

Plan

- Field theory : a short reminder
 - free fields (KG details, Fermion)
 - Scattering matrix in perturbation
 - Wick theorem to Feynman rules
- Why Monte-Carlo/automated tools?
- Lagrangian to the Feynman rules
 - Model file : Parameters, fields, gauge group and Lagrangian
 - Running FeynRules
- Demo

Why automated tools

- Algorithmic

- Less error prone

- Long $f^{abc} G_{\mu\nu}^a G^{b\nu\rho} G_\rho^{c\mu} \ni$ 4 gluons vertex

$$\begin{aligned}
 & 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_1^{\mu_4} p_2^{\mu_3} \eta_{\mu_1, \mu_2} - 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_1^{\mu_3} p_2^{\mu_4} \eta_{\mu_1, \mu_2} + 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_1^{\mu_3} p_3^{\mu_4} \eta_{\mu_1, \mu_2} + \\
 & 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_2^{\mu_3} p_3^{\mu_4} \eta_{\mu_1, \mu_2} + 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_1^{\mu_4} p_4^{\mu_3} \eta_{\mu_1, \mu_2} + 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_2^{\mu_4} p_4^{\mu_3} \eta_{\mu_1, \mu_2} - \\
 & 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} \eta_{\mu_3, \mu_4} p_1 \cdot p_3 \eta_{\mu_1, \mu_2} - 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} \eta_{\mu_3, \mu_4} p_1 \cdot p_4 \eta_{\mu_1, \mu_2} - 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} \eta_{\mu_3, \mu_4} p_2 \cdot p_3 \eta_{\mu_1, \mu_2} - \\
 & 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} \eta_{\mu_3, \mu_4} p_2 \cdot p_4 \eta_{\mu_1, \mu_2} + 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_1^{\mu_2} p_2^{\mu_4} \eta_{\mu_1, \mu_3} + 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_1^{\mu_4} p_3^{\mu_2} \eta_{\mu_1, \mu_3} - \\
 & 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_2^{\mu_4} p_3^{\mu_2} \eta_{\mu_1, \mu_3} - 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_1^{\mu_2} p_3^{\mu_4} \eta_{\mu_1, \mu_3} - 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_1^{\mu_4} p_4^{\mu_2} \eta_{\mu_1, \mu_3} + \\
 & 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_3^{\mu_4} p_4^{\mu_2} \eta_{\mu_1, \mu_3} - 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_1^{\mu_2} p_2^{\mu_3} \eta_{\mu_1, \mu_4} - 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_1^{\mu_3} p_3^{\mu_2} \eta_{\mu_1, \mu_4} + \\
 & 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_1^{\mu_3} p_4^{\mu_2} \eta_{\mu_1, \mu_4} - 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_2^{\mu_3} p_4^{\mu_2} \eta_{\mu_1, \mu_4} - 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_1^{\mu_2} p_4^{\mu_3} \eta_{\mu_1, \mu_4} - \\
 & 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_3^{\mu_3} p_4^{\mu_2} \eta_{\mu_1, \mu_4} - 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_1^{\mu_4} p_2^{\mu_1} \eta_{\mu_2, \mu_3} - 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_1^{\mu_4} p_3^{\mu_1} \eta_{\mu_2, \mu_3} + \\
 & 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_2^{\mu_4} p_3^{\mu_1} \eta_{\mu_2, \mu_3} - 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_2^{\mu_1} p_4^{\mu_3} \eta_{\mu_2, \mu_3} - 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_2^{\mu_4} p_4^{\mu_1} \eta_{\mu_2, \mu_3} - \\
 & 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_3^{\mu_4} p_4^{\mu_1} \eta_{\mu_2, \mu_3} + 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_1^{\mu_3} p_2^{\mu_1} \eta_{\mu_2, \mu_4} - 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_2^{\mu_3} p_3^{\mu_1} \eta_{\mu_2, \mu_4} - \\
 & 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_1^{\mu_3} p_4^{\mu_1} \eta_{\mu_2, \mu_4} + 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_2^{\mu_3} p_4^{\mu_1} \eta_{\mu_2, \mu_4} - 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_2^{\mu_1} p_4^{\mu_3} \eta_{\mu_2, \mu_4} + \\
 & 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_3^{\mu_1} p_4^{\mu_3} \eta_{\mu_2, \mu_4} + 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_1^{\mu_2} p_3^{\mu_1} \eta_{\mu_3, \mu_4} + 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_2^{\mu_1} p_3^{\mu_2} \eta_{\mu_3, \mu_4} + \\
 & 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} p_1^{\mu_2} p_4^{\mu_1} \eta_{\mu_3, \mu_4} + 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_3^{\mu_2} p_4^{\mu_1} \eta_{\mu_3, \mu_4} + 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} p_2^{\mu_1} p_4^{\mu_2} \eta_{\mu_3, \mu_4} - \\
 & 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} p_3^{\mu_1} p_4^{\mu_2} \eta_{\mu_3, \mu_4} + 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} \eta_{\mu_1, \mu_4} \eta_{\mu_2, \mu_3} p_1 \cdot p_2 - 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} \eta_{\mu_1, \mu_3} \eta_{\mu_2, \mu_4} p_1 \cdot p_2 + \\
 & 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} \eta_{\mu_1, \mu_4} \eta_{\mu_2, \mu_3} p_1 \cdot p_3 + 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} \eta_{\mu_1, \mu_3} \eta_{\mu_2, \mu_4} p_1 \cdot p_4 + 6ig_s f_{a_1, a_4, a} f_{a_2, a_3, a} \eta_{\mu_1, \mu_3} \eta_{\mu_2, \mu_4} p_2 \cdot p_3 + \\
 & 6ig_s f_{a_1, a_3, a} f_{a_2, a_4, a} \eta_{\mu_1, \mu_4} \eta_{\mu_2, \mu_3} p_2 \cdot p_4 + 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} \eta_{\mu_1, \mu_4} \eta_{\mu_2, \mu_3} p_3 \cdot p_4 - 6ig_s f_{a_1, a_2, a} f_{a_3, a_4, a} \eta_{\mu_1, \mu_3} \eta_{\mu_2, \mu_4} p_3 \cdot p_4
 \end{aligned}$$

Many diagrams

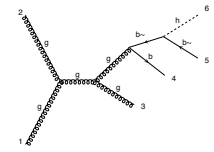


diagram 7 QCD=3, QED=1

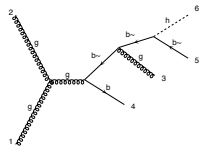


diagram 8 QCD=3, QED=1

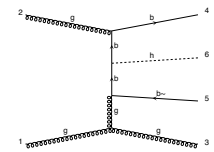


diagram 9 QCD=3, QED=1

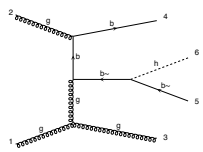


diagram 10 QCD=3, QED=1

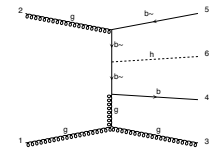


diagram 11 QCD=3, QED=1

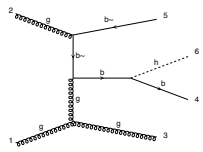


diagram 12 QCD=3, QED=1

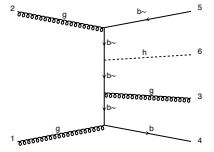


diagram 19 QCD=3, QED=1

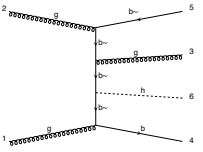


diagram 20 QCD=3, QED=1

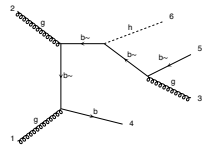


diagram 21 QCD=3, QED=1

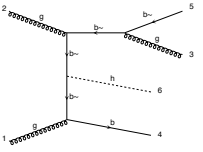


diagram 22 QCD=3, QED=1

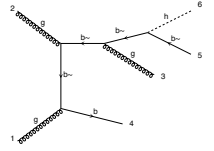


diagram 23 QCD=3, QED=1

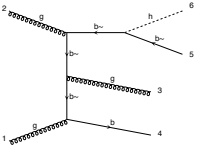


diagram 24 QCD=3, QED=1

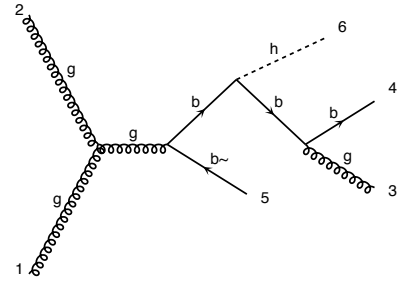


diagram 1 QCD=3, QED=1

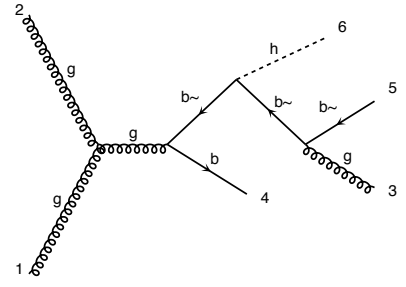


diagram 3 QCD=3, QED=1

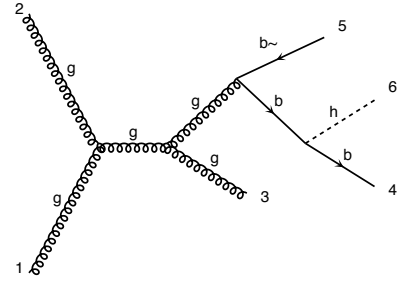


diagram 5 QCD=3, QED=1

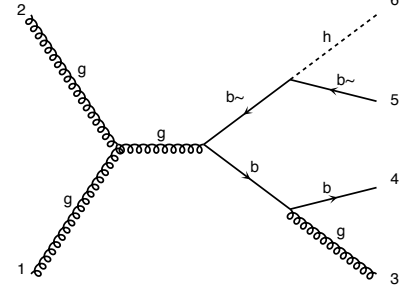


diagram 2 QCD=3, QED=1

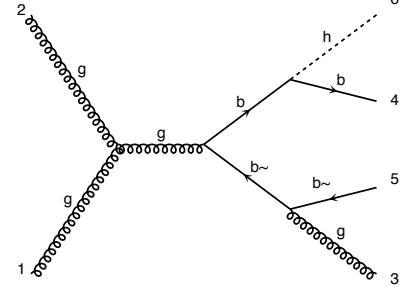


diagram 4 QCD=3, QED=1

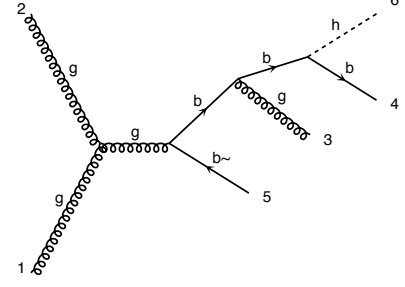


diagram 6 QCD=3, QED=1

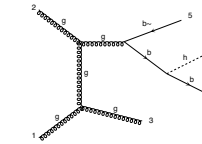


diagram 13 QCD=3, QED=1

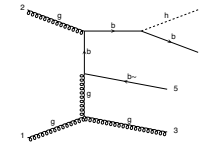


diagram 14 QCD=3, QED=1

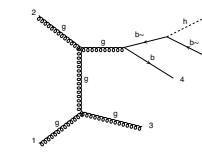


diagram 15 QCD=3, QED=1

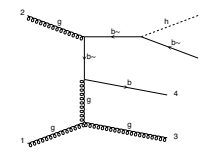


diagram 16 QCD=3, QED=1

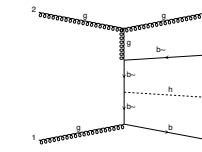


diagram 17 QCD=3, QED=1

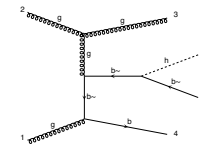


diagram 18 QCD=3, QED=1

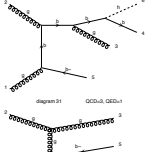


diagram 31 QCD=3, QED=1

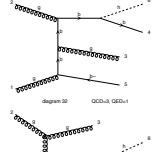


diagram 32 QCD=3, QED=1

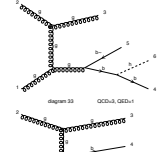


diagram 33 QCD=3, QED=1

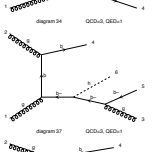


diagram 34 QCD=3, QED=1

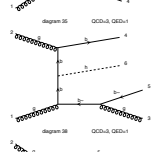


diagram 35 QCD=3, QED=1

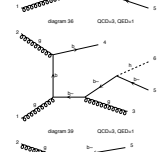


diagram 36 QCD=3, QED=1

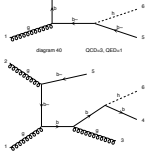


diagram 37 QCD=3, QED=1

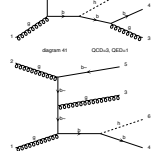


diagram 38 QCD=3, QED=1

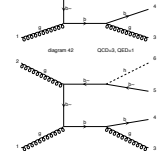


diagram 39 QCD=3, QED=1

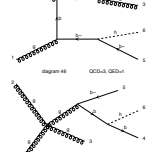


diagram 40 QCD=3, QED=1

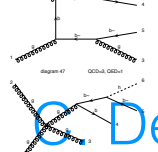


diagram 41 QCD=3, QED=1

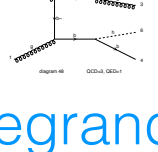
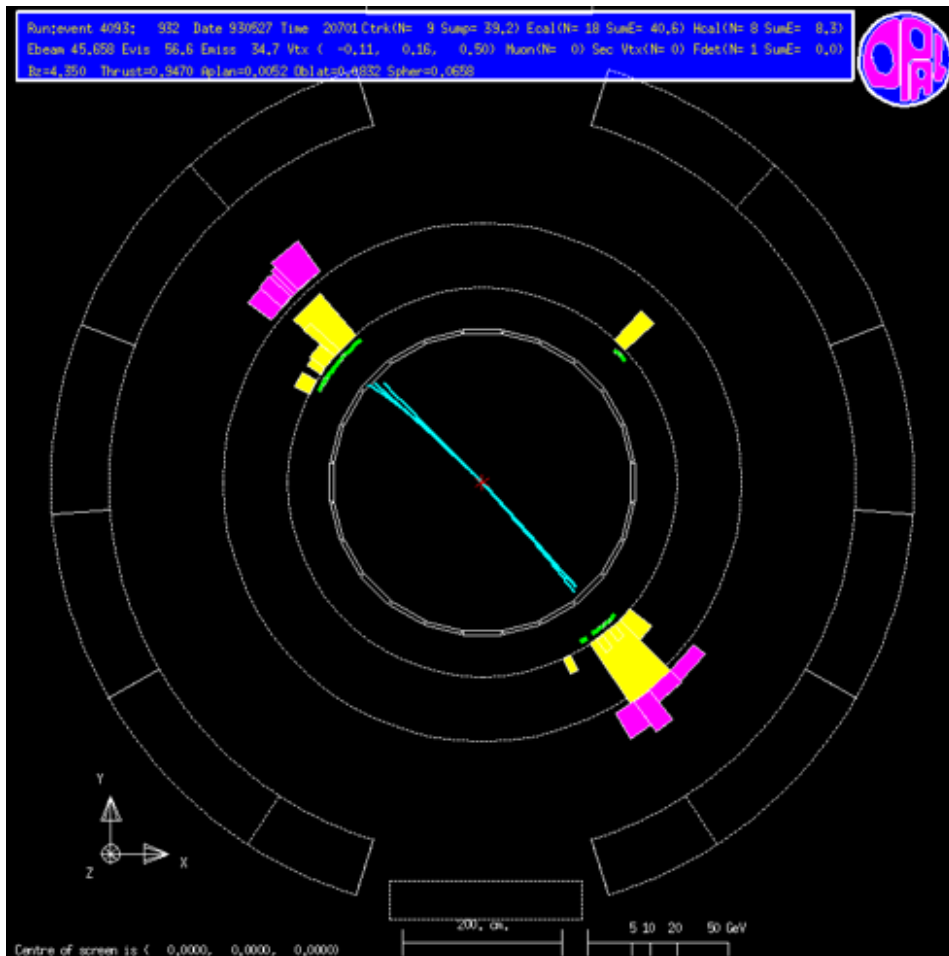


diagram 42 QCD=3, QED=1

Hadron colliders

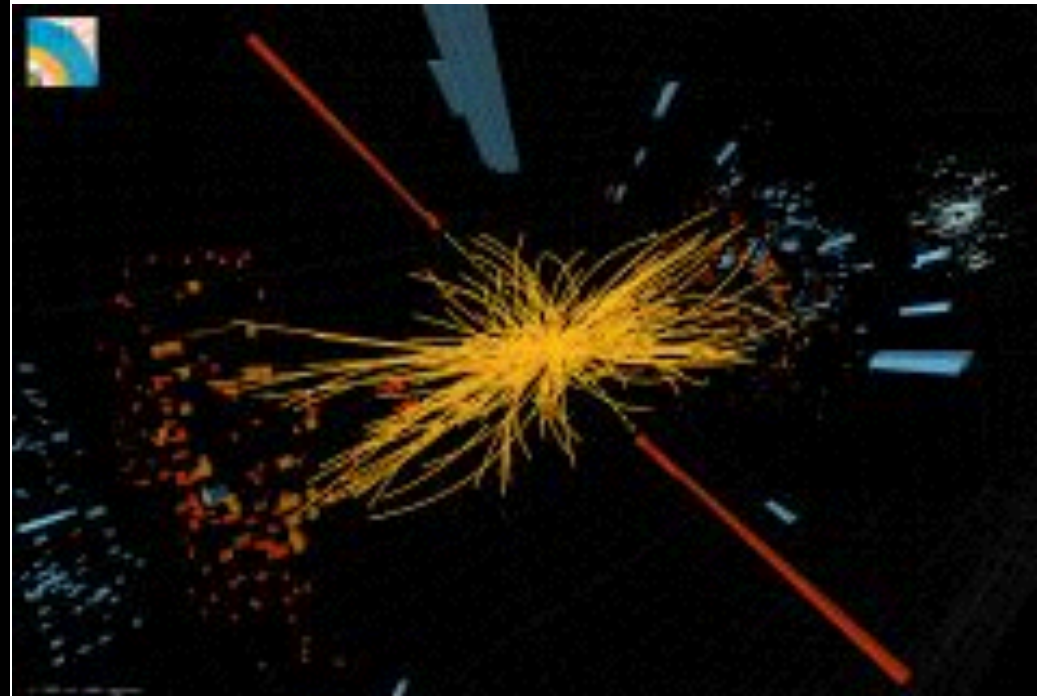
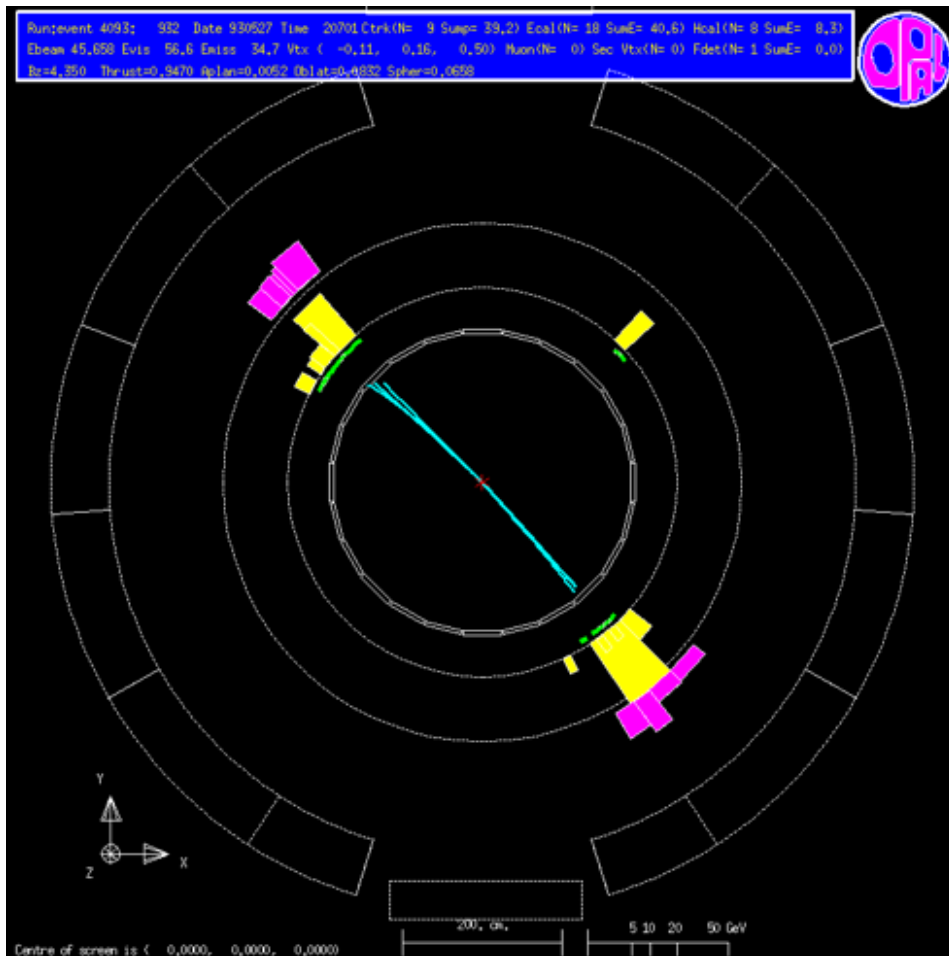
LEP



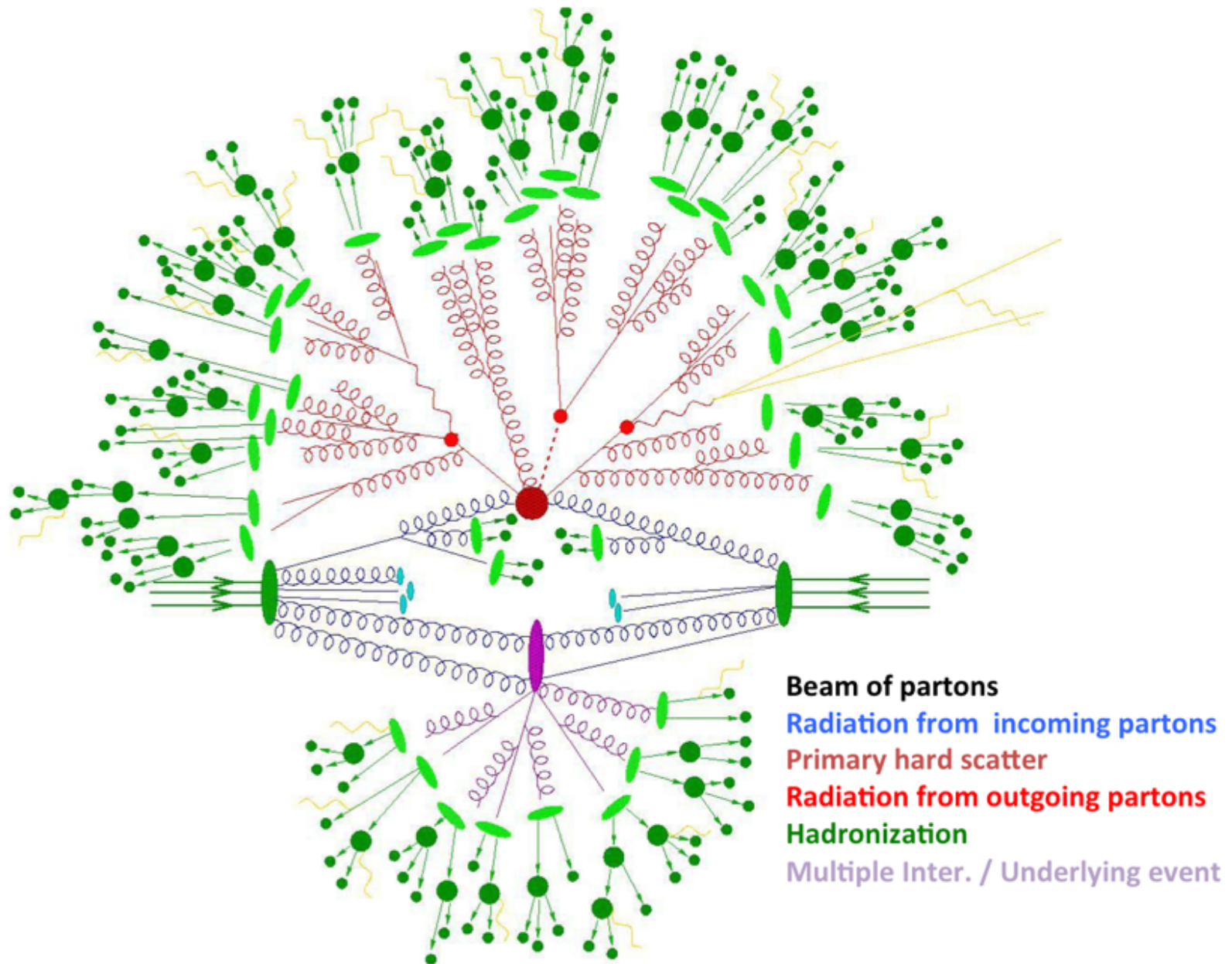
Hadron colliders

LEP

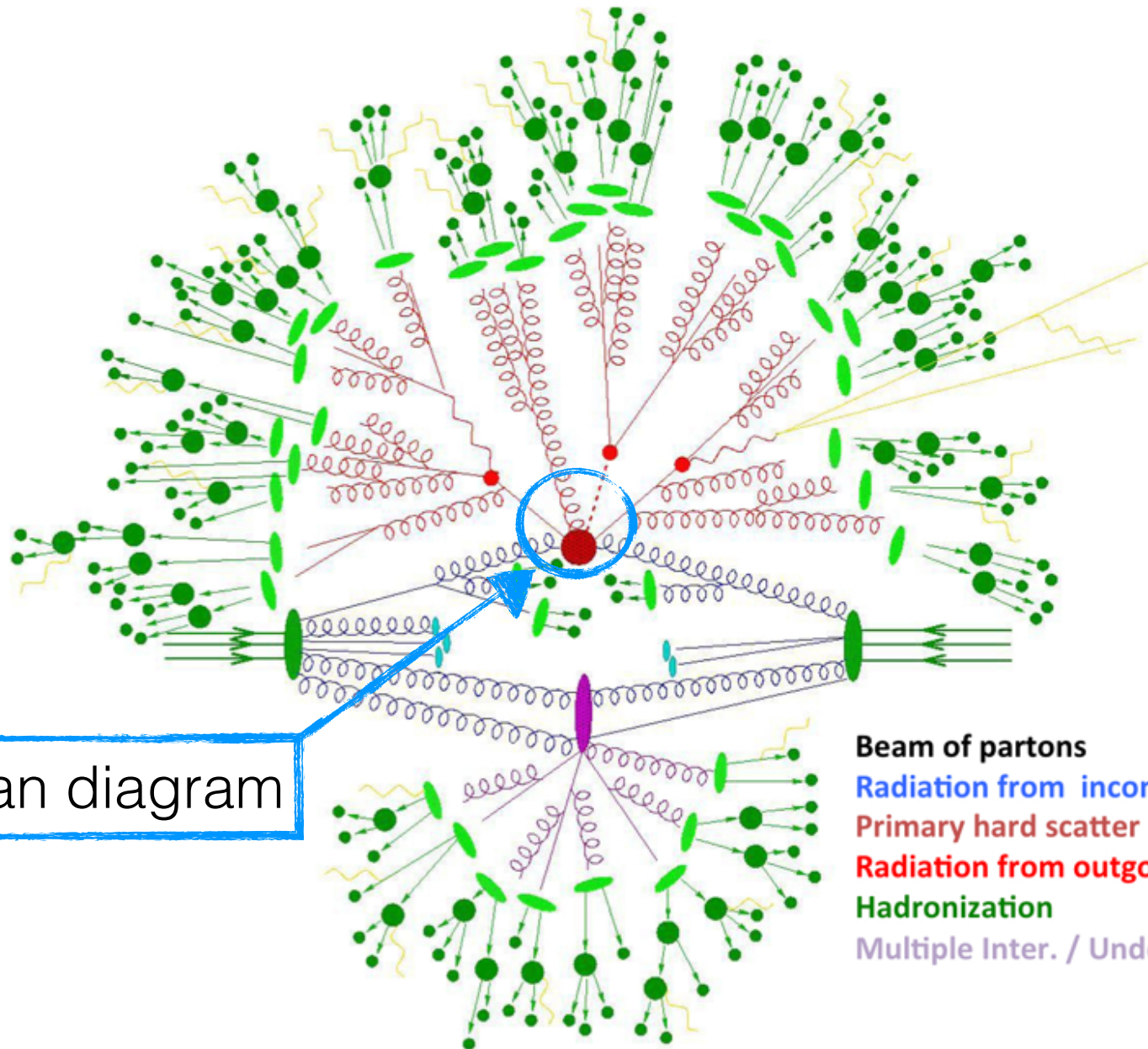
LHC



Hadron collider event



Hadron collider event

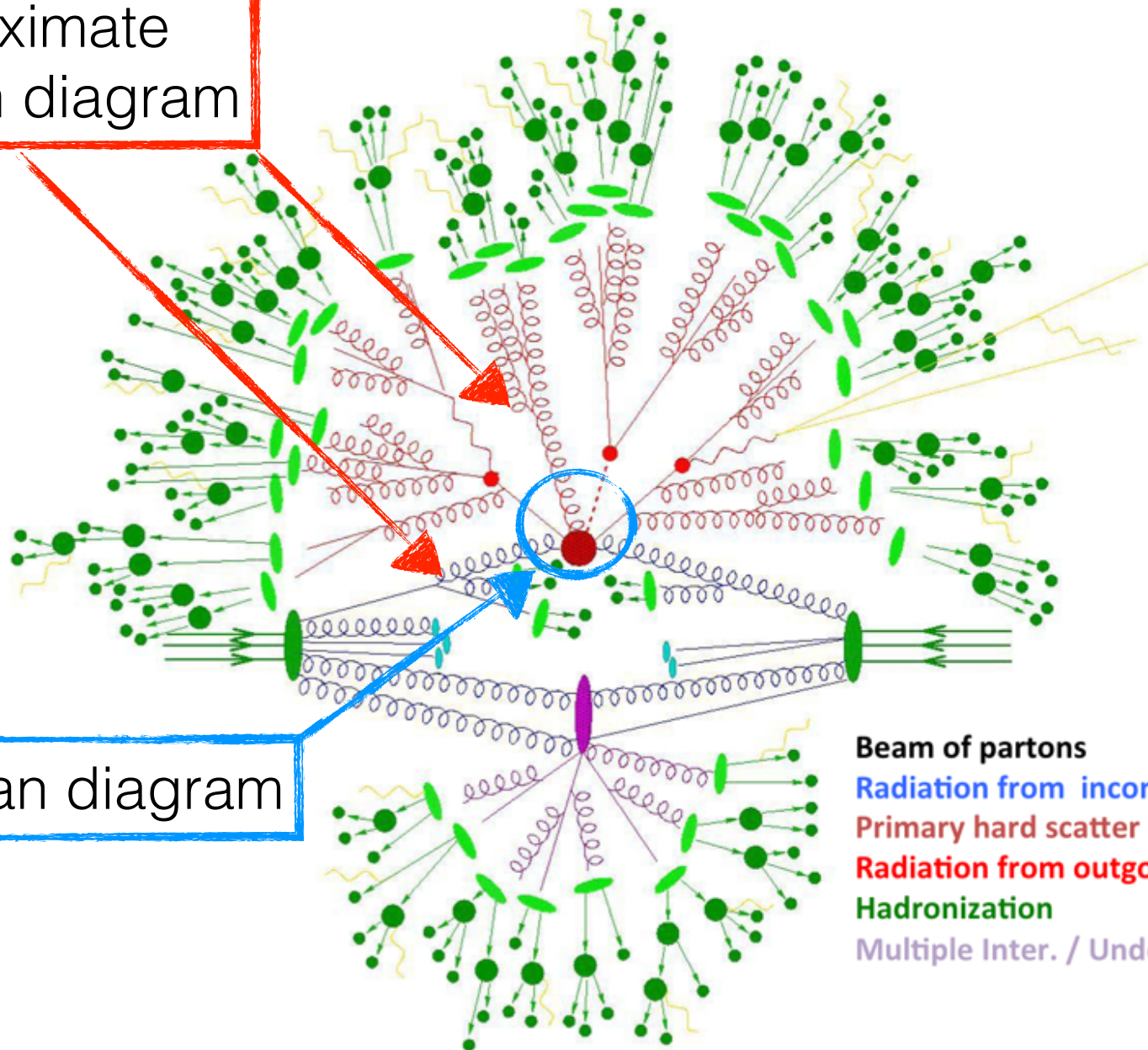


Feynman diagram

- Beam of partons
- Radiation from incoming partons
- Primary hard scatter
- Radiation from outgoing partons
- Hadronization
- Multiple Inter. / Underlying event

Hadron collider event

Approximate
Feynman diagram



Feynman diagram

Beam of partons

Radiation from incoming partons

Primary hard scatter

Radiation from outgoing partons

Hadronization

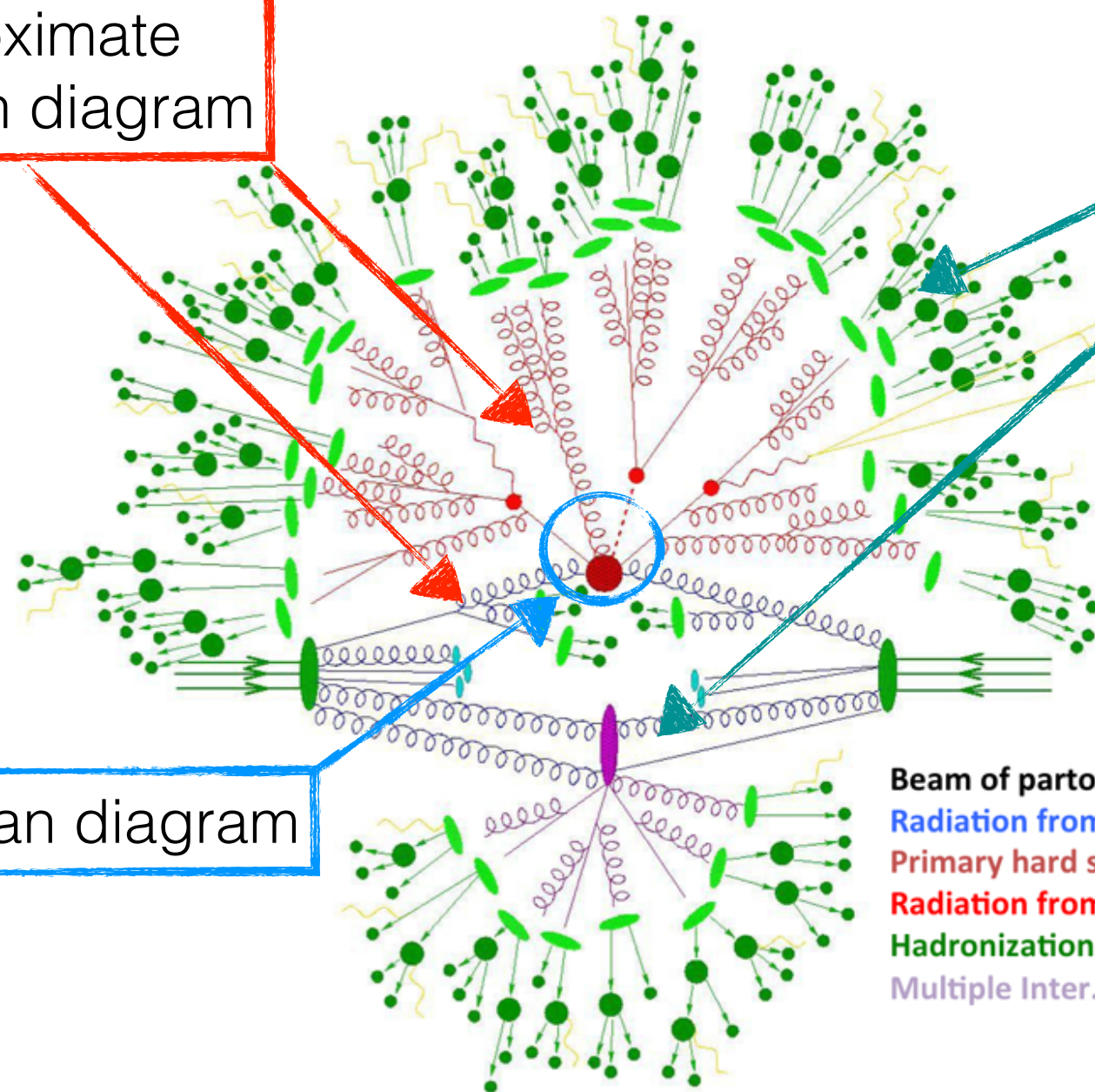
Multiple Inter. / Underlying event

Hadron collider event

Approximate Feynman diagram

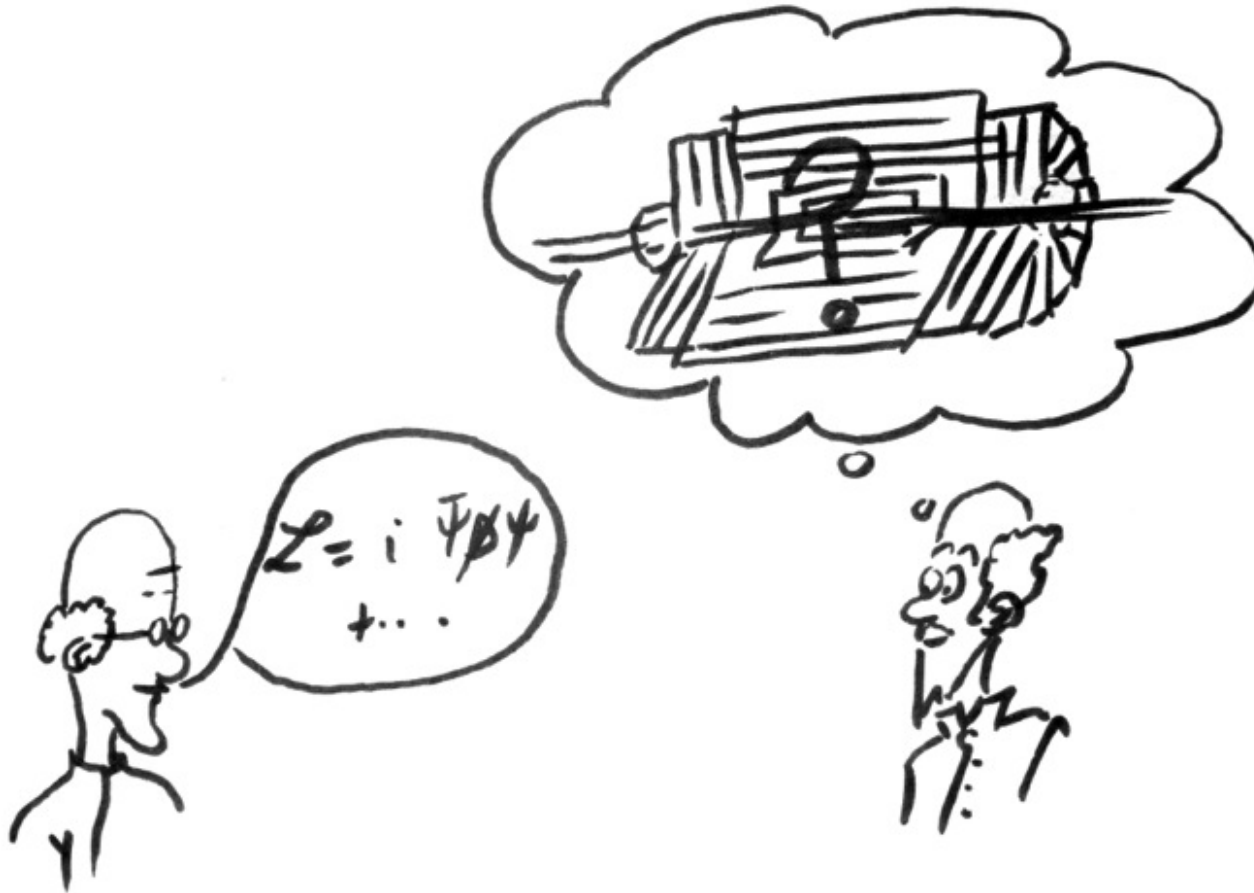
Models

Feynman diagram



- Beam of partons
- Radiation from incoming partons
- Primary hard scatter
- Radiation from outgoing partons
- Hadronization
- Multiple Inter. / Underlying event

Why BSM simulation?

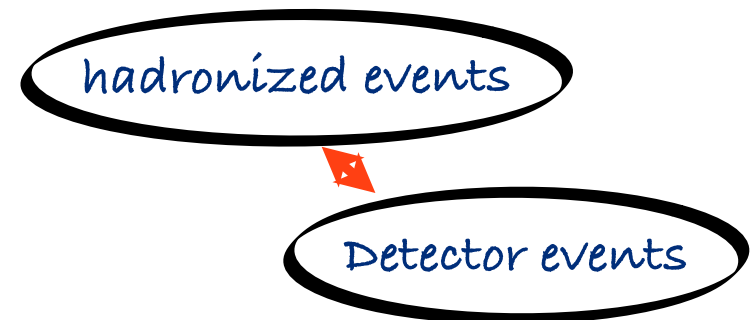


BSM simulation

Detector events

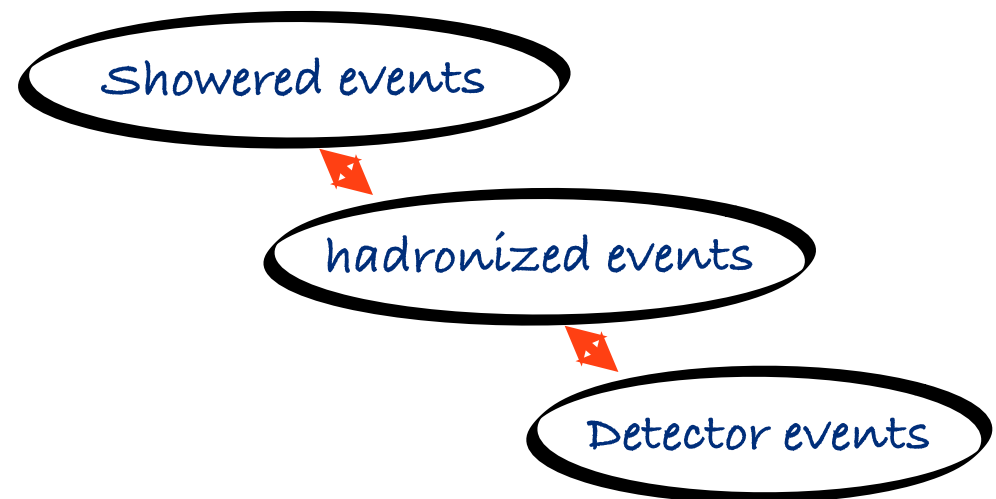
C. Degrande

BSM simulation

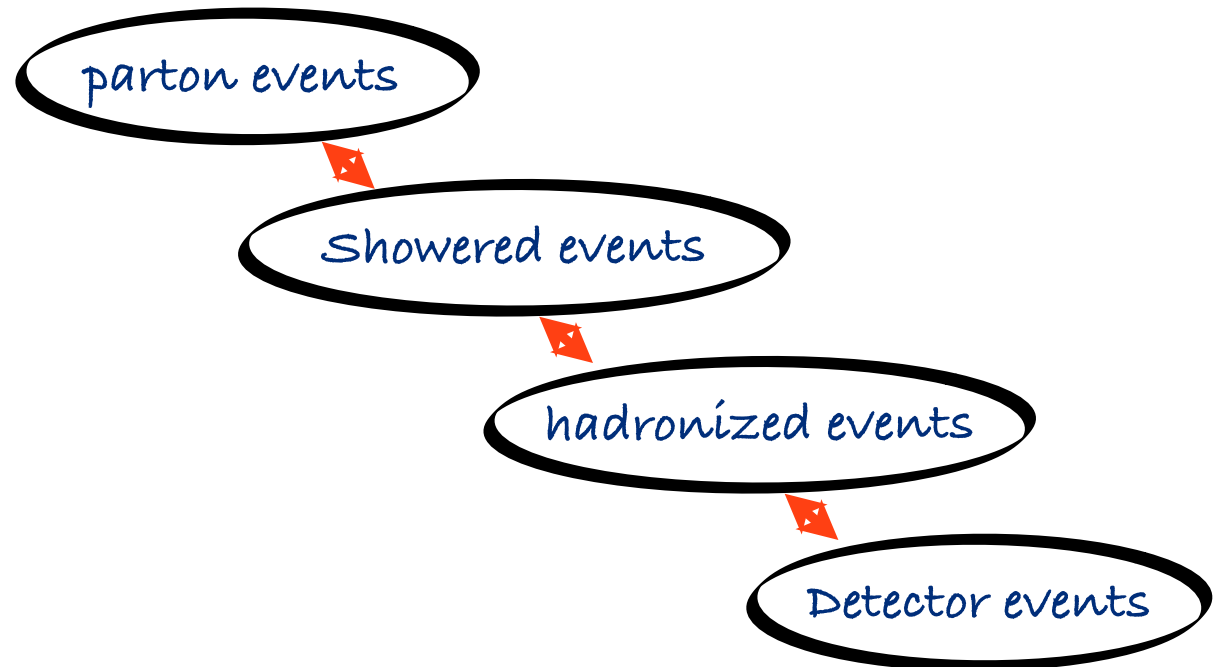


C. Degrande

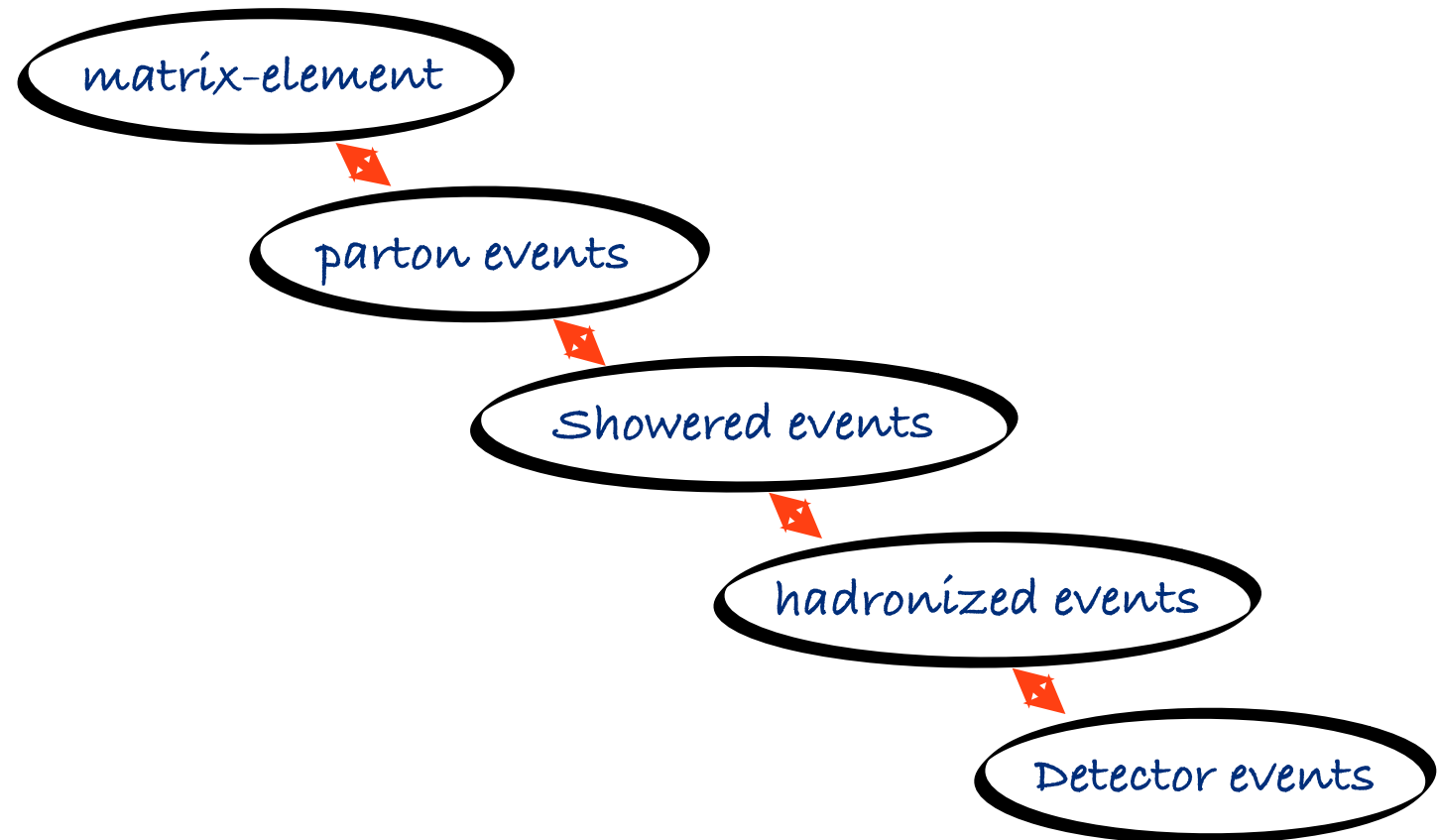
BSM simulation



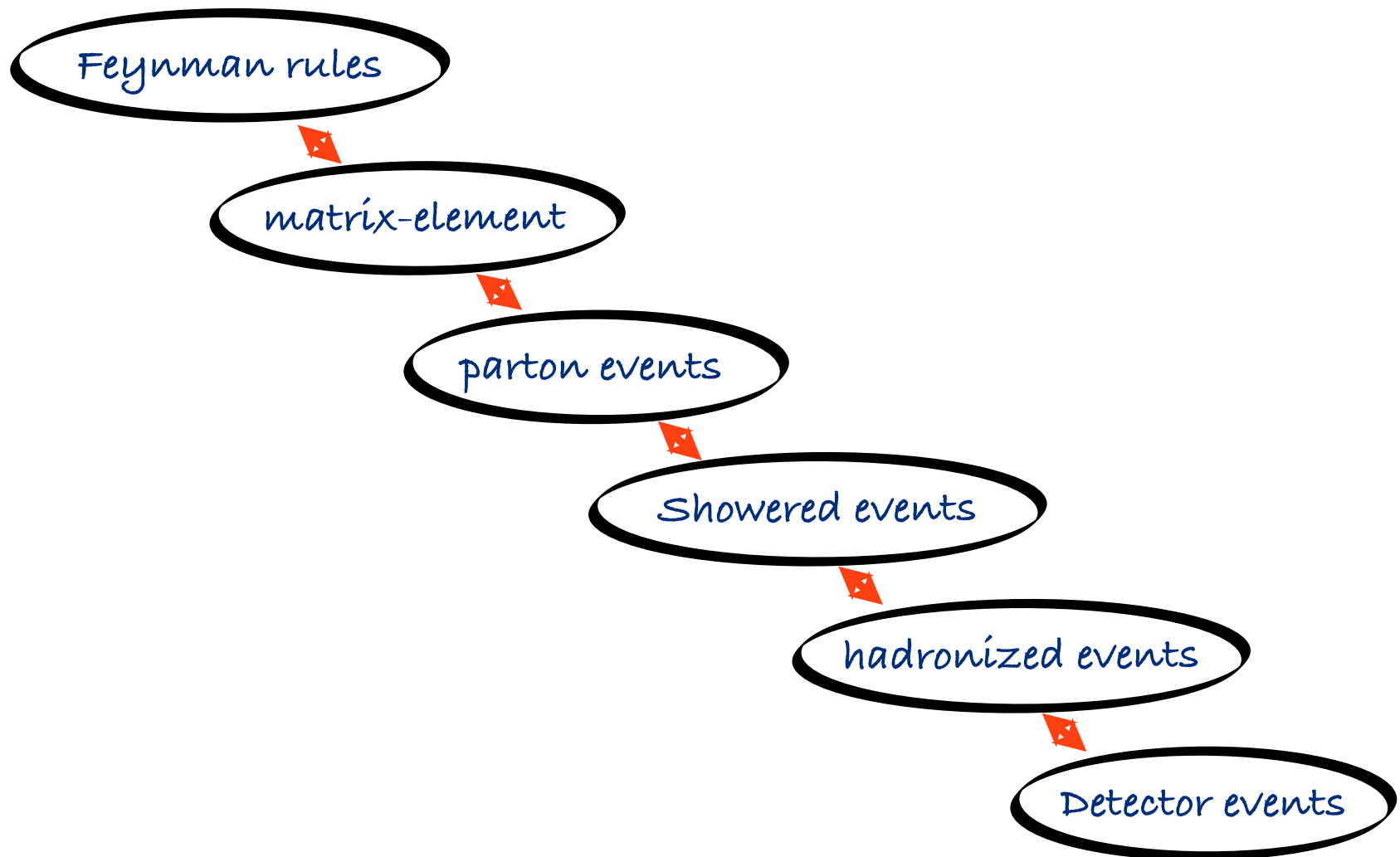
BSM simulation



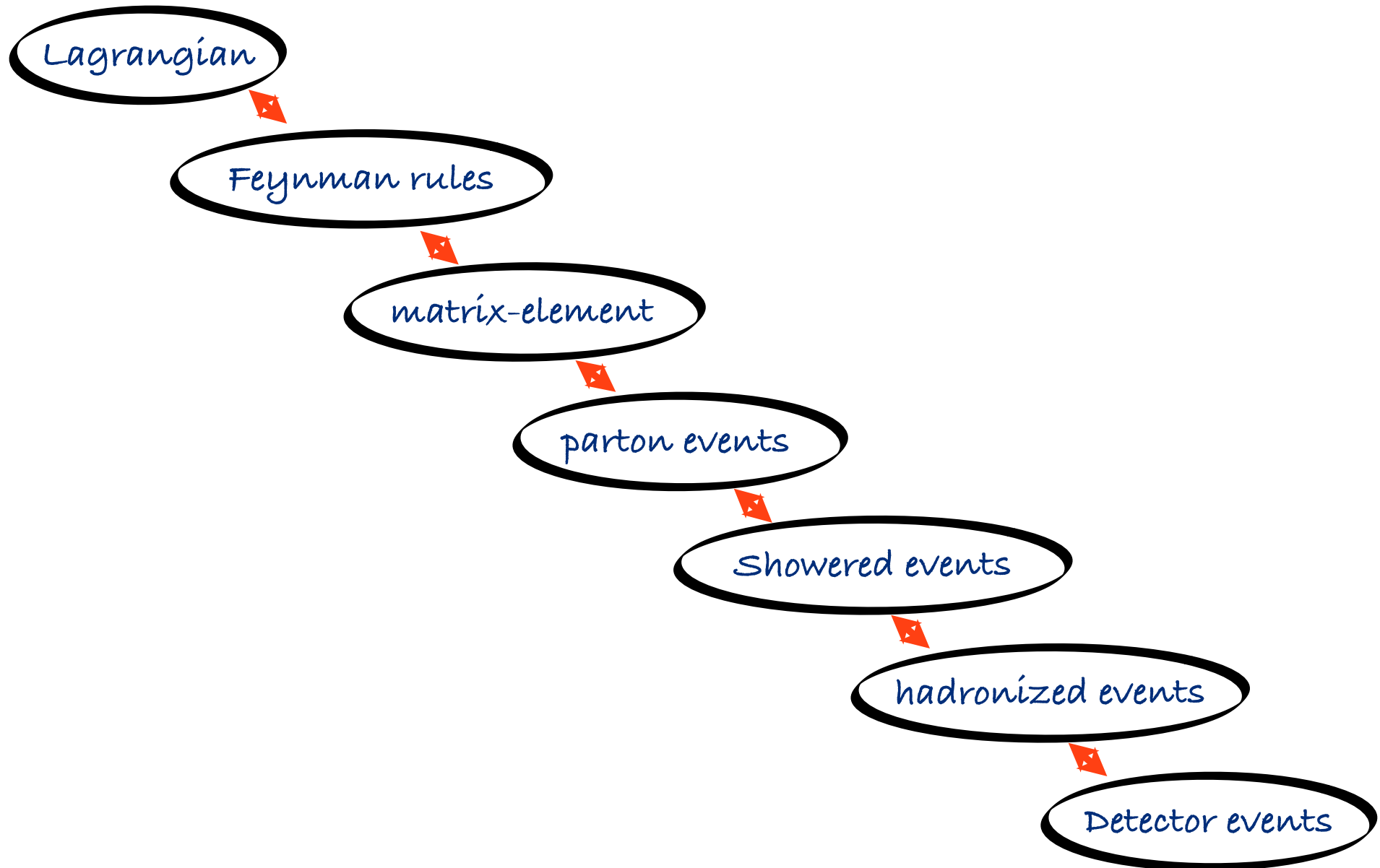
BSM simulation



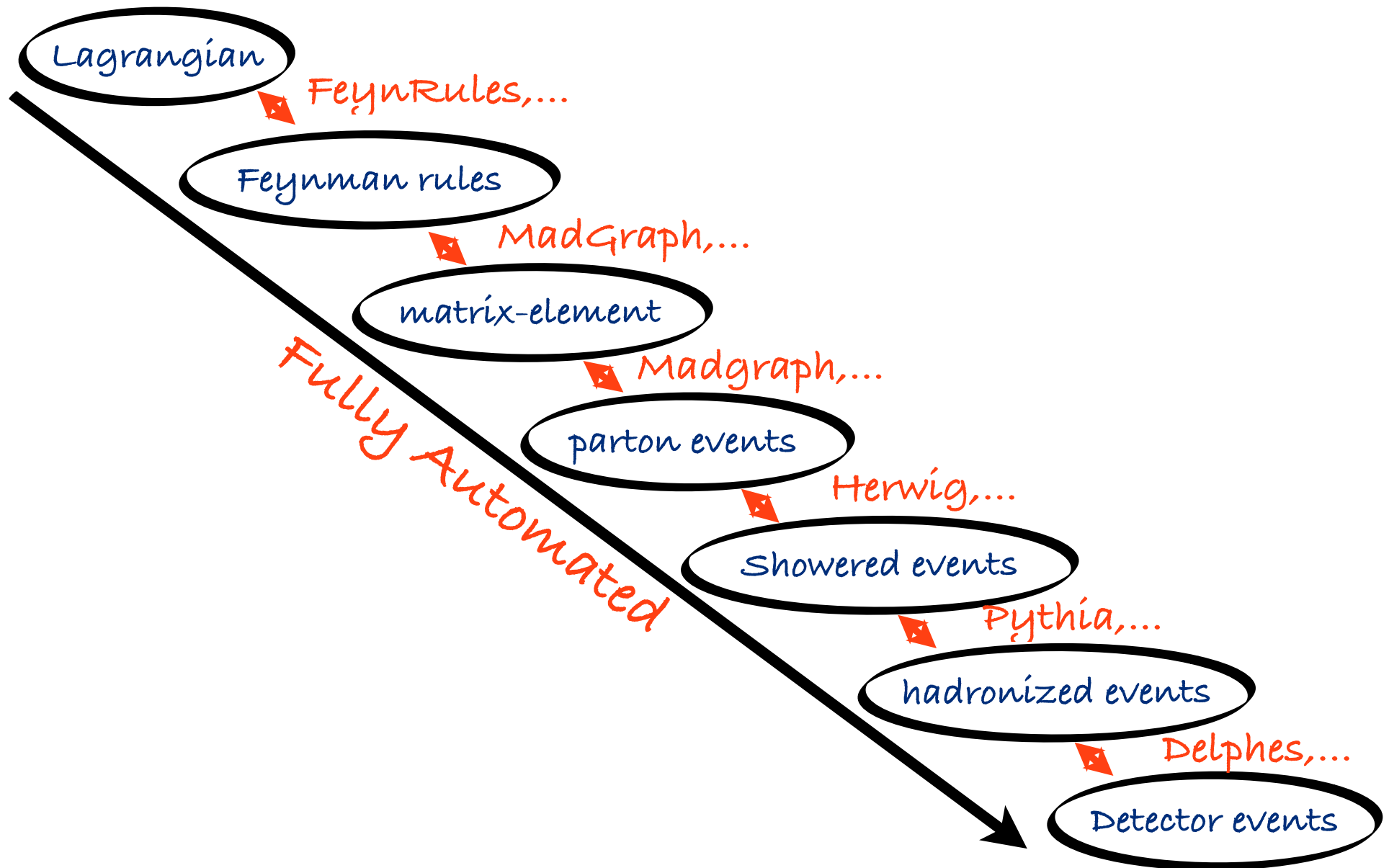
BSM simulation



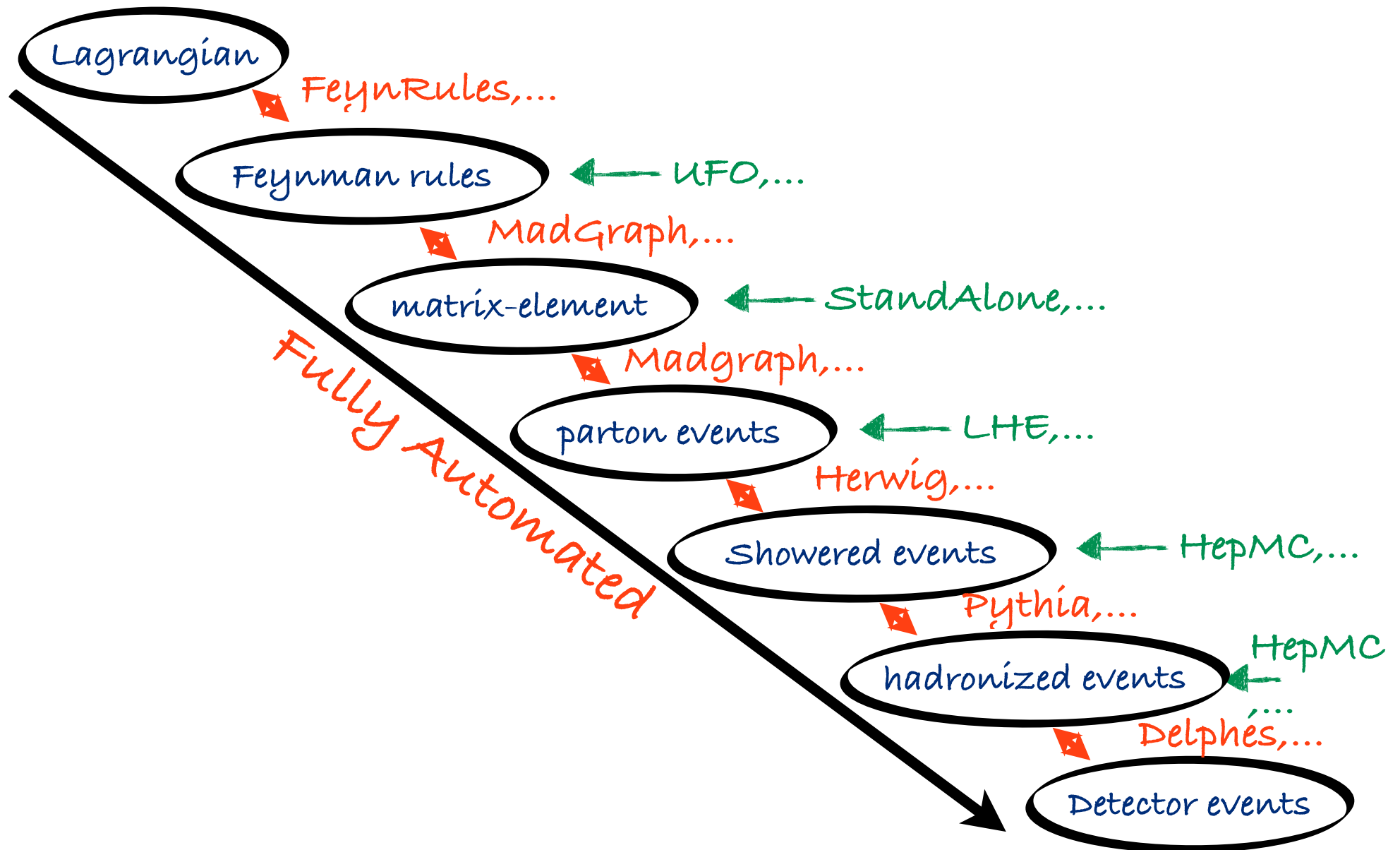
BSM simulation



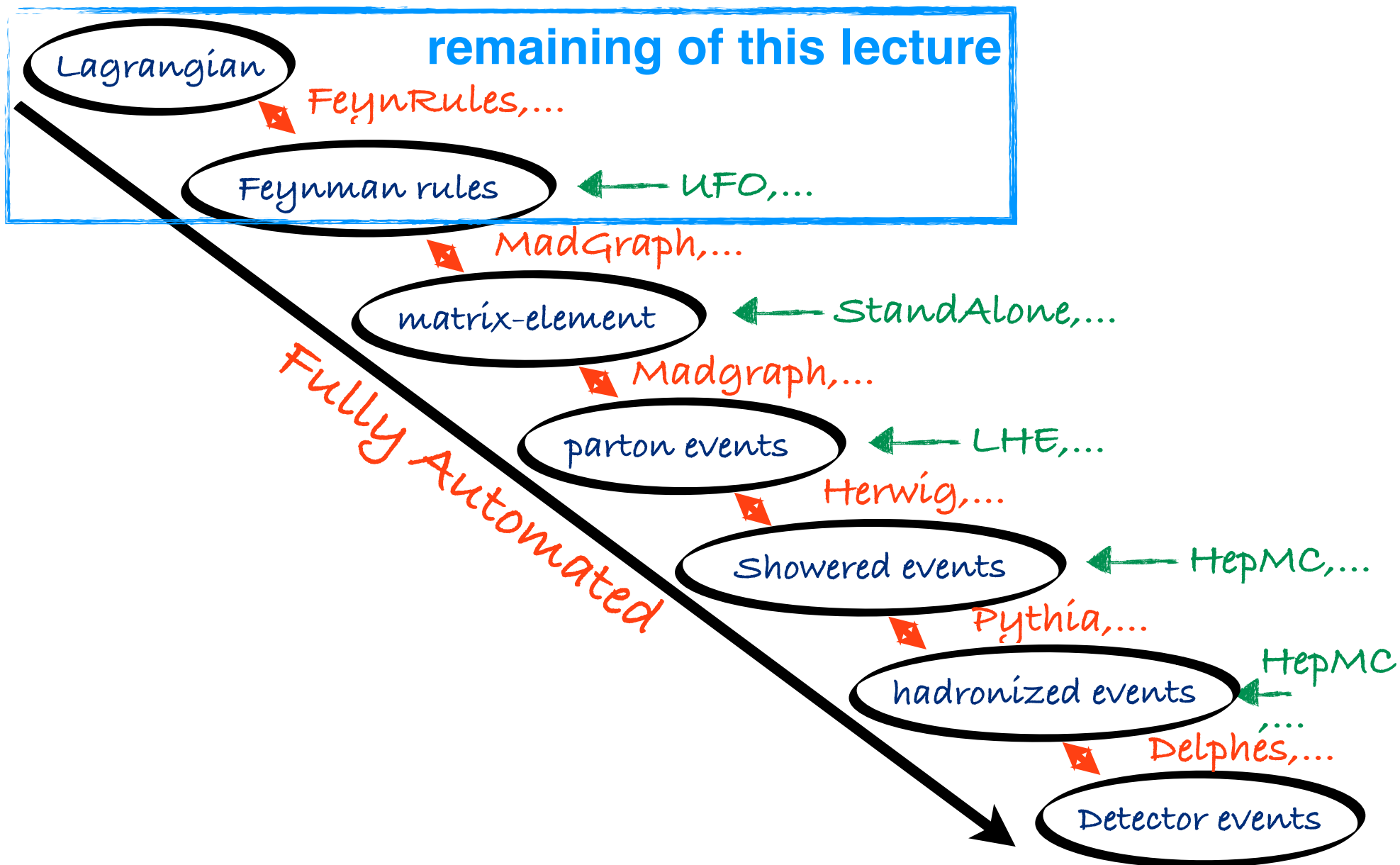
BSM simulation



BSM simulation



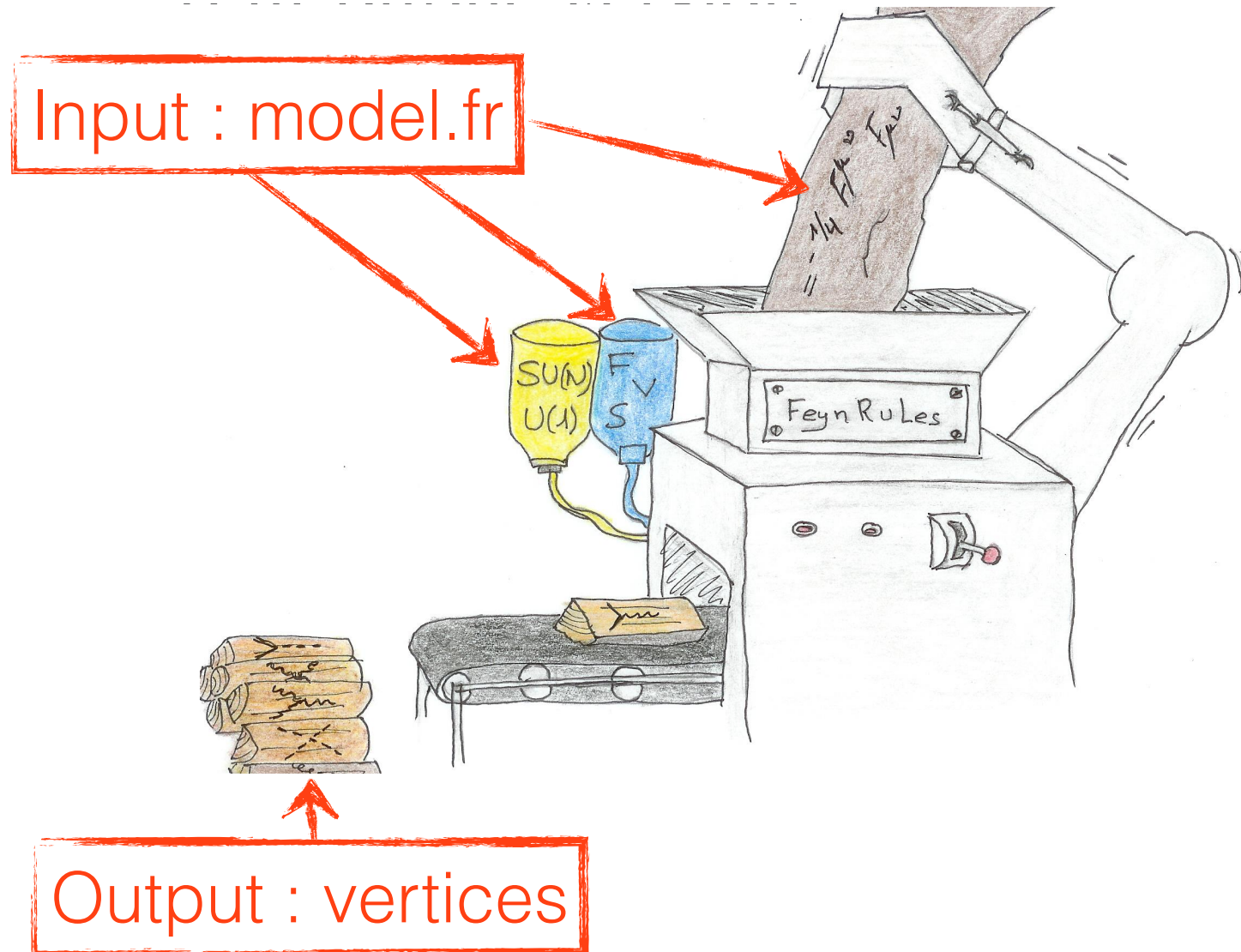
BSM simulation



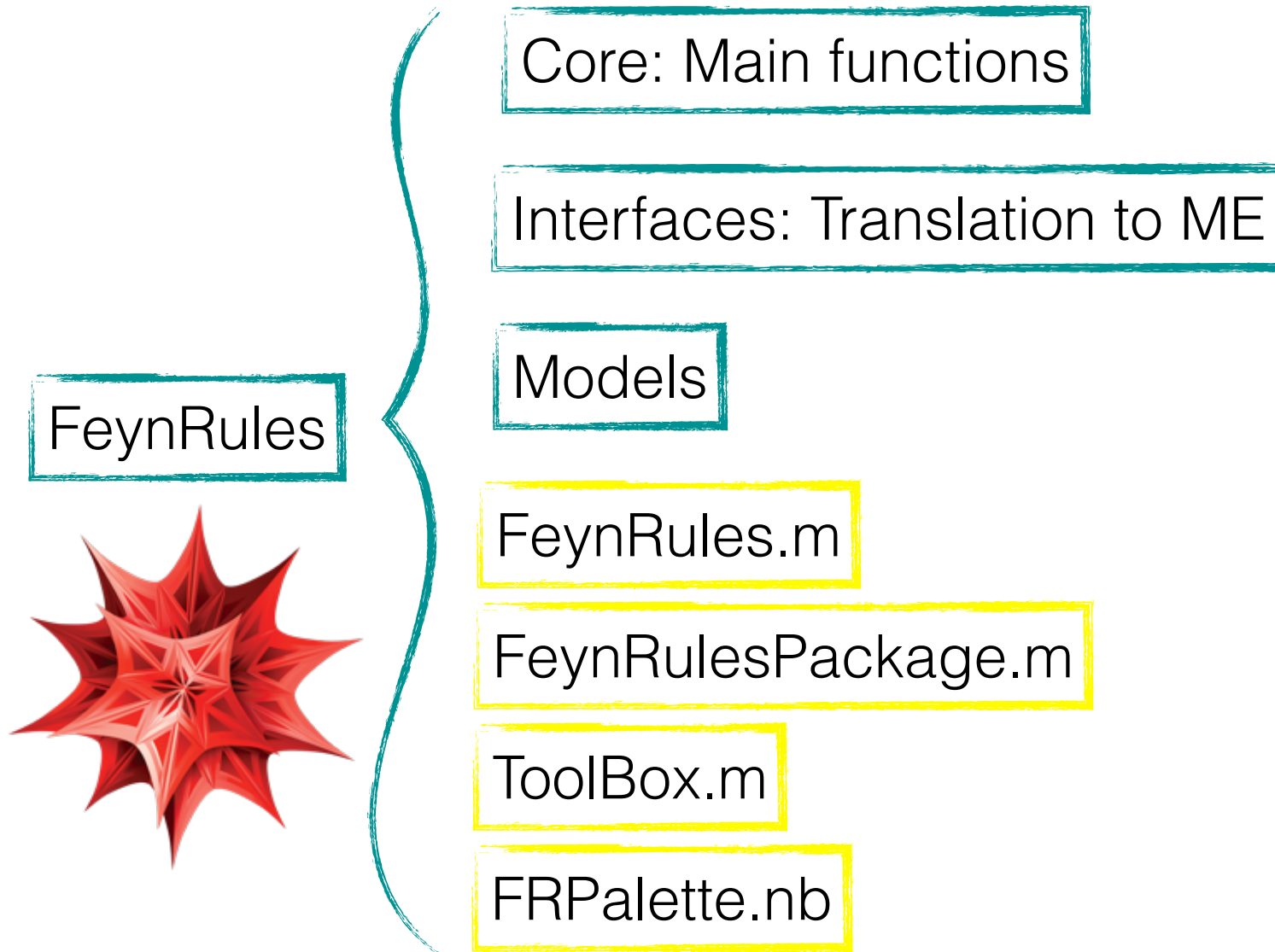
Plan

- Field theory : a short reminder
 - free fields (KG details, Fermion)
 - Scattering matrix in perturbation
 - Wick theorem to Feynman rules
- Why Monte-Carlo/automated tools?
- Lagrangian to the Feynman rules
 - Model file : Parameters, fields, gauge group and Lagrangian
 - Running FeynRules
- Demo

