When this application is used as actual parameter for an OPTIMIZE command these redundancies are removed before the actual optimization process starts. Likewise, an ALGSTRUCTR application only leads to identification of repeatedly occuring sub-structures in its input. ALGSTRUCTR, ALGHORNER, see the next subsection, and ALGOPT all apply the same output production strategy, i.e. it might be necessary to restore the previous algebraic mode status by applying the function RESTOREALL.

```
OFF EXP, PERIOD$
```

MATRIX a(2,2);

```
a:=mat((x+y+z,x*y),((x+y)*x*y,(x+2*y+3)^3-x));
```

```
 a := \begin{bmatrix} x + y + z & x*y & ] \\ [ & & ] \\ a := \begin{bmatrix} & 3 & ] \\ [(x + y)*x*y & (x + 2*y + 3) & -x] \end{bmatrix}
```

```
GSTRUCTR <<a;b:=(x+y)^2;c:=(x+y)*(y+z);d:=(x+2*y)*(y+z)*(z+x)^2>> NAME v$
```

```
a(1,1) := v1
a(1,2) := x*y
a(2,1) := v2*x*y
a(2,2) := v4
      2
b := v2
c := v2*v5
        2
d := v6*v7 *v5
  where
     v7 := x + z
     v6 := x + 2*y
     v5 := y + z
             3
     v4 := v3 - x
     v3 := x + 2*y + 3
     v2 := x + y
     v1 := x + y + z
ALGSTRUCTR({a,b=(x+y)^2,c=(x+y)*(y+z),d=(x+2*y)*(y+z)*(z+x)^2},v);
\{a(1,1)=x + y + z,
a(1,2)=x*y,
v2=x + y,
 a(2,1)=v2*x*y,
                  3
 a(2,2)=(x + 2*y + 3) - x,
    2
 b=v2 ,
 v5=y + z,
 c=v2*v5,
                   2
 d=(x + 2*y)*(x + z) *v5
```

```
RESTORABLES;
```

{a}

ARESTORE a\$

```
alst:=
 ALGOPT(ALGSTRUCTR({a,b=(x+y)^2,c=(x+y)*(y+z),d=(x+2*y)*(y+z)*(z+x)^2},v),s);
*** a declared operator
alst := \{s5=x + z,
         a(1,1)=s5 + y,
         a(1,2)=x*y,
         v2=x + y,
         a(2,1)=a(1,2)*v2,
         s6=x + 2*y,
         s4=s6 + 3,
                  3
         a(2,2)=s4 - x,
             2
         b=v2 ,
         v5=y + z,
         c=v2*v5,
             2
         d=s5 *s6*v5}
% ----
\% The above delivered warning is caused by the decloupling of a and its
\% status as matrix. Therefore a(1,2) can function in the rhs of a(2,1).
% After an ARESTORE instruction a can restart its life as matrix_id.
```

% ---

a; a ARESTORE a\$ a; [ x + y + z x\*y ] [ ] [ 3 ] [(x + y)\*x\*y (x + 2\*y + 3) - x]

## 4.3 Horner-rules: GHORNER and ALGHORNER

Horner-rule based expression modification is a SCOPE facility, called GHORNER. The syntax of the command is similar to the GSTRUCTR syntax:

<REDUCE\_command> ::= ··· | GHORNER <stat\_group> [VORDER <id\_seq>];

The VORDER part is optional. Application of a (generalized) Horner-rule assumes an identifier ordering. The syntax of the identifier sequence is:

 $\langle id\_seq \rangle ::= \langle id \rangle [, \langle id\_seq \rangle].$ 

We assume the rhs's in the stat\_group to be polynomials in the identifiers, partly or completely given in the id\_seq. The left-to-right ordering of this sequence replaces the existing system identifier ordering. Identifiers, omitted from the **vorder** sequence have a lower preference and follow the existing system ordering. The rewritten rhs's are presented as a side-effect. FORTRAN notation is of course permitted. It is simply an extended print facility.

The PSOPFN-type variant of the GHORNER command is called ALGHORNER. Its syntax is:

<function\_application> ::= ··· | ALGHORNER (<arg\_list> [,{<id\_seq>}])

The syntax for the arg\_list can be found in subsection 4.2. The result is presented in the form of an algebraic mode list. An ALGHORNER application can be used as (part of an) actual parameter of either an OPTIMIZE command

or an ALGOPT application.

#### Example 16

We illustrate the Horner-facilities by rewriting the expression of example 2, before optimizing it. Observe that application of ALGHORNER in the default algebraic mode setting is useless. Due to the algebraic mode regime the rewritten expression is expanded again. We also show some Taylor-series remodelling.

#### ON EXP\$

```
z:=a<sup>2</sup>*b<sup>2</sup>+10*a<sup>2</sup>*m<sup>6</sup>+a<sup>2</sup>*m<sup>2</sup>+2*a*b*m<sup>4</sup>+2*b<sup>2</sup>*m<sup>6</sup>+b<sup>2</sup>*m<sup>2</sup>;
      2 2
                  2 6
                          2 2
                                                 2 6
                                                          2 2
                                         4
z := a *b + 10*a *m + a *m + 2*a*b*m + 2*b *m + b *m
GHORNER z:=z VORDER a;
         2 6
                  2 2
                                           2
                                   4
                                                    6
                                                          2
z := (2*b *m + b *m) + a*(2*b*m + a*(b + 10*m + m))
GHORNER z:=z VORDER b;
          2 6 2 2
                                         2
                                  4
                                                  6
                                                          2
z := (10*a *m + a *m ) + b*(2*a*m + b*(a + 2*m + m ))
hlst:={z=z}
ALGHORNER(hlst, {a,b,m});
               2 6 2 2
    2 2
                                       4
                                             2 6 2 2
{z=a *b + 10*a *m + a *m + 2*a*b*m + 2*b *m + b *m }
OPTIMIZE ALGHORNER(hlst, {a,b,m}) INAME s;
s1 := m*m
s0 := s1*s1
s2 := b*b
s4 := 2*s0
z := a*(a*(s2 + s1*(10*s0 + 1)) + s4*b) + s2*s1*(s4 + 1)
OPTIMIZE ALGHORNER(hlst, {b,m}) INAME s;
s2 := m*m
```

```
s0 := s2*s2
s1 := a*a
s4 := 2*s0
z := b*(b*(s1 + s2*(s4 + 1)) + s4*a) + s2*(s1 + 10*s1*s0)
```

% Hornering Taylor-series: PROCEDURE taylor(fx,x,x0,n); sub(x=x0,fx)+(FOR k:=1:n SUM(sub(x=x0,df(fx,x,k))\*(x-x0)^k/factorial(k)))\$ hlst2:={f1=taylor(e^x,x,0,4),f2=taylor(cos x,x,0,6)}; 4 3 2 x + 4\*x + 12\*x + 24\*x + 24 hlst2 := {f1=-----, 24 4 2 6 - x + 30\*x - 360\*x + 720 f2=-----} 720 OPTIMIZE ALGHORNER(hlst2,{x}); 24 + x\*(24 + x\*(12 + x\*(4 + x)))f1 := -----24 g7 := x\*x 720 + g7\*(g7\*(30 - g7) - 360)f2 := -----720 ON ROUNDED\$ hlst2:=hlst2; 3 2 4 hlst2 := {f1=0.04166666666667\*x + 0.16666666666667\*x + 0.5\*x + x + 1, 4 2 6 f2= - 0.001388888888889\*x + 0.041666666666667\*x - 0.5\*x + 1 } OPTIMIZE ALGHORNER(hlst2, {x});

```
f1 := 1 + x*(1 + x*(0.5 + x*(0.04166666666667*x + 0.166666666667)))
g9 := x*x
f2 := 1 + g9*(g9*(0.0416666666667 - 0.0013888888889*g9) - 0.5)
```

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# 5 File Management and Optimization Strategies

Both the OPTIMIZE command and the ALGOPT function accept input from file(s). Obviously, this input ought to obey the usual syntactical rules, as introduced in the previous (sub)sections.

The OPTIMIZE command is designed as a syntactical extension of REDUCE itself, i.e. the meaning of its actual parameters is understood from the tokencontext in the command. However, an ALGOPT application requires one, two or three actual parameters without additional provisions or conditions. The ALGOPT facility is added to provide a simple, user friendly, algebraic mode tool. Therefore -in contrast with the OPTIMIZE command- it does not allow to direct output to a file; the default REDUCE features for dealing with output files can be applied. The previously given syntax requires some extensions:

<SCOPE\_application> ::= <OPTIMIZE command> | <ALGOPT application> <OPTIMIZE command> ::=

OPTIMIZE <object\_seq> [IN <file\_id\_seq>] [OUT <file\_id>] [INAME <cse\_prefix>] |

OPTIMIZE [<<br/>object\_seq>] IN <file\_id\_seq> [OUT <file\_id>] [INAME <cse\_prefix>]

<ALGOPT application> ::=

```
ALGOPT(<a_object_list>[,<string_id_list>][,<cse_prefix>]) |
ALGOPT([<a_object_list>,]<string_id_list>[,<cse_prefix>])
```

The different variations for the object\_seq and the a\_object\_list and the meaning of cse\_prefix are introduced in the subsections 3.1 and 3.2. The syntax of the file handling features is:

<file_id_seq></file_id_seq>	::=	$<$ file_id> [, <file_id_seq>]</file_id_seq>
<file_id></file_id>	::=	$<$ id $>$   $<$ string_id $>$
$<$ string_id_list $>$	::=	$<$ string_id>   { <string_id_seq>}</string_id_seq>
$<$ string_id_seq $>$	::=	$<$ string_id> [, $<$ string_id_seq>]
<string_id></string_id>	::=	$\ll d \gg \parallel \ll d > . < f_extension > $

The differences in input-file management are introduced for practical reasons. As stated above, the ALGOPT function can have up to three arguments. To be able to distinguish the optional second argument from the first and the last requires file-names to be given in the form of strings. The OPTIMIZE command follows the ordinary REDUCE rules for file names.

File management can be used as a tool for input partioning. If m > 1 then  $N^m > \sum_{i=1}^k n_i^k$  for positive integers N and  $n_i$ , such that  $N = \sum_{i=1}^k n_i$ . In view of the time-complexity of the optimization algorithm, it may be worth the effort to partition SCOPE input of size N in k partitions, of sizes  $n_i$ , i = 1, ..., k. We can start optimizing the contents of file fi.1, containing the initial  $n_1$ -sized piece of code, and store the result of this operation in file fo.1. Consecutive steps provide an optimization of the combined contents of the files fo.i and fi.(i+1), i=1,..., k-1. During this iterative process, or during variations of this strategy, it is better not to perform a finishing touch. The switch AGAIN, which is normally OFF, can be used, when set ON, to avoid this. The switch serves an additional purpose. When switched ON storage of partly optimized code in a file will include all relevant information, needed to restore the required status of system generated sub-expression names.

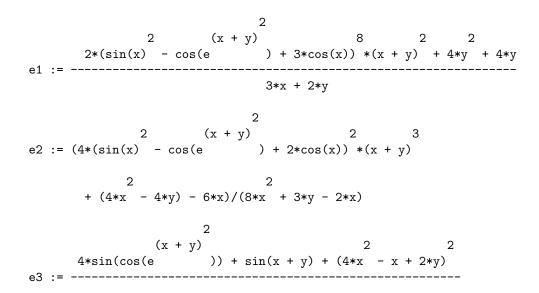
We illustrate SCOPE's file management facilities with example.

