D.C.P.L. - A Distributed Control Programming Language

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D. C. P. L.

A DISTRIBUTED CONTROL PROGRAMMING LANGUAGE

by

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ABSTRACT*

In this thesis, a computation is considered a system of asynchronously cooperating "independent" programs (coroutines) linked by paths of information along which messages are sent.

A programming language called DCPL, a Distributed Control Programming Language, in which such computations may be expressed, and which may be considered as a system-oriented programming language, is presented. A tree-structured representation and a very dynamic binding give to a DCPL program the flexibility of the highest level programming languages together with the potential of concurrency of the asynchronous computational structures.

The locality of references which is exhibited in any DCPL program allows a new computer organization using sequential storage devices with large transfer rate instead of random-access storage devices with relatively low transfer rate. Moreover, the computer is expected to achieve a large throughput by taking the parallelism into account.

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CHAPTER I

INTRODUCTION

I.1 Introduction.

In this thesis, a computation is considered a system of asynchronously cooperating "independent" programs (coroutines) linked by paths of information along which messages are sent.

A programming language called DCPL, a Distributed Control Programming Language, in which such computations may be expressed, and which may be considered as a system-oriented programming language, is presented. A tree structured representation and a very dynamic binding give to a DCPL program the flexibility of the highest level programming languages together with the potential of concurrency of the asynchronous computational structures.

The locality of references which is exhibited in any DCPL program allows a new computer organization using (inexpensive) sequential storage devices with large transfer rate instead of (expensive) random access storage devices with relatively low transfer rate. Moreover, the computer is expected to achieve large throughput by taking the parallelism into account.

I.2 Variables in computing processes.

In his paper "On certain basic concepts of programming languages" [27] Niklaus Wirth wrote:

"The elementary concepts of computing processes are:

- o There exist certain quantities, to be called "values" and elementary classes or types (possibly only one) of values among whose elements given elementary relationships hold. These relationships or mappings are represented in a computer by its operations which generate a new value (called result) which has the specified relationship to the given value(s) (called operands).
- o There exist cells (usually called "variables") which are able to contain a value, and which have a name. That name serves to refer to the contained value.
- o There exists an operator for the assignment of a new value to a cell.

... "

These concepts are widely accepted today, and they underlie any actual implementation of a conventional programming language.

In DCPL we support the first part of the quotation: there is a universe of values structured in classes or types, and mappings from some classes to some possibly different classes which are actualized by operations. There may be, for instance, in our universe, integer and logical values forming the classes I and L, and the operations:

operations	operators	mappings
addition	$+$	$I \times I \to I$
disjunction	$\vee$	$L \times L \to L$
negation	$\sim$	$L \to L$
equality	$=$	$I \times I \to L$

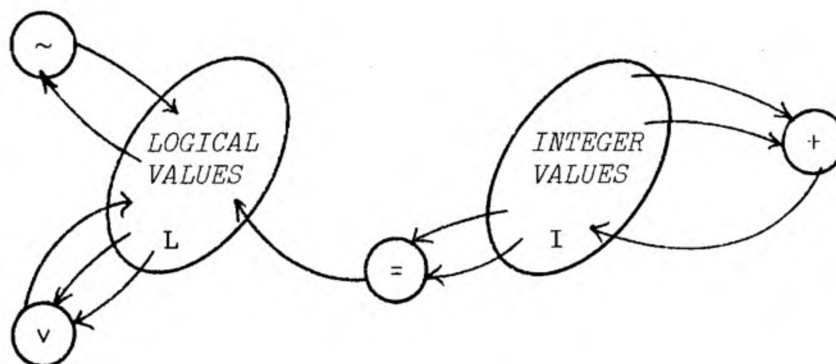

Fig. I-1

In fact, it is not our intention to impose any restriction on our universe. Consequently, we are leaving the list of types and operations open-ended. For this reason, and since syntax has received a rather speedy treatment, the emphasis being placed upon semantics, it would be proper to consider DCPL as describing a family of languages rather than defining completely one specific language.

DCPL, however, does not support the second part of the quotation: one variable in DCPL has no meaning by itself; a system of mutually bound variables defines communication paths.

I.2.1 An example.

Let us consider the evaluation of the following simple expression:

(1) $(2 + 3 \rightarrow x;\ (6-1)\rightarrow y;\ x + y \rightarrow x\ ;\ y + 1 \rightarrow y\ ;(2 \times y) - x\ )$

in a conventional programming language (fig. I-2):

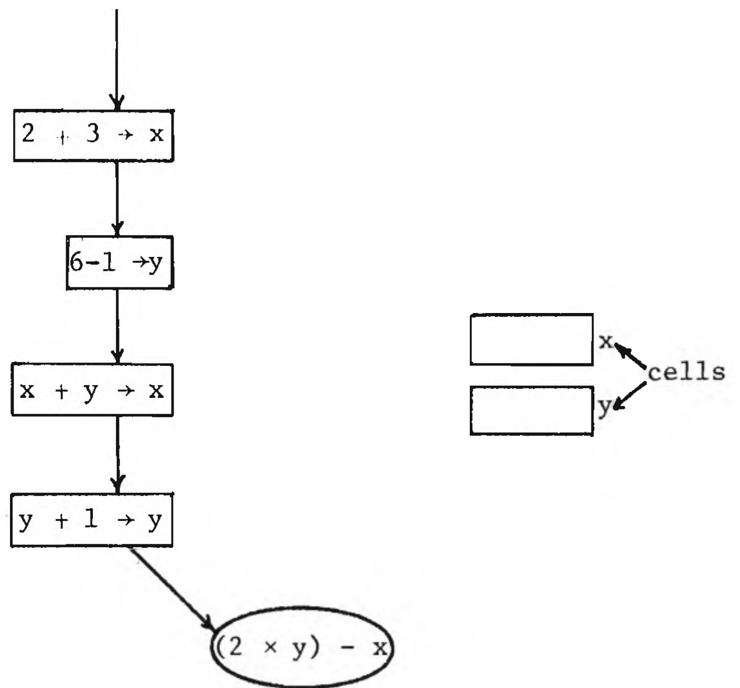

Fig. I-2

- there are two cells x and y ;

- the expression is viewed as a sequence of statements. The execution of any statement is a simple process involving some cells: whenever the statement $x + y \to x$, for instance, is executed, the values contained in x and y are retrieved, and their sum is then stored in x;

- the statements are executed serially one after the other. The last item is not a statement; its value is considered to be the value of the expression.

The same expression may be interpreted in a quite different way in DCPL. In order to study gradually the notions involved, we start by interpreting a much simpler expression without variables:

(2) $(2 + 3 + 2) + 2 + 1$

Association having to be done on the right, the expression may be represented as the tree of fig. I-5a. We may view each node as a simple automaton, and each edge as a channel of information. Whenever a node represents an integer, it sends up spontaneously along the edge its own value, and then vanishes (fig. I-3).

Whenever a node represents an addition, it waits until it receives a value from both the right and the left side; then it adds the two values and sends the result up along the channel, and vanishes (fig. I-4). The evaluation of expression (2) is displayed in fig. I-5.

Let us consider again expression (1). The expression may be represented as a syntax tree (fig. I-6).

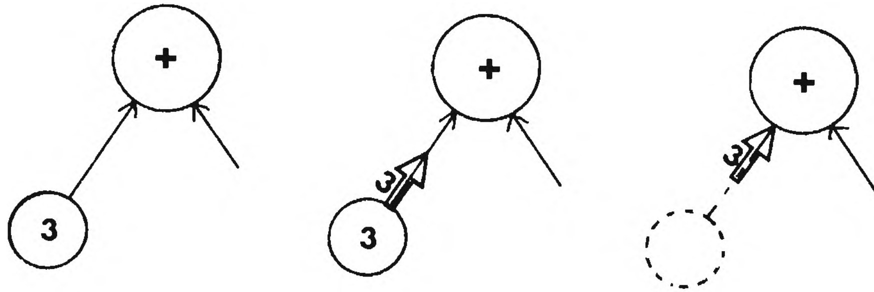

Fig. I-3

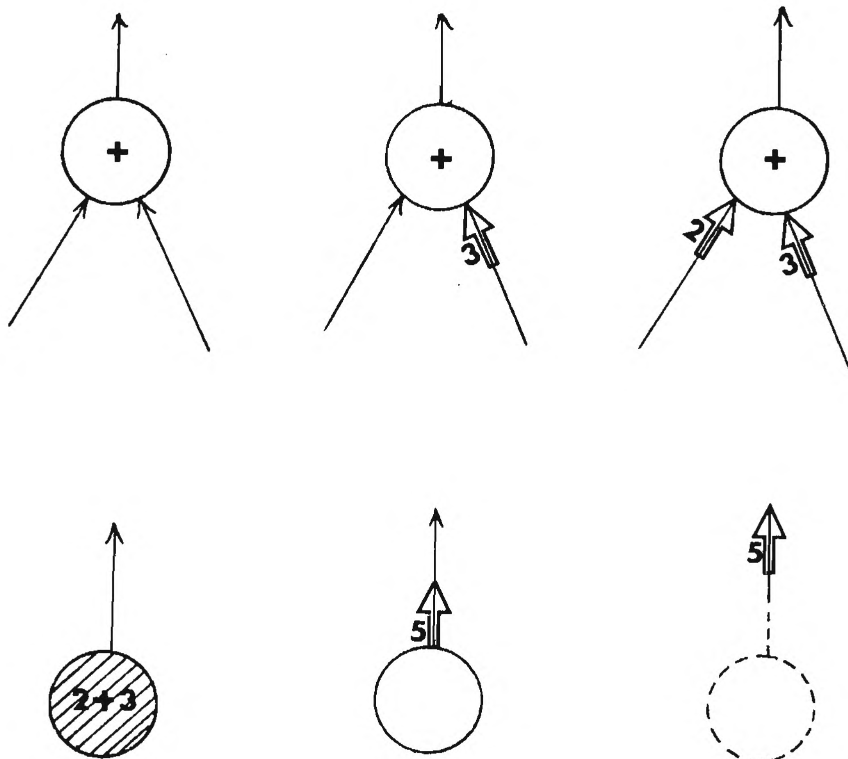

Fig. I-4

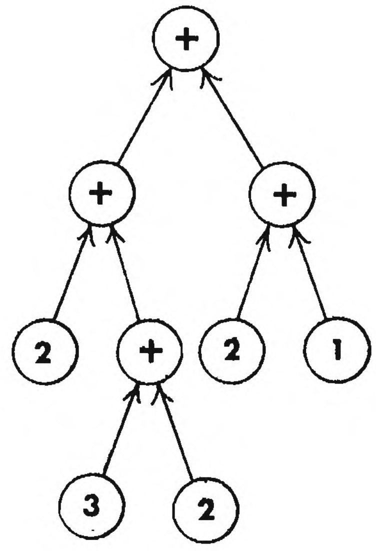

Fig. I-5a

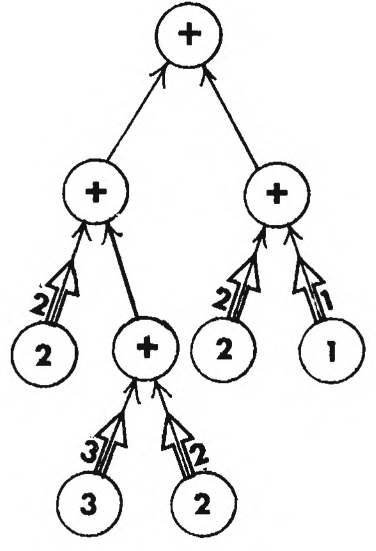

Fig. I-5b

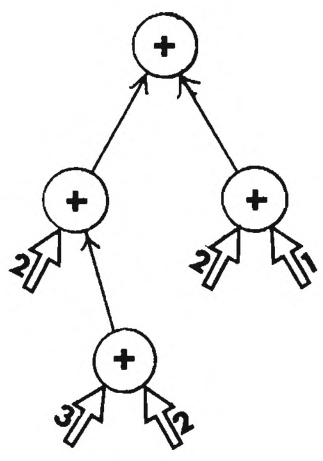

Fig. I-5c

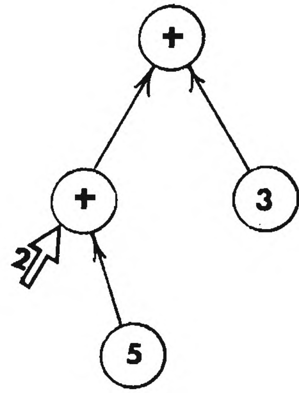

Fig. I-5d

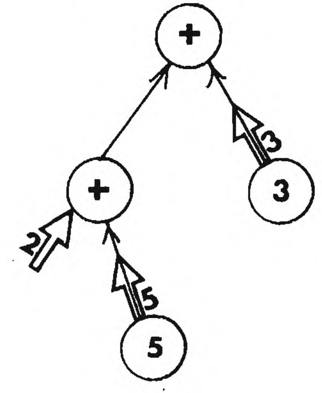

Fig. I-5e

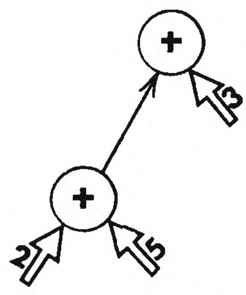

Fig. I-5f

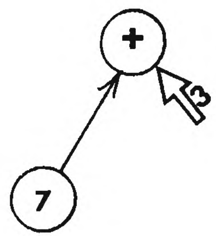

Fig. I-5g

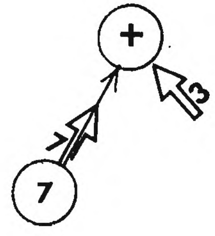

Fig. I-5h

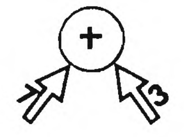

Fig. I-5i

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Fig. I-5j

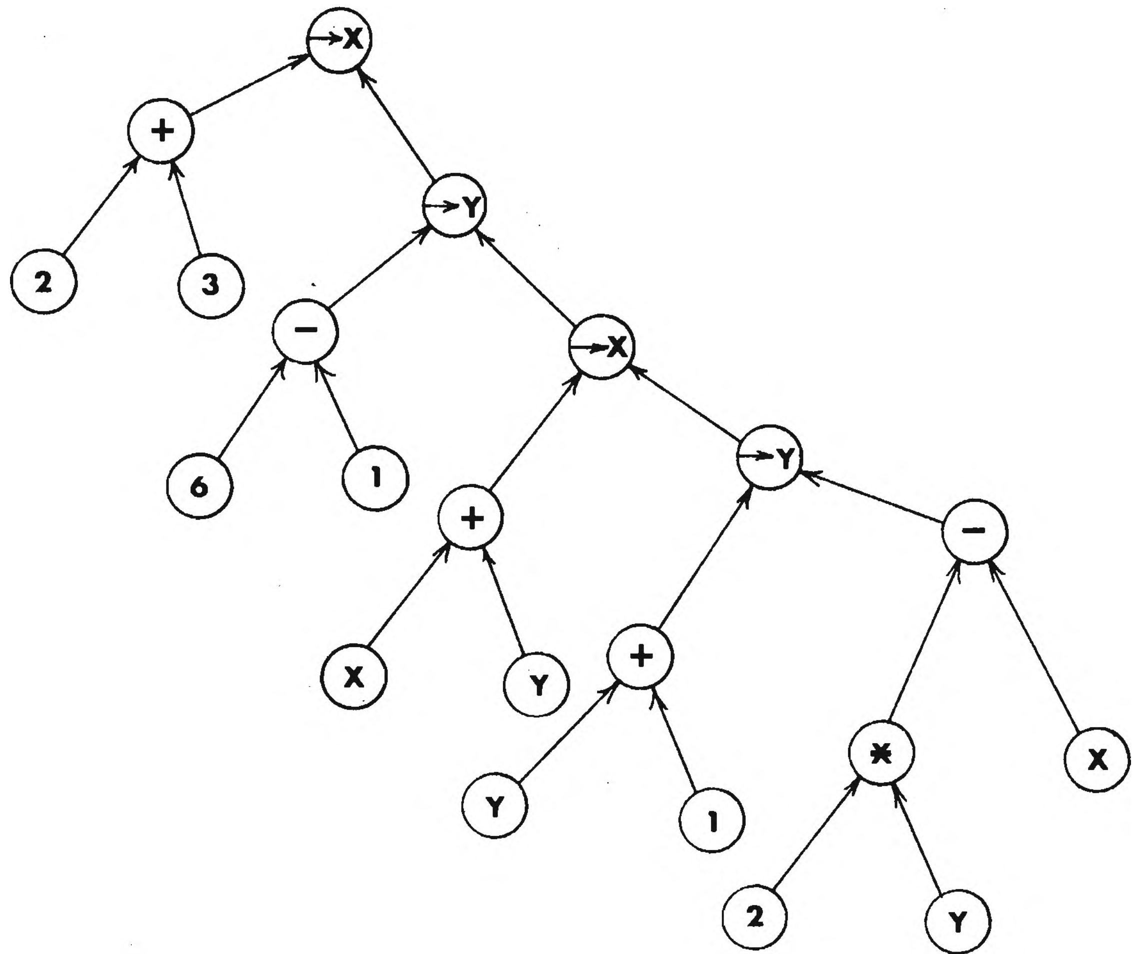

Fig. I-6

Each node is here again considered as being an automaton which has an address in some address space. Whenever a node A knows the address of another node B, the former may send a message to B (fig. I-7). Moreover any node may send a message up the tree; any node receiving such a message may either pick it up or pass it along upward.

The nodes representing values or operations have the same behavior as before. Whenever a node represents some variable x,y, etc..., it sends spontaneously up the tree its name and its address. Such a message can be picked up by a binder (→x , →y are binders) containing the same variable name only if the message reaches the binder by the right. The binder possesses then the address of the node which has originated the message (fig. I-8).

The various nodes which act spontaneously are completely independent from one another. However, in order to have figures

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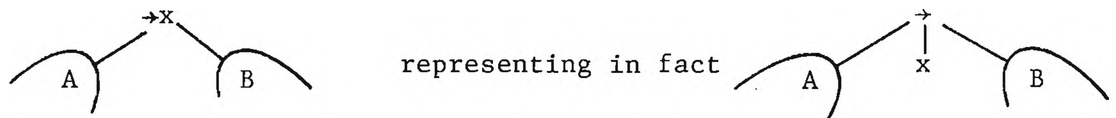
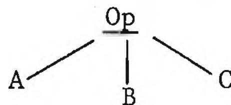
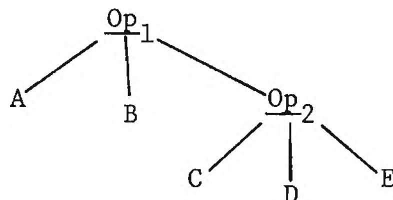
<u>Note</u> about the syntax: → is considered as being a triadic operator (operator of degree 3),



Whenever an operator <u>Op</u> is triadic, A <u>Op</u> B;C represents:



The association being done on the left, A $\underline{Op}_1$ B;C $\underline{Op}_2$ D;E represents:



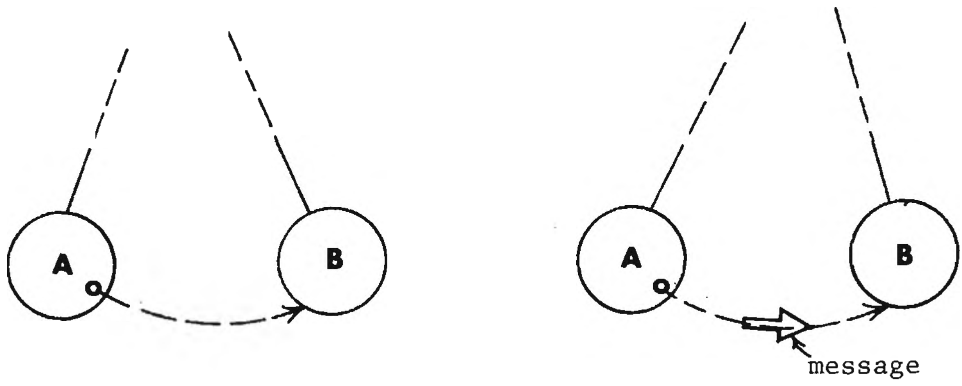


Fig. I-7

address of the originating node X

Fig. I-8

more readable and to display the binding in a more visible way, the binding operations are performed first.

Whenever a binder has received a value on its left, and addresses of variable nodes on its right, it sends the value received to each variable node whose address it has received (fig. I-9).

Whenever a binder receives a value from its right, it passes this value up along the tree and vanishes (fig. I-10).

The description of the evaluation is given in fig. I-11.

I.2.2 Variables as defining paths of information.

If we view a computation as a sequence of statements, variables are to be considered as denoting cells in which some results may be stored for subsequent use. Conversely if beforehand variables are considered as denoting cells, it is necessary to be assured that no attempt will be made to use a value before it has been produced, or to overwrite a value which is still to be used. As a result some sequencing has to be done, sequencing whose viscosity will decrease the amount of possible parallelism.

If we view a computation as occurring in space and in time, we will use variables whenever we want some information to be transferred from one place to some other place(s). Whenever a variable is free, it has no meaning by itself. A system of mutually bound variables is used as a symbolic device defining paths of information along which information flows toward computation. Whenever the needed information reaches an operator node, the simple process it represents would be activated asynchronously. Its completion may result in sending some information to some other nodes.

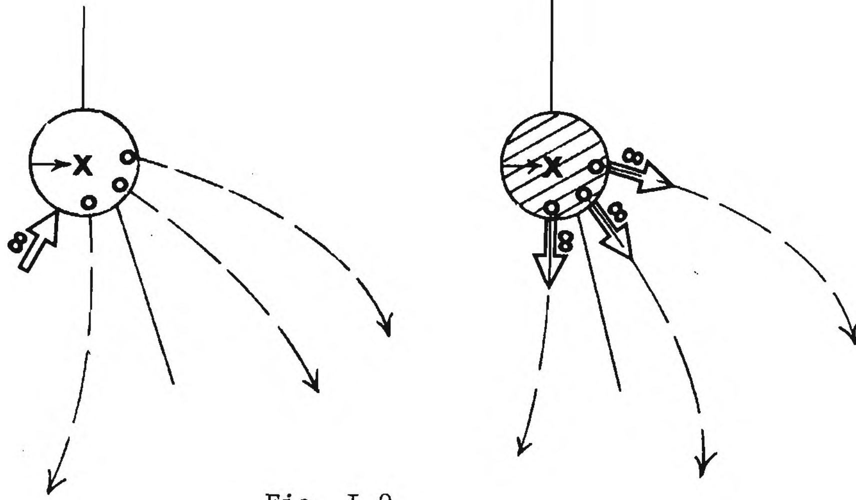


Fig. I-9

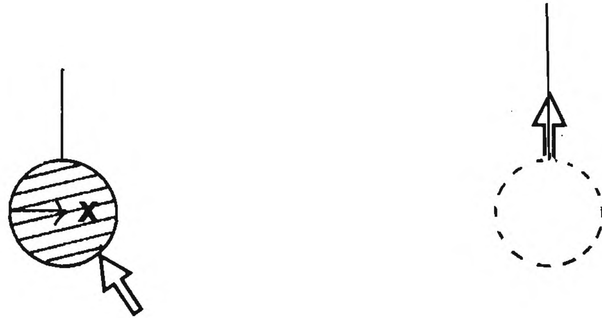


Fig. I-10

Fig. I-11a

Fig. I-11b

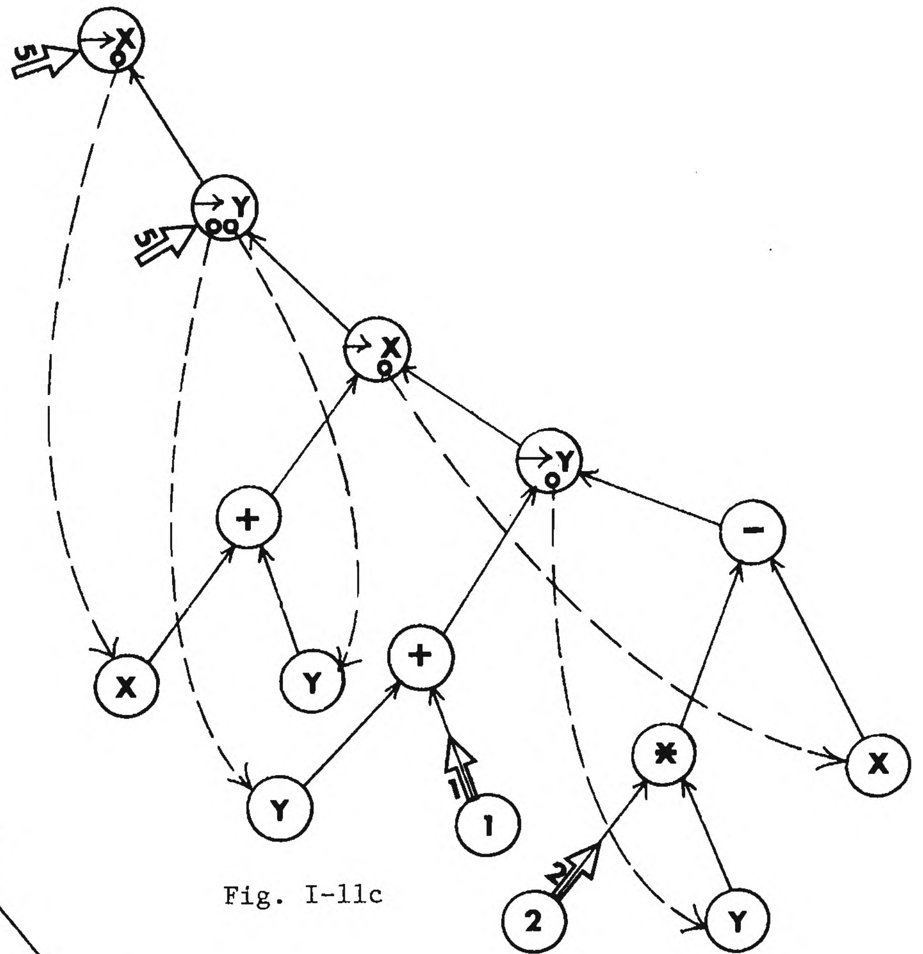


Fig. I-11c

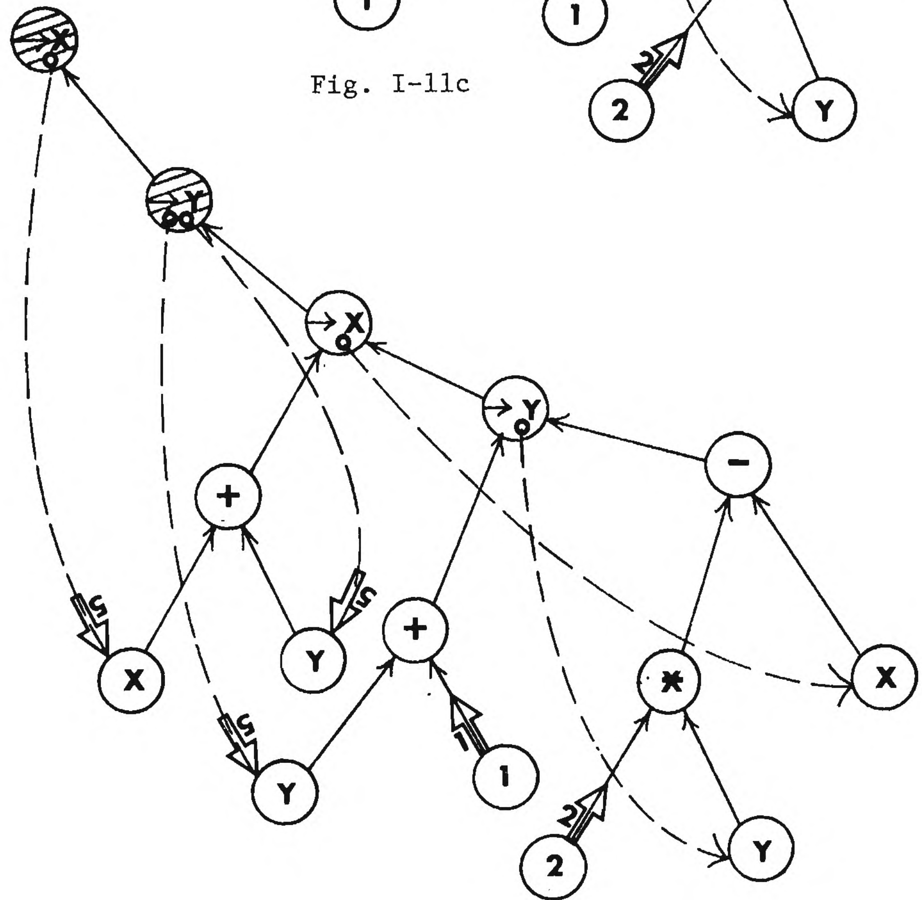


Fig. I-11d

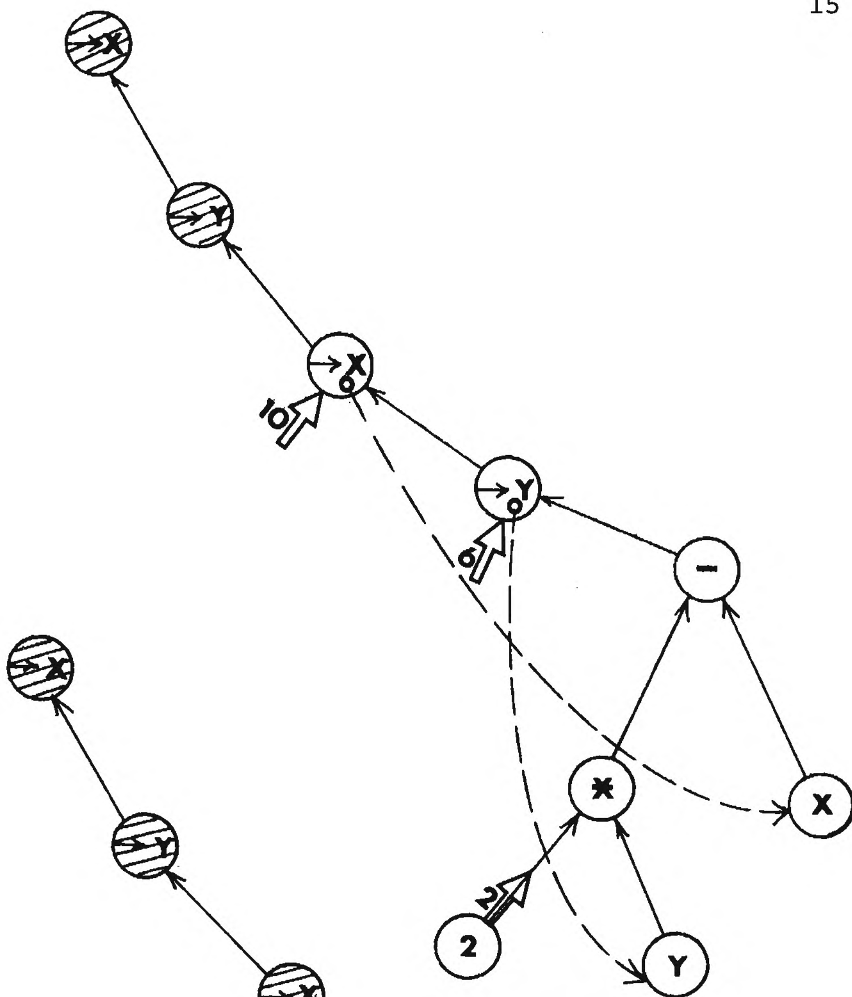


Fig I-11e

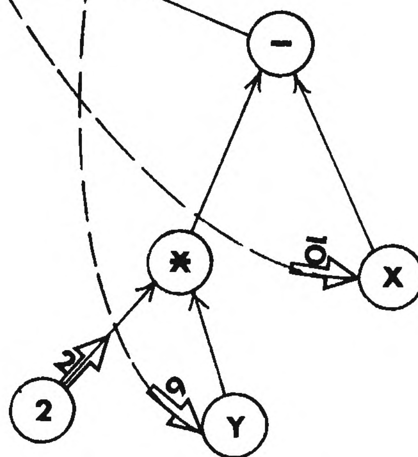


Fig. I-11f

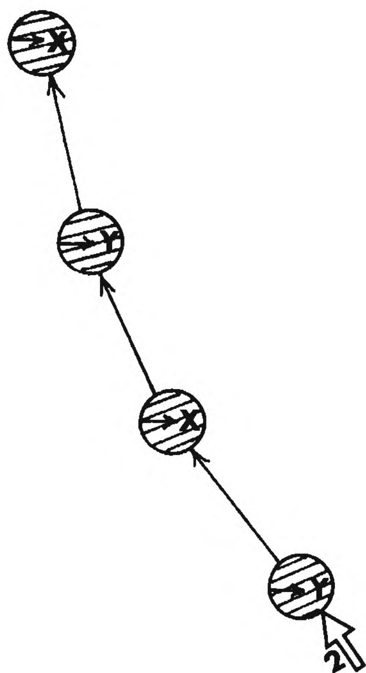


Fig. I-11g

Fig. I-11h

Fig. I-11i

In DCPL a computation may be regarded as an object composed of small automata which react to one another. Since a variable may receive a computation as value, representing for instance a procedural argument, the object representing a computation may expand and shrink with a behavior which may lead us to think of Von Neumann's self reproducing automata [24]. We may notice however that our structures are not to be implemented in some cellular space but programmed on a storage device.

The binding of variables superimposes to the tree structure a graph structure similar to a program graph. Fig. I-12, for instance displays the graph of our previous example. Such a program graph accounts for all the possible parallelism (or preferably concurrency) which may occur in the computation.

I.3 DCPL as a programming language.

As a programming language, DCPL has much in common with languages emphasizing expressions (rather than statements) and having to some extent the lambda calculus as background machine (McCarthy [18]; Landin [12,13,14]; A. Evans [7]). Any computation is a structured object whose evaluation produces a value. Moreover we can have procedural arguments: a procedure may be constructed in some place, produced as a value and sent to some other places where copies of the procedure are created (implemented in space).

However, DCPL presents many peculiarities: a computation, viewed as an object, may produce explicit side-effects on the environment in which the computation is embedded. Together with a binder 'lambda' which binds the variables which are to receive an argument from the environment (as in some extent "value" in

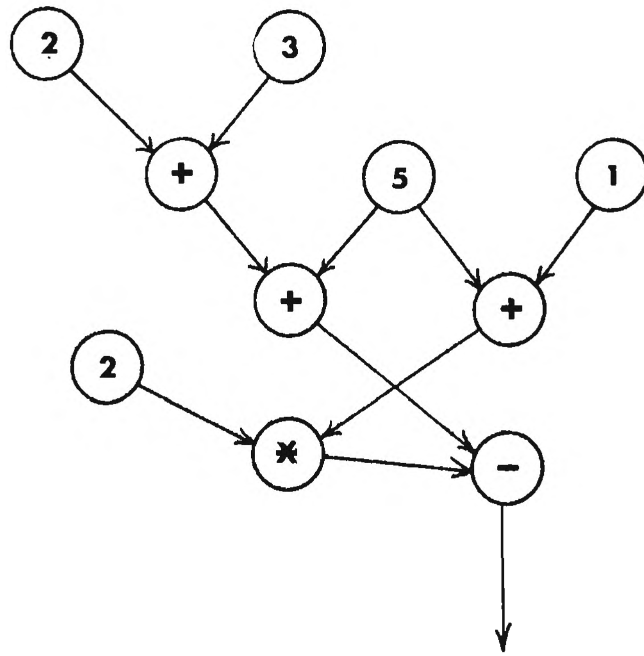


Fig. I-12

Algol 60), there is a binder 'mu' which binds the variables which are to send an argument to the environment (fig. I-13). 'Mu' , as a binder, is in some way similar to "result" in Algol W [26] with this important difference: an argument may be sent to the environment before the computation has been completed or even while the computation is actively worked out; as a result a computation may ask to the environment how to pursue the process or, some special conditions having occurred, if the computation is not to be cancelled , etc ...

As it has been already mentioned, DCPL is a programming language implicitly displaying parallelism to a large extent. Indeed, there is no need for special devices such as fork, join, parbegin/parend etc ... which determine parallelism explicitly.

DCPL is a system-oriented programming language, this aspect being discussed in the next section.

I.4 DCPL as a system-oriented programming language.

"System" is regarded here as denoting a group of interacting procedures constituting a collective entity. A sophisticated industrial organization, an administration, a hospital, are systems: a number of departments are services interacting to one another. If a computer is to be used integrated in such an environment, it is likely it should look like an information network; moreover programming should reflect such an organization.

In DCPL we are able to write asynchronously cooperating "independent" programs (coroutines) linked by paths of information along which messages are sent, and to write them recursively, i.e. any one of the previous programs may itself be a construct of cooperating independent subprograms (fig. I-14). It is possible to embed in DCPL sequential programs and to master their synchro-

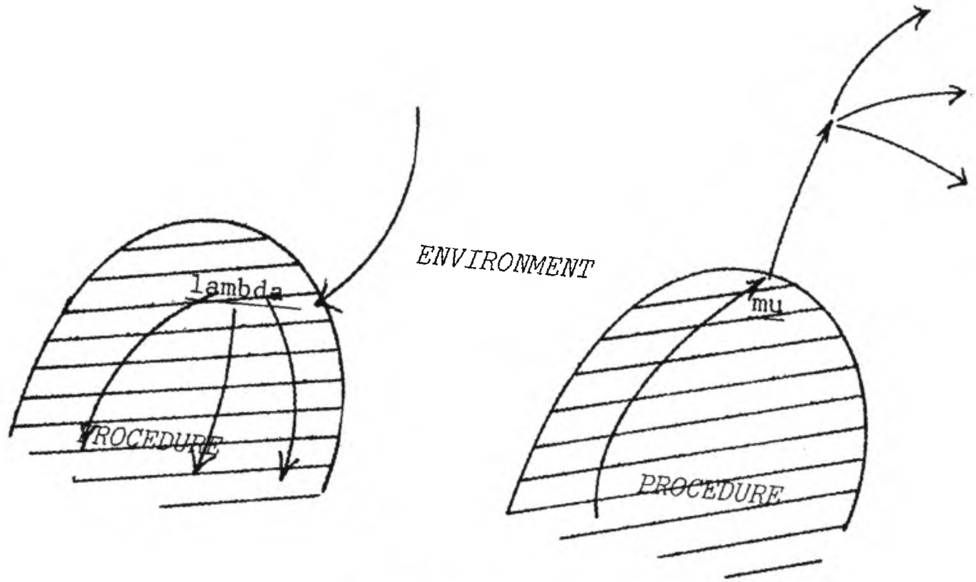


Fig. I-13

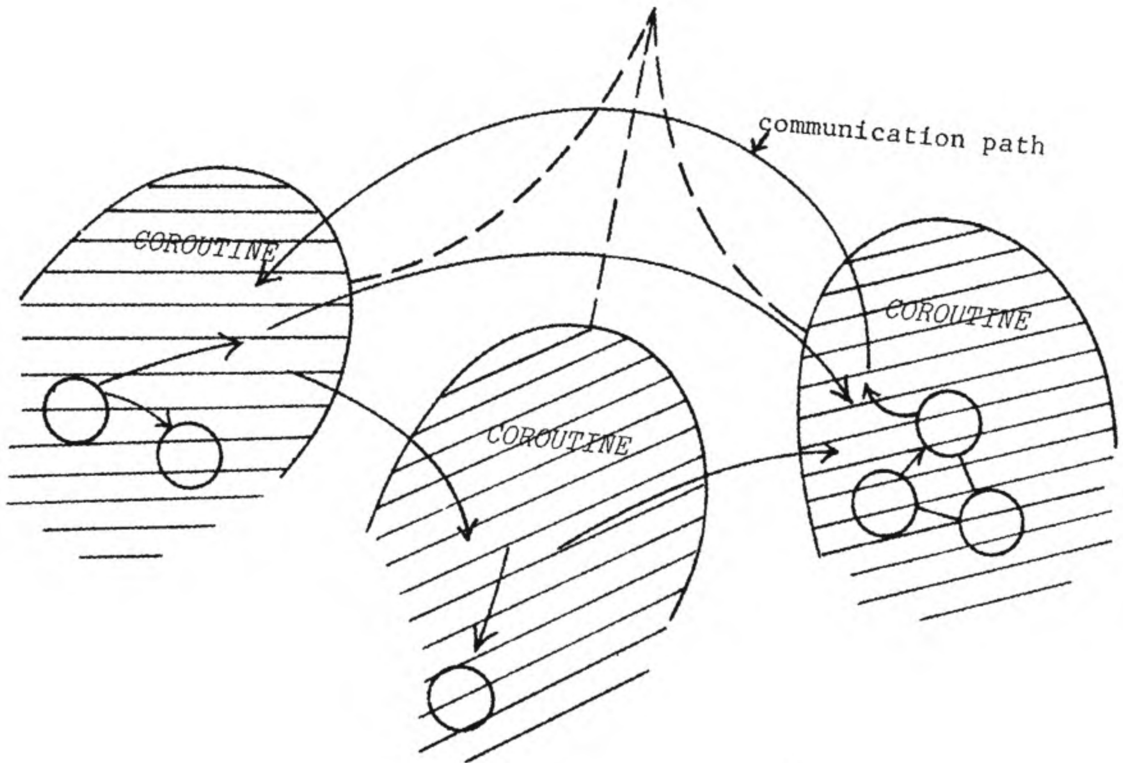


Fig. I-14

nization. For such programs DCPL looks like a host system . In fact it is possible to write in DCPL a hierarchy of host/guest systems.

Paths of information permit as well to have full programming generality: the same program may be debugged in a testing environment, made available in a program shop and put by some user in his own environment without the need of any surgery (Krutar [8]) . In conventional programming languages, such a programming generality may not be available for two reasons: 1. A procedure in general contains the names of some other procedures to call; as a result the former procedure is bound to the environment which contains the latter ones. The situation is better when procedural arguments are allowed. 2. Input/output operations are performed with particular instructions (read, write etc ...); thus it is not possible to debug a procedure in a testing environment with I/O devices simulated by some programs. In DCPL inputs and outputs are considered as paths of information coming into and going from the procedures. Such paths may be connected as well to I/O devices as to programs.

I.5 Machine organization.

DCPL gives to machine organization a new perspective.

Whenever a program is expressed in a current programming language, the control may jump from one place to any other one, a same cell may be accessed from quite different places. This results in a serious lack of locality. This would not be of any importance if today's computers were still Von Neumann type machines: one processor has access to a random access memory whose cells may be considered as being all "equidistant" from the processor.

Processors becoming faster and faster (and cheaper and cheaper) the trend in machine organization is to hierarchies of memories. However, unless many iterative computations are expected to occur in the fastest level, it is necessary to have at any level of the hierarchy a transfer rate large enough to "feed" the processor.

A large tranfer rate may be obtained by taking at each level a large block as unit of transferable information (the slower the level of memory, the larger the block).

One may believe, however, that only a few words in such blocks would be really used. For this reason Jack Dennis suggests in [4] that information should be moved on demand with the word as information unit, a large transfer rate being assured by performing many computations in parallel.

In DCPL it is possible to consider a program as a construct of "simply-connected" computations which, once triggered, could be brought in the fastest level of memory and completed without the need for any additional information.

Moreover it is possible to replace random access memory by sequential rotative memory. This will be discussed extensively in part III of this thesis.

In part II DCPL is presented. Part one is concerned with preliminaries and discussions which the author believes to be relevant to the subject and important. Some readers might prefer to skip them and go directly to part two (beginning p. 86).

PART ONE

PRELIMINARIES

CHAPTER II

SOME CONSIDERATIONS ABOUT PROCESSES

This chapter is intended to present some notions about simple processes . We are indebted to Holt [9] and Shapiro [22] for some of the concepts presented here.

II.1 A process as a sequence of transformations .

Some simple processes may be considered as leading from one object to some other object through a sequence of elementary transformations.

II.1.1 The object-flow model .

The process illustrated in fig. II-1 may be visualized by considering each node as being a processor able to perform the corresponding elementary transformation, and considering each arrow as being a communication path. The object enters on the left. It then advances along the path. On reaching a processor, it triggers the corresponding transformation. Then the resulting object continues to advance along the path. When the process is completed the result emerges at the right end.

II.1.2 Control signals .

Another point of view may be adopted. There are places where objects may be stored and retrieved. A processor does not receive on the path the object to be transformed, but a <u>control</u>

Fig. II-1

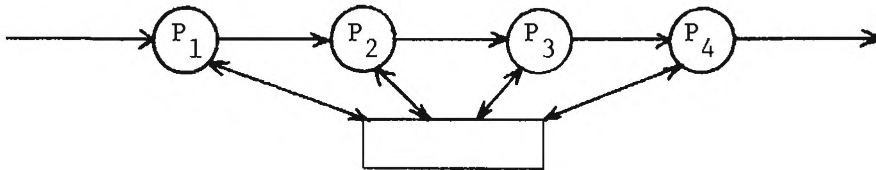


Fig. II-2

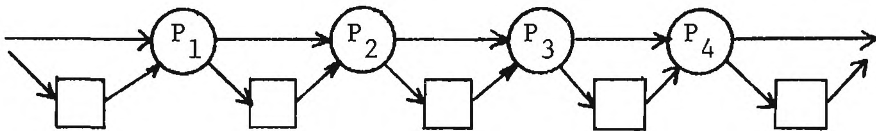


Fig. II-3

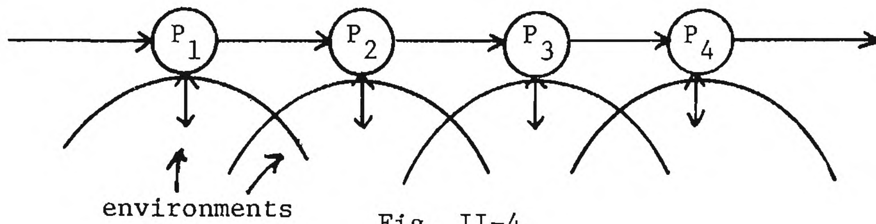


Fig. II-4

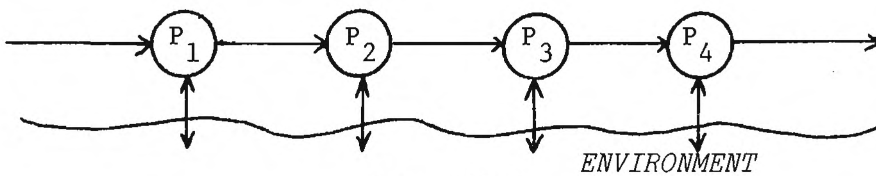


Fig. II-5

signal. On receiving it, the processor expects to find the object to be transformed at some specific place. When the transformation has been completed, it puts the resulting object at some (possibly different) place and sends a control signal to the next processor to operate. In fig. II-2 one single place is used; a cascade of places is used in fig. II-3.

The latter situation may be somewhat abstracted. To each processor is associated an environment (environments may overlap one another). Each processor is able to perform a specific transformation on its environment. The process is controlled by a control signal as before (fig. II-4). In fig. II-5 there is one common environment accessible by each processor.

In the light of these latter interpretations the arrows in fig. II-1 appear to have two purposes: 1. They order in time the occurences of the different transformations, carrying an implicit control signal (which is the object itself); 2. They specify, for each processor, on which object the corresponding transformation is to be done.

II.1.3 Production line.

Up to now we were interested in transforming one object into another one. Let us suppose we want to apply a process to a sequence of incoming objects, fig. II-1 being interpreted as representing a production line.

For instance we may consider a row of objects advancing on the communication path. Any processor performs repetitively its transformation on each incoming object. If the various processors operate at different rates, it becomes necessary for each arrow to act as a first-in-first-out

queue (fig. II-6).

In some instances, it may be preferable to just have one queue before the process. It is then necessary to synchronize in some way the various processors.

A straightforward solution is to process only one object at a time, one processor at most being at work at any time: P_1 is not to accept any new object before P_4 has completed the transformation of the current one. As shown in fig. II-7 , a backward path links P_4 to P_1. On completion of its task P_4 sends a control signal on this path. On receiving it, P_1 initiates the processing of a new object.

Let us abstract the situation in the following way. We represent an object as well as a control signal by a token placed on the corresponding arrow. An elementary transformation is triggered whenever all the incoming arrows contain a token (fig. II-8b): the tokens are then removed from these arrows and the transformation is in progress (fig. II-8c). On completion of the transformation, a token is placed on each outgoing arrow (fig. II-8d). The processor then stays idle until there is again a token on each incoming arrow (fig. II-8a).

The process is described in figures II-9a to II-9j .

A similar scheme may be used whenever the process uses places to store objects as in fig. II-2 and in fig. II-3 (fig. II-10 and II-11). There is one token representing a control signal, which performs a loop (the servicing of the input and the output in these examples is not discussed here).

