#### **Lattice Grammar**

The grammar is context free<sup>1</sup> and the text is in free-format<sup>2</sup> style. However, the length of a statement is limited to 132 characters.

# {<text>} or /\*<text>\*/ **Statements** A statement may run over several lines and is terminated by ";": <statement>; **Constants** Constants are defined by: <constant name> = <arithmetic expression>; Lattice File The lattice file contains: define lattice; ringtype = $<1 \mid 0>$ ; Energy=<beam energy [GeV/c²] (for classical radiation, and quantum fluctuations)>; codeps= $<\epsilon$ for the closed orbit finder (Newton search)>; $dp = <\Delta p/p>$ ; <element definition> <element definition>

anywhere between tokens.

**Comments** 

<sup>&</sup>lt;sup>1</sup> Chomsky type 2: grammars that only allow one symbol on the left hand side of a production, e.g. Algol-60 (specified by the Bachus-Nauer Form), Pascal, C or C++ (but not Fortran, i.e. DATA statements).

<sup>2</sup> White space (space, tab, and newline) has no significance (except in strings), i.e., it can be inserted

```
<section definition>
<section definition>
cell: <element | section> {,<element | section>}, symmetry=1;
end;
                                     Elements
Marker
<element name>: Marker;
Example:
                 M1: Marker;
Beam Position Monitor
<element name>: Beam Position Monitor;
Example:
         BPM1: Beam Position Monitor;
Drift
<element name>: drift,
         L=<length [m]>;
Example:
         L1: Drift, L=0.30;
Corrector
<element name>: Corrector,
         [L=<length [m], default: 0 (thin kick)>,]
         <direction: 'horizontal' | 'vertical'>;
Example:
```

COH: Corrector, horizontal;

## **Dipole**

<element name>: Sextupole,

```
<element name>: Bending,
          L=<length (\rho [m] · \phi [rad])>,
          [Roll=<Design roll angle [°]>,]
           T= dend angle [°]>,
          [T1=<entrance angle [°]>,]
           [T2 = \langle exit angle [°] \rangle,]
          [Gap = \langle gap [m] \rangle,
          [K=< b<sub>2</sub> [m<sup>-2</sup>] (gradient)>,
           N=<no of integration steps>,
           Method=<integration method: '0' | '2' | '4'), default: '2' (matrix style, 2<sup>nd</sup> and
                    4<sup>th</sup> order symplectic integrator>;
           [HOM=<list of systematic higher order multipole errors (random errors
                    are assigned in the input file): n, bn, an, ...(order, integrated skew- and
                    normal multipole strengths)>;]
Example
          B: Bending, L=0.70, T=10.0, T1:=5.0, T2:=5.0, K=-1.0, N=8, Method=2;
Quadrupole
<element name>: Quadrupole,
          L=<length [m]>,
           [Roll<sup>3</sup>=<Design roll angle [°]>,]
           K=\langle b_2 [m^{-2}] (gradient) \rangle,
           N=<no of integration steps>,
          Method=<integration method: '0' | '2' | '4'), default: '2' (matrix style, 2<sup>nd</sup> and
                    4<sup>th</sup> order symplectic integrator>;
           [HOM=<list of systematic higher order multipole errors (random errors
                    are assigned in the input file): n, bn, an, ...(order, integrated skew- and
                      normal multipole strengths)>;]
Example
QF: Quadrupole, L=0.5, K=2.2134, N=4, Method=4;
Sextupole
```

```
[L=<length [m], default 0 (thin kick)>,]
[Roll³=<Design roll angle [°]>,]
K=<b₃ [m⁻³] (sextupole strength)>;
N=<no of integration steps>,
Method=<integration method: '0' | '2' | '4'), default: '2' (matrix style, 2<sup>nd</sup> and 4<sup>th</sup> order symplectic integrator>;
[HOM=<list of systematic higher order multipole errors (random errors are assigned in the input file): n, bn, an, ...(order, integrated skew- and normal multipole strengths)>;]
```

#### Example

```
SF : Sextupole, K=-10.2363;
```

## Multipole

```
<element name>: Multipole,
    L=<length [m]>,
    [Roll³=<Design roll angle [°]>,]
    [T=<bend angle [°]>,]
    [T1=<entrance angle [°]>,]
    [T2=<exit angle [°]>,]
    N=<no of integration steps>,
    Method=<integration method: '0' | '2' | '4'), default: '2' (matrix style, 2<sup>nd</sup> and 4<sup>th</sup> order symplectic integrator>;
    [HOM=<list of systematic higher order multipole errors (random errors are assigned in the input file): n, bn, an, ...(order, integrated skew- and normal multipole strengths)>;]
```

## Examples

```
B: multipole, L=0.70, T=10.0, T1=5.0, T2=5.0, HOM=(2, -1.0, 0), N=8, Method=2;
QF: multipole, L=0.70, HOM=(2, 2.50, 0.0, 4, 1.01e7, 0.0), N=8, Method=2;
```

## Wiggler

```
<name>: Wiggler,
L=<length [m]>,
BoBrho=<B/Brho [m<sup>-1</sup>]>,
Lambda=<period [m]>,
kx=<[m]>,
```

N=<no of integration steps>, Method=<integration method: '0' | '2' | '4'), default: '2' (matrix style, 2<sup>nd</sup> and 4<sup>th</sup> order symplectic integrator>;

## Example

U143: wiggler, L=4.80, BoBrho=0.5, Lambda=0.15, N=20, Method=0;

## **Cavity**

## Example

CAV : Cavity, Frequency = 499.95e6, Voltage=1.22e6, HarNum=328;

#### **Sections**

<section name>: <element name> {, <element name>};

## **Operations**

#### Reversal

inv(<section name>)

Note, this command reverses the list of elements, i.e., it does not change the parameters of the elements. In particular, the resulting list is only mirror symmetric if the individual elements are mirror symmetric.