#### Example 17

We assume to have three files, called f1, f2 and f3. Each file contains only one assignment. We simply show different variations of the use of these files. With ON INPUTC the contents of the files is made visible.

#### ON INPUTC\$

OPTIMIZE IN f1,f2,f3 INAME s;



3\*y + f(x,g( - cos(x)))

```
s3 := sin(x)
s20 := x + y
s6 := s20*s20
    s6
s4 := cos(e)
s8 := cos(x)
s31 := s3*s3 - s4
s2 := s31 + 3*s8
s44 := s2*s2
s43 := s44*s44
s36 := 4*y
s34 := 2*y
s10 := s34 + 3*x
      s36 + s36*y + 2*s6*s43*s43
e1 := -----
               s10
s13 := s31 + 2*s8
s33 := 4*x*x
s30 := s33 - x
s35 := 3*y
```

We repeat the same process. However, this time we apply input partitioning. The switch AGAIN is turned ON. Output is redirected to the output file fo.1 in an OFF NAT fashion and ended with the required ;end; closure, thus made ready for re-use during a next step. The default mode of operation is OFF AGAIN and ON NAT. If the switch NAT is turned OFF file output is automatically ended by ;end;.

Due to the ON INPUTC effect we can also observe that the identifiers gsym and cses are apparently used to store relevant information about cse names.

ON AGAIN, INPUTC\$

```
2 2
+ (4*x - 4*y) - 6*x)/(8*x + 3*y - 2*x)
OFF AGAIN$
ALGOPT({"fo.2","f3"},u);
gsym := g0002
cses := t23 + t11 + t26 + t7 + t17 + t19
t19 := x + y
      2
t17 := t19
t7 := cos(x)
         2 t17
t26 := sin(x) - cos(e)
t11 := 3*x + 2*y
                8
           2
     4*y + 4*y + 2*(t26 + 3*t7) *t17
e1 := -----
                t11
      2
t23 := x
                                2
     4*t23 - 2*t11 + 4*t19*(t26 + 2*t7) *t17
e2 := -----
              8*t23 - 2*x + 3*y
                    2
                           2 2
             (x + y)
4*sin(cos(e )) + sin(x + y) + (4*x - x + 2*y)
e3 := -----
                 3*y + f(x,g( - cos(x)))
```

```
2
t26=u5 - u6,
u33=2*y,
t11=u33 + 3*x,
u10=t26 + 3*t7,
    2
u46=u10 ,
     2
u45=u46 ,
u35=4*y,
         2
  2*u20*u45 + u35*y + u35
e1=-----,
          t11
   2
t23=x ,
u13=t26 + 2*t7,
u36=4*t23,
u31=u36 - x,
u34=3*y,
               2
   - 2*t11 + 4*u13 *u20*u23 + u36
e2=-----,
          2*u31 + u34
u24=u31 + u33,
                       2
  sin(u23) + 4*sin(u6) + u24
e3=-----}
     f(x,g( - t7)) + u34
```

Observe that the initial characters of the sub-expression names indicate their moment of generation. We used **f** and **g** as operators. Therefore, a warning was produced ahead of the ALGOPT output. Since an OPTIMIZE command produces output as a side-effect these warnings were not given earlier.

# 6 Generation of Declarations

GENTRAN's DECLARE statement can be used as an optional extension of the OPTIMIZE command, and as illustrated in example 18. The syntax of such an extension is in accordance with the GENTRAN rules:

```
<OPTIMIZE command> ::=
    OPTIMIZE <object_seq> [IN <file_id_seq>] [OUT <file_id>]
        [INAME <cse_prefix>] [DECLARE <declaration_group>] |
    OPTIMIZE [<object_seq>] IN <file_id_seq> [OUT <file_id>]
        [INAME <cse_prefix>] [DECLARE <declaration_group>]
```

The syntax of the declaration\_group is:

$<$ declaration_group>	::=	$<$ declaration $>$   $\ll$ $<$ declaration_list $>$ $\gg$
$<$ declaration_list $>$	::=	$<$ declaration $>$ [; $<$ declaration_list>]
<declaration $>$	::=	<range_list>: IMPLICIT <type>   <id_list>:<type></type></id_list></type></range_list>
$<$ range_list $>$	::=	<range>[,<range_list>]</range_list></range>
<range></range>	::=	$\langle id \rangle   \langle id \rangle - \langle id \rangle$
<id_list></id_list>	::=	$<$ id>[, $<$ id_list>]
<type></type>	::=	integer   real   complex   real*8   complex*16

The symbol table features of GENTRAN are used. During the subtask R (see subsection 2.3) of an OPTIMIZE command evaluation, all typing information is installed in the symbol table. Once optimization is ready all relevant information for completing the declarations ought to be known, i.e. the contents of the symbol table and the result of the optimization operations, collected in prefix form in a list, called prefixlist. This prefixlist is employed do decide which not vet typed identifiers and system selected cse names have to be entered in the symbol table. We make use of earlier provided information, delivered via the DECLARE option, (sub)expression structure and the normal hierarchy in data types. The strategy to achieve this form of dynamic typing is based on chapter 6 of [1]. Once the table is completed a list of declarations is produced and precedes the other SCOPE output. SCOPE output is by default given in REDUCE notation. Therefore such lists of declarations are also given in REDUCE text. Incomplete initial typing information can lead to overtyping after optimization, such as complex in stead of real, for instance. It can therefore lead to erroneous results and even to an error message. A safe procedure is to use the DECLARE option of the <code>OPTIMIZE</code> command for typing all identifiers, occuring in the input set  $\mathrm{E}_0$ 

Alternative output can be obtained via an application of the function OPTLANG. This function accepts one argument from the set {fortran, c, ratfor, pascal<sup>1</sup>, f90, nil}. The fortran(77) choice can also be made by turning ON the switch FORT. The nil option is necessary if one wants to switch back to the usual REDUCE output. not yet generally available. The output modules of GENTRAN are used for producing formatted code in the user selected target language. The f90 option, for the production of fortran90 code, is not yet provided by the standard GENTRAN version [3].

Especially the above given syntax rules for typing require some additional explanation:

- The corresponding types in Fortran are integer, real, complex, double precision and complex\*16.
- The GENTRAN switch DOUBLE is automatically turned ON, when a type real\*8 or type complex\*16 is introduced in a DECLARE option. The same mode of operation is introduced when floating point numbers appear in SCOPE input. Fixed floats do not produce this side effect.
- When generating fortran code we have to be aware of a possibly existing statement length limitation. If one is afraid that a declaration statement will become too long, for instance due to a huge number, dynamically added cse-names, it may be better to use IMPLICIT typing.
- C neither supports IMPLICIT types nor has the types complex and complex\*16. The remaining types are denoted by int, float and double, respectively.
- Array and/or matrix definitions are also considered to be id's in id\_list's in declarations. However, we have to be aware of the instantaneous replacement of array- and/or matrix entries, when expressions are simplified. Therefore, we have to use operators, functioning as array and/or matrix names in code we want to optimize. We return to this question in the sections 7 and 8.

When the ON/OFF AGAIN strategy is applied we have to be aware of the above outlined declaration strategy. The last OPTIMIZE command, executed

<sup>&</sup>lt;sup>1</sup>The **pascal** module of GENTRAN is not error free. Especially the template file features do not function correctly.

directly after choosing  ${\tt OFF}$  AGAIN, has to be extended with the  ${\tt DECLARE}$  option.

Array and/or matrix names only occur in literally parsed information. In all other situations we have to make use of REDUCE operators. Normally, function applications inside SCOPE input are instantaneously replaced by newly selected cse names after putting them in the function table. Usually array and/or matrix entries are considered to be function applications. However, when due to a DECLARE option array and/or matrix names are known via the contents of the symbol table, such entries are substituted back before SCOPE produces output.

### Example 18

A simple OPTIMIZE command, extended with a DECLARE option, is executed for the various output options of GENTRAN, including the f90 alternative.

```
OPTLANG fortran$
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real; b(5):integer>>$
      INTEGER B(5), I, S10, S9
      REAL A(4,4), X(4), Y(5)
      S10=I+1
      S9=I-1
      X(S10) = A(S10, S9) + B(I)
      Y(S9) = A(S9, S10) - B(i)
OPTLANG ratfor$
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real; b(5):integer>>$
integer b(5),i,s10,s9
real a(4,4),x(4),y(5)
{
    s10=i+1
    s9=i-1
    x(s10)=a(s10,s9)+b(i)
    y(s9)=a(s9,s10)-b(i)
}
```

OPTLANG c\$

```
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real; b(5):integer>>$
int b[6],i,s10,s9;
float a[5][5],x[5],y[6];
{
    s10=i+1;
   s9=i-1;
   x[s10]=a[s10][s9]+b[i];
    y[s9]=a[s9][s10]-b[i];
}
OPTLANG pascal$
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real; b(5):integer>>$
var
    s9,s10,i: integer;
   b: array[0..5] of integer;
    y: array[0..5] of real;
    x: array[0..4] of real;
    a: array[0..4,0..4] of real;
begin
    s10:=i+1;
    s9:=i-1;
    x[s10]:=a[s10,s9]+b[i];
    y[s9]:=a[s9,s10]-b[i]
end;
OPTLANG nil$
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real; b(5):integer>>$
integer b(5),i,s10,s9
real a(4,4), x(4), y(5)
s10 := i + 1
s9 := i - 1
x(s10) := a(s10, s9) + b(i)
y(s9) := a(s9, s10) - b(i)
```

```
OPTLANG fortran$
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real*8; b(5):integer>>$
      INTEGER B(5), I, S10, S9
      DOUBLE PRECISION A(4,4), X(4), Y(5)
      S10=I+1
      S9=T-1
      X(S10) = A(S10, S9) + B(I)
      Y(S9) = A(S9, S10) - B(I)
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << x(4),y(5):real; b(5):complex>>$
***** Type error:
real x(4), y(5)
complex b(5)
(integer all) s9
integer s5,i
real := complex(all)
***** Wrong typing
Cont? (Y or N)
We can restart REDUCEand rerun the example with the Fortran90 version
of SCOPE. It results in:
LOAD_PACKAGE scope90$
OPTLANG f90$
OPTIMIZE x(i+1):=a(i+1,i-1)+b(i),y(i-1):=a(i-1,i+1)-b(i)
INAME s
DECLARE << a(4,4),x(4),y(5):real; b(5):integer>>$
      REAL, DIMENSION (4,4)::A
      INTEGER,DIMENSION(5)::B
      INTEGER:: I,S10,S9
      REAL, DIMENSION(4)::x
      REAL, DIMENSION(5)::y
      S10=I+1
      S9=I-1
      X(S10) = A(S10, S9) + B(I)
      Y(S9) = A(S9, S10) - B(I)
```

## 6.1 Coefficient Arithmetic and Precision Handling

REDUCE knows a variety of coefficient domains, as presented in subsection 9.11 of the REDUCE 3.6 manual [17], entitled *Polynomial Coefficient* Arithmetic. As stated in subsection 3.1 SCOPE supports integer and real coefficients. By turning ON the switch ROUNDED we introduce float arithmetic for coefficients. The operator PRECISION can be applied to change the default, machine dependent *precision*. Internally, REDUCE uses floating point numbers up to the precision supported by the underlying machine hardware, and so-called *bigfloats* for higher precision. The internal precision is two decimals greater than the exernal precision to guard against roundoff inaccuracies. Rounded numbers are normally printed to the specified precision. If the user wishes to print such numbers with less precision, the printing precision can be set by the command PRINT\_PRECISION. If a case arises where use of the machine arithmetic leads to problems, a user can force REDUCE to use the bigfloat representation by turning ON the switch ROUNDBF, which is normally OFF. GENTRAN, and thus SCOPE as well, support bigfloat notation. However the precision is a responsibility of the user. A possibility is to use the PRINT\_PRECISION command, both for algebraic mode output and for output in a selected target language, like fortran77. SCOPE uses the given precision for selecting cse's. Although complex arithmetic is not supported in SCOPE, a simple alternative is provided. When using float arithmetic in REDUCE the protected name I can be used to denote  $\sqrt{-1}$ . If the I is included in a declaration list as an identifier of type complex(!\*), its assumed value is automatically put ahead of the resulting optimized code. We illustrate the different possibilities in example 19. Comment is included.

