

FormCalc 7

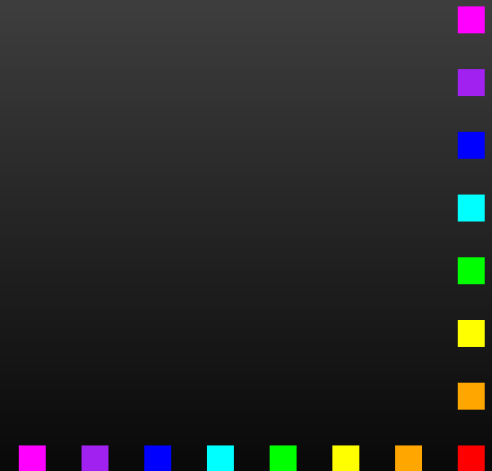
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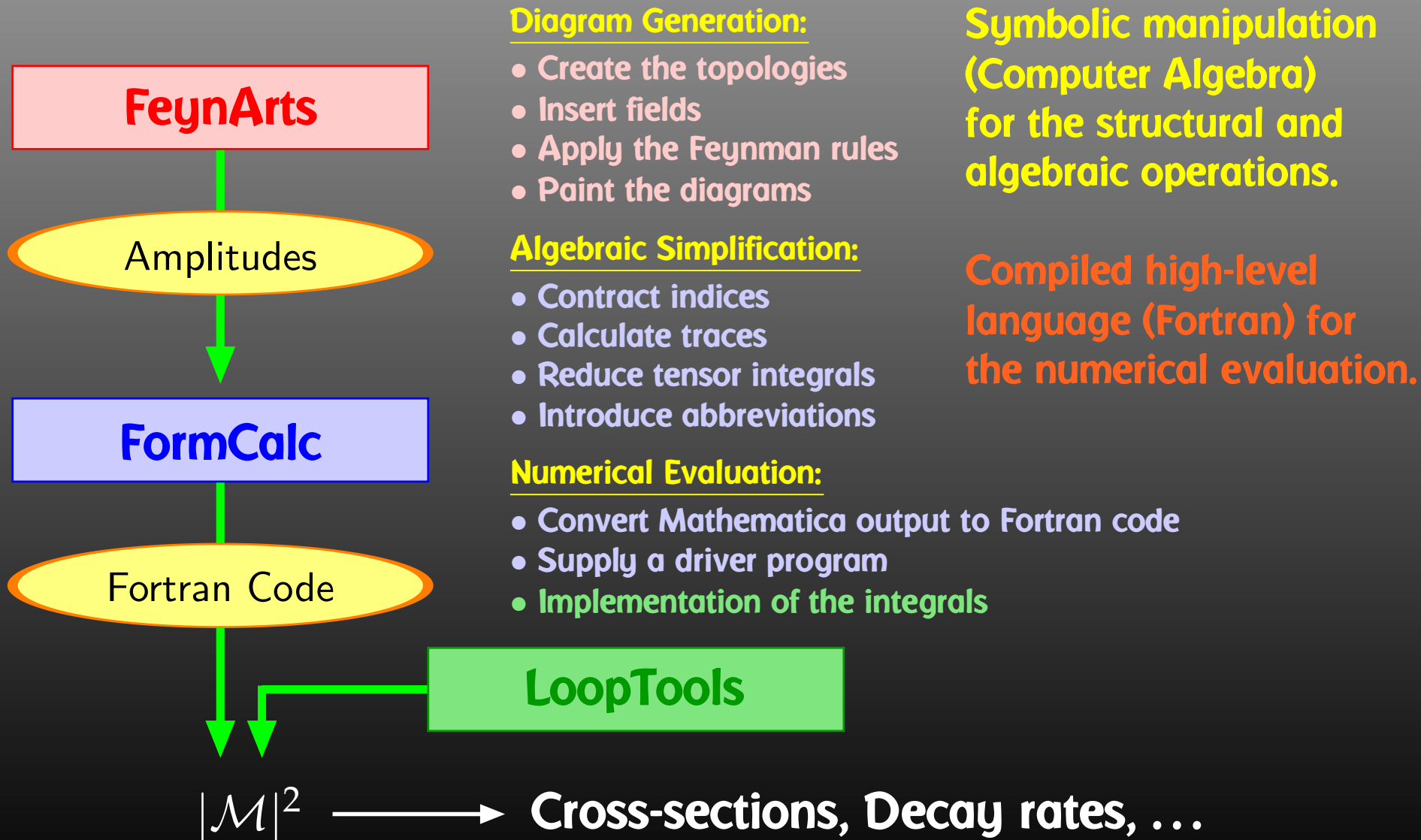
From microsoft.com/en-us/windows7:

Why get Version 7?

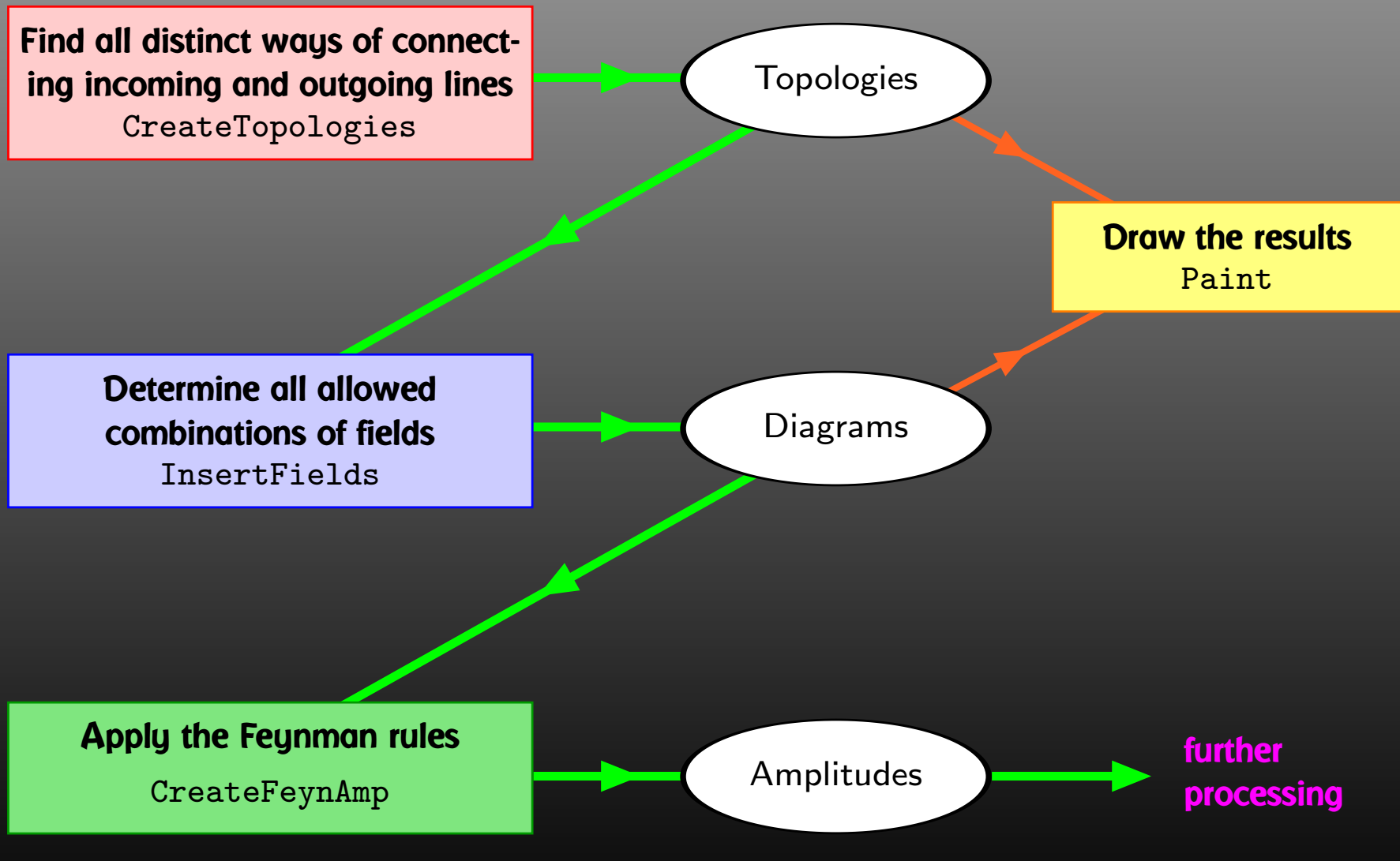
- To simplify everyday tasks
- To work the way you want
- To do new things



Automated Diagram Evaluation



FeynArts



Algebraic Simplification

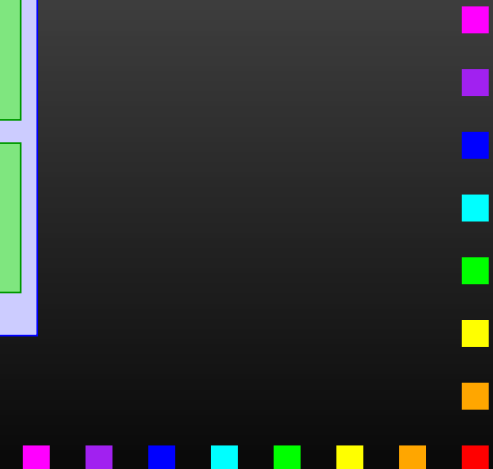
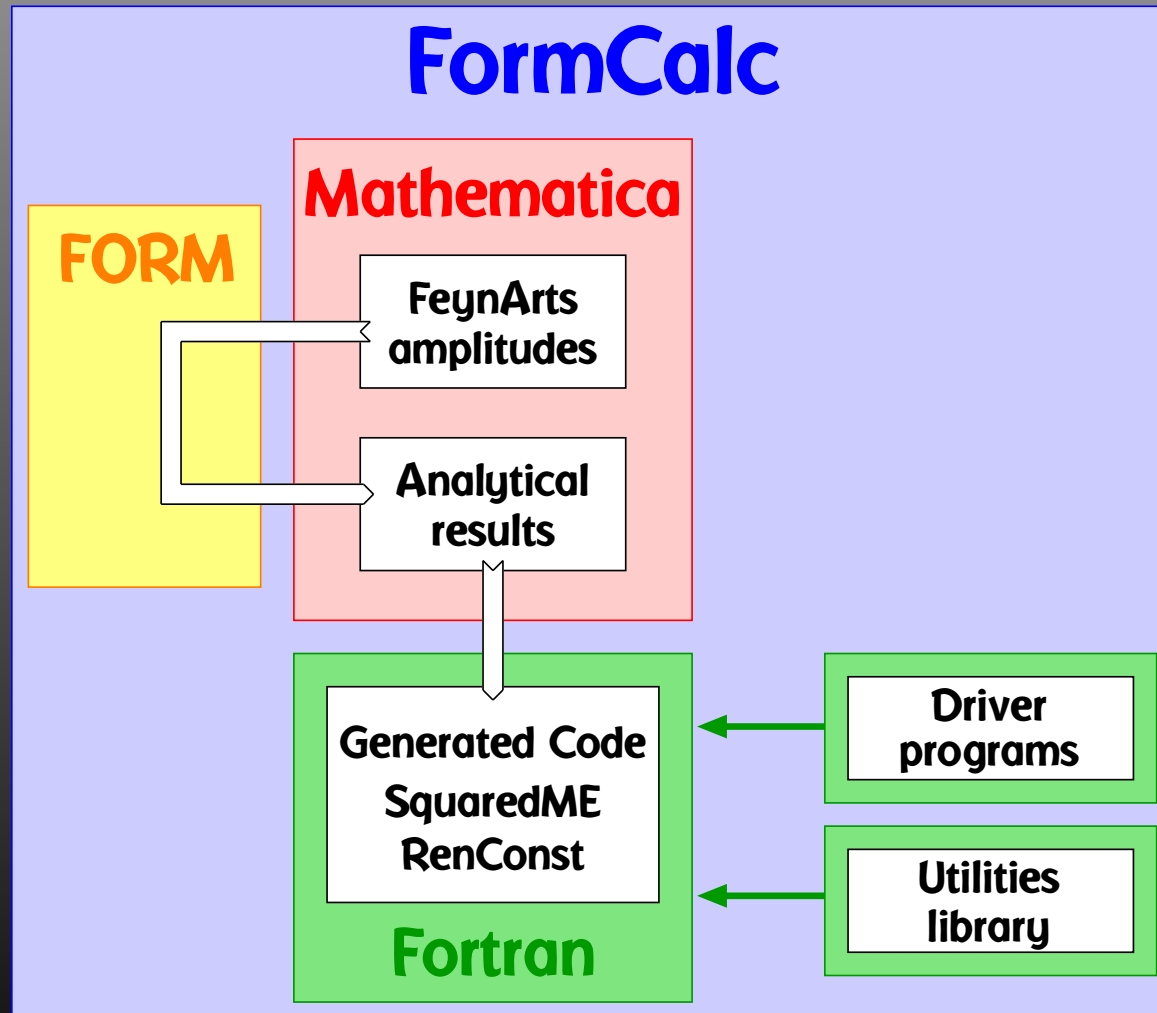
The amplitudes of `CreateFeynAmp` are in **no good shape for direct numerical evaluation.**