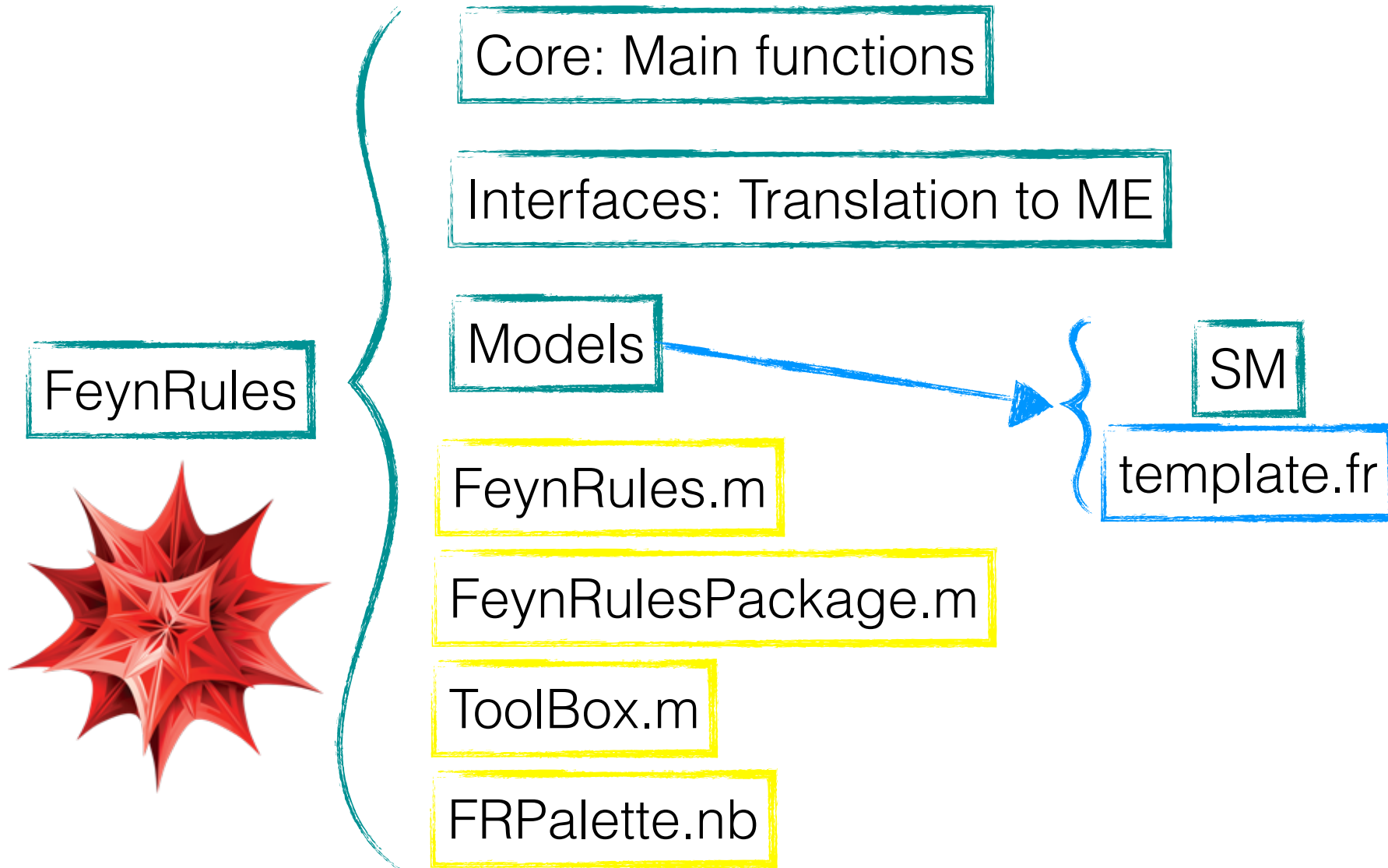
FeynRules in a nutshell



FeynRules structure



FeynRules structure



model file

```
(***** This is a template model file for FeynRules *****)
```

```
(***** Index definition *****)
```

```
IndexRange[ Index[Generation] ] = Range[3]
```

```
IndexFormat[Generation, f]
```

```
(***** Parameter list *****)
```

```
M$Parameters = {  
}  
(***** Gauge group list *****)
```

Definition of variables
in Mathematica syntaxe

```
M$GaugeGroups = {  
}  
(***** Particle classes list *****)
```

```
M$ClassesDescription = {  
}
```

Model information

```
M$ModelName = "my_new_model";
```

```
M$Information = {  
  Authors      -> {"Mr. X", "Ms. Y"},  
  Institutions -> {"UC Louvain"},  
  Emails       -> {"X@uclouvain.be", "Y@uclouvain.be"},  
  Date         -> "01.03.2013",  
  References   -> {"reference 1", "reference 2"},  
  URLs         -> {"http://feynrules.irmp.ucl.ac.be"},  
  Version      -> "1.0"  
};
```

Good practice for credit, issue(s) tracking

Indices definition

**Used in parameters, gauge groups
and fields**

Tells FR to remplace
summed indices by
the explicite sum

```
IndexRange[ Index[Colour] ] = Range[3];  
IndexRange[ Index[SU2W] ] = Unfold[ Range[3] ];  
IndexRange[ Index[Gluon] ] = NoUnfold[ Range[8] ];
```

Indices definition

Used in parameters, gauge groups
and fields

```
IndexRange[ Index[Colour] ] = Range[3];  
IndexRange[ Index[SU2W] ] = Unfold[ Range[3] ];  
IndexRange[ Index[Gluon] ] = NoUnfold[ Range[8] ];
```

Tells FR to remplace
summed indices by
the explicite sum

Tells FA/FC **not** to
remplace summed
indices by the
explicite sum

Indices definition

Used in parameters, gauge groups
and fields

```
IndexRange[ Index[Colour] ] = Range[3];  
IndexRange[ Index[SU2W] ] = Unfold[ Range[3] ];  
IndexRange[ Index[Gluon] ] = NoUnfold[ Range[8] ];
```

Tells FR to remplace
summed indices by
the explicite sum

Format:

```
IndexStyle[ Colour, i ];  
IndexStyle[ Gluon, a ];
```

Tells FA/FC **not** to
remplace summed
indices by the
explicite sum

Indices definition

Used in parameters, gauge groups
and fields

```
IndexRange[ Index[Colour] ] = Range[3];  
IndexRange[ Index[SU2W] ] = Unfold[ Range[3] ];  
IndexRange[ Index[Gluon] ] = NoUnfold[ Range[8] ];
```

Tells FR to remplace summed indices by the explicite sum

Format:

```
IndexStyle[ Colour, i ];  
IndexStyle[ Gluon, a ];
```

Tells FA/FC **not** to remplace summed indices by the explicite sum

Predefined indices: Lorentz, Spin, Spin1, Spin2

Parameters definition

```
M$Parameters = {  
    param1 == { options1 },  
    param2 == { options2 },  
    ...  
};
```

```
aEWM1 == {  
    ParameterType -> External,  
    BlockName     -> SMINPUTS,  
    OrderBlock    -> 1,  
    Value         -> 127.9,  
    InteractionOrder -> {QED,-2},  
    Description    -> "Inverse of the EW coupling constant at the Z  
pole"  
},
```

Compulsory!

Numerical value

Parameters definition

```
M$Parameters = {  
  param1 == { options1 },  
  param2 == { options2 },  
  ...  
};
```

```
MW == {  
  ParameterType -> Internal,  
  Value -> Sqrt[MZ^2/2+Sqrt[MZ^4/4-Pi/Sqrt[2]*aEW/  
Gf*MZ^2]],  
  TeX -> Subscript[M,W],  
  Description -> "W mass"  
},
```

Expression

Parameters definition


```
M$Parameters = {  
    param1 == { options1 },  
    param2 == { options2 },  
    ...  
};
```

```
aEWM1 == {  
    ParameterType -> External,  
    BlockName     -> SMINPUTS,  
    OrderBlock    -> 1,  
    Value         -> 127.9,  
    InteractionOrder -> {QED,-2},  
    Description    -> "Inverse of the EW coupling constant at the Z  
pole"  
},
```

Parameters definition

```
M$Parameters = {  
    param1 == { options1 },  
    param2 == { options2 },  
    ...  
};
```

```
aEWM1 == {  
    ParameterType -> External,  
    BlockName     -> SMINPUTS,  
    OrderBlock    -> 1,  
    Value         -> 127.9,  
    InteractionOrder -> {QED,-2},  
    Description    -> "Inverse of the EW coupling constant at the Z  
pole"  
},
```

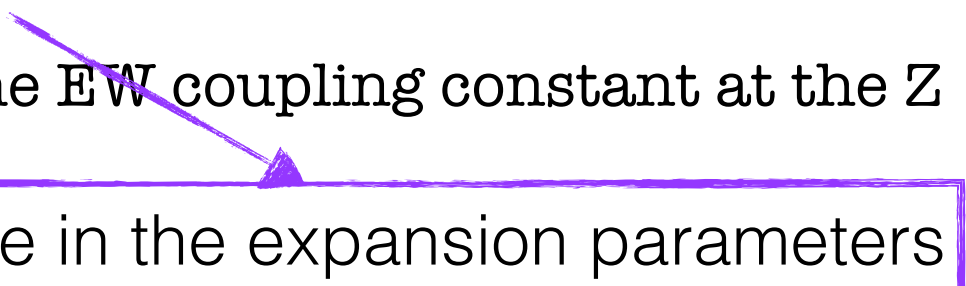


For the LHA cards

Parameters definition

```
M$Parameters = {  
    param1 == { options1 },  
    param2 == { options2 },  
    ...  
};
```

```
aEWM1 == {  
    ParameterType -> External,  
    BlockName     -> SMINPUTS,  
    OrderBlock    -> 1,  
    Value         -> 127.9,  
    InteractionOrder -> {QED,-2},  
    Description    -> "Inverse of the EW coupling constant at the Z  
pole"  
},
```



Dependence in the expansion parameters

Parameters definition

```
M$Parameters = {  
    param1 == { options1 },  
    param2 == { options2 },  
    ...  
};
```

```
aEWM1 == {  
    ParameterType -> External,  
    BlockName     -> SMINPUTS,  
    OrderBlock    -> 1,  
    Value         -> 127.9,  
    InteractionOrder -> {QED,-2},  
    Description    -> "Inverse of the EW coupling constant at the Z  
pole"  
},
```

Parameters definition

Tensor parameters

```
CKM == {  
  ParameterType -> Internal,  
  Indices       -> {Index[Generation], Index[Generation]},  
  ComplexParameter -> True,  
  Unitary       -> True,  
  Value         -> {CKM[1,1] -> Cos[cabi], CKM[1,2] -> Sin[cabi],  
CKM[1,3] -> 0,  
                  CKM[2,1] -> -Sin[cabi], CKM[2,2] -> Cos[cabi],  
CKM[2,3] -> 0,  
                  CKM[3,1] -> 0, CKM[3,2] -> 0, CKM[3,3] -> 1 },  
  TeX          -> Superscript[V,CKM],  
  Description -> "CKM-Matrix" }
```

Parameters definition

Tensor parameters

```
CKM == {  
  ParameterType -> Internal,  
  Indices       -> {Index[Generation], Index[Generation]},  
  ComplexParameter -> True,  
  Unitary       -> True,  
  Value         -> {CKM[1,1] -> Cos[cabi], CKM[1,2] -> Sin[cabi],  
CKM[1,3] -> 0,  
                  CKM[2,1] -> -Sin[cabi], CKM[2,2] -> Cos[cabi],  
CKM[2,3] -> 0,  
                  CKM[3,1] -> 0, CKM[3,2] -> 0, CKM[3,3] -> 1},  
  TeX          -> Superscript[V,CKM],  
  Description   -> "CKM-Matrix"}
```

Parameters definition

Tensor parameters

```
CKM == {  
  ParameterType -> Internal,  
  Indices       -> {Index[Generation], Index[Generation]},  
  ComplexParameter -> True,           Default: Tensor is True, scalar is  
  Unitary       -> True,               False  
  Value         -> {CKM[1,1] -> Cos[cabi], CKM[1,2] -> Sin[cabi],  
CKM[1,3] -> 0,  
                CKM[2,1] -> -Sin[cabi], CKM[2,2] -> Cos[cabi],  
CKM[2,3] -> 0,  
                CKM[3,1] -> 0, CKM[3,2] -> 0, CKM[3,3] -> 1 },  
  TeX          -> Superscript[V,CKM],  
  Description -> "CKM-Matrix" }
```

Parameters definition

Tensor parameters

```
CKM == {  
  ParameterType -> Internal,  
  Indices       -> {Index[Generation], Index[Generation]},  
  ComplexParameter -> True,  
  Unitary       -> True,  
  Value         -> {CKM[1,1] -> Cos[cabi], CKM[1,2] -> Sin[cabi],  
CKM[1,3] -> 0,  
                  CKM[2,1] -> -Sin[cabi], CKM[2,2] -> Cos[cabi],  
CKM[2,3] -> 0,  
                  CKM[3,1] -> 0, CKM[3,2] -> 0, CKM[3,3] -> 1 },  
  TeX          -> Superscript[V,CKM],  
  Description  -> "CKM-Matrix"}
```


Interaction order

In the SM : QCD the power of g_s
QED the power of e

```
aEWM1 == { ...  
  InteractionOrder -> {QED,-2},  
  Description      -> "Inverse of the EW coupling constant at the Z pole"  
},
```

```
vev == {...  
  InteractionOrder -> {QED,-1},  
  Description      -> "Higgs vacuum expectation value"  
},
```

Interaction order

In the SM : QCD the power of g_s
 QED the power of e

aEWM1 == { ...

InteractionOrder -> {QED,-2},

Description -> "Inverse of the EW coupling constant at the Z pole"

},

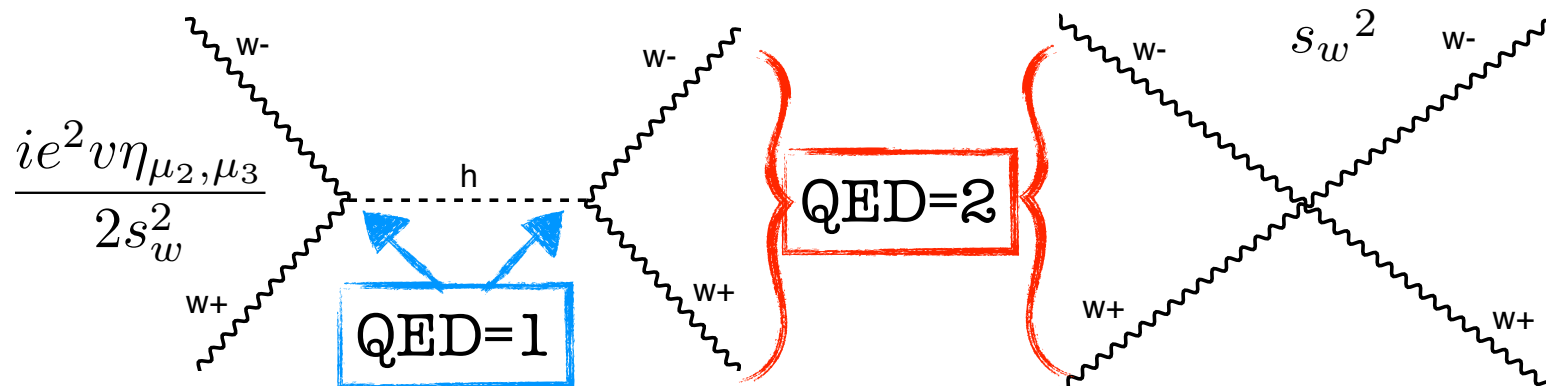
vev == {...

InteractionOrder -> {QED,-1},

Description -> "Higgs vacuum expectation value"

},

$$\frac{ie^2 (\eta_{\mu_1, \mu_4} \eta_{\mu_2, \mu_3} + \eta_{\mu_1, \mu_3} \eta_{\mu_2, \mu_4} - 2\eta_{\mu_1, \mu_2} \eta_{\mu_3, \mu_4})}{s_w^2}$$



Interaction order

In the SM : QCD the power of g_s
QED the power of e

```
vev == {...  
  InteractionOrder -> {QED,-1},  
  Description      -> "Higgs vacuum expectation value"  
},
```

```
yu == {...  
  InteractionOrder -> {QED, 1},  
  Description      -> "Up-type Yukawa couplings"  
},
```

Such that masses have QED=0

Interaction order

In the SM : QCD the power of g_s
QED the power of e

```
vev == { ...  
  InteractionOrder -> {QED, -1},  
  Description      -> "Higgs vacuum expectation value"  
},
```

```
yu == { ...  
  InteractionOrder -> {QED, 1},  
  Description      -> "Up-type Yukawa couplings"  
},
```

Such that masses have QED=0

However y_t is not a small parameter!

