

```
In[1]:= dir = NotebookDirectory[];
```

```
In[2]:= SetDirectory[dir];
```

```
In[3]:= << DynamicSystems.m
```

```
In[4]:= Chemostat
```

Chemostat

$$N'(t) = \alpha_1 \frac{C(t)}{1+C(t)} N(t) - N(t)$$

$$C'(t) = \alpha_2 - C(t) - \frac{C(t)}{1+C(t)} N(t)$$

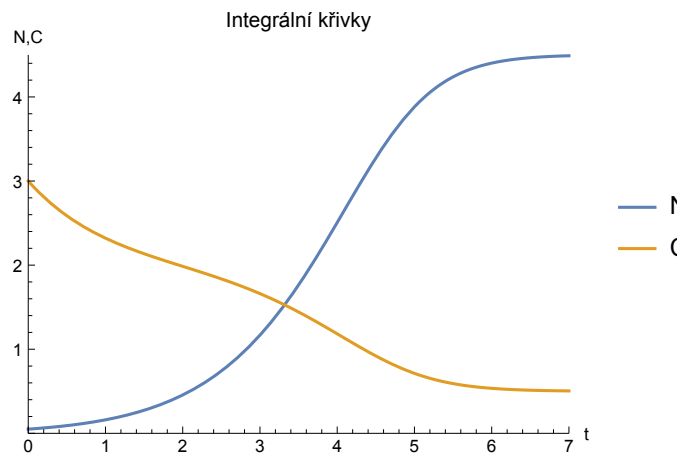
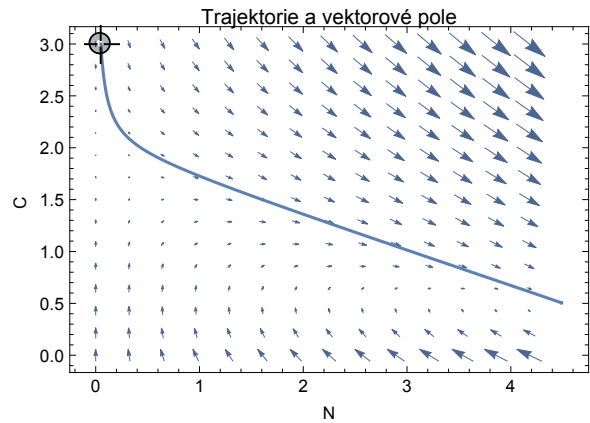
α_1

α_2

t_1

$N_{\max}$

Out[4]= $C_{\max}$

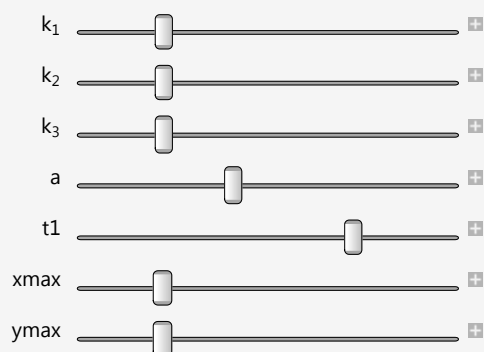


In[5]:= BZ

Bělousov–Žabotínského reakce

$$x'(t) = a k_1 x(t) - k_2 x(t)y(t)$$

$$y'(t) = k_2 x(t)y(t) - k_3 y(t)$$

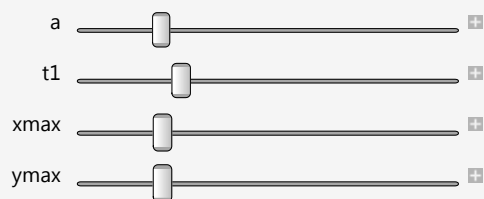


In[6]:= Sch

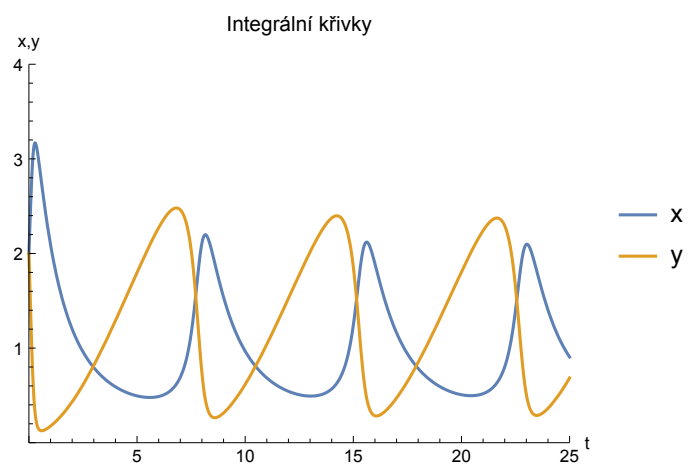
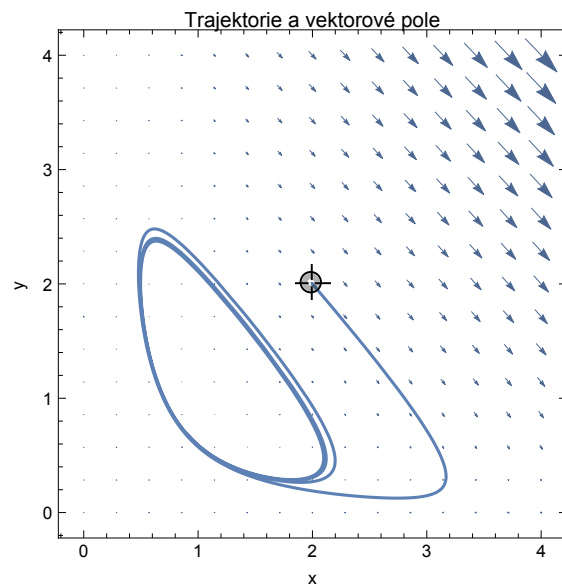
Schnakenbergův systém

