## **Exercises 12**

1. Write a turtle command sequence which generates a balls-and-sticks molecule model of

formic acid (structural formula: 
$$\begin{bmatrix} 0 \\ H \end{bmatrix}$$
 ).

The atoms shall be represented by spheres with different sizes and colours (depending on the chemical element), and the bonds by cylinders (**F(...)** command of the turtle). The double bond shall be represented by a thicker cylinder.

Hint: The colour of an atom  $\mathbf{x}$  can be specified as in the following example:

Test your solution with GroIMP.

- 2. (a) Write an L-system which simulates the primary growth of a plant in annual steps. The *annual shoots* of the vertical main axis (stem) shall all have the same length. The uppermost annual shoot shall bear an *apical bud* (= a red sphere) and a *lateral bud* (= a green sphere). The apical bud is supposed to produce a new annual shoot of the main axis next year, and from the lateral bud shall grow a shorter *lateral shoot* with a branching angle of 45°, which will terminate its growth next year (i.e., there are no buds at the lateral shoots). The positions of the lateral branches are alternating (left-right-left-right-...) along the stem. The simulation shall start with an apical bud.
- (b) Modify the model by introducing a trend: Assume that the annual shoots get 10 per cent shorter each year.
- (c) Assume additionally that the apical bud produces a flower ( = a large blue cone) after 7 years, and that the plant then stops to grow.

Test your solutions with GroIMP.

Remarks: By M(-s) you can cause the turtle to move back along the main axis by stepsize s. Cone (h, r) stands for a cone with height h and radius r.

3. Open the example "Molecules" in GroIMP's built-in example portfolio ("File" / "Show Examples").

Make several model runs by clicking on the buttons "Run run" and "Reset", and observe what happens.

Now modify the model in the following ways:

- (a) Increase the number of atoms from 10 to 20.
- (b) Switch off the output of text to the console.
- (c) Double the distance threshold for formation of a bond between two atoms.