Fig. II-6

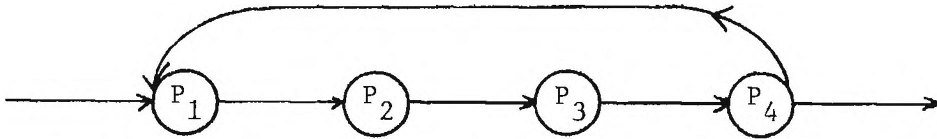


Fig. II-7

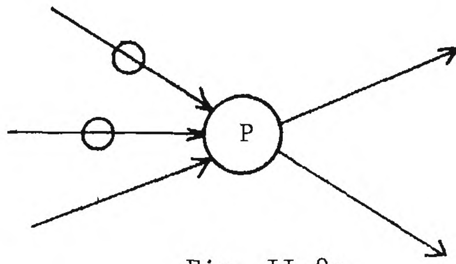


Fig. II-8a

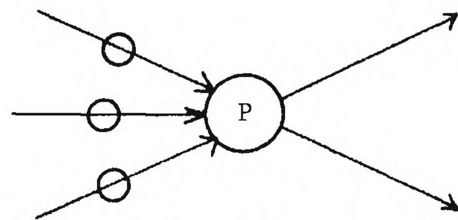


Fig. II-8b

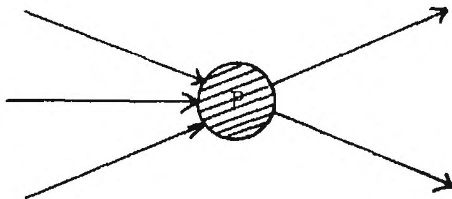


Fig. II-8c

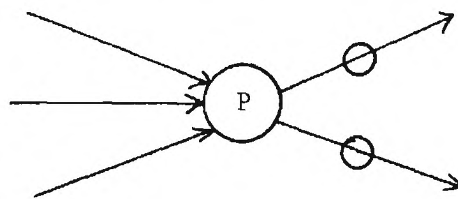


Fig. II-8d

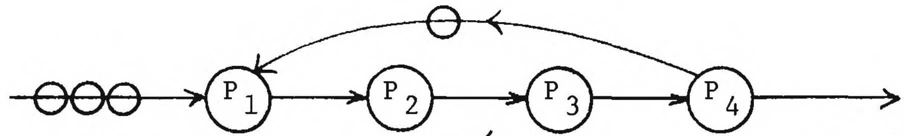


Fig. II-9a

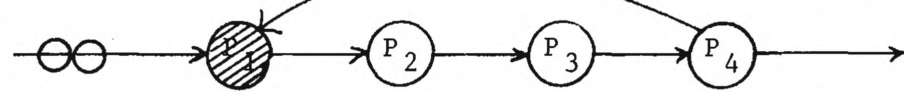


Fig. II-9b

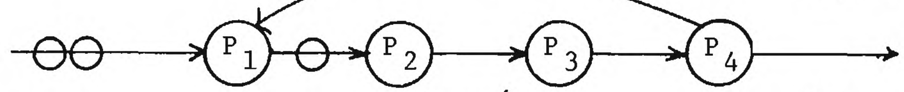


Fig. II-9c

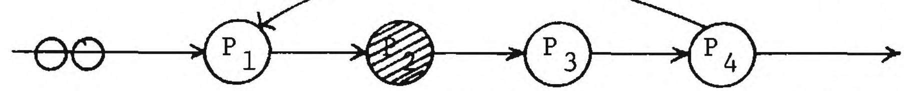


Fig. II-9d

Fig. II-9e

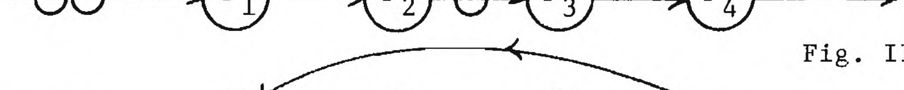


Fig. II-9f

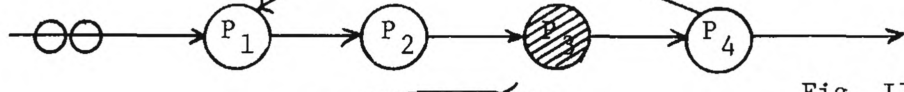


Fig. II-9g

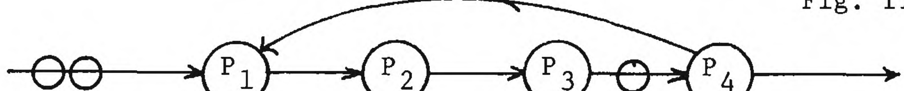


Fig. II-9h

Fig. II-9i

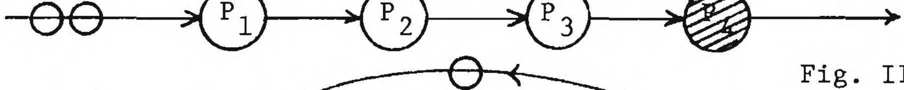


Fig. II-9j

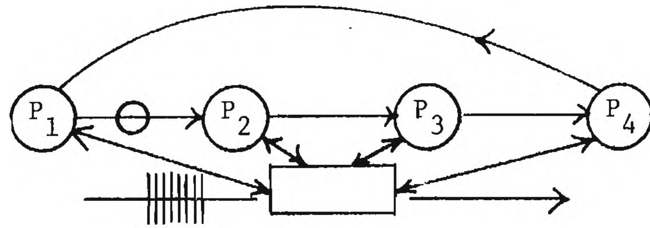


Fig. II-10

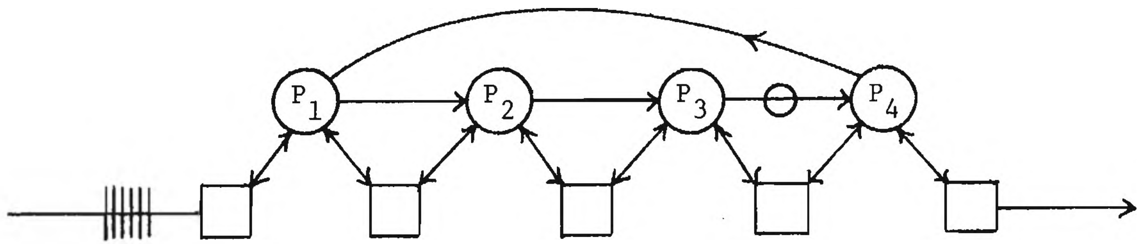


Fig. II-11

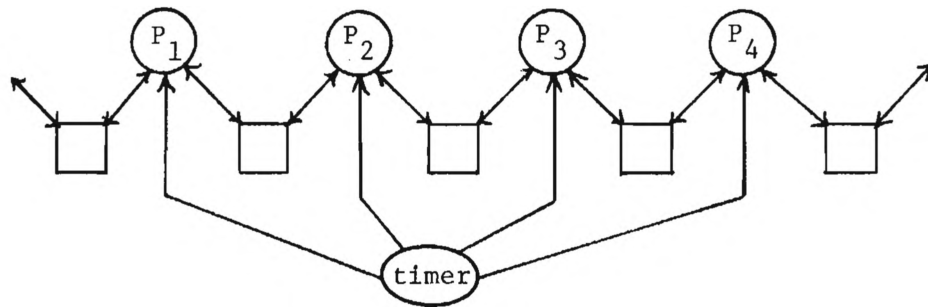


Fig. II-12

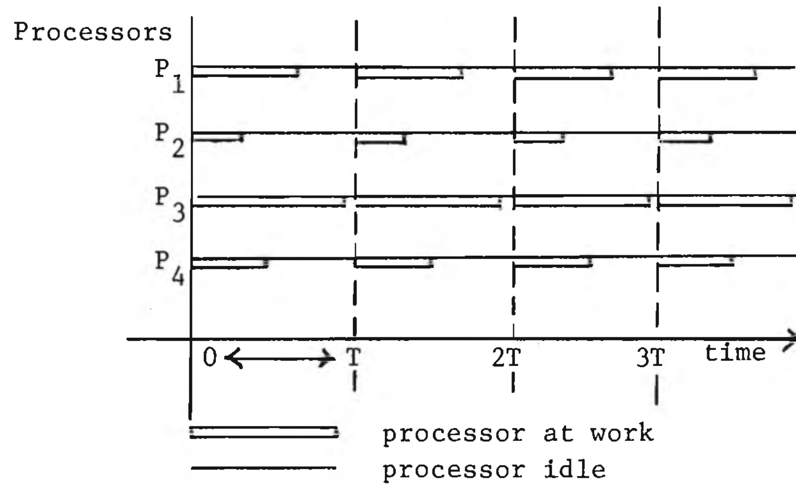


Fig. II-13

II.1.4 Pipe-line.

If a large throughput (number of objects processed per unit of time) is desired, such straightforward solutions are not satisfactory: at any time only one processor is working; we would prefer to have the various processors working concurrently on different objects.

If the time required by one processor to transform one object is independent of the object and if this time is known, synchronization may be performed with a central timer. Every T (an amount of time during which any processor may perform its task on one object) the timer triggers all the processors by sending them a control signal (fig. II-12 and II-13).

A more general solution may be achieved by replacing the backward path in fig. II-7 by a sequence of backward paths as displayed in fig. II-14. Fig. II-15a-d describe the process.

The processes we have just described are called pipe-lines. In fig. II-12 the pipe-line is synchronous; in fig. II-14 the pipe-line is asynchronous.

II.1.5 Petri-net.

The situation of fig. II-14 may be modeled by a Petri-net (fig. II-16).

The behavior of a Petri-net is very close to the behavior described in fig. II-8. A Petri-net is made out of transitions (bars in fig. II-16), and places (circles in fig. II-16). An arrow may lead from a place to a transition, or from a transition to a place; a place in the former case is called an input place of the transition, and in the latter case, an output place of the transition

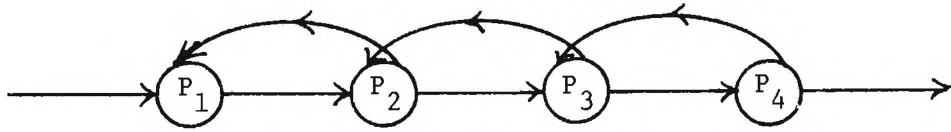


Fig. II-14

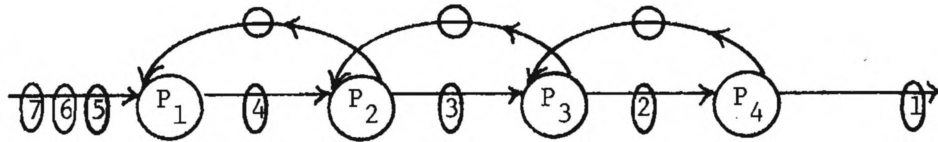


Fig. II-15a

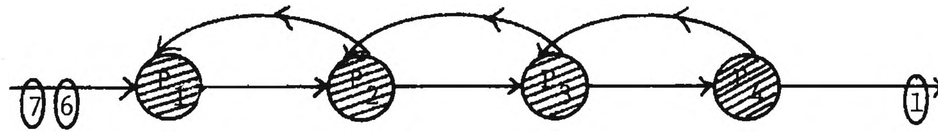


Fig. II-15b

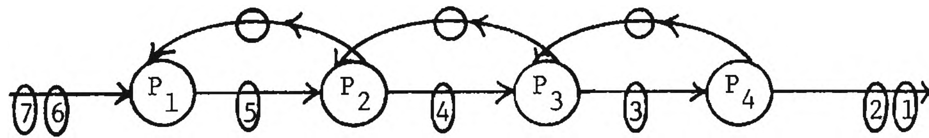


Fig. II-15c

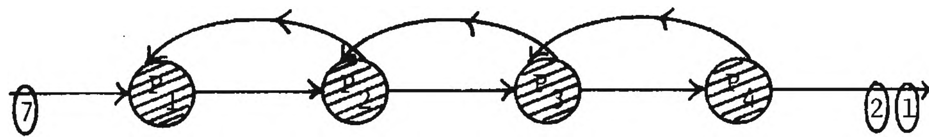


Fig. II-15d

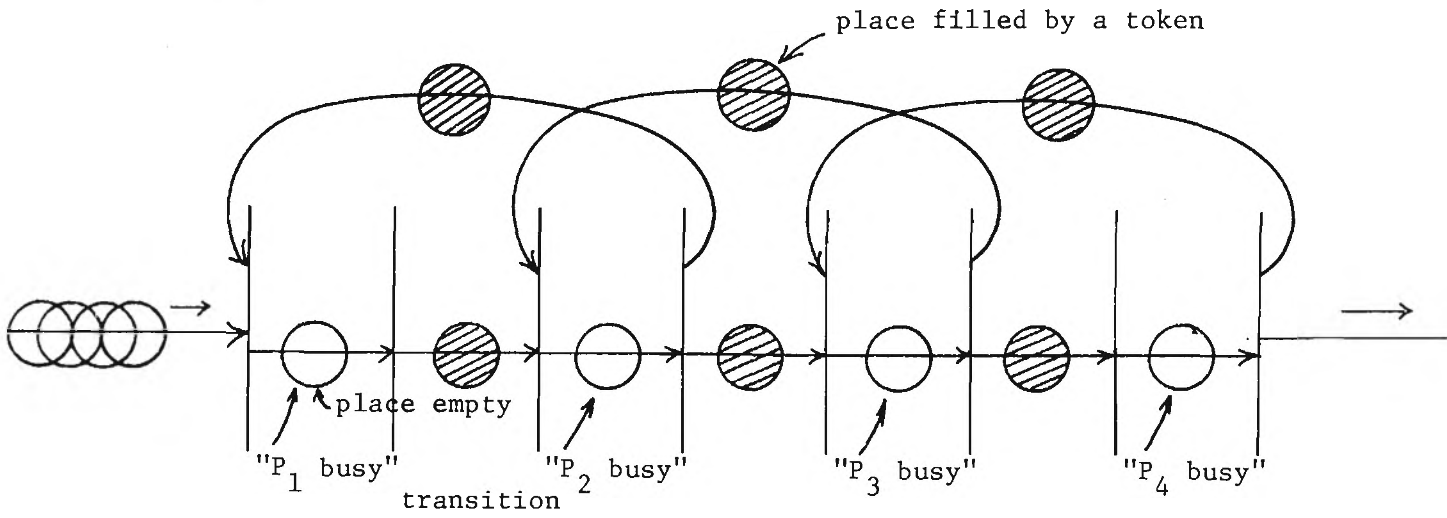


Fig. II-16

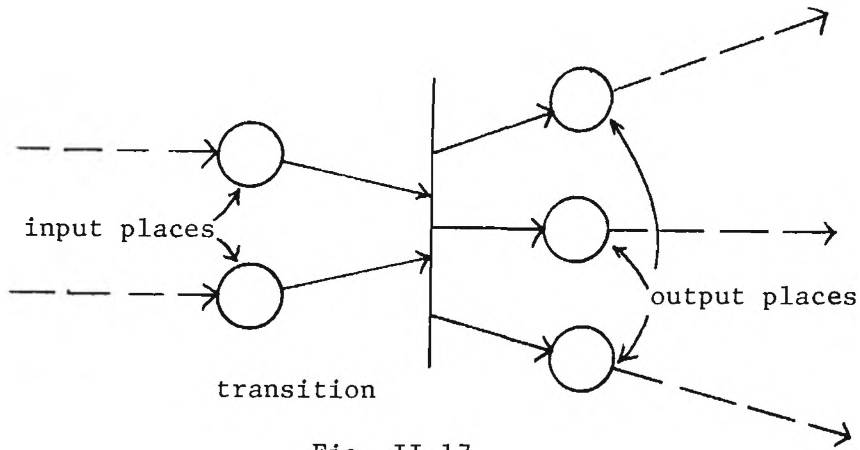


Fig. II-17

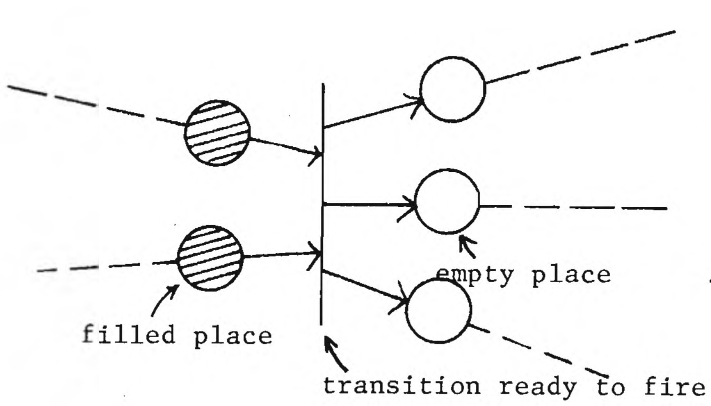


Fig. II-18a

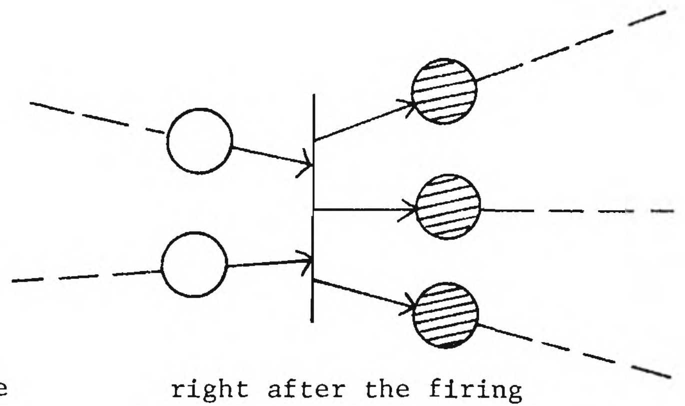


Fig. II-18b

(fig. II-17). A place is either empty or filled by a token. A transition is ready to fire whenever all its input places are filled. (It is supposed that the Petri-net is such that all the output places are empty at the time of firing. For a detailed discussion, please see Holt [9]). The firing of a transition may be viewed as a spontaneous and instantaneous operation: each input place is emptied, a token is placed in each output place of the transition (fig. II-18a-b).

It may be interesting to note that a place may be the input of several transitions. Several of these transitions may be ready to fire at the same time; however only one transition among them may fire at a given time (fig. II-19a-b).

Remark.

A given situation may be modeled in quite different ways depending on the emphasis which is to be placed on various conditions. For instance, fig. II-20 and fig. II-16 account quite differently for the same situation.

II.2 A process as a system of transformations.

In the previous section we discussed simple processes whose elementary transformations were applied sequentially to some object.

In general an elementary process may be applied to one or several inputs producing one or several outputs (fig. II-21a). We require of any would-be elementary process that it may only be applied when all its inputs are available and that its outputs are available only after the process has been completed (fig. II-21b).

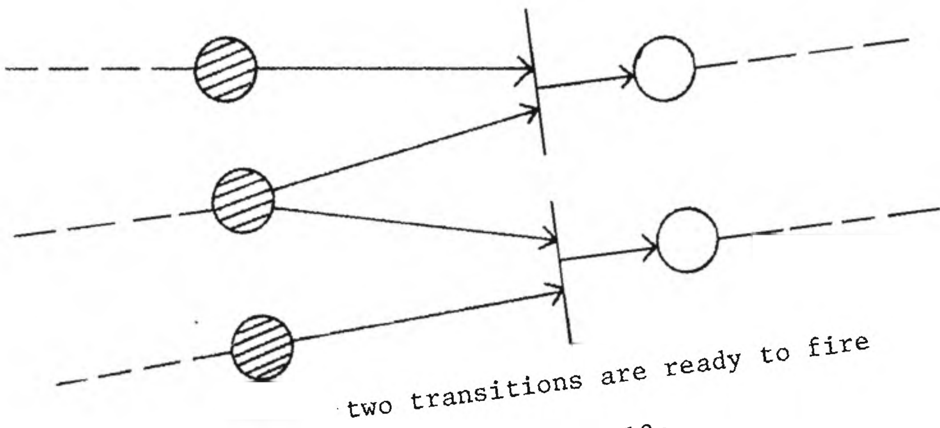


Fig. II-19a

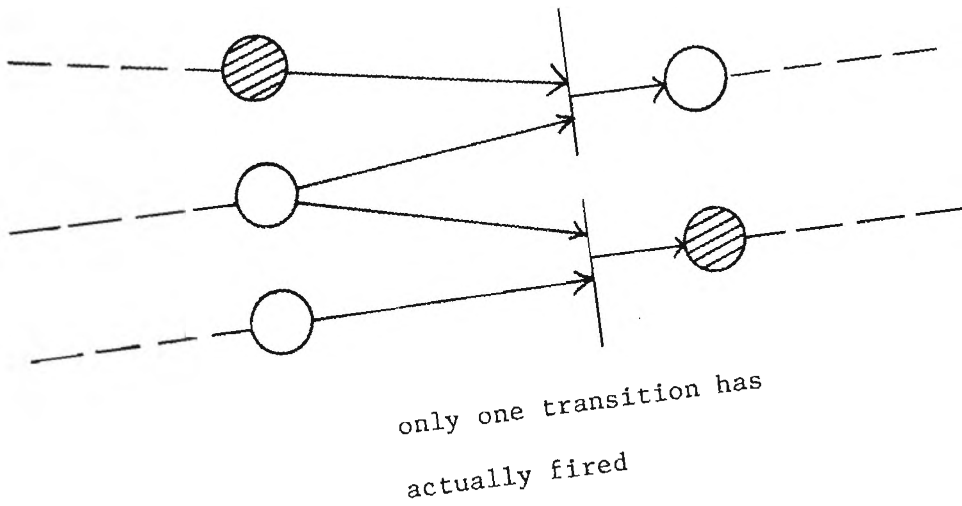


Fig. II-19b

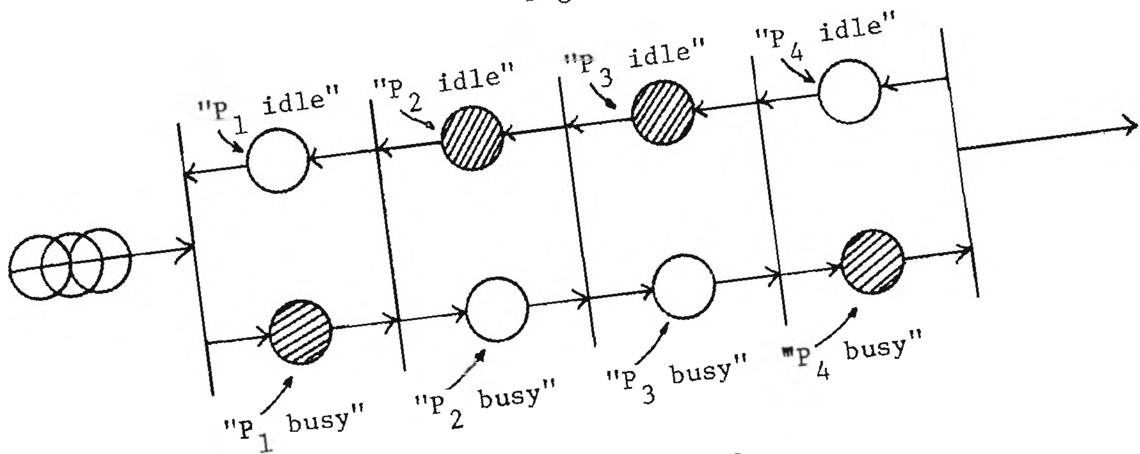


Fig. II-20

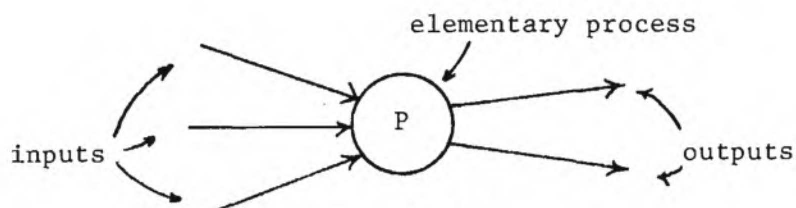


Fig. II-21a

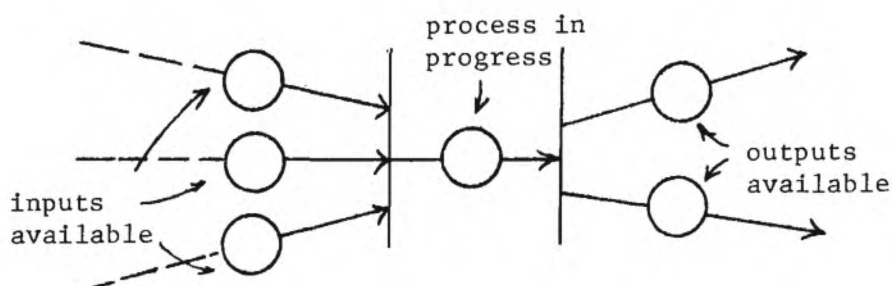


Fig. II-21b

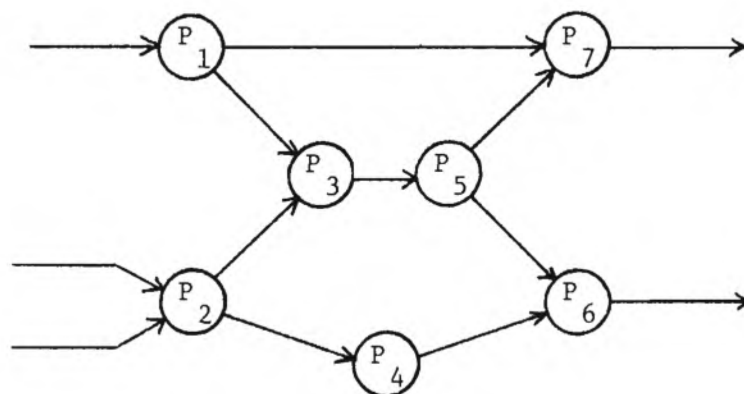


Fig. II-22

II.2.1 The object-flow model.

Let us now consider a process leading from a set of objects to another set of objects by applying different elementary processes $P_1, P_2, \ldots$ as indicated in fig. II-22.

By and large, our previous discussion may be applied here again. Each node may be considered as being a processor. Whenever an object is produced it advances on the corresponding path until it reaches the next processor. A processor is triggered whenever there is an object on each of its input lines.

II.2.2 Control signals.

In another implementation there are places in which and from which objects may be stored and retrieved. Control signals may travel on the arrows. A processor is triggered whenever it has received a control signal on each of its input lines. The processor retrieves the objects stored in its input places, performs the associated elementary process, stores the resulting objects in its output places, and sends a control signal on each of its output lines. Fig. II-23 and II-24 present two possible configurations.

II.2.3 Determinacy.

We may abstract such a situation by considering each processor as being able to perform a transformation on its environment after it has been triggered by some control signals (fig. II-25). However it is important to notice that the result of the process may depend on the order in which elementary processes have been applied if the different environments are interdependent or if there is one global environment. A process is said to be <u>deterministic</u> or

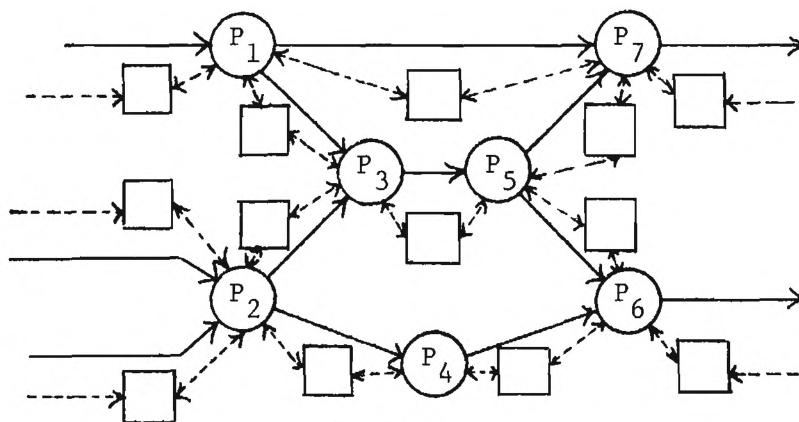


Fig. II-23

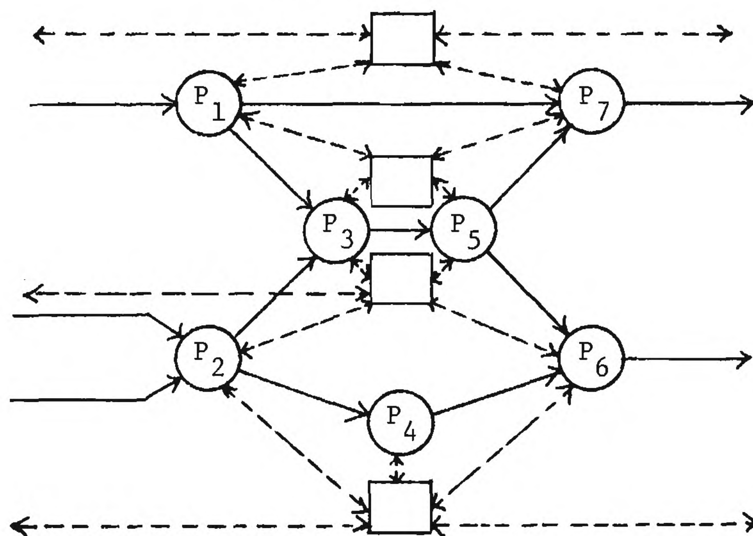


Fig. II-24

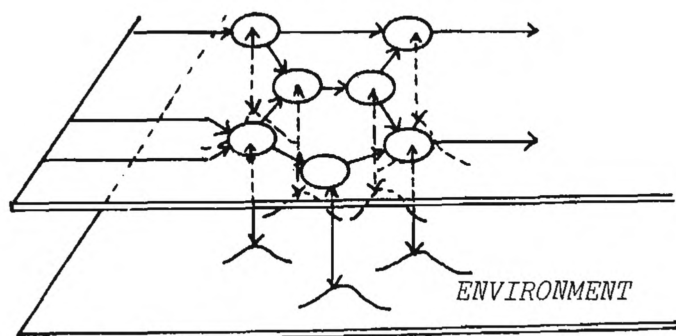


Fig. II-25

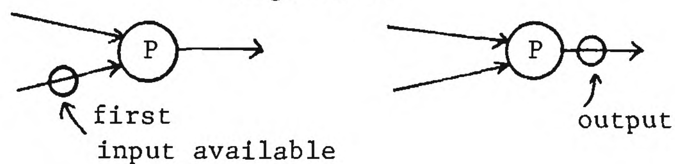


Fig. II-26

output-functional if the resulting objects or the resulting environments do not depend on the order in which elementary processes have been applied [15].

A process is said to be completely functional[a] if, for any elementary process P_i, the objects or the local environment to which P_i is applied do not depend on the order in which elementary processes are applied, at the time when P_i is triggered.

The process of fig. II-22 with objects flowing on arrows is determinate if the elementary processes are output-functional; we owe this result to the requirement we made about elementary processes (fig. II-21). However the process would not be determinate if we had allowed races to occur with, for instance, an elementary process having as an output the first input received (fig. II-26).

The processes of fig. II-22 and II-23 are determinate; however, some other configurations of places would have given non-determinate processes.

II.2.4 Service-on-demand.

A same elementary process may occur at different locations in a process. Let us assume that the process of fig. II-27 is to be carried out by a team of four workers, each worker being able to perform a specific elementary process A,B,C or D.

The flow chart of the process is displayed on a table with a light bulb on each node and a place on each arrow. Whenever all

---

a) The term "determinate" is often used instead of deterministic. However determinate is sometimes opposed to deterministic and means then completely functional. Such opposition is not a very important one from the point of view of this thesis, and we shall use "determinate" in general. However, whenever the opposition is meaningful we will use deterministic and completely functional.

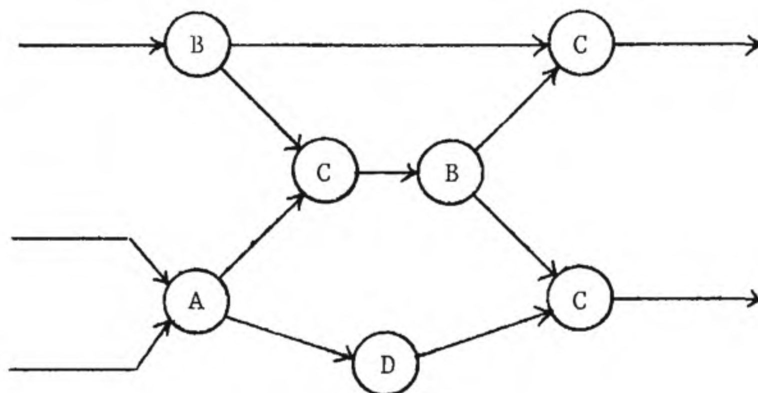


Fig. II-27

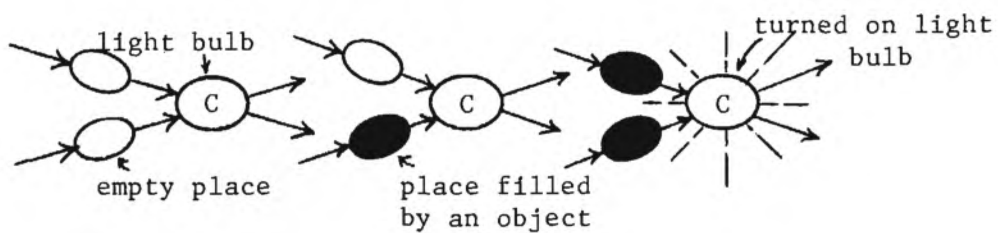


Fig. II-28a Fig. II-28b Fig. II-28c

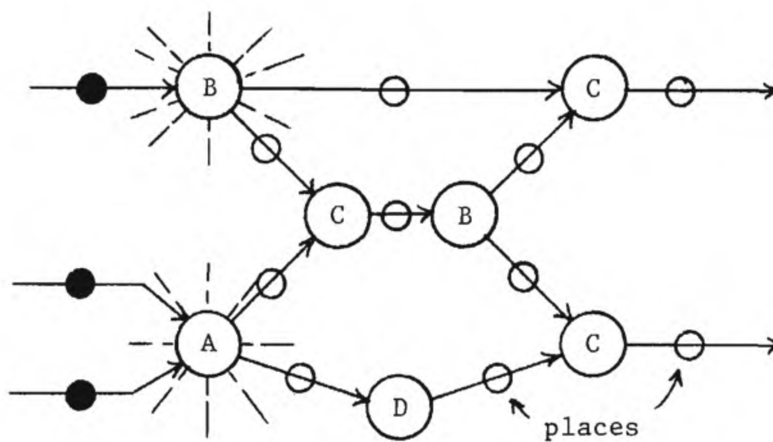


Fig. II-29

input places of a node are occupied by objects, the bulb on this node is automatically turned on (fig. II-28).

At the beginning the objects to be processed are put on the input places (fig. II-29).

Whenever a worker is idle, he looks for a turned-on bulb representing an elementary process he is able to perform. On finding one, he takes the objects from the input places, performs the corresponding elementary process and puts the resulting objects on the output places; then the worker is idle again.

In such an organization, the workers are said to <u>service on demand</u> the elementary processes. The demands are issued locally whenever some local condition occurs. The same scheme may be used if one worker is able to perform more than one elementary process. However, if several workers were able to perform a same elementary process and if they could have access concurrently to the flow chart, a system of arbiters(assigning to each idle worker an elementary process to be performed) would be necessary (see Patil [20]).

Let us consider again fig. II-22. Two elementary processes are said to be <u>ordered</u> if there is a path of arrows leading from one to the other. The executions of two such elementary processes are then necessarily ordered in time. Two elementary processes are said to be <u>concurrent</u> if they are not ordered. When the whole process is carried out, two such elementary processes may be performed either concurrently or one after the other, according to considerations which are not relevant to the level of organization in which we are interested here. In the situation of fig. II-25, a sufficient condition for the process to be deterministic is that any pair

of concurrent elementary processes (considered as operators on the environment) commute. A sufficient condition for the process to be completely functional is that any pair of concurrent elementary processes have "independent" environments (fig. II-30). This latter condition holds in the situations displayed in fig. II-23 and II-24.

II.2.5 Sequential execution.

Let us now suppose that the process of fig. II-27 is to be performed by one worker executing the different elementary processes one after the other. At any time there is a set of pairwise concurrent elementary processes which are ready to be executed (in our previous scheme these processes are those whose bulbs are turned on).
Our worker may choose at random one process in this set and performs it. The occurence of this elementary process is no longer in the set, but at its completion, new elementary processes might be in the set. Then the worker repeats the same sequence of actions until the whole process is completed.

If the process is planned to be executed by only one worker, it may be described as a sequence of elementary processes. Obviously different sequences are possible (fig. II-31 and II-32). Together with the sequence we have to know the inputs and the outputs of each elementary process and how one is related to another. In fig. II-33 corresponding inputs and outputs are linked together. In fig. II-34 there are four places; to each input (resp. output) is associated a place from which (resp. in which) the object is to be taken (resp. stored). These figures display the same topologies as fig. II-32 and fig. II-35.

Fig. II-30

Fig. II-31

Fig. II-32

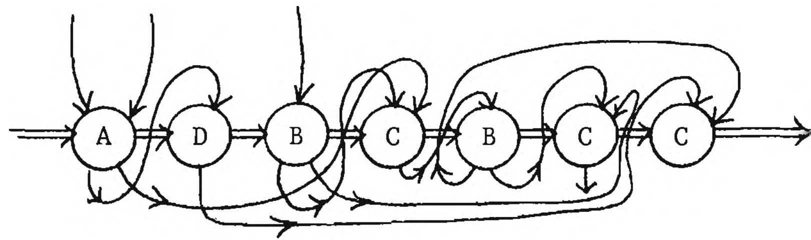


Fig. II-33

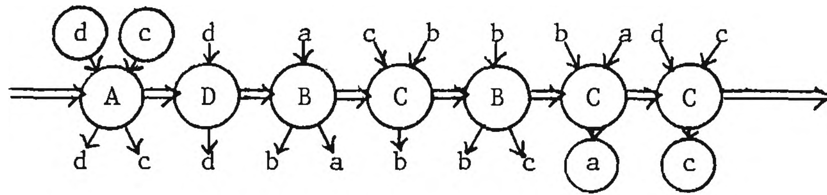


Fig. II-34

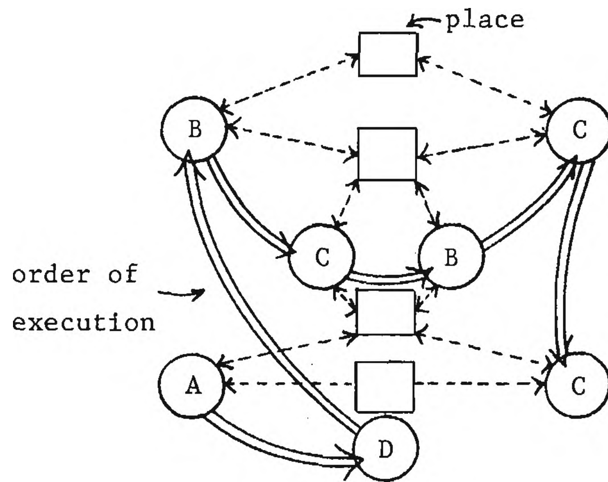


Fig. II-35

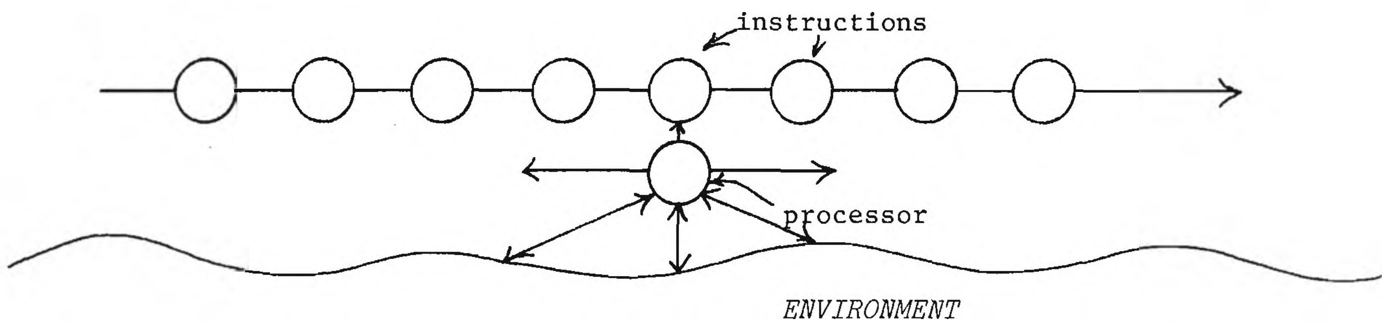


Fig. II-36

II.2.6 Centralized vs. distributed control machine.

Abstracting the situation, we may consider the description of the process as a sequence of instructions. One processor scans such a sequence performing one at a time the corresponding elementary transformations on the environment (fig. II-36).

We call a machine with one or several processors working in such a way a centralized control machine. On the other hand we call a distributed control machine a machine whose special purpose processors act on a service-on-demand basis: any action is triggered by local conditions occurring asynchronously.

Although such a distinction may appear rather artificial under certain circumstances, the notion of a distributed control machine will be helpful in the sequel.

CHAPTER III

A COMPUTATION AS THE REALIZATION OF A FORMAL OBJECT

So far, we have not particularized our discussion to any special object or to any special elementary process. In this chapter we consider formal objects called obs which may be built up, and then realized relatively to our universe of values and operations.

III.1 Combinations and obs.

The following notions are due to Curry [3]. They are very similar to Landin's notion of applicative structure [12]. In many other works, such notions are expressed from a syntactic point of view.

An _ob_ is a formal object which is either _primitive_ and called an _atom_, or built up by combining already constructed obs according to some elementary processes called _combinations_. A combination is a particular elementary process: an ob being given, we know at once what is its structure; we know whether or not the ob is an atom, and if the ob is not an atom we know how it has been constructed: by applying which combination to which obs. It is therefore understood that obs constructed by different processes are different as obs.

As a result, concatenation of symbols and strings, into strings, is not a combination. In Lisp, forming the "cons" of two S-expressions is a combination. In the same way a sentence

which is syntactic relatively to a phrase structure grammar is not an ob; however, any syntax tree of the same sentence may be considered as an ob.

A symbol denoting a combination is called a combinator. The degree of a combination or of a combinator is the number of obs to which the combinator is applied. An atom is considered as a combination of degree zero.

Fig. III-1 represents an ob O, with C_0 , C_2 and C_3 being respectively combinators of degree zero, two and three. The figure displays the "topology" of O, i.e., the different obs, components occurring in the construction of O.

III.2 Realization of a process or of an object: categories and functors.

Let us consider an elementary process with input and output places (upper part of fig. III-2). When realizing such a process, each place is mapped into a value of the universe of values and operations (see introduction), each elementary process into an operation, such that the value associated to the output place is obtained by applying the operation (associated to the elementary process) to the values associated to the input places (fig. III-2).

It is possible to realize an ob by assigning a value to each component and an operation to each combination with the same rule as previously: the value associated to the ob, result of the combination, is the value obtained by applying the operation (associated to the combination) to the values associated to the obs, operands of the combination (fig. III-3).

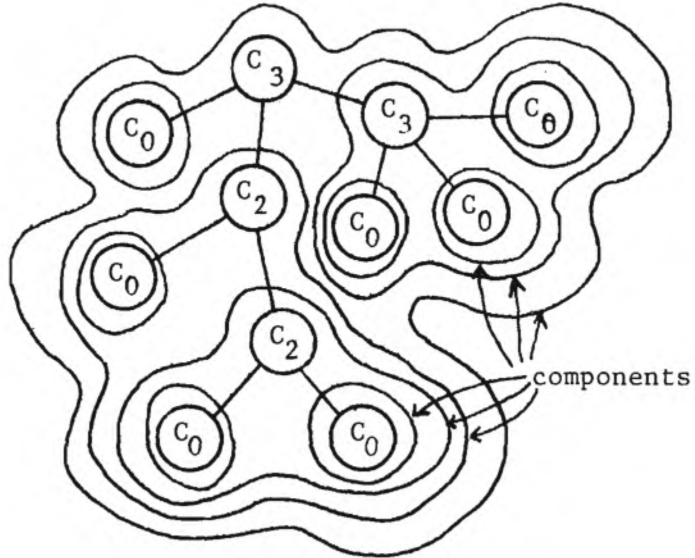


Fig. III-1

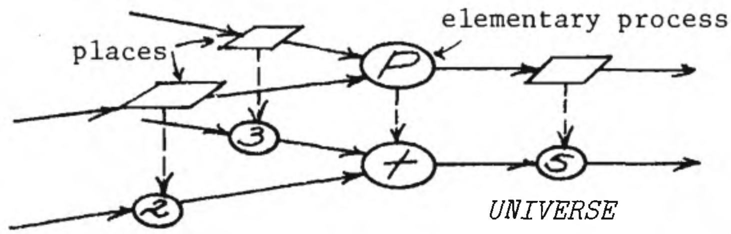


Fig. III-2 (to be viewed in space)

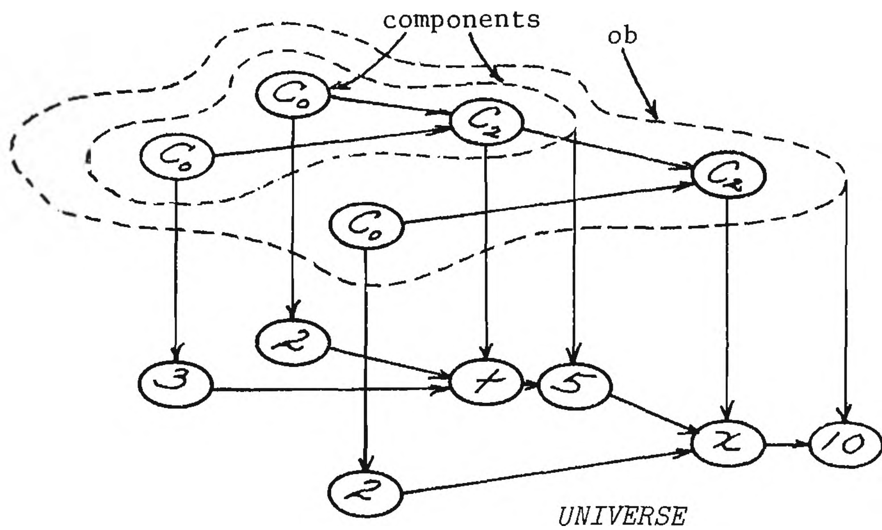


Fig. II-3 (to be viewed in space)

It is useful to abstract slightly such situations by introducing a variation of the concept of categories and functors (see Mitchell [17] ; as defined here, categories may be extended in a trivial way to be mathematical categories).

A <u>category</u> is a set of objects among which some given relations, which we shall suppose to be mappings, hold. So, for instance, the previous universe is a category. Any ob is also a category: the components of the ob are the objects of the category, and the relations are here the mappings actualized by the combinations.

Let C and C' be two categories. A <u>functor</u> is a mapping $T: C \to C'$ which associates:

- to each object O in C, an object T(O) in C'
- to each mapping R in C, a mapping T(R) in C'

such that whenever

$$O_1 , O_2 , \ldots , O_n \xrightarrow{R} O'_1 , O'_2 , \ldots , O'_p \quad \text{in } C$$

we have either

$$T(O_1),T(O_2), \ldots ,T(O_n) \xrightarrow{T(R)} T(O'_1),T(O'_2), \ldots ,T(O'_p)$$

and T is called a covariant functor (fig. III-4a), or

$$T(O'_1),T(O'_2), \ldots ,T(O'_p) \xrightarrow{T(R)} T(O_1),T(O_2), \ldots ,T(O_n)$$

and the function is called a contravariant functor (fig. III-4b).

III.3 Iterative and recursive processes.

The previous definitions are significant because covariant and contravariant functors lead to two different computational schemata. We shall see that they correspond to the intuitive distinction of an iterative versus a recursive computation, when applied to sequential processes.

T is a covariant functor:

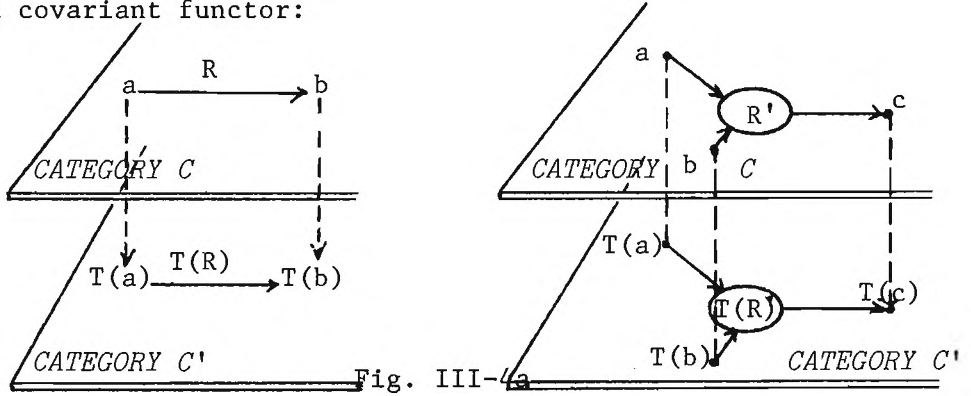


Fig. III-4a

T is a contravariant functor:

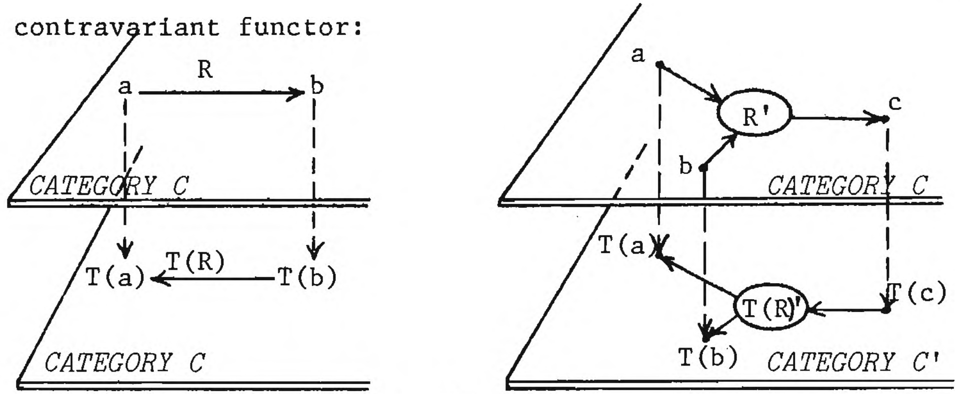


Fig. III-4b

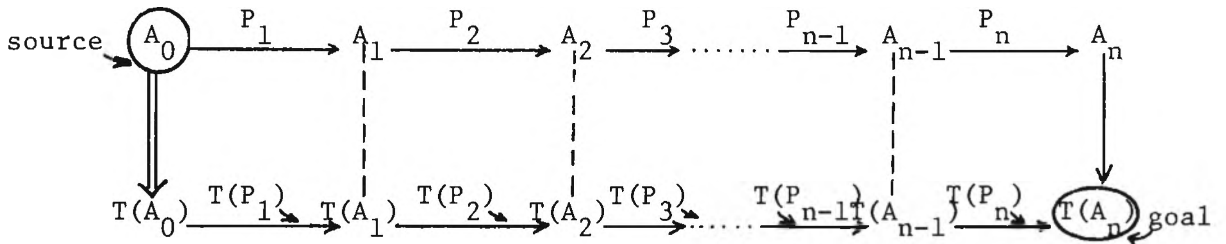


Fig. III-5

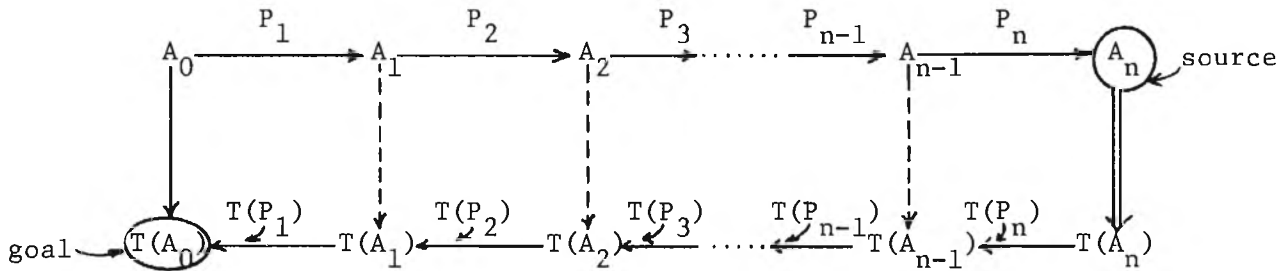


Fig. III-6

To any mapping R is associated an operation T(R) (which may be primitive or not) in the universe of values and operations, and the correspondance $R \longrightarrow T(R)$ is given beforehand. For some objects a_i , called <u>sources</u>, the values $T(a_i)$ are given in the universe of values. For some b_j the values $T(b_j)$,called <u>goals</u>, are to be determined.

