

```
0 #
1 #
2 #
3 #
4 #
5 #
6 #
```

## IDAS integrator

We solve a system  $\dot{x}(t)=f(x(t),y(t),t)$   $z_0=g(x(t),y(t),t)$

```
15 from casadi import *
16 from numpy import *
17 from pylab import *
```

We solve the following simple dae system that describes the dynamics of a pendulum:

$$\begin{aligned} x' &= u, \quad y' = v, \quad u' = \lambda x, \quad v' = \lambda y - g \\ \text{s.t. } x^2 + y^2 &= L \end{aligned}$$

We retain g and L as parameters

[http://en.wikipedia.org/wiki/Differential\\_algebraic\\_equation#Examples](http://en.wikipedia.org/wiki/Differential_algebraic_equation#Examples)

```
26 L = SX.sym("L")
27 g = SX.sym("g")
```

differential states

```
30 x=SX.sym("x")
31 y=SX.sym("y")
32 u=SX.sym("u")
33 v=SX.sym("v")
```

algebraic states

```
36 lambda=SX.sym("lambda")
```

All states and parameters

```
39 x_all = vertcat(x,u,y,v)
40 z_all = lambda
41 p_all = vertcat(L,g)
```

the initial state of the pendulum

```
45 P_ = [5,10] # parameters
46 X_ = [3,-1.0/3,4,1.0/4] # differential states
47
48 XDOT_ = [-1.0/3,1147.0/240,1.0/4,-653.0/180] # state derivatives
49
50 Z_ = [1147.0/720] # algebraic state
```

We construct the DAE system

```
50 ode = vertcat(u, lambda*x, v, lambda*y+g)
51 alg = x**2+y**2-L**2
52 dae = {'x':x_all, 'z':z_all, 'p':p_all, 'ode':ode, 'alg':alg}
```

```
53 f = Function('f', [x_all, z_all, p_all], [ode, alg], ['x', 'z', 'p'], ['ode', 'alg'])
```

Let's check we have consistent initial conditions:

```
56 res = f(p=P_, x=X_, z=Z_)
57 print(res['ode']) # This should be same as XDOT_
```

```
[-0.333333, 4.77917, 0.25, 16.3722]
```

```
59 print(res['alg']) # This should be all zeros
```

```
0
```

Let's check our jacobian  $\frac{dg}{dy}$ :

```
62 j = jacobian(alg, lambda)
63 print(j)
```

```
00
```

Note that the jacobian is not invertible: it is not of DAE-index 1

This system is not solvable with idas, because it is of DAE-index 3. It is impossible to lambda from the last element of the residual.

We create a DAE system solver

```
70 I = integrator('I', 'idas', dae, {'calc_ic':False, 'init_xdot':XDOT_})
```

This system is not solvable with idas, because it is of DAE-index 3. It is impossible obtain lambda from the last element of the residual.

```
78 try:
79     I(p=P_, x0=X_, z0=Z_)
80 except Exception as e:
81     print(e)
82
```

```
.../casadi/interfaces/sundials/idas_interface.cpp:560: IDASolve returned "
IDA_CONV_FAIL". Consult IDAS documentation.
```

**Error:**

```
CasADi - 2018-02-11 02:38:00 WARNING("I:daeF failed: NaN detected for output ode, at
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CasADi - 2018-02-11 02:38:00 WARNING("I:daeF failed: NaN detected for output ode, at
CasADi - 2018-02-11 02:38:00 WARNING("I:daeF failed: NaN detected for output ode, at
At t = 0 and h = 2.06366e-19, the corrector convergence failed repeatedly or with |h|
```

We construct a reworked version of the DAE (index reduced), now it is DAE-index 1

```
81 ode = vertcat(u, lambda*x)
82 alg = vertcat(x**2+y**2-L**2, u*x+v*y, u**2-g*y+v**2+L**2*lambda)
83 x_all = vertcat(x,u)
84 z_all = vertcat(y,v,lambda)
85 dae = {'x':x_all, 'z':z_all, 'p':p_all, 'ode':ode, 'alg':alg}
```

```

86 f = Function('f', [x_all, z_all, p_all], [ode, alg], ['x', 'z', 'p'], ['ode',
    'alg'])
    the initial state of the pendulum
90 P_ = [5,10] # parameters
91
92 X_ = [3,-1.0/3] # differential states
93
94 XDOT_ = [-1.0/3,1147.0/240] # state derivatives
95
96 Z_ = [4,1.0/4,1147.0/720] # algebraic state
    Let's check we have consistent initial conditions:
97
98 res = f(p=P_, x=X_, z=Z_)
99 print(res['ode']) # This should be the same as XDOT_
100
101 [-0.333333, 4.77917]
102
103 print(res['alg']) # This should be all zeros
104
105 [0, 0, 0]
    Let's check our jacobian:
106 J = f.factory('J', f.name_in(), ['jac:alg:z'])
107 res = J(p=P_, x=X_, z=Z_)
108 print(array(res["jac_alg_z"]))
109
110 [[ 8.    0.    0. ]
111 [ 0.25  4.    0. ]
112 [-10.   0.5  25. ]]
    $frac{dg}{dy}$ is invertible this time.
    We create a DAE system solver
113
114 l = integrator('I', 'idas', dae, {'t0':0, 'tf':1, 'init_xdot':XDOT_})
115 res = l(p=P_, x0=X_, z0=Z_)
116 print(res['xf'])
117
118 [4.68624, 2.34688]

```

Although this initialisation is consistent, it coincides with a singular point.

## Possible problems

If you would initialize with:

```

115 P_ = [5,10] # parameters
116
117 X_ = [5,0] # states

```

You will get an error:

```

121 try:
122     l(p=P_, x0=X_, z0=Z_)
123 except Exception as e:
124     print(e)

```

```

.../casadi/interfaces/sundials/idas_interface.cpp:560: IDASolve returned "
IDA_TOO_MUCH_WORK". Consult IDAS documentation.

```

**Error:**

At t = 6.02906e-09, , mxstep steps taken before reaching tout.