# **Documentation**

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## **Types**

```
\# Main.Branch — Type.
```

A series of Solutions. One can push, unshift, pop, and shift new/old Solutions to/from it.

```
# Main.Solution — Type.
```

A Solution comprises a single solution vector and its parent

```
\# Main.ContinuationMethod — Type.
```

Continuation method implementations extend ContinuationMethod and the corresponding show and step. Responsible for changing the project itself

```
\# Main.PC — Type.
```

An "implementation" of a ContinuationMethod using a predictor-corrector-method with step-size adaption. Look up the theoretical documentation for more details.

```
\# Main.SystemCore — Type.
```

An abstract type that requires to implement the basic functions needed for path following: Homotopy H, its jacobian J (and show to display a GUI)

```
\# Main.Galerkin — Type.
```

"Implementation" of SystemCore; Only requires the ode f and its derivative f', the respective homotopy and its jacobian are derived. Again, look up the theoretical documentation for more detail.

```
\# Main.Project — Type.
```

Contains all found Branches

# Main.Session — Type.

Comprises everything needed for path following: Project, SystemCore, ContinuationMethod, and Visualization. Sessions can be managed via create, save, and load.

## **Utility Functions**

#### Differentiation

```
\# mbNewton.forwardDifference — Function.
forwardDifference(homotopy, v, epsilon)
forwardDifference(homotopy, epsilon)
Returns the difference quotient in v or an approximate jacobian using this
# mbNewton.centralDifference — Function.
centralDifference(homotopy, v, epsilon)
centralDifference(homotopy, epsilon)
Same as forwardDifference, but with the symmetric difference quotient.
# mbNewton.broyden — Function.
broyden(homotopy, jacobian)
Creates an approximate jacobian based on broyden-updates and maintainance
of an internal state. Consider this for performance improvements.
all jacobians can be used in
# mbNewton.newton — Function.
newton(homotopy, jacobian, v , pred[, init, callback, useOpt])
Newton's method. For example: v1 = newton(H, J, v0, predEps(0.001)).
```

### Galerkin

Returns all points from the transientIterations-times numerical integration of H with fixed step-size transientStepSize, and all points from the steadyStateIterations-times integration with fixed step-size steadyStateStepSize (assuming it hit the steady state part), and the periodicity of the found cycle.

```
# Main.findCyclePoincare — Function.
```

```
findCyclePoincare(F, y[, plane, clusterRating, nIntersections,
    maxCycles, sampleSize, transientIterations, transientStepSize,
    steadyStateStepSize])
extracts a single cycle of the steady state of ode F using poincare cuts through
the plane.
\# Main.prepareCycle — Function.
prepareCycle(data, h, P[, fac])
cut single cycle of length P*fac from data, resample, shift s.t. X(0) 0, Fourier
transform.
\# mbInterpolate.interpolateLanczos — Function.
interpolateLanczos(V, a::Integer)
simple periodic (!) Lanczos interpolation
\# mbInterpolate.interpolateTrigonometric — Function.
interpolateTrigonometric(a, a, b)
Returns trigonometric polynomial. Use with 2a,-2b and divide by 2m+1 to use
with rfft coefficients.
GUI
# Main.mkControlGrid — Function.
mkControlGrid(D, C)
Creates a grid of controls with labels, handlers and encapsulated storage (Dic-
tionary D) c in C is Tuple (name::String, ::DataType, init, v...)
\# Main.ctrl — Function.
ctrl(D, x)
Dictionary D, Tuple x=(name::String, ::DataType, init, v...); used by
mkControlGrid
Session Control
# Main.create — Function.
create(homotopy, jacobian, projection)
# Main.save — Function.
save(filename, session[, overwrite])
\# Main.load — Function.
```

load(filename, homotopy, jacobian, projection)

### ODE

```
# mbRK.rk — Function.
rk(butcherTableau)
returns a runge-kutta method using the respective tableau:
function(f, t0, y0, h, pred[, init, callback])
e.g. rk1, or rk4. Examines the ode f starting from t0, y0 with fixed stepsize h until pred evalutes to false.
```

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