

Introduction to FORM

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- ▶ A powerful symbolic manipulation system specialized for high-energy physics
- ▶ Development started by Jos Vermaseren at Nikhef
- ▶ First released in 1989, with roots in Schoonschip (Martinus Veltman's system from the 1960s)
- ▶ Designed to handle extremely large algebraic expressions in quantum field theory
- ▶ Distinguishes itself by its speed and memory efficiency for massive calculations
- ▶ Used in numerous advanced calculations in particle physics
- ▶ Open source since 2000, with ongoing development by the community

- ▶ In 1960s Martinus J. G. Veltman developed first CAS - SCHOONSCHIP.

- ▶ Later on he was writing:

... on the basis of this program [SCHOONSCHIP], the commercially successful software program Mathematica was developed by Wolfram. While most theoreticians were getting lost in all kinds of formulas, I could just get things calculated. That gave me an enormous advantage.

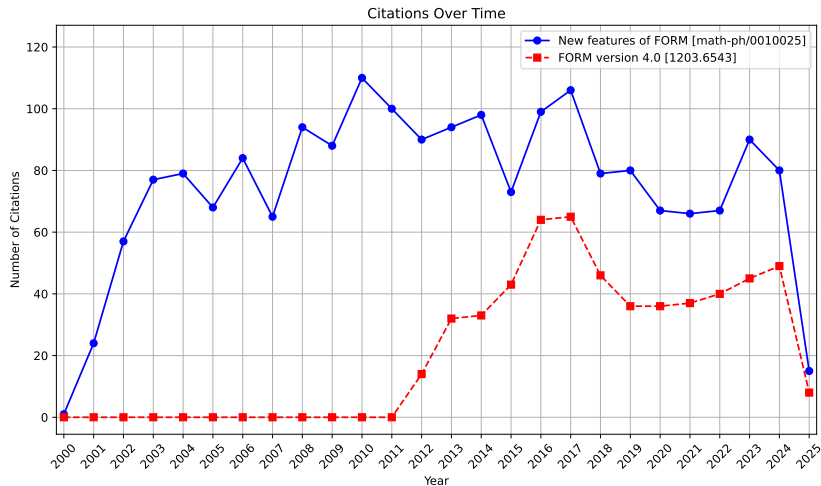


M. Veltman

- ▶ Many other computer algebra systems were released afterwards, which are more efficient and user friendly than SCHOONSCHIP.
- ▶ Until now the core principle remains unchanged:
 - ▶ **The user** is expected to map the natural language to artificial language constructed by some developer
- ▶ But still, not all of CASs are created equally...



- ▶ "FORM is an ancient tool no longer relevant in modern physics"
- ▶ "It's not actively maintained and there are a lots of bugs"
- ▶ "Modern CAS systems have made FORM obsolete"
- ▶ "FORM is not user-friendly with its steep learning curve"
- ▶ "It's only useful for specialized calculations that nobody does anymore"
- ▶ "The syntax is too cryptic and difficult to learn"
- ▶ "It's not compatible with modern computational workflows"
- ▶ "Only a handful of experts can actually use it effectively"
- ▶ "There's no community support and you are left alone with your problems"





- ▶ Free and open-source software with an active development community
- ▶ High efficiency in handling large polynomial expressions
- ▶ Transparent memory management and algorithmic approaches
- ▶ Highly portable across different computing environments and platforms
- ▶ Minimalistic design focused on performance
- ▶ Numerous relevant public libraries are implemented in FORM
- ▶ Parallel processing capabilities for modern multi-core systems
- ▶ Specialized features designed specifically for quantum field theory
- ▶ Continuous improvements with new releases and updates respecting the original design philosophy



- ▶ Demonstrate core functionality through practical examples
- ▶ Assume zero prior knowledge
- ▶ Highlight common pitfalls
- ▶ Show efficient techniques for large calculations
- ▶ Equip you with workarounds for limitations

Today:

- ▶ Basic functionalities, basic exercises

Tomorrow:

- ▶ Advanced features, advanced exercises



- ▶ The historical homepage of FORM: <https://www.nikhef.nl/~form>
- ▶ The actual development/discussions: <https://github.com/vermaseren/form>
- ▶ The Reference Manual:
<https://www.nikhef.nl/~form/maindir/documentation/reference/reference.html>
- ▶ The latest official release: v4.3.1
- ▶ The actual development version: v5.0.0-beta.1
- ▶ Installation:
 - ▶ From official github repository
 - ▶ Using a helper script: <https://github.com/magv/hepware>



hello_world.frm:

```
1  #-  
2  #message Hello World  
3  .end
```

form hello_world.frm:

```
FORM 5.0.0-beta.1 (Mar  7 2025, v5.0.0-beta.1-122-g638f84b)  Run: Mon Mar 2  
  #-  
~~~Hello World  
  0.00 sec out of 0.00 sec
```



- ▶ The code is organized in *modules*.
- ▶ The commonly used modules:
 - ▶ `.sort`
The general end-of-module. Causes execution of all active expressions, and prepares them for the next module.
 - ▶ `.end`
Executes all active expressions and terminates the program.
 - ▶ `.global`
No execution of expressions. It just saves declarations made thus far from being erased by a `.store` instruction.
 - ▶ `.store`
Executes all active expressions. Then it writes all active global expressions to an intermediate storage file and removes all other non-global expressions. Removes all memory of declarations except for those that were made before a `.global` instruction.

- ▶ Modules consist of statements
- ▶ Statements should always appear in the following order:
Declarations
....
Specifications
....
Definitions
...
Statements
...
Output control statement
- ▶ FORM manual lists for each keyword which category does it belong to

Example of a valid FORM code



```
1  #-
2  * Declarations
3  Symbols m,n;
4  * Specifications
5  Off statistics;
6  * Definitions
7  Local A = (m+n)^2;
8  * Output
9  Print;
10 .sort
11 * Declarations
12 Symbol z;
13 * Specifications
14 Drop;
15 * Definitions
16 Local B = A+z;
17 * Statement
18 identify n = 3;
19 * Output
20 Print;
21 .end
```

Design choice:

- ▶ Expressions are always expanded
- ▶ Operations inside the module are applied term-by-term

Consequences:

- ▶ Possible proliferation of terms to be processed
- ▶ Effective memory management
- ▶ Difficult to implement anything beyond single term based operations

```
1  Symbols a,b,c;  
2  Local F = 14*a-23*b;  
3  Print "1: %t";  
4  id a = b+c;  
5  Print "2: %t";  
6  Print;  
7  .end
```

Output:

```
#-  
1:  + 14*a  
2:  + 14*b  
2:  + 14*c  
1:  - 23*b  
2:  - 23*b
```

```
F =  
14*c - 9*b;
```

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Compare

```
1  Symbols a,b,c,d,e;  
2  
3  Local test = (a+b+c+d)^30;  
4  id a = -b+5;  
5  
6  id c = -d+e;  
7  Print;  
8  .end
```

vs.

```
1  Symbols a,b,c,d,e;  
2  
3  Local test = (a+b+c+d)^30;  
4  id a = -b+5;  
5  .sort  
6  id c = -d+e;  
7  Print;  
8  .end
```



- Two fundamental types of functions:

```
1  #-  
2  CFunctions A,B,C;  
3  Functions  F,G,H;  
4  
5  L test = G(25)*C(1,2)*A(3,4)*B(6,5)*H(1)*F(4);  
6  
7  Print;  
8  .end
```

Output:

```
test =  
      G(25)*H(1)*F(4)*A(3,4)*B(6,5)*C(1,2);
```

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- Substitutions are made using `identify (id)` statement:

```
1  #-  
2  CFunction F;  
3  Symbols a,b;  
4  
5  L test = F(1,a);  
6  id F(a?,b?) = F(b,a,a+b);  
7  
8  Print;  
9  .end
```

Output:

```
test =  
  F(a,1,1 + a);
```

0.00 sec out of 0.00 sec

- ▶ `a?`, `b?` are wildcards
- ▶ They match the objects depending their declaration
- ▶ One might have complicated restrictions on wildcards, e.g. `a?{,>-2}`, match symbols which are larger than `-2`.
- ▶ The symbols starting with `?` are ranged wildcards, matching with everything until the end range:

```
1  #-  
2  CFunction F;  
3  
4  L test = F(37,28);  
5  id F(?a) = F(1,?a);  
6  
7  Print;  
8  .end
```

What are these programs going to produce?