A number of steps have to be done analytically:

- **contract indices as far as possible,**
- **evaluate fermion traces,**
- **perform the tensor reduction,**
- **add local terms arising from \mathcal{D} -(divergent integral) (dim reg + dim red),**
- **simplify open fermion chains,**
- **simplify and compute the square of $SU(N)$ structures,**
- **“compactify” the results as much as possible.**



FormCalc Internals



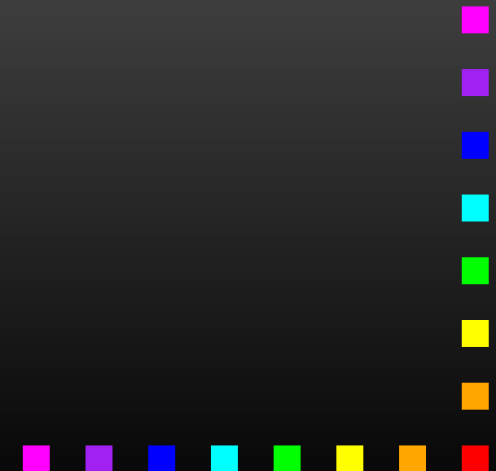
FormCalc 7

New Features:

- Analytic tensor reduction,
- Unitarity methods (OPP),
- Improved code generation,
- Command-line parameters for model initialization, MSSM (SM) initialization via FeynHiggs.

Cuba:

- Built-in Parallelization available.



Analytic Tensor Reduction

Work done in collaboration with S. Agrawal.

Passarino-Veltman reduction is still useful. So far:

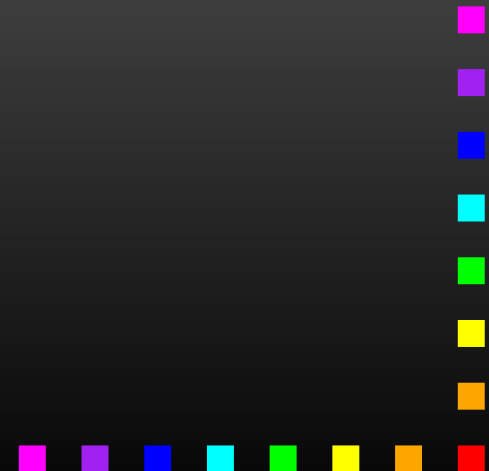
- introduction of tensor coefficients in FormCalc, e.g.

$$\int d^4q \frac{q_\mu q_\nu}{D_0 D_1} \sim B_{\mu\nu} = g_{\mu\nu} B_{00} + p_\mu p_\nu B_{11}$$

- complete reduction to scalars only numerically in LoopTools.

Available now: Analytic Reduction in FormCalc.

```
CalcFeynAmp[... , PaVeReduce -> True]
```



Analytic Tensor Reduction

Reduction formulas from Denner & Dittmaier, hep-ph/0509141.
Not straightforward to implement in FORM.

Apart from analytic considerations, this is useful e.g. for low-energy observables, where small momentum transfer may lead to **numerical instabilities in numerical reduction**, as in:

$$B_\mu = p_\mu B_1 \quad \text{for} \quad p \rightarrow 0$$

Unless FormCalc finds a way to cancel it immediately, the **inverse Gram determinant appears wrapped in IGram** in the output, so is available for further modifications.



Unitarity Methods

Work done in collaboration with E. Mirabella.

