

The Secret Life of the N-Matrix Gradient Descent Inverse Kinematics Revisited

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Motivation



- Robots should move!
- ▶ But describing robot motions in terms of joints is cumbersome



There are quite a lot of different IK approaches. But most real-world robots use geometric IKs.

So did we...

What do we want?



- Tell the robot where to move (not how)
 - Move left foot to position p_{left_foot}
 - ► Move right foot to position p_{right_foot}
 - While having the feet in a "natural" orientation
- Tell the robot how to move (constraints)
 - Keep the COM somewhere "safe"

We build tasks for all this



Joint:

Pretty self-explanatory (something elementary that moves)

Value:

The angle of a rotation-**joint** or the stroke of a piston-**joint**, etc.

Configuration/Pose/Posture:

A vector containing all joint-values

▶ Task:

Something the robot shall acomplish, eg:

- Move a body part to a certain target
- Orient a body part in a direction
- Have a certain value at a joint

Tasks generate Jacobians and errors

► Target/Method:

The representation of the solution space for a task eg:

- ▶ Point move a bodypart to a position $\rightarrow dim(taskspace) = 0$
- Line move a bodypart onto a line → dim(taskspace) = 1
- Plane move a bodypart onto a plane → dim(taskspace) = 2
- Space for highdimensional robots you could solve for n-dim spaces!



Demo

python eins.py

Derivation



Here comes the math!

Yey!



Define a loss function:

$$\mathcal{L}(\Delta q) = ||\phi(q + \Delta q) - y^*||_C^2 + ||\Delta q||_W^2$$
 (1)

W and C are weighting matrices. We really care about the solution of our Task $\phi(q+\Delta q)-y^*\to 0$ so we set $C\to \infty$ W scales the joint-changes. Since we treat every joint equally we choose W=I

with
$$\phi(q + \Delta q) = \phi(q) + J * \Delta q$$
 (local linearization) we get:
$$\mathcal{L}(\Delta q) = \underbrace{||\phi(q) + J * \Delta q - y^*||_C^2}_{\text{move to target}} + \underbrace{||\Delta q||_W^2}_{\text{be lazy}}$$



You know what time it is!

It's derive and set zero-time!



$$\frac{\delta}{\delta \Delta q} \mathcal{L}(\Delta q) = 0^{T} = ||\phi(q) + J * \Delta q - y^{*}||_{C}^{2} + ||\Delta q||_{W}^{2}$$

$$0^{T} = 2(\phi(q) - y^{*} + J\Delta q)^{T}CJ + 2\Delta q^{T}W$$

$$0^{T} = (J\Delta q - \vec{e})^{T}CJ + \Delta q^{T}W$$

$$0 = J^{T}C^{T}(J\Delta q - \vec{e}) + W^{T}\Delta q$$

$$0 = -J^{T}C^{T}\vec{e} + J^{T}C^{T}J\Delta q + W^{T}\Delta q$$

$$J^{T}C^{T}\vec{e} = (J^{T}C^{T}J + W^{T})\Delta q$$

$$(2)$$

$$(J^{T}C^{T}J + W^{T})^{-1}J^{T}C^{T}\vec{e} = \Delta q$$

use the Woodbury identity:

$$\Delta q = W^{T^{-1}} J^T (J W^{T^{-1}} J^T + C^{T^{-1}})^{-1} \vec{e}$$
 (3)



with

$$\lim_{C\to\infty} \Delta q = W^{T^{-1}} J^T (JW^{T^{-1}} J^T + C^{T^{-1}})^{-1}$$
 and $W = I$ we get:

$$J^{\#} = J^{\mathsf{T}} (JJ^{\mathsf{T}} + \epsilon I)^{-1}$$

that's the Moore-Penrose-inverse a.k.a. **pseudo-inverse**!!!!

Sweet!

(4)

Quick Recap



Where are we?

▶ We can calculate a change in configuration for a task!

What do we need?

Jacobians and error vectors?



 Errors are the difference of the current value in target space to the target in target space
 eq:

Where do I want to put the finger minus where is my finger now.

Jacobians are local linearization of the task space given infinitesimal changes in configuration space.

eg:

How does my finger move if I change the value of my elbow (and my wrist)



A Jacobian(-matrix) is the derivative of the **local linearization** of the forward kinematics of a **task**. eg:

How does my finger (**end-effector**) move if I change the value of my elbow (and my wrist).

Jacobians are dependant on the robots posture.