Let us consider a sequence of objects, each object (the first one excepted) being obtainable by applying a particular elementary process to the previous object in the sequence. To each elementary process P_i is associated an operation $T(P_i)$, T being a covariant functor. The first object A_0 is the source; the value associated to the last object A_n is the goal (fig. III-5).

The computation of $T(A_n)$ may be performed straightforwardly with one storage place whose contents will be called the current value. At the beginning $T(A_o)$ is the current value; the control then proceeds sequentially from the source to the last object of the sequence; whenever a mapping P_i is encountered, the operation $T(P_i)$ is applied to the current value, and the result is taken as the new current value. When the control reaches A_n , the current value is the goal $T(A_n)$.

In pseudo-Algol, the process may be described by means of an iteration:

```
... currentvalue := T(Ao) ;
    for i := 1 step 1 until n do
       currentvalue :=T(Pi)(currentvalue) ;
    .........
```

We consider now the same sequence of objects, T being a contravariant functor. A_n is the source and $T(A_o)$ the goal (fig. III-6).

It may seem that $T(A_n)$ being given, the computation of $T(A_o)$ is completely similar to our previous computation $T(A_0)$ and $T(A_n)$ being permuted: $T(P_n)$, $T(P_{n-1})$, ... , $T(P_1)$ would be applied sequentially to $T(A_n)$. However, in general, this reverse sequence of mappings is not known beforehand.

The control goes as before from A_o to A_n : whenever a mapping P_i is encountered, the corresponding operation $T(P_i)$ is placed at the top of a last-in-first-out queue (a stack). After the control has reached A_n , the computation proceeds as in our previous example, the operations being retrieved, now, from the stack.

The arrows on the diagram may serve as a built-in stack. Let us consider as an example the factorial function:

(1) fact(n) = [<u>if</u> n = 0 <u>then</u> 1 <u>else</u> n × fact(n-1)]

Let us compute fact(3). The goal is fact(3), the source is 0 with fact(0) being 1. The diagram is displayed in fig. III-7.

We can interpret the program as a diagram modification scheme. The arrows $T(P_i)$ are constructed as the control proceeds along the arrows P_i. When the control reaches the source the two sequences are bound. Then the control proceeds towards the goal, performing the operations $T(P_i)$. (Fig. III-8).

We may wonder what interpretation an iterative factorial may have.

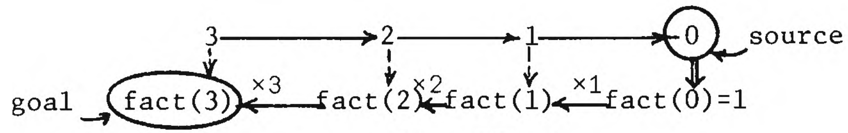


Fig. III-7

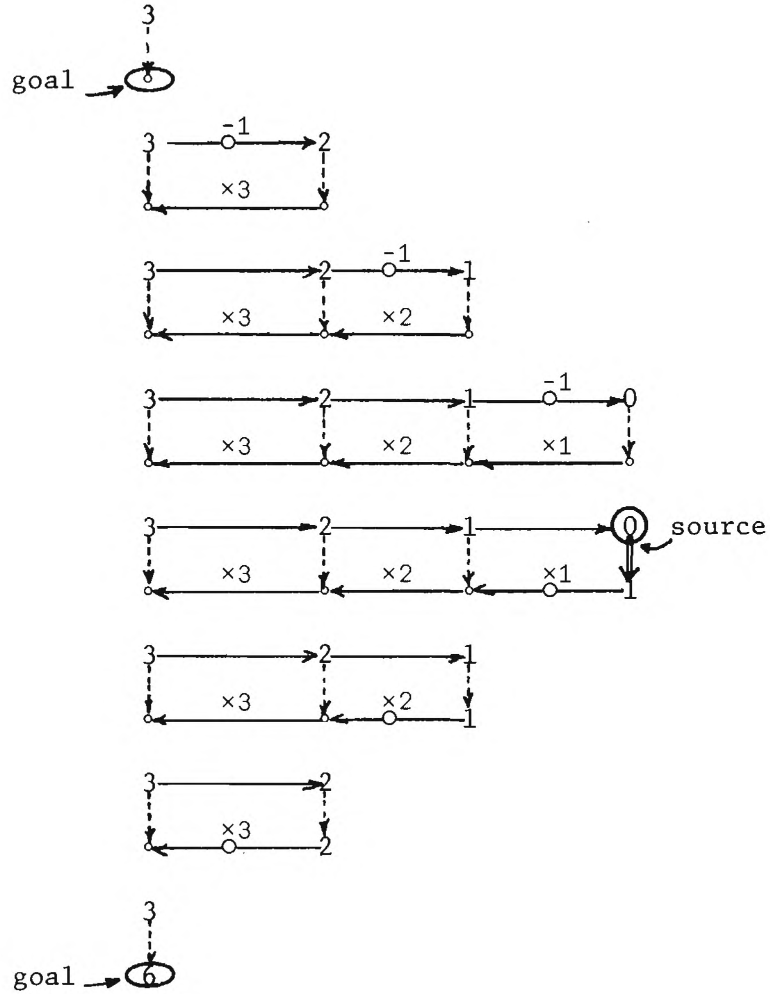


Fig. III-8

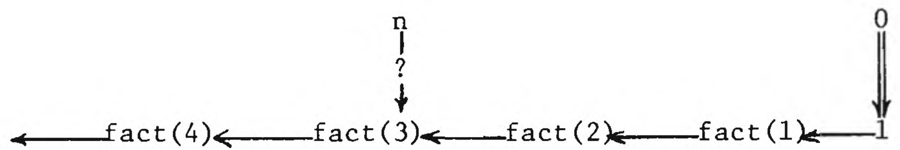


Fig. III-9

A "while-iterative" factorial: (2) [x := 1 ; y := 1 ;

```
while x ≤ n do begin y:= x × y ;
                     x:= x + 1 end ;
y ]
```

A "for-iterative" factorial: (3) [y:= 1;

```
for i:=1 step 1 until n do y:= i×y ;
y ]
```

In these programs, the programmer knows beforehand what the sequence $T(P_n)$, $T(P_{n-1})$, ... is. As a result fact(0) being known, fact(1), fact(2), ... may be computed iteratively, one after the other (fig. III-9). In the first program, at each step, a test determines whether the goal is reached or not. In the second program, the programmer knows beforehand that the goal will be reached whenever the control has passed three arrows.

As a result we can consider algorithm (1) as the most general. Algorithm (2) uses a particular situation. Algorithm (3) uses a still more particular situation.

In the previous discussion elementary processes were restricted to transform one object into another one. The general case where elementary processes are applied on several input objects and produce several output objects may receive a rather similar treatment.

Objects and mappings relating one object to another constitute a category (fig. III-10). Let T be a functor, associating to each elementary process P_i an operation in the universe $T(P_i)$. If T is covariant (fig. III-11) we can take the input objects a,i,j as sources, the goals being the values of the output objects c and m. The computation is performed with a distributed control. If T is

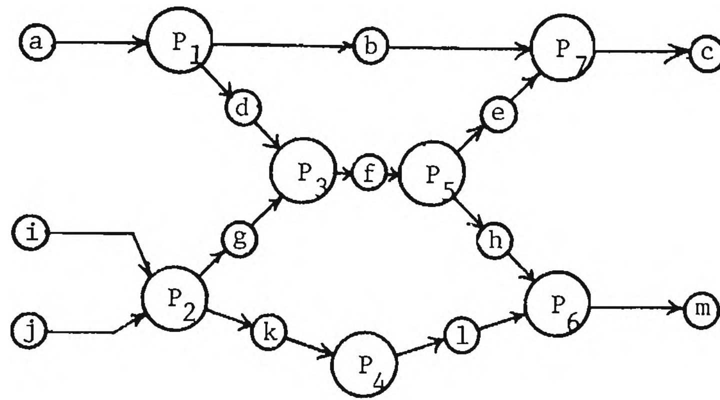
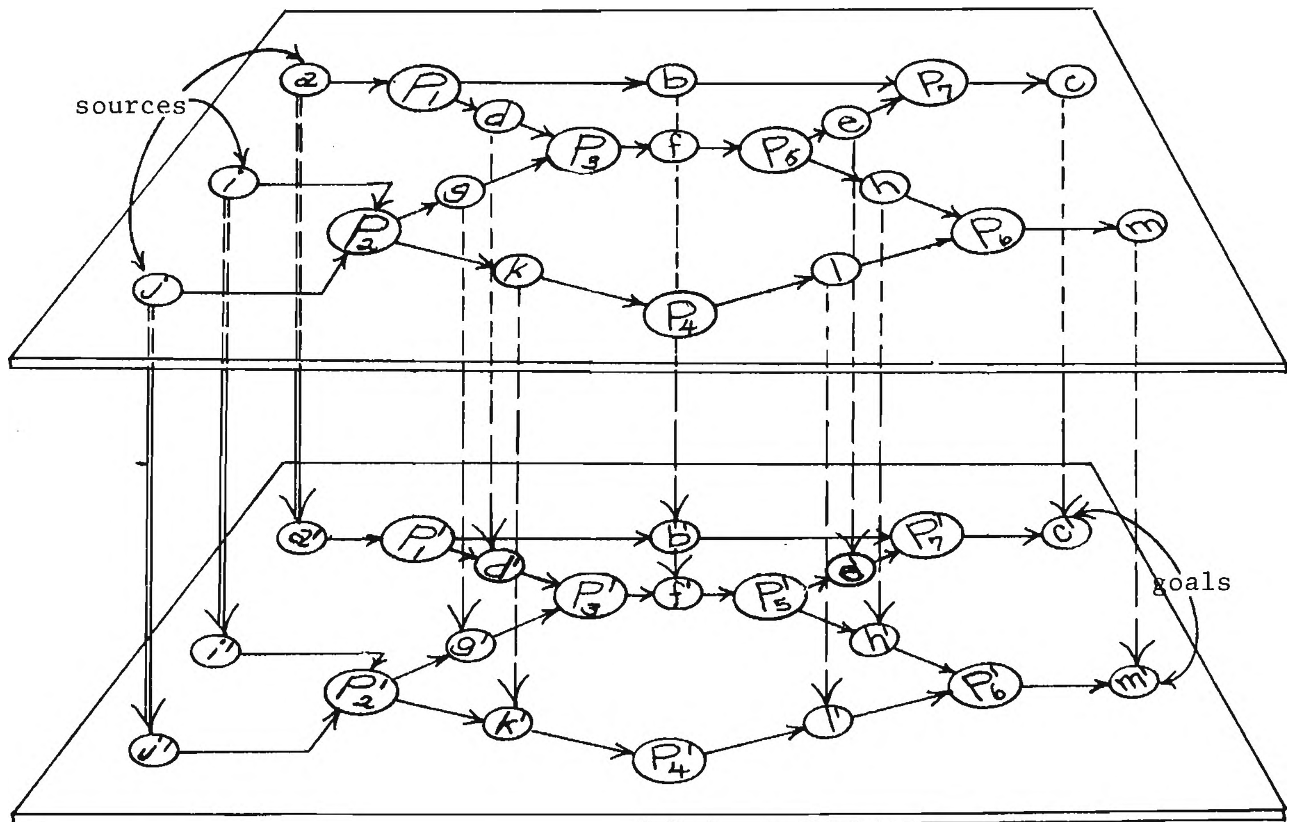


Fig. III-10



T is a covariant functor (in this figure X' denotes T(X) for any symbol X)

Fig. III-11

contravariant (fig. III-12) the output objects are taken as sources, the goals being the values of the input objects T(a), T(i), T(j). The arrows representing the mappings $T(P_i)$ are used as a first-in-last-out-like storage device. We may notice that a stack,being an essentially sequential storage device, cannot be used here.

III.4 Synthetic and inherited attributes.

III.4.1 Definitions.

When applied to an ob, covariant and contravariant functors lead to the notions of synthetic and inherited attributes. Indeed, it is sometimes interesting to associate to each component of an ob an attribute, these attributes being structurally related to one another. In our terminology, the components of an ob are considered to constitute a category, which is mapped by some functor into the universe of values and operations. If the functor is covariant, the attribute of the given ob may be computed whenever the attributes of the atoms (primitive obs) are known. Such attributes are called synthetic attributes. If the functor is contravariant, the attribute of an atom may be computed whenever the attribute of the given ob is known. Such attributes are called inherited attributes (see Knuth [11] where a syntactic point of view has been taken).

III.4.2 The value of an expression as a synthetic attribute.

The expression (2 +3 +2) + 2 + 1 (see section I.2.1) may be regarded as an ob. Its evaluation requires only one synthetic attribute, the value.

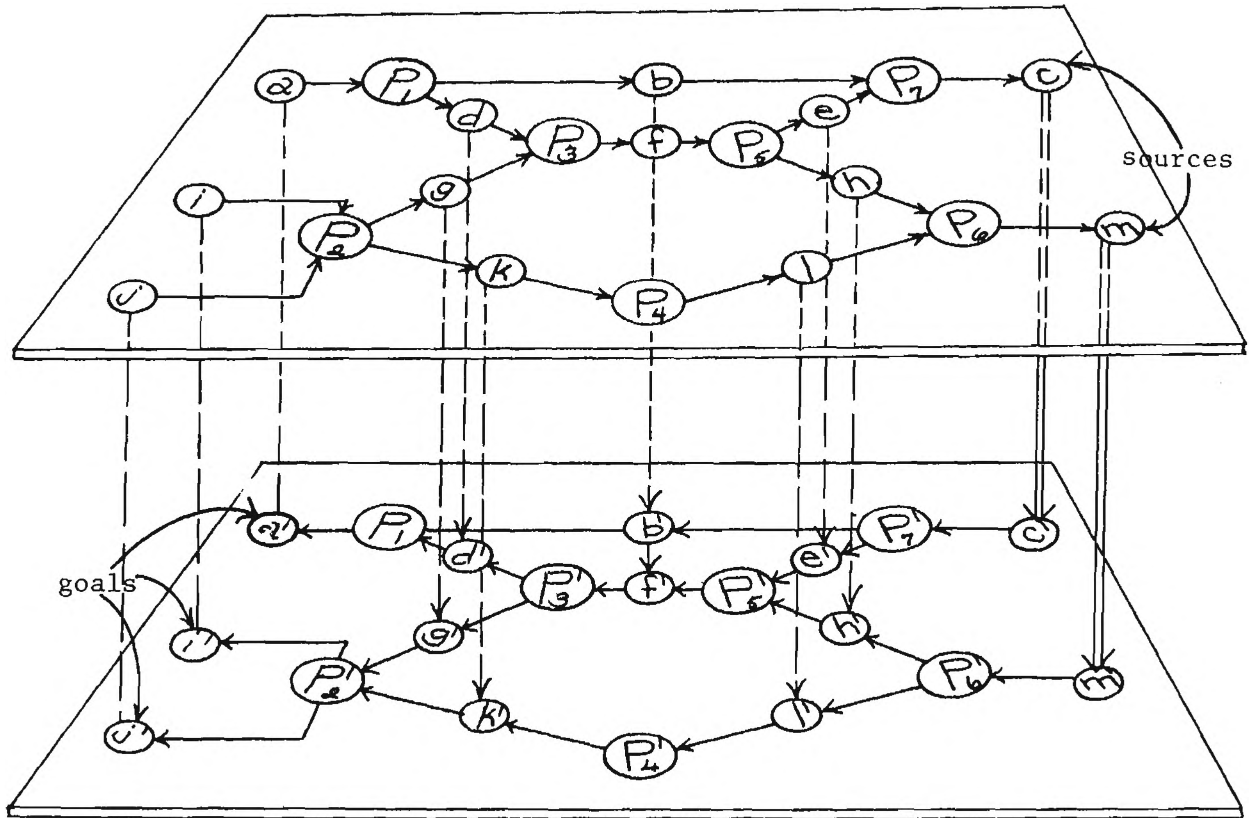


Fig. III-12

Whenever an ob like $(x + y + x) + x + 1$, containing formal variables, is to be evaluated, it is necessary that the environment determines the value of each variable. The realization of the ob is a covariant functor (fig. III-13).

III.4.3 The environment as an inherited attribute.

In the previous example the environment was considered as being global relatively to the expression, and as being accessible by any variable. We may also consider the environment as an inherited attribute.[a] The evaluation may then be viewed as having two phases: the binding and the evaluation proper (fig. III-14).

In the example of section I.2.1,

$(2 + 3 \to x \; ; \; 5 \to y \; ; \; x + y \to x \; ; \; y + 1 \to y \; ; \; 2 \times y - x)$,

the $\to x$ and $\to y$ may be considered as being operators on environments (fig. III-15).

---

a) To each node is associated an attribute: a local environment. As a result, to two occurences of a same variable in a computation may correspond two different local environments which may associate to these variable-nodes two different values.

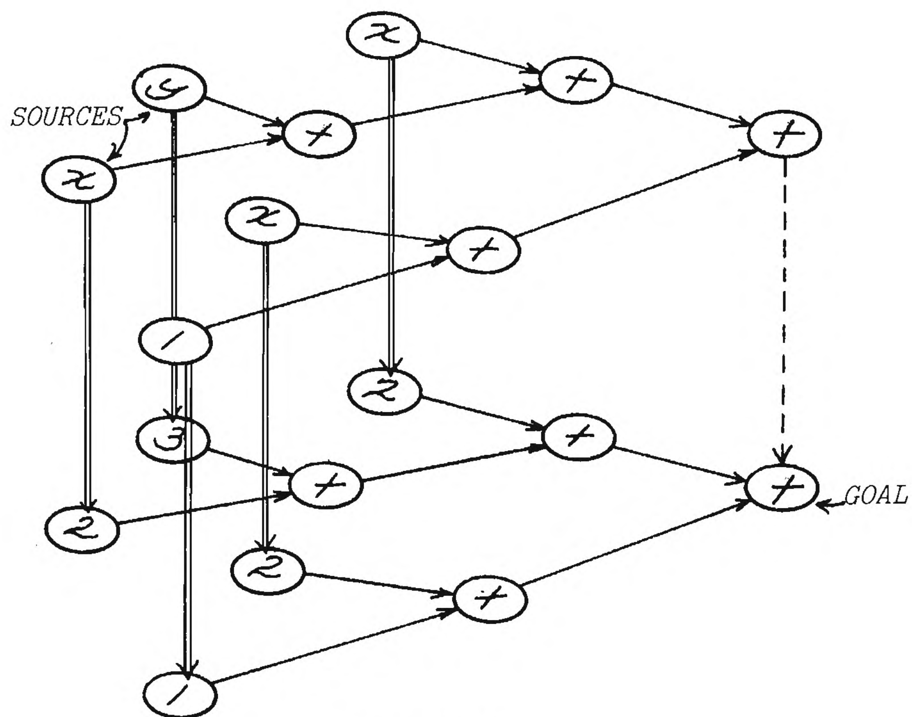


Fig. III-13

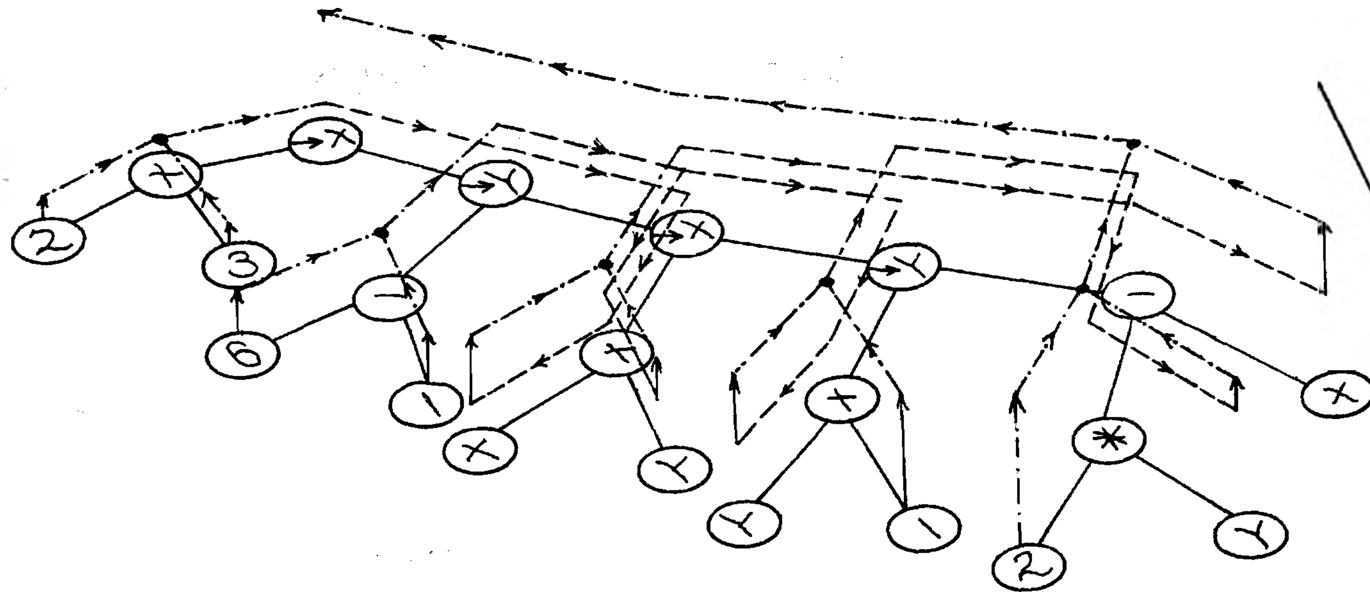
binding

evaluation

proper

goal

Fig. III-14



—·—·—·— value (synthetic attribute)

— — — — — piece of environment (inherited attribute)

Fig. III-15

CHAPTER IV

SOME COMPUTATION MODELS

IV.1 Introduction

In this chapter, two different approaches of modelling computations are discussed. In both of these approaches, the models present in some way concurrency and distribution of control.

In the first approach, the directed graph is taken as a model of computation [6,10,10a,16,21,23] . In section IV.2 we will discuss the "computation model with data flow sequencing" due to Adams [1]. Such a representation is of interest for us since a DCPL program may appear to be a data flow model after the binding has been performed.

In the second approach, a functional computation is considered as an expression; the evaluation of such an expression may be carried out by some abstract machine. The lambda-calculus provides a machine whose elementary operations are replacement and substitution (section IV.3).

Curry's combinators allow to use a much simpler machine, using replacement only. Strikingly enough, variables are not used at all in this representation. As a result of the simplicity of the machine and of the extreme locality of control, the representation of an expression may appear rather complex (section IV.4).

The second part of this chapter (sect. IV.3 and IV.4)

contains some technicalities. Some readers might prefer to skip it.

IV.2 The Adams' computation model with data flow sequencing.[a)]

The directed graph is used as a model of the computation.

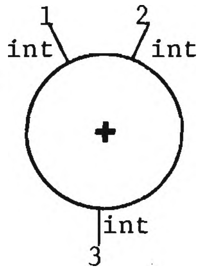
"The nodes of the graph represent computation steps and the edges represent transmission paths for data and control. An edge may be thought of as a queue of data produced by one node and waiting to be consumed by another. A computation step may be initiated whenever each edge directed into that node of the graph contains the amount of data required for the node to execute properly."

Fig. IV-1 displays some nodes which are activated whenever there is an input value on each input edge. Then, output values to be put on the output edges are computed according to a function f associated with the node. "φ" is a notational device: placing "φ" on an edge means placing nothing on this edge.

Fig. IV-2 displays some nodes having a more involved behavior. Relatively to a node, an input edge may be in any one of two status: it is either locked or unlocked. In fig. IV-2, the status of input edges is given, for instance, as ULL , which means that port 1 is unlocked and that ports 2 and 3 are locked. The computation step of a node is initiated whenever there is an input value on each unlocked input edge. Output values are computed as before according to an associated function f; moreover, a new status for each input edge is determined according to another function g associated with the node. The blocking capabilities of these nodes permit the computations to be determinate.

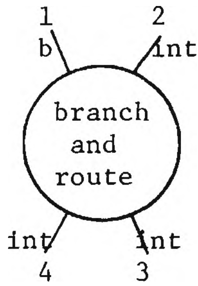
---

a) We have taken the freedom to quote some passages of, and to reproduce some figures from Adams [1].



f: $V_3 \leftarrow V_1 + V_2$;

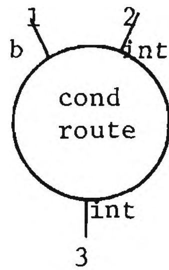
Fig. IV-1a



f: $V_3 \leftarrow$ if V_1 = false then V_2 else φ;

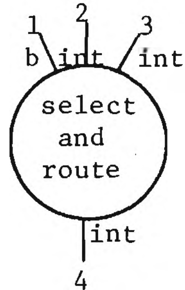
$V_4 \leftarrow$ if V_1 = true then V_2 else φ;

Fig. IV-1b



f: $V_3 \leftarrow$ if V_1 = true then V_2 else φ;

Fig. IV-1c



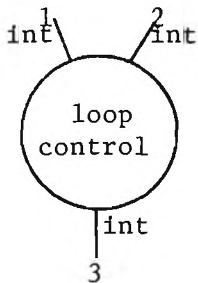
g: ULL → LUL if V_1 = true f: $V_4 \leftarrow \varphi$;

ULL → LLU if V_1 = false $V_4 \leftarrow \varphi$;

LUL → ULL $V_4 \leftarrow V_2$;

LLU → ULL $V_4 \leftarrow V_3$;

Fig. IV-2a



g: UL → LU f: $V_3 \leftarrow V_1$;

LU → LU $V_3 \leftarrow V_2$;

Fig. IV-2b (from Adams [1])

Fig. IV-3 and IV-4 represent an iterative [1, p. 31] and a recursive [1,p.32] factorial respectively. The latter figure witnesses to the recursive character of a graph procedure:

"When a node in a graph procedure represents a recursive call upon the procedure of which it is a part, a copy of the called graph procedure is created. Thus, during the execution of a graph program, an auxiliary graph referred to as the executing graph will be constructed. Initially the executing graph will consist of the main graph procedure G, with the initial data for the program placed on the edges of G. The initial data must be of the same type as the edges on which it is placed. Whenever a procedure node in the executing graph is ready for execution, a copy of the defining graph procedure will be created and added to the executing graph [a)]; and when the procedure terminates, the created copy will be deleted. The executing graph can thus expand and contract dynamically during the execution of the program."

Remark. Such a situation may be regarded as an instance of the realization of a contravariant functor (compare III-8 and IV-5).

In DCPL we will find the same notion of implementing in space successive generations of a recursive procedure.

IV.3 Functional representation: the lambda-calculus [b)]

The discussion in this section and in the next one is based on Curry [3]. The lambda-calculus provides a framework in which functional expressions may be represented and evaluated.

---

a) Such a creation and addition of a graph procedure to the executing graph is referred to in this thesis as an implementation in space.

b) There are different lambda-calculi, each one having its particular idea about what objects represent the same function. However since in programming languages the emphasis is put to the application of a function to an argument, and not to a function itself, only one calculus is generally used (β-conversion lambda-calculus).

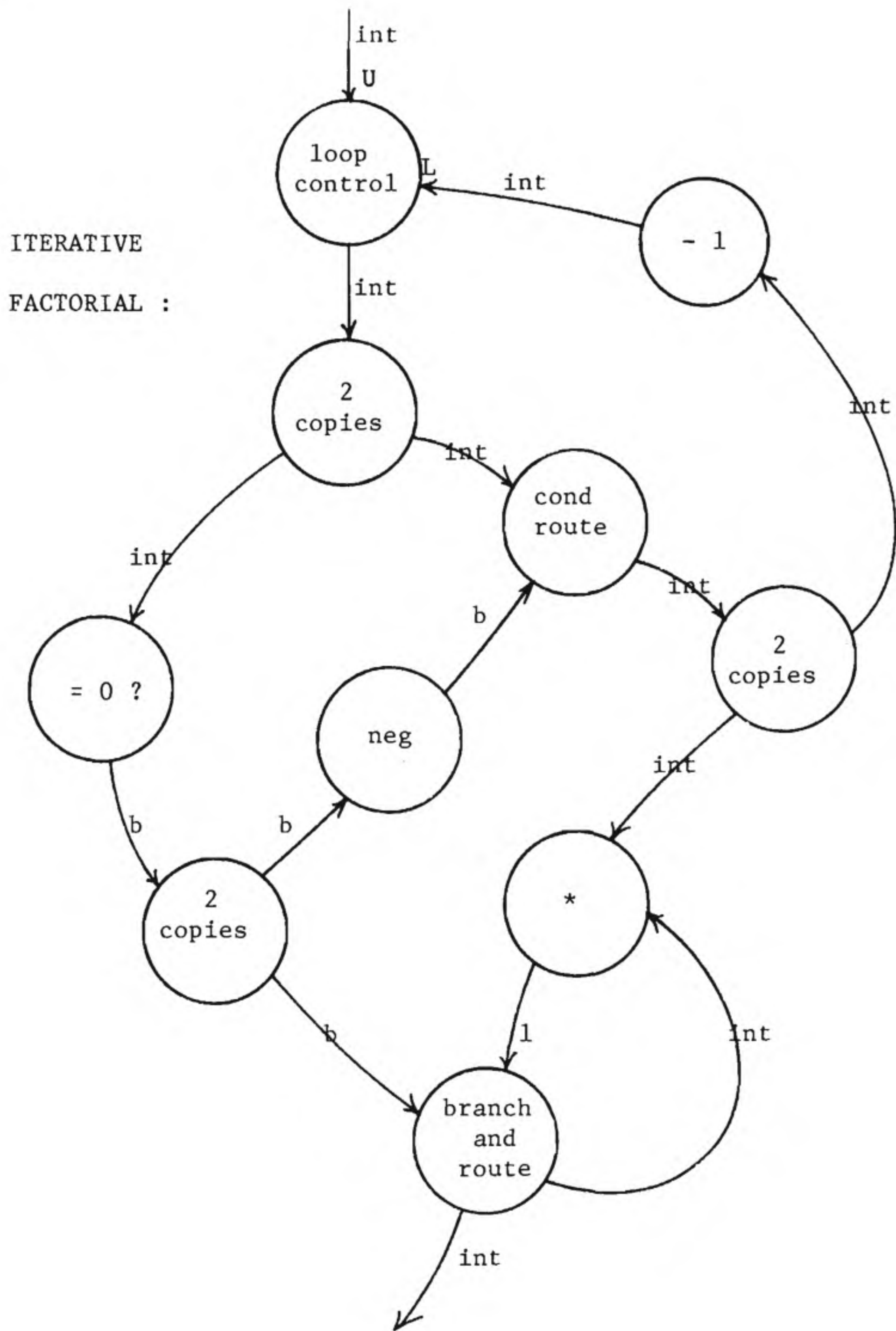


Fig. IV-3 (Adams [1],p.31)

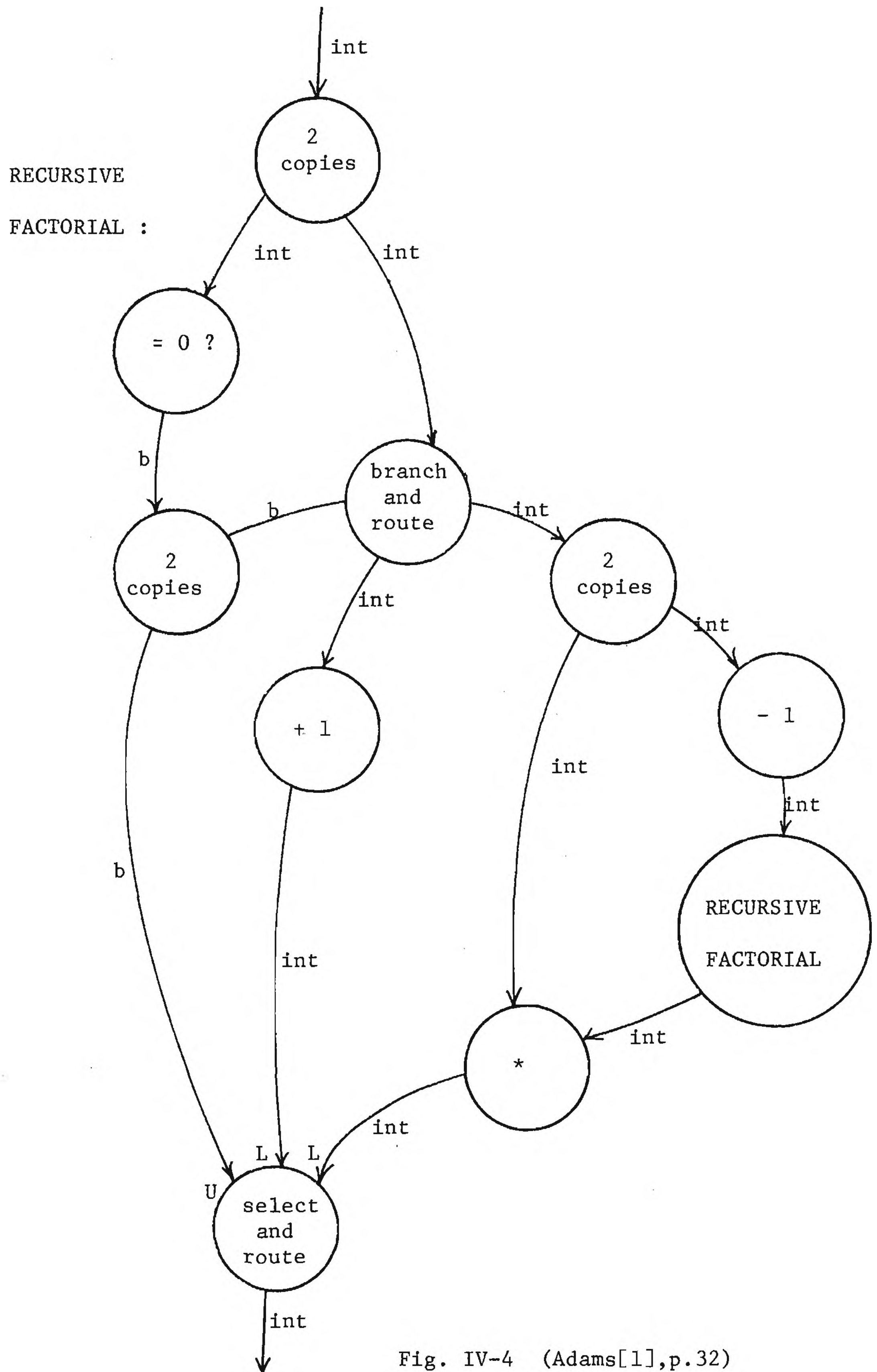


Fig. IV-4 (Adams[1],p.32)

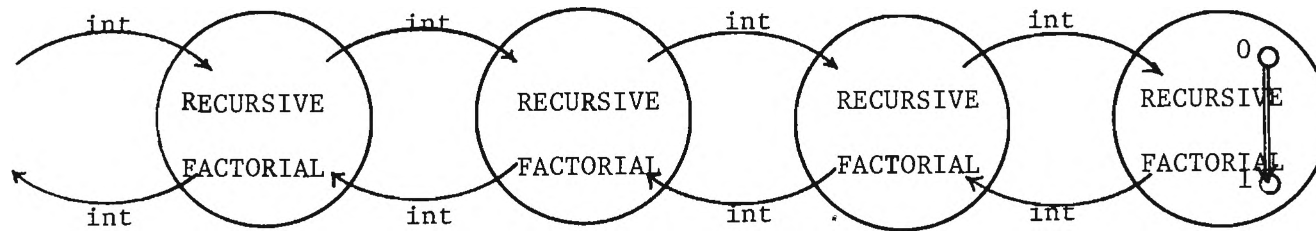


Fig. IV-5

Such evaluations present possibilities of concurrency and distribution of control. Since Mc Carthy [18] , the lambda-calculus has often been considered in connection with programming languages: Landin [12,13,14], A. Evans [7], Morris [19].

IV.3.1 Variables in mathematics.

Let us consider with Curry the following mathematical statements:

(1) $$(x + 1)^2 = x^2 + 2\,x + 1$$

(2) $$\frac{d}{dx}\,x^2 = 2\,x$$

(3) $$\int_o^3 x^2\,dx = 9$$

In statement (1) the variable x may be considered as having an intuitive meaning: for any integer, for instance a, we have the relation $(a + 1)^2 = a^2 + 2a + 1$.

We cannot interpret the use of the variable x in such a simple way in (2) and (3): they do not enunciate any statement about some object for which x stands. In fact, (2) and (3) state some properties about a function, the square function. The use of variables in these statements may be considered as only a notational device. The use of variables may even be more explicit with Church's Lambda notation (Church [2]); $\lambda x.x^2$ denotes then the square function. If D and I_a^b denote respectively derivation and integration between a and b, the statements (2) and(3) become:

(2') $$D(\lambda\,x.x^2) = \lambda x.2x$$

(3') $$I_o^3(\lambda\,x.x^2) = 9$$

The lambda notation allows us to consider a function, at least conceptually, as an object: functions become part of the universe of discourse, and statements about functions may be formulated.

IV.3.2 Functional representations in programming languages.

Functions appear naturally in programming languages whenever some object is to be evaluated in some realization (see chapter III). Generally statements about functions which are considered in programming languages are very limited; with the exception, may be, of some symbolic manipulation oriented programming languages, functions are considered as far as they will be applied to some arguments at some time. There is a notation to represent functions and another notation for the application of a function to some arguments, and there is a background machine which may evaluate functional expressions. If such a background machine is to be an abstract (or formal) one, a notation is not sufficient, it is necessary to be able to represent a functional expression. By representation we mean:

1. Any function may be represented as an object, better as an ob.
2. There is a combination called application which allows to represent in an explicit way the ob obtained by combining,with the application, a function to its arguments.
3. There is an abstract machine which can evaluate the previous ob and produce the representation of the resulting object.

IV.3.3 An abstract reducing machine.

We are interested here in abstract machines performing reductions since these machines are particularly simple: when applied to some object, they look for any component correspon- ding to a given pattern and replace it by an associated com- ponent . A macro-processor, an interpreter using a simple prece- dence grammar (Wirth [28]) might be considered as reducing machines.

IV.3.3.1 Substitution versus replacement.

We present these notions pictorially: in fig. IV-6 an ob X is substituted for an atom x in an ob A; the result is an ob B.(We may notice that

1. x is supposed to be an atom
2. each occurrence of x is replaced by an occurrence of X.).

In fig. IV-7, a component X of an ob A is replaced by an ob Y; the result is an ob B. (Note:

1. X is not necessarily an atom, but an ob
2. Even if there are several occurrences of X in A, only the considered occurrence is replaced.).

IV.3.3.2 Reduction rules.

Let us consider two obs A and B and the reduction rule $A \twoheadrightarrow B$. Let X be an ob having A as a component. We say that X may be reduced in Y, and we note $X \gtrdot Y$ if Y is obtained by replacing, in X, the component A by B. We note $X \doteq Y$ whenever $X \gtrdot Y$ or $Y \gtrdot X$ and $\geq$ and $=$ represent respectively the quasi-ordering (symmetric and transitive) and the equivalence generated by $\gtrdot$ and $\doteq$.

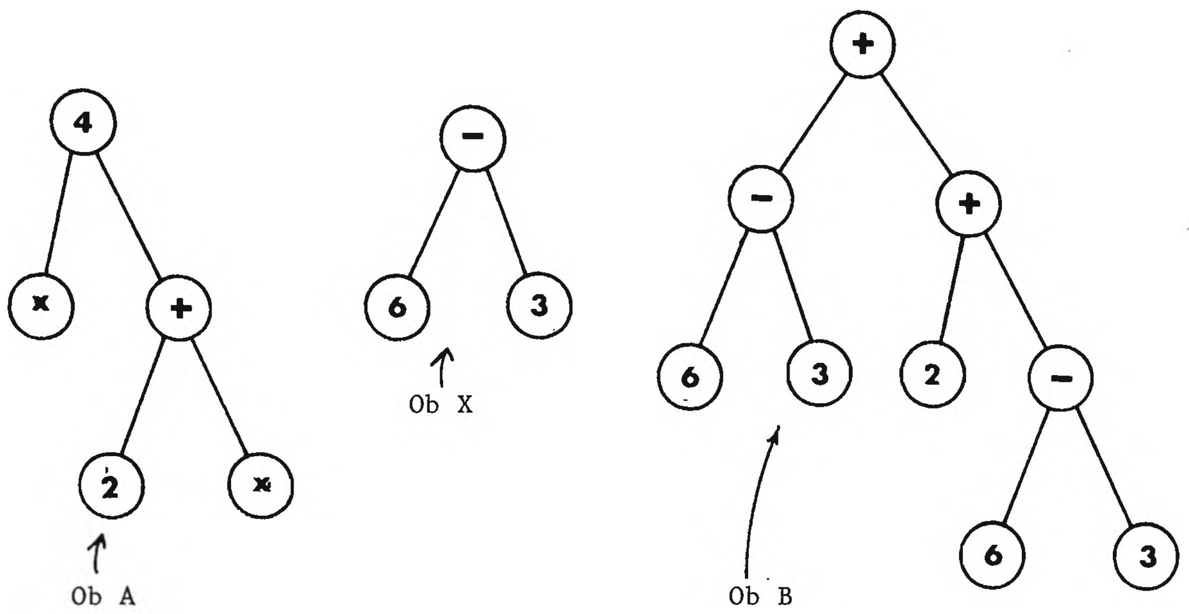


Fig. IV-6

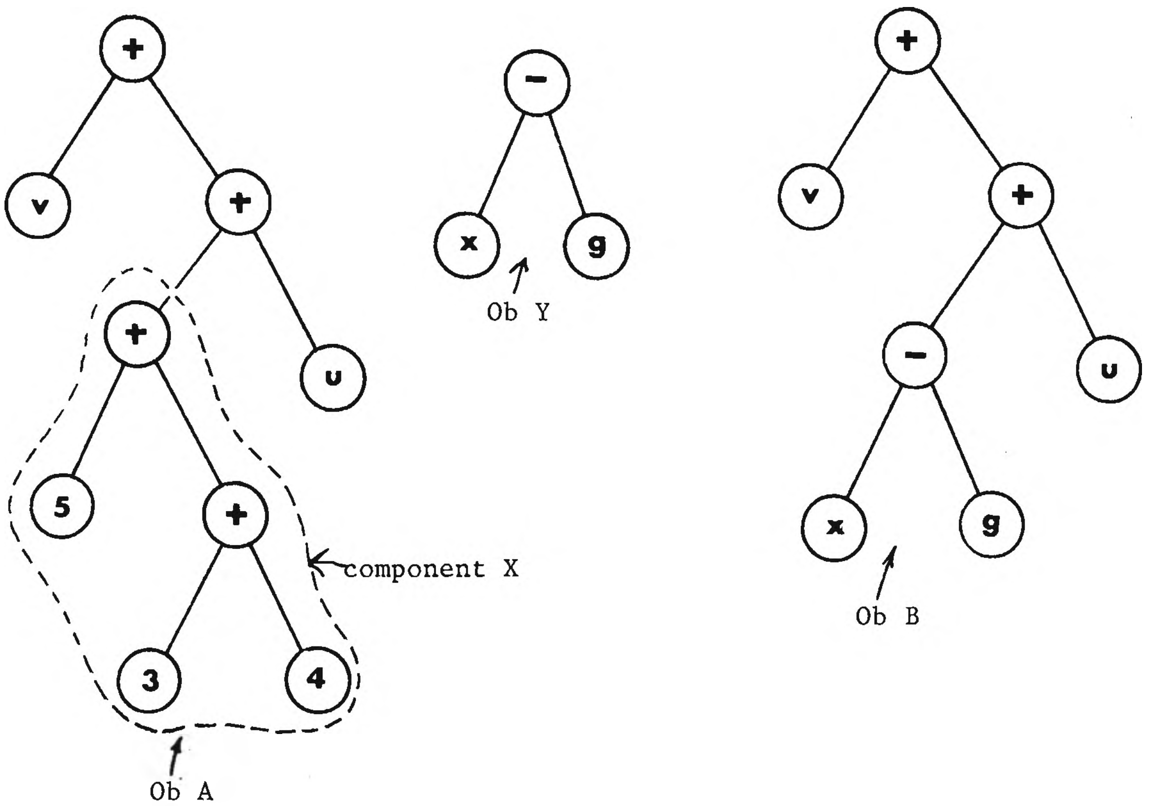


Fig. IV-7

The same definitions may be given with a set of reduction rules $\left\{ A_i \rightarrow B_i \right\}_{i \in I}$. In summary:

$X \cdot\!> Y \iff$ there is in the ob X a component A_i whose replacement by the corresponding B_i produces the ob Y

$X \doteq Y \iff$ either $X \cdot\!> Y$ or $Y \cdot\!> X$.

$X \geq Y \iff$ there is a sequence of objects $X_o, X_1, \ldots X_n$, $X_o \equiv X$ and $X_n \equiv Y$ such that $X_o \cdot\!> X_1 \ldots \cdot\!> X_n$ ($\equiv$ denotes the identity of obs)

$X = Y \iff$ there is a sequence of objects $X_o, X_1, \ldots X_n$ such that $X_o \equiv X$ and $X_n \equiv Y$ and $X_o \doteq X_1 \doteq \ldots \doteq X_n$

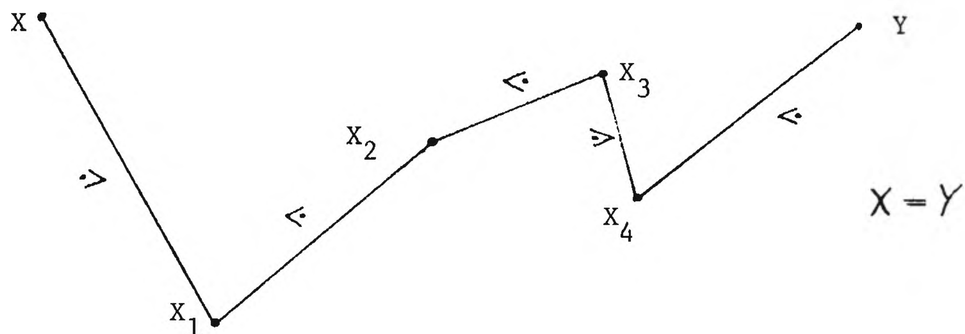


Fig. IV-8

IV.3.3.3 Reducing machine.

Whenever a set of reduction rules is given we can consider the following reduction process, to be applied to any ob X:

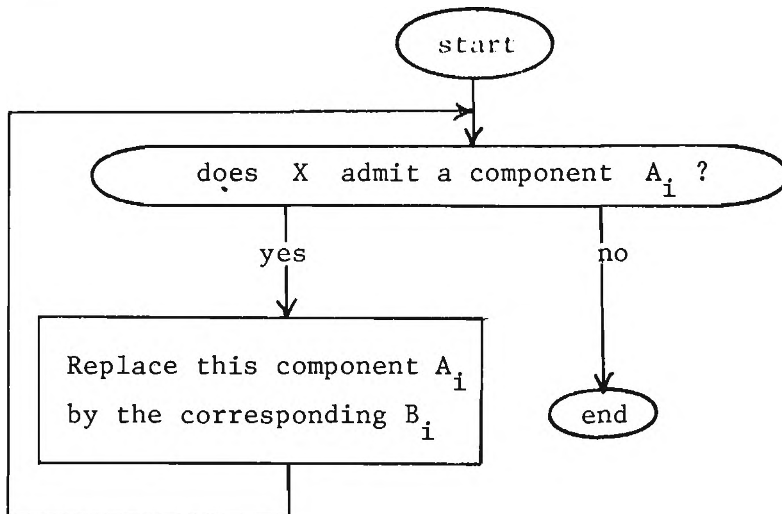


Fig. IV-9

Obviously, such a process is not in general deterministic. Moreover, applied to an ob X it may or may not stop. If it does stop, producing an ob X_o , there is no reduction which may be performed on X_o: X_o is said to be a <u>normal</u> <u>form</u> of X.

We are interested in abstract reducing machines which are deterministic to some extent, namely in machines verifying the following condition (Church-Rosser): if $X = Y$ there is an ob Z such that $X \geq Z$ and $Y \geq Z$.

