

## Lab 5 – Sage Tutorial

### 1. Linearization

**What?:** Linearization refers to the local approximation of a nonlinear function by a linear one. Here, 'local' refers to the fact that the approximation is 'good' only in a (small) neighborhood of a reference point.

**Why?:** Linear problems are always explicitly solvable (think of algebraic or differential equation) while nonlinear problems aren't. A good linear approximation will thus provide valuable insights into the more complicated nonlinear 'geometry'.

**How?:** The linear approximation is constructed by using the first terms of a Taylor expansion  $f(x) = f(x^*) + f'(x^*)(x - x^*) + HOT$  (i.e., by discarding Higher Order Terms)). For vector-valued functions of several variables,  $f'(x^*)$  is replaced by its high-dimensional counterpart, the Jacobi matrix (or Jacobian).

**Sage** does provide a specialized command for computing the linearization/Jacobian

```
x, y = var('x y')
f = (2*x-x^2-x*y, -y+x*y)
jacobian(f, (x,y))
```

```
Out[70]: [-2*x - y + 2      -x]
          [                y      x - 1]
```

Notice the fact that this is a (matrix-valued) function (of  $x$  and  $y$ ).

After inserting (different) numerical values for  $x$  and  $y$  one obtains (different) matrices

```
In [7]: jac = jacobian(f, (x,y))
A1 = jac.subs(x=0,y=0)
A2 = jac.subs(x=2,y=0)
A3 = jac.subs(x=1,y=1)
A1, A2, A3
```

```
Out[7]: (
  [ 2  0]  [-2 -2]  [-1 -1]
  [ 0 -1], [ 0  1], [ 1  0]
)
```

These can be analyzed with standard Linear Algebra methods (eigenvalues, eigenvectors, etc.).

## 2. Plotting direction fields and orbits

One of the most powerful methods for understanding nonlinear (planar dynamics) is based on representing the direction(vector) field which generates the dynamics.

**Geometrically:** the systems orbits are tangent to the generating vector field.

**Sage** has inbuilt commands for computing and plotting both direction fields and orbits.

For **plotting the direction field** we first define the field (RHS of the system) to be plotted ( $f$  has two components,  $f[0]$  and  $f[1]$  in **Sage**)

```
In [71]: x, y = var('x y')
         f = (2*x-x^2-x*y, -y+x*y)
```

then use the appropriate plot command (with good limits for the  $x$  and  $y$  ranges)

```
In [57]: # the direction field
         plot_vector_field((f[0],f[1]), (x,-1,2.2),(y,-1,1.2))
```

In order to **compute orbits** (as most systems are **not** explicitly solvable in terms of simple functions) one resorts to a numerical ODE solver

```
sol = desolve_odeint(f, ics, times, [x, y])
line(zip(sol[:,0],sol[:,1]))
```

Here, the arguments of the solve command are: ' $f$ ' the nonlinear RHS of the equation, the initial conditions ' $ics$ ', ' $times$ ' a 3-tuple containing the initial time, final time and time step.

The ' $line$ ' command is used in generating the orbit (ODE solver only generates a discrete sequence of points).

Finally, one can superimpose orbits over a direction field (using a for loop to generate multiple orbits). Mind the ' $+=$ ' syntax inside the for loop and the fact that