### Example 19

```
OPTLANG fortran$
ON ROUNDED, DOUBLE$
% ----
% We start with precision 6. The returned value is the internal
% precision supported by the underlying machine hardware.
% ----
PRECISION 6;
```

```
OPTIMIZE x1:= 2
                    *a + 10
                                     * b.
        x2:= 2.00001 *a + 10
                                      * b,
        x3:= 2  *a + 10.00001  * b,
        x4:= 6
                  *a + 30
                                      * b,
        x5:= 2.0000001*a + 10.000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, X1, X2, X3, X4, X5
     S1=10*B
     X1=S1+2*A
     X2=S1+2.00001D0*A
     X3=X1
     X4=3*X1
     X5=X1
% ---
\% Explanation: X1 is a cse of X3, X4 and X5, but not of X2, because
\% the coefficient 2.00001 is given in 6 decimal digits.
% Increase in precision will show this.
% ---
PRECISION 7$
OPTIMIZE x1:= 2
                    *a + 10
                                     * b,
        x2:= 2.00001 *a + 10
                                      * b,
        x3:= 2  *a + 10.00001  * b,
        x4:= 6 *a + 30
                                      * b,
        x5:= 2.0000001*a + 10.000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.00001D0*A
     X3=S1+1.000001D1*B
     X4=3*X1
     X5=X1
PRECISION 8$
```

```
OPTIMIZE x1:= 2 *a + 10
                                    * b,
       x2:= 2.00001 *a + 10
                                     * b,
        x3:= 2 *a + 10.00001 * b,
        x4:= 6
                     *a + 30
                                     * b.
        x5:= 2.0000001*a + 10.000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.00001D0*A
     X3=S1+1.000001D1*B
     X4=3*X1
     X5=2.000001D0*A+1.0000001D1*B
% ---
% All rhs's were taken literally. Let us now increase precision and
\% simplify the rhs's before optimization. It is in fact a repetition
\% of the examples above, this time with a larger precision.
% ---
PRECISION 20$
OPTIMIZE x1:=:2
                                   *a + 10
                                                                * b,
                                                                * b,
        x2:=:2.00000000000000000 *a + 10
                                   x3:=:2
                                                                * b,
                                   *a + 30
        x4:=:6
                                                                * b,
        x5:=:2.000000000000000001*a + 10.0000000000000000000 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.00000000000000001D0*A
     X3=S1+1.0D1*B
     X4=3*X1
     X5=1.0D0*X1
PRECISION 21$
```

```
OPTIMIZE x1:=:2
                                   *a + 10
                                                                * b,
        x2:=:2.00000000000000000 *a + 10
                                                                * b,
                                   x3:=:2
                                                                * b,
        x4:=:6
                                   *a + 30
                                                                * b.
        x5:=:2.000000000000000001*a + 10.00000000000000000000 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.00000000000000001D0*A
     X3=S1+1.000000000000000001D1*B
     X4=3*X1
     X5=2.0D0*A+1.0D1*B
PRECISION 22$
                                                                * b,
OPTIMIZE x1:=:2
                                   *a + 10
        x2:=:2.00000000000000000 *a + 10
                                                                * b,
                                   x3:=:2
                                                                * b,
        x4:=:6
                                   *a + 30
                                                                * b.
        x5:=:2.000000000000000001*a + 10.00000000000000000000 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.00000000000000001D0*A
     X3=S1+1.000000000000000001D1*B
     X4=3*X1
     X5=2.0000000000000000000000001D0*A+1.0D1*B
% ---
% However, we can observe some differences in both modes of operation, when
% selecting a precision around the precision supported by the undelying
% machine hardware. Then the switch ROUNDBF can better be turned ON.
% ---
OFF ROUNDBF$
PRECISION 12$
```

```
OPTIMIZE x1:= 2.00
                         *a + 10.00
                                         * b.
                                         * b,
        x2:= 2.0000000001*a + 10
        x3:= 2  *a + 10.00000001* b,
        x4:= 6
                         *a + 30
                                        * b,
        x5:= 2.00000000001*a + 10.00000000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.0000000001D0*A
     X3=S1+1.000000001D1*B
     X4=3*X1
     X5=X1
OPTIMIZE x1:=:2.00 *a + 10.00
                                          * b,
        x2:=:2.0000000001*a + 10
                                         * b,
        x3:=:2
                       *a + 10.00000001* b,
        x4:=:6
                         *a + 30
                                        * b,
        x5:=:2.00000000001*a + 10.00000000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.0D0*A
     X3=S1+10.0D0*B
     X4=3*X1
     X5=X1
% ----
% Observe that simplification prior to optimization leads to internal
\% roundings, which differ from the rounding used for literally taken
\% coefficients. This difference disappeares with ON ROUNDBF.
% ---
ON ROUNDBF$
```

```
OPTIMIZE x1:= 2.00
                         *a + 10.00
                                         * b,
        x2:= 2.0000000001*a + 10
                                         * b.
        x3:= 2
                       *a + 10.00000001* b,
        x4:= 6
                        *a + 30
                                       * b.
        x5:= 2.00000000001*a + 10.00000000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.0000000001D0*A
     X3=S1+1.000000001D1*B
     X4=3*X1
     X5=X1
OPTIMIZE x1:=:2.00
                        *a + 10.00
                                        * b,
        x2:=:2.0000000001*a + 10
                                         * b,
        x3:=:2
                       *a + 10.00000001* b,
        x4:=:6
                        *a + 30
                                        * b,
        x5:=:2.00000000001*a + 10.00000000001 * b
INAME s
DECLARE <<x1,x2,x3,x4,x5,a,b: real>>$
     DOUBLE PRECISION A, B, S1, S2, X1, X2, X3, X4, X5
     S1=2*A
     S2=10*B
     X1=S2+S1
     X2=S2+2.0000000001D0*A
     X3=S1+1.000000001D1*B
     X4=3*X1
     X5=X1
% ----
% Complex arithmetic is not supported in SCOPE. However the Fortan equivalent
% of I, a protected name in REDUCE, is automatically created, ahead of the
% optimized code, whenever I is included in the declaration as a type complex
% or a type complex*16 identifier.
% ---
OPTIMIZE a:=b+c
INAME s
DECLARE <<a,b,i,c: complex>>;
```

## 6 GENERATION OF DECLARATIONS

COMPLEX\*16 B,I,C,A I=(0.0D0, 1.0D0) A=B+C

OFF DOUBLE\$

OPTIMIZE a:=b+c INAME s DECLARE <<a,b,i,c: complex>>;

> COMPLEX B,I,C,A I=(0.0, 1.0) A=B+C

## 7 Dealing with Data Dependencies

SCOPE is designed to optimize blocks B of straight line code, i.e. sequences of n assignment statements  $S_i$  of the form  $\lambda_i := \rho_i$ , where  $i = 1, \dots, n$ . If an identifier occurs in  $\lambda_i$ , it is said to be *defined* in  $S_i$ . All identifiers occuring in  $\rho_i$  are said to be used in  $S_i$ . The set  $\text{DEF}(S_i)$  is formed by the identifiers defining  $S_i$ , usually only one. The set  $USE(S_i)$  is formed by the identifiers, which are used in  $S_i$ . The relation  $\text{DEF}(S_i) \in \text{USE}(S_i)$ , for  $1 \leq i < j \leq n$ , is called a *flow dependency* and denoted by  $S_i \to S_j$ . The relation  $\text{DEF}(S_i) \in \text{USE}(S_j)$ , for  $1 \leq j \leq i \leq n$  is called an *anti* dependency and denoted by  $S_i \not\rightarrow S_j$ . The set of inputs of B, denoted by I(B), consists of identifiers, which are used in B, before being defined, if defined at all. The set of outputs of B, denoted by O(B), consists of the set of all last definitions of identifiers, occuring in B. So a block of straight line code can be introduced as a triple  $B = \{S, I, O\}$ , where S stands for the sequence  $S_1; S_2; \dots; S_{n-1}; S_n$ , and where I and O define the inputs and outputs, respectively. When optimizing source code defined by B, i.e. the sequence S, the intention is to mechanically produce an equivalent, but computationally less complex sequence, preserving the relation between inputs and outputs. Due to anti dependencies, i.e. redefinitions of the rules for computing identifier values or stepwise computing such values, |O(B)| <n is possible. But that in turn implies that some of the used identifiers, although being literally identical, represent different values. Therefore, a mechanical search for cse's can only be maintained if these critical identifiers are adequately renamed internally before the optimization process itself is started. As long as the relation between I(B) and O(B) is preserved it is even allowed to partly maintain these additional names, when presenting the results of an optimization operation. Furthermore it is worth noting that *dead code* can be left out, when ever occuring. Such code can be introduced through redundant assignments. The SCOPE features for dealing with data dependencies are illustrated, using the following artificial piece of code:

$$\begin{array}{rclcrcrcrcrcrc} S_1 & :& a_{1,x+1} & := & g+h.r^f\\ S_2 & :& b_{y+1} & := & a_{1,2.x+1}.(g+h.r^f)\\ S_3 & :& c1 & := & h.r.a_{2,1+x}/g\\ S_4 & :& c2 & := & c1.a_{1,x+1}+sin(d)\\ S_5 & :& a_{1,x+1} & := & c1+2\\ S_6 & :& d & := & b_{y+1}.a_{1,x+1}\\ S_7 & :& a_{1,1+2.x} & := & a_{1,x+1}.b_{y+1}.c/(d.g^2)\\ S_8 & :& b_{y+1} & := & a_{1.x+1}+b_{y+1}+sin(d)\\ S_9 & :& a_{1,x+1} & := & b_{y+1}.c+h/(g+sin(d))\\ S_{10} & :& d & := & k.e+d.(a_{1,1+x}+3)\\ S_{11} & :& e & := & d.(a_{1,1+x}+3)+sin(d)\\ S_{12} & :& f & := & d.(a_{1,1+x}+3)+sin(d)\\ S_{13} & :& g & := & d.(3+a_{1,1+x})+f \end{array}$$