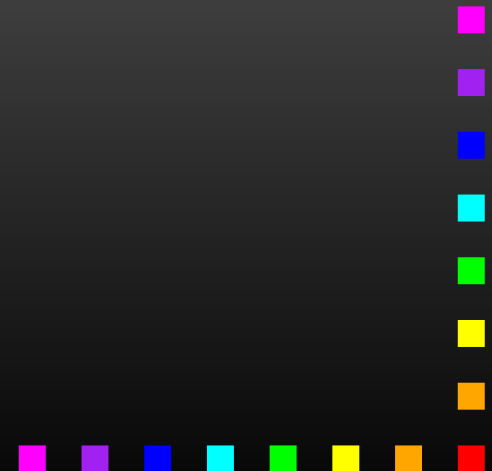
We employ the **OPP (Ossola, Papadopoulos, Pittau)** methods as implemented in the two libraries **CutTools** and **Samurai**.

Instead of introducing tensor coefficients, the **numerator is put into a subroutine** which is **sampled by the OPP function**, as in:

$$\varepsilon_1^\mu \varepsilon_2^\nu B_{\mu\nu}(p, m_1^2, m_2^2) = B_{\text{cut}}(2, N, p, m_1^2, m_2^2)$$

where

$$N(q_\mu) = (\varepsilon_1 \cdot q) (\varepsilon_2 \cdot q)$$



Unitarity Methods

So far tested on a handful of $2 \rightarrow 2$ and $2 \rightarrow 3$ processes, get **agreement to about 10 digits.**

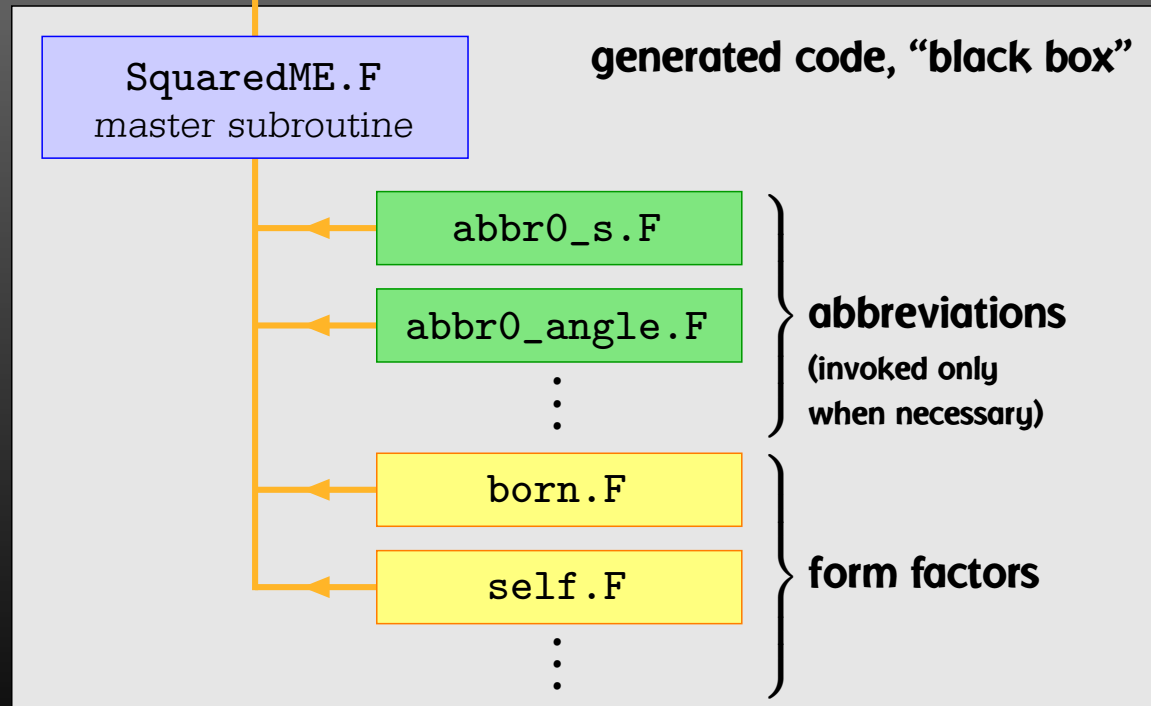
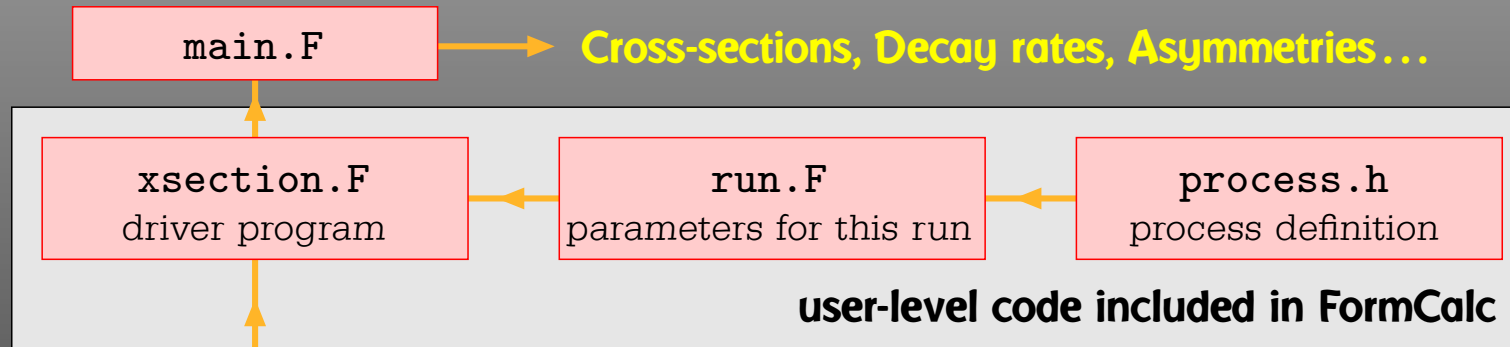
Performance somewhat wanting as of now, Passarino-Veltman beats OPP hands down in the processes we looked at.

