

# Introduction to grace system

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## Contents:

1. Introduction
2. Structure of grace
3. Installation
4. Sample calculation

# 1. Introduction

- Problem:
  - Too many processes in high energy collider physics
  - Too many Feynman graphs for a process with many final particles.

Example:

- $pp \Rightarrow W + 3jet$ 
    - 31 processes
    - 736 Tree Feynman Graphs
    - 20915 One-loop Feynman Graphs
  - $pp \Rightarrow W + 4jet$ 
    - 96 processes
    - 9716 Tree Feynman Graphs
    - 414906 One-loop Feynman Graphs
- (initial 1 generation, final 2 generations)

- Solution

Automatize perturbative calculation on computers.

- Generation of Feynman graphs
- Generation of code for calculation
- Without approximation
- Generation of event generator

- Status of grace
  - Full automatic for tree processes: ( $2 \Rightarrow 6, 7, 8, \dots$ )
    - \* Electro-weak
    - \* SUSY (MSSM)
    - \* QCD with Parton shower/PDF

This tutorial is for tree processes.

- Automatic code generation for 1-loop processes:  
 $(2 \Rightarrow 2, 3)$

- \* Electro-weak
  - \* SUSY (MSSM)

Calculation requires large CPU time for  $2 \Rightarrow 3, \dots$

Need to write many scripts to run jobs.

- Automatic generation NLO QCD processes

- \* Generation of Matrix Elements
  - \* NLO parton shower
  - \* Hadronization

In progress

- Event generators
  - \* `grc4f`
  - \* `susy23`
  - \* `GR@PPA`
  - \* Used in KORALW

Joint to grace project!

## 2. Structure of grace

- **input**

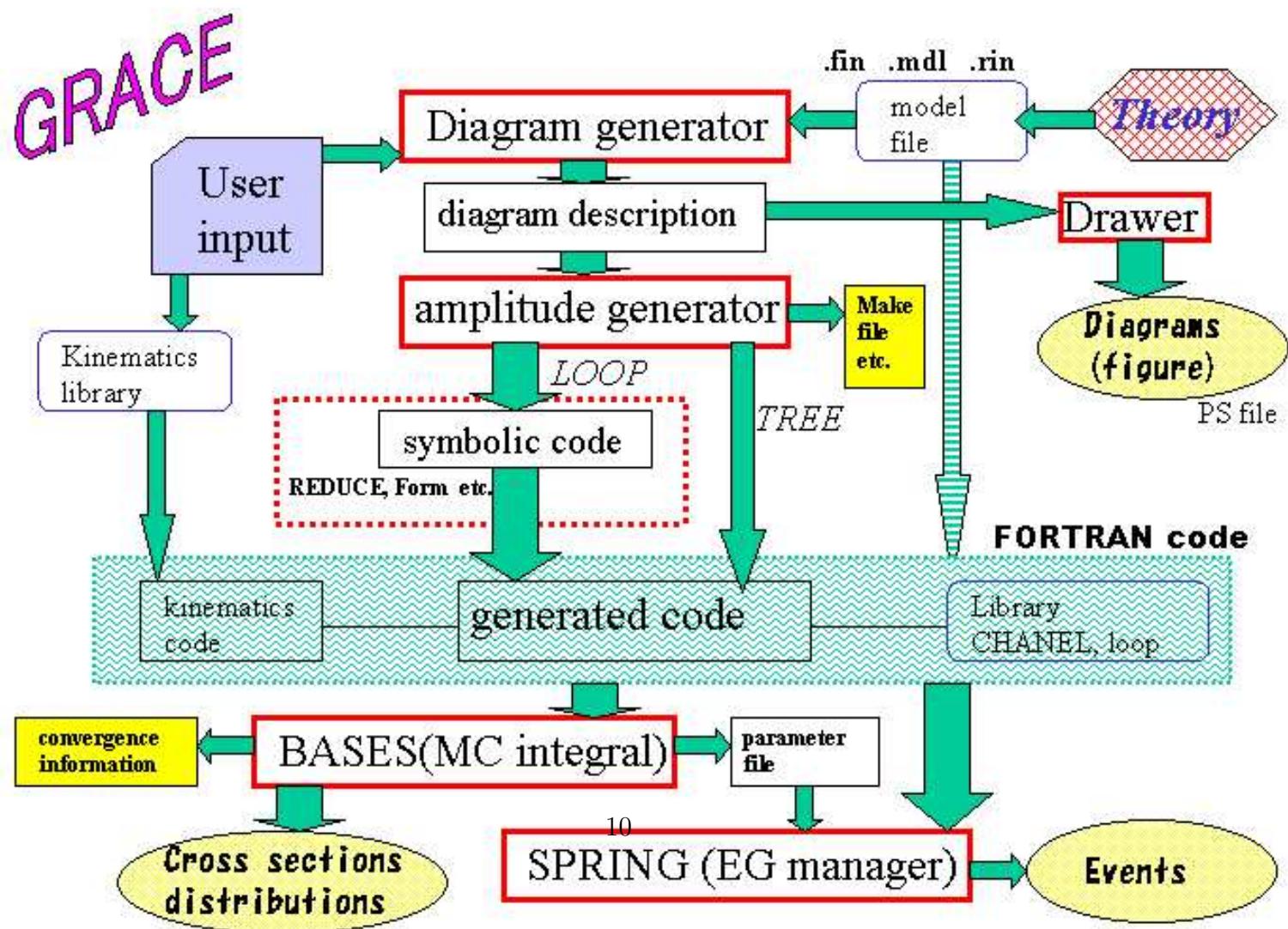
- `in.prc` : Description of processes
- `*.mdl` : Description of the model
- `*.fin` : Model dependent run-time routines