The different flow and anti dependencies can be vizualized in the following way:

The identifiers, occuring in the piece of code given above, can be defined, can be used or can play both roles. Identifiers, used in one or more of the  $\lambda_i, i = 1 \cdots n$  figure in subscript expressions. The set notation  $\{\cdots\}$  is used to explicitly describe USE sets. Since all DEF sets consist of one element only, we omitted the set notation for the DEF sets. This provides a simple tool to distinguish flow and anti dependencies:

 $: \lambda_1 \to \{\rho_3, \rho_4\} \not\to \lambda_5 \to \{\rho_6, \rho_7, \rho_8\} \not\to \lambda_9 \to \{\rho_{10}, \rho_{11}, \rho_{12}, \rho_{13}\}$  $a_{1,x+1}$ :  $\{\rho_1, \rho_2, \rho_3, \rho_7, \rho_9\} \not\rightarrow \lambda_{13}$ g:  $\{\rho_1, \rho_2, \rho_3, \rho_9\}$ hr:  $\{\rho_1, \rho_2, \rho_3\}$ f :  $\{\rho_1, \rho_2\} \not\rightarrow \lambda_{12} \rightarrow \rho_{13}$ :  $\lambda_2 \rightarrow \{\rho_6, \rho_7, \rho_8\} \not\rightarrow \lambda_8 \rightarrow \{\rho_9\}$  $b_{y+1}$ :  $\{\rho_2\} \not\rightarrow \lambda_7$  $a_{1,2,x+1}$ :  $\lambda_3 \rightarrow \{\rho_4, \rho_5\}$ c1:  $\{\rho_3\}$  $a_{2,1+x}$ c2:  $\lambda_4$ :  $\{\rho_4\} \not\rightarrow \lambda_6 \rightarrow \{\rho_7, \rho_8, \rho_9, \rho_{10}\} \not\rightarrow \lambda_{10} \rightarrow \{\rho_{11}, \rho_{12}, \rho_{13}\}$ dc $: \{\rho_7, \rho_9\}$ :  $\{\rho_{10}\}$ k:  $\{\rho_{10}\} \not\rightarrow \lambda_{12}$ e

We observe that

$$I(B) = \{x, y, g, h, r, f, a_{1,2,x+1}, a_{2,1+x}, d, c, k, e\},\$$
$$O(B) = \{a_{1,x+1}, g, f, b_{y+1}, a_{1,2,x+1}, c1, c2, d, e\}$$

and thus, that

$$I(B)\cap O(B)\neq \emptyset.$$

The identifiers  $a_{1,1+x}$  and d are redefined twice and the identifier  $b_{y+1}$  once. Only the input occurrences and the last output definitions need to be preserved.

We also observe that some of the identifiers are subscripted. In our example the subscript expressions are constructed with input identifiers only. More general situations are conceivable. The set of subscript expressions contains cse's. Since our optimization strategy is based on an all level approach expressions, like 1 + x and y + 1 are certainly discovered as cse's.

An OPTIMIZE command can be extended with a DECLARE option, indicating that both a and b are array names. Their subscript expressions are optimized. In addition, the a and b entries are considered to be array entries, and not as function applications. The latter will happen when the DECLARE option is omitted. Vector architectures make often use of machine specific optimizing compilers. Therefore it may be better not to optimize the subset of subscript expressions. We implemented some facilities to take such machine specific limitations into account. When turning ON the switch VECTORC

the finishing touch is omitted and subscript expressions are not optimized. The function

VECTORCODE <a\_or\_m\_id\_seq>

can be used to operate more selectively. The a\_or\_m\_id\_seq consists of one or more array and/or matrix names, separated by comma's. Only the actual parameters of VECTORCODE are assumed to be names of arrays. So only their subscript expressions are disregarded during an optimization process. We can undo the effect of the VECTORCODE setting with a command of the form:

VCLEAR <a\_or\_m\_id\_(sub)seq>

The actual parameters are supposed to be taken from the sequence of actual parameters of its counterpart, the function VECTORCODE.

The different settings are now illustrated in:

#### Example 20

The above given block of code  $B = \{S, I, O\}$  is optimized in five different situations. To start with we hand it over to SCOPE, using an OPTIMIZE command, without using the DECLARE option. We observe, looking at the results, that some renaming of non significant identifier definitions have been performed. The first occurrence of  $a_{1,1+x}$  is replaced by s34, the second by s4. The first occurrence of  $b_{y+1}$  is replaced by s3, but the occurrences of d, a scalar identifier, are maintained.

Especially the role of the scalar d is quite interesting. The first definition of d is given in  $S_6$  In  $S_7$  d is used twice: explicitly in the denominator and in a hidden way in the numerator as well. The optimized version of  $S_7$  shows a factor d/d. The reason is that d locally replaces an internally created cse name, substituted for  $\rho_6$  and for part of the numerator of  $\rho_7$ . Like illustrated in example 13 a second SCOPE application can further simplify  $\rho_7$ . The scalar d is also used as argument of the *sin*-function in  $S_4, S_8, S_9, S_{11}$  and  $S_{12}$ . The d-values in  $S_4$ , in  $\{S_8, S_9\}$  and in  $\{S_{11}, S_{12}\}$  are all different, due to the redefinitions in  $S_6$  and in  $S_{10}$ . Therefore we recognize two different cse's, containing sin(d): s24 and e. The role of  $a_{1,x+1}$  is of course very similar, albeit less transparant, due to the renamings.

The second run showes that adding a DECLARE option does not influence the form of the output in this particular case. Besides the production of declarations, the result of both runs is identical. All relevant input and output names are preserved in both runs.

```
OPTIMIZE a(1,x+1) := g+h*r^f,
        b(y+1) := a(1,2*x+1)*(g+h*r^f),
               := (h*r)/g*a(2,1+x),
:= c1*a(1,x+1) + sin(d),
        c1
        c2
        a(1,x+1) := c1^{(5/2)},
               := b(y+1)*a(1,x+1),
        d
        a(1,1+2*x):= (a(1,x+1)*b(y+1)*c)/(d*g^2),
        b(y+1) := a(1,1+x)+b(y+1) + sin(d),
        a(1,x+1) := b(y+1)*c+h/(g + sin(d)),
                := k*e+d*(a(1,1+x)+3),
        d
                 := d*(a(1,1+x)+3) + sin(d),
        е
        f
                 := d*(3+a(1,x+1)) + sin(d),
                 := d*(3+a(1,x+1))+f
        g
INAME s$
s0 := x + 1
f
s34 := r *h + g
s2 := 1 + y
s6 := 2*x + 1
s3 := s34*a(1,s6)
              r*h
c1 := a(2,s0)*-----
              g
c2 := sin(d) + s34*c1
s4 := sqrt(c1)*c1*c1
d := s4*s3
            d*c
a(1,s6) := -----
           g*g*d
s24 := sin(d)
b(s2) := s4 + s3 + s24
            h
a(1,s0) := ----- + b(s2)*c
          g + s24
s29 := 3 + a(1,s0)
d := s29*d + e*k
s33 := s29*d
e := s33 + sin(d)
f := e
g := s33 + f
```

```
OPTIMIZE a(1,x+1) := g+h*r^f,
         b(y+1) := a(1,2*x+1)*(g+h*r^f),
         c1 := (h*r)/g*a(2,1+x),
c2 := c1*a(1+x+1) + sin
         c2
                  := c1*a(1,x+1) + sin(d),
         a(1,x+1) := c1^{(5/2)},
         d := b(y+1)*a(1,x+1),
         a(1,1+2*x):= (a(1,x+1)*b(y+1)*c)/(d*g^2),
         b(y+1) := a(1,1+x)+b(y+1) + sin(d),
         a(1,x+1) := b(y+1)*c+h/(g + sin(d)),
                   := k*e+d*(a(1,1+x)+3),
         d
         е
                  := d*(a(1,1+x)+3) + sin(d),
         f
                  := d*(3+a(1,x+1)) + sin(d),
                   := d*(3+a(1,x+1))+f
         g
INAME s
DECLARE <<a(5,5),b(7),c,c1,c2,d,f,g,h,r:real*8; x,y:integer>>$
      INTEGER X,Y,S0,S2,S6
      DOUBLE PRECISION C, H, R, S34, S3, C1, C2, S4, S24, B(7), A(5,5), S29, K, D, S33
     . ,E,F,g
      S0=X+1
      S34=R**F*H+G
      S2=1+Y
      S6=2*X+1
      S3=S34*A(1,S6)
      C1=A(2,S0)*((R*H)/G)
      C2=DSIN(D)+S34*C1
      S4=DSQRT(C1)*C1*C1
      D=S4*S3
      A(1,S6)=(D*C)/(G*G*D)
      S24=DSIN(D)
      B(S2) = S4 + S3 + S24
      A(1,S0)=H/(G+S24)+B(S2)*C
      S29=3+A(1,S0)
      D=S29*D+DEXP(1.0D0)*K
      S33=S29*D
      E=S33+DSIN(D)
      F=DEXP(1.0D0)
      G=S33+F
```

#### OFF FORT\$

Observe the differences in the f and F assignments. When generating fortran77 code all right hand side occurrences of e are automatically consid-

ered as appearances of the base of the natural logarithm and are translated accordingly.

The third run is influenced by turning ON the switch VECTORC. We observe that all array references are substituted back, without having replaced repeatedly occuring identical subscript expressions by internally selected cse names.

The fourth and the last run are governed by the functions VECTORCODE and VCLEAR, after having turned OFF the switch VECTORC.