Currently optimizing performance:

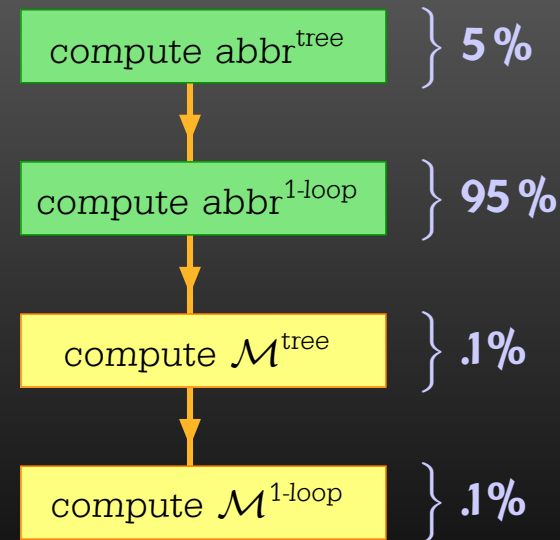
- Option to specify the N in N -point up to which Passarino-Veltman is used, above OPP
- Minimizing OPP calls to reduce sampling effort - work in progress.
- Already looked into tweaking caching of loop integrals, but pointless: lower- N integrals also needed by OPP.



Numerical Evaluation in Fortran 77



CPU-time (rough)



Code generation

Currently: Output in Fortran 77.

Code generator is rather sophisticated by now, e.g.

- **Expressions too large** for Fortran are split into parts, as in

```
var = part1  
var = var + part2  
...
```

- **High level of optimization**, e.g. common subexpressions are pulled out and computed in temporary variables.
- **Many ancillary functions** make code generation versatile and highly automatable, such that the resulting code needs few or no changes by hand.

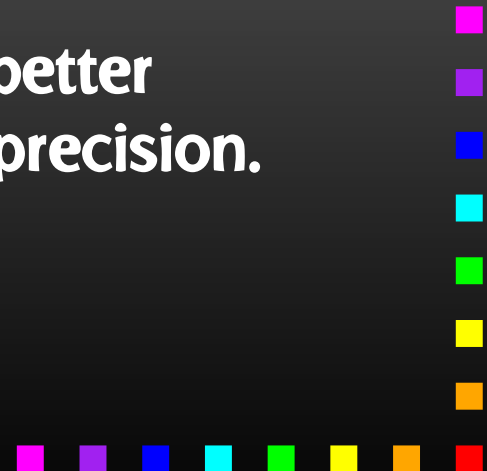


Improvements in Code Generation

- Working on **Output in C**, makes integration into C/C++ codes easier and allows for GPU programming.
- Main subroutine SquaredME.F is now **sectioned by comments**, to aid **automated substitution** e.g. with sed, for example:

```
* BEGIN VARDECL  
...  
* END VARDECL
```

- Introduced **data types** Real **and** Complex for better abstraction, can e.g. be changed to different precision.