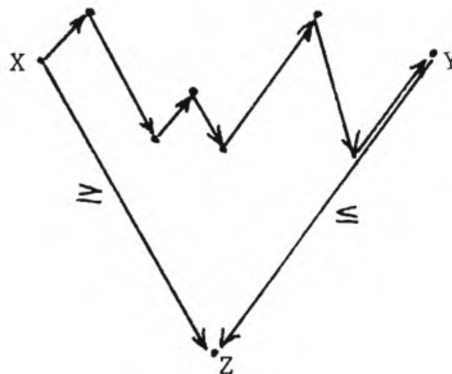


Fig. IV-10

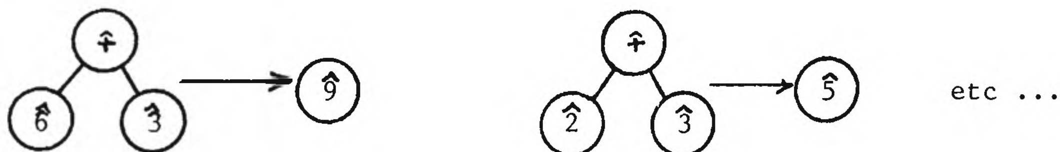
Indeed such a condition guarantees that whenever the machine stops when applied to an ob X, it produces the same normal form (however in some simulation the machine may not stop). Such a normal form may be considered as the "value" of X, the reduction machine being therefore an evaluating machine.

IV.3.3.4 A trivial example.

Let us consider a trivial representation of simple computations without variables:

- to each value v in the universe corresponds a combinator of degree zero $\hat{v}$ and to each operator of degree n, opr, a combinator of degree n, $\hat{opr}$. We call CL_o the class of obs which may be constructed with these combinators. For any operator opr and for any values v_1 , v_2 , ... v_n producing $opr(v_1,v_2, \ldots ,v_n)$ we have the reduction rule :

For instance if there is one operation, the addition, and integers as values, the reduction rules will be of the form:



Any sum will be trivially evaluated by our machine (fig. IV-11). The Church-Rosser condition is here obviously verified; the machine

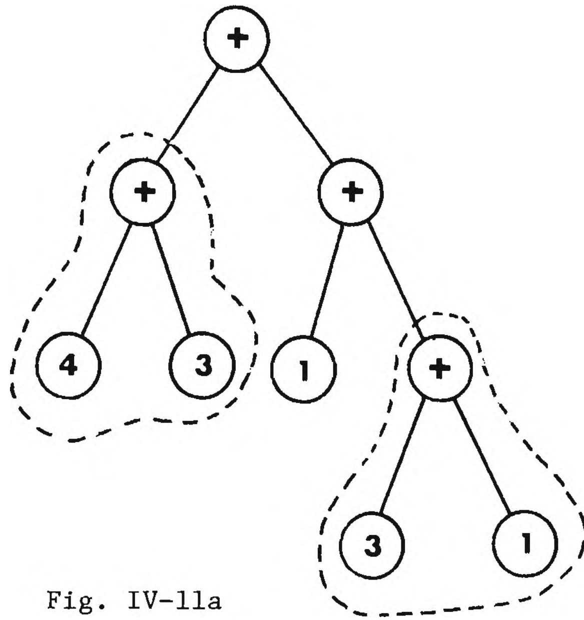


Fig. IV-11a

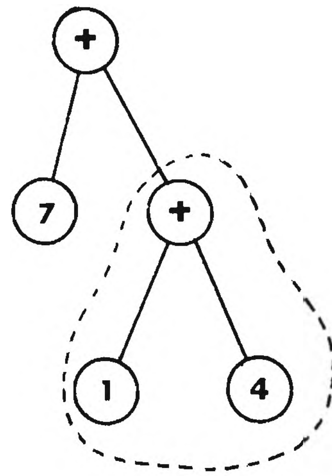


Fig. IV-11b

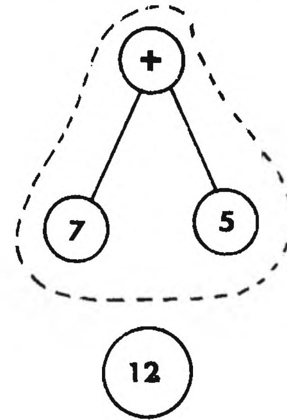


Fig. IV-11c

Fig. IV-11d

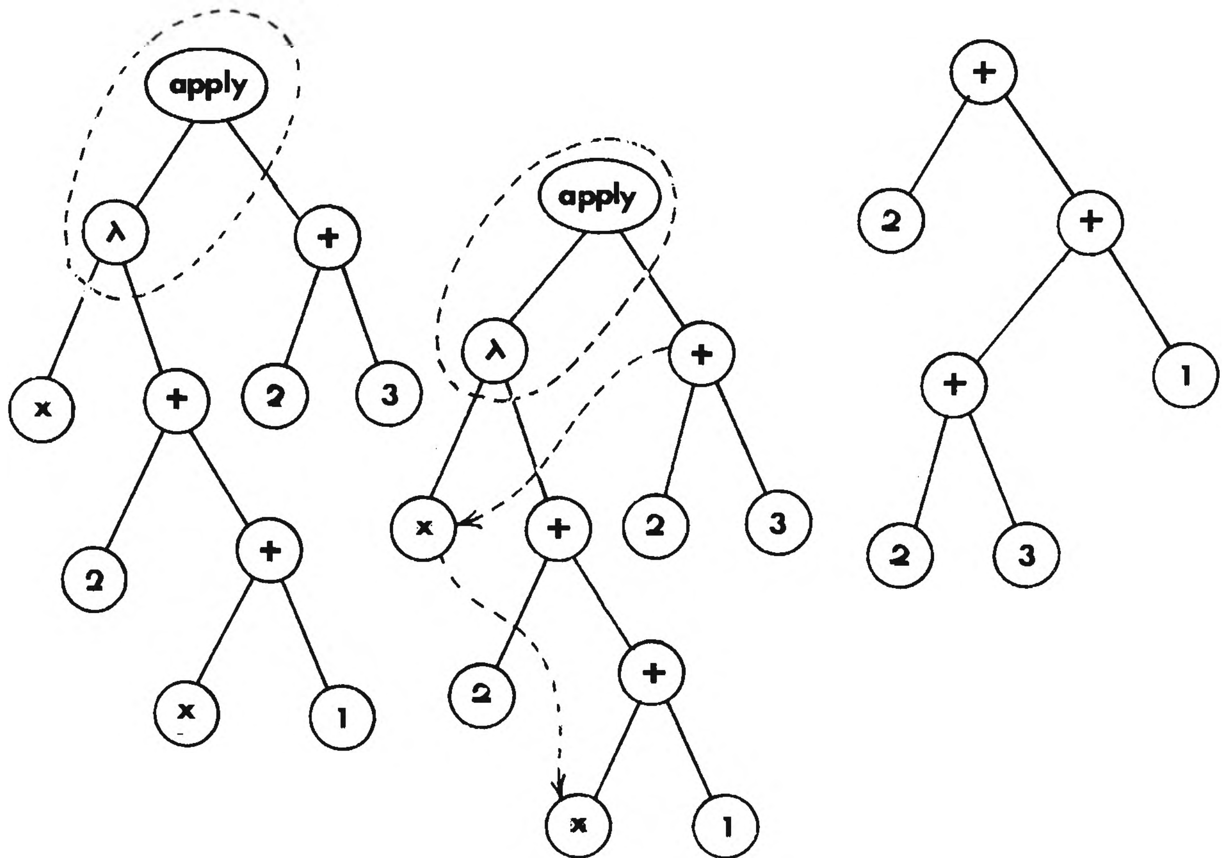


Fig. IV-13

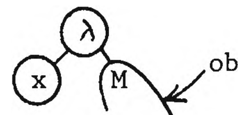
is deterministic and has a distributed control. The result $\widehat{12}$, as a combinator, is the normal form of the given expression and may be considered as being the value of the expression. The relation of equality we have defined corresponds to the traditional meaning of equality of sums.

IV.3.4 Evaluation of λ -expressions.

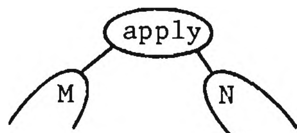
Let us consider, together with the previous combinators, a new atom called a formal variable x. Any ob M containing x may be considered as a function $CL_o \longrightarrow CL_o$ (CL_o is the class of obs obtained with the combinators associated to values and operations of the universe): indeed, to any ob A in CL_o we may associate the ob obtained by substituting A for x in M, which may be denoted by $[A/x]M$ and the function $A \longrightarrow [A/x]M$ may be denoted by $\lambda x.M$ ([2]). The operation $M \longrightarrow \lambda x.M$ is called abstraction relatively to the formal variable x.

We can <u>represent</u> such a function as an ob in the following way:

1. There is a combination of degree 2, called <u>abstraction</u> and denoted by a combinator λ. Whenever abstraction is applied to two obs, the first one must be a formal variable. A written notation for such a combination is $\lambda x.M$; the ob is



2. There is a combination of degree 2 called <u>application</u>, and denoted by the combinator '<u>apply</u>'. A written notation for



is MN .

3. There is an abstract reducing machine defined by the reduction rules generated by the following reduction scheme: Whenever M and N are two obs and x is a formal variable, $(\lambda x.M)N \rightarrow [N/x]M$, or graphically:

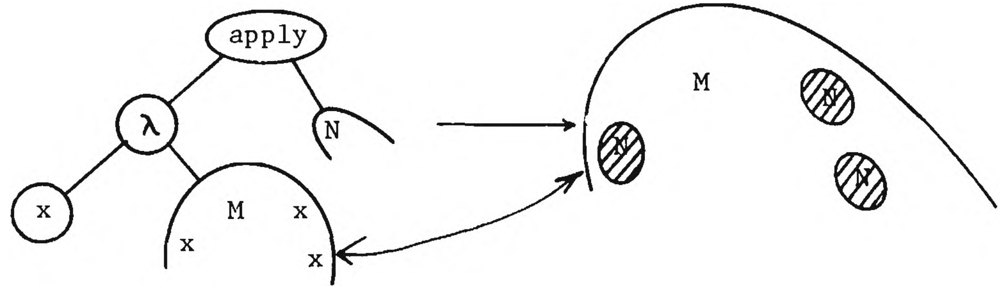


Fig. IV-12

It is proven that these reductions verify the Church-Rosser condition (Church-Rosser theorem, see Curry's [3,chapter 4])

We now have a machine which can evaluate expressions such as:

$(\lambda x.(2 + (x + 1)))(2 + 3)$ (fig. IV-13)

or: $(\lambda x.x + 1)((\lambda x.2 + x)\ 3)$ (fig. IV-14).

In fact, very general computations may be expressed with such a model (see Landin [12]).

Remark 1: The reduction process may not stop when applied to some peculiar obs like: $Y \equiv \lambda f.[(\lambda h.f(hh))(\lambda h.f(hh))]$ called "paradoxical combinator" (see Morris [19]).

Remark 2. The reduction machine is not in a strict sense deterministic: we can only state that whenever the machine stops, it produces the same value; however some simulation may produce a value

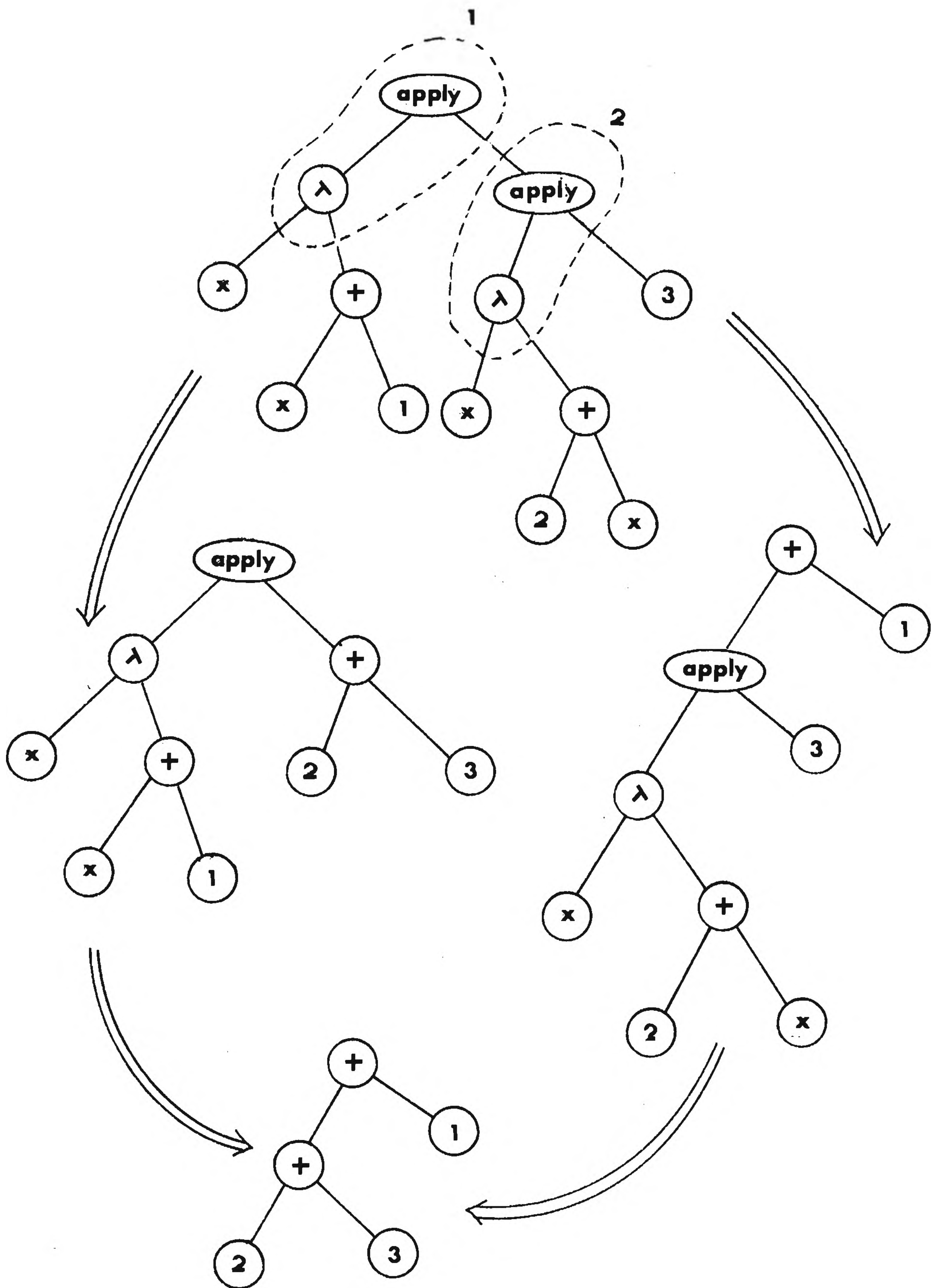
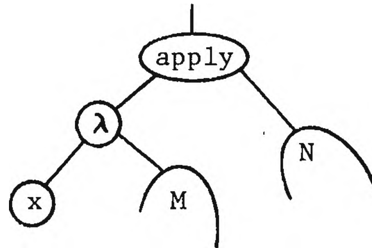


Fig. IV-14

and some other simulation may go for ever. Example: any expression $(\lambda x.2)A$, A being any ob, may be reduced to 2 since, x not occurring in 2, $[A/x]2 \equiv 2$. However if $A \equiv Y$ (the paradoxical combinator), the machine may undefinitely reduce Y. Therefore a reduction of $(\lambda x.2)Y$ may not terminate.

The machine may be modified in order to be deterministic: a reduction such as the one of



may only be applied when the argument N is not reducible any more. Then $(\lambda x.2)Y$ (Y being the paradoxical combinator) becomes a computation which does not terminate in any case.

In Mc Carthy [18] and Landin [12] a sequential machine is used (the order of the reduction is uniquely determined). The evaluation is therefore deterministic.

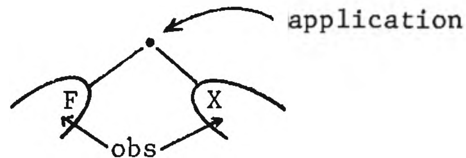
IV.4 Another functional representation: Curry's combinators.

In the previous section, an abstract machine evaluating λ-expressions was supposed to be able to perform replacements and substitutions. We consider a replacement as a simple and basic transformation; however we view a substitution as a much more complex transformation. The question then arises whether it is possible to represent functional computations with an abstract machine which is only able to perform replacements. This is possible with Curry's combinators and moreover formal variables are not needed anymore in the representation. Such a property is concep-

tually important enough to legitimate this section of the thesis. In fact the study of combinatory logic has been a point of departure from which many ideas expressed in this dissertation have emerged.

IV.4.1 Notations and representations.

In this section (and in this section only), we suppose that there is only one combination called application. Whenever F and X are two obs, we represent the ob obtained by applying F to X by FX; in a graphical notation:



Whenever parentheses are not used, association is to be performed on the left (in this section only). So FXYZ ≡ (FX)YZ ≡ (FXY)Z

Any combination different from the application is represented by an atom; for instance, if A denotes addition, A 2 3 represents (A 2) 3 ; A may be considered as an operation of degree 2, (A 2) as an operation of degree 1 . Fig. IV-15 represents the ob:

A (A 2 (A 3 5)) 1

IV.4.2 The combinators K,I,ϕ .

Let us suppose that together with the combinators associated with values and operations we have the following combinators: K,I,ϕ. To each of these combinators is associated a reduction scheme:

For any obs X,Y,Z,T we have:

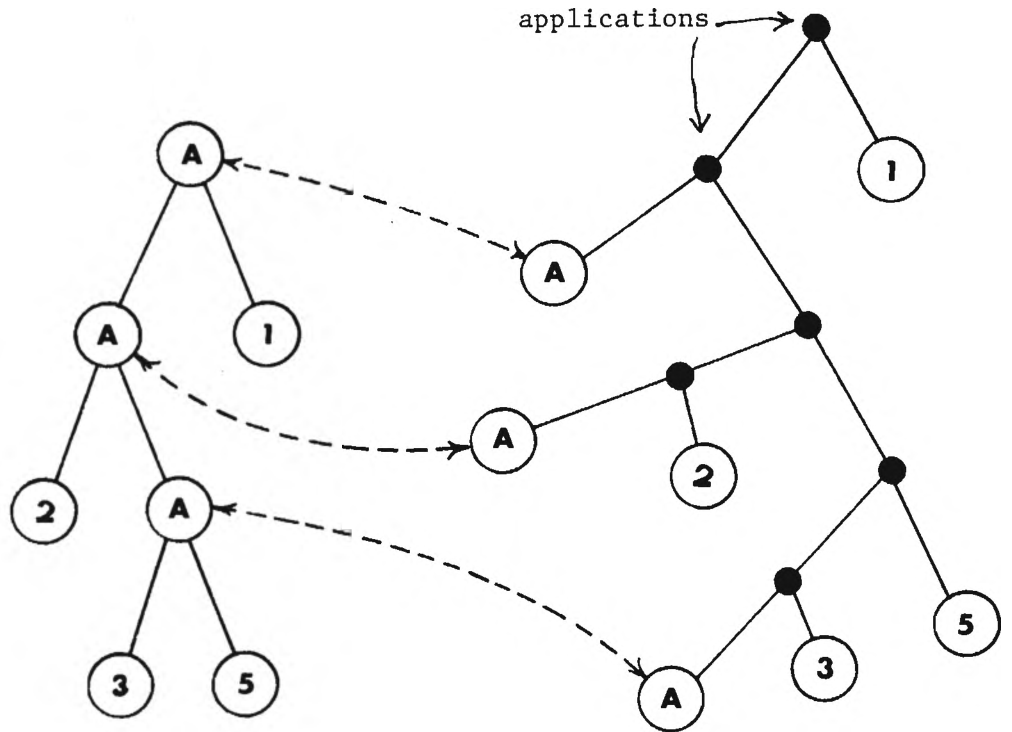
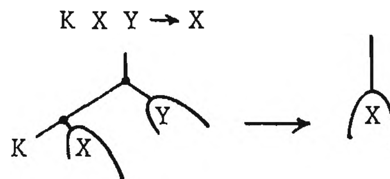
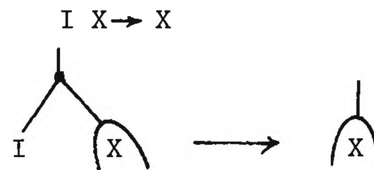


Fig. IV-15

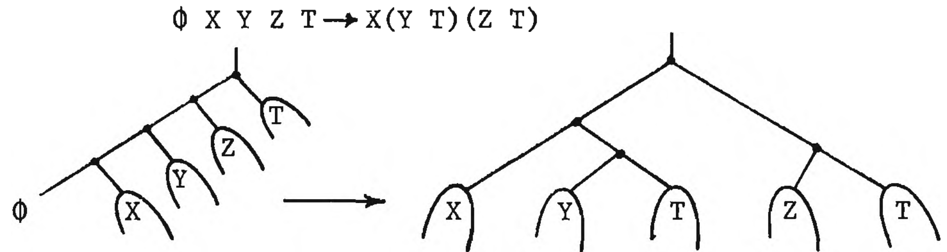
rule (K) : K X Y → X



rule (I) : I X → X



rule (Φ) : Φ X Y Z T → X(Y T)(Z T)



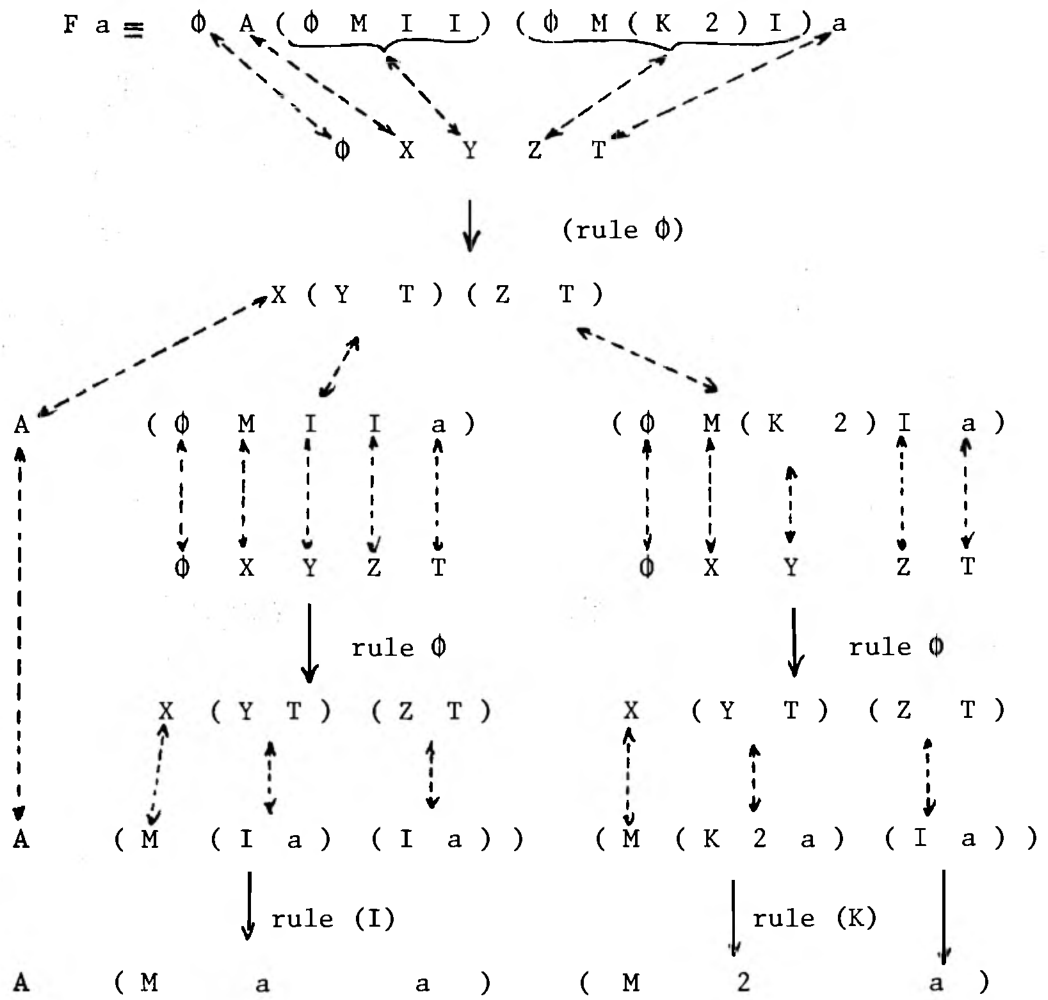
As we have seen in section IV.3 these rules define an abstract reducing machine.

IV.4.3 An example.

A and M denoting the addition and the multiplication, let us consider the ob:

F ≡ Φ A (Φ M I I) (Φ M (K 2) I)

Let us reduce the ob F a , with 'a' standing for any integer:



We have F a ≥ A (M a a) (M 2 a) . So, F is a function which, applied to any integer a, produces A (M a a) (M 2 a) . F is therefore an ob representing λ x . A (M x x) (M 2 x) or, with usual notations, $\lambda x.(x \times x) + (2 \times x)$ and F does not contain any variable.

IV.4.4 An algorithm.

We may wonder if every function may be represented in this way: Let us consider an ob M which may contain a formal variable x. Is there an ob F which does not contain any variable, representing λx.M ? The answer is yes. In order to simplify the discussion, let us assume that M only contains operations of degree 2. Let us call [x]M the ob F we are looking for. We have the following algorithm:

<u>if</u> M does not contain x <u>then</u> [x]M ≡ K M
<u>else</u> <u>if</u> M is x <u>then</u> [x]M ≡ I
<u>else</u> ,

M has the form:
<operator> <left operand> <right operand>

and:

[x]M ≡ Φ <operator>([x]<left operand>)([x]<right operand>)

This algorithm, applied to λx.A(Mxx)(Mxx) produces in fact F.

IV.4.5 B,C,W,S,K.

If M contains several variables x, y, z , then λx y z . M is obtained by forming [x] ([y] ([z] M)) .

We have particularized our discussion with operations of degree 2. However any function may be represented with:

B	(B):	B f g x ≥ f(g x)
C	(C):	C f x y ≥ f(y x)
W	(W):	W f x ≥ f x x

or, alternatively, with:

S	(S)	$S\ f\ g\ x \geq f\ x(g\ x)$
K	(K)	$K\ c\ x \geq c$

With any of these combinators the abstract machine verifies the Church-Rosser condition.The remarks at the end of IV.3 are still valid here.

Remark. Each of the previous combinators may be considered as a lambda-expression. For instance, we can take:

$$B \equiv \lambda f\ g\ x\ .\ f(g\ x)$$

$$C \equiv \lambda f\ x\ y\ .\ f(y\ x)$$

$$W \equiv \lambda f\ x.\ f\ x\ x$$

Then B,C,W may be considered as a "base", any lambda-expression being expressible as an ob constructed with B,C,W.

This discussion is the last of our preliminaries. We are now going to study DCPL.

PART TWO

D C P L

A DISTRIBUTED CONTROL PROGRAMMING LANGUAGE

CHAPTER V

D.C.P.L.

A DISTRIBUTED CONTROL PROGRAMMING LANGUAGE

V.1 The general frame.

V.1.1 A computation as a formal object.

In DCPL, any computation is considered as a formal object, an ob (see section III.1). In the computing machine, such an ob is actualized as a tree whose nodes, representing combinators, are automata, and whose edges are directed channels of information, directed from "son" to "father" (fig. V-1). A node may have zero, one or several sons; the number of sons it has is called its degree (fig. V-2).

The edges being directed channels of information, each node may send a message up the tree; such a message is called a notice (fig. V-3). Whenever a node receives a notice, it may either pick it up or pass it along upward.

The tree is considered to be in some space in which each node has an address. There is a communication system allowing any node to send a message to any other node whose address it knows; such a message is called a reply. A reply contains explicitly the address of the addressee.

Whenever a node A wants some information from ancestor B, A sends up a notice with its own address (fig. V-4); such a notice is called a request. A request is a notice containing the address

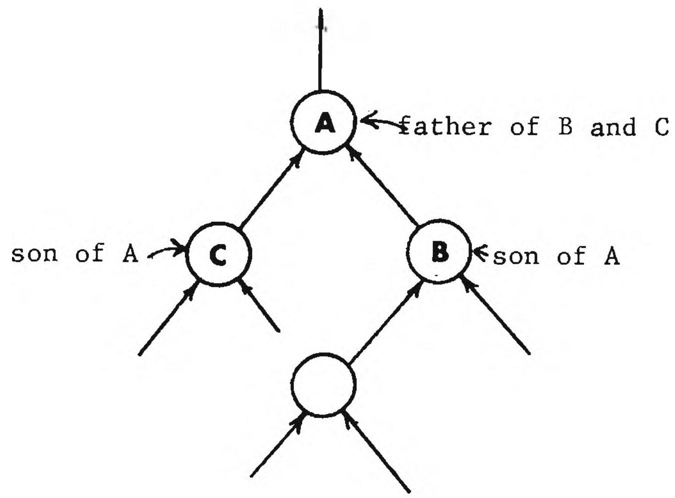


Fig. V-1

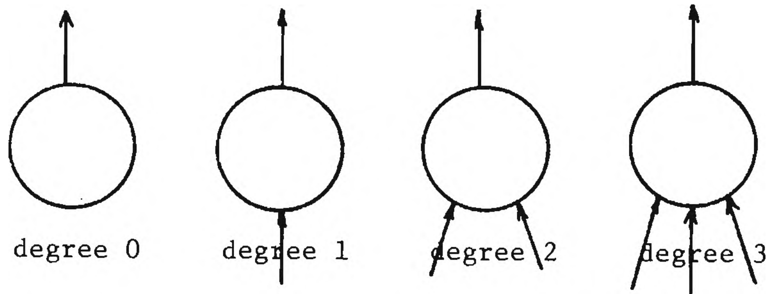


Fig. V-2

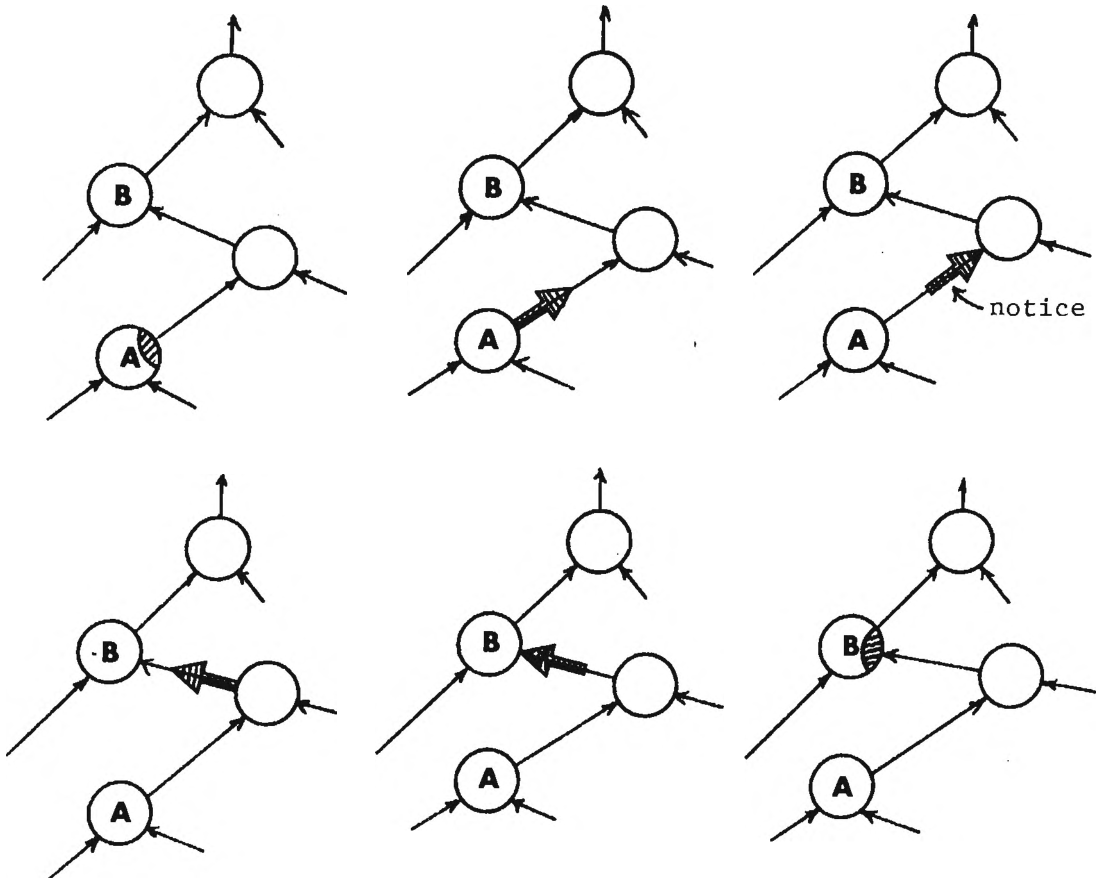


Fig. V-3

B
address of A
A

B
request
A

B
A

B
A

B
address of A
A

Fig. V-4

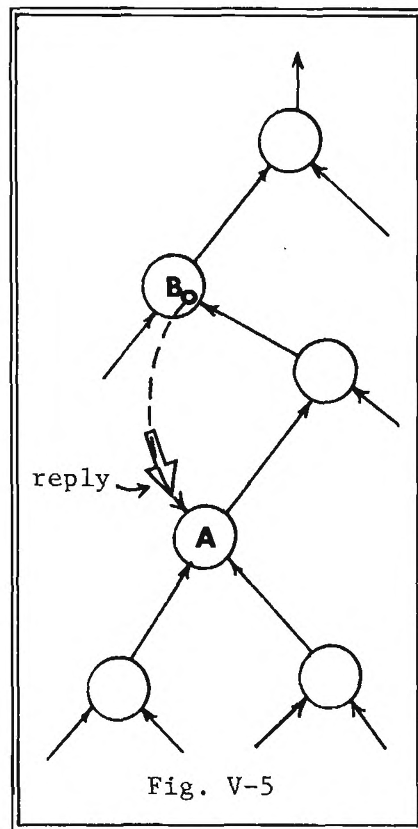


Fig. V-5

of the sender to which some reply is to be made. A notice which is not a request is a simple notice. On receiving the request sent by node A, the node B may send back the required information (fig. V-5).

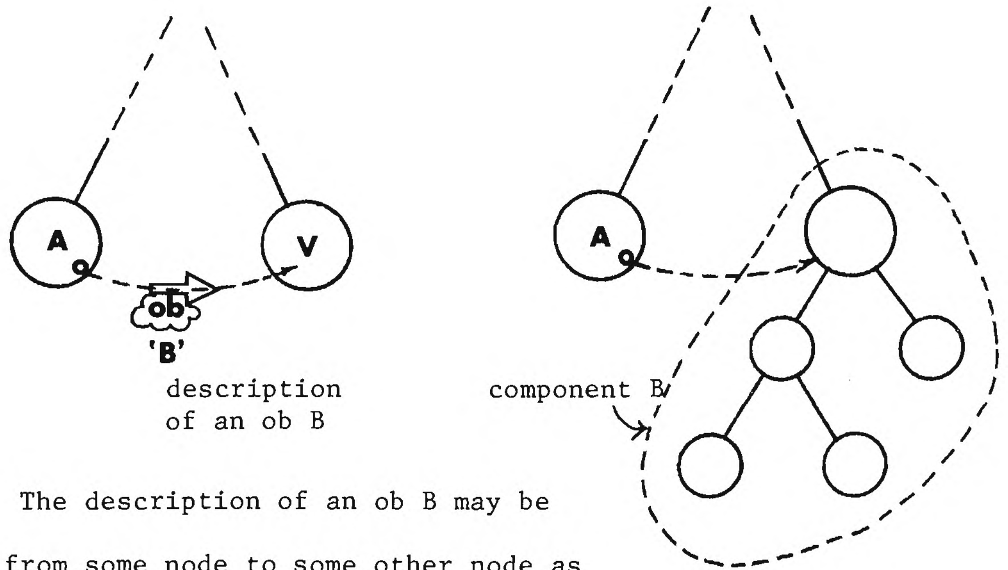
V.1.2 Values and operations: their representation.

As it has been already noted in the introduction (chapter I), we assume that there is a universe of values structured in classes or types, and mappings from some classes to some possibly different classes which are actualized by operations, and we leave the list of types and operations open-ended. We require however that:

1. to each value in the universe corresponds a specific combinator, represented by a value-node. Spontaneously a value-node sends up the tree a simple notice containing the value it represents; the node then disappears (fig. I-3, chapter I).

2. to each operation in the universe corresponds a specific operator represented by an operation-node, which has the degree of the operation it represents. An operation-node waits until it receives a value from each of its sons; it performs, then, the operation it represents on these values.The result is sent up the tree, and the operation-node disappears(fig. I-4). Fig. I-5, in the introduction, shows how a simple expression without variable may be evaluated.

3. to each ob corresponds a value in the universe,which may be considered as a description of the ob. As a result, a description of an ob may be sent as a message from some part of a computation to some other part, where the described ob may be built and implemented in space, extending the tree (fig. V-6).



The description of an ob B may be sent from some node to some other node as a value. When received the ob.B may be implemented in space.

Fig. V-6

As a result of these behaviors, the tree representing a computation may grow and shrink during its life.[a)]

V.1.3 The other combinators.

Together with the combinators associated to the values or operations, DCPL uses many other specific combinators:

Triadic combinators (degree 3):

'→'	(setdown)
'←'	(setup)
'*IF*' or '|'	(if)

Dyadic combinators (degree 2);

'*NEW*'	(new)
'*LAMBDA*'	(lambda)
'*MU*'	(mu)
'↑'	(sendup)
'↓'	(sendown)
'□→'	(trigger)
'*IMPL*'	(implement)
'∘'	(compose)
'*OR*'	(or-event)
'*AND*'	(and-event)

---

a) Such a behavior is similar to the behavior of the executing graph used in the Adams' model (IV.2). Instead of graphs, the model described here manipulates trees.

<u>Degree zero combinators</u>:

∅ (null ob)

a whole class of variables: *X*, *Y*, *JOHN*, *VAR* , ...

a whole class of alpha-variables: α*X*, α*Y*, α*JOHN*, α*VAR* , ...

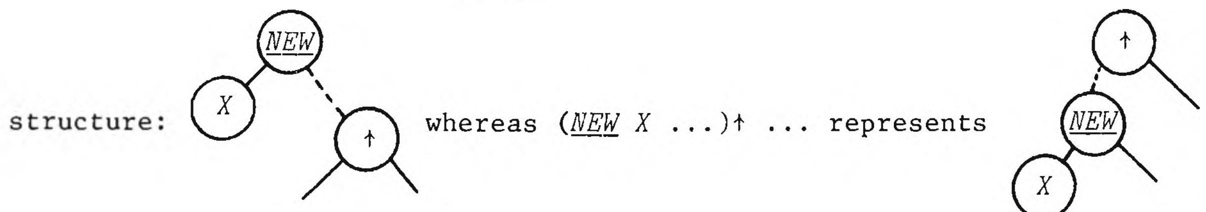
The function and the behavior of each of these nodes will be progressively presented in the sequel.

V.2 About the syntax.

The tree-structured representation is fundamental in DCPL. However, it is useful to have a linear representation easier to manipulate. Rather than a straight parenthesization or a heavy syntactic apparatus, we prefer the use of a few replacement rules: programs gain a familiar appearance without any syntactic freezing. The following rules determine a one-to-one mapping between the tree-structured and the linear representation.

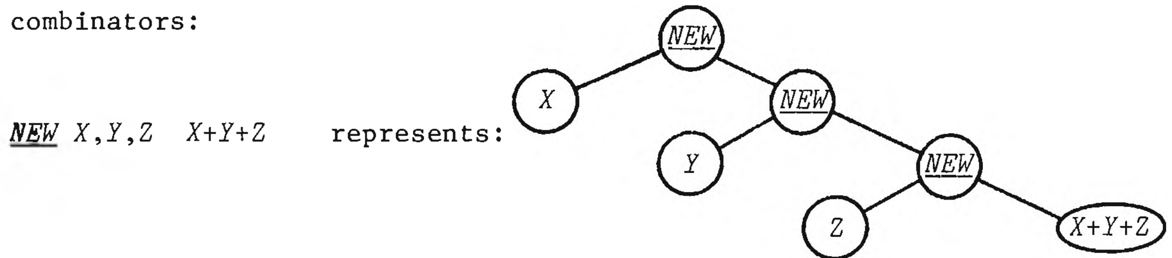
<u>Rule 1</u>. Parentheses may be used freely and no distinction is made between parentheses and brackets.

<u>Rule 2</u>: Any combinator of degree 2 may be infixed. Association is to be done to the right, unless explicitly specified through the use of parentheses. For instance $X+Y+Z$ is to be structured as $(X+(Y+Z))$. As a consequence, whenever two combinators of degree 2 are not ordered by parentheses, the latter is, on the tree structure, a descendant of the former: <u>*NEW*</u> *X* .... ↑ .... has the following tree structure: whereas (<u>*NEW*</u> *X* ...)↑ ... represents



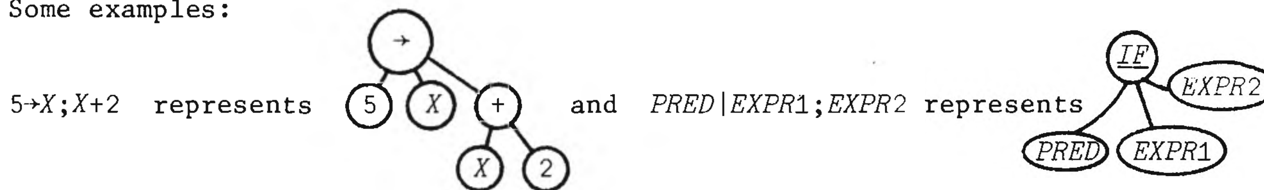
<u>Rule 3</u>. Much like in Algol where <u>integer</u> X,Y,Z; is used instead of <u>integer</u> X; <u>integer</u> Y; <u>integer</u> Z; , one can factorize combinators:

<u>*NEW*</u> *X*,*Y*,*Z* *X*+*Y*+*Z* represents:

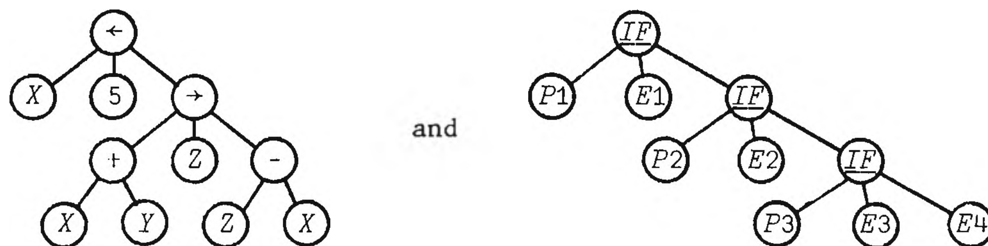


<u>Rule 4</u>. Combinators of degree 3 may be infixed between their first two arguments, a semi-colon separating the second argument from the third. Rule 2 is then applicable to such combinators. Some examples:

5→*X*;*X*+2 represents [tree] and *PRED*|*EXPR*1;*EXPR*2 represents [tree]

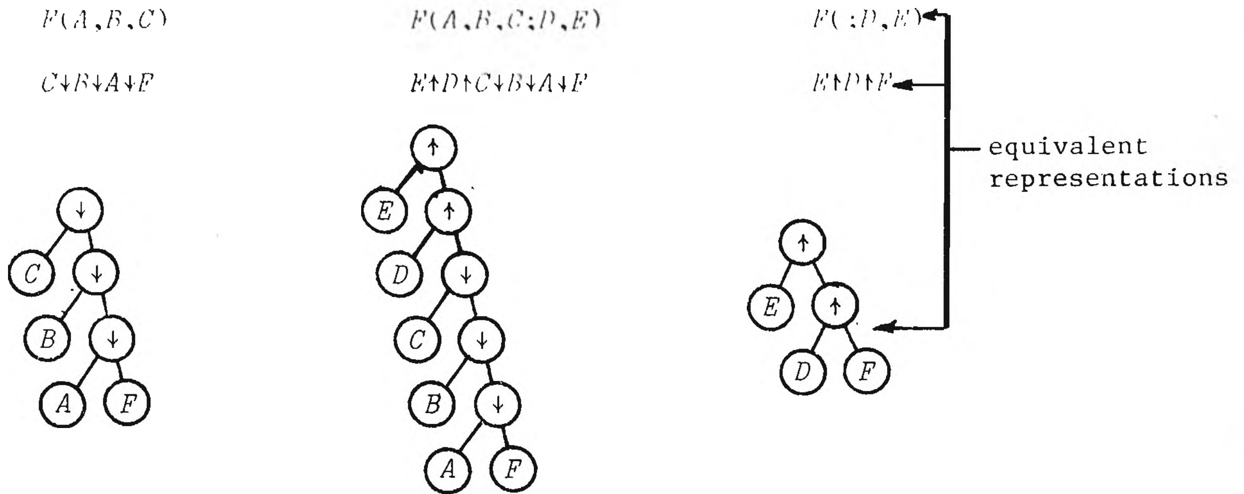


The interest of this rule is sharpened when it is used repetitively on a nested structure:



may be represented respectively by *X*←5;(*X*+*Y*)→*Z*;*Z*-*X* and by *P*1|*E*1;*P*2|*E*2;*P*3|*E*3;*E*4 or, with a non-necessary pair of parentheses, (*X*←5;(*X*+*Y*)→*Z*;*Z*-*X*) and (*P*1|*E*1;*P*2|*E*2;*P*3|*E*3;*E*4) .

<u>Rule 5</u>. As it will be discussed in the sequel, the application of a function to an argument or to a parameter is represented through the use of '↓' ('sendown') or '↑' ('sendup'). However, in order to conform to traditional notations, we allow expressions such as *F*(*A*,*B*,*C*) , *F*(*A*,*B*,*C*;*D*,*E*) , *F*(;*D*,*E*) :



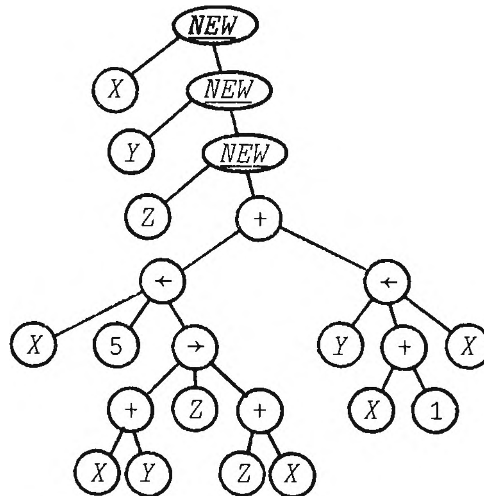
Remark. Application of rule 5 may be misleading: $F(A,B,C)$ is an ob; (A,B,C) is a notational convenience which has no meaning by itself, it does not represent any ob.

An example:

According to these rules, the expression

$\underline{NEW}$ X,Y,Z $(X\leftarrow 5;(X+Y)\rightarrow Z;Z+X)+(Y\leftarrow X+1;X)$

is the following ob:

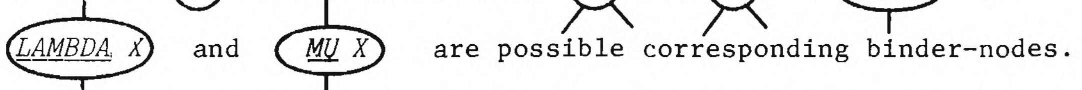


This is as much as is necessary to know about the "syntax" of DCPL.

V.3 The binding process.

In DCPL a system of mutually bound variables determines communication paths. The process which builds these paths is called the binding process. The binding process superimposes a graph structure to the tree structure which represents a computation.

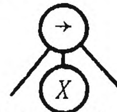
A communication path leads from a "binder-node" to a variable-node. If (X) is a variable-node, (→X) a), (X←), (NEW X), (LAMBDA X) and (MU X) are possible corresponding binder-nodes.

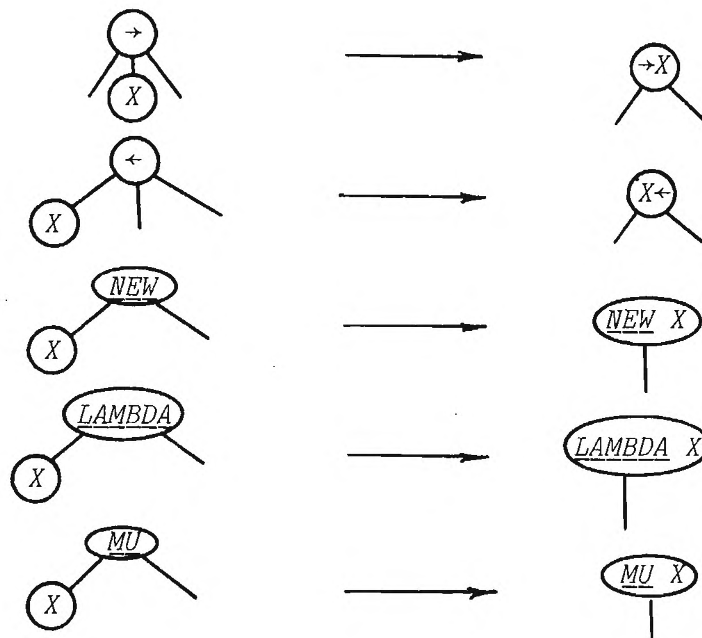


Any variable-node sends spontaneously a request containing

---

a) Note: '→' is a combinator of degree 3, whose second argument must be a variable. However, the construct

is considered as being one node  . In fact, in the following list, the constructs on the left are considered to represent the constructs on the right:

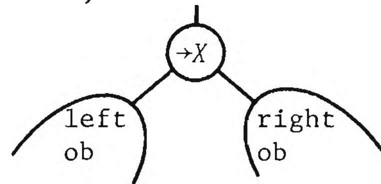


its address and its name (for instance 'X'), up the tree (fig. V-7). Such a request is picked up by the first corresponding binder-node encountered. The variable-node is then bound by this binder-node: a communication path leads from the binder-node to the variable-node.