ON VECTORC\$

```
OPTIMIZE a(1,x+1) := g+h*r^f,
        b(y+1) := a(1,2*x+1)*(g+h*r^f),
        c1
                := (h*r)/g*a(2,1+x),
                 := c1*a(1,x+1) + sin(d),
        c2
        a(1,x+1) := c1^{(5/2)},
                  := b(y+1)*a(1,x+1),
         d
         a(1,1+2*x):= (a(1,x+1)*b(y+1)*c)/(d*g^2),
        b(y+1) := a(1,1+x)+b(y+1) + sin(d),
         a(1,x+1) := b(y+1)*c+h/(g + sin(d)),
         d
                 := k*e+d*(a(1,1+x)+3),
                  := d*(a(1,1+x)+3) + sin(d),
         е
        f
                  := d*(3+a(1,x+1)) + sin(d),
                   := d*(3+a(1,x+1))+f
         g
INAME s$
              f
a(1,x + 1) := r *h + g
b(y + 1) := a(1,x + 1)*a(1,2*x + 1)
                 r*h
c1 := a(2,x + 1)*----
g
c2 := sin(d) + a(1,x + 1)*c1
                         2
a(1,x + 1) := sqrt(c1)*c1
d := a(1, x + 1) * b(y + 1)
```

```
d*c
a(1,1 + 2*x) := -----
                 2
                g *d
s20 := sin(d)
b(y + 1) := a(1,x + 1) + b(y + 1) + s20
                h
a(1,x + 1) := ----- + b(y + 1)*c
              g + s20
s25 := a(1, x + 1) + 3
d := s25*d + e*k
s28 := s25*d
e := s28 + sin(d)
f := e
g := s28 + f
OFF VECTORC$
VECTORCODE a,b$
OPTIMIZE a(1,x+1) := g+h*r^f,
        b(y+1) := a(1,2*x+1)*(g+h*r^f),
                 := (h*r)/g*a(2,1+x),
        c1
        c2
                  := c1*a(1,x+1) + sin(d),
        a(1,x+1) := c1^{(5/2)},
        d
                 := b(y+1)*a(1,x+1),
        a(1,1+2*x):= (a(1,x+1)*b(y+1)*c)/(d*g^2),
        b(y+1) := a(1,1+x)+b(y+1) + sin(d),
        a(1,x+1) := b(y+1)*c+h/(g + sin(d)),
                 := k*e+d*(a(1,1+x)+3),
        d
        е
                  := d*(a(1,1+x)+3) + sin(d),
        f
                 := d*(3+a(1,x+1)) + sin(d),
                 := d*(3+a(1,x+1))+f
        g
INAME s$
              f
a(1,x + 1) := r *h + g
b(y + 1) := a(1,x + 1)*a(1,2*x + 1)
                 r*h
c1 := a(2,x + 1)*----
                  g
c2 := sin(d) + a(1,x + 1)*c1
a(1,x + 1) := sqrt(c1)*c1*c1
d := a(1, x + 1) * b(y + 1)
                 d*c
```

```
a(1,1 + 2*x) := -----
              g*g*d
s22 := sin(d)
b(y + 1) := a(1,x + 1) + b(y + 1) + s22
            h
a(1,x + 1) := ----- + b(y + 1)*c
              g + s22
s27 := 3 + a(1, x + 1)
d := s27*d + e*k
s30 := s27*d
e := s30 + sin(d)
f := e
g := s30 + f
VCLEAR b$
OPTIMIZE a(1,x+1) := g+h*r^f,
        b(y+1) := a(1,2*x+1)*(g+h*r^f),
             := (h*r)/g*a(2,1+x),
:= c1*c(1
        c1
        c2
                := c1*a(1,x+1) + sin(d),
        a(1,x+1) := c1^{(5/2)},
                 := b(y+1)*a(1,x+1),
        d
        a(1,1+2*x):= (a(1,x+1)*b(y+1)*c)/(d*g^2),
        b(y+1) := a(1,1+x)+b(y+1) + sin(d),
        a(1,x+1) := b(y+1)*c+h/(g + sin(d)),
                := k*e+d*(a(1,1+x)+3),
        d
                 := d*(a(1,1+x)+3) + sin(d),
        е
                 := d*(3+a(1,x+1)) + sin(d),
        f
                  := d*(3+a(1,x+1))+f
        g
INAME s$
             f
a(1,x + 1) := r *h + g
s1 := y + 1
s2 := a(1,x + 1)*a(1,1 + 2*x)
                 r*h
c1 := a(2, 1 + x) * - - - - -
                 g
c2 := sin(d) + a(1,x + 1)*c1
a(1,x + 1) := sqrt(c1)*c1*c1
d := a(1, x + 1) * s2
                 d*c
a(1,1 + 2*x) := -----
```

# 8 A Combined Use of GENTRAN and SCOPE 1.5

As already stated in subsection 2.4 each GENTRAN command is evaluated separately, implying that the symbol table, required for the production of declarations, is fresh at the beginning of a GENTRAN command evaluation. Turning ON the switch GENTRANOPT leads to the optimization of the arithmetic code, defined in the GENTRAN command, obeying the new GENTRANOPT regime. In addition, we can observe that each separate GEN-TRAN command can produce its own declarations.

To increase flexibility we introduced facilities for making declarations, associated with a group of GENTRAN commands and for the optimization of the arithmetic in such a group as well. We implemented two function pairs of parameter-less functions:

(DELAYDECS, MAKEDECS)

and

(DELAYOPTS, MAKEOPTS).

Both pairs function as "brackets" around groups of statements. The OPTS pair can be used (repeatedly) inside a DECS pair. Both pairs achieve an alterned GENTRAN mode of operation. All GENTRAN productions between such a pair are collected in an internal format, say if\_list. The DELAY... functions initialize the modified mode of operation. The MAKE... functions restore the previous mode of operation in combination with the production either of declarations, associated with the contents of if\_list, or of an optimized representation of the contents of if\_list.

Example 21 serves to introduce a variety of approaches to code generation, based on the use of these pairs of brackets and of the switch GENTRANOPT. We illustrate a more realistic use in example 22: generation of optimized fortran77 code for the computation of the entries of the inverse of a symmetric (3,3) matrix. It is a continuation of example 8 in subsection 3.1 and example 11 in subsection 3.2. It was also presented in [9], albeit in a slightly different form.

#### Example 21

The output of combined GENTRAN and SCOPE operations is by default given in fortran77 notation. We illustrate the different possibilities in the form of small pieces of code. • The pair (DELAYDECS, MAKEDECS) encloses four GENTRAN commands. The first is needed to initialize the symbol table. The literal translation in internal format of the last three commands is stored in the if\_list. The application of MAKEDECS leads to the restoration of the default GENTRAN regime, applied to the if\_list and leading to the result, presented in the form of fortran77 code.

```
DELAYDECS$
```

```
GENTRAN DECLARE <<a,b,c,d,q,w:real>>$
GENTRAN a:=b+c$
GENTRAN d:=b+c$
GENTRAN <<q:=b+c;w:=b+c>>$
```

MAKEDECS\$

REAL A,B,C,D,Q,W A=B+C D=B+C Q=B+C W=B+C

• We repeat the previous commands, but execute them in a slightly different setting by turning ON the switch GENTRANOPT. This time the arithmetical rules in each of the three last GENTRAN commands are optimized separately. This is illustrated by the form of the output. The last piece of code contains the cse B+C, which is presented under the name Q in the fortran77 output.

ON GENTRANOPT\$

DELAYDECS\$

```
GENTRAN DECLARE <<a,b,c,d,q,w:real>>$
GENTRAN a:=b+c$
GENTRAN d:=b+c$
GENTRAN <<q:=b+c;w:=b+c>>$
```

MAKEDECS\$

```
REAL B,C,A,D,Q,W
A=B+C
D=B+C
Q=B+C
W=Q
```

OFF GENTRANOPT\$

• We can improve the optimization strategy by using the function pair (DELAYOPTS, MAKEOPTS) in stead of the pair (DELAYDECS, MAKEDECS). All blockwise combinable arithmetic is collected. These blocks of straight line code are optimized as separate input sets  $E_{\rm in}$ , when activating MAKEOPTS.

DELAYOPTS\$

```
GENTRAN a:=b+c$
GENTRAN d:=b+c$
GENTRAN <<q:=b+c;w:=b+c>>$
```

MAKEOPTS\$

- A=B+C D=A Q=A W=A
- We can further improve the optimization strategy by using the function pair (DELAYOPTS, MAKEOPTS) inside the pair (DELAYDECS, MAKEDECS). All blockwise combinable arithmetic is collected. These blocks of straight line code are optimized as separate input sets  $E_{\rm in}$ , when activating MAKEOPTS. But this time the results are added in internal format to the if\_list version, being maintained, as to obey the DELAYDECS regime.

DELAYDECS\$

GENTRAN DECLARE <<a,b,c,d,q,w:real>>\$

DELAYOPTS\$

GENTRAN a:=b+c\$
GENTRAN d:=b+c\$
GENTRAN <<q:=b+c;w:=b+c>>\$

MAKEOPTS\$

MAKEDECS\$

REAL B,C,A,D,Q,W A=B+C D=A Q=A

W=A

• A slightly more realistic example suggests how easily optimized code for sets of matrix entries can be obtained. We use two identical matrices **a** and **d**. The latter is not introduced at the REDUCE level, but simply used inside a GENTRAN command. The entries of a are initialized inside a double for-loop. After each initialization a GEN-TRAN command is applied, using the special assignment operator ::=: for correctly combining entry names and entry values. The RE-DUCE algebraic mode assignments are again used, when identifying the matrix **d** with the matrix **a**, applying the special assignment operator :=: in a separate GENTRAN command. The latter command is internally expanded into separate GENTRAN commands for each entry assignment. By using the pair (DELAYOPTS, MAKEOPTS) one block of straigt line code is constructed and optimized. It consists of two sets of assignments, one for the entries of the matrix **a** and another for the entries of the matrix d. The presented output shows that both matrices are indeed identical.

MATRIX a(4,4);

DELAYDECS\$

GENTRAN DECLARE <<a(4,4),d(4,4),b,c:real>>\$

DELAYOPTS\$

FOR i:=1:4 DO FOR j:=1:4 DO <<	a(i,j):=(i+j)*(b+c)+i*b+j*c;
	GENTRAN a(i,j)::=:a(i,j)>>\$
GENTRAN d:=:a\$	

MAKEOPTS\$

MAKEDECS\$

```
REAL B,C,G56,G54,G57,G55,A(4,4),D(4,4)
A(1,1)=3.0*(B+C)
G56=5.0*C
A(1,2)=G56+4.0*B
G54=5.0*B
G57=7.0*C
A(1,3)=G57+G54
A(1,4)=6.0*B+9.0*C
A(2,1)=G54+4.0*c
```

```
A(2,2)=2.0*A(1,1)
G55=7.0*B
A(2,3) = G55 + 8.0 * C
A(2,4)=2.0*A(1,2)
A(3,1)=G56+G55
A(3,2)=G57+8.0*B
A(3,3)=3.0*A(1,1)
A(3,4)=10.0*B+11.0*C
A(4,1)=9.0*B+6.0*C
A(4,2)=2.0*A(2,1)
A(4,3)=11.0*B+10.0*C
A(4,4)=4.0*A(1,1)
D(1,1)=A(1,1)
D(1,2)=A(1,2)
D(1,3)=A(1,3)
D(1,4)=A(1,4)
D(2,1)=A(2,1)
D(2,2)=A(2,2)
D(2,3)=A(2,3)
D(2,4)=A(2,4)
D(3,1)=A(3,1)
D(3,2)=A(3,2)
D(3,3)=A(3,3)
D(3,4)=A(3,4)
D(4,1)=A(4,1)
D(4,2)=A(4,2)
D(4,3)=A(4,3)
D(4,4) = A(4,4)
```

• Finally, and again only for illustrative purposes the fifth piece of code is again executed in an almost identical manner. We omit the declarations and replace the instruction GENTRAN d:=:a\$ by the command GENTRAN a:=:a\$. Hence the matrix a is simply redefined. As stated in section 7 redundant assignments — producing dead code, for instance by copying previous assignments — are automatically removed. as part of the optimization process. Therefore the optimized entry values of the matrix a are presented only once.

