

The MGv5 model structure

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1 particles.py

Each particle is defined as a dictionary, *e.g.* for the electron,

```
particles[0] = {
    'name':'e-',
    'antiname':'e+',
    'spin':2,
    'color':1,
    'mass':'ME',
    'width':'WE',
    'texname':'e^-',
    'antitexname':'e^+',
    'line':'S',
    'charge':-1
}
```

All attributes are strings, except for **spin**, **color**, and **charge**. The number for the spin corresponds to $(2s + 1)$, for color it is the representation. All particles can have a additional optional attribute that tags the particle as non propagating,

```
'propagating':False
```

The default value is **True**, in which case the attribute can be omitted.

2 parameters.py

Each parameter is defined as a dictionary, *e.g.* for the strong coupling,

```
parameters[0] = {
    'name':'gs',
    'nature':'external' / 'internal',
    'type':'real' / 'complex',
    'value': ...
    'texname':'g_s'
```

```
}
```

For external parameters, `value` is a number, whereas for internal parameters it is an algebraic expression. For external parameters there are two more attributes linking the parameter to the corresponding entry in the `param_card.dat`,

```
'lhablock': 'SMINPUTS',  
'lhacode': [1,2,3,...],
```

3 vertices.py

Each vertex is represented as a dictionary, *e.g.* for the u, \bar{u}, g vertex,

```
vertices[0] = {  
    'particles': [u, u, g],  
    'color': [ 'T(a3,i2,i1)', ...],  
    'lorentz': [ L1, L2, ...],  
    'couplings': [ (0,0):'g1', (0,1):'g2', ...],  
    'orders': ['QCD', ...]  
}
```

`color` is a list of all color structures appearing in the vertex. Similarly, `lorentz` is a list containing all Lorentz structures appearing in the vertex. Note that '`L1`', '`L2`', ... denote the classes of HELAS routines (See below). Finally, `couplings` is a list linking the couplings to the color and Lorentz structures, *i.e.*, `(0,0):'g1'` represents $g_1 * T(a_3, i_2, i_1) * L_1$, *etc.*

4 couplings.py

```
couplings[0] = {  
    'name': 'g1',  
    'expression': ...  
}
```

5 lorentz.py

```
buildingblocks[0] = {  
    'name': 'B1',  
    'expression': ...  
}
```

```
lorentz[0] = {  
    'name': 'L1',
```

```

'particles':[2, 2, 1],
    # spin structure
'onshell': [{'L10': ...}]
    # there are two entries for fermions
'offshell': [ [ {'L11': ... }], [ {'L12': ... }], ....]
    # L11 is the Lorentz structure L1 with particle 1 off shell, etc.
}

```