A Jacobian(-matrix) is the derivative of the **local linearization** of the forward kinematics of a **task**.

$$J = \frac{\delta}{\delta q} \phi(q) = \begin{pmatrix} \frac{\delta}{\delta q_1} \phi(q)_1 & \frac{\delta}{\delta q_2} \phi(q)_1 & \dots & \frac{\delta}{\delta q_n} \phi(q)_1 \\ \frac{\delta}{\delta q_1} \phi(q)_2 & \frac{\delta}{\delta q_2} \phi(q)_2 & \dots & \frac{\delta}{\delta q_n} \phi(q)_2 \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\delta}{\delta q_1} \phi(q)_d & \frac{\delta}{\delta q_2} \phi(q)_d & \dots & \frac{\delta}{\delta q_n} \phi(q)_d \end{pmatrix}$$
(5)

- *1 : Where goes the endeffector when we change the first joint
- *2 : Where goes the endeffector when we change the second joint
- *3: Where goes the endeffector when we change the n-th joint

Jacobians



- ▶ Jacobians are nxd-matrixes
 - one column for each joint
 - one row for each target-space dimension



```
class RotationJoint:
    def getLocationDerivative(self, location):
        return np.matrix([[-location[1, 0]],[location[0, 0]]])

    def getOrientationDerivative(self, orientation):
        return np.matrix([[-orientation[1, 0]],[orientation[0, 0]]])

class PistonJoint:
    def getLocationDerivative(self, location):
        return np.matrix([[1],[0]])

    def getOrientationDerivative(self, orientation):
```

return np.matrix([[0],[0]])





```
class LocationTask(PathedTask):
 def init (self, robot, path):
   PathedTask. init (self, robot, path)
 def getlacobian (self):
   transform = np.matrix(np.eye(3, 3))
   jacobian = np.matrix(np.zeros((2, self.dof)))
   for node, direction in self.path.path:
     subtransform = np.matrix(np.eye(2, 2))
     if node numbor != 0:
      idx = self.robot.getIndexOfActiveNode(node.name)
      if direction == Direction.FROM CHILD:
        jacobian[:,idx] = node.getLocationDerivative(transform[0:2,2])
      elif direction == Direction.FROM PARENT:
        [acobian[:,idx] = -node.getLocationDerivative(transform[0:2,2])
     if direction == Direction.FROM CHILD:
      subtransform = node.getTransform()[0:2,0:2]
      transform = node.qetTransform() * transform
     elif direction == Direction.FROM PARENT or direction == Direction.LINK:
      subtransform = node.getBackTransform()[0:2,0:2]
      transform = transform * node.getBackTransform()
     jacobian = subtransform * jacobian
   return jacobian
 def getError(self):
   return self.method.getTarget() — self.method.transform(self.getCurrentValue())
```



Demo

python linearlK1.py

Basic Solver



```
def solve_simple(robot, task, epsilon=0):
    helper = J * J.transpose()
    J_pinv = J.transpose() * inv(helper * epsilon * np.matrix(np.eye(helper.shape[0])))
    dq = J_pinv * task.getError()
    return dq
```

Multiple Tasks



But wait there is more!

what if we want to do several things simultaneously?

Multiple Tasks



Simultaneous tasks:

"stack" the jacobians and the error vectors into a big-jacobian and a big-error



Demo

python linearIK2.py



```
def solve_simple(robot, tasks, epsilon=0):
    numCols = robot.getDOF()
    bigJacobian = np.matrix(np.zeros((0, numCols)))
    bigError = np.matrix(np.zeros((0, 1)))
    for task in tasks:
        bigJacobian = np.concatenate((bigJacobian, task.getJacobian()))
        bigError = np.concatenate((bigJacobian, task.getJacobian()))
        bigFror = np.concatenate((bigJacobian, task.getError()))
        helper = bigJacobian * bigJacobian.transpose()
        J_pinv = bigJacobian.transpose() * inv(helper * epsilon * np.matrix(np.eye(helper.shape[0])))
        dq = J_pinv * bigError
        return da
```



But wait there is even more!!!