DELAYOPTS\$

```
A(1,1)=3.0*(B+C)
G111=5.0*C
A(1,2) = G111 + 4.0 * B
G109=5.0*B
G112=7.0*C
A(1,3) = G112 + G109
A(1,4)=6.0*B+9.0*C
A(2,1)=G109+4.0*C
A(2,2)=2.0*A(1,1)
G110=7.0*B
A(2,3)=G110+8.0*C
A(2,4)=2.0*A(1,2)
A(3,1) = G111 + G110
A(3,2)=G112+8.0*B
A(3,3)=3.0*A(1,1)
A(3,4)=10.0*B+11.0*C
A(4,1)=9.0*B+6.0*C
A(4,2)=2.0*A(2,1)
A(4,3)=11.0*B+10.0*C
A(4,4)=4.0*A(1,1)
```

Example 22

Let us now again look at the symmetric (3,3) matrix m, already used in the examples 8 and 11. For completeness we begin by showing the entry values. We generate optimized fortran77 code for the inverse mnv of m and store it in a file, named inverse.code, using the function pair (GENTRANOUT, GENTRANSHUT). Inside this pair we apply the pair (DELAYDECS, MAKEDECS). The latter pair encloses in turn the pair (DELAYOPTS, MAKEOPTS). Inside these innermost brackets four different sections of code can be distinguished. The first section consists of three LITERAL commands, for inserting comment in the target code. The second is formed by a double for-loop. Essential are the applications of the GENTRAN functions tempvar and markvar for assigning internal names to matrix entry values. These applications are very similar to the approach, chosen in example 11. The third section is again formed by LITERAL commands and the last orders the creation of the entries of the inverse matrix mnv. Before introducing the file inverse.code we selected SO as initial cse name, using the function INAME, and turned ON the switches ACINFO and DOUBLE.

Observe that directly after the MAKEOPTS instruction two sets of tables are printed with information about the arithmetic complexity of the two blocks of code, which are generated in the second and the last section. We activated ACINFO to show this effect. The tables are printed as a side effect. The output itself is directed to the file inverse.code. Looking at the contents of this file, given below, shows three different kinds of internally generated names. A number of S-names is created during the optimization of the first block of straight line code, created with the second section. In this piece of code we also notice T-names, generated by tempvar applications. The initial T character is the default internal GENTRAN selection for (temporarily needed) names. And finally G-names, introduced by applications of gensym, during the second optimization operation for reducing the arithmetic complexity of the entries of mnv. Because the code splitting is internally performed, the second SCOPE application is missing its INAME initialisation, thus leading to the application of gensym. Observe as well that INAME can be used as a separate facility as well.

```
OFF EXP$
```

MATRIX m(3,3),mnv(3,3)\$ IN "matrix.M"\$ m(1,1) := - ((j30y - j30z + 9\*m30\*p )\*sin(q3) 2 2 - 18\*cos(q2)\*cos(q3)\*m30\*p - j10y - j30y - m10\*p 2 - 18\*m30\*p ) 2 m(2,1) := m(1,2) := - ((j30y - j30z + 9\*m30\*p)\*sin(q3))2 - 9\*cos(q2)\*cos(q3)\*m30\*p - j30y - 9\*m30\*p ) 2 m(3,1) := m(1,3) := - 9\*sin(q2)\*sin(q3)\*m30\*p 2 2 2 m(2,2) := - ((j30y - j30z + 9\*m30\*p )\*sin(q3) - j30y - 9\*m30\*p ) m(3,2) := m(2,3) := 0

```
2
m(3,3) := j30x + 9*m30*p
INAME sO$
ON ACINFO, DOUBLE$
GENTRANOUT "inverse.code"$
DELAYDECS$
 GENTRAN DECLARE <<mnv(3,3),p,m30,j30y,j30z,q3,m10,q2,j10y,j30x:real>>$
 DELAYOPTS$
   GENTRAN LITERAL "C",tab!*," ",cr!*$
   GENTRAN LITERAL "C",tab!*," -- Computation of relevant M-entries --",cr!*$
   GENTRAN LITERAL "C",tab!*," ",cr!*$
   FOR j:=1:3 DO FOR k:=j:3 DO
    IF m(j,k) NEQ O THEN
     << s:=tempvar('real); markvar s;
        GENTRAN eval(s):=:m(j,k);
        m(j,k):=m(k,j):=s
     >>$
    GENTRAN LITERAL "C",tab!*," ",cr!*$
   GENTRAN LITERAL "C",tab!*," -- Computation of the inverse of M --",cr!*$
   GENTRAN LITERAL "C",tab!*," ",cr!*$
   GENTRAN mnv:=:m^(-1)$
 MAKEOPTS$
Number of operations in the input is:
Number of (+/-) operations
                            : 17
Number of unary - operations : 4
Number of * operations
                         : 30
Number of integer ^ operations : 14
Number of / operations : 0
Number of function applications : 9
```

Number of operations after optimization is:

```
Number of (+/-) operations : 11
Number of unary - operations : 1
Number of * operations
                             : 12
Number of integer \hat{} operations : 0
Number of / operations : 0
Number of function applications : 4
Number of operations in the input is:
Number of (+/-) operations
                             : 20
Number of unary - operations : 4
Number of * operations : 45
Number of integer ^ operations : 20
Number of / operations
                              : 9
Number of function applications : 0
Number of operations after optimization is:
Number of (+/-) operations
                            : 4
Number of unary - operations : 2
Number of * operations : 11
Number of integer ^ operations : 0
Number of / operations : 4
Number of function applications : 0
MAKEDECS$
GENTRANSHUT "inverse.code"$
The contents of the file inverse.code is:
     DOUBLE PRECISION P,M30,J30Y,J30Z,Q3,M10,Q2,J10Y,J30X,S0,S7,S5,S4,
     . S13, S11, T0, T3, T1, T2, T4, G153, G152, G151, G147, G155, G156, MNV(3, 3)
С
С
      -- Computation of relevant M-entries --
С
     SO=DSIN(Q3)
     S7=P*P
     S5=S7*M30
     S4=S5*DCOS(Q3)*DCOS(Q2)
     S13=9.0D0*S5
     S11=(S13+J30Y-J30Z)*S0*S0
     T0=J30Y+J10Y+18.0D0*(S4+S5)+S7*M10-S11
```

```
T3=S13+J30Y-S11
      T1=T3+9.0D0*S4
      T2=-(S13*DSIN(Q2)*S0)
      T4=S13+J30X
С
С
       -- Computation of the inverse of M --
С
      G153=T2*T2
      G152=T1*T1
      G151=T0*T4
      G147=G151*T3-(G153*T3)-(G152*T4)
      G155=T4/G147
      MNV(1,1)=G155*T3
      MNV(1,2) = -(G155 * T1)
      G156=T2/G147
      MNV(1,3) = -(G156*T3)
      MNV(2,1) = MNV(1,2)
      MNV(2,2)=(G151-G153)/G147
      MNV(2,3) = G156 * T1
      MNV(3,1) = MNV(1,3)
      MNV(3,2) = MNV(2,3)
      MNV(3,3)=(T0*T3-G152)/G147
```

Let us now repeat the generation process in a slightly different setting. We leave out the comment generating instructions, thus creating only one block of straight line code to be optimized. We choose for an S-name selection based on tempvar applications and for T-names for cse's. This time the default use of gensym is not necessary.

The contents' of both output files illustrate quotient optimization. All denominators, being the determinant of the matrix m, are identical. The set of rational entries of MNV contains the cse's G155 (T45) and G156 (T46).

```
TEMPVARNAME!*:='s$
TEMPVARNUM!*:=0$
INAME tO$
```

GENTRANOUT "inverse.code2"\$

DELAYDECS\$

GENTRAN DECLARE <<mnv(3,3),p,m30,j30y,j30z,q3,m10,q2,j10y,j30x:real>>\$
DELAYOPTS\$

```
FOR j:=1:3 D0 FOR k:=j:3 D0
IF m(j,k) NEQ 0 THEN
<< s:=tempvar('real); markvar(s);
    GENTRAN eval(s):=:m(j,k);
    m(j,k):=m(k,j):=s
>>$
```

```
GENTRAN mnv:=:m<sup>(-1)</sup>$
```

#### MAKEOPTS\$

Number of operations in the input is:

Number	of	(+/-) operations	:	37
Number	of	unary - operations	:	8
Number	of	* operations	:	75
Number	of	integer ^ operations	:	34
Number	of	/ operations	:	9
Number	of	function applications	:	9

Number of operations after optimization is:

```
Number of (+/-) operations: 15Number of unary - operations: 3Number of * operations: 23Number of integer ^ operations: 0Number of / operations: 4Number of function applications: 4
```

#### MAKEDECS\$

```
GENTRANSHUT "inverse.code2"$
```

The contents of the file inverse.code2 is:

```
DOUBLE PRECISION P,M30,J30Y,J30Z,Q3,M10,Q2,J10Y,J30X,T9,T40,T32,
. T31,T49,T47,S0,S3,S1,S2,S4,T39,T38,T36,T30,T45,T46,MNV(3,3)
T9=DSIN(Q3)
T40=P*P
T32=T40*M30
T31=T32*DCOS(Q3)*DCOS(Q2)
T49=9.0D0*T32
T47=(T49+J30Y-J30Z)*T9*T9
S0=J30Y+J10Y+18.0D0*(T31+T32)+T40*M10-T47
```

```
S3=T49+J30Y-T47
S1=S3+9.0D0*T31
S2=-(T49*DSIN(Q2)*T9)
S4=T49+J30X
T39=S2*S2
T38=S1*S1
T36=S0*S4
T30=T36*S3-(T39*S3)-(T38*S4)
T45=S4/T30
MNV(1,1) = T45 * S3
MNV(1,2) = -(T45 * S1)
T46=S2/T30
MNV(1,3) = -(T46 * S3)
MNV(2,1) = MNV(1,2)
MNV(2,2)=(T36-T39)/T30
MNV(2,3)=T46*S1
MNV(3,1) = MNV(1,3)
MNV(3,2)=MNV(2,3)
MNV(3,3)=(S0*S3-T38)/T30
```

A comparison between the arithmetic complexities given here and in example 11 shows that computing the entries of  $MNV(=M^{-1})$  instead of the value of the determinant of M, only requires 2 extra additions, 1 extra negation, 6 extra multilications and 4 extra divisions.

Other examples of this combined use of GENTRAN and SCOPE can be found in [31, 7]. The symbolic-numeric strategy discussed in [7] also relies on the ALGOPT facilities, which were introduced earlier.