Interaction order

$\text{InteractionOrderHierarchy} = \{ \{ \text{QCD}, 1 \},$
 $\{ \text{QED}, 2 \} \};$

$$g_s \sim e^2$$

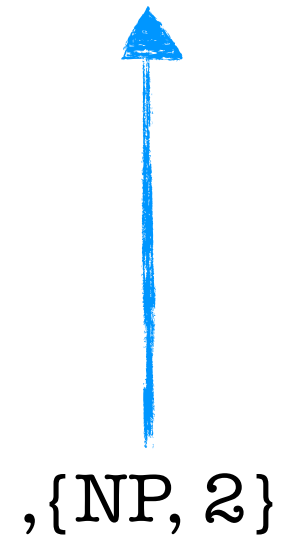
Interaction order

$\text{InteractionOrderHierarchy} = \{ \{ \text{QCD}, 1 \}, \{ \text{QED}, 2 \} \};$

$$g_s \sim e^2$$

$$\mathcal{L} = \mathcal{L} + \sum_i \frac{1}{\Lambda^2} \mathcal{O}_i + \mathcal{O}(\Lambda^{-4})$$

NP the power of Λ^{-2}



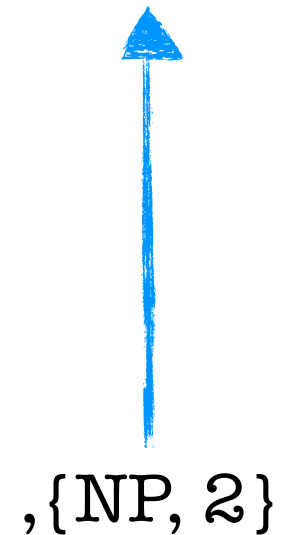
Interaction order

$\text{InteractionOrderHierarchy} = \{ \{ \text{QCD}, 1 \}, \{ \text{QED}, 2 \} \};$

$$g_s \sim e^2$$

$$\mathcal{L} = \mathcal{L} + \sum_i \frac{1}{\Lambda^2} \mathcal{O}_i + \mathcal{O}(\Lambda^{-4})$$

NP the power of Λ^{-2}



$\text{InteractionOrderLimit} = \{ \{ \text{NP}, 1 \} \};$

Max power per diagram of Λ^{-2} is 1

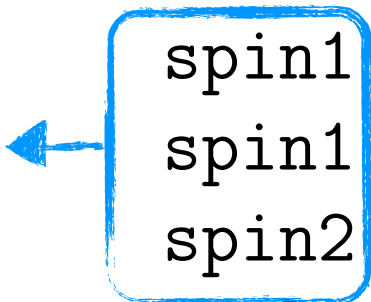
Fields definition I

```
M$ClassesDescription = {  
  spin1[1] == { options1 },  
  spin1[2] == { options2 },  
  spin2[1] == { options3 },  
  ...}
```


Fields definition I

S	0
W,F	1/2
V	1
RW,R	3/2
T	2
U	-1

```
M$ClassesDescription = {  
  spin1[1] == { options1 },  
  spin1[2] == { options2 },  
  spin2[1] == { options3 },  
  ... }
```



Fields definition I

S	0
W,F	1/2
V	1
RW,R	3/2
T	2
U	-1

```
M$ClassesDescription = {  
  spin1[1] == { options1 },  
  spin1[2] == { options2 },  
  spin2[1] == { options3 },  
  ... }
```

Unique Id

Fields definition I

ClassName->..., SelfConjugate->...,
 Indices->..., QuantumNumbers->...,
 FlavorIndex->..., ClassMembers,
 Mass->..., Width->..., PDG->...,
 Definitions->..., Unphysical->...,
 Chirality->..., MajoranaPhase->...,
 WeylComponents->...,
 Goldstone->..., Ghost->..., ...(Format)

S	0
W,F	1/2
V	1
RW,R	3/2
T	2
U	-1

```

M$ClassesDescription = {
  spin1[1] == { options1 },
  spin1[2] == { options2 },
  spin2[1] == { options3 },
  ... }
  
```

Unique Id

Fields definition II

```
F[3] == {  ClassName      -> uq,
          ClassMembers   -> {u, c, t},
          Indices        -> {Index[Generation], Index[Colour]},
          FlavorIndex    -> Generation,
          SelfConjugate  -> False,
          Mass           -> {Mu, {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},
          Width          -> {0, 0, {WT,1.50833649}},
          QuantumNumbers -> {Q -> 2/3},
          PDG            -> {2, 4, 6},
          ...
}
```

Fields definition II

Spin index



```
F[3] == {  ClassName      -> uq,  
          ClassMembers  -> {u, c, t},  
          Indices        -> {Index[Generation], Index[Colour]},  
          FlavorIndex    -> Generation,  
          SelfConjugate  -> False,  
          Mass           -> {Mu, {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},  
          Width          -> {0, 0, {WT,1.50833649}},  
          QuantumNumbers -> {Q -> 2/3},  
          PDG            -> {2, 4, 6},  
          ...  
}
```

Fields definition II

Generation index distinguishes
the class members

```
F[3] == {  ClassName    -> uq,  
         ClassMembers -> {u, c, t},  
         Indices      -> {Index[Generation] Index[Colour]},  
         FlavorIndex  -> Generation,  
         SelfConjugate -> False,  
         Mass         -> {Mu, {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},  
         Width        -> {0, 0, {WT,1.50833649}},  
         QuantumNumbers -> {Q -> 2/3},  
         PDG          -> {2, 4, 6},  
         ...  
}
```

Fields definition II

```
F[3] == {  ClassName      -> uq,  
          ClassMembers  -> {u, c, t},  
          Indices       -> {Index[Generation], Index[Colour]},  
          FlavorIndex   -> Generation,  
          SelfConjugate -> False,  
          Mass          -> {Mu, {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},  
          Width        -> {0, 0, {WT,1.50833649}},  
          QuantumNumbers -> {Q -> 2/3},  
          PDG          -> {2, 4, 6},  
          ...  
}
```

Same representation

Fields definition II

```
F[3] == {  ClassName      -> uq,  
          ClassMembers  -> {u, c, t},  
          Indices       -> {Index[Generation], Index[Colour]},  
          FlavorIndex   -> Generation,  
          SelfConjugate  -> False,  
          Mass          -> {Mu, {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},  
          Width         -> {0, 0, {WT,1.50833649}},  
          QuantumNumbers -> {Q -> 2/3},  
          PDG           -> {2, 4, 6},  
          ...  
}
```

External parameters

Fields definition II

```
F[3] == {  ClassName      -> uq,
          ClassMembers  -> {u, c, t},
          Indices        -> {Index[Generation], Index[Colour]},
          FlavorIndex    -> Generation,
          SelfConjugate  -> False,
          Mass           -> {Mu {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},
          Width          -> {0, 0, {WT,1.50833649}},
          QuantumNumbers -> {Q -> 2/3},
          PDG            -> {2, 4, 6},
          ...
}
```

External parameters

Generic label

```
Mass -> {MW, Internal}
Mass -> {MZ, 91.188}
Mass -> {{MU,0}, {MC,0}, {MT, 174.3}}
Mass -> {Mu, {MU, 0}, {MC, 0}, {MT, 174.3}}
```

Fields definition II

```
F[3] == {  ClassName      -> uq,  
          ClassMembers  -> {u, c, t},  
          Indices        -> {Index[Generation], Index[Colour]},  
          FlavorIndex    -> Generation,  
          SelfConjugate  -> False,  
          Mass           -> {Mu, {MU, 2.55*^-3}, {MC,1.27}, {MT,172}},  
          Width          -> {0, 0, {WT,1.50833649}},  
          QuantumNumbers -> {Q -> 2/3},  
          PDG            -> {2, 4, 6},  
          ...  
}
```

Not used in FR but by
following codes

Fields definition III

Interaction eigenstates

```
V[12] == {  
  ClassName    -> Wi,  
  Unphysical   -> True,  
  SelfConjugate -> True,  
  Indices      -> {Index[SU2W]},  
  FlavorIndex  -> SU2W,  
  Definitions  -> { Wi[mu_,1] -> (Wbar[mu]+W[mu])/Sqrt[2],  
    Wi[mu_,2] -> (Wbar[mu]-W[mu])/(I*Sqrt[2]), Wi[mu_,3] -> cw  
    Z[mu] + sw A[mu]}  
}
```

FR does not export
them to matrix
element code

Physical fields

Fields definition IV

```
U[11] == {  
  ClassName    -> ghB,  
  Unphysical   -> True,  
  SelfConjugate -> False,  
  Ghost        -> B,  
  Definitions  -> { ghB -> -sw ghZ + cw ghA }  
},
```

ClassName of the
boson

```
S[2] == {  
  ClassName    -> GO,  
  SelfConjugate -> True,  
  Goldstone    -> Z,  
  ...  
},
```

Gauge Groups

```
M$GaugeGroups = {  
  U1Y == {  
    Abelian      -> True,  
    CouplingConstant -> g1,  
    GaugeBoson   -> B,  
    Charge       -> Y  
  },...  
  SU3C == {  
    Abelian      -> False,  
    CouplingConstant -> gs,  
    GaugeBoson   -> G,  
    StructureConstant -> f,  
    Representations -> {T,Colour},  
    SymmetricTensor -> dSUN  
  }  
};
```

Gauge Groups

```
M$GaugeGroups = {  
  U1Y == {  
    Abelian      -> True  
    CouplingConstant -> g1,  
    GaugeBoson   -> B,  
    Charge       -> Y  
  },...  
  SU3C == {  
    Abelian      -> False,  
    CouplingConstant -> gs,  
    GaugeBoson   -> G,  
    StructureConstant -> f,  
    Representations -> {T,Colour},  
    SymmetricTensor -> dSUN  
  }  
};
```

Gauge Groups

```
M$GaugeGroups = {  
  U1Y == {  
    Abelian      -> True,  
    CouplingConstant -> g1,  
    GaugeBoson   -> B,  
    Charge       -> Y  
  },...  
  SU3C == {  
    Abelian      -> False,  
    CouplingConstant -> gs,  
    GaugeBoson   -> G,  
    StructureConstant -> f,  
    Representations -> {T,Colour},  
    SymmetricTensor -> dSUN  
  }  
};
```

Gauge Groups

```
M$GaugeGroups = {  
  U1Y == {  
    Abelian      -> True,  
    CouplingConstant -> g1,  
    GaugeBoson   -> B,  
    Charge       -> Y  
  },...  
  SU3C == {  
    Abelian      -> False,  
    CouplingConstant -> gs,  
    GaugeBoson   -> G,  
    StructureConstant -> f,  
    Representations -> {T,Colour}  
    SymmetricTensor -> dSUN  
  }  
};
```

Generator label

Associated index

Gauge groups

- Predefined strength tensor

$$\text{FS}[A, \mu, \nu, a] \begin{array}{l} \text{abelian} \end{array} \Rightarrow F_{\mu\nu}^a = \partial_\mu A_\nu^a - \partial_\nu A_\mu^a + gf^a_{bc} A_\mu^b A_\nu^c$$

- Predefined covariant derivative (from the **Indices** and **QuantumNumbers** of the fields)

$$\text{DC}[\phi, \mu] \Rightarrow D_\mu \phi = \partial_\mu \phi - ig A_\mu^a T_a \phi$$

Lagrangian

$$\mathcal{L}^{QCD} \equiv -\frac{1}{4} G_a^{\mu\nu} G_{\mu\nu}^a + i\bar{d}\not{D}d$$

$$L = -1/4 \text{FS}[G, \text{mu}, \text{nu}, \text{a}] \text{FS}[G, \text{mu}, \text{nu}, \text{a}] \\ + \text{I} \text{dqbar}.\text{Ga}[\text{mu}].\text{DC}[\text{dq}, \text{mu}]$$

FeynRules creates the “anti”-particle name

Dot to avoid commuting the fermions

$\text{dqbar}.\text{Ga}[\text{mu}].\text{T}[\text{a}].\text{dq}$

$\rightarrow \text{Ga}[\text{mu}, \text{s}, \text{r}] \text{T}[\text{a}, \text{i}, \text{j}] \text{dqbar}[\text{s}, \text{f}, \text{i}].\text{dq}[\text{r}, \text{f}, \text{j}]$

FeynRules restores the indices internally

Running FeynRules

In Mathematica :

Loading Feynrules

```
$FeynRulesPath = SetDirectory[ <the address of the package> ];  
<< FeynRules`
```

Loading the model

```
LoadModel[ < file.fr >, < file2.fr >, ... ]
```

Extracting the Feynman rules

```
vertsQCD = FeynmanRules[ LQCD ];
```

Checking the Lagrangian

```
CheckKineticTermNormalisation[ L ]  
CheckMassSpectrum[ L ]
```

Running FeynRules

In Mathematica :

Loading Feynrules

```
$FeynRulesPath = SetDirectory[ <the address of the package> ];  
<< FeynRules`
```

Loading the model

```
LoadModel[ < file.fr >, < file2.fr >, ... ]
```

All the model files should
be loaded at once

Extracting the Feynman rules

```
vertsQCD = FeynmanRules[ LQCD ];
```

Checking the Lagrangian

```
CheckKineticTermNormalisation[ L ]  
CheckMassSpectrum[ L ]
```

Running FeynRules

In Mathematica :

Loading Feynrules

```
$FeynRulesPath = SetDirectory[ <the address of the package> ];  
<< FeynRules`
```

Loading the model

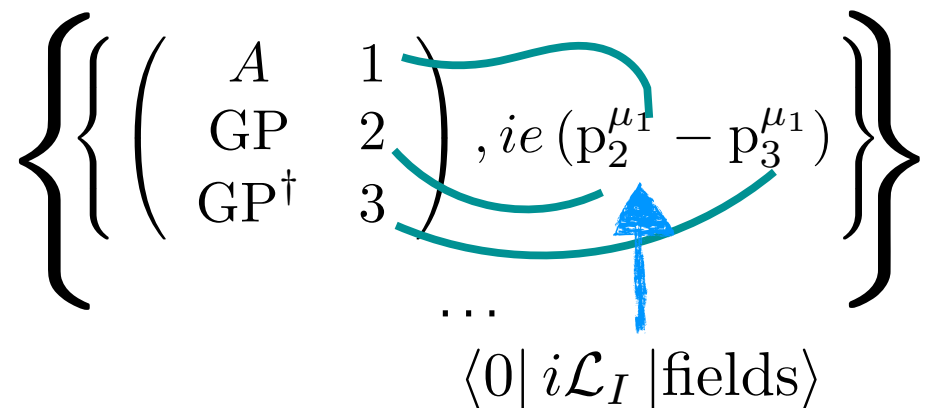
```
LoadModel[ < file.fr >, < file2.fr >, ... ]
```

Extracting the Feynman rules

```
vertsQCD = FeynmanRules[ LQCD ];
```

Checking the Lagrangian

```
CheckKineticTermNormalisation[ L ]  
CheckMassSpectrum[ L ]
```



All momenta are incoming

Running FeynRules

In Mathematica :

Loading Feynrules

```
$FeynRulesPath = SetDirectory[ <the address of the package> ];  
<< FeynRules`
```

Loading the model

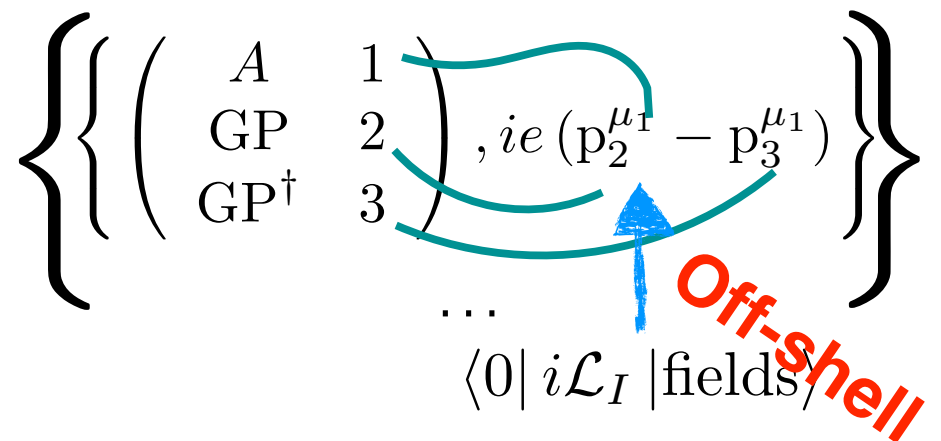
```
LoadModel[ < file.fr >, < file2.fr >, ... ]
```