A binder-node may bind several variable-nodes. Whenever a binder-node receives an argument, it sends a copy of this argument to each variable-node it binds (fig. V-8).

V.3.1 The binder-node setdown: '→'.

X being an identifier, $\to X$ combines two components:



$\to X$ may only pick up the requests reaching it from the right (fig. V-9).

On receiving a value simple notice on the left, setdown sends a copy of this value to every variable-node it binds, on the right. The scope of $\to X$ is therefore limited to the right component; $\to X$ may be considered as assigning the value produced on the left to the variable-nodes X in its scope (fig. V-10).

On receiving a value simple notice on the right, $\to X$ passes it along up the tree and disappears (fig. V-11).

Example 1:

The evaluation of the expression $(5 \to X; X+2)$ is displayed in fig. V-12.

Example 2:

The evaluation of the expression $(2 \to X; 3 \to Y; X+Y)$ is displayed in fig. V-13.

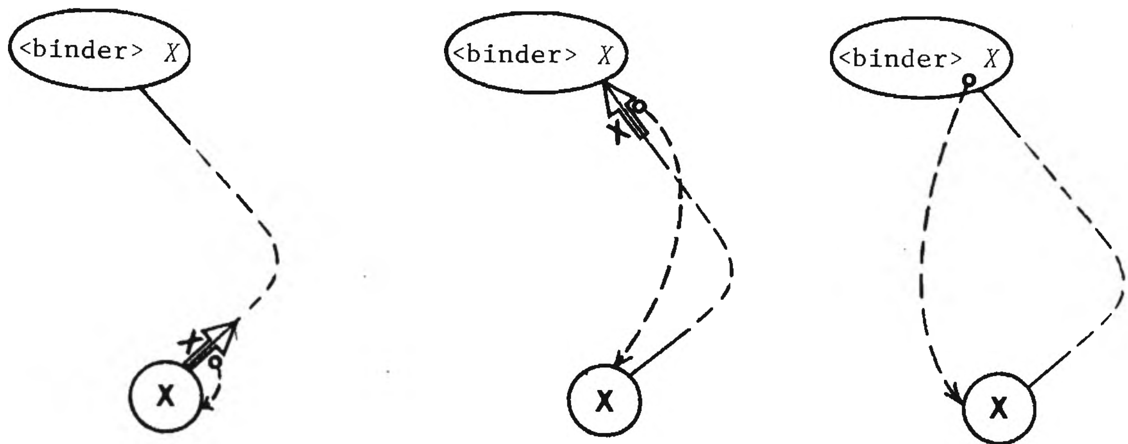


Fig. V-7

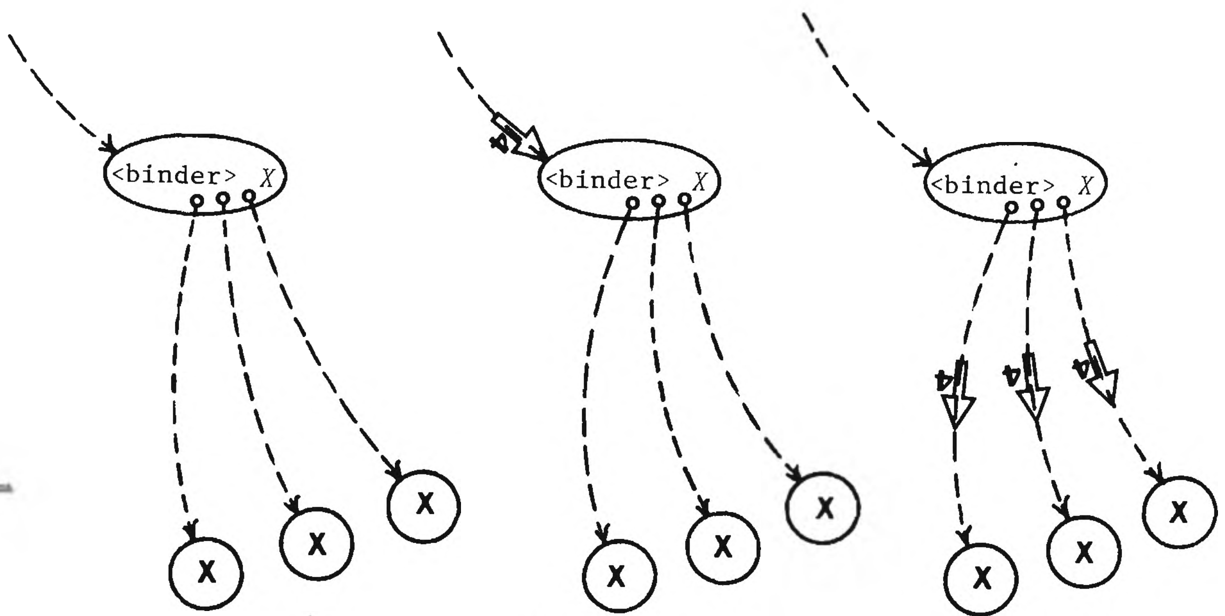


Fig. V-8

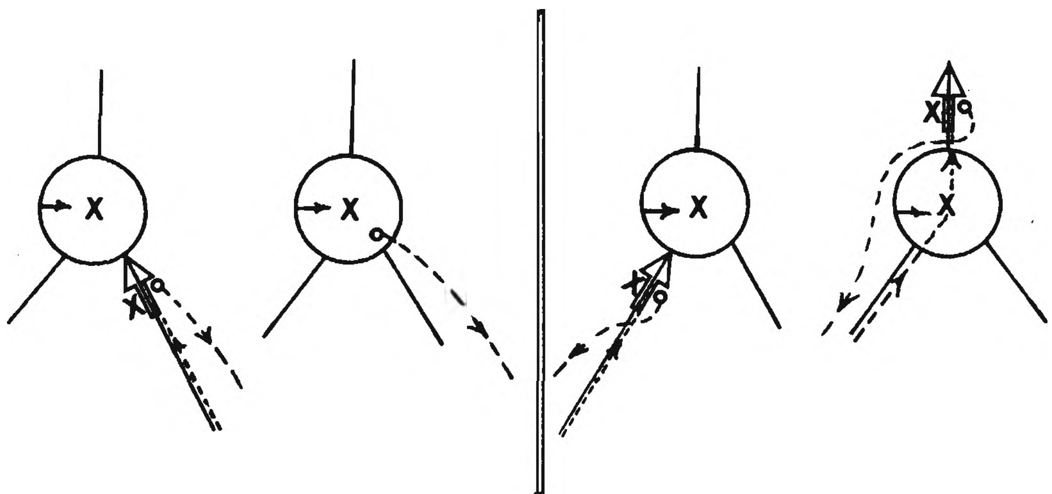


Fig. V-9

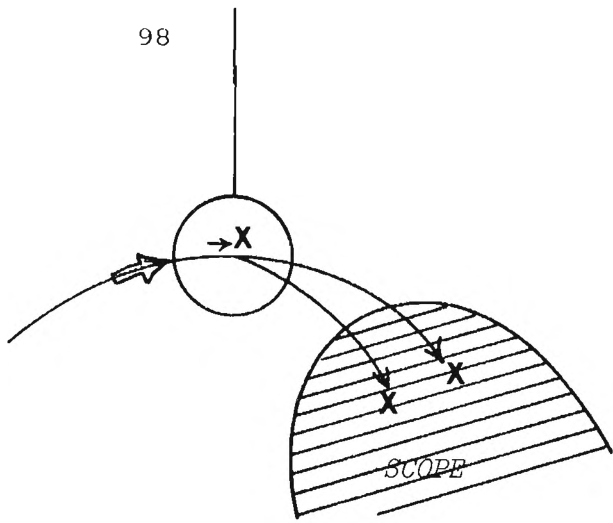


Fig. V-10

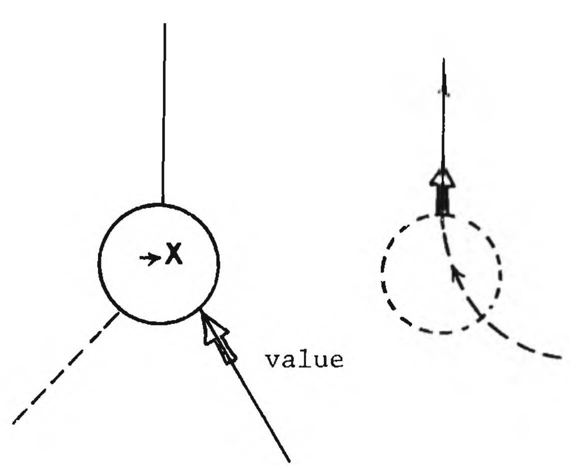
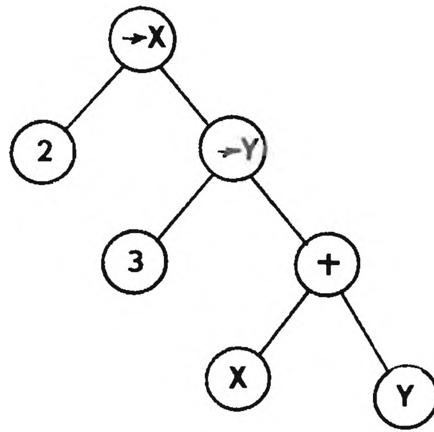


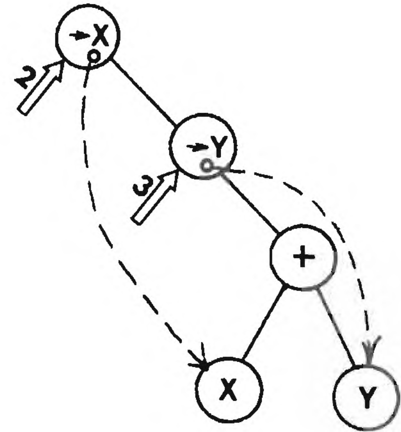
Fig. V-11

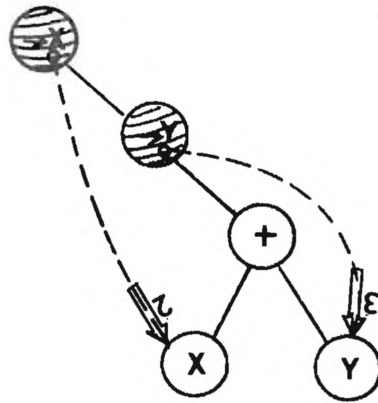
(5→X;X+2)

Fig. V-12

(2→X;3→Y;X+Y)







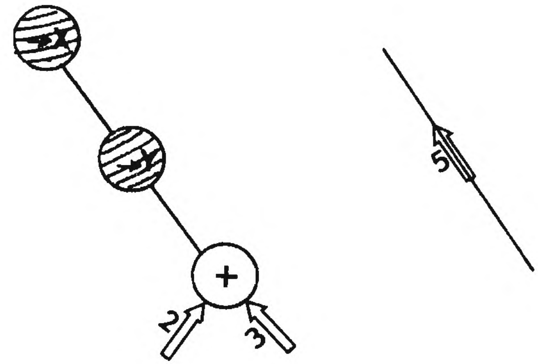


Fig. V-13

Example 3:

The evaluation of the expression

$$[(2+3)\rightarrow X;5\rightarrow Y;(X+Y)\rightarrow X;(Y+1)\rightarrow Y;(2\times Y)-X]$$

has already been studied in the first chapter (fig. I-6 and I-11).

V.3.2 The binder-nodes setup '←', and '*NEW*'.

As already noted, the scope of $\rightarrow X$ is limited to the right component. $X\leftarrow$ allows to send a value in the right component and in the environment as well (fig. V-14). In order to delimit exactly the scope of $X\leftarrow$, *NEW* X is used as a "top-binder": the scope of $X\leftarrow$ is entirely under the corresponding *NEW* X (fig. V-15).

$\rightarrow X$ and $X\leftarrow$ bind exactly in the same way any variable-node X in their right component. However, contrarily to $\rightarrow X$, $X\leftarrow$ sends up the tree a request with its name which is picked up by the first *NEW* X the request encounters. As a reply, *NEW* X sends to $X\leftarrow$ the address of any variable-nodes whose requests it has received (fig. V-16).

Whenever $X\leftarrow$ receives a value on the left, it sends a copy of it to each variable-node whose address it possesses (fig. V-17).

$X\leftarrow$, as $\rightarrow X$, disappears whenever it receives a value simple notice on the right. *NEW* X disappears whenever it receives a value simple notice.

Example 1:

Fig. V-18 displays the evaluation of the expression:

$$\underline{NEW}\ V\ (2\rightarrow U;V\leftarrow U+1;U+V)\ +\ (3\rightarrow U;U+V)$$

in the left argument of the addition, a value V is produced; V is used in the evaluation of both arguments.

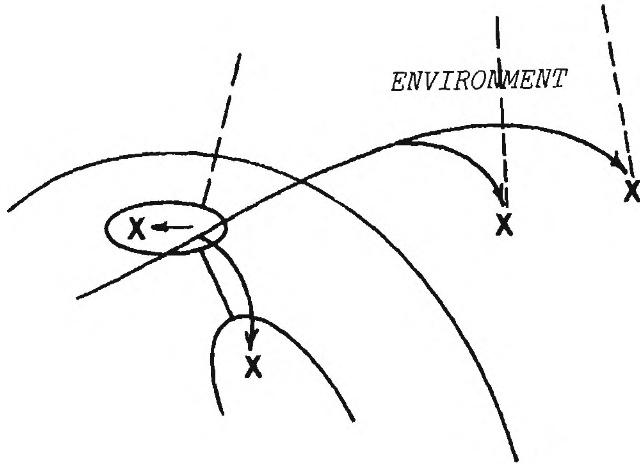


Fig. V-14

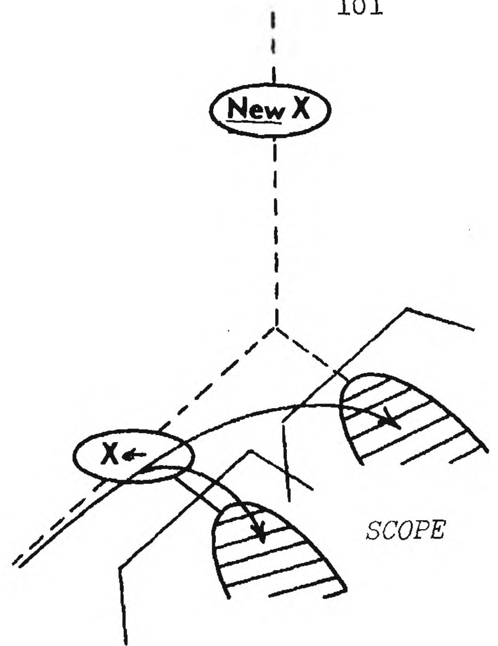


Fig. V-15

Fig. V-16

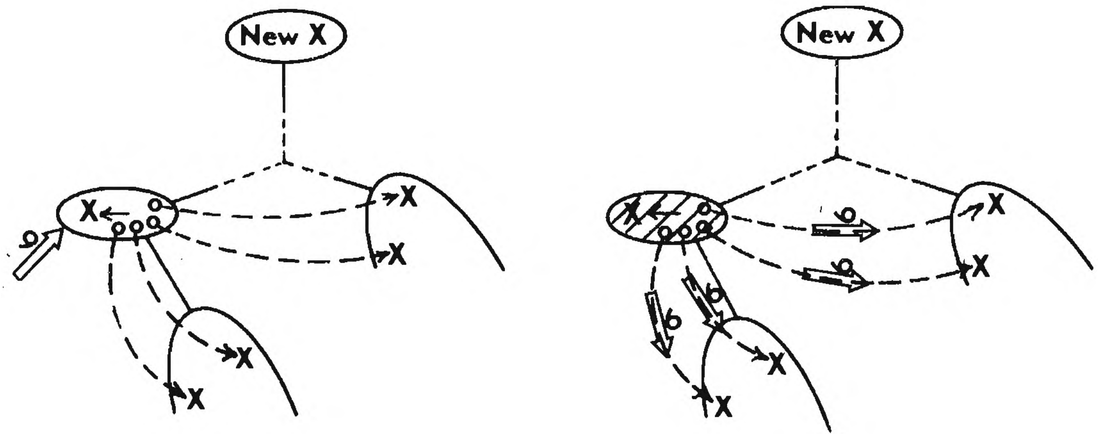


Fig. V-17

$\underline{NEW}\ V\ (2\rightarrow U; V\leftarrow U+1; U+V) + (3\rightarrow U; U+V)$

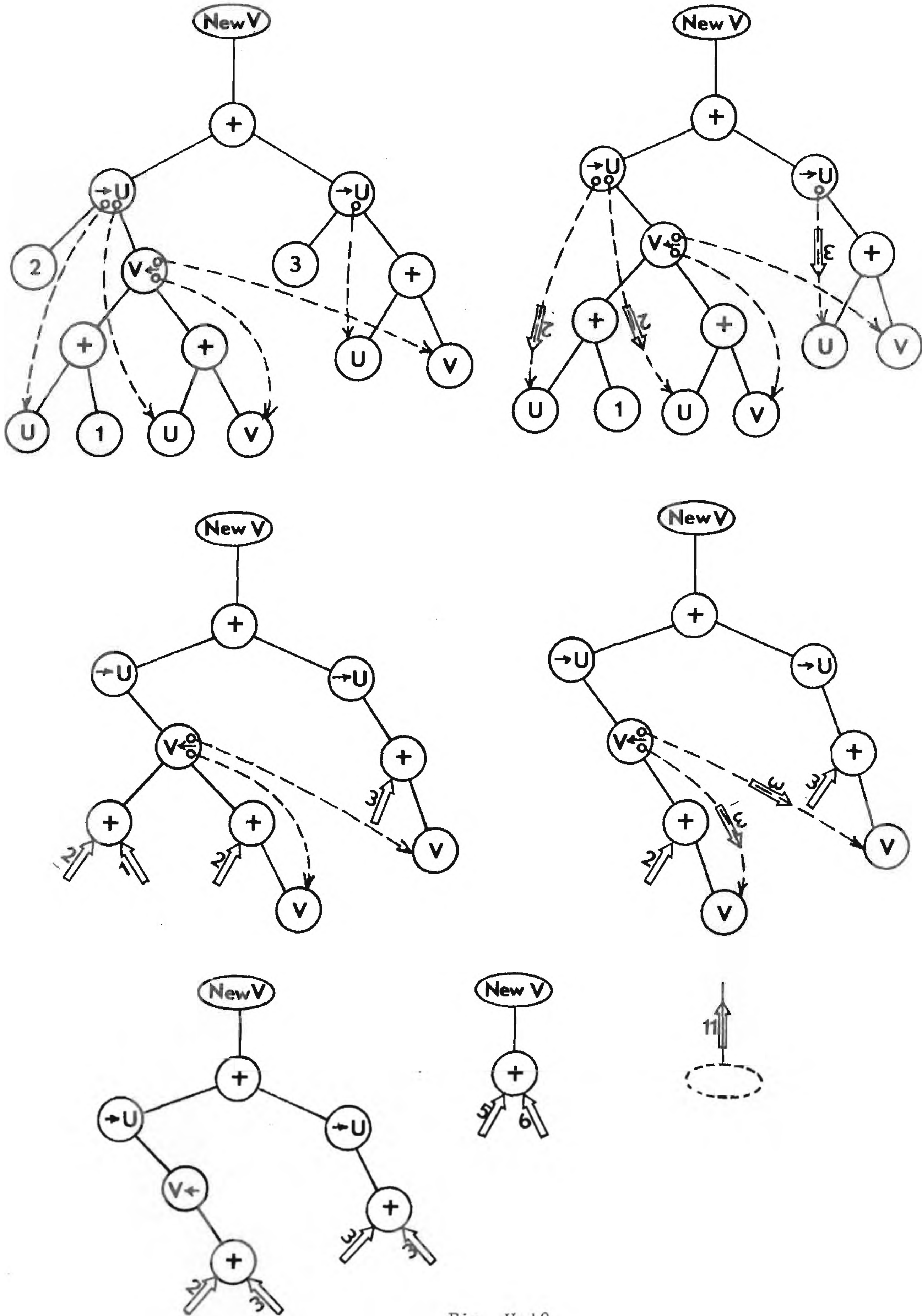


Fig. V-18

Example 2:

Fig. V-19 displays the evaluation of the expression:

$$\underline{\underline{NEW}}\ U,V \quad (U \leftarrow 5;[U+V] \rightarrow Z;Z+U) + (V \leftarrow U+1;U) \ .$$

Such an expression is remarkable in that it cannot be evaluated by a sequential machine: the evaluation of the first argument requires the value of V which is produced in the second argument whose evaluation requires the value U which is produced in the first argument. In DCPL there is no deadlock since the two arguments are evaluated concurrently.

V.3.3 The binding of procedures.

A procedure is a computation ob which is intended for implementation and usage in various parts of a computation program. In order to ensure generality of utilization, the set of variables used inside a procedure must not conform to the set of variables used in the environments in which the procedure is to be implemented (please, see fig. V-20).

Whenever a variable in a procedure is to receive an argument from the environment, the variable is bound inside the procedure to a *LAMBDA* binder-node; the procedure itself is interfaced with its environment through a node sendown '↓' (fig. V-21a).

Whenever a variable in a procedure is to send an argument to the environment, the variable is bound inside the procedure to a *MU* binder-node; the procedure itself is interfaced with its environment through a node sendup '↑' (fig. V-21b).

$\underline{NEW}\ U,V\ (U\leftarrow 5;[U+V]\rightarrow Z;Z+U)\ +\ (V\leftarrow U+1;U)$

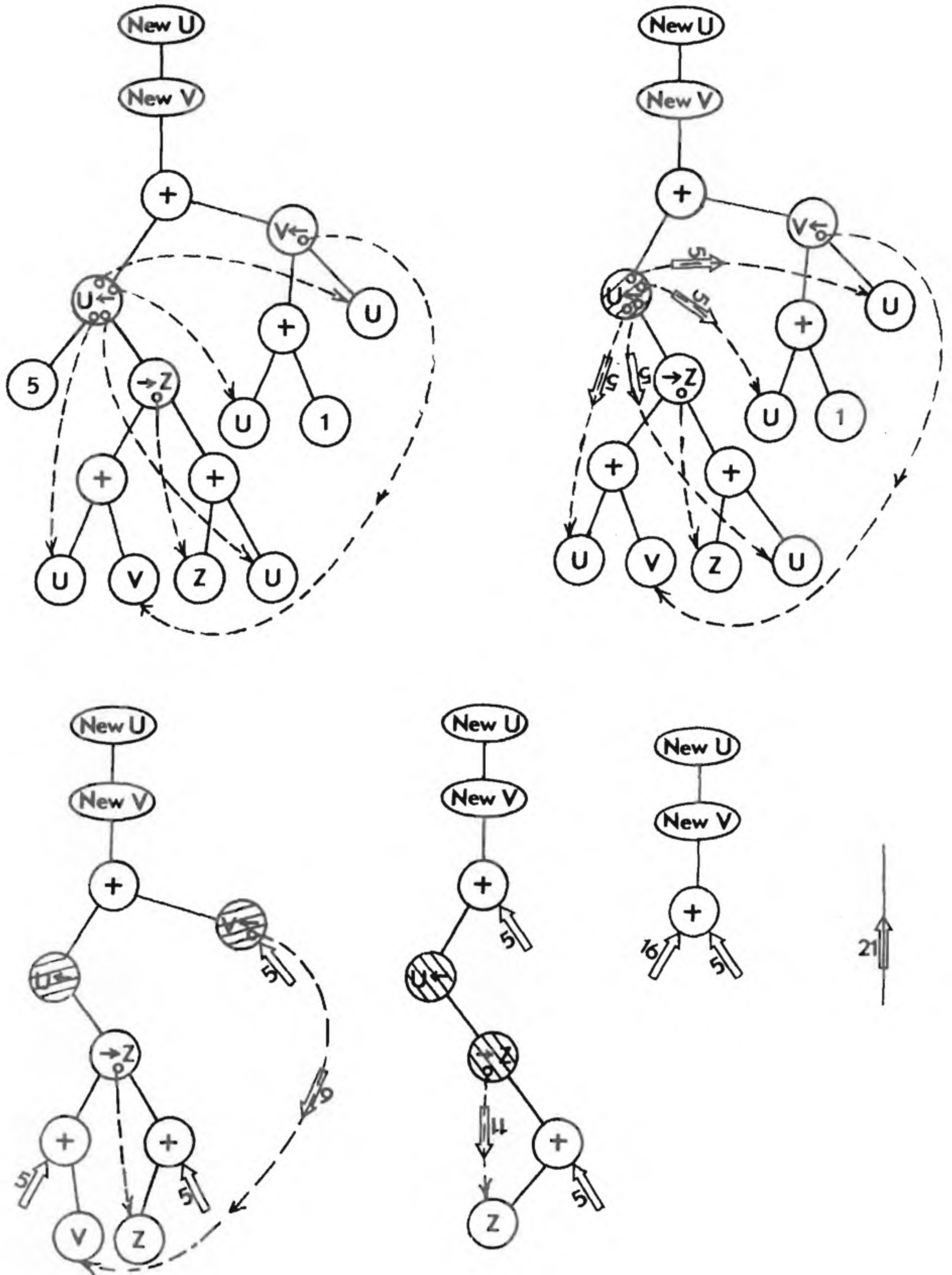


Fig. V-19

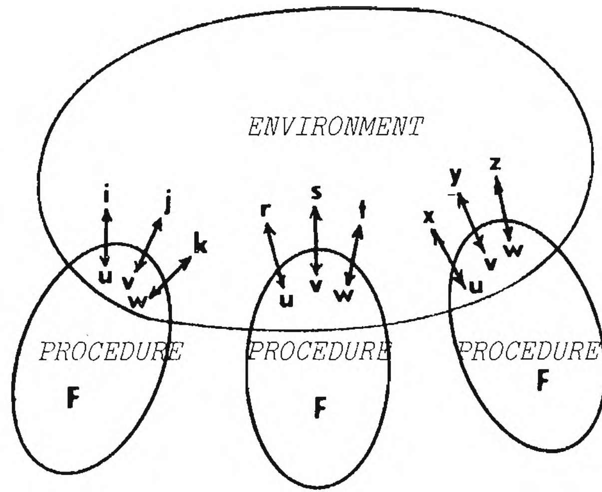


Fig. V-20

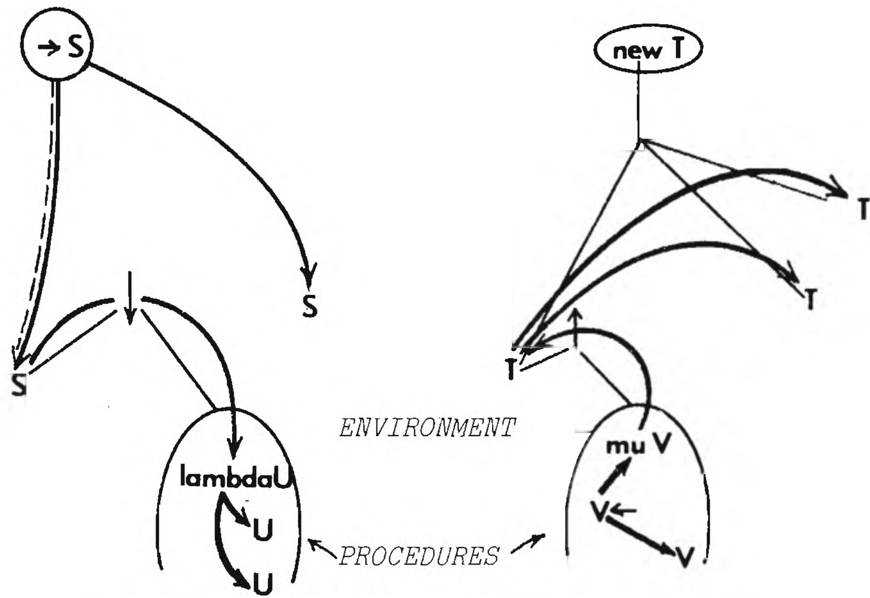
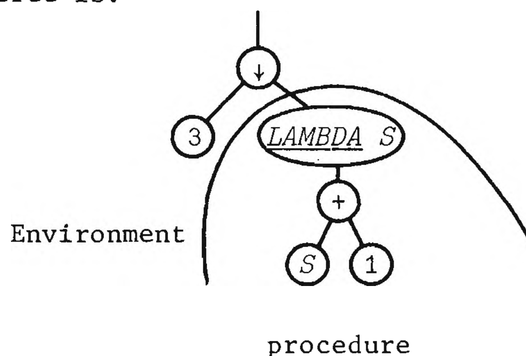


Fig. V-21a

Fig. V-21b

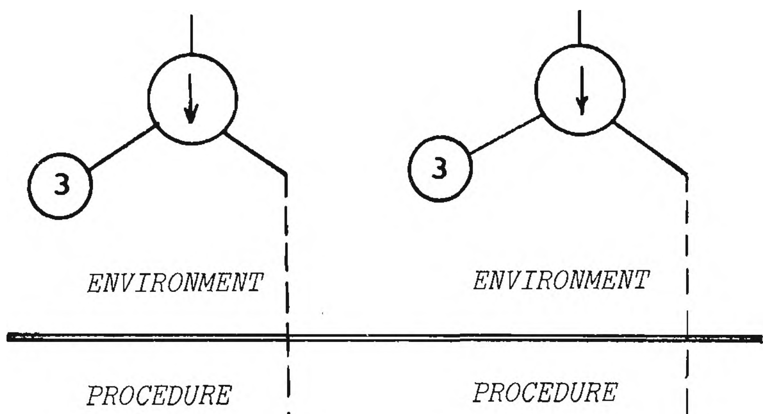
'LAMBDA' and sendown '↓'.

An example: LAMBDA S (S+1) is a procedure with one lambda-variable S. [LAMBDA S (S+1)] (3) represents that procedure applied to one argument 3. According to the (syntax) replacement rule number 5, the computation tree is:



In fig. V-22a and V-22b the variable-node S is bound to the lambda-node: the former sends up the tree a request which is picked up by the latter. This binding occurs inside the procedure itself and may be performed before the implementation in space of the procedure, i.e., the physical embedding of the procedure in its environment. Then the procedure is bound to its environment: the lambda-node sends up the tree a request with its address which is picked up by the sendown node (fig. V-22c and V-22d). The argument may flow down, from the environment into the procedure, and reach the variable-node (remainder of figure V-22).

Whenever a procedure contains several lambda-variables, it is necessary to have the arguments matched with the lambda-nodes (fig. V-23). The matching is performed dynamically in the following way: when a lambda-node sends up the tree a request with its address, this request contains a count set to zero. This count is incremented by one whenever the demand encounters another lambda-node, and decremented by one whenever the demand encounters a sendown node



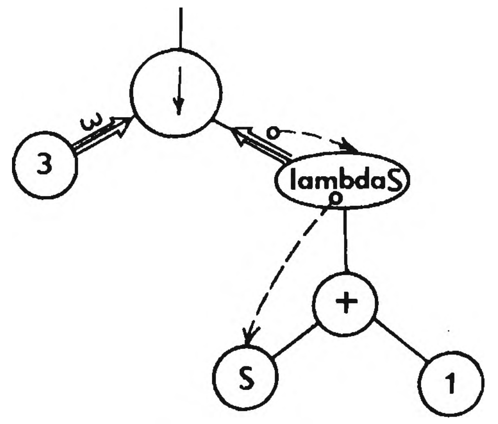
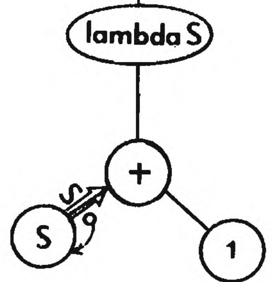


Fig. V-22c



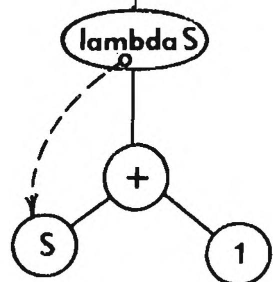
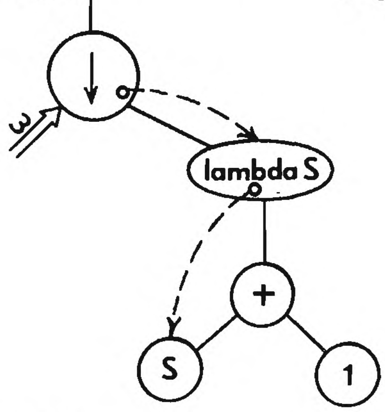
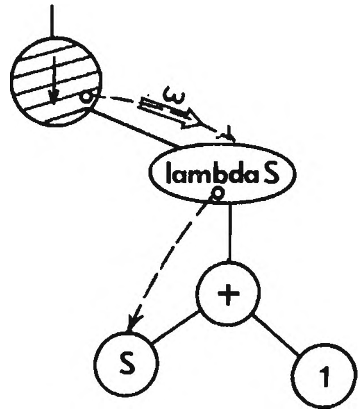


Fig. V-22a

Fig. V-22b





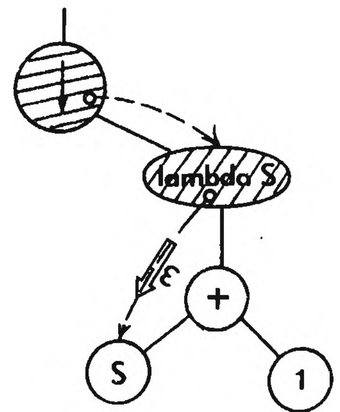
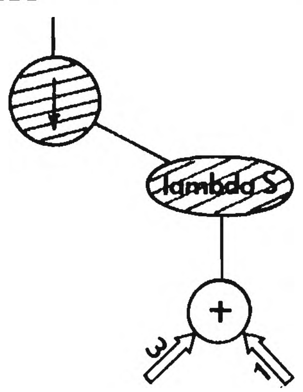


Fig. V-22d



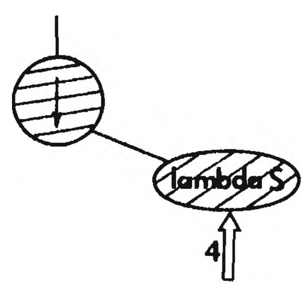


Fig. V-22

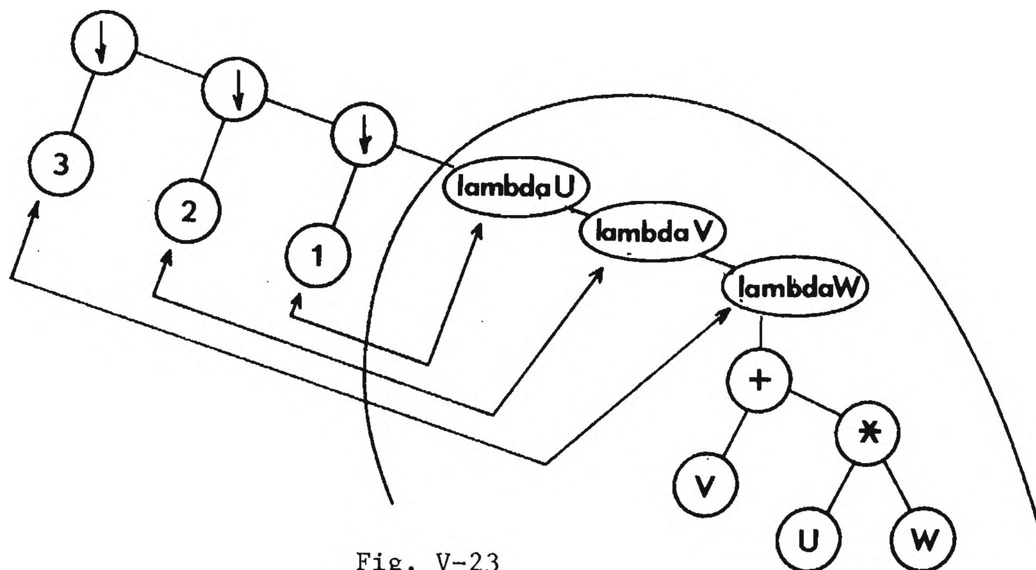


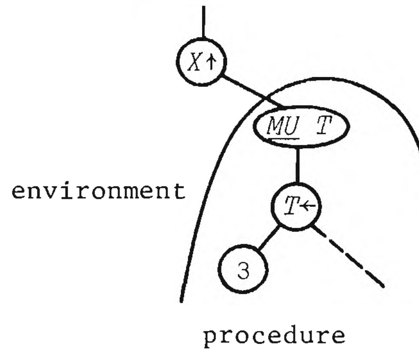
Fig. V-23

Fig. V-24

with a non-zero count. The demand is picked up by the first sendown node which receives it with a count equal to zero (fig. V-24).

'MU' and sendup '↑'.

An example: ..... [MU T ... T←3 ...] (;X) ...



The binding inside the procedure and the binding of the procedure to the environment are performed exactly as for lambda and sendown. However, instead of receiving an argument, the variable in the procedure receives the address in the environment to which the produced value is to be sent (fig. V-25).

Let us consider again the example of fig. V-19:

NEW U,V (U←5;[U+V]→Z;Z+U) + (V←U+1;U)

It may be expressed as

NEW U,V $F(V;U)$ + $G(U;V)$

with F ≡ MU U LAMBDA V (U←5;[U+V]→Z;Z+U)

and G ≡ LAMBDA U MU V (V←U+1;U) . In order to make it clear that U and V are bound in F and G, we can take

F ≡ MU R LAMBDA S (R←5;[R+S]→T;T+R)

and G ≡ LAMBDA R MU S (S←R+1;R) (fig. V-26).

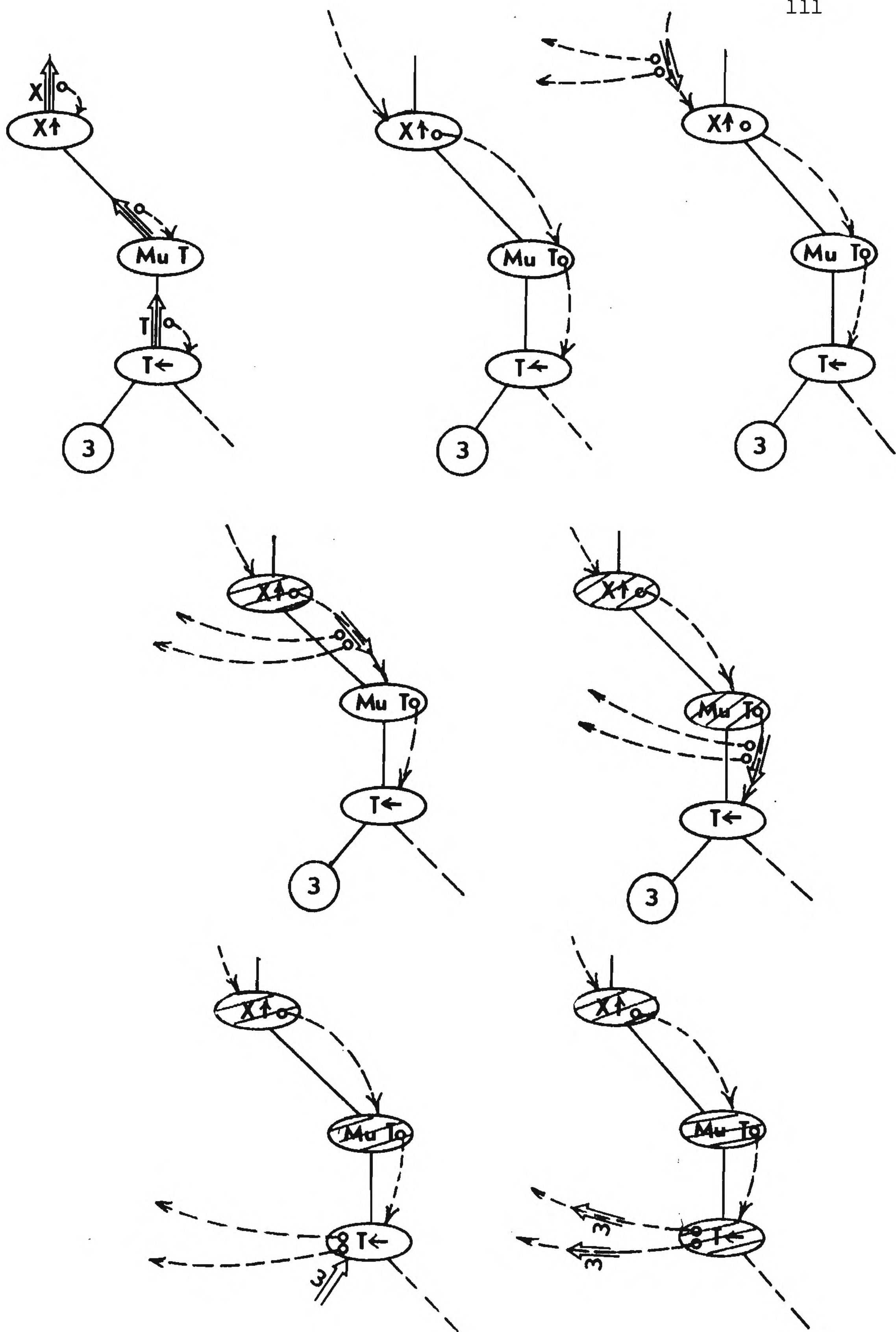


Fig. V-25

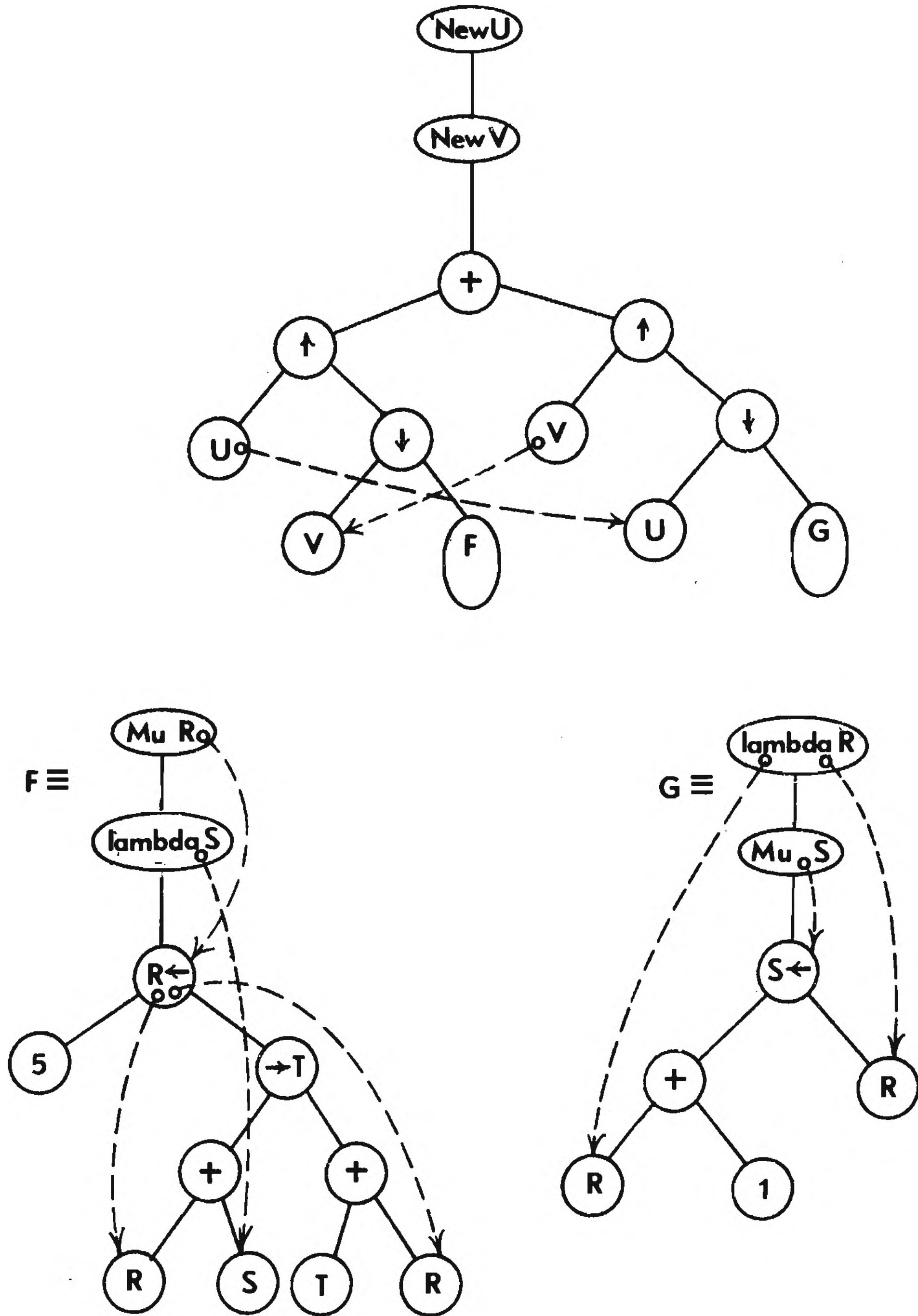


Fig. V-26a

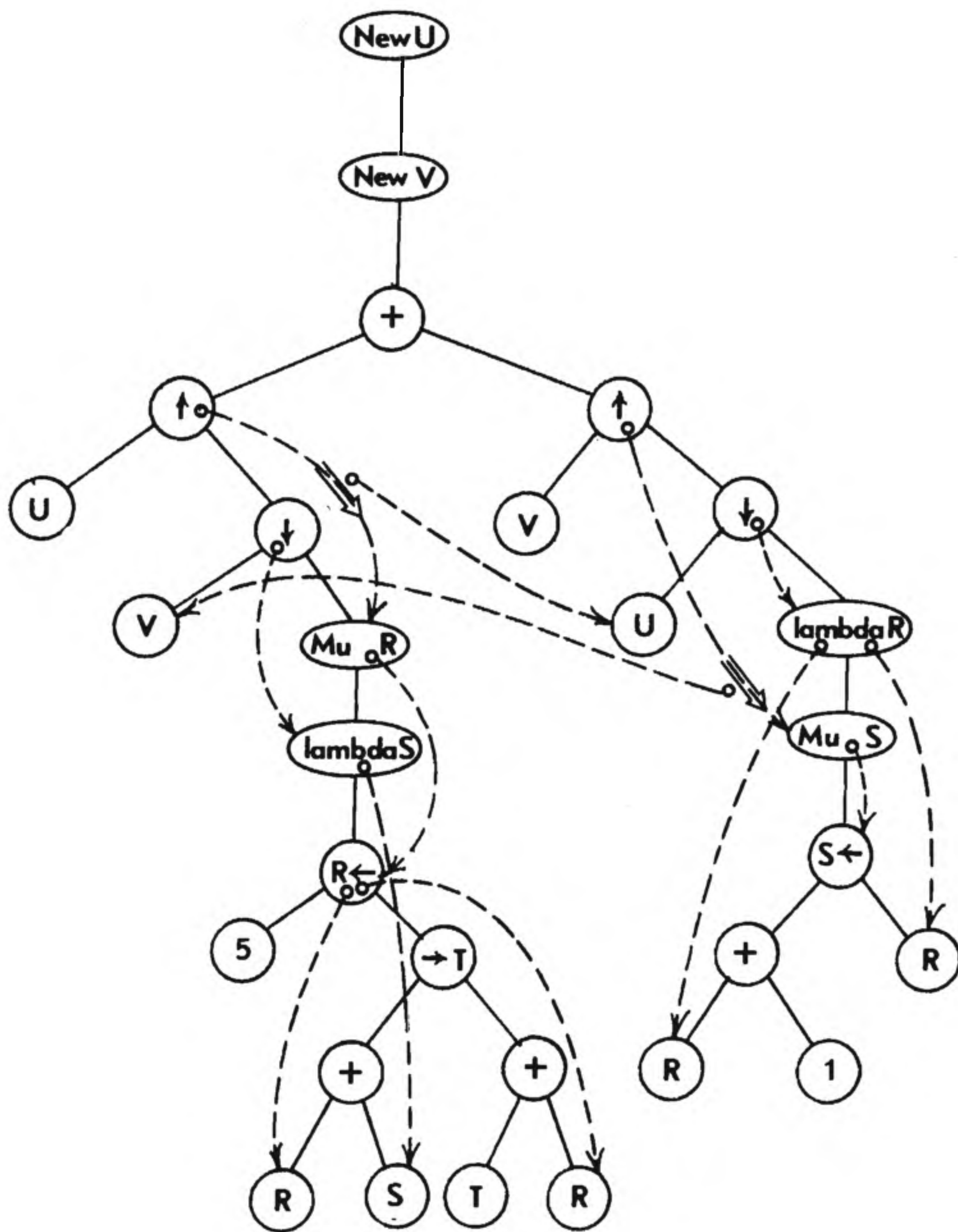


Fig. V-26b

V.4 Procedures as arguments.

V.4.1 Pseudo-values.

As stated in section V.1.2, any ob has its description as a value in the universe of values. Such a value is called a pseudo-value: "*LAMBDA* S (S+1)" is a pseudo-value representing the ob *LAMBDA* S (S+1).

Whenever a variable receives a pseudo-value as an argument, it implements the ob described by the pseudo-value at its own place. As a result a procedure may be received as an argument and implemented if the corresponding pseudo-value is given beforehand (fig. V-27).

Let us however consider the following example:

.... *LAMBDA* S,T $(S+T) \rightarrow F$ .... $F(5) \rightarrow G$ .... $G(4)$ ....

In this example the procedure *LAMBDA* S,T $(S+T)$ is first implemented in the environment A and receives the argument 5; then the resulting procedure is to be sent and implemented in the place of G in the environment B. This may only occur if the ob [*LAMBDA* S,T $(S+T)$](5) is able to produce a pseudo-value representing itself. We will see in the sequel how we may implement such a pattern of behavior.