# 9 Symbolic Mode Use of SCOPE 1.5

Both the OPTIMIZE command and the ALGOPT function are transformed into the same symbolic mode function, called SYMOPTIMIZE. It is this function, which governs the optimization process as a whole, delivering the results of an optimization run as a side effect, for instance by making it visible on a screen or by storing it in a file. Using SYMOPTIMIZE is straighforward, once the syntax for its five actual parameters is known. If we set ON INTERN a SYMOPTIMIZE application will deliver a list, containing the correctly ordered results of an optimization operation in the form of assignment statements in prefix form in Lisp notation. The thus provided results can function as one of the five actual parameters for SYMOPTIMIZE as well. This simple feature helps avoiding file traffic when stepwise optimizing code and as illustrated earlier in example 17 in section 5. Before illustrating that in example 23 we present the syntax of the actual parameters for SYMOPTIMIZE:

<SCOPE\_application> ::= ··· | SYMOPTIMIZE(<ssetq\_list>, <infile\_list>, <outfile\_name>, <cse\_prefix>

,

<declaration\_list>)

$<$ ssetq_list $>$	::=	$(< ssetq\_seq >)$
$<$ ssetq_seq $>$	::=	$<$ ssetq_stat $>$ [ $<$ ssetq_seq $>$ ]
$<$ ssetq_stat $>$	::=	(setq <lhs_id> <rhs>)   (rsetq <lhs_id> <rhs>)  </rhs></lhs_id></rhs></lhs_id>
		$(lsetq < lhs_id > < rhs >)   (lrsetq < lhs_id > < rhs >)$
<lhs_id></lhs_id>	::=	<id>   <subscripted_id></subscripted_id></id>
$<$ subscripted_id>	::=	( <id> <s_subscript_seq>)</s_subscript_seq></id>
$<$ s_subscript_seq>	::=	$[ ]$
$<$ s_subscript $>$	::=	<integer>   <integer prefix_expression=""></integer></integer>
< rhs >	::=	<prefix_expression></prefix_expression>
$<$ infile_list $>$	::=	$(< infile\_seq >)$
$<$ infile_seq $>$	::=	$<$ infile_name> [ $<$ infile_seq>]
$<$ infile_name $>$	::=	<string_id></string_id>
$< outfile\_name >$	::=	<string_id></string_id>
$<$ declaration_list $>$	::=	$(< declaration\_seq>)$
$<$ declaration_seq>	::=	$<$ declaration $>$ $<$ declaration_seq>
<declaration $>$	::=	( <type> <lhs_id_seq>)</lhs_id_seq></type>
$<$ lhs_id_seq>	::=	<lhs_id> <lhs_id_seq></lhs_id_seq></lhs_id>

The above given syntax requires some explanation:

• The presented ssetq syntax is incomplete. The prefix equivalent of any object, introduced in subsection 3.1 and of any a\_object, defined in subsection 3.2, is accepted as ssetq item. Such prefix equivalents can be obtained quite easily by using the function show:

# SYMBOLIC PROCEDURE show u; prettyprint u\$ SYMBOLIC OPERATOR show\$

The explicit presentation of a subset of the syntax rules for ssetq is given to suggest that local simplification in symbolic mode can be brought in easily by using the assignment operators lsetq, lrsetq and rsetq. The algebraic mode equivalent of these operators is ::=, ::=: and :=:, respectively. Their effect on simplification is discussed in subsection 2.4 and already shown in a number of examples. In addition it is worth noting that any (sub)expression in a lhs\_id or a rhs may contain any number of calls to eval. These calls lead to simplification of their arguments, prior to optimization. Details about the use of eval are presented in the GENTRAN manual [12].

• Since we operate in symbolic mode the last four formal parameters have possibly to be replaced by quoted actual parameters. This is illustrated in example 23.

- The infile\_seq consists of file names in string notation. The contents' of such input files may contain any form of infix input, obeying the syntax rules for objects and/or a\_objects, as introduced in the subsections 3.1 and 3.2, respectively.
- The single output file name outfile\_name ought to be given in string notation as well. The outfile\_name is properly closed. The default output is REDUCE infix in an ON NAT fashion. Alternatives are discussed above: ON AGAIN or OFF NAT, both leading to re-readable output, or an application of OPTLANG for a non nil argument.
- The declaration list presents declarations in prefix notation. The list is used to initialize the symbol table prior to optimization. This information is used for dynamically typing the result of an optimization process. In addition it is used to determine wether subscripted names denote array elements or a function call. The latter is replaced by a cse name in the presented output, whereas the former is not.
- The five parameters of SYMOPTIMIZE correspondent with optional extensions of the OPTIMIZE command. When part of these options remains unused, nil has to taken as value for the corresponding actual parameters.

We illustrate the symbolic mode variant of the OPTIMIZE command by repeating example 17 from section 5, albeit in a modified setting.

#### Example 23

The script explains itself.

LISP\$

ON ACINFO, INPUTC, INTERN, AGAIN\$

prettyprint(prefixlist:=SYMOPTIMIZE(nil,'("f1" "f2"),nil,'c,nil))\$

 $e1 := \frac{2 (x + y) 8 2 2}{3*x + 2*y}$   $e2 := (4*(\sin(x) - \cos(e)) + 3*\cos(x)) *(x + y) + 4*y + 2*y +$ 

```
2
             2
       + (4*x - 4*y) - 6*x)/(8*x + 3*y - 2*x)
Number of operations in the input is:
Number of (+/-) operations
                              : 16
Number of unary - operations
                               : 0
Number of * operations
                               : 16
Number of integer ^ operations : 11
Number of / operations
                        : 2
Number of function applications : 8
Number of operations after optimization is:
Number of (+/-) operations
                              : 16
Number of unary - operations : 0
Number of * operations
                              : 16
Number of integer ^ operations : 6
Number of / operations
                         : 2
Number of function applications : 4
((setq gsym c23)
   (setq cses (plus c18 c10 c20 c8 c6 c14))
   (setq c14 (plus x y))
   (setq c6 (expt c14 2))
   (setq c8 (cos x))
   (setq c20 (plus (expt (sin x) 2) (minus (cos (expt e c6)))) )
   (setq c10 (plus (times 3 x) (times 2 y)))
   (setq e1
      (quotient
         (plus
            (times 4 y)
            (times 4 (expt y 2))
            (times 2 c6 (expt (plus c20 (times 3 c8)) 8)))
        c10))
   (setq c18 (expt x 2))
   (setq e2
      (quotient
         (plus
            (times 4 c18)
            (minus (times 2 c10))
            (times 4 c6 c14 (expt (plus c20 (times 2 c8)) 2)))
         (plus (times 8 c18) (minus (times 2 x)) (times 3 y)))) )
```

```
OFF INTERN, AGAIN, PERIOD$
ON DOUBLE, FORT$
SYMOPTIMIZE(prefixlist,'("f3"), '"f7",'d,'((real e1 e2 e3 x y)))$
gsym := c23
cses := c18 + c10 + c20 + c8 + c6 + c14
c14 := x + y
       2
c6 := c14
c8 := cos(x)
          2
              c6
c20 := sin(x) - cos(e)
c10 := 3*x + 2*y
            2
                      8
     4*y + 4*y + 2*c6*(c20 + 3*c8)
e1 := -----
                c10
      2
c18 := x
                                   2
     4*c18 - 2*c10 + 4*c6*c14*(c20 + 2*c8)
e2 := -----
             8*c18 - 2*x + 3*y
                   2
             (x + y)
                                          2
                                  2
     4*\sin(\cos(e)) + \sin(x + y) + (4*x - x + 2*y)
e3 := -----
                       -----
                3*y + f(x,g( - cos(x)))
Number of operations in the input is:
Number of (+/-) operations : 23
Number of unary - operations : 1
```

```
: 20
Number of * operations
Number of integer ^ operations : 9
Number of / operations
                               : 3
Number of function applications : 11
Number of operations after optimization is:
Number of (+/-) operations
                               : 15
Number of unary - operations
                             : 1
Number of * operations
                               : 24
Number of integer ^ operations : 0
Number of / operations
                               : 3
Number of function applications : 8
The contents of the output file f7 is:
      DOUBLE PRECISION X,Y,D19,D16,C8,D1,D2,C20,D29,C10,D6,D38,D37,D31,
     . E1,C18,D9,D32,D27,D30,E2,D20,E3
     D19=X+Y
     D16=D19*D19
      C8=DCOS(X)
      D1=DSIN(X)
      D2=DCOS(DEXP(D16))
      C20=D1*D1-D2
      D29=2*Y
      C10=D29+3*X
      D6=C20+3*C8
     D38=D6*D6
     D37=D38*D38
     D31=4*Y
     E1=(D31+D31*Y+2.0D0*D16*D37*D37)/C10
      C18=X*X
      D9=C20+2*C8
      D32=4*C18
      D27=D32-X
      D30=3*Y
      E2=(D32-(2.0D0*C10)+4.0D0*D9*D9*D19*D16)/(D30+2.0D0*D27)
      D20=D29+D27
      E3=(4.0D0*DSIN(D2)+DSIN(D19)+D20*D20)/(D30+F(X,G(-C8)))
```

We especially designed these symbolic mode facilities for our joint research with Delft Hydraulics concerning code generation for an incompressible Navier-Stokes problem [15].

A final remark: The ON PREFIX mode of operation, in both algebraic and symbolic mode causes the results of a SCOPE application to be presented in the form of an association list, called **Prefixlist**. The pairs are formed by lhs\_id's and rhs values in prefixform. This lisp S-expression can be used to create an alternative version of the optimization results, in whatever target language the user prefers to choose.

#### Example 24

We show the ON PREFIX effect. When switching to symbolic mode (command 5) we can again obtain the output, assigned as value to the global identifer prefixlist. The ON PREFIX facility allows storage in a file for later use. When working in symbolic mode it is of course possible to apply ON INTERN in stead and to remove the setq extensions from the provided output value, if desired.

```
REDUCE 3.7, 15-Apr-1999 ...
1: LOAD_PACKAGE nscope$
2: OPTIMIZE a:=b+c*sin(x), d:=c*sin(x)*cos(y);
g7 := sin(x)*c
a := g7 + b
d := g7*cos(y)
3: ON PREFIX$
4: input 2;
Prefixlist:=
((g3 times (sin x) c) (a plus g3 b) (d times g3 (cos y)))
5: LISP$
6: prettyprint prefixlist$
((g3 times (sin x) c) (a plus g3 b) (d times g3 (cos y)))
7: caar prefixlist;
g3
7: cdar prefixlist;
```

(times (sin x) c)

9: BYE;

# 10 A Syntax Summary of SCOPE 1.5

REDUCE is extended with some commands, designed to apply the facilities offered by SCOPE in a flexible way. The syntactical rules, defining how to activate SCOPE in both algebraic and symbolic mode, are given in subsection 10.1. A short overview of the set of additional functions is given in subsection 10.2 and the relevant switches are again presented in subsection 10.3.

#### 10.1 SCOPE's Toplevel Commands

We assume the syntax of id's, integer's and the like to be already known. Hence we do not present an exhaustive description of the rules.

```
<REDUCE_command> ::= ··· | <SCOPE_application>
<GSTRUCTR_application> | <GHORNER_application> |
<SCOPE_application> ::= <OPTIMIZE command> |
<ALGOPT application> | <SYMOPTIMIZE application>
<OPTIMIZE command> ::=
    OPTIMIZE <object_seq> [IN <file_id_seq>] [OUT <file_id>]
        [INAME <cse_prefix>] [DECLARE <declaration_group>] |
    OPTIMIZE [<object_seq>] IN <file_id_seq> [OUT <file_id>]
        [INAME <cse_prefix>] [DECLARE <declaration_group>] |
    OPTIMIZE [<object_list>[,<string_id_list>][,<cse_prefix>]) |
    ALGOPT(<a_object_list>,[<string_id_list>][,<cse_prefix>]) |
    ALGOPT([<a_object_list>,]<string_id_list>[,<cse_prefix>])) |
    ALGOPT([<a_object_list>,]<string_id_list>[,<cse_prefix>])) |
    <SYMOPTIMIZE application> ::=
    SYMOPTIMIZE application> ::=
    SYMOPTIMIZE(<ssetq_list>, <infile_list>, <outfile_name>, <cse_prefix>,
        <declaration_list>)
```