- **output**

- `*.f` etc. : Generated code
- `Histograms` : Value of cross section
- `Events` : Generated events

For tree processes

1. Graph generation (`grc`: all order)
2. Graph drawing (`gracefig`, `grcdraw`)
3. Generation of Fortran code (`grcfort`)
4. Library of helicity amplitude (`chanel`)
5. Kinematics library (`dbkinem`)
6. Phase space integration (`bases`)
7. Event generation (`spring`)



## Confirmation of the results

- **Gauge invariance** at one phase space point.
    - ~ 15 digits in double precision
    - ~ 30 digits in quadruple precision
- ⇒ logical consistency of generated code.
- ⇒ correctness of small contributing terms.  
⇒ may give large contributions for different parameters or conditions.
- ⇒ check numerical stability.

## 4. Installation

### 1. System Requirement

- (a) ANSI C compiler
- (b) Fortran 77 compiler
- (c) Standard Unix commands
- (d) X Window system : optional for drawing graphs
- (e) Motif Toolkit or its clone ([OpenMotif](#), [lesstif](#)) :  
optional for drawing graphs

2. Getting source code:

`http://minami-home.kek.jp/`

3. Expand source code:

`tar xvzf grace.220.yyyy.mmdd.tgz`

4. Configuration:

`cd grace.2.2.0  
./Config.sh`

## 5. Compilation:

make

make install

## 6. Check system:

make test-install

for  $e^+e^- \rightarrow W^+W^-\gamma$  in SM:

tail testinst/sm/eewwa/fort/gauge.out

and for  $e^+e^- \rightarrow \gamma\tilde{\chi}_1^+\tilde{\chi}_1^-$  in MSSM:

tail testinst/mssm/asw1SW1/fort/gauge.out

## 4. Sample calculation

### 1. Add commands to search PATH

```
setenv PATH $PATH:$HOME/grace.2.2.0/bin  
rehash
```

in "csh" or "tcsh", or

```
PATH=$PATH:$HOME/grace.2.2.0/bin  
export PATH
```

in "sh" or "bash".

### 2. Preparation of input file

(a) Create working directory

cd

mkdir grcwork

cd grcwork

(b) Copy sample input file “in.prc”

cp ~/grace.2.2.0/testinst/sm/eewwa/input/in.prc .

### (c) Edit input file “in.prc”

```
%%%%%%%%
Model="sm.mdl";
%%%%%%%
Process;
    ELWK=3;
    Initial={electron, positron};
    Final  ={photon W-plus, W-minus};
    Expand=Yes;
    OPI=No;
    Kinem="2302";
Pend;
```

- `Model="sm.mdl";`  
Calculate in standard model  
Try to find in `grace.2.2.0/lib/model/sm.mdl`
- `ELWK=3;`  
The order of Electro-Weak coupling constants is 3
- `Initial={...}`  
Initial particles
- `Final={...}`  
Final particles
- `Kinem="2302";`  
Select kinematics : see `grace.2.2.0d/lib/dbkinem/doc`

## (d) Names of particles

Try

```
grep Particle grace.2.2.0/lib/model/sm.mdl
```

For the standard model:

particle	anti-p.	type	charge	color
W-plus	W-minus	vector	1	1
Z	Z	vector	0	1
Photon	Photon	vector	0	1
Gluon	Gluon	vector	0	1

particle	anti-p.	type	charge	color
Higgs	Higgs	scalar	0	1
chi-plus	chi-minus	scalar	1	1
chi-3	chi-3	scalar	0	1

particle	anti-p.	type	charge	color
nu-e	nu-e-bar	fermion	0	1
nu-mu	nu-mu-bar	fermion	0	1
nu-tau	nu-tau-bar	fermion	0	1
electron	positron	fermion	-1	1
muon	anti-muon	fermion	-1	1
tau	anti-tau	fermion	-1	1

particle	anti-p.	type	charge	color
u	u-bar	fermion	2/3	3
c	c-bar	fermion	2/3	3
t	t-bar	fermion	2/3	3
d	d-bar	fermion	-1/3	3
s	s-bar	fermion	-1/3	3
b	b-bar	fermion	-1/3	3

particle	anti-p.	type	charge	color
c-plus	c-plus-bar	ghost	1	1
c-minus	c-minus-bar	ghost	-1	1
c-z	c-z-bar	ghost	0	1
c-a	c-a-bar	ghost	0	1
c-g	c-g-bar	ghost	0	8

### 3. Feynman graphs generation:

**grc**

This command reads "in.prc" and creates a file "out.grf".

You can read "out.grf"

4. Draw Feynman graphs:

`gracefig`

or for creation of ps file:

`gracefig -p`

5. Fortran code generation

`grcfort`

The generated code is compiled by

`make make integ make spring`

or

`make all`

## 6. Gauge invariance check

`./gauge`

For the confirmation of the values of parameters: **uncomment** the line in "gauge.f"

`CC call prmass`

## 7. Integration over phase space

`./integ`

You should carefully examine the convergence behavior of the integration. If the convergence is not so good,

you must increase the number of sampling points or  
replace kinematics routines.

## 8. Event generation

`./spring`