V.4.2 Pseudo-argument.

There is a special message $\emptyset$ called pseudo-argument: whenever a variable X receives a pseudo-argument, it sends up the tree the pseudo-value representing itself, "X" (fig. V-29).

Whenever the combinator representing an operation receives the proper number of input values, it computes the resulting value and sends it up the tree. If the same combinator receives at least

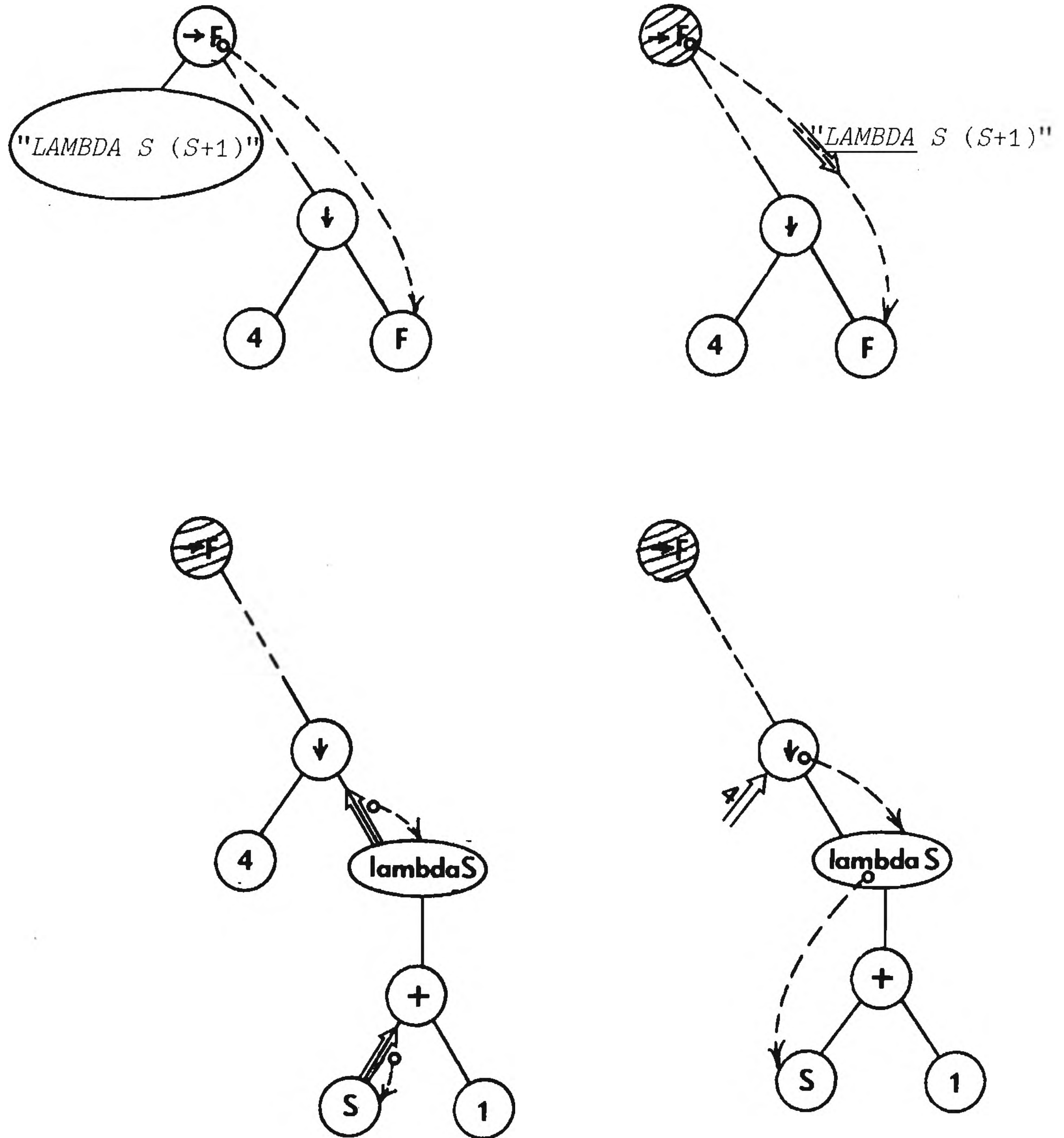


Fig. V-27

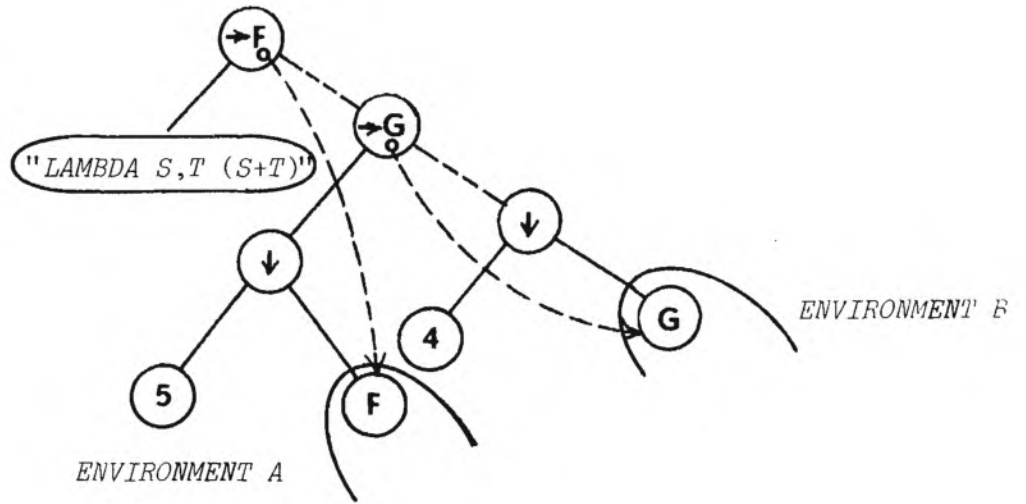


Fig. V-28

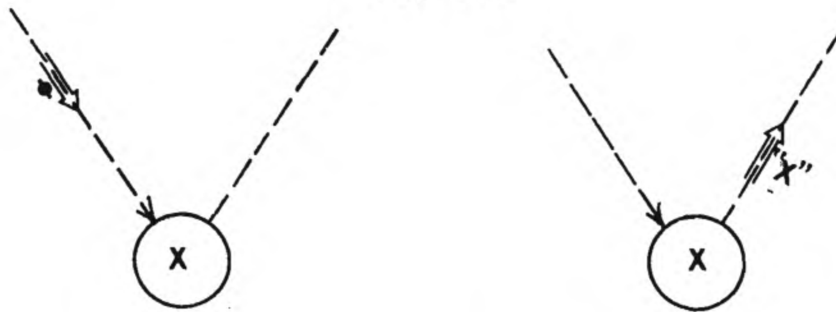


Fig. V-29

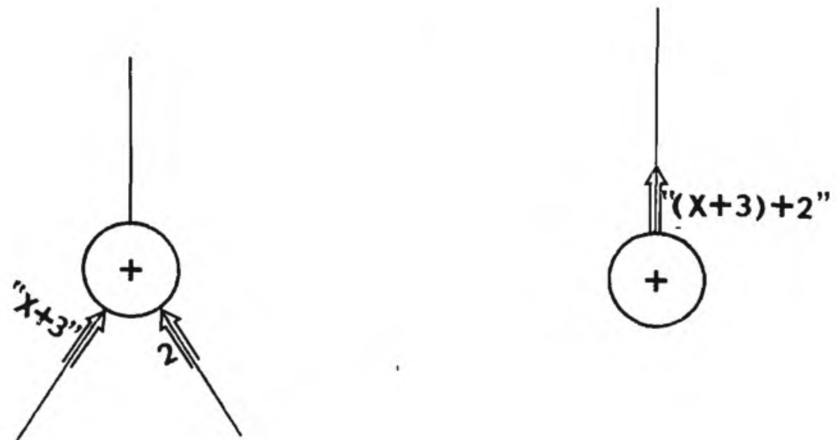


Fig. V-30

one pseudo-value representing an ob, it produces a pseudo-value which is sent up the tree: for instance, in fig. V-30 the combinator '+' receives a pseudo-value "X+3" and a value 2; it sends up the tree the pseudo-value "(X+3)+2".

An ob may have some of its variables receiving a value and some other variables receiving a pseudo-value. As a result, this ob is partly evaluated into an ob whose associated pseudo-value is produced and sent up the tree (fig. V-31).

V.4.3 Procedural arguments.

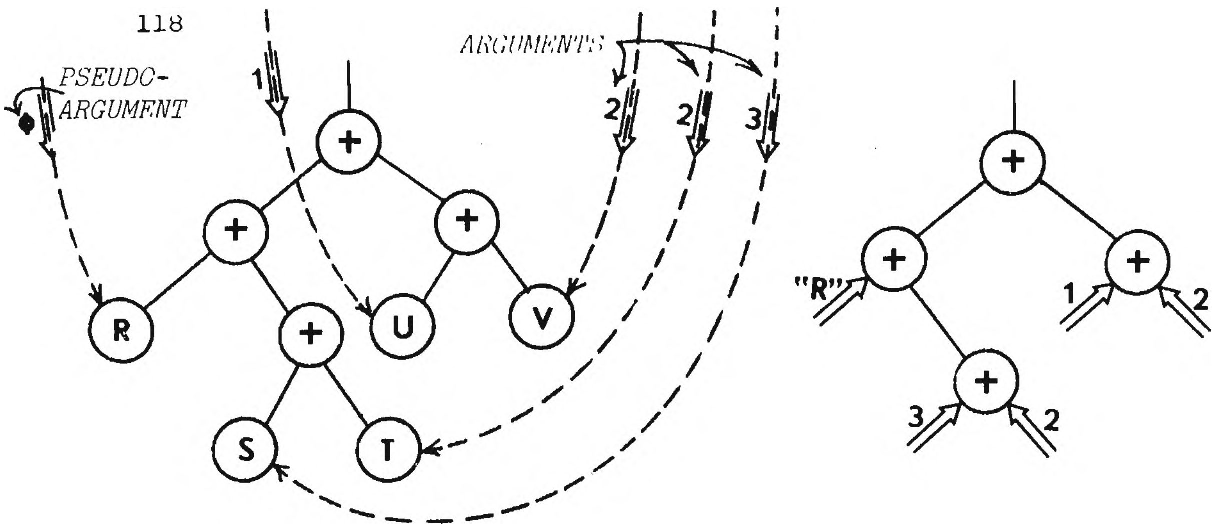
We are now able to handle the example of fig. V-28: whenever the node $\to G$ receives from the left a request sent by a lambda-node, it sends as a reply a pseudo-argument (fig. V-32). As a result the ob for which F(5) stands may be partially evaluated, the corresponding pseudo-value is produced and sent to the variable G (fig. V-33).

The problem of having procedures as arguments is not yet completely solved: we have seen how a combinator representing an operation reacts whenever it receives a pseudo-value; however we do not know yet how other combinators (for instance →,↓, ...) are to react. Let us study the two following examples:

Example 1: ([*LAMBDA* X ([*LAMBDA* Y (Y+1)]→G;G(X))]→F;F(2)) (fig. V-34).

Example 2: ([*LAMBDA* X ([*LAMBDA* Y (Y+X)]→G;G(X))]→F;F(2)) (fig. V-35).

In example 1, the component ([*LAMBDA* Y (Y+1)]→G;G(X)) may be handled as previously (fig. V-36):



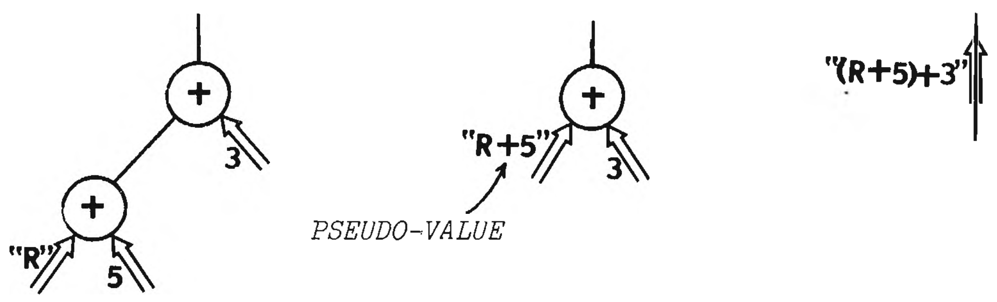


Fig. V-31

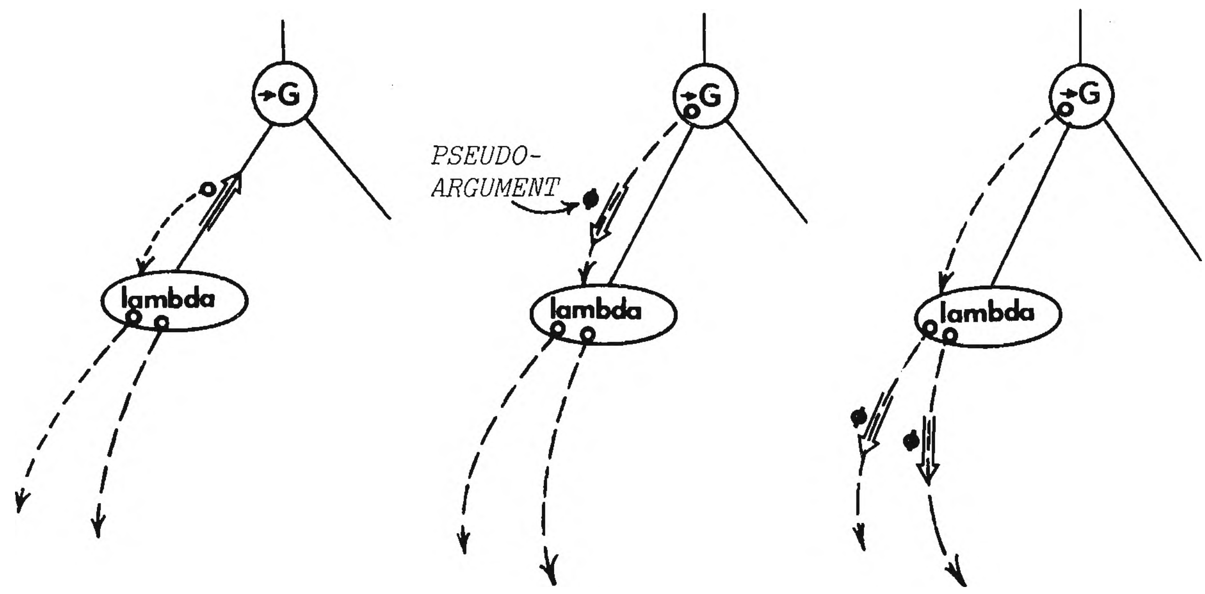
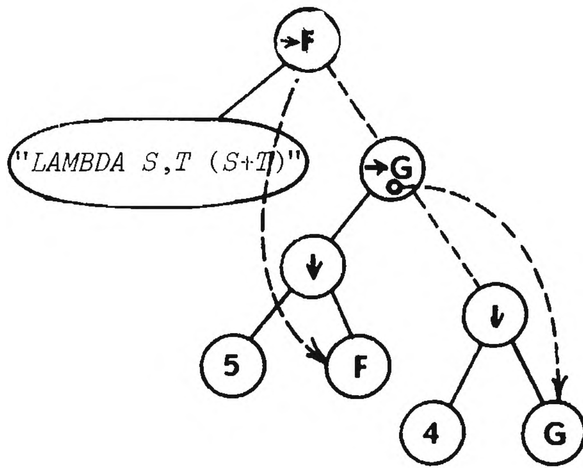
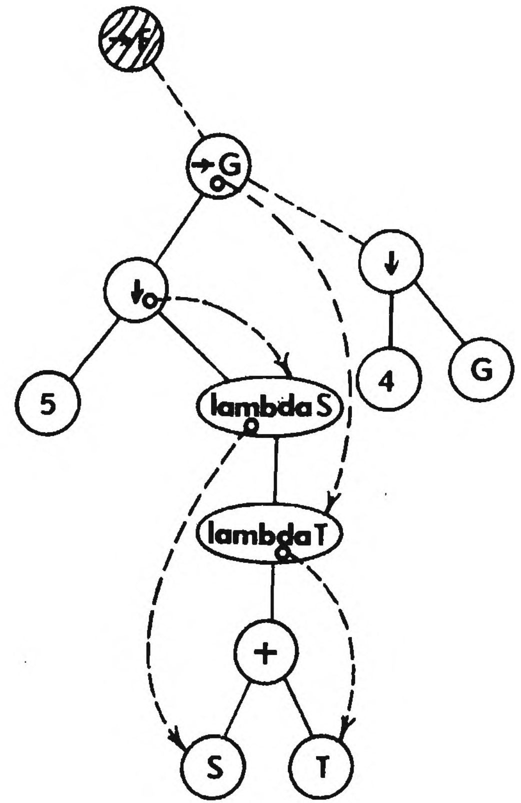
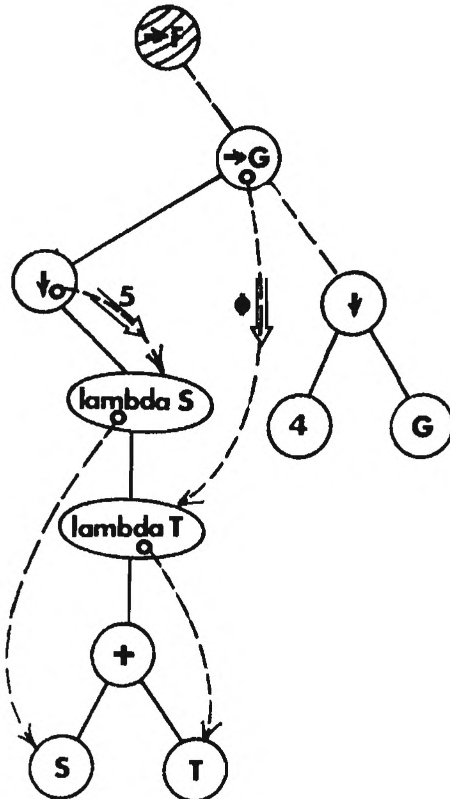


Fig. V-32







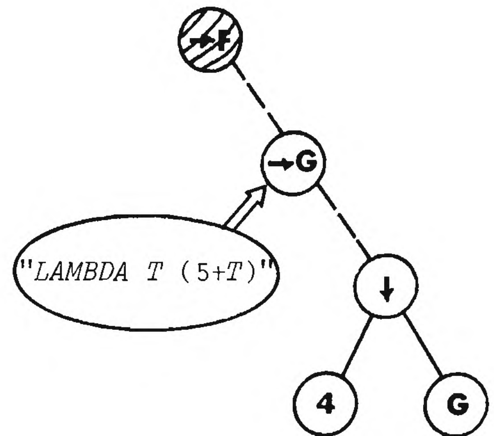


Fig. V-33

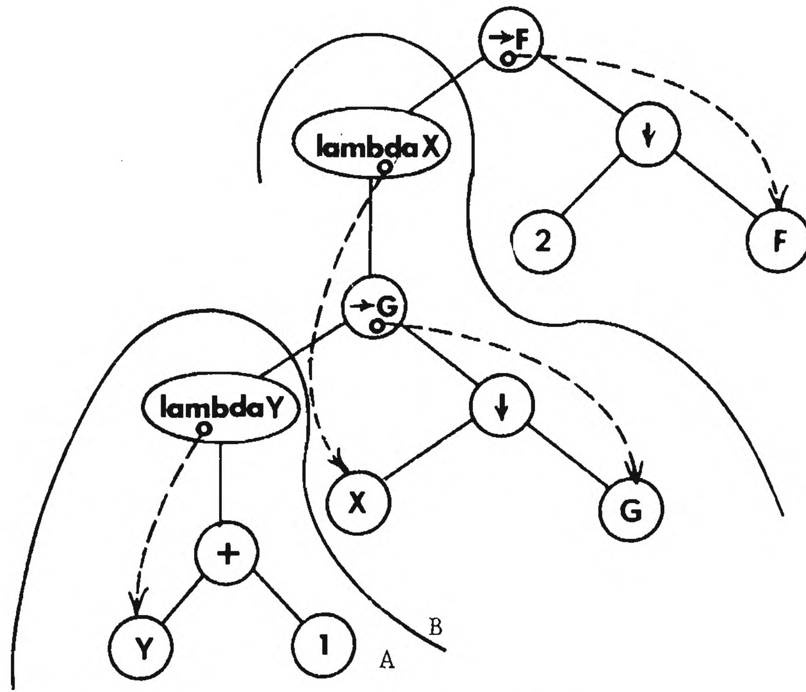


Fig. V-34

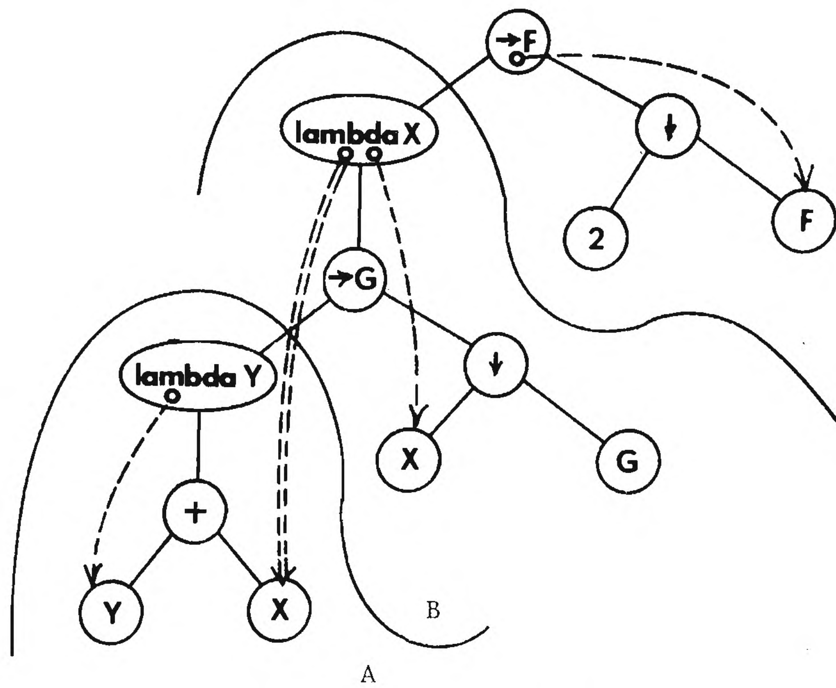


Fig. V-35

1. $\rightarrow G$ sends to the lambda-node <u>LAMBDA</u> Y a pseudo-argument ϕ.

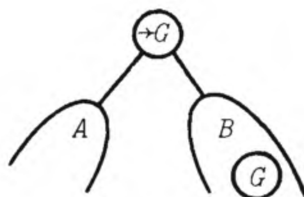
2. on receiving the pseudo-value "<u>LAMBDA</u> Y $(Y+1)$", $\rightarrow G$ passes it along to the variable-node G where the procedure is to be implemented.

The variable-node X (in $G(X)$) receives a pseudo-argument sent by $\rightarrow F$ via <u>LAMBDA</u> X ; it sends up the tree the pseudo-value "X" which is received, on the left, by the node sendown,$\downarrow$. On receiving this pseudo-value, sendown sends a pseudo-argument to the procedure on the right, and waits for a pseudo-value from this procedure (fig. V-37). Then $\downarrow$ combines the two pseudo-values into a pseudo-value it sends up the tree. The completion of the evaluation of example 1 is displayed on fig. V-38.

In example 2 the procedure <u>LAMBDA</u> Y $(Y+X)$ cannot be transmitted and implemented in the place of G for X is externally bound: it is indeed our policy to implement an argument only after its complete evaluation; we require (see fig. V-35) that the implementation of B in the place of F occurs before the implementation of A in the place of G .

As a result $\rightarrow G$ must have here another behavior: on receiving a pseudo-value on the left it sends a pseudo-argument to each variable-node it binds on the right. On receiving a pseudo-value on its right $\rightarrow G$ combines the two pseudo-values and sends the resulting pseudo-value up the tree (fig. V-39).

In summary, in the following situation



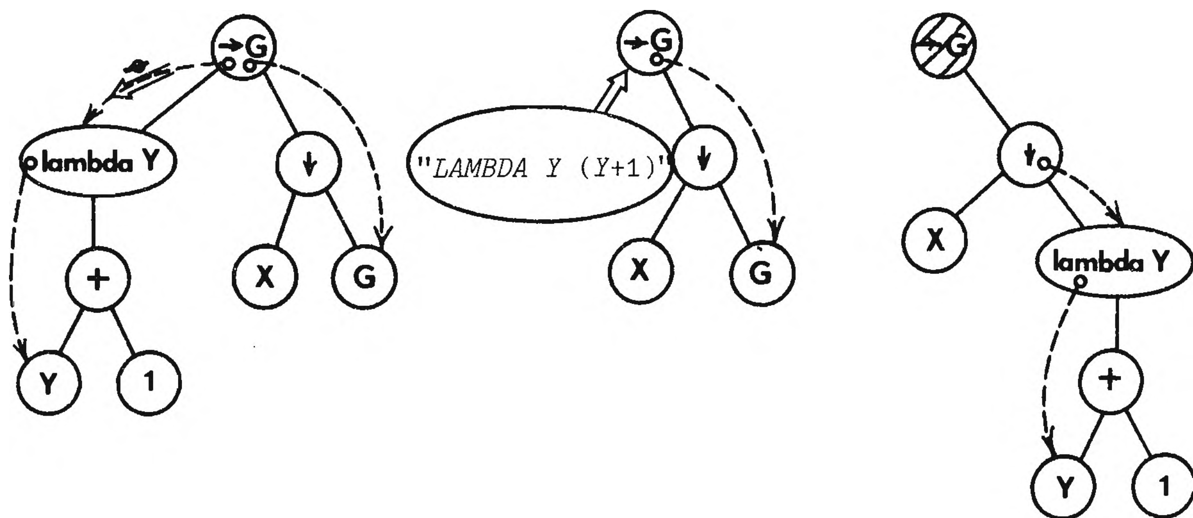


Fig. V-36

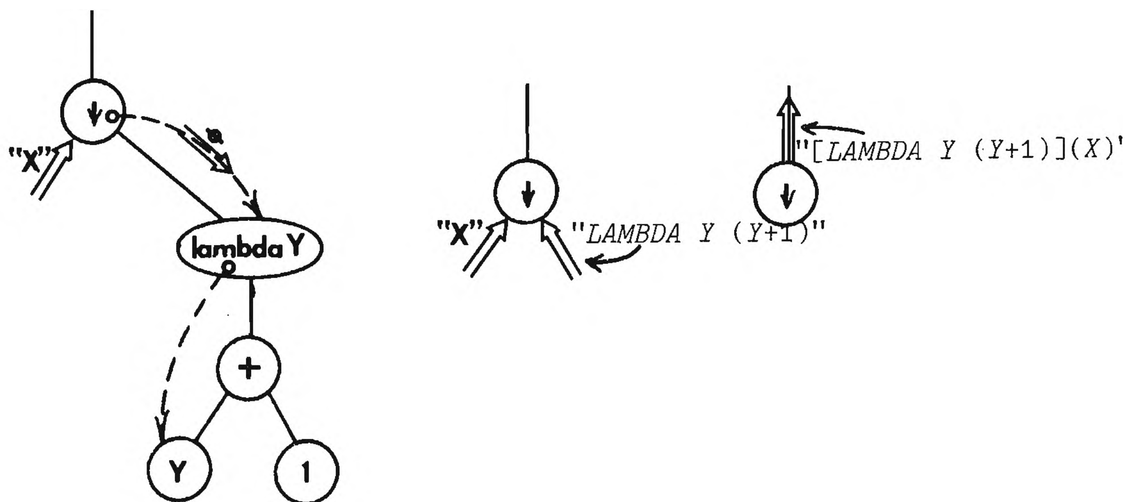
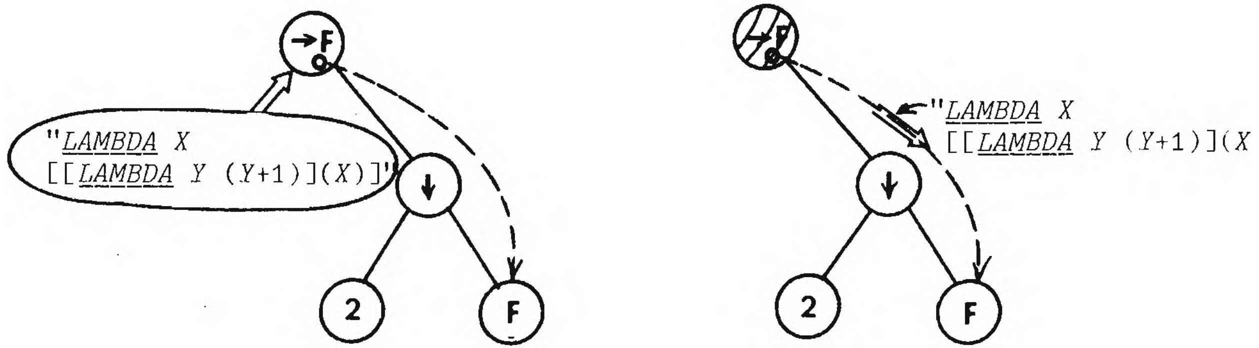


Fig. V-37



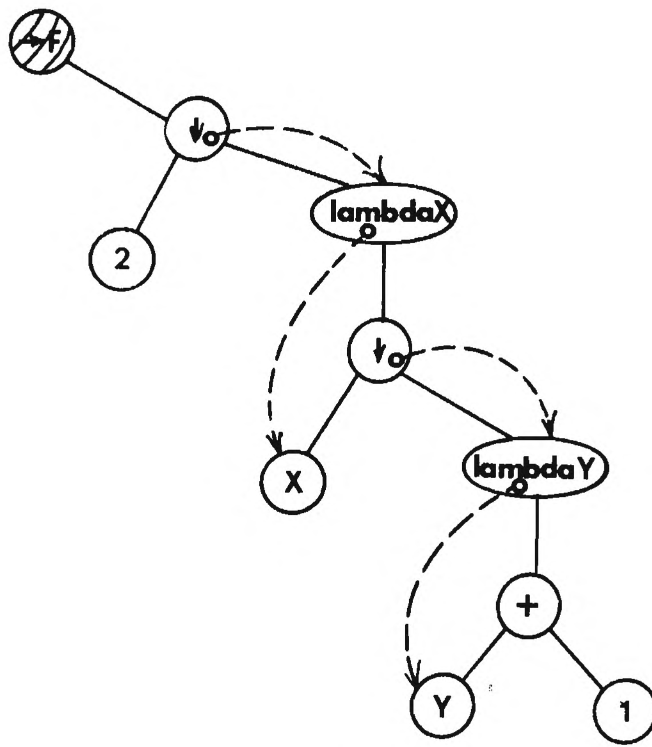


Fig. V-38

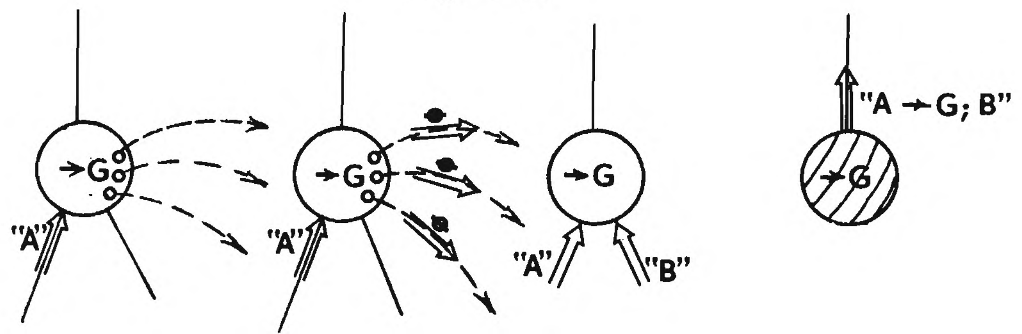


Fig. V-39

the ob *A* may or may not be to receive some argument from the environment.

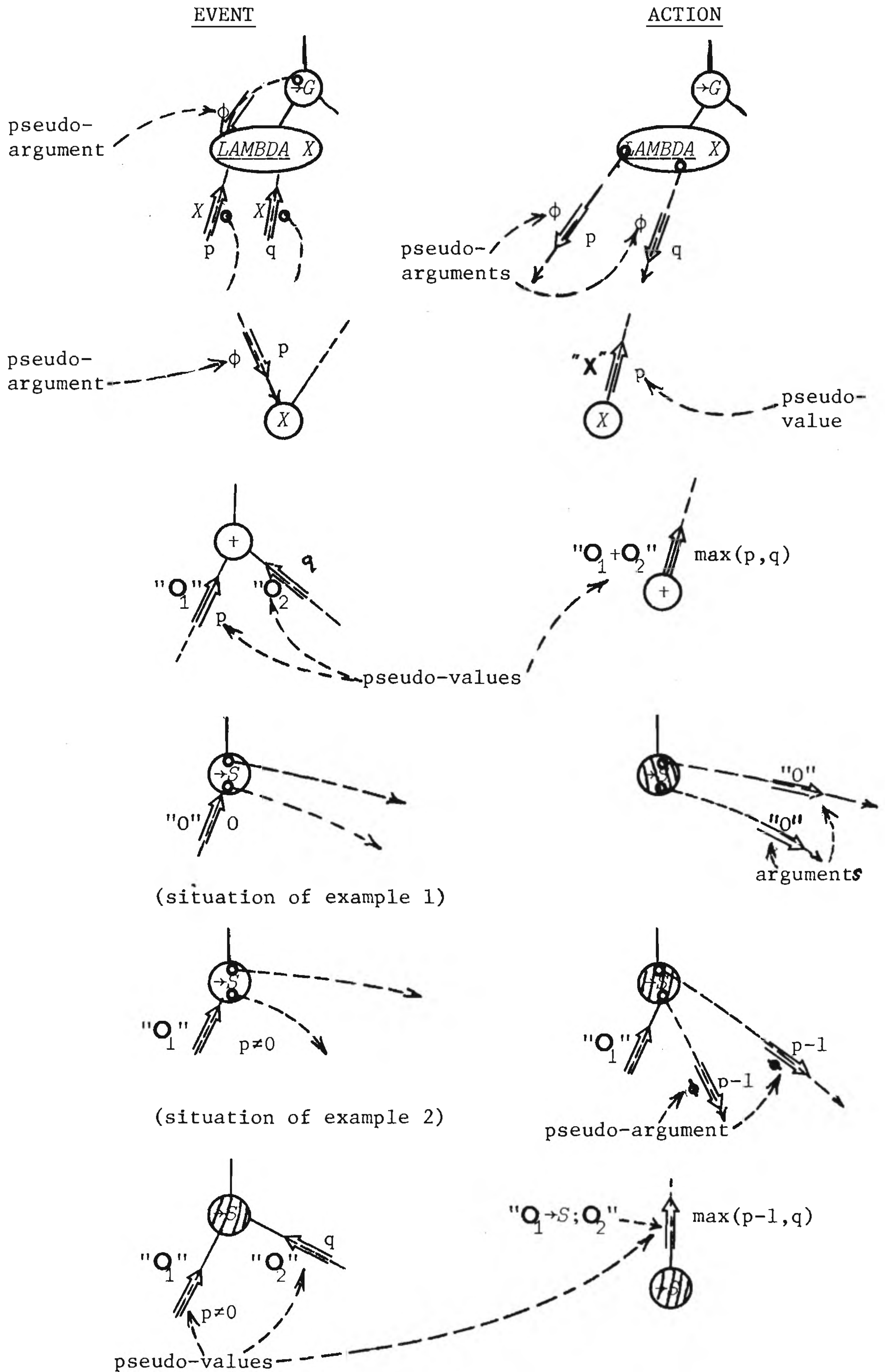
In the latter situation (fig. V-40) and if *A* does receive these arguments, *A* produces a pseudo-value "*A*", after having received a pseudo-argument ϕ from $\to G$ (fig. V-41). This pseudo-value "*A*" is sent by $\to G$ to the variable-node *G* where *A* is implemented.

If *A* receives at least one pseudo-argument from the environment (fig. V-42), the evaluation of *A* is not completed and "*A*" is not to be sent and implemented in the place of *G* . Instead, on receiving the pseudo-value "*A*", $\to G$ sends to the variable node *G* a pseudo-argument. On receiving the pseudo-value "*B*", $\to G$ sends up the tree the pseudo-value "$A\to G;B$" (fig. V-39).It is therefore necessary that $\to G$ is able to distinguish between these two situations.

We implement such a behavior in the following way:

Whenever a variable-node sends up the tree a request with its own address, such a request contains an integer called count, equal to zero at the beginning. While the request flows up the tree, the count is incremented by one every time the request reaches on the left one of the combinators setdown '→' , setup '←' , or sendown '↓' .

Such a count is associated to pseudo-arguments and pseudo-values as shown in the following table:


EVENT
ACTION
pseudo-argument
LAMBDA X
X
p
q
pseudo-arguments
pseudo-value
max(p,q)
pseudo-values
"O"
0
(situation of example 1)
arguments
p≠0
(situation of example 2)
p-1
pseudo-argument
max(p-1,q)
pseudo-values

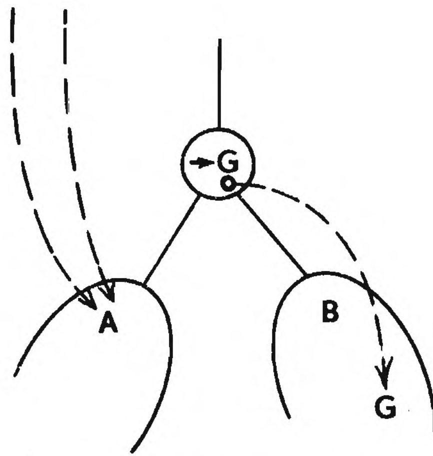


Fig. V-40

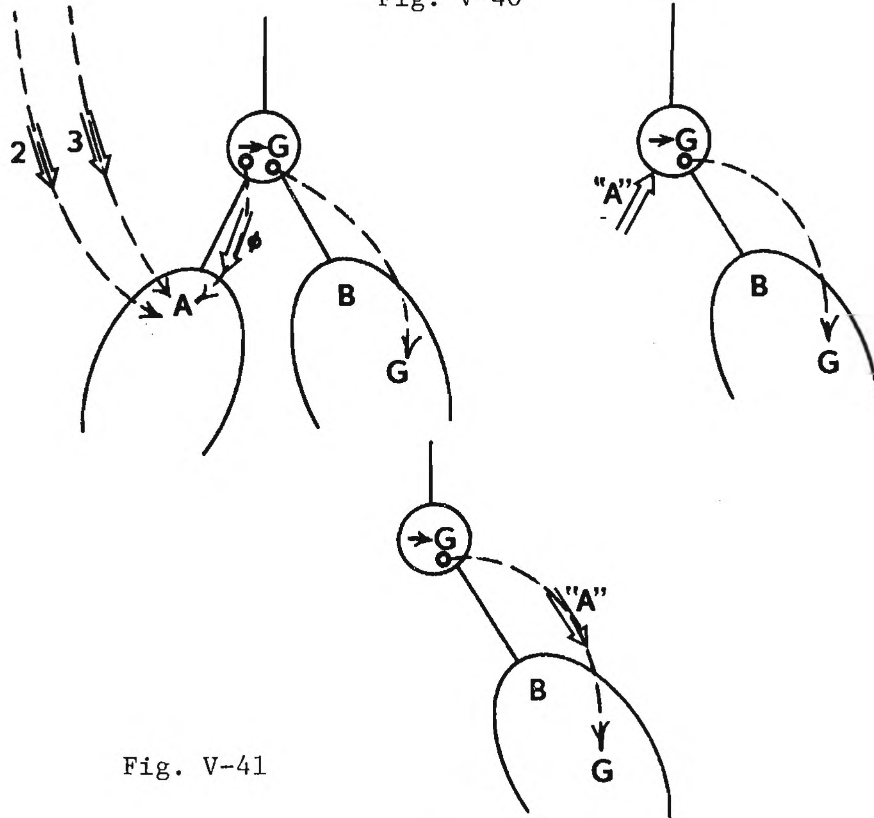


Fig. V-41

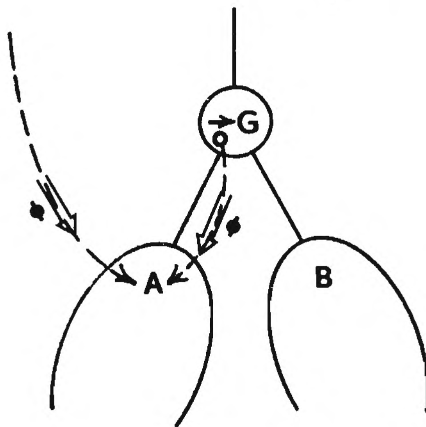


Fig. V-42

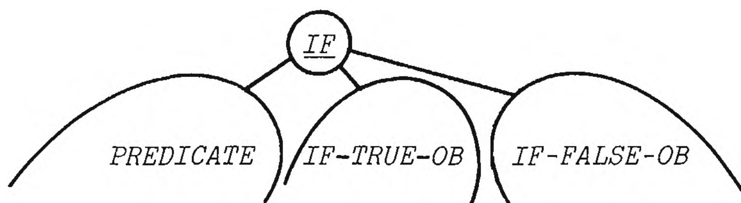
Remark.

Whenever a combinator (→S) receives from the left a pseudo-value with a count of zero, it is assured (by these behaviors) that the left ob has not received a pseudo-argument from above (→S). As a result, the pseudo-value received may be implemented (see example 1).

The behaviors of (←) and (↓) are similar to the one described for (→).

V.5 Conditional expressions.

'*IF*' (or '|') is a combinator of degree three:



The ob "predicate" is assumed to produce a logical value: *FALSE* or *TRUE* . In the former case the ob "if-true-ob" is deleted from the computation tree; in the latter case it is the ob "if-false-ob" which is deleted from the computation tree (fig. V-43a and V-43b).

As shown in section V.3, any ob may exchange information with its environment; as a result, messages can freely flow down from the environment into the obs "predicate", "if-true-ob", "if-false-ob" which may be considered as being evaluated concurrently.

However, it is important to prevent an ob which eventually would be cancelled from sending messages to its environment. For this reason, the messages which are sent to the environment by the obs "if-true-ob" and "if-false-ob" are picked up and kept in two different packages at the '*IF*'-combinator level (fig. V-44a). When the value of the ob "predicate" reaches the combinator, one of the two packages is destroyed

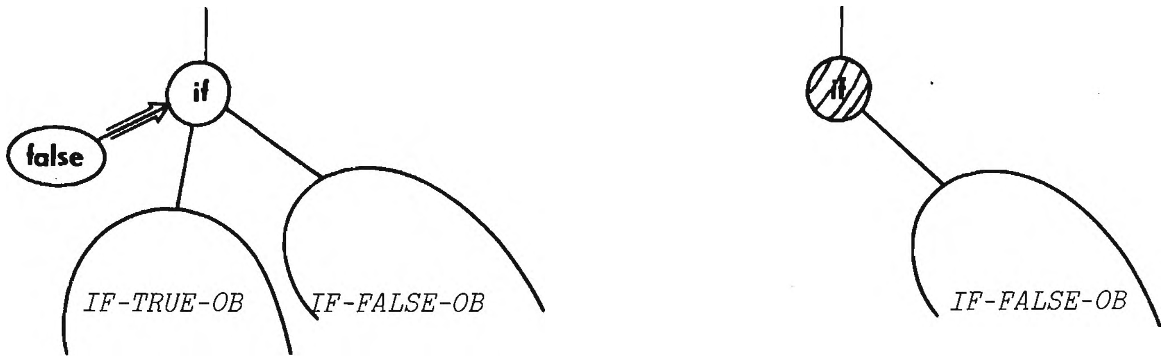


Fig. V-43a

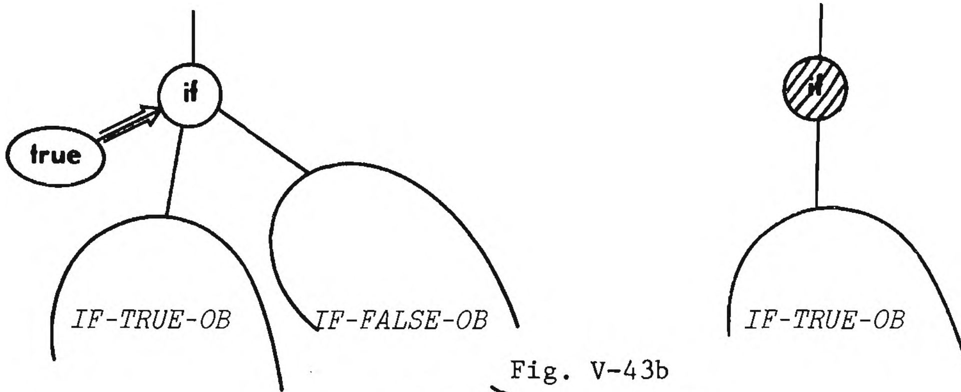


Fig. V-43b

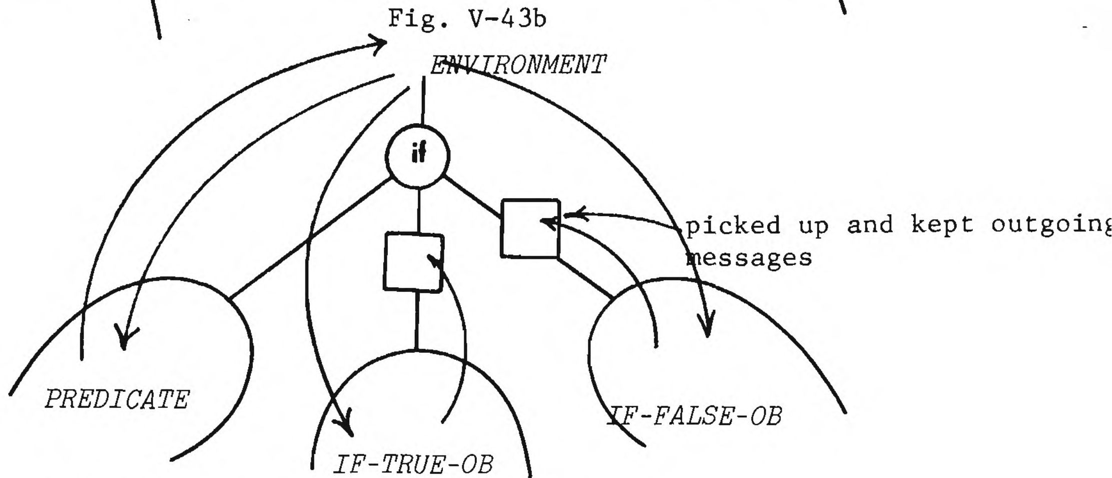


Fig. V-44a

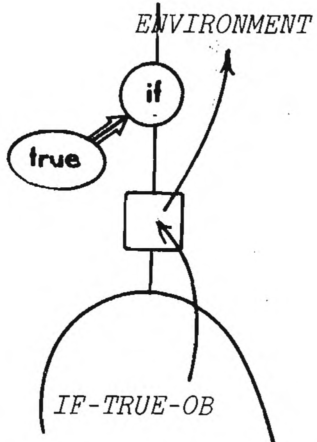


Fig. V-44b

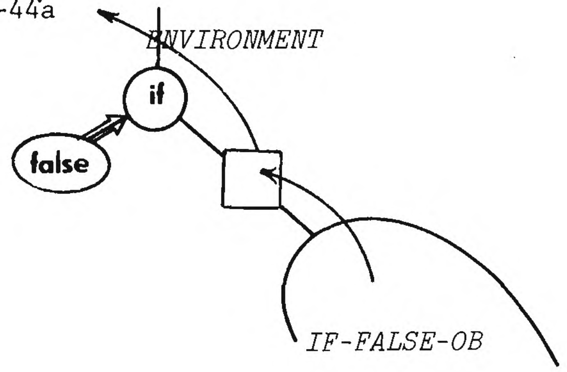


Fig. V-44c

and the messages contained in the other one are freed (fig. V-44b and V-44c).

V.6 DCPL as a system-oriented programming language.

In the previous section of this chapter, we have seen how simple DCPL programs may be considered as expressions which may be evaluated. This section sustains our claim that DCPL is a system-oriented programming language: a computation is considered a system of asynchronously cooperating "independent" programs (coroutines) linked by paths of information along which messages are sent.

V.6.1 Asynchronous events and sequential processes.

Sequential processes which are triggered by the occurrence of some asynchronous event can be embedded in DCPL. Whenever some actions are to be synchronized[a] in some way, such sequential processes may be used (these processes may be very small and perform just one elementary action).

Any value may be considered as an event whenever we are interested in knowing whether the value has been received or not, disregarding the value itself.

We may operate on events with the operators '*AND*' and '*OR*', which are not to be confused with the logical operators '$\wedge$' and '$\vee$'. The nodes which correspond to these operators consider any value they receive as an event. An *AND*-node produces an event whenever it receives an event from both of its sons. An *OR*-node produces an event whenever it receives an event from one of its two sons (fig. V-45). Thus it is possible to have an expression producing an event (fig. V-46).

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a) The word "synchronize" is here a poor choice. We mean that these actions must be safeguarded from one another as they would be with semaphores (Dijkstra [5]).