$<$ GSTRUCTR_application $>$	::=	$\texttt{GSTRUCTR} < \texttt{stat\_group} > [\texttt{NAME} < \texttt{cse\_prefix} > ]$
$<$ stat_group $>$	::=	$\ll < \text{stat_list} \gg$
<stat_list></stat_list>	::=	$\langle \text{gstat} \rangle [; \langle \text{stat\_list} \rangle]$
$\langle gstat \rangle$	::=	$<$ name> := $<$ expression> $ $ $<$ matrix_id>
$<$ GHORHER_application $>$	::=	$\texttt{GHORNER} < \!\! \texttt{stat\_group} > [\texttt{VORDER} < \!\! \texttt{id\_seq} \!\! > ]$
<id_seq></id_seq>	::=	$<$ id>[, $<$ id_seq>]

<object_seq> <object> <stat> <assignment operator=""> <alglist> <eq_seq> <alglist_production> <name> <a_subscript_seq> <a_subscript> <cse_prefix></cse_prefix></a_subscript></a_subscript_seq></name></alglist_production></eq_seq></alglist></assignment></stat></object></object_seq>	::= ::= ::= ::= ::= ::= ::= ::=	<pre><stat>   <alglist>   <alglist_production> <name> <assignment operator=""> <expression> :=   ::=   ::=:   :=: {<eq_seq>} <name> = <expression>[,<eq_seq>] <name>   <function_application></function_application></name></eq_seq></expression></name></eq_seq></expression></assignment></name></alglist_production></alglist></stat></pre>
<a_object_list> <a_object_seq> <a_object></a_object></a_object_seq></a_object_list>	::= ::= ::=	<a_object>   {<a_object>[,<a_object_seq>]} <a_object>[,<a_object_seq>] <id>   <alglist>   <alglist_production>   {<a_object>}</a_object></alglist_production></alglist></id></a_object_seq></a_object></a_object_seq></a_object></a_object>
$<$ function_application>	::=	$\begin{array}{l} \texttt{ALGSTRUCTR} \ (<\!\!\operatorname{arg\_list}>[, <\!\!\operatorname{cse\_prefix}>]) \mid \\ \texttt{ALGHORNER} \ (<\!\!\operatorname{arg\_list}>[, \{<\!\!\operatorname{id\_seq}>\}]) \mid \\ & \ldots \end{array}$
<arg_list> <arg_seq> <arg> <arg_list_name></arg_list_name></arg></arg_seq></arg_list>	::= ::= ::=	
<file_id_seq> <file_id> <string_id_list> <string_id_seq> <string_id></string_id></string_id_seq></string_id_list></file_id></file_id_seq>	::= ::= ::= ::=	$<$ string_id>   { $<$ string_id_seq>} $<$ string_id> [, $<$ string_id_seq>]
<declaration_group> <declaration_list> <declaration> <range_list> <range> <id_list> <type></type></id_list></range></range_list></declaration></declaration_list></declaration_group>	::= ::= ::= ::= ::=	$\langle id \rangle$   $\langle id \rangle$ - $\langle id \rangle$
<ssetq_list> <ssetq_seq> <ssetq_stat></ssetq_stat></ssetq_seq></ssetq_list>		<pre>(<ssetq_seq>) <ssetq_stat> [<ssetq_seq>] (setq <lhs_id> <rhs>)   (rsetq <lhs_id> <rhs>)   (lsetq <lhs_id> <rhs>)   (lrsetq <lhs_id> <rhs>)</rhs></lhs_id></rhs></lhs_id></rhs></lhs_id></rhs></lhs_id></ssetq_seq></ssetq_stat></ssetq_seq></pre>

<lhs_id></lhs_id>	::=	$<$ id>   $<$ subscripted_id>
$<$ subscripted_id>	::=	$( )$
$<$ s_subscript_seq>	::=	$<$ s_subscript>[ $<$ s_subscript_seq>]
$<$ s_subscript $>$	::=	<integer>   <integer prefix_expression=""></integer></integer>
$\langle rhs \rangle$	::=	<pre>cprefix_expression&gt;</pre>
$<$ infile_list $>$	::=	$(< infile_seq >)$
$<$ infile_seq $>$	::=	$<$ infile_name> [ $<$ infile_seq>]
$<$ infile_name $>$	::=	<string_id></string_id>
$<$ outfile_name $>$	::=	<string_id></string_id>
$<$ declaration_list $>$	::=	$(< declaration\_seq>)$
$<$ declaration_seq>	::=	$<$ declaration> $<$ declaration_seq>
<declaration $>$	::=	$(< type > < lhs_id_seq >)$
$<$ lhs_id_seq $>$	::=	<lhs_id> <lhs_id_seq></lhs_id_seq></lhs_id>

# 10.2 Additional SCOPE-functions

Fifteen additional functions can be used. We shortly summarize their name and use:

Name(s)	Introduced in	See the examples:
SCOPE_SWITCHES	3	1
SETLENGTH, RESETLENGTH	3.1	2, 5, 12  and  13
ARESULTS, RESTORABLES, ARESTORE, RESTOREALL	3.1	9, $12 \text{ and } 15$
OPTLANG	6	18
VECTORCODE, VCLEAR	7	20
DELAYDECS, MAKEDECS, DELAYOPTS, MAKEOPTS	8	21  and  22
INAME	8	22

## 10.3 The relevant REDUCE, GENTRAN and SCOPE switches

We also shortly summarize the use of the switches, which were introduced in section 3 in example 1:

## 10 A SYNTAX SUMMARY OF SCOPE 1.5

Name	Origin	Illustrated in the examples:
ACINFO	SCOPE	8, 11, 22 and 23
AGAIN	SCOPE	17 and 23
DOUBLE	GENTRAN	23
EVALLHSEQP	REDUCE	11
EXP	REDUCE	8, 13, 16 and 22
FORT	REDUCE	4, 6, 8, 11, 20  and  23
FTCH	SCOPE	4 and 5

# 10 A SYNTAX SUMMARY OF SCOPE 1.5

Name	Origin	Illustrated in the examples:
GENTRANOPT	GENTRAN	21
INPUTC	SCOPE	3, 6, 7, 13, 17  and  23
INTERN	SCOPE	23
NAT	REDUCE	8 and 17
PERIOD	REDUCE	4
PREFIX	SCOPE	24
PRIALL	SCOPE	2
PRIMAT	SCOPE	13
POUNDBF	REDUCE	19
ROUNDED	REDUCE	7 and 22
SIDREL	SCOPE	13
VECTORC	SCOPE	20

## 11 SCOPE 1.5 Installation Guide

SCOPE 1.5 is easily installed. The usual code compilation facilities of RE-DUCE can be applied. In view of the frequent interaction between SCOPE and GENTRAN a compiled version of GENTRAN is required during the creation of the load module for SCOPE 1.5. The compilation process itself is vizualized below.

```
faslout "~infhvh/mkscope/scope_15";
lisp in "~infhvh/mkscope/cosmac.red"$
lisp in "~infhvh/mkscope/codctl.red"$
lisp in "~infhvh/mkscope/codopt.red"$
lisp in "~infhvh/mkscope/codad1.red"$
lisp in "~infhvh/mkscope/codad2.red"$
lisp in "~infhvh/mkscope/coddec.red"$
lisp in "~infhvh/mkscope/coddec.red"$
lisp in "~infhvh/mkscope/coddpri.red"$
lisp in "~infhvh/mkscope/codpri.red"$
lisp in "~infhvh/mkscope/codpri.red"$
lisp in "~infhvh/mkscope/codpri.red"$
lisp in "~infhvh/mkscope/codpri.red"$
lisp in "~infhvh/mkscope/codbr.red"$
```

faslend ;
end;

The subdirectory mkscope in the author's directory system contains the files with the source code of SCOPE 1.5. The order in which the files are read in is irrelevant except the first and the last. The file cosmac.red contains one module, named cosmac, which consists of a set of smacro procedures, designed to simplify access to the lower levels of the expression data base, employed during optimization. These smacro's are used in all other code sections. The last file in optional and usually executed to include new patches into a recompilable version of the package. Once it is stored in the fas1 directory of the local REDUCE system it is available as a load\_package.

In short, the files contain the following code sections:

• cosmac.red contains the module cosmac, consisting of smacro proce-

dures, allowing access to the expression data base.

- codctl.red consists of the three modules codctl, restore and minlenght. The first is a large module, containing the optimization process managing facilities. The second module is added to regulate the interplay with the REDUCE simplifier, when entering optimizer output in algebraic mode using functions like ARESULTS. The last module serves to vary the minimal length of cse's using SETLENGTH and RESETLENGTH.
- codmat.red contains the module codmat, defining the parsing process and the expression data base setup and access facilities.
- codopt.red's content is formed by the module codopt. It is the kernel of the optimization process, the implementation of the extended Breuer algorithm.
- codad1.red contains the module codad1, formed by additional facilities for improving the lay-out of the overall result, for information migration between different sections of the expression data base and for the application of the distributive law to remodel and compactify (sub)expression structure at any level.
- codad2.red contains the module codad2. This module defines five different possible activities during the optimization process. The first four regulate the so called *finishing touch*. The last one is a new section, defining how to optimize *rational forms* as part of the overall optimization process.
- coddec.red covers the declaration facilities, presented in section 6, collected in the module coddec and based on chapter 6 of [1]. The symbol table setup of GENTRAN is used.
- codpri.red is also formed by one module, called codpri. It covers all printing facilities. The first section is applied when the switch PRIMAT is turned ON. The latter is used to produce an internal list of pairs, consisting of the left hand side and the right hand side of assignment statements in prefix notation, and defining the output of the optimization process in sequential order. This prefix list is delivered to GENTRAN or REDUCE to make the results visible in the user prefered form. The intial version of this list, created directly after the optimization process, is improved using a collection of functions, also grouped together in the module codpri. These improvements may for instance be necessary to remove temporarily introduced names, internally employed as a result of a data dependency analysis.

- codgen.red consists of the module codgen. The interface between GENTRAN and SCOPE 1.5, introduced in section 8 of this manual is defined in this module.
- codhrn.red is formed by the module ghorner. It defines the facilities, presented in section 4.3 of this manual.
- codstr.red contains the module gstructr. This module defines the possibilities for expression structure recognition, as presented in section 4.2 of this manual.
- coddom.red finally, consists of one module, called coddom. It covers additional coefficient domain functions, needed to make the extended Breuere algorithm and the additional functions, collected in the modules codad1 and codad2 for instance, applicable for (multiple presicion) floating point coefficients.

A few additional remarks:

- GENTRAN plays an important role when creating declarations and output. The package is automatically loaded when executing one of the first instructions in the module codct1. Hence it may be necessary to look critically to the load instruction in codct1 before installing SCOPE 1.5. By changing this load instruction we easily created a fortran90 compatable SCOPE version [3]. At present it is only available for our own internal and experimental use.
- We believe the code to be almost version independent. Over the past years all uses of nil have been critically reviewed. However the coddom module may require version based maintenance when installing SCOPE 1.5.
- The present size of the source code is given in the table below. Comment is included in these figures

File naam	number of lines	number of characters
cosmac.red	172	4761
codctl.red	1439	61466
codmat.red	1488	72733
codopt.red	1243	68809
codad1.red	801	39175
codad2.red	1314	59217
coddec.red	928	41069
codpri.red	1371	62600
codgen.red	214	9120
codhrn.red	752	30549
codstr.red	308	11199
coddom.red	204	6638

• The modules **ghorner** and **gstructr** can be left out without harming the other facilities, presented in this manual.

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