$$x'(t) = x^2(t)y(t) - x(t)$$

$$y'(t) = a - x^2(t)y(t)$$



Out[6]=

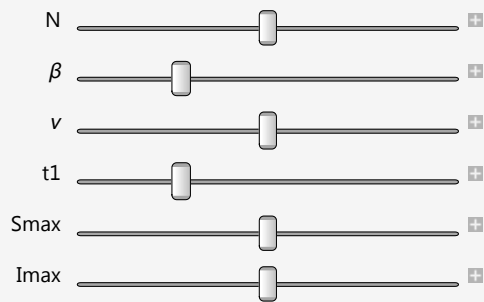


In[7]:= **SIR****SIR model**

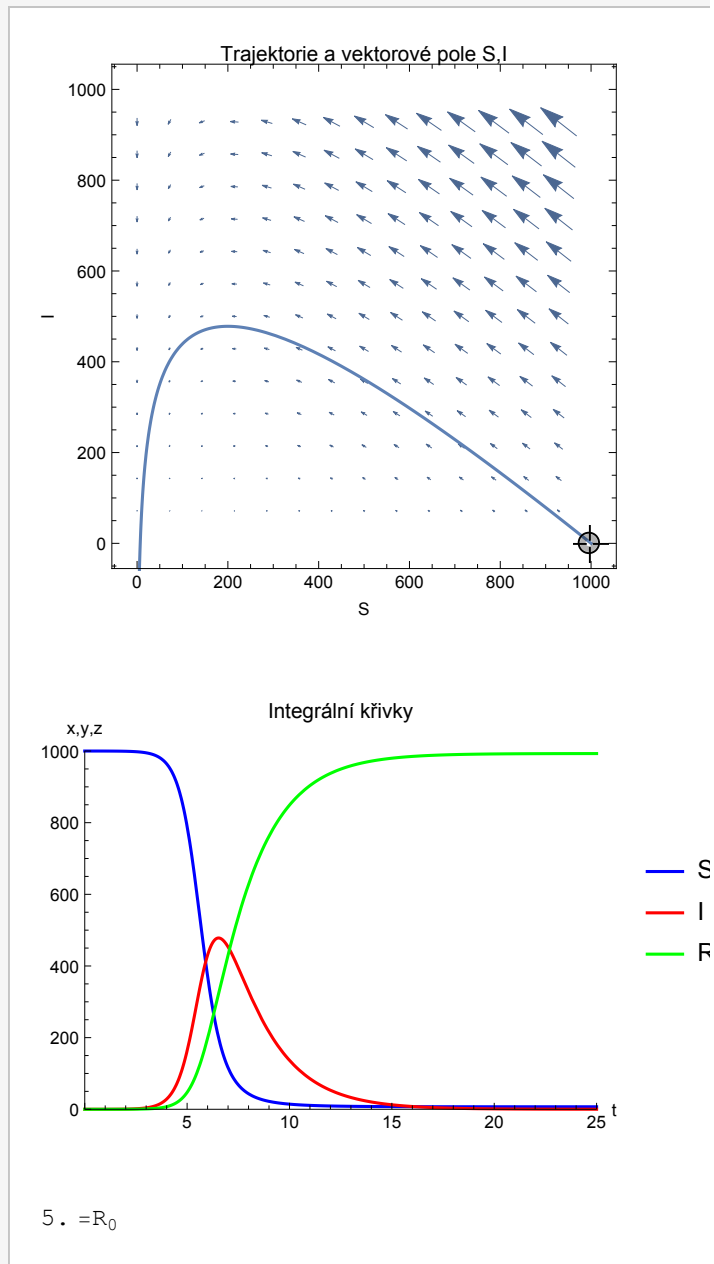
$$S'(t) = -\beta S(t) I(t)$$

$$I'(t) = \beta S(t) I(t) - \nu I(t)$$

$$R'(t) = \nu I(t)$$



Out[7]=



In[8]:= SIRS

SIR model

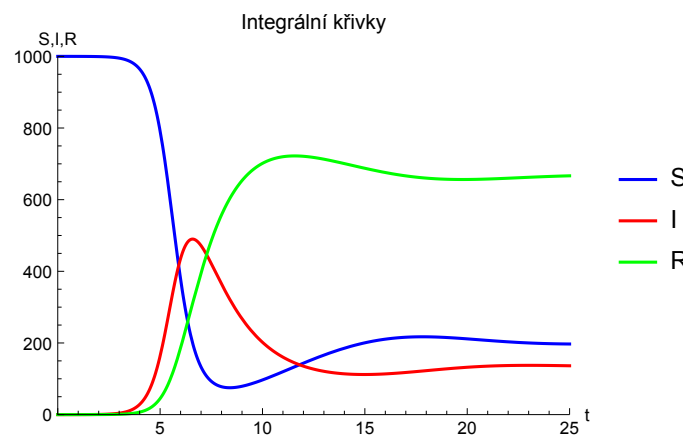
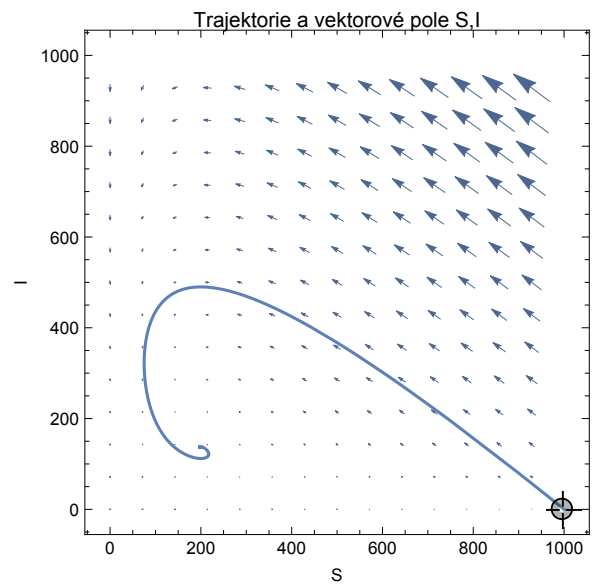
$$S'(t) = -\beta S(t) I(t)$$

$$I'(t) = \beta S(t) I(t) - \nu I(t)$$

$$R'(t) = \nu I(t)$$

N β ν γ t_1 Smax Imax

Out[8]=

5. $=R_0$