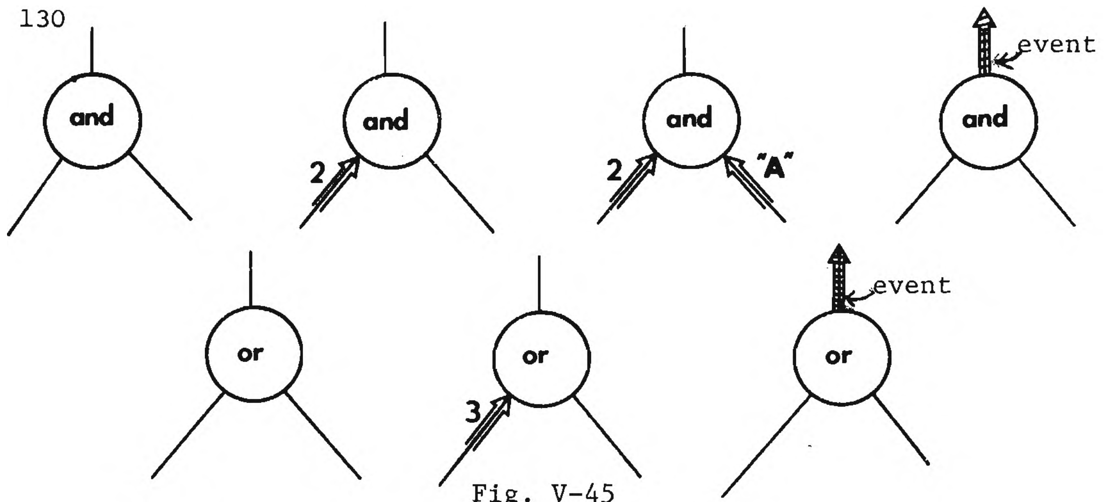
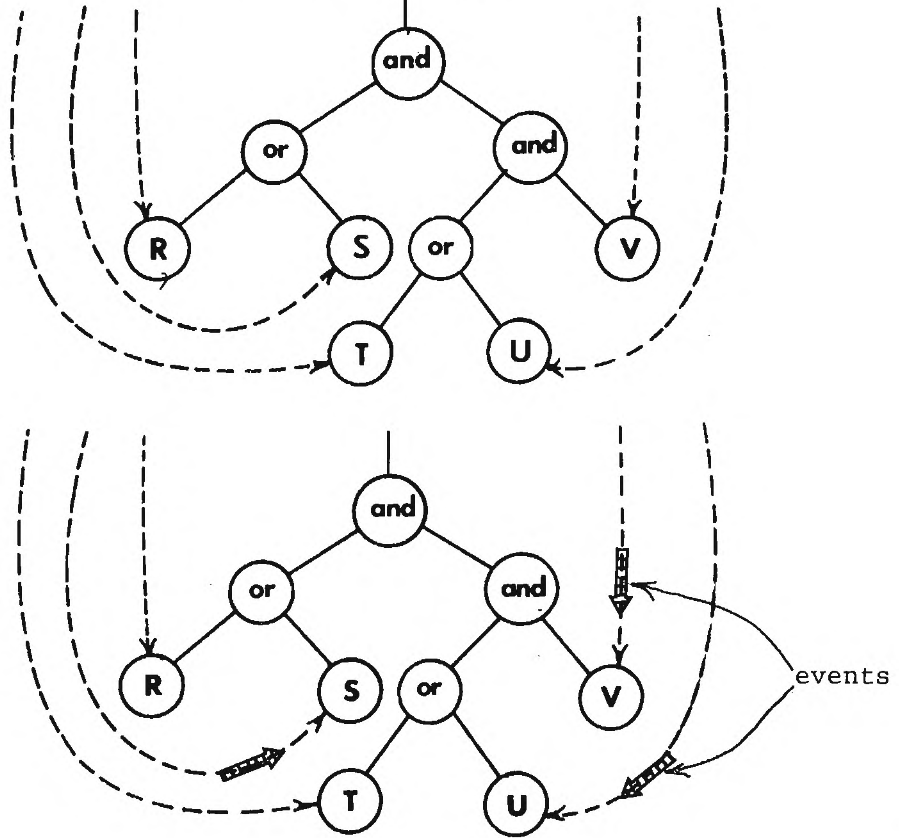


Fig. V-45



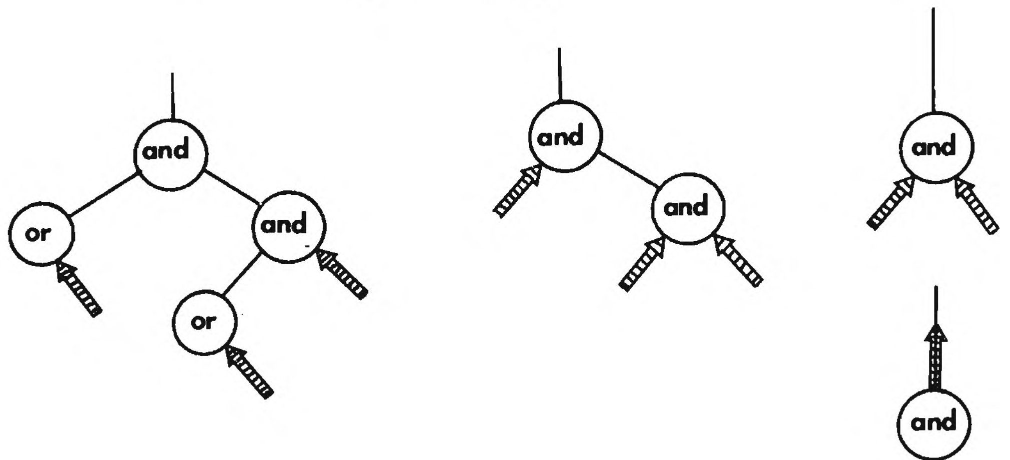
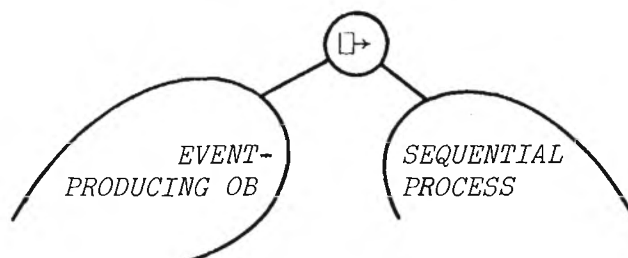


Fig. V-46

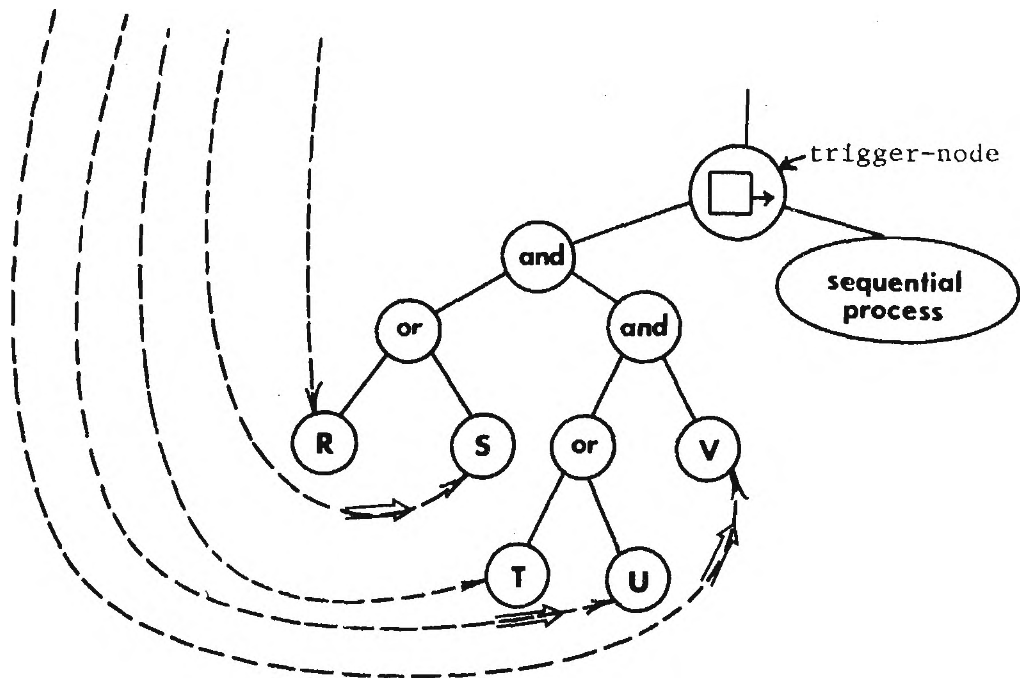
Such an expression may be used to trigger a sequential process with the combinator trigger, '⊡→':



Whenever the trigger-node receives an event from the left, it triggers the sequential process on the right (fig. V-47).

A sequential process is considered as a sequence of statements. Whenever such a sequential process is triggered, the statements are executed sequentially. A sequential process may be connected to its DCPL environment by communication paths (fig. V-48). Some statements may, when executed, send some values on these paths; some other statements may only be executed after they have received a value from such a path. If the control reaches such a statement before the arrival of the required value, the control stays pending in this statement until the value does arrive: the servicing of the communication paths in a sequential process may be viewed as some kind of input/output operations.

A sequential process may be DCPL-like; in this case the binding of variables is performed as already described (each variable-node sends a request which is picked up by the corresponding binder, etc ...). However it may be useful to embed in DCPL some subset of a usual sequential programming language; the programs in these languages being considered as sequential processes, the DCPL environment serves as a host system. In this case a special binding process



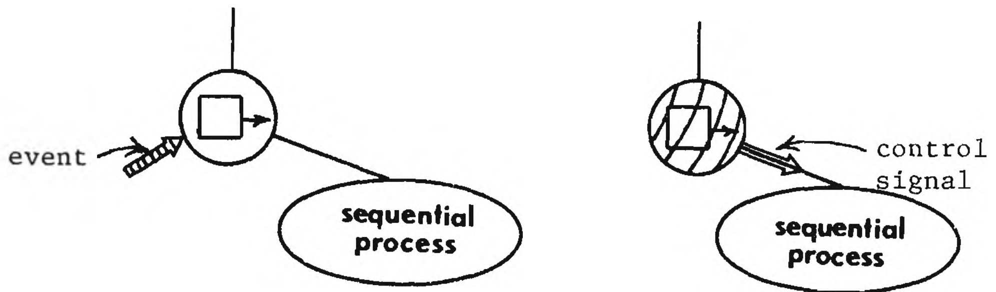


Fig. V-47

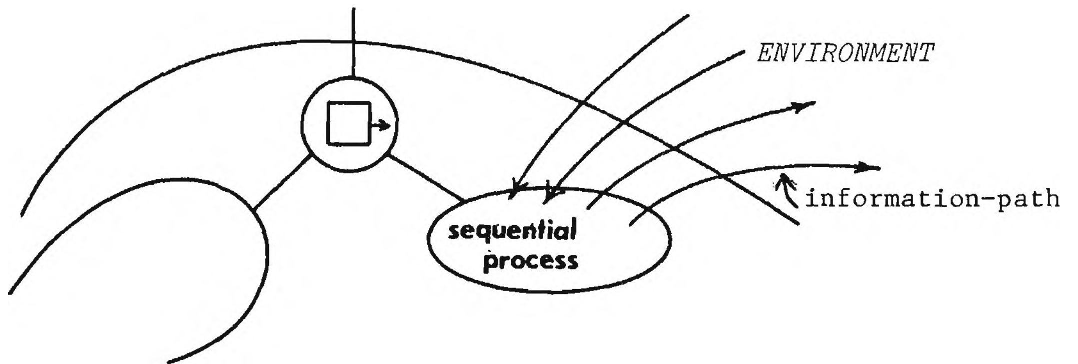


Fig. V-48

must be worked out in order to bind the sequential processes to their environments.

V.6.2 Cells.

Simple or structured cells may be used in DCPL. A cell is declared with its type: *INTEGER X* , *LOGICAL Y* etc ... . Each cell is implemented as a node in a computation tree (fig. V-49). The scope of a cell declaration is limited to the portion of the subtree under the declaration node which is not superseded by another cell declaration with the same name. In order to assure that no attempt of retrieving a value from a cell will occur before the value is stored, these actions are to be ordered and synchronized. For this reason, in DCPL a cell can only be accessed from a sequential process. There is an assignment operator ':=' to assign a value to a cell; the value is retrieved whenever the name of the cell is referred to in an expression.

V.6.3 Recursive procedures.

Let us try to implement a recursive procedure like the factorial function. For instance we may suggest for n! something like

[LAMBDA *N*[(*N*=0)|1;*N*×*FACT*(*N*-1)[a]]]

or more specifically we may suggest for 3!

("[LAMBDA *N* [(*N*=0)|1;*N*×*FACT*(*N*-1)]]"→*FACT*;*FACT*(3))

In this program, several occurrences of the ob assigned to *FACT* are to be embedded one in another (fig. V-50). At some point the

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a) '*IF*' and '|' denote the same combination.

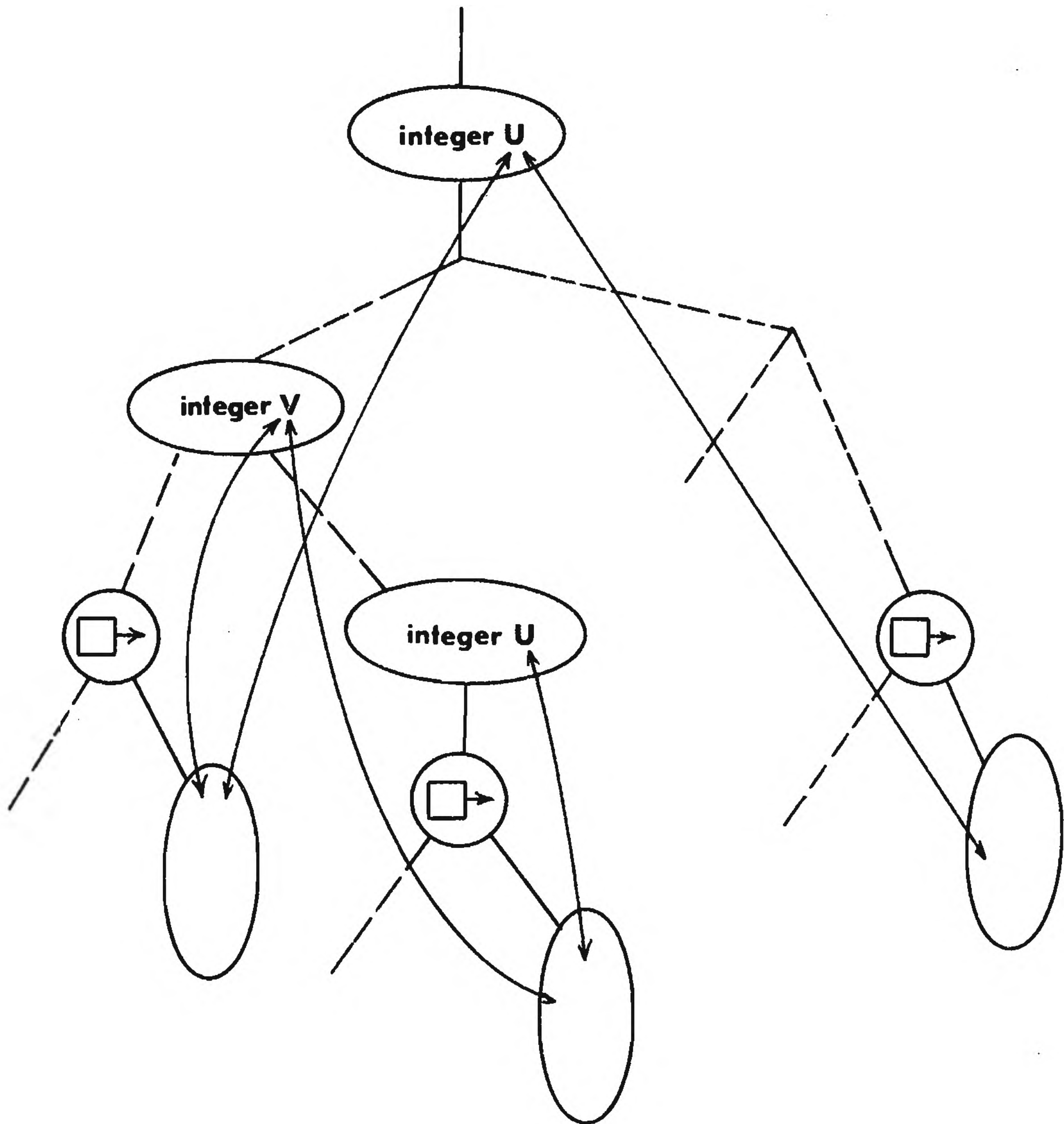


Fig. V-49

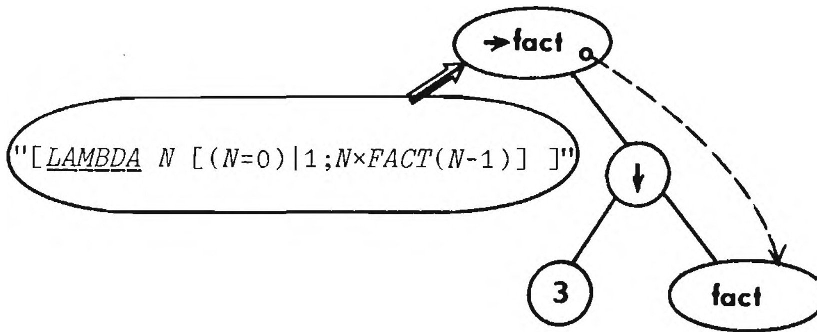


Fig. V-50a

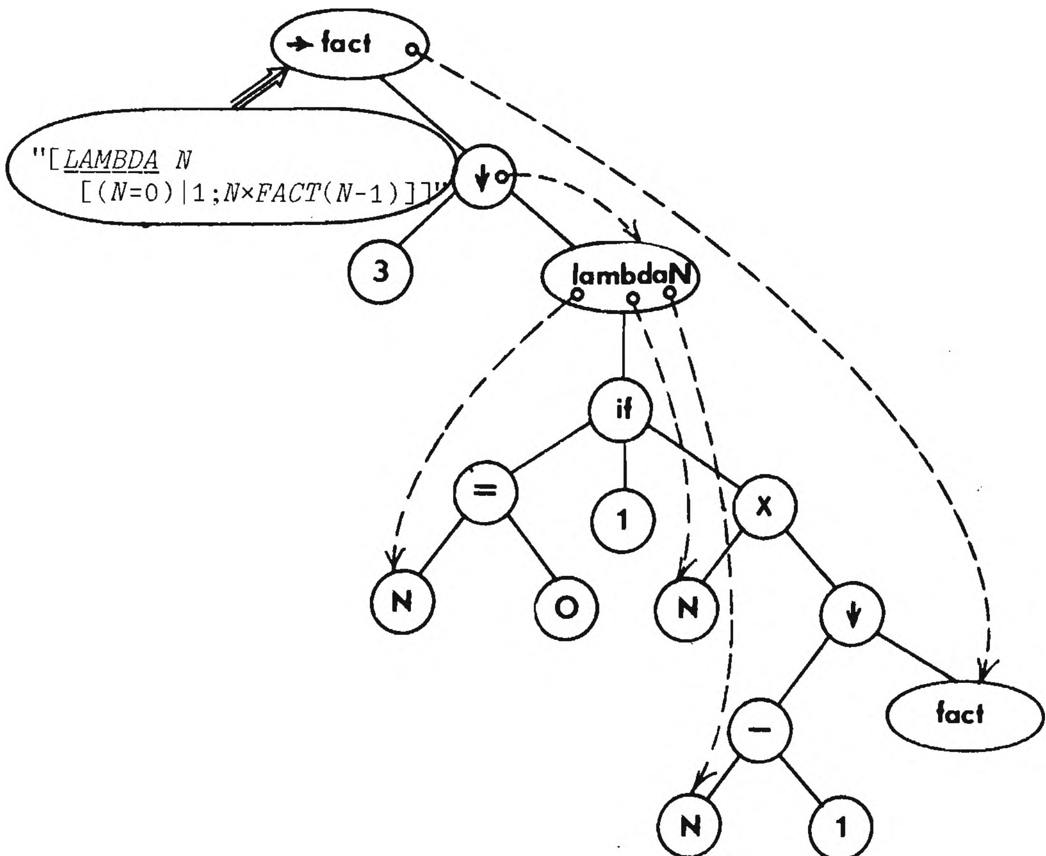


Fig. V-50b

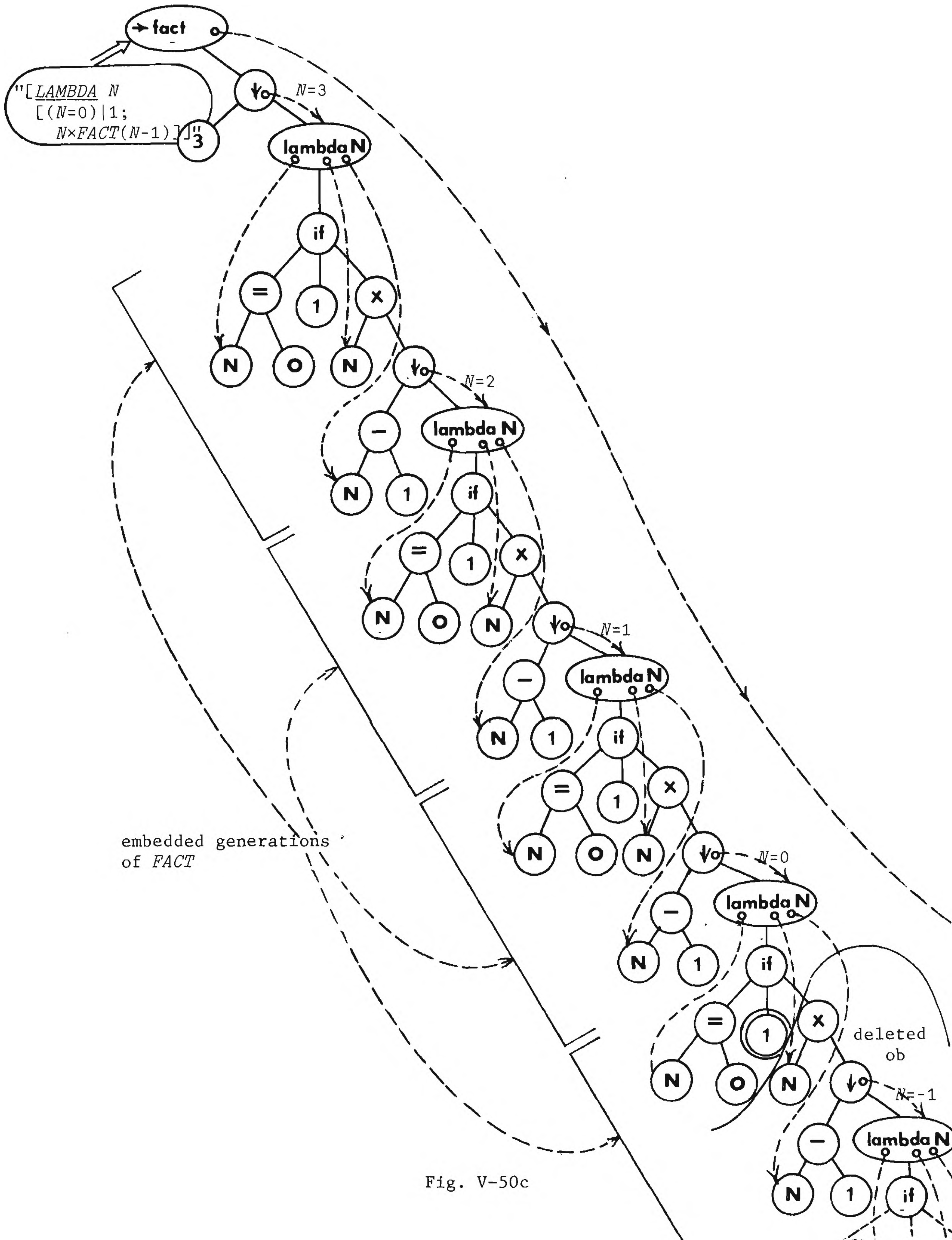


Fig. V-50c

variable *N* will be locally equal to zero and the right subtree *N*×*FACT*(*N*-1) of the corresponding '*IF*' combinator will be disregarded.

However such a program presents a major shortcoming: the implementations of the different occurrences of *FACT* are in no way synchronized with the computation itself; thousands of generations of *FACT* may be needlessly implemented.

With the combinator '*IMPL*' (implement) we can master the implementation of a procedure. '*IMPL*' stops any request reaching it on the right. Whenever it receives an event on the left, it frees the stopped requests. The variable-nodes on the right may then receive their arguments (fig. V-51).

In the program

("[*LAMBDA* *N* [(*N*=0)|1;*N*×(*N* *IMPL* *FACT*)(*N*-1)]]"→*FACT*; *FACT*(3))

the implementation of *FACT* is synchronized with the availability of the variable *N*. In the following program an implementation of *FACT* is only performed if the corresponding test produces a *FALSE* value

("[*LAMBDA* *N* *NEW* *T* [[(*N*=0)|*TRUE*;(*T*←1;*FALSE*)]|1;*N*×(*T* *IMPL* *FACT*)(*N*-1)]]"→*FACT*; *FACT*(3))

Some other programs would permit a fixed amount of "look-ahead".

V.6.4 Paths of information: alpha-variables.

Alpha-variables (as for instance α*X* , α*Y* , α*JOHN* etc...) allow the construction of alpha-paths along which an indeterminate number of arguments may be sent. As far as ob-construction and binding are concerned, alpha-variables have exactly the same behavior as variables (there is one exception, see the "cyclic behavior" subsection). However, passing an alpha-value never results in the

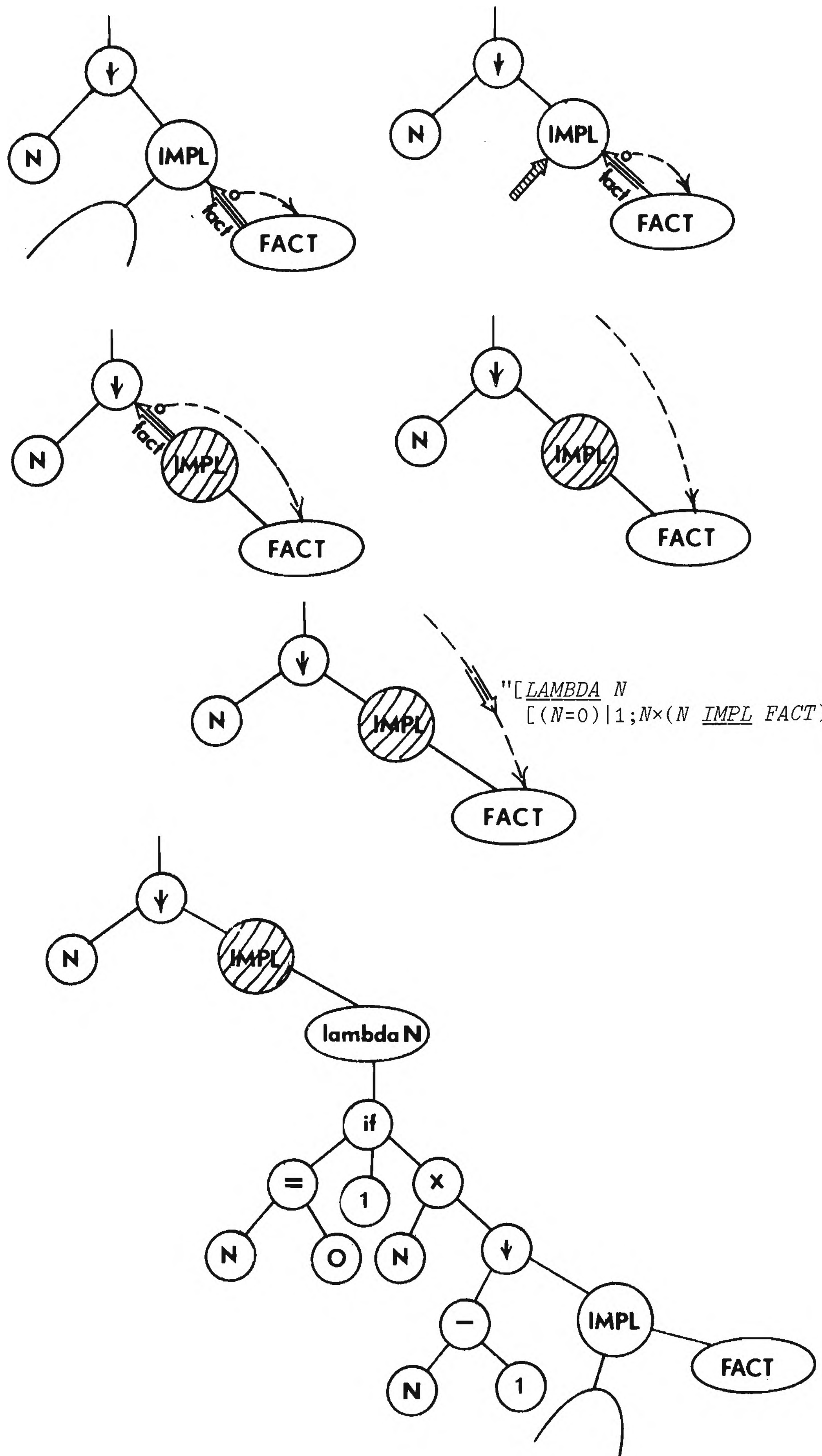


Fig. V-51

deletion of a node; as a result, alpha-variables allow a computation tree to carry out repetitively a same computation.

An alpha-variable may successively receive arguments which it passes along, up the tree, as alpha-values (fig. V-52).

An operation-node may successively receive an alpha-value on each incoming path and produces each time a resulting alpha-value (fig. V-53). If an operation-node has an incoming alpha-path and an incoming simple-path, it uses the value received on the simple path iteratively with the successive values received on the alpha-path (fig. V-54).

Cyclic behavior.

Iterative and cyclic behaviors may be modeled with alpha-paths. In particular a setdown-node may bind a setup-node constituting a recirculating path. For instance, the program

$$\underline{MU}\ \alpha U\ \ [1\rightarrow\alpha V;\ \alpha U\leftarrow\alpha V;\ \alpha V\leftarrow\alpha V+1;\ \phi]$$

is an infinite loop sending out on the path αU the successive integers starting at 1 (fig. V-55).

Interlinked coroutines.

Alpha-paths may interconnect coroutines. In the following example, the procedure *PROC*1 sends messages to procedures *PROC*2 and *PROC*3:

$\underline{NEW}\ \alpha U$ ... *PROC*1(;αU) ... *PROC*2(αU) ... *PROC*3(αU)   (fig. V-56).

In order to associate together procedures it is useful to have a particular dyadic combinator compose '∘' ; whenever compose receives a value from one of its two sons, it produces this value and deletes the other son. Our previous example may be written:

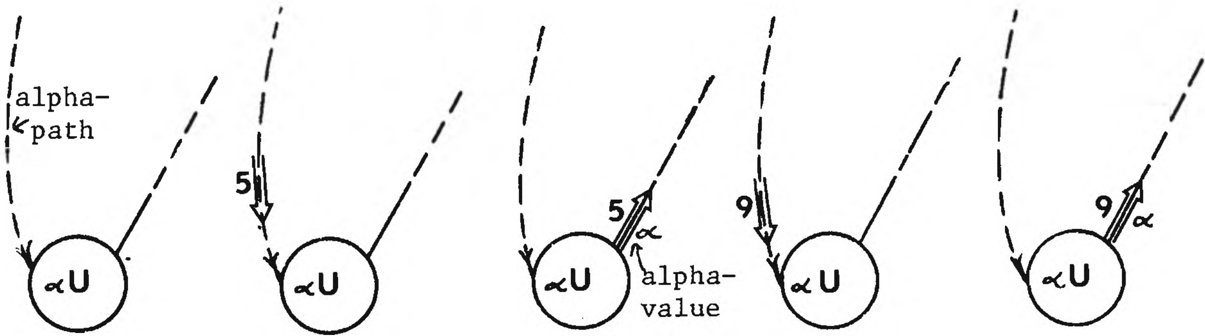


Fig. V-52

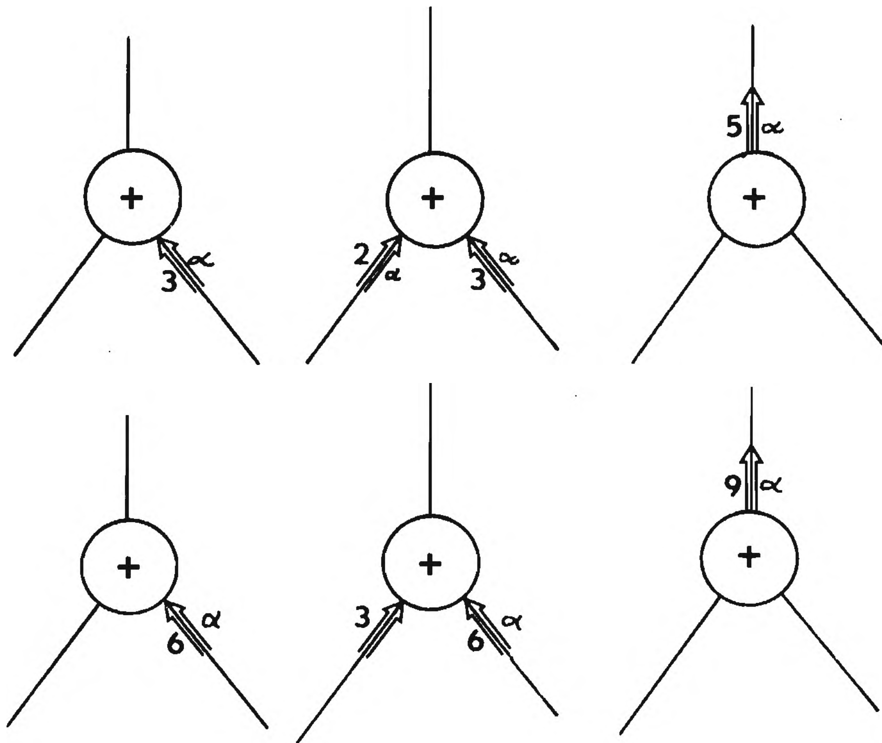


Fig. V-53

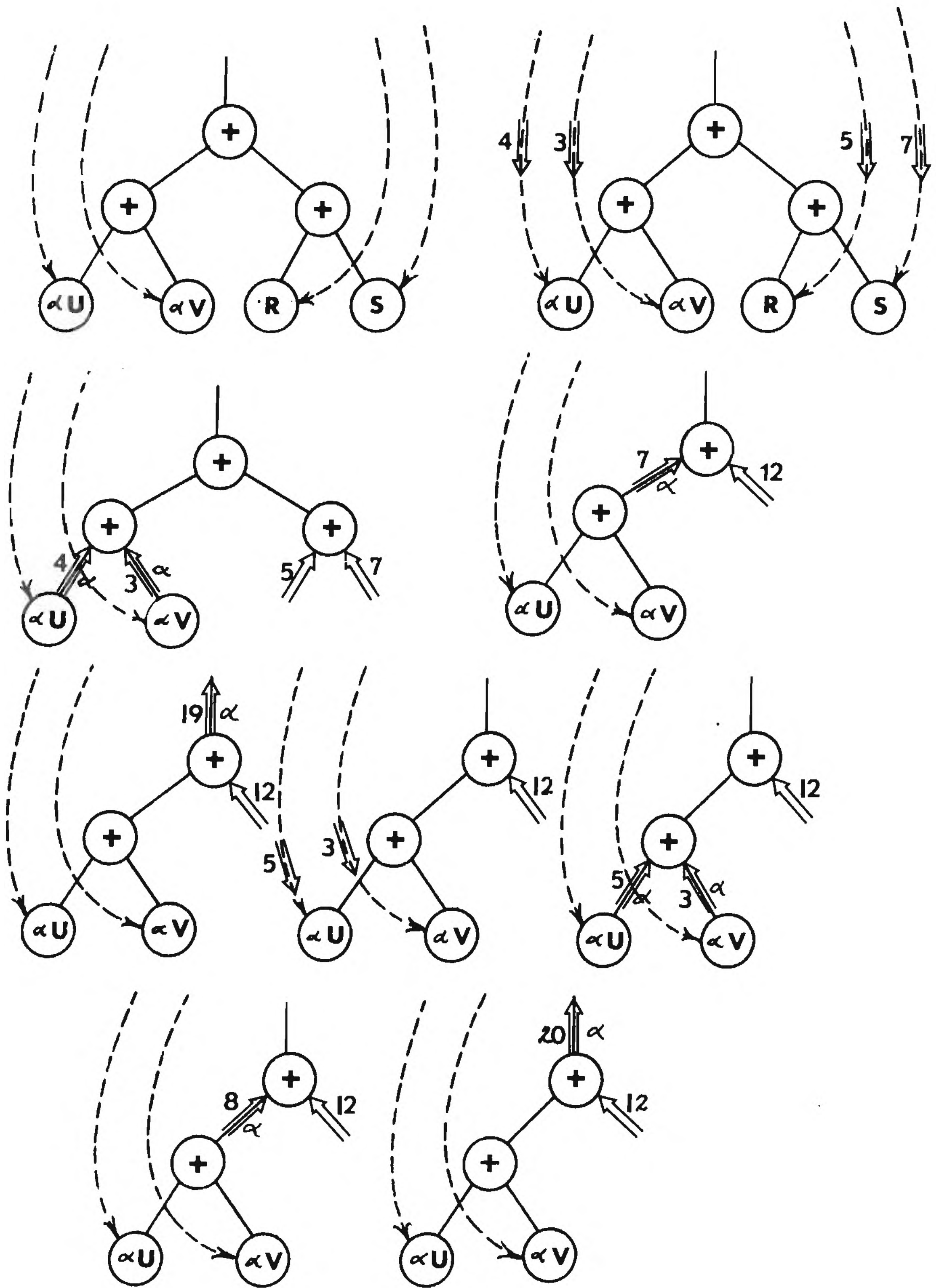
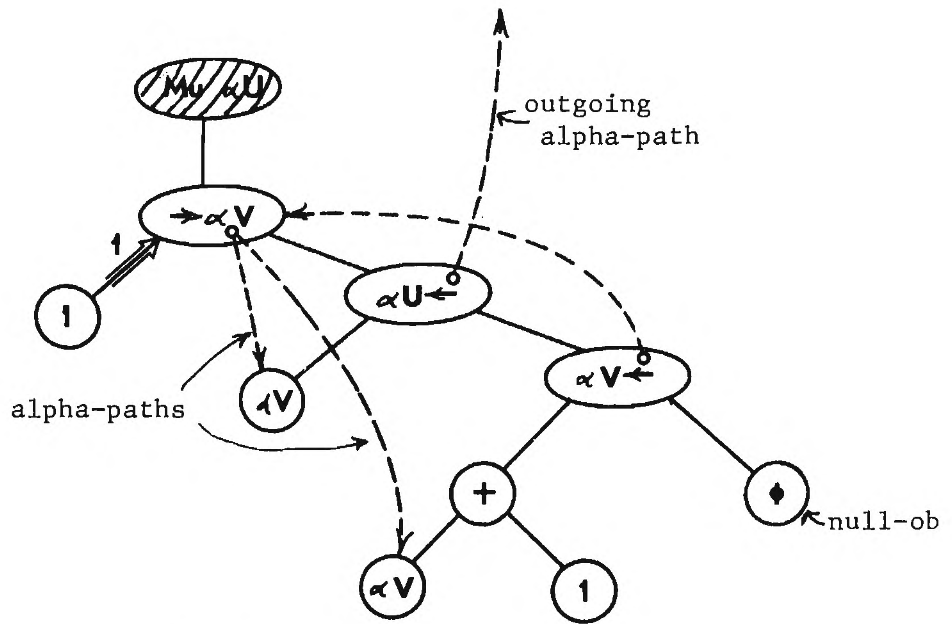
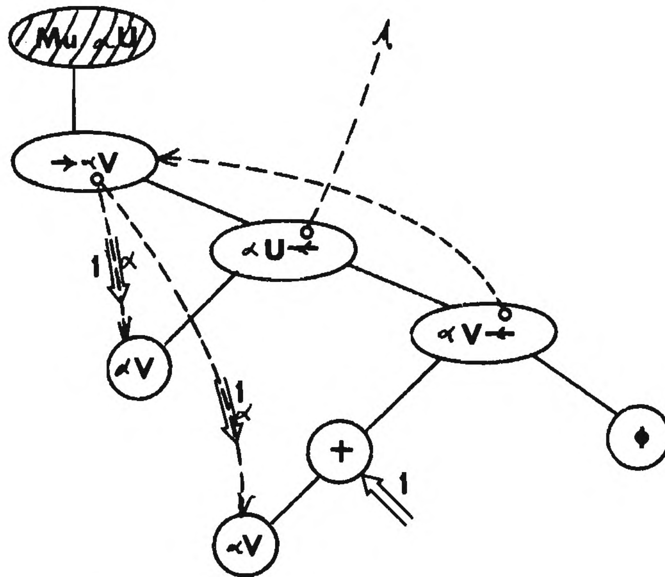


Fig. V-54





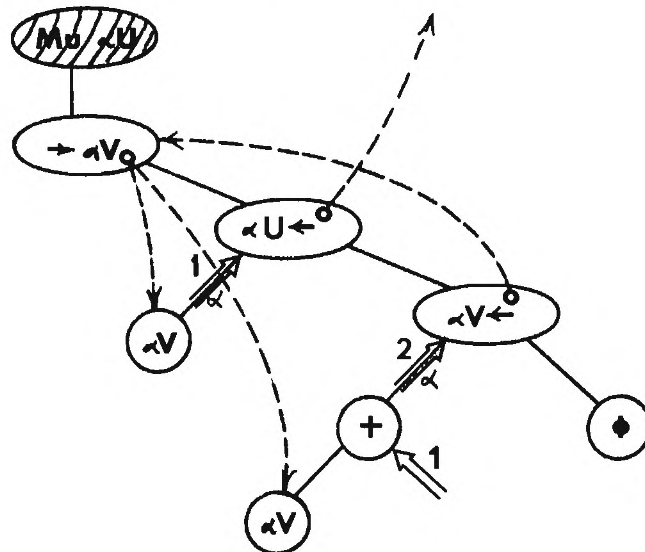
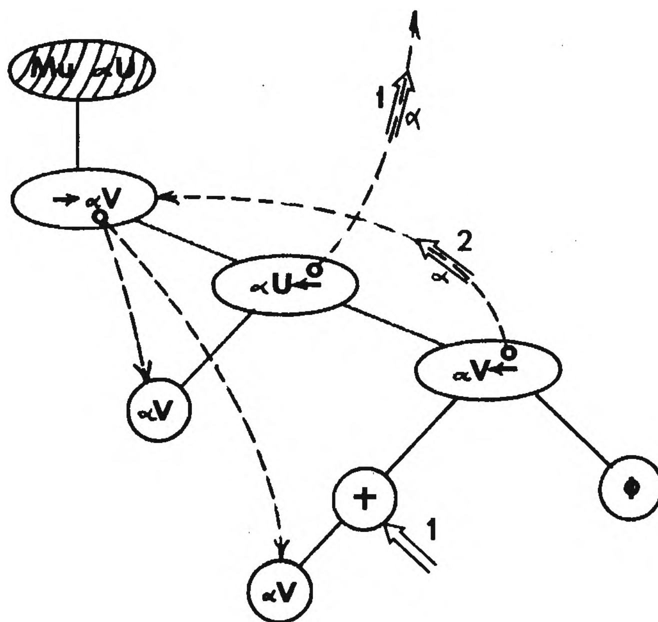


Fig. V-55a,b,c



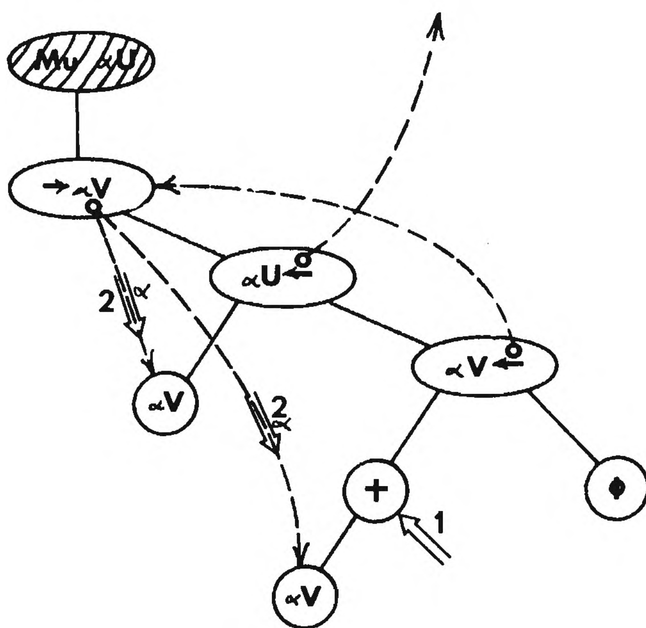


Fig. V-55d,e

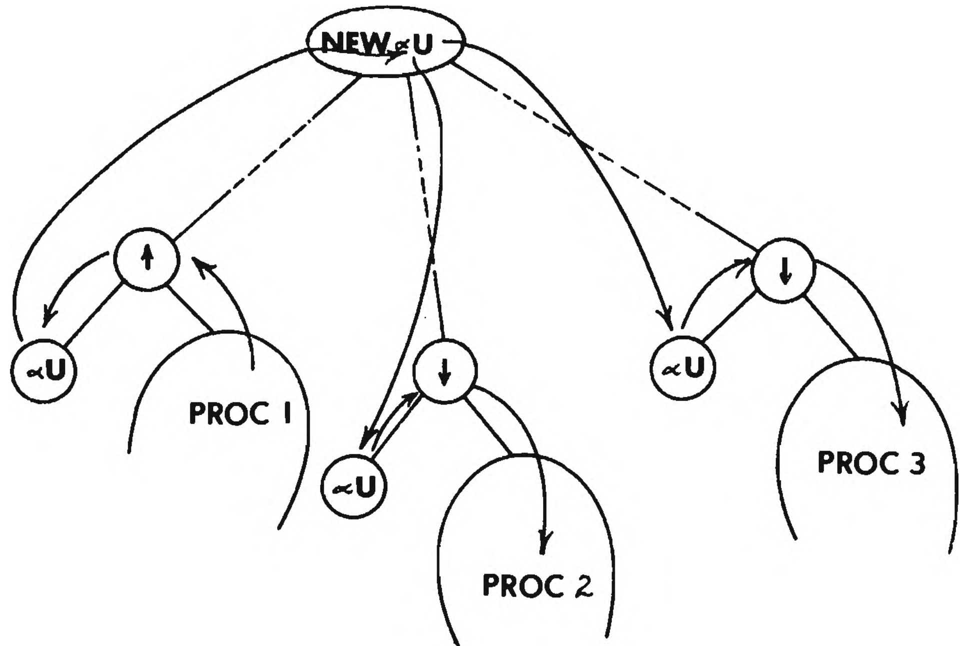


Fig. V-56

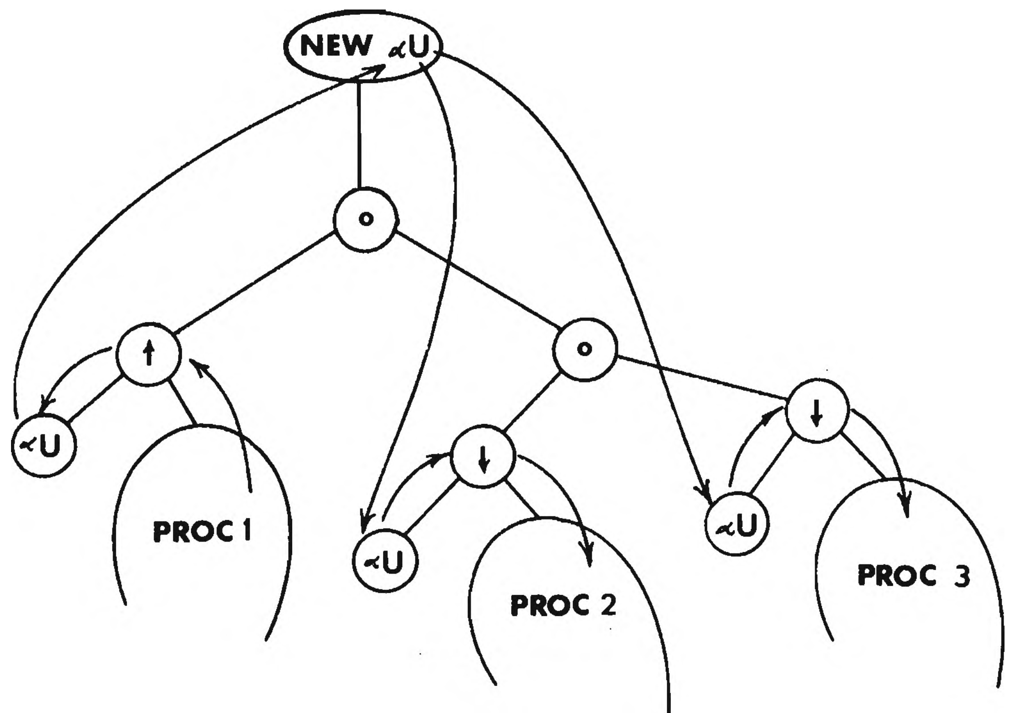


Fig. V-57

[<u>*NEW*</u> αU *PROC*1(;αU)∘*PROC*2(αU)∘*PROC*3(αU)] (fig. V-57).

In the following example, any of the procedures *PROCU*,*PROCV*, *PROCW* can send information which is duplicated and sent to the two other procedures:

[<u>*NEW*</u> αU,αV,αW *PROCU*(αV,αW;αU)∘*PROCV*(αU,αW;αV)∘*PROCW*(αU,αV;αW)]
(fig. V-58).

Whenever a procedure is interfaced with its environment, an alpha-variable in the environment may be bound with another alpha-variable or with a simple variable.

In the former case the path in the environment is connected to the path in the procedure; any number of arguments may use these paths.

In the latter case the procedure must be recursive, the path in the environment is connected successively to each generation of the recursion[a] (fig. V-59). If the path goes into the procedure the incoming arguments are queued on reaching sendown, constituting a First-In-First-Out queue. Whenever such a sendown receives a request from a lambda-node, the first element in the queue is popped out and sent as a reply to the lambda-node (fig. V-60).

<u>Queue declaration</u>.

