idas

November 7, 2023

This file is part of CasADi.

CasADi -- A symbolic framework for dynamic optimization.

Copyright (C) 2010-2023 Joel Andersson, Joris Gillis, Moritz Diehl,

KU Leuven. All rights reserved.

Copyright (C) 2011-2014 Greg Horn

CasADi is free software; you can redistribute it and/or modify it under the terms of the GNU Lesser General Public License as published by the Free Software Foundation; either version 3 of the License, or (at your option) any later version.

CasADi is distributed in the hope that it will be useful, but WITHOUT ANY WARRANTY; without even the implied warranty of MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE. See the GNU Lesser General Public License for more details.

You should have received a copy of the GNU Lesser General Public License along with CasADi; if not, write to the Free Software Foundation, Inc., 51 Franklin Street, Fifth Floor, Boston, MA 02110-1301 USA

1 IDAS integrator

We solve a system $\dot{x}(t) = f(x(t), y(t), t) \ 0 = g(x(t), y(t), t)$

```
[1]: from casadi import *
from numpy import *
from pylab import *
```

We solve the following simple dae system that describes the dynamics of a pendulum: x' = u, y' = v, u' = lambda * x, v' = lambda * y - g s.t. $x^{2+y}2 = L$

We retain g and L as parameters http://en.wikipedia.org/wiki/Differential_algebraic_equation#Examples

```
[2]: L = SX.sym("L")
g = SX.sym("g")
```

differential states

```
[3]: x=SX.sym("x")
y=SX.sym("y")
u=SX.sym("u")
v=SX.sym("v")
```

algebraic states

```
[4]: lambd=SX.sym("lambda")
```

All states and parameters

```
[5]: x_all = vertcat(x,u,y,v)
z_all = lambd
p_all = vertcat(L,g)
```

the initial state of the pendulum

```
[6]: P_{-} = [5,10] # parameters X_{-} = [3,-1.0/3,4,1.0/4] # differential states XDOT_{-} = [-1.0/3,1147.0/240,1.0/4,-653.0/180] # state derivatives Z_{-} = [1147.0/720] # algebraic state
```

We construct the DAE system

Let's check we have consistent initial conditions:

```
[8]: res = f(p=P_, x=X_, z=Z_)
print(res['ode']) # This should be same as XDOT_
print(res['alg']) # This should be all zeros
```

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

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

```
[9]: j = jacobian(alg,lambd)
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

```
[10]: 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.

```
[11]: try:
    I(p=P_, x0=X_, z0=Z_)
    except Exception as e:
        print(e)
```

```
Error in Function::call for 'I' [IdasInterface] at .../casadi/core/function.cpp:1401:
Error in Function::call for 'I' [IdasInterface] at .../casadi/core/function.cpp:330: .../casadi/interfaces/sundials/idas_interface.cpp:596: IDASolve returned "IDA_CONV_FAIL". Consult IDAS documentation.
```

At t = 0 and h = 5.40977e-14, the corrector convergence failed repeatedly or with |h| = hmin.

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

```
[12]: ode = vertcat(u,lambd*x)
    alg = vertcat(x**2+y**2-L**2, u*x+v*y,u**2-g*y+v**2+L**2*lambd)
    x_all = vertcat(x,u)
    z_all = vertcat(y,v,lambd)
    dae = {'x':x_all, 'z':z_all, 'p':p_all, 'ode':ode, 'alg':alg}
    f = Function('f', [x_all, z_all, p_all], [ode, alg], ['x', 'z', 'p'], ['ode', u o'alg'])
```

the initial state of the pendulum

```
[13]: P_ = [5,10] # parameters

X_ = [3,-1.0/3] # differential states

XDOT_ = [-1.0/3,1147.0/240] # state derivatives

Z_ = [4,1.0/4,1147.0/720] # algebraic state
```

Let's check we have consistent initial conditions:

```
[14]: res = f(p=P_, x=X_, z=Z_)
print(res['ode']) # This should be the same as XDOT_
print(res['alg']) # This should be all zeros
```

```
[-0.333333, 4.77917]
[0, 0, 0]
```

Let's check our jacobian:

```
[15]: J = f.factory('J', f.name_in(), ['jac:alg:z'])
res = J(p=P_, x=X_, z=Z_)
print(array(res["jac_alg_z"]))
```

 $\frac{dg}{du}$ is invertible this time.

We create a DAE system solver

```
[16]: I = integrator('I', 'idas', dae, {'t0':0, 'tf':1, 'init_xdot':XDOT_})
res = I(p=P_, x0=X_, z0=Z_)
print(res['xf'])
```

[4.68624, 2.34688]

CasADi - 2023-11-07 01:54:38 WARNING("The options 't0', 'tf', 'grid' and 'output_t0' have been deprecated.

The same functionality is provided by providing additional input arguments to the 'integrator' function, in particular:

- * Call integrator(..., t0, tf, options) for a single output time, or
- * Call integrator(..., t0, grid, options) for multiple grid points.

The legacy 'output_t0' option can be emulated by including or excluding 't0' in 'grid'.

Backwards compatibility is provided in this release only.") [.../casadi/core/integrator.cpp:515]

2 Possible problems

If you would initialize with:

```
[17]: P_ = [5,10] # parameters
X_ = [5,0] # states
```

You will get an error:

```
[18]: try:
    I(p=P_, x0=X_, z0=Z_)
    except Exception as e:
        print(e)
```

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