- ► Tasks might not fully define the robot
 - For the remaining degrees of freedom we want to solve other tasks
 - ► Those other tasks must not interfere with the primary tasks!



when rank(J) ≠ numDOF:
J does not fully utilize the robot

$$J^{\#}J \tag{6}$$

is the (orthogonal) range-projector of J

$$I - J^{\#}J = \mathcal{N} \tag{7}$$

is the (orthogonal) **nullspace-projector** of J



• any vector we multiply with $\mathcal N$ will be within the **nullspace** of $\mathcal J$ This is what we are looking for: $\mathbf 0 = \mathcal J \mathcal N \vec{\mathbf V}$



Utilizing the \mathcal{N} -matrix:

$$\Delta q = \Delta q_1 + \Delta q_2$$

$$\Delta q_1 = J_1^\# \vec{e_1}$$

$$\Delta q_2 = \mathcal{N}_1 J_2^\# \vec{e_2}$$
(8)

This is pretty much exactly what you'll find in textbooks



```
def solve simple(robot, tasks, epsilon=0):
 numCols = robot.getDOF()
 da = np.matrix(np.zeros((numCols. 1)))
 Nv = np.matrix(np.eve(numCols, numCols))
 reallyBiglacobian = np.matrix(np.zeros((0, numCols)))
 for taskGroup in taskGroups:
   biglacobian = np.matrix(np.zeros((0. numCols)))
   bigError = np.matrix(np.zeros((0, 1)))
   for task in taskGroup:
     biglacobian = np.concatenate((biglacobian, task.getlacobian()))
     bigError = np.concatenate((bigError, task.getError()))
   iacobian pinv = pinv(bigJacobian, epsilon)
   dq += Ny * jacobian pinv * bigError
   reallyBiglacobian = np.concatenate((reallyBiglacobian, biglacobian))
   Nv = np.matrix(np.eve(numCols. numCols)) - pinv(reallvBiglacobian. nullspaceEpsilon) *
      reallyBiglacobian
 return da
```



Demo

python linearlK3.py python linearlK3_bad.py



What happened?

▶ The calculation for Δq_2 does not yield optimal solutions!



What happened?

- ▶ The calculation for Δq_2 does not yield optimal solutions!
- ▶ Looks like we're utilizing the N-matrix incorrecty!



Lets look again at:

$$\mathcal{L}(\Delta q) = ||\phi(q + \Delta q) - y^*||_C^2 + ||\Delta q||_W^2$$
(11)

$$\Delta q = W^{T^{-1}} J^T (J W^{T^{-1}} J^T + C^{T^{-1}})^{-1} \vec{e}$$
 (12)

- C weights the error in task-space (which is important)
- W weights the error in joint-space
 - we need a matrix that punishes usage of joints that are already utilized
 - $W = \mathcal{N}^{-1}$ does exactely what we need!



$$\Delta q_i = \mathcal{N}_i J_i^T (J_i \mathcal{N}_i J_i^T + \epsilon I)^{-1} \vec{e}_i \qquad (13)$$

with:

$$\hat{J}_{0,i-1} = \begin{pmatrix} J_{0,0} \\ \vdots \\ J_{0,0n} \\ J_{1,0} \\ \vdots \\ J_{1,0n} \end{pmatrix} \text{ jacobians from task-group 1}$$

$$\vdots \\ J_{i-1,0} \\ \vdots \\ J_{i-1,i-1n} \\ \end{pmatrix} \text{ jacobians from task-group i-1}$$

$$\mathcal{N}_i = I - \hat{J}_{0,i-1}^{\theta} \hat{J}_{0,i-1}$$

(14)

With \mathcal{N} as W^{-1} we get:

$$\Delta q_2 = \mathcal{N} J_{2\mathcal{N}_1}^{\#} \vec{e}_2$$

$$= \begin{pmatrix} 0 & 0 \\ 0 & 0 \\ 1 & 0 \\ 0 & 1 \end{pmatrix} \vec{e}_2$$
(15)



Demo

python linearlK4.py python linearlK4_better.py



But wait there is even more again!!!!

- ► The robot still converges weirdly
 - We calculate Δq as

$$\Delta q = \sum_{i}^{N} \Delta q_{i} \tag{16}$$

▶ But we could already know q when we calculate Δq_i !

$$q_i = q_{i-1} + \Delta q_i \tag{17}$$

read this as:

• q_i is the robots posture after applying Δq_i



Demo

python linearIK5.py python nice_robot.py



This presentation and the code can be found here:





- Marc Toussaint
 Robotics Course
 - Michael Gienger and Marc Toussaint and Christian Goerick
 Whole-body Motion Planning Building Blocks for Inteligent Motion
 Planning for Humanoid Robots
 online
 - Gunüter Schreiber, Christian Ott, Gerd Hirzinger Interactive Redundant Robotics: Control of the Inverted Pendulum with Nullspace Motion