It is possible to declare the end of a path as a queue: αU <u>*QUEUE*</u> *V*; the scope of such a declaration is the subtree under it. Whenever *V* is referred to in a sequential process, the first of the queue is popped out; it may then be forwarded to the environment of the sequential process (fig. V-61).

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a) each generation receives one argument from the path.

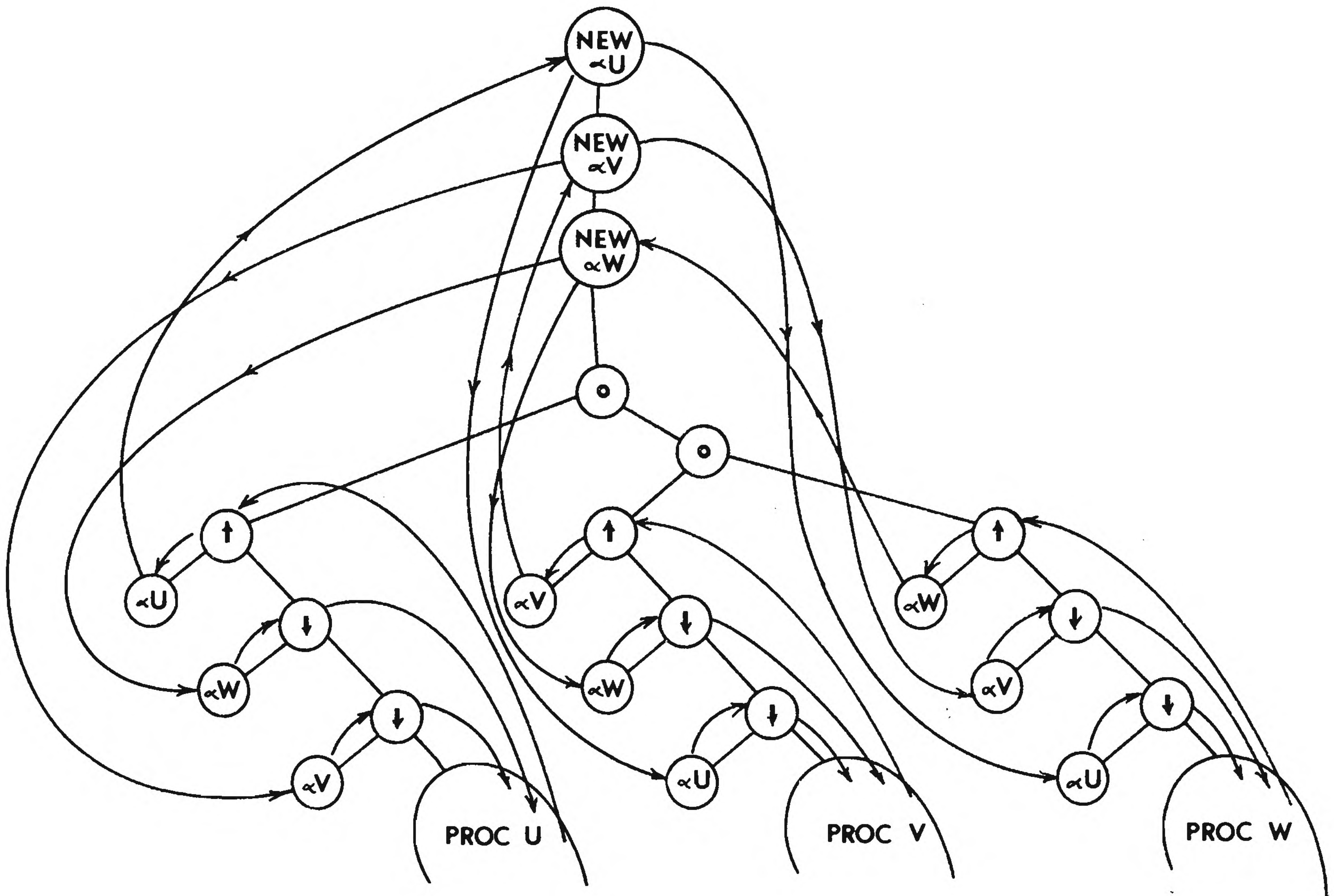

Fig. V-58

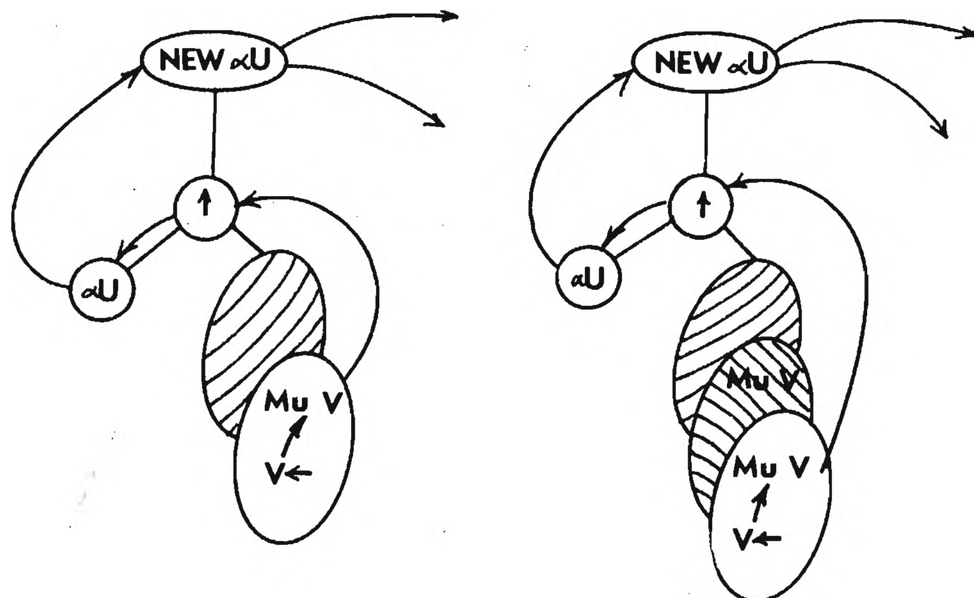

Fig. V-59

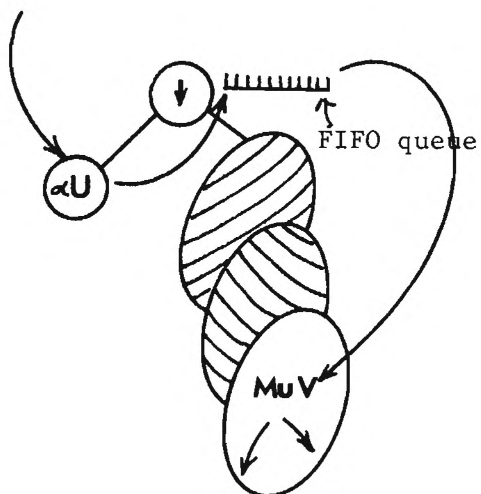

Fig. V-60

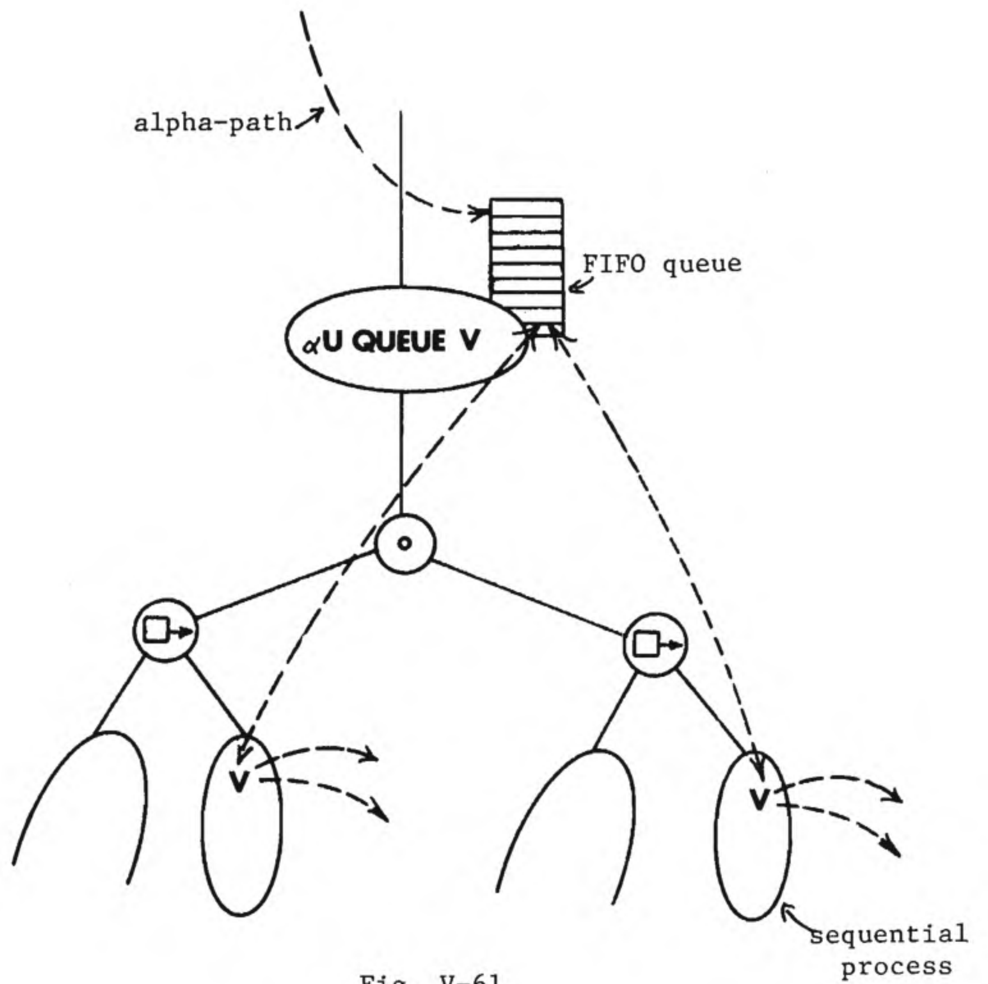

Fig. V-61

V.7 Summary.

Let us review the main notions encountered in this chapter.

1. A computation is viewed as a tree whose nodes are automata and whose edges are directed channels of information.

2. Each node has an address in a particular space; any node may send a message to any other node whose address it knows.

3. The universe of values and operations is left open-ended. To each value corresponds a value-node and to each operation an operation-node.

4. A few replacement rules allow to have a linear representation of any DCPL computation tree.

5. A binding process superimposes a graph structure to the computation tree: each variable-node sends a request (containing its address) up the tree. This request is picked up by the corresponding binder.

6. Setdown '→' allows to send a value "down the tree". Setup '←' allows to send a value "down the tree" and to the environment; its scope is delimited by the binder *NEW*.

7. A variable in a procedure may either receive an argument from the environment or send an argument to the environment. In the former case, the variable is bound inside the procedure with a *LAMBDA* binder-node and the procedure is interfaced with its environment through a node sendown '↓'.In the latter case, the variable is bound inside the procedure with a *MU* binder-node and the procedure is interfaced with its environment through a node sendup '↑'.

8. The description of any computation tree may be transmitted as a pseudo-value. Whenever a variable receives a pseudo-value,

the corresponding computation tree is implemented in the place of the variable. "X+1" is a pseudo-value describing the tree X+1.

9. A procedure may be modified through interaction with its environment, and then be transmitted as a pseudo-value in order to be implemented in some other place.

10. The different parts of a conditional expression are evaluated concurrently. However, as long as the result of the test is not available, the outgoing messages of the alternate subtrees are picked up and retained.

11. With the nodes *AND* and *OR* it is possible to have an expression produce an event.

12. '□→' , trigger, is used to trigger sequential processes, and *IMPL* to master the synchronization of the implementation of the different generations of a recursive procedure.

13. Cells may be declared; they may only be accessed from a sequential process.

14. Alpha-variables allow the construction of alpha-paths on which an indeterminate number of arguments can be sent. They permit cyclic behavior, interlinked coroutines and implicit or explicit FIFO queue servicing.

It should be noted that the concepts described here do not constitute a set of primitive concepts.

Moreover, the behavior of the nodes might be modelled in very different ways. We consider the general notions discussed as more important than the described schemes themselves.

PART THREE

MACHINE ORGANIZATION

CHAPTER VI

IMPLEMENTING DCPL : MACHINE ORGANIZATION

VI.1 Introduction.

Any machine which is intended to actualize and to carry out a DCPL program must respect and be able to express the following properties:

1. A computation in DCPL is represented as a tree whose nodes are automata and whose edges are directed channels of information.
2. Each node has an address in a certain address space; any node may send a message to any other node whose address it knows.
3. During its life, a computation tree evolves, growing and shrinking.

The third requirement may appear to be the most stringent one. We would like to have some host physical structure which can grow and shrink as the guest computation it contains evolves, synthesizing itself from the components available in the surrounding milieu, like a DNA molecule governs its own reproduction. In [24], Von Neumann used a similar analogy, comparing a computing process to a self-reproducing organism. He used as supporting structure a cellular space. Such an approach for designing a machine for DCPL may be valid from a theoretical point of view, however as a result of

concentration phenomena[a], it would turn out to be too impractical.

Different kinds of machine organization are described in the sequel. In particular it is shown how large transfer rate sequential memories may be used. Moreover an organization allowing swapping in advance is discussed.

VI.2 Cellular structure using busses as communication paths.

DCPL computation trees may be implemented in the supporting physical structure shown in fig. VI-1: autonomous "active" cells may communicate to one another through a system of busses; each cell has an address and may send a message via a bus to any other cell. Each node of the computation tree is actualized by one or several cells (the amount of storage a node may need is not bound: there is indeed no limit to the number of pointers a node may contain). The edges of the tree are actualized with pointers.
At any time there are cells which are free: they do not participate to any computation tree. Whenever a computation tree grows (or shrinks), free cells are allocated to that tree (or deallocated and freed). If the busses may be used concurrently, such an organization allows concurrency and distribution of control. However, since every cell is "active", such an organization would be prohibitively expensive if large computation trees were to be carried out. Moreover, since there is a lack of locality in such an organi-

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a) As any reproducing organism, trees may grow exponentially: their embedding in a n-cellular space may therefore be problematic.

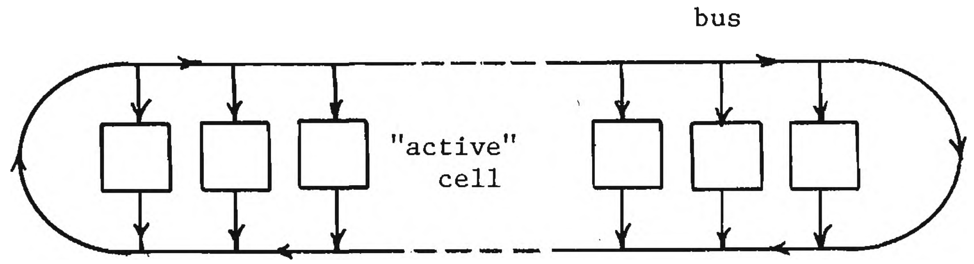

Fig. VI-1

Fig. VI-2

zation, such a communication system would probably clog rapidly whenever the size of a computation would grow. In the organization displayed in fig. VI-2, we can take advantage of the locality a program may present: each cell has a hierarchy of neighborhoods. However, with this organization the cell-allocation problem may become more difficult. It should be noted that such a cellular organization with busses may be very attractive if used as the fastest level of a "memory" hierarchy: the main computation tree is implemented in some other type of organization, small parts of a computation with many local interactions being swapped into the cellular structure to be carried out.

VI.3 A machine organization with the nodes stored on a random access storage medium.

In the previous section, a node was actualized by an "active" cell. This is a straightforward approach when considering a node as an automaton. However, we may view a node behavior in the following way: a node is quiescent until it receives a message; the interaction of a node with a message addressed to it results in a new quiescent state of the node and possibly messages for some other nodes.[a] From this interpretation derives the following machine organization:

The nodes are placed on a storage medium, each node having a physical address on this storage. At any time, the addresses of the nodes for which some messages are waiting form a queue to which

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a) Whenever a node acts spontaneously in DCPL, this action occurs right after the creation of the node; it may then be considered that a correspondent message is addressed to the node at its creation.

is associated the set of corresponding messages. One or several processors may access the nodes on the storage medium, the queue of addresses and the corresponding set of messages. Any processor may service the queue: it takes the first message, accesses the node to which the message is addressed, "computes" the interaction, stores the new content of the node, and places at the end of the queues the addresses of the nodes to which the resulting messages (if any) are to be sent.

Fig. VI-3 displays such an organization with one processor; in fig. VI-4 several processors may access different parts of the storage medium; to each processor is associated a queue; any processor may access the queue associated to the other processors and place messages in it.

VI.4 A machine organization with a sequential rotative memory.

The computation trees may be implemented on a sequential rotative memory. For instance, in fig. VI-5, a processor controls a read/write head; the addresses of the node for which some messages are waiting, and the corresponding messages, are queued; however, such a queue is not serviced on a FIFO basis: the addresses in the queue are ordered according to the order of appearance of the corresponding nodes under the head. Thus, the top of the queue contains the address of the first node to appear under the head.

One processor may service several sets of tracks, nodes being implemented on each of these sets (fig. VI-6).

Several processors may serve different sets of tracks (fig. VI-7) or may serve the same track at different locations

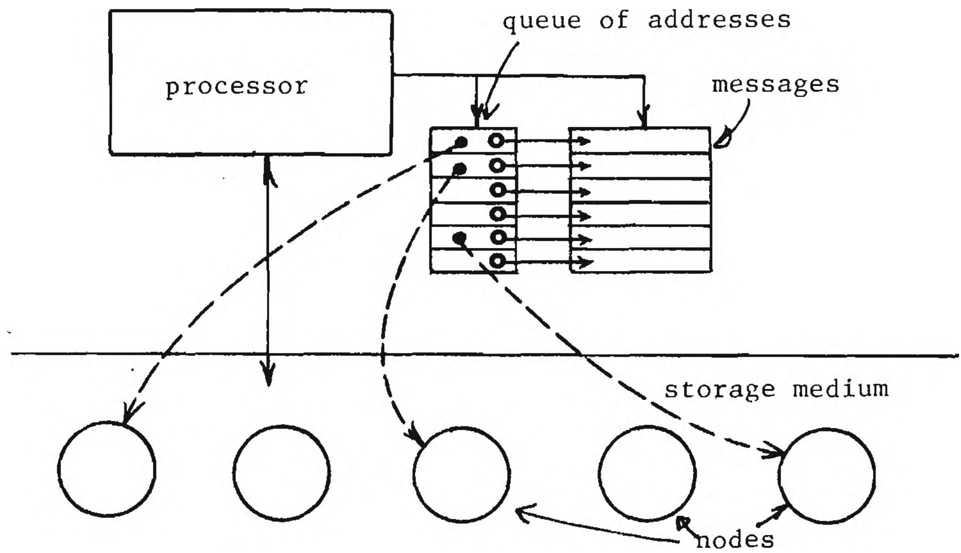

Fig. VI-3

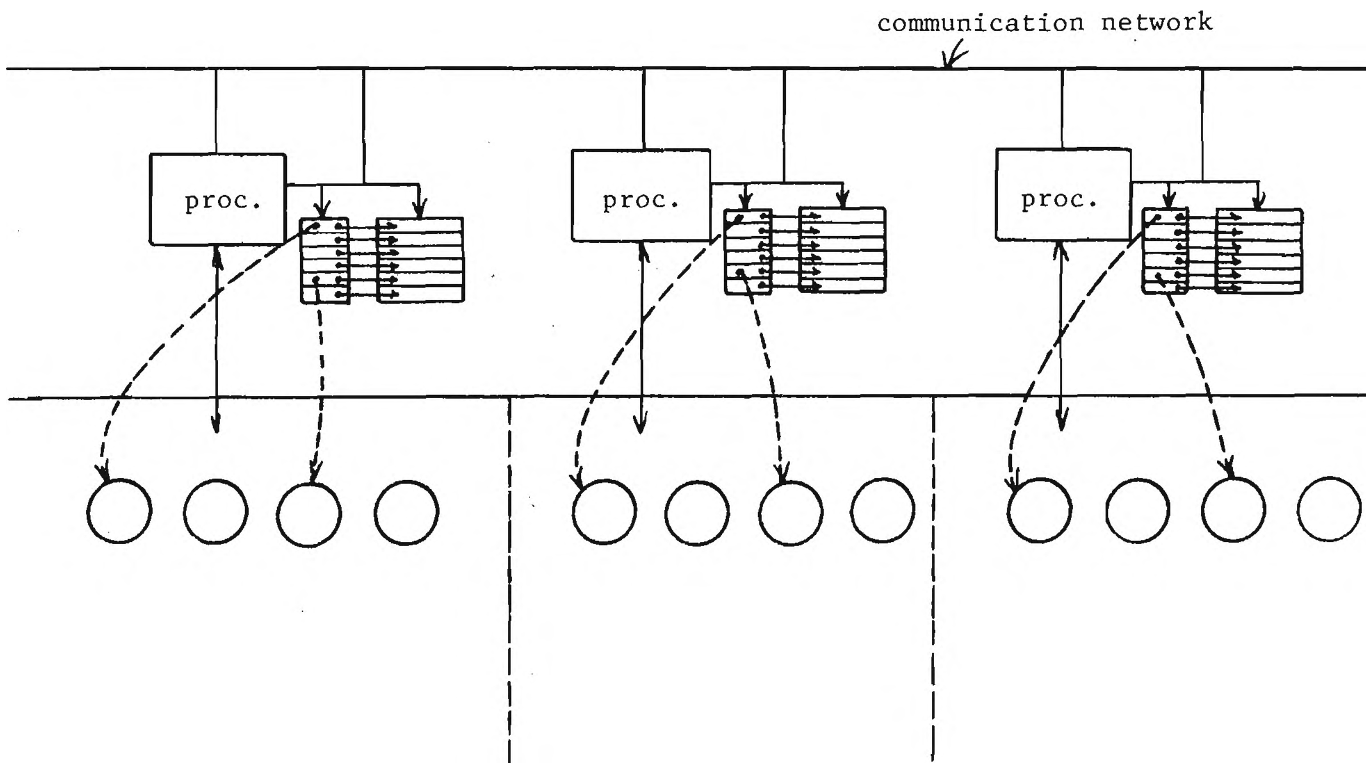

Fig. VI-4

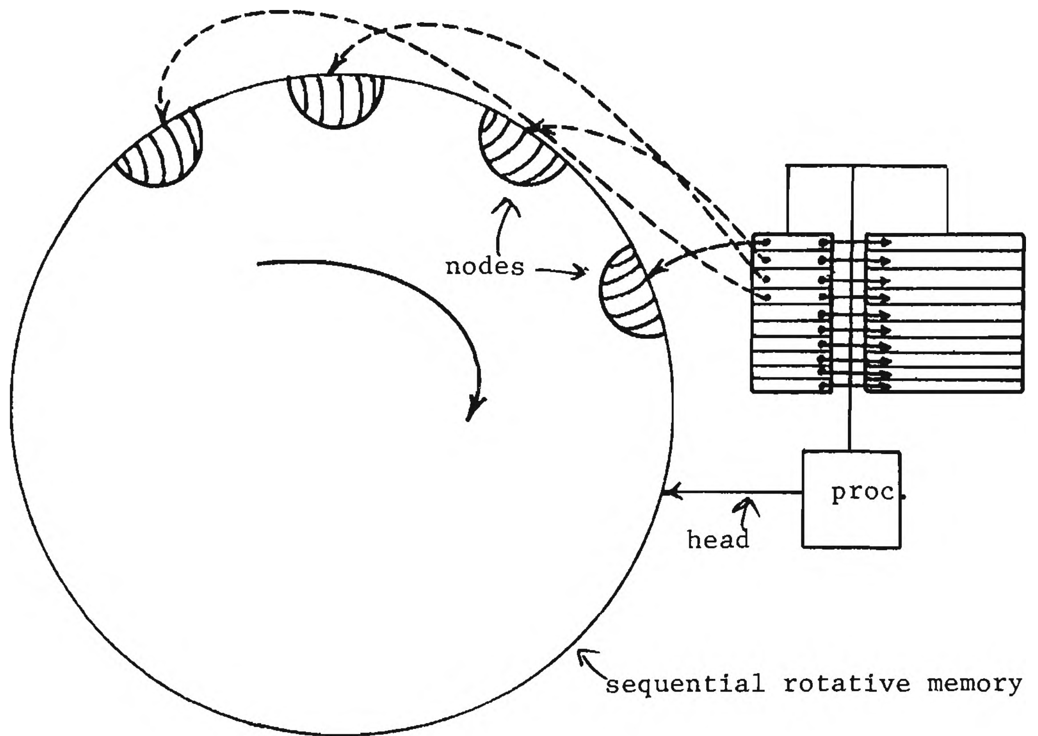

Fig. VI-5

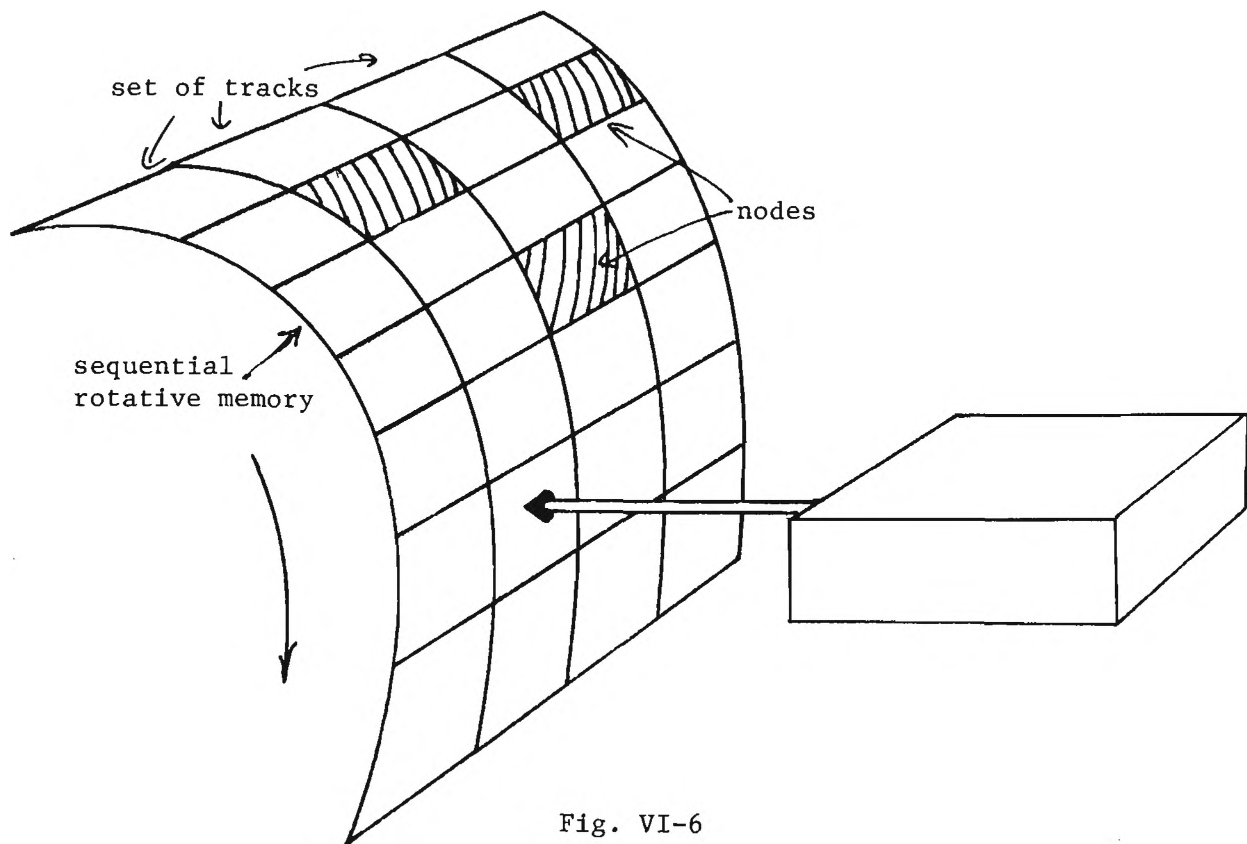

Fig. VI-6

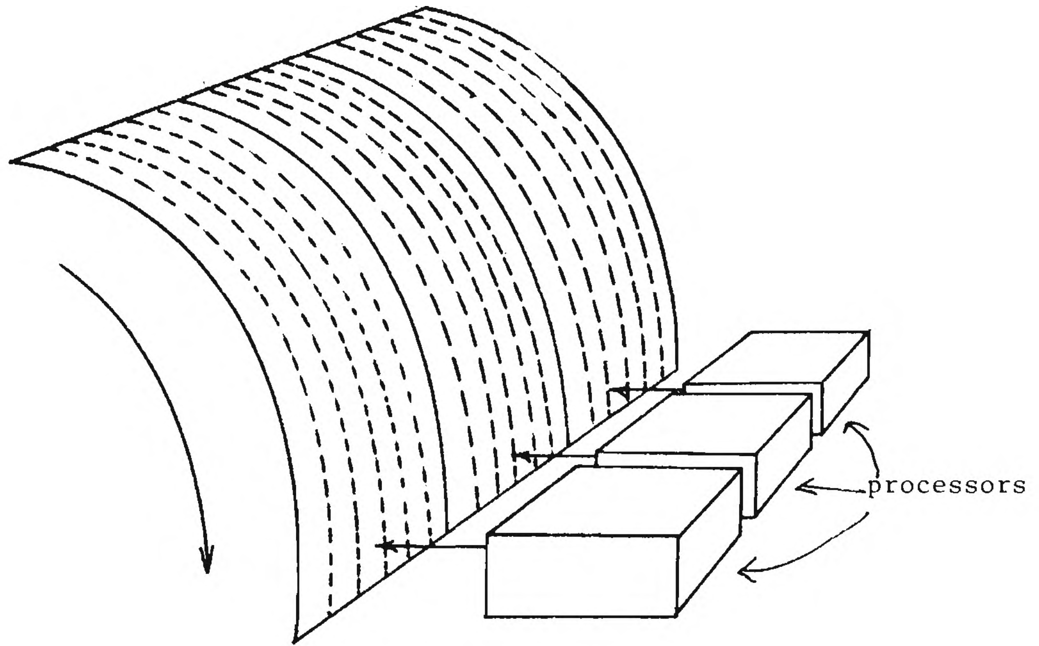

Fig. VI-7

processors

rotative
sequential
memory

communication
system

Fig. VI-8

(fig. VI-8). In both cases a communication system must link each processor to the queues associated to the other processors.

A DCPL computation tree may be implemented on a sequential rotative memory so that a son always appears under the head before its father. As a result, any request may be sent from son to father up the tree in a fraction of the rotation time (fig. VI-9). In the same way, structured cells may be implemented in order to be accessed and searched in a fraction of a rotation time (fig. VI-10).

We consider these machine organizations as being of special interest: not only sequential memories are generally rather cheap, but they also allow very large transfer rates. It becomes then possible to take advantage of the possibility of having very fast processors (for instance, a processor designed on a wafer with a cycle time of 20 ns.).

VI.5 Hierarchy of memories.

Whenever it is desired to take advantage of a very fast processor, a hierarchy of random-access memories is used (fig. VI-11): the smaller a level, the faster its memory([25]).When a memory hierarchy is used, the processor works in the fastest level, programs being swapped back and forth between the various levels. It is hoped that the processor would feel that the whole memory is as fast as the fastest one. Unless many iterative computations are expected to occur in the fastest level, it is necessary to have between any two levels a transfer rate large enough to "feed" the processor.

A large transfer rate may be obtained by taking at each level a large block as a unit of transferable information (the slower the level, the larger the block).

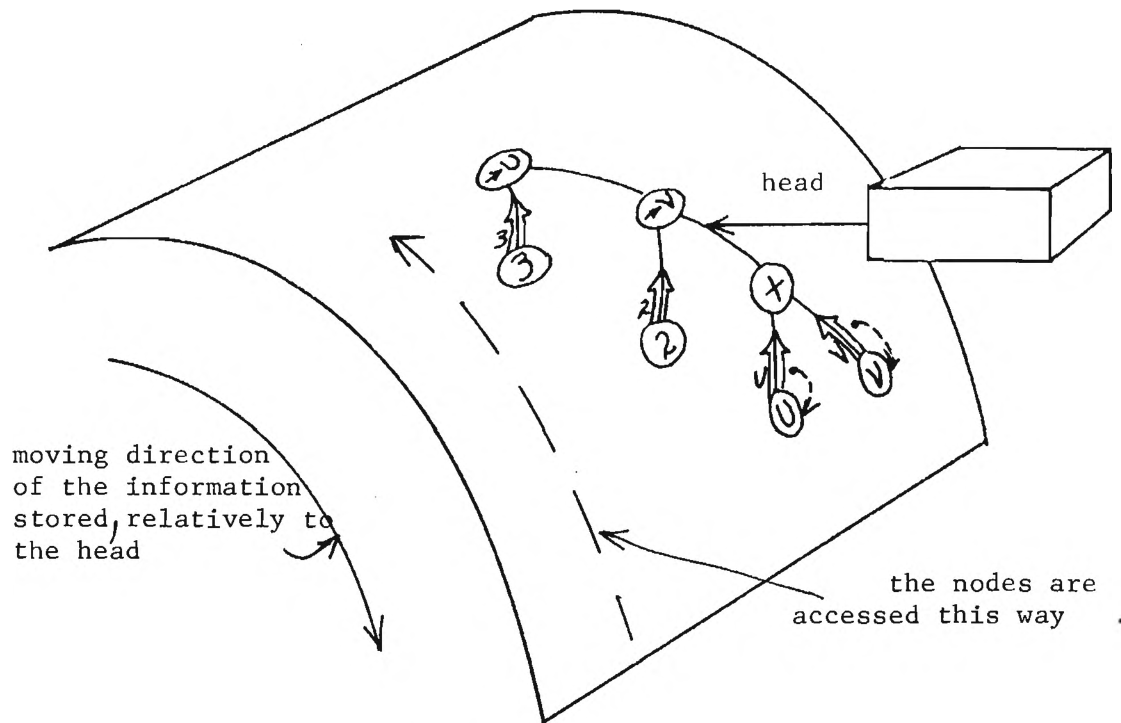

Fig. VI-9

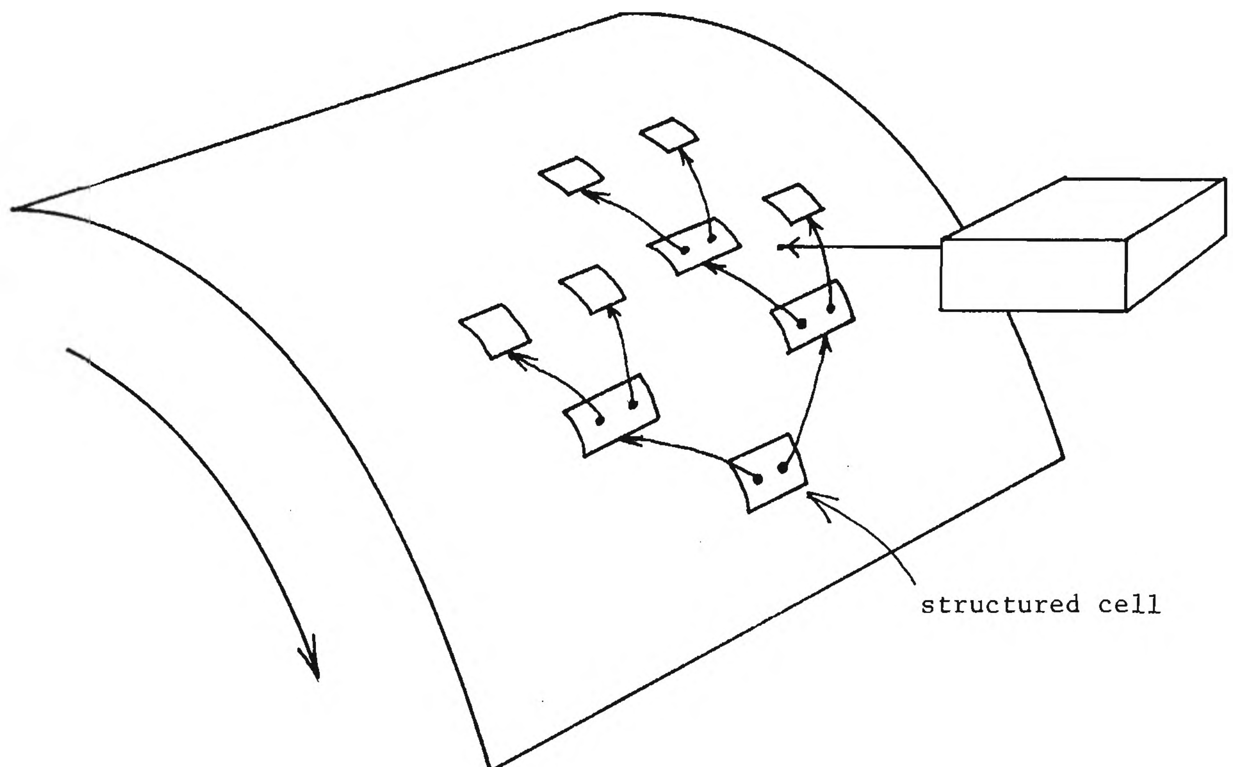

Fig. VI-10

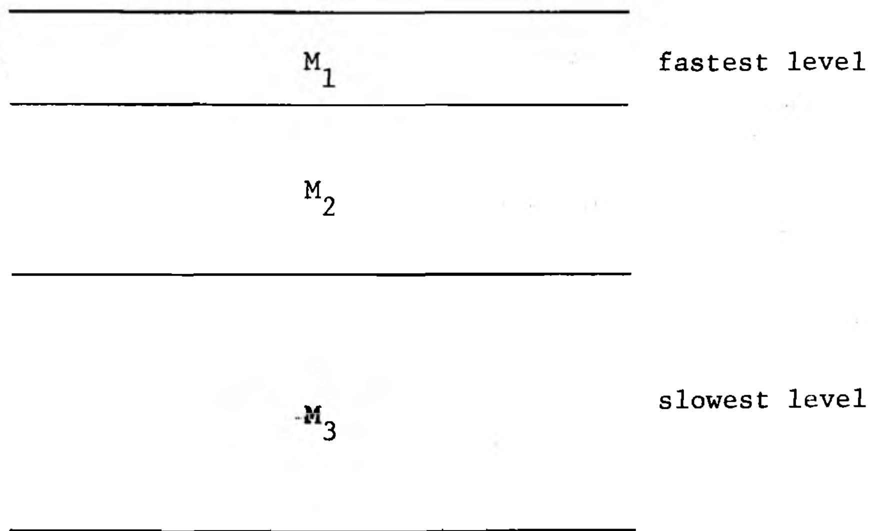

Fig. VI-11: A memory hierarchy

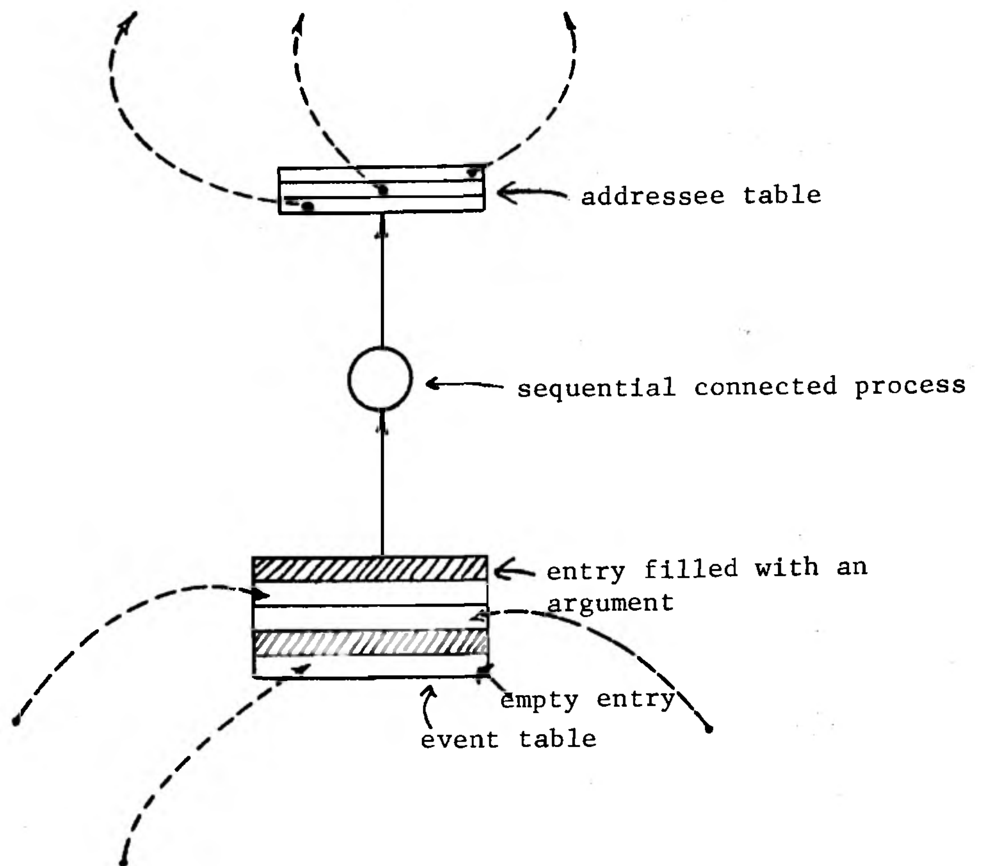

Fig. VI-12: A sequential connected program element

One may believe, however, that only a few words in such blocks would be really used. For this reason Jack Dennis suggests in [4] that information should be moved on demand with the word as information unit, a large transfer rate being assured by performing many computations in parallel.

In order to be able to use efficiently large units of transferable information, we introduce the notion of connected program. A computation is said to be connected if:

1. it can be activated whenever a certain set of arguments is available, and then completed without the need of any external information;
2. no partial result must be supplied to its environment before the completion of the computation.

As a result, a connected computation may be brought in the fastest level of a memory hierarchy and completed without the need for any information from some other level; moreover, there is no advantage in breaking the connected computation into pieces, carrying out separately the various pieces.

The notion of simple connectedness arises because a connected computation tree may be too large to be swapped in the fastest level of the memory hierarchy and have some of its parts not being connected.

A computation is said to be simply connected if it is connected and if any subcomputation it contains is connected. As a result, a simply connected computation may be brought by parts in the fastest level.

Consider the following examples in DCPL:

1. the computation tree (fig. V-18):

   NEW V (2→U; V←U+1; U+V) + (3→U; U+V)

   is simply connected.

2. the computation tree (fig. V-19):

   NEW U,V (U←5; [U+V]→Z; Z+U) + (V←U+1; U)

   is connected, but not simply: the ob (U←5; [U+V]→Z; Z+U) is not connected since the argument V can only come after the argument U has been sent to the second expression.

In DCPL it is easy to determine simply connected parts by taking for instance the largest parts which do not contain any setup '←'(these simply connected parts are by no means maximal).

Whenever a computation tree is simply connected, it is irrelevant for its environment whether the computation is carried out with a distributed or a sequential control.

VI.6 A computation as a network of sequential connected programs.

A sequential connected program element may be described by giving (fig. VI-12):

1. an event table, each entry of which may receive an argument from the environment,
2. a sequential connected process which may be applied to the arguments contained in the event table when this table is filled.
3. addresses of some other event table entries to which a produced value is to be sent (addressee table).

Remark: Any reference in a sequential connected process is local

to its program element. Any binding with the environment involves only the event and the addressee tables.

Whenever an argument is sent to an entry of the event table, this entry is filled with it. The sequential process is triggered when all the entries of the event table are filled. At the completion of the computation, the produced values are sent to some entries of some other program elements, according to the addressee table.

A computation may be considered as a network of sequential connected program elements (fig. VI-13). Fig. VI-14 displays a machine organization which can carry out such a computation. The configuration is similar to the one of fig. VI-5: the network of sequential connected program elements is stored on a sequentially accessed rotative memory; at any time there are messages waiting for an entry in an event table; the addresses of these entries are ordered in the order of appearance of these entries under the head. Whenever an event table containing an entry for which a message is waiting passes under the head, the table is examined. If, in addition to the entry for which there is a message, a non-filled entry remains, the message is just placed in the event table at the proper entry. On the contrary if, with the exception of the entry for which there is a message, the event table is filled, the corresponding program element is swapped into the scratch-pad memory (see fig. VI-14) and executed there; at its completion, the addresses of the event tables to which the messages are to be sent are placed together with the corresponding messages, at the proper places in the queue.

Let us suppose now that some of the sequential processes

Fig. VI-13

Fig. VI-14

are stored in a library, their addresses appearing only in the program elements (fig. VI-15). A computation looks at any time like the network of fig. VI-16. Whenever there is a message for an entry in an event table which will be filled with this message, a request for the proper sequential process in the library will be made; the elementary program can only be performed when this process will be available. We can already know to which entries in which event tables we will have messages to send: if these entries pass under the head while the required program has not yet been provided, the messages to be sent to these entries are not yet available. However we may see whether the event table would be filled if the message were available; if so, we may already request from the library the sequential process whose address is contained in this last elementary program. The scheme may be performed again with this new elementary program. For instance, in fig. VI-17, one message triggers the request for five sequential processes from the library.

We have seen that it is possible to single out from the computation network a subgraph of elementary programs which can be performed as soon as the corresponding sequential processes will be available from the library.

Whenever a message is conditional, we cannot know in advance whether this message will be sent or not to the associated entry in the corresponding event table. It is then a question of policy to extend or not the scheme to such event tables.

We can now propose another machine organization with a hierarchy of sequential rotative memories (fig. VI-18). For instance,

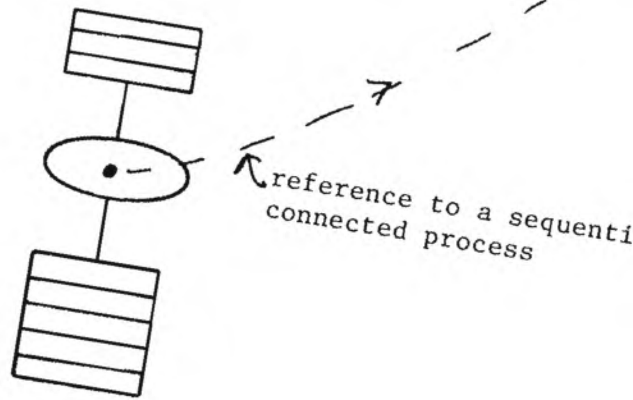

Fig. VI-15

library

Fig. VI-16

queue of addresses

messages

subgraph to be swapped, singled out by the current message

Fig. VI-17

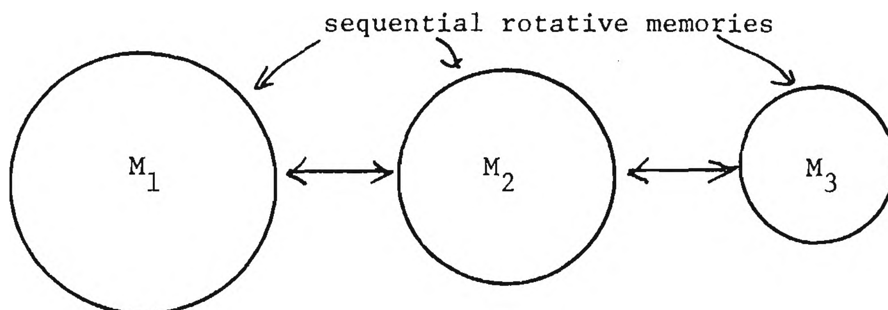

Fig. VI-18

each memory might have the same transfer rate and the capacity might be almost proportional to the rotation time:

	M1	M2	M3
rotation time	16 ms	1 ms	64 μs
capacity	256K	16K	256 words

Subgraphs of computation networks may be swapped between these memory levels using the previous scheme.

A sequential memory has a privileged direction: consequently, an implementation of a network will be more efficient if a greater number of messages are to be sent in the privileged direction.

The internal binding of a DCPL procedure may be performed before the procedure is placed in a library. This binding superimposes to the procedure a graph structure which may be expressed as a network of sequential connected program elements. Moreover, this network may be implemented contiguously in a logical space (for instance, a segment). Whenever this procedure is implemented in a computation tree, this segment is stored on the rotative memory in a way which keeps (relatively to the privileged direction) the topology of the network. Then the procedure is bound to its environment, in the computation tree, according to the DCPL binding rules. The convention we have taken previously (a son appears under the head before its father) assures a non-optimal but to a large extent satisfying solution.

A program in DCPL exhibits, to a large extent, the flow of information, the possible concurrency, and the "topology" of the computation structure. This allows new machine organizations which, we believe, would permit to obtain a larger throughput with less resources.

CHAPTER VII

CONCLUSION

Notions which are relevant to both programming language design and machine organization have been discussed in this thesis.

Progresses made in these two domains occurred generally independently; since these developments were to be compatible, artificial restrictions have been imposed in order to define a common frame of reference. We believe such restrictions can only be removed by comprehensive approaches.

Today's programming languages view computers as if they were still simple von Neumann's type machines. A computation is mainly considered as a sequence of instructions, which can modify the contents of some cells. As a result, almost any possibility of having concurrency and distribution of control in computation structures is lost.

Today's machine and system organizations consider programming languages as if they were unable to grasp the information and control topology of the computations which are expressed with them. As a result, possibilities to plan in advance the computation process are lost, producing, in our opinion, less efficient systems.

A DCPL program exhibits concurrency, distribution of control and locality of references. We believe that this will permit to have a more efficient machine organization with less expensive resources:

with a traditional organization, a very fast processor (a "processor on a chip" will be very inexpensive in just a few years) would not be able to be fully used unless a prohibitively expensive very fast memory were to be used, or small iterative computations were expected to occur very often. The machine organization proposed here is believed to be able to "feed" such a processor by using relatively not expensive, very large transfer rate, sequential memories.

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VITA

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