Extracting the Feynman rules

```
vertsQCD = FeynmanRules[ LQCD ];
```

Checking the Lagrangian

```
CheckKineticTermNormalisation[ L ]  
CheckMassSpectrum[ L ]
```



All momenta are incoming

Checks

CheckHermiticity[L, options]

CheckDiagonalKineticTerms[L, options]

CheckDiagonalMassTerms[L, options]

CheckDiagonalQuadraticTerms[L, options]

CheckKineticTermNormalisation[L, options]

$$\begin{array}{lll} \frac{1}{2}\partial_\mu\phi\partial^\mu\phi - \frac{1}{2}m^2\phi^2 & \frac{1}{2}\bar{\lambda}i\not{\partial}\lambda - \frac{1}{2}m\bar{\lambda}\lambda & -\frac{1}{4}F_{\mu\nu}F^{\mu\nu} - \frac{1}{2}m^2A_\mu A^\mu \\ \partial_\mu\phi^\dagger\partial^\mu\phi - m^2\phi^\dagger\phi & \bar{\psi}i\not{\partial}\psi - m\bar{\psi}\psi & -\frac{1}{2}F_{\mu\nu}^\dagger F^{\mu\nu} - m^2A_\mu^\dagger A^\mu \end{array}$$

CheckMassSpectrum[L, options]

Toolbox

`ExpandIndices[L, options]`

`GetKineticTerms[L, options]`

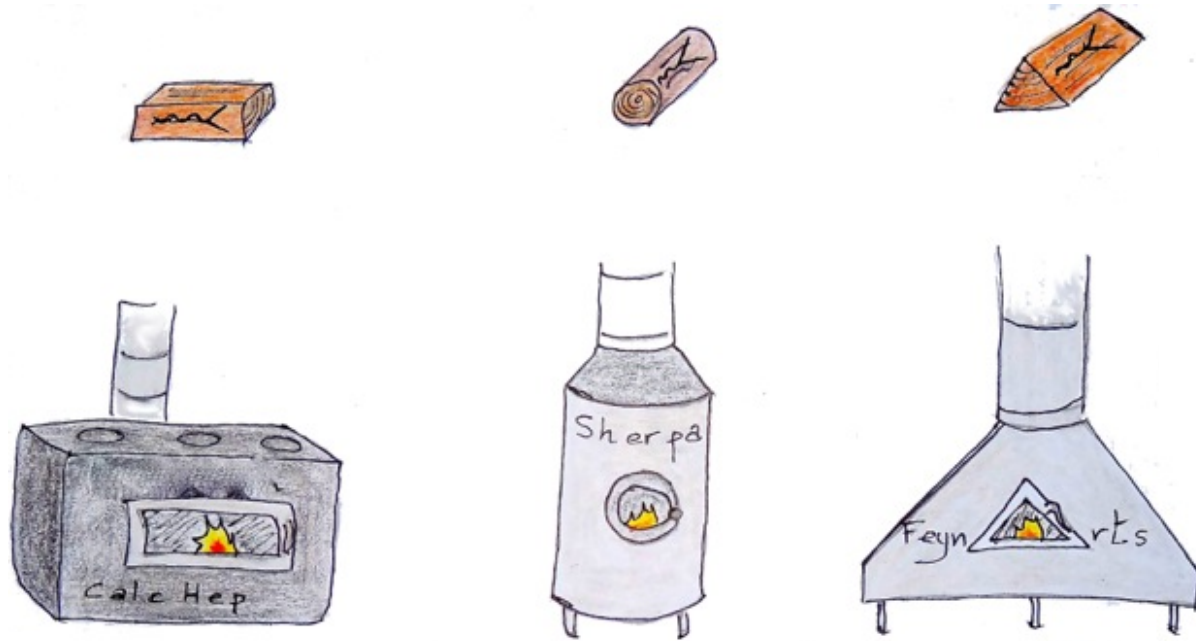
`GetMassTerms[L, options]`

`GetQuadraticTerms[L, options]`

`GetInteractionTerms[L, options]`

`SelectFieldContent[L, list]`

FeynRules outputs



FeynRules
outputs can be
used directly by
event generators

UFO : output with
the full information
used by several
generators



Conventions

- FeynRules does not care which symbol is used
- However there are conventions for the translation to matrix element computation tools
 - α_s is hardcoded in most code (running) as well as other SM parameters (α_{EW}^{-1}, G_F)
 - SU(3) representations
 - PDG numbering scheme, LHA block, ...

FeynRules outputs

Generating the output

WriteCHOutput[L]

WriteFeynArtsOutput[L]

WriteSHOutput[L]

WriteWOutput[L]

WriteUFO[L]

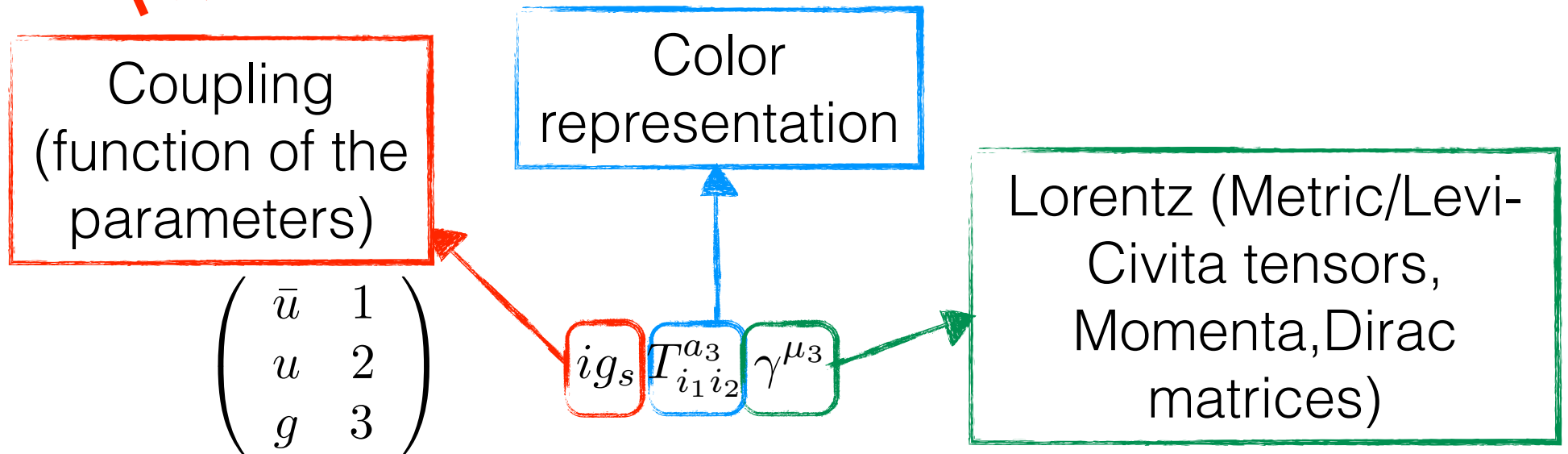
FeynRules takes care of all the conventions

FeynRules outputs

Generating the output

```
WriteCHOutput[ L ]  
WriteFeynArtsOutput[ L ]  
WriteSHOutput[ L ]  
WriteWOutput[ L ]  
WriteUFO[ L ]
```

FeynRules takes care of all the conventions

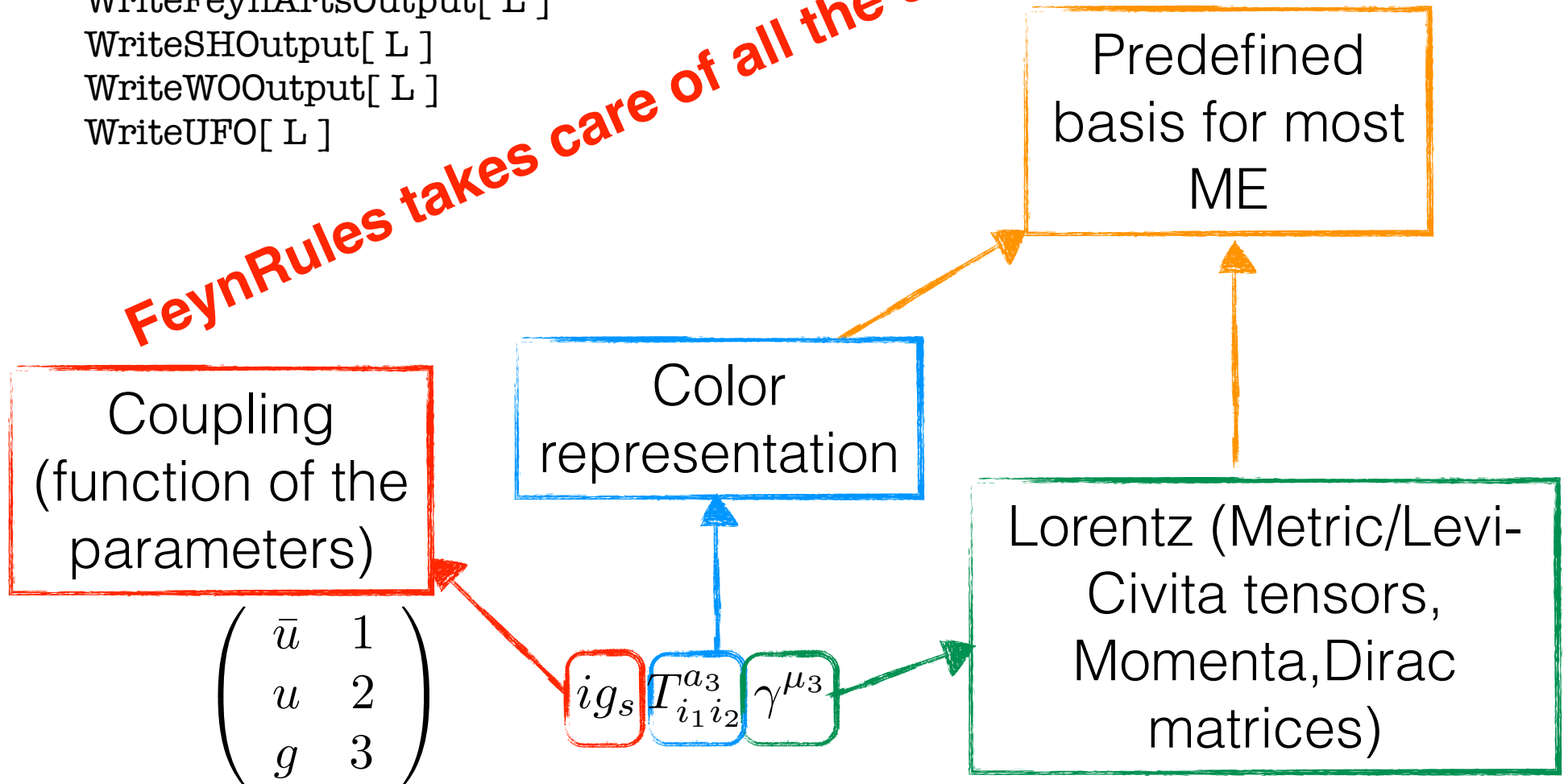


FeynRules outputs

Generating the output

```
WriteCHOutput[ L ]  
WriteFeynArtsOutput[ L ]  
WriteSHOutput[ L ]  
WriteWOutput[ L ]  
WriteUFO[ L ]
```

FeynRules takes care of all the conventions



Sherpa output

[feynrules.dat](#): A static file, setting up the model in Sherpa.

[Particle.dat](#): The list of all particles together with their properties.

[param_card.dat](#): LH-like file defining the numerical values of the external parameters.

[ident_card.dat](#): File linking the entries in [param_card.dat](#) to the variables used in the Sherpa code.

[param_definition.dat](#): File containing analytical expressions for all the internal parameters.

[Interactions.dat](#): File defining all the interaction vertices with their couplings.

Sherpa output

feynrules.dat: A static file, setting up the model in Sherpa.

Particle.dat: The list of all particles together with their properties.

param_card.dat: LH-like file defining the numerical values of the external parameters.

ident_card.dat: File linking the entries in param_card.dat to the variables used in the Sherpa code.

param_definition.d

the internal para

Interactions.dat:

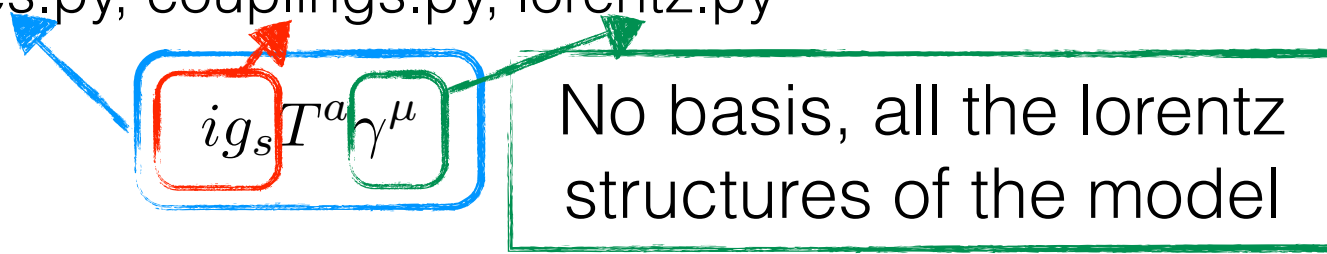
couplings.

```
VERTEX 5 22 5 # b a b
1 -0.3333333333333333*ee*(0,1) # right-handed coupling
2 -0.3333333333333333*ee*(0,1) # left-handed coupling
3 D[1,3] # colour structure
4 FFV # Lorentz structure
```

$$-\frac{i}{3}e\delta_{ij}\gamma^\mu\gamma_+ - \frac{i}{3}e\delta_{ij}\gamma^\mu\gamma_-$$

Color or Lorentz structures of the SM and MSSM only

UFO

- Generic output with the **full** model information
 - coupling_orders.py, parameters.py, particles.py, write_param_card.py, __init__.py,
 - vertices.py, couplings.py, lorentz.py
- 

No basis, all the lorentz structures of the model
- decays.py
- CT_vertices.py, CT_couplings.py (For NLO)
- Python module used in MadGraph, Herwig, Gosam(, Sherpa)

UFO

vertices.py

```
V_135 = Vertex(name = 'V_135',  
               particles = [ P.u__tilde__, P.u, P.g ],  
               color = [ 'T(3,2,1)' ],  
               lorentz = [ L.FFV1 ],  
               couplings = {(0,0):C.GC_11})
```

Lorentz.py

```
FFV1 = Lorentz(name = 'FFV1',  
               spins = [ 2, 2, 3 ],  
               structure = 'Gamma(3,2,1)')
```

couplings.py

```
GC_11 = Coupling(name = 'GC_11',  
                 value = 'complex(0,1)*G',  
                 order = {'QCD':1})
```

Plan

- Field theory : a short reminder
 - free fields (KG details, Fermion)
 - Scattering matrix in perturbation
 - Wick theorem to Feynman rules
- Why Monte-Carlo/automated tools?
- Lagrangian to the Feynman rules
 - Model file : Parameters, fields, gauge group and Lagrangian
 - Running FeynRules
- Demo