```
1  #-  
2  CFunction F;  
3  Symbols a,b;  
4  
5  L test = 1/F(1,a);  
6  id F(a?,b?) = F(b,a,a+b);  
7  
8  Print;  
9  .end
```

or

```
1  #-  
2  CFunction F;  
3  Symbols a,b;  
4  
5  L test = F(1,a);  
6  id a = b;  
7  
8  Print;  
9  .end
```

Operate on the arguments of a function:

```
1  #-  
2  CFunction F;  
3  Symbols a,b;  
4  
5  L test = F(1,a);  
6  Argument F;  
7      id a = b;  
8  EndArgument;  
9  
10 Print;  
11 .end
```

Output:

```
test = F(1,b);
```

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```
1  Symbols a,b,c,d,e;
2  CFunction pow;
3
4  Local test = pow(a+b+c+d,30);
5  Argument pow;
6  id a = -b+5;
7  id c = -d+e;
8  EndArgument;
9
10 id pow(?k,a?) = exp_(?k,a);
11 Print;
12 .end
```

Time =	0.00 sec	Generated terms =	31
	test	Terms in output =	31
		Bytes used =	1068

Compare the timing/memory usage with the previous approach.

- `dd_` is a totally symmetric tensor represented as a sum of symmetrized product of Kronecker deltas (`d_`):

```
1  #-  
2  AutoDeclare Vector v;  
3  Local F = dd_(v1,v2,v3,v4);  
4  Print;  
5  .end
```

Output:

```
F =  
v1.v2*v3.v4 + v1.v3*v2.v4 + v1.v4*v2.v3;
```

Fun quest:

$$\langle \Omega | T \{ \phi(x_1) \dots \phi(x_n) \} | \Omega \rangle = \frac{\langle 0 | T \left\{ \phi_0(x_1) \dots \phi_0(x_n) e^{i \int d^4x \mathcal{L}_i[\phi_0]} \right\} | 0 \rangle}{\langle 0 | T \left\{ e^{i \int d^4x \mathcal{L}_i[\phi_0]} \right\} | 0 \rangle}$$

Demonstrate that in the S -matrix vacuum bubble diagrams cancel at any order of a perturbation theory.

► `e_` is a Levi-Civita Tensor

```
1  #-  
2  Dimension 3;  
3  Indices i,j,k,p,q,r;  
4  Local test = e_(i,j,k) * e_(p,q,r);  
5  *Wrong  
6  *id r = k;  
7  
8  contract;  
9  Print;  
10 .end
```

How can we make a replacement $r \rightarrow k$ to obtain the following output?

```
test =  
    d_(i,p)*d_(j,q) - d_(i,q)*d_(j,p);
```

What would you do if we don't know how many indices are inside `e_`?

- `g_` objects represent the Dirac gamma matrices

```
1  #-  
2  Dimension D;  
3  Index i1,i2;  
4  Local test = g_(1, i1)*g_(1,i2);  
5  tracen 1;  
6  Print;  
7  .end
```

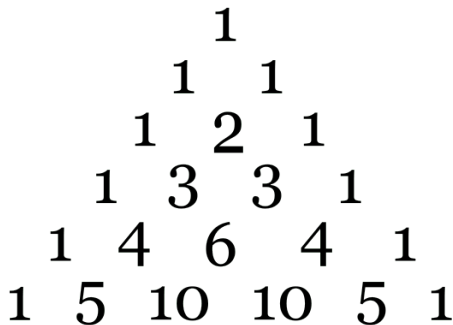
Produces:

```
test =  
4*d_(i1,i2);
```

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Show that:

$$\text{tr}\left(\gamma^\mu \gamma^\nu \gamma^\rho \gamma^\sigma \gamma^5\right) = -4i\epsilon^{\mu\nu\rho\sigma}.$$



- ▶ Build the Pascal's triangle from top to bottom
- ▶ Define a procedure which can extract an element at the depth i with the column number j
- ▶ Calculate the sum of all elements at the depth i .
- ▶ i and j are integers between 100 and 1000.

Thanks for your attention!