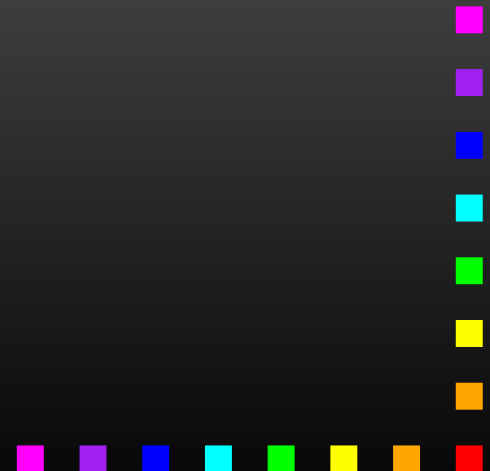
Command-line parameters for model initialization

Extension of command-line argument parsing:

```
run :arg1 :arg2 ... uuuu 0,1000
```

The ':'-arguments are **passed to model initialization code.**

Internal routines in `xsection.F` accordingly have additional parameters `argv, argc`.

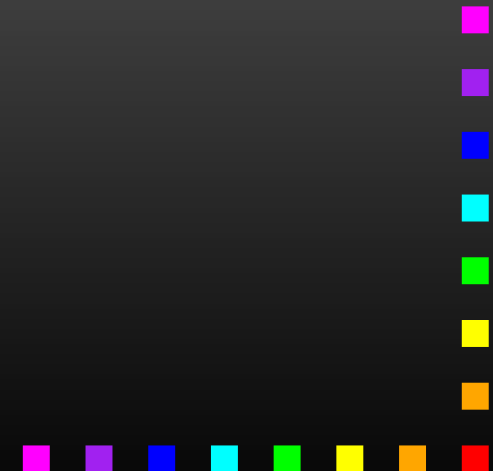


Model Initialization through FeynHiggs

- `model_fh.F` uses **FeynHiggs as Frontend for FormCalc-generated code:**

```
run :fhparameterfile :fhflags uuuu 0,1000
```

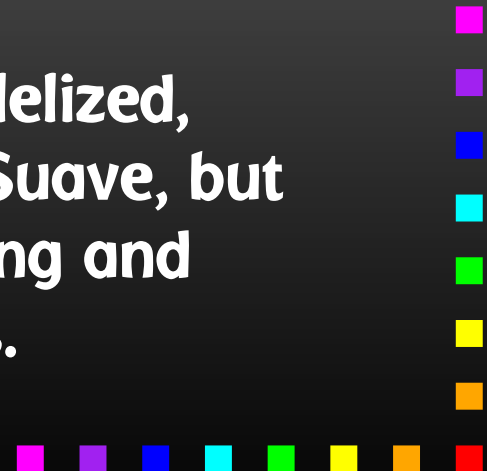
- FeynHiggs initializes MSSM (SM) parameters and passes them to FormCalc code.
- No duplication of initialization code.
- Parameters consistent between Higgs-mass computation and cross-section calculation.
- Needs FeynHiggs 2.8.1 or above.



Cuba Parallelization

New version Cuba 3 features built-in parallelization:

- Just **set environment variable** CUBACORES to the number of cores one wishes to utilize.
- Uses `fork/wait`, with back-communication of results through pipes, thus **no need to write reentrant integrand function**, works identically in Fortran and C/C++.
- Effective on multi-core machines only, no parallelization across network.
- Only the sampling function is presently parallelized, which is the optimal solution for Vegas and Suave, but sub-optimal Divonne and Cuhre. Benchmarking and optimization for the latter is work in progress.



Summary

New Features in FormCalc 7:

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- Analytic tensor reduction in CalcFeynAmp,
- Unitarity (OPP) methods using either the Samurai or CutTools library,
- Improved code generation,
- Command-line parameters for model initialization,
- Initialization of MSSM parameters via FeynHiggs.

Cuba:

feynarts.de/cuba

- Built-in Parallelization available simply by setting an environment variable.