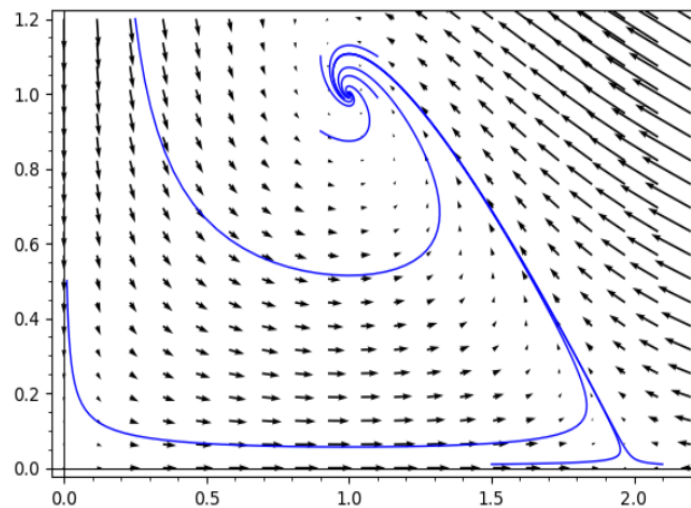
the 'show' command must remain outside the for loop (otherwise multiple plots are generated).

```
In [65]: # this is the solve block
times = srange(0,10,0.1)
ics = (-1.5, 3)
sol = desolve_odeint(f, ics, times, [x, y])

# the direction field
p_tot = plot_vector_field((f[0],f[1]), (x,0,2.2),(y,0,1.2))

# orbits are added to the same plot
for ics in [(0.01, 0.5), (2.1,0.01), (1.5,0.01), (0.9,1.1), (1.1, 1.1), (1.1, 0.99), (0.9,0.9), (0.25, 1.2)]:
    sol = desolve_odeint(f, ics, times, [x, y])
    p_tot += line(zip(sol[:,0],sol[:,1]))

show(p_tot)
```



### 3. Representing level curves using contour plots

First integrals of (nonlinear) planar systems can be represented/visualized in terms of their level curves, just as geographical landscapes are plotted on maps (this is why sometimes mathematicians use expressions like 'energy landscape').

**Sage** has a predefined command, called 'contour\_plot' for plotting level curves.

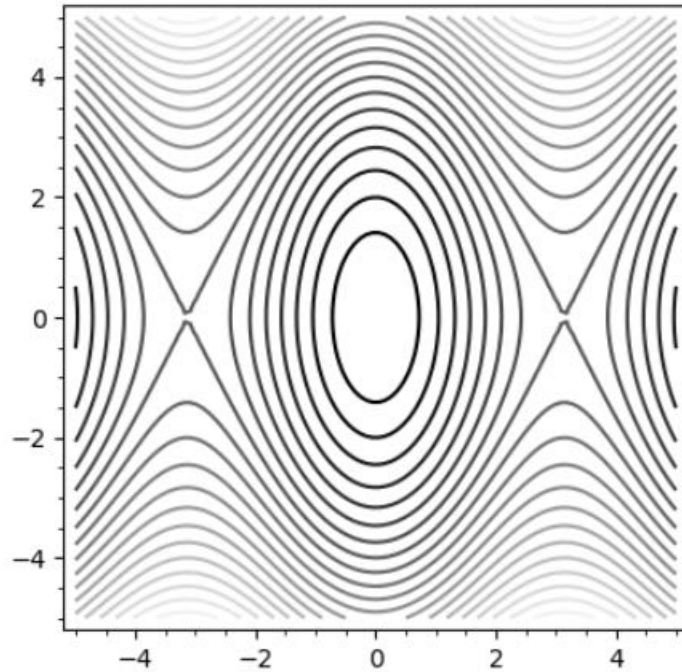
The arguments of the command include: the function to be plotted, the x and y ranges, the number of level curves to be plotted as well as visual options such as 'fill'.

Obviously, the scalar-valued function (energy) to be plotted must be defined first.

```
In [3]: x, y = var('x y')
        H(x,y) = y^2 - 8*cos(x)

        # we can plot the level curves of H
        contour_plot(H, (-5,5), (-5,5), fill=False, contours=20)
```

Out[3]:



Changing the graphical appearance options can be done in numerous ways

```
In [7]: x, y = var('x y')
        H(x,y) = y^2 - 8*cos(x)

        # we can plot the level curves of H
        contour_plot(H, (-5,5), (-5,5), contours=20)
```

Out[7]:

