## Hartley - Zisserman reading club

## Part I: Hartley and Zisserman Appendix 6:

 Iterative estimation methodsPart II: Zhengyou Zhang:

A Flexible New Technique for Camera Calibration

Presented by Daniel Fontijne

## HZ Appendix 6: Iterative estimation methods

Topics:

- Basic methods: Newton, Gauss-Newton, gradient descent.
- Levenberg-Marquardt.
- Sparse Levenberg-Marquardt.
- Applications to homography, fundamental matrix, bundle adjustment.
- Sparse methods for equations solving.
- Robust cost functions.
- Parameterization.

Lecture notes which I found useful
(methods for non-linear least squares problems):
http://www2.imm.dtu.dk/pubdb/views/edoc_download.php/3215/pdf/imm3215.pdf

## Iterative estimation methods

Problem: how to find minimum of non-linear functions?

## Iterative estimation methods

Problem: how to find minimum of non-linear functions?
Examples of HZ problems: -homography estimation.
-fundamental matrix estimation.
-multiple image bundle adjustment.
-camera calibration (Zhang paper).

## Iterative estimation methods

Problem: how to find minimum of non-linear functions?
Examples of HZ problems:
-homography estimation.
-fundamental matrix estimation.
-multiple image bundle adjustment.
-camera calibration (Zhang paper).
Examples of my recent problems:
-optimization of skeleton geometry given marker data.
-optimization of skeleton pose given marker data.

## Iterative estimation methods

Problem: how to find minimum of non-linear functions?
Examples of HZ problems:
-homography estimation.
-fundamental matrix estimation.
-multiple image bundle adjustment.
-camera calibration (Zhang paper).
Examples of my recent problems:
-optimization of skeleton geometry given marker data.
-optimization of skeleton pose given marker data.
Central approach of Appendix 6: Levenberg-Marquardt.
Questions: Pronunciation? Why LM?

## Newton iteration

UNIVERSITEIT
VAN
Amsterdam

Goal: minimize $X=f(P)$ for $P$.
X is the measurement vector.
$\mathbf{P}$ is the parameter vector.
f is some non-linear function.

## Newton iteration

UNIVERSITEIT
VAN
AMSTERDAM

Goal: minimize $X=f(P)$ for $P$.
$\mathbf{X}$ is the measurement vector.
$\mathbf{P}$ is the parameter vector.
f is some non-linear function.
In other words:
Minimize $\epsilon=\mathbf{X}-\mathbf{f}(\mathbf{P})$.

## Newton iteration

Goal: minimize $X=f(P)$ for $P$.
X is the measurement vector.
$\mathbf{P}$ is the parameter vector.
f is some non-linear function.
In other words:
Minimize $\boldsymbol{\epsilon}=\mathbf{X}-\mathbf{f}(\mathbf{P})$.
We assume $\mathbf{f}$ is locally linear at each $\mathbf{P}_{i}$, then $\mathbf{f}\left(\mathbf{P}_{i}+\Delta_{i}\right)=\mathbf{f}\left(\mathbf{P}_{i}\right)+\mathrm{J}_{i} \Delta_{i}$,
where matrix $\mathrm{J}_{i}$ is the Jacobian $\partial \mathbf{f} / \partial \mathbf{P}$ at $\mathbf{P}_{i}$.

## Newton iteration

Goal: minimize $X=f(P)$ for $P$.
X is the measurement vector.
$\mathbf{P}$ is the parameter vector.
f is some non-linear function.
In other words:
Minimize $\boldsymbol{\epsilon}=\mathbf{X}-\mathbf{f}(\mathbf{P})$.
We assume $\mathbf{f}$ is locally linear at each $\mathbf{P}_{i}$, then $\mathbf{f}\left(\mathbf{P}_{i}+\Delta_{i}\right)=\mathbf{f}\left(\mathbf{P}_{i}\right)+\mathrm{J}_{i} \Delta_{i}$,
where matrix $\mathrm{J}_{i}$ is the Jacobian $\partial \mathbf{f} / \partial \mathbf{P}$ at $\mathbf{P}_{i}$.
So we want to minimize $\left\|\epsilon_{i}+\mathrm{J}_{i} \Delta_{i}\right\|$ for some vector $\Delta_{i}$.

## Newton iteration

Goal: minimize $X=f(P)$ for $P$.
$\mathbf{X}$ is the measurement vector.
$\mathbf{P}$ is the parameter vector.
f is some non-linear function.
In other words:
Minimize $\boldsymbol{\epsilon}=\mathbf{X}-\mathbf{f}(\mathbf{P})$.
We assume $\mathbf{f}$ is locally linear at each $\mathbf{P}_{i}$, then
$\mathbf{f}\left(\mathbf{P}_{i}+\Delta_{i}\right)=\mathbf{f}\left(\mathbf{P}_{i}\right)+\mathrm{J}_{i} \Delta_{i}$,
where matrix $\mathrm{J}_{i}$ is the Jacobian $\partial \mathbf{f} / \partial \mathbf{P}$ at $\mathbf{P}_{i}$.
So we want to minimize $\left\|\epsilon_{i}+\mathrm{J}_{i} \Delta_{i}\right\|$ for some vector $\Delta_{i}$.
Find $\Delta_{i}$ either using normal equations: $\mathrm{J}_{i}^{\mathrm{T}} \mathrm{J}_{i} \Delta=-\mathrm{J}_{i}^{\mathrm{T}} \epsilon_{i}$
or using pseudo-inverse: $\Delta_{i}=-\mathrm{J}_{i}^{+} \epsilon_{i}$.
Iterate until convergence . . .

## Gauss-Newton method

UNIVERSITEIT
VAN
AMSTERDAM

Suppose we want to minimize some cost function
$g(\mathbf{P})=\frac{1}{2}\|\boldsymbol{\epsilon}(\mathbf{P})\|^{2}=\frac{1}{2} \boldsymbol{\epsilon}(\mathbf{P})^{\mathrm{T}} \boldsymbol{\epsilon}(\mathbf{P})$.

## Gauss-Newton method

Suppose we want to minimize some cost function
$g(\mathbf{P})=\frac{1}{2}\|\boldsymbol{\epsilon}(\mathbf{P})\|^{2}=\frac{1}{2} \boldsymbol{\epsilon}(\mathbf{P})^{\mathrm{T}} \boldsymbol{\epsilon}(\mathbf{P})$.
We may expand in a Taylor series up to second degree $g(\mathbf{P}+\Delta)=g+g_{\mathbf{P}} \Delta+\Delta^{\mathrm{T}} g_{\mathbf{P P}} \Delta / 2$, where subscript P denotes differentiation.

## Gauss-Newton method

Suppose we want to minimize some cost function
$g(\mathbf{P})=\frac{1}{2}\|\boldsymbol{\epsilon}(\mathbf{P})\|^{2}=\frac{1}{2} \boldsymbol{\epsilon}(\mathbf{P})^{\mathrm{T}} \boldsymbol{\epsilon}(\mathbf{P})$.
We may expand in a Taylor series up to second degree
$g(\mathbf{P}+\Delta)=g+g_{\mathbf{P}} \Delta+\Delta^{\mathrm{T}} g_{\mathbf{P P}} \Delta / 2$,
where subscript P denotes differentiation.
Differentiating w.r.t. $\Delta$, setting to zero results in $g_{\mathbf{P P}} \Delta=-g_{\mathbf{P}}$. Using this equation we could compute $\Delta$ if we knew $g_{\mathbf{P P}}$ and $g_{\mathbf{P}}$.

## Gauss-Newton method

Suppose we want to minimize some cost function
$g(\mathbf{P})=\frac{1}{2}\|\boldsymbol{\epsilon}(\mathbf{P})\|^{2}=\frac{1}{2} \boldsymbol{\epsilon}(\mathbf{P})^{\mathrm{T}} \boldsymbol{\epsilon}(\mathbf{P})$.
We may expand in a Taylor series up to second degree
$g(\mathbf{P}+\Delta)=g+g_{\mathbf{P}} \Delta+\Delta^{\mathrm{T}} g_{\mathbf{P P}} \Delta / 2$,
where subscript P denotes differentiation.
Differentiating w.r.t. $\Delta$, setting to zero results in $g_{\mathbf{P P}} \Delta=-g_{\mathbf{P}}$. Using this equation we could compute $\Delta$ if we knew $g_{\mathbf{P P}}$ and $g_{\mathbf{P}}$.

Gradient vector: $g_{\mathrm{P}}=\epsilon_{\mathrm{P}}^{\mathrm{T}} \boldsymbol{\epsilon}=\mathrm{J}^{\mathrm{T}} \boldsymbol{\epsilon}$. Intuition?
Hessian: $g_{\mathbf{P P}}=\boldsymbol{\epsilon}_{\mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon}_{\mathbf{P}}+\boldsymbol{\epsilon}_{\mathbf{P} \mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon} \approx \mathrm{J}^{\mathrm{T}} \mathrm{J}$. Assume linear again . . .

## Gauss-Newton method

Suppose we want to minimize some cost function
$g(\mathbf{P})=\frac{1}{2}\|\boldsymbol{\epsilon}(\mathbf{P})\|^{2}=\frac{1}{2} \boldsymbol{\epsilon}(\mathbf{P})^{\mathrm{T}} \boldsymbol{\epsilon}(\mathbf{P})$.
We may expand in a Taylor series up to second degree
$g(\mathbf{P}+\Delta)=g+g_{\mathbf{P}} \Delta+\Delta^{\mathrm{T}} g_{\mathbf{P P}} \Delta / 2$,
where subscript $\mathbf{P}$ denotes differentiation.
Differentiating w.r.t. $\Delta$, setting to zero results in $g_{\mathbf{P P}} \Delta=-g_{\mathbf{P}}$. Using this equation we could compute $\Delta$ if we knew $g_{\mathbf{P P}}$ and $g_{\mathbf{P}}$.

Gradient vector: $g_{\mathrm{P}}=\boldsymbol{\epsilon}_{\mathrm{P}}^{\mathrm{T}} \boldsymbol{\epsilon}=\mathrm{J}^{\mathrm{T}} \boldsymbol{\epsilon}$. Intuition?
Hessian: $g_{\mathbf{P P}}=\boldsymbol{\epsilon}_{\mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon}_{\mathbf{P}}+\boldsymbol{\epsilon}_{\mathbf{P} \mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon} \approx \mathrm{J}^{\mathrm{T}} \mathrm{J}$. Assume linear again . . .
Putting it all together we get $\mathrm{J}^{\mathrm{T}} \mathrm{J} \Delta=-\mathrm{J}^{\mathrm{T}} \epsilon$.
So we arrive at the normal equations again.
(So what was the point?)

## Gradient descent

Universiteit
VAN
AMSTERDAM

Gradient descent or steepest descent searches in the direction of most rapid decrease $-g_{\mathbf{P}}=-\epsilon_{\mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon}$.

## Gradient descent

Gradient descent or steepest descent searches in the direction of most rapid decrease $-g_{\mathbf{P}}=-\epsilon_{\mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon}$.

So we take steps $\lambda \Delta=-g_{\mathbf{P}}$ where $\lambda$ controls the step size and is found through line search.

## Gradient descent

Gradient descent or steepest descent searches in the direction of most rapid decrease $-g_{\mathbf{P}}=-\epsilon_{\mathbf{P}}^{\mathrm{T}} \boldsymbol{\epsilon}$.

So we take steps $\lambda \Delta=-g_{\mathbf{P}}$ where $\lambda$ controls the step size and is found through line search.

A problem is zig-zagging which can cause slow convergence:


## Levenberg-Marquardt

UNIVERSITEIT
VAN
Amsterdam

Levenberg-Marquardt is a blend of Gauss-Newton and gradient descent. Update equation:
$\left(\mathrm{J}^{\mathrm{T}} \mathrm{J}+\lambda \mathrm{I}\right) \Delta=-\mathrm{J}^{\mathrm{T}} \boldsymbol{\epsilon}$.

## Levenberg-Marquardt

Levenberg-Marquardt is a blend of Gauss-Newton and gradient descent. Update equation:
$\left(\mathrm{J}^{\mathrm{T}} \mathrm{J}+\lambda \mathrm{I}\right) \Delta=-\mathrm{J}^{\mathrm{T}} \boldsymbol{\epsilon}$.
Algorithm:

- Initially set $\lambda=10^{-3}$.
- Try update equation.
- If improvement: divide $\lambda$ by 10. I.e., shift towards Gauss-Newton.
- Else: multiply $\lambda$ by 10. I.e., shift towards gradient descent.


## Levenberg-Marquardt

Levenberg-Marquardt is a blend of Gauss-Newton and gradient descent. Update equation:
$\left(\mathrm{J}^{\mathrm{T}} \mathrm{J}+\lambda \mathrm{I}\right) \Delta=-\mathrm{J}^{\mathrm{T}} \boldsymbol{\epsilon}$.
Algorithm:

- Initially set $\lambda=10^{-3}$.
- Try update equation.
- If improvement: divide $\lambda$ by 10. I.e., shift towards Gauss-Newton.
- Else: multiply $\lambda$ by 10. I.e., shift towards gradient descent.

The idea is (?):
-take big gradient descent steps far away from minimum.
-take Gauss-Newton steps near (hopefully quadratic) minimum.

## Sparse Levenberg-Marquardt 1/2

UNIVERSITEIT
VAN
AMSTERDAM

In many estimation problems, the Jacobian is sparse. One should this to lower the time complexity (sometimes even from $O\left(n^{3}\right)$ to $O(n)$ ).

## Sparse Levenberg-Marquardt 1/2

In many estimation problems, the Jacobian is sparse.
One should this to lower the time complexity (sometimes even from $O\left(n^{3}\right)$ to $O(n)$ ).

In the example, the parameters are partitioned into two blocks:
$\mathbf{P}=\left(\mathbf{a}^{T}, \mathbf{b}^{T}\right)^{T}$
The Jacobian then has the form $\mathrm{J}=[A \mid B]$, with
$A=[\partial \hat{\mathbf{X}} / \partial \mathbf{a}], \quad B=[\partial \hat{\mathbf{X}} / \partial \mathbf{b}]$.

## Sparse Levenberg-Marquardt 1/2

In many estimation problems, the Jacobian is sparse.
One should this to lower the time complexity (sometimes even from $O\left(n^{3}\right)$ to $O(n)$ ).

In the example, the parameters are partitioned into two blocks:
$\mathbf{P}=\left(\mathbf{a}^{T}, \mathbf{b}^{T}\right)^{T}$
The Jacobian then has the form $\mathrm{J}=[A \mid B]$, with
$A=[\partial \hat{\mathbf{X}} / \partial \mathbf{a}], \quad B=[\partial \hat{\mathbf{X}} / \partial \mathbf{b}]$.
Using $A$ and $B$, the normal equations $\left(\mathrm{J}^{\mathrm{T}} \mathrm{J}\right) \Delta=-\mathrm{J}^{\mathrm{T}} \boldsymbol{\epsilon}$ take on the the form

$$
\left[\begin{array}{c|c}
A^{T} A & A^{T} B \\
\hline B^{T} A & B^{T} B
\end{array}\right]\binom{\boldsymbol{\delta}_{\mathbf{a}}}{\hline \boldsymbol{\delta}_{\mathbf{b}}}=\binom{A^{T} \boldsymbol{\epsilon}}{\hline B^{T} \boldsymbol{\epsilon}} .
$$

## Sparse Levenberg-Marquardt 2/2

If the normal equations are written as (what's with the *?)

$$
\left[\begin{array}{cc}
\mathrm{U}^{*} & \mathrm{~W} \\
\mathrm{~W}^{T} & \mathrm{~V}^{*}
\end{array}\right]\binom{\boldsymbol{\delta}_{\mathrm{a}}}{\boldsymbol{\delta}_{\mathrm{b}}}=\binom{\boldsymbol{\epsilon}_{A}}{\boldsymbol{\epsilon}_{B}},
$$

we can rewrite this to

$$
\left[\begin{array}{cc}
\mathrm{U}^{*}-\mathrm{WV}^{*-1} \mathrm{~W}^{T} & 0 \\
\mathrm{~W}^{T} & \mathrm{~V}^{*}
\end{array}\right]\binom{\boldsymbol{\delta}_{\mathrm{a}}}{\boldsymbol{\delta}_{\mathrm{b}}}=\binom{\boldsymbol{\epsilon}_{A}-\mathrm{WV}^{*-1} \boldsymbol{\epsilon}_{B}}{\boldsymbol{\epsilon}_{B}}
$$

by multiplying on the left by $\left[\begin{array}{cc}\mathrm{I} & \mathrm{WV}^{*-1} \\ 0 & \mathrm{I}\end{array}\right]$.
Now first solve the top half, then the lower half using back-substitution.

Robust cost functions 1/5

UNIVERSITEIT
VAN
Amsterdam


## Robust cost functions 2/5

UNIVERSITEIT
VAN
AMSTERDAM

Squared-error is not usable unless outliers are filtered out.

## Robust cost functions 2/5

Squared-error is not usable unless outliers are filtered out. Alternatives:

- Blake-Zisserman: outliers are given a constant cost.

Disadvantages: not a PDF, not convex.

## Robust cost functions 2/5

Squared-error is not usable unless outliers are filtered out. Alternatives:

- Blake-Zisserman: outliers are given a constant cost.

Disadvantages: not a PDF, not convex.

- Corrupted Gaussian: blend two Gaussians, one for inliers and one for outliers.

Disadvantages: not convex.

## Robust cost functions 2/5

Squared-error is not usable unless outliers are filtered out. Alternatives:

- Blake-Zisserman: outliers are given a constant cost.

Disadvantages: not a PDF, not convex.

- Corrupted Gaussian: blend two Gaussians, one for inliers and one for outliers.

Disadvantages: not convex.

- Cauchy: (?).
disadvantages: not convex.


## Robust cost functions 2/5

## Squared-error is not usable unless outliers are filtered out.

 Alternatives:- Blake-Zisserman: outliers are given a constant cost.

Disadvantages: not a PDF, not convex.

- Corrupted Gaussian: blend two Gaussians, one for inliers and one for outliers.

Disadvantages: not convex.

- Cauchy: (?).
disadvantages: not convex.
- L1: absolute error (not squared).

Disadvantages: not differentiable at 0 , minimum is not at a single point when summed.

## Robust cost functions 2/5

## Squared-error is not usable unless outliers are filtered out.

 Alternatives:- Blake-Zisserman: outliers are given a constant cost.

Disadvantages: not a PDF, not convex.

- Corrupted Gaussian: blend two Gaussians, one for inliers and one for outliers.

Disadvantages: not convex.

- Cauchy: (?).
disadvantages: not convex.
- L1: absolute error (not squared).

Disadvantages: not differentiable at 0 , minimum is not at a single point when summed.

- Huber cost function: like L1, but 'rounded'.

Disadvantages: non-continuous derivative from $2^{n d}$ order and up.

## Robust cost functions 2/5

## Squared-error is not usable unless outliers are filtered out.

 Alternatives:- Blake-Zisserman: outliers are given a constant cost.

Disadvantages: not a PDF, not convex.

- Corrupted Gaussian: blend two Gaussians, one for inliers and one for outliers. Disadvantages: not convex.
- Cauchy: (?).
disadvantages: not convex.
- L1: absolute error (not squared).

Disadvantages: not differentiable at 0 , minimum is not at a single point when summed.

- Huber cost function: like L1, but 'rounded'.

Disadvantages: non-continuous derivative from $2^{n d}$ order and up.

- Pseudo Huber: like Huber, but with continuous derivatives.

Robust cost functions 3/5

UNIVERSITEIT VAN
AMSTERDAM

Figure A.6.5






Robust cost functions 4/5

UNIVERSITEIT VAN

Figure A.6.6





## Robust cost functions 5/5

Summary:

- Squared-error cost function is very susceptible to outliers.


## Robust cost functions 5/5

Summary:

- Squared-error cost function is very susceptible to outliers.
- The non-convex functions (like L1 and corrupted Gaussian) may be good, but they have local minima. So do not use them unless already close to true minimum.


## Robust cost functions 5/5

Summary:

- Squared-error cost function is very susceptible to outliers.
- The non-convex functions (like L1 and corrupted Gaussian) may be good, but they have local minima. So do not use them unless already close to true minimum.
- Best: Huber and Pseudo-Huber.


## Fooling LM implementations

Universiteit
VAN
AMSTERDAM

Most implementations of Levenberg-Marquardt use the squared error cost function. What if you want a different cost function $C$ instead?

## Fooling LM implementations

Most implementations of Levenberg-Marquardt use the squared error cost function. What if you want a different cost function $C$ instead?

Replace the each difference $\delta_{i}$ with a weighted version

$$
\delta_{i}^{\prime}=w_{i} \delta_{i}
$$

such that

$$
\left\|\delta_{i}\right\|^{2}=w_{i}^{2}\left\|\delta_{i}\right\|^{2}=C\left(\left\|\delta_{i}\right\|\right)
$$

## Fooling LM implementations

Most implementations of Levenberg-Marquardt use the squared error cost function. What if you want a different cost function $C$ instead?

Replace the each difference $\delta_{i}$ with a weighted version

$$
\delta_{i}^{\prime}=w_{i} \delta_{i}
$$

such that

$$
\left\|\delta_{i}\right\|^{2}=w_{i}^{2}\left\|\delta_{i}\right\|^{2}=C\left(\left\|\delta_{i}\right\|\right)
$$

Thus

$$
w_{i}=\frac{\sqrt{C\left(\left\|\delta_{i}\right\|\right)}}{\left\|\delta_{i}\right\|}
$$

## Parameterization for Levenberg-Marquardt

A good parameterization for use with LM is singularity free (at least in area visited during optimization). This means:

- continuous,
- differentiable,
- one-to-one.


## Parameterization for Levenberg-Marquardt

A good parameterization for use with LM is singularity free (at least in area visited during optimization). This means:

- continuous,
- differentiable,
- one-to-one.

So latitude-longitude is not suitable to parameterize sphere.

## Parameterization for Levenberg-Marquardt

A good parameterization for use with LM is singularity free (at least in area visited during optimization). This means:

- continuous,
- differentiable,
- one-to-one.

So latitude-longitude is not suitable to parameterize sphere.
And Euler angles are not suitable to parameterize rotations.

## Parameterization for Levenberg-Marquardt

A good parameterization for use with LM is singularity free (at least in area visited during optimization). This means:

- continuous,
- differentiable,
- one-to-one.

So latitude-longitude is not suitable to parameterize sphere.
And Euler angles are not suitable to parameterize rotations.
Gauge freedom?

## Parameterization for Levenberg-Marquardt

A good parameterization for use with LM is singularity free (at least in area visited during optimization). This means:

- continuous,
- differentiable,
- one-to-one.

So latitude-longitude is not suitable to parameterize sphere.
And Euler angles are not suitable to parameterize rotations.
Gauge freedom?
Variance?

## Parameterization of 3-D rotations

UnIVERSITEIT
VAN

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.

## Parameterization of 3-D rotations

Universiteit
VAN

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.
Observations: (these are mostly just general observations about $\log$ (rotation))

- Identity rotation: $\mathrm{t}=\mathbf{0}$.


## Parameterization of 3-D rotations

UNIVERSITEIT
VAN

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.
Observations: (these are mostly just general observations about $\log$ (rotation))

- Identity rotation: $\mathbf{t}=\mathbf{0}$.
- Inverse rotation: -t.


## Parameterization of 3-D rotations

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.
Observations: (these are mostly just general observations about $\log$ (rotation))

- Identity rotation: $\mathbf{t}=\mathbf{0}$.
- Inverse rotation: -t.
- Small rotation: the rotation matrix is $\mathrm{I}+[\mathbf{t}]_{\times}$.


## Parameterization of 3-D rotations

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.
Observations: (these are mostly just general observations about $\log$ (rotation))

- Identity rotation: $\mathbf{t}=\mathbf{0}$.
- Inverse rotation: -t.
- Small rotation: the rotation matrix is $\mathrm{I}+[\mathbf{t}]_{\times}$.
- For small rotations: $\mathrm{R}\left(\mathrm{t}_{1}\right) \mathrm{R}\left(\mathrm{t}_{2}\right) \approx \mathrm{R}\left(\mathrm{t}_{1}+\mathrm{t}_{2}\right)$.


## Parameterization of 3-D rotations

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.
Observations: (these are mostly just general observations about $\log$ (rotation))

- Identity rotation: $\mathbf{t}=\mathbf{0}$.
- Inverse rotation: -t.
- Small rotation: the rotation matrix is $I+[\mathbf{t}]_{\times}$.
- For small rotations: $\mathrm{R}\left(\mathrm{t}_{1}\right) \mathrm{R}\left(\mathrm{t}_{2}\right) \approx \mathrm{R}\left(\mathrm{t}_{1}+\mathrm{t}_{2}\right)$.
- All rotations can be represented by t with $\|\mathrm{t}\| \leq \pi$. When $\|\mathbf{t}\|=n 2 \pi$, ( $n$ positive integer) you get identity rotation again (singularity).


## Parameterization of 3-D rotations

3-D rotation matrix: 9 elements, only 3 degrees of freedom.
Angle-axis (3-vector) representation: 3 elements, 3 d.o.f.
Observations: (these are mostly just general observations about $\log$ (rotation))

- Identity rotation: $\mathbf{t}=\mathbf{0}$.
- Inverse rotation: -t.
- Small rotation: the rotation matrix is $I+[\mathbf{t}]_{\times}$.
- For small rotations: $\mathrm{R}\left(\mathrm{t}_{1}\right) \mathrm{R}\left(\mathrm{t}_{2}\right) \approx \mathrm{R}\left(\mathrm{t}_{1}+\mathrm{t}_{2}\right)$.
- All rotations can be represented by t with $\|\mathrm{t}\| \leq \pi$. When $\|\mathbf{t}\|=n 2 \pi$, ( $n$ positive integer) you get identity rotation again (singularity).
- Normalization: stay away from $\|\mathbf{t}\|=2 \pi$.


## Parameterization of homogeneous vectors

UNIVERSITEIT
VAN

Let $\mathbf{v}$ be a $n$-D-vector (already stripped of 'extra' homogeneous coordinate?).

Then parameterize it as $n+1$ vector:
$\bar{v}=\left(\operatorname{sinc}(\|\mathbf{v}\| / 2) \mathbf{v}^{T}, \cos (\|\mathbf{v}\| / 2)\right)^{T}$.

## Parameterization of the $n$-sphere

UNIVERSITEIT
VAN
Amsterdam

How to parameterize unit vectors $\mathbf{x}$ ?
Compute Householder matrix (reflection) such that $\mathrm{H}_{\mathbf{v}(\mathrm{x})} \mathbf{x}=(0, \ldots, 0,1)^{\mathrm{T}}$.

## Parameterization of the $n$-sphere

UNIVERSITEIT
VAN

How to parameterize unit vectors x ?
Compute Householder matrix (reflection) such that $\mathrm{H}_{\mathbf{v}(\mathbf{x})} \mathbf{x}=(0, \ldots, 0,1)^{\mathrm{T}}$.
(i) $f(\mathbf{y})=\hat{\mathbf{y}} /\|\hat{\mathbf{y}}\|$ where $\hat{\mathbf{y}}=\left(\mathbf{y}^{T}, 1\right)^{T}$, (?)
(ii) $f(\mathbf{y})=\left(\operatorname{sinc}(\|\mathbf{y}\| / 2) \mathbf{y}^{T}, \cos (\|\mathbf{y}\| / 2)\right)^{T}$ (?). both have a Jacobian $\partial f / \partial \mathbf{y}=[\mathrm{I} \mid \mathbf{0}]^{T}$.

## Parameterization of the $n$-sphere

How to parameterize unit vectors x ?
Compute Householder matrix (reflection) such that $H_{\mathbf{v}(\mathbf{x})} \mathbf{x}=(0, \ldots, 0,1)^{\mathrm{T}}$.
(i) $f(\mathbf{y})=\hat{\mathbf{y}} /\|\hat{\mathbf{y}}\|$ where $\hat{\mathbf{y}}=\left(\mathbf{y}^{T}, 1\right)^{T}$, (?)
(ii) $f(\mathbf{y})=\left(\operatorname{sinc}(\|\mathbf{y}\| / 2) \mathbf{y}^{T}, \cos (\|\mathbf{y}\| / 2)\right)^{T}$ (?).
both have a Jacobian $\partial f / \partial \mathbf{y}=[\mathrm{I} \mid \mathbf{0}]^{T}$.
So 'constrained' Jacobian can be computed

$$
\mathrm{J}=\frac{\partial C}{\partial \mathbf{y}}=\frac{\partial C}{\partial \mathbf{x}} \frac{\partial \mathbf{x}}{\partial \mathbf{y}}=\frac{\partial C}{\partial \mathbf{x}} \mathrm{H}_{\mathbf{v}(\mathbf{x})} \mathbf{x}[\mathrm{I} \mid \mathbf{0}]^{T} .
$$

## Zhang Paper

UNIVERSITEIT VAN

## Zhengyou Zhang

A Flexible New Technique for Camera Calibration (1998)

## Zhang Paper

# Zhengyou Zhang <br> A Flexible New Technique for Camera Calibration (1998) 

As implemented for:
Matlab The Camera Calibration Toolbox for Matlab
C++ Intel OpenCV

## Internal Camera Calibration 1/4

UNIVERSITEIT
VAN

Primary use of the Zhang algorithm is internal camera calibration. It computes:

- focal center $c_{x}$ and $c_{y}$.
- focal length $f_{x}$ and $f_{y}$.
- skew $s$ (optional).


## Internal Camera Calibration 1/4

Primary use of the Zhang algorithm is internal camera calibration. It computes:

- focal center $c_{x}$ and $c_{y}$.
- focal length $f_{x}$ and $f_{y}$.
- skew $s$ (optional).

In short, the camera intrinsic matrix:

$$
\mathbf{A}=\left[\begin{array}{ccc}
f_{x} & s & c_{x} \\
0 & f_{y} & c_{y} \\
0 & 0 & 1
\end{array}\right]
$$

## Internal Camera Calibration 2/4

The Zhang algorithm also computes radial lens distortion parameters $\left[k_{1}, k_{2}, k_{3}, k_{4}\right]$.

The original paper uses

$$
\begin{aligned}
& x_{d}=x+x\left(k_{1}\left(x^{2}+y^{2}\right)+k_{2}\left(x^{2}+y^{2}\right)^{2}\right), \\
& y_{d}=y+y\left(k_{1}\left(x^{2}+y^{2}\right)+k_{2}\left(x^{2}+y^{2}\right)^{2}\right),
\end{aligned}
$$

where $x$ and $y$ are normalized image coordinates and $x_{d}$ and $y_{d}$ are the distorted coordinates.

## Internal Camera Calibration 2/4

The Zhang algorithm also computes radial lens distortion parameters $\left[k_{1}, k_{2}, k_{3}, k_{4}\right]$.

The original paper uses

$$
\begin{aligned}
& x_{d}=x+x\left(k_{1}\left(x^{2}+y^{2}\right)+k_{2}\left(x^{2}+y^{2}\right)^{2}\right), \\
& y_{d}=y+y\left(k_{1}\left(x^{2}+y^{2}\right)+k_{2}\left(x^{2}+y^{2}\right)^{2}\right),
\end{aligned}
$$

where $x$ and $y$ are normalized image coordinates and $x_{d}$ and $y_{d}$ are the distorted coordinates.

But the implementations use a more complex model

$$
\begin{aligned}
& x_{d}=x+x\left(k_{1}\left(x^{2}+y^{2}\right)+k_{2}\left(x^{2}+y^{2}\right)^{2}\right)+x_{t d} \\
& y_{d}=y+y\left(k_{1}\left(x^{2}+y^{2}\right)+k_{2}\left(x^{2}+y^{2}\right)^{2}\right)+y_{t d}
\end{aligned}
$$

where

$$
\begin{aligned}
& x_{t d}=2 k_{3} x y+k_{4}\left(3 x^{2}+y^{2}\right), \\
& y_{t d}=2 k_{4} x y+k_{3}\left(x^{2}+3 y^{2}\right) .
\end{aligned}
$$

## Internal Camera Calibration 3/4

Example of internal camera calibration parameters.
Camera: PixeLINK A741, 2/3 inch CMOS sensor, 1280x1024.
Lens: Cosmicar 8.5 mm fixed focal length.
$f_{x}=1272.872$ pixels $=8.528 \mathrm{~mm}$
$f_{y}=1272.988$ pixels $=8.529 \mathrm{~mm}$
$c_{x}=632.740$
$c_{y}=507.648$
$k_{1}=-0.204$
$k_{2}=0.171$
$k_{3}=-0.00074896$
$k_{4}=0.00008878$

## Internal Camera Calibration 4/4

UNIVERSITEIT VAN
AMSTERDAM

Show lens distortion in DASiS video viewer. . .


## External Camera Calibration

The Zhang algorithm may also be used for external camera calibration.

Camera rotation and translation are computed as side-product of internal calibration.

If two cameras see the same calibration pattern at the same time, their relative position and orientation may be computed.

## Overall approach

UNIVERSITEIT
VAN
AMSTERDAM

- Measure projected position of points in a plane (e.g., checkerboard).


## Overall approach

UNIVERSITEIT
VAN
AMSTERDAM

- Measure projected position of points in a plane (e.g., checkerboard).
- Do so for at least two different camera orientations.


## Overall approach

- Measure projected position of points in a plane (e.g., checkerboard).
- Do so for at least two different camera orientations.
- Setup equations in order to estimate camera intrinsics.


## Overall approach

- Measure projected position of points in a plane (e.g., checkerboard).
- Do so for at least two different camera orientations.
- Setup equations in order to estimate camera intrinsics.
- Given camera intrinsics, estimate extrinsics.


## Overall approach

- Measure projected position of points in a plane (e.g., checkerboard).
- Do so for at least two different camera orientations.
- Setup equations in order to estimate camera intrinsics.
- Given camera intrinsics, estimate extrinsics.
- Estimate radial distortion.


## Overall approach

- Measure projected position of points in a plane (e.g., checkerboard).
- Do so for at least two different camera orientations.
- Setup equations in order to estimate camera intrinsics.
- Given camera intrinsics, estimate extrinsics.
- Estimate radial distortion.
- Use Levenberg-Marquardt to optimize initial estimates.


## Basic Equations

Universiteit
VAN
Amsterdam

Plane ('checkerboard') is at $Z=0$.

## Basic Equations

Universiteit
VAN
AMSTERDAM

Plane ('checkerboard') is at $Z=0$.
Homogeneous 2-D image point: $\widetilde{\mathbf{m}}$.
Homogeneous 3-D world point: $\widetilde{\mathrm{M}}=\left[\begin{array}{llll}X & Y & 0 & 1\end{array}\right]^{T}$.

## Basic Equations

UNIVERSITEIT
VAN

Plane ('checkerboard') is at $Z=0$.
Homogeneous 2-D image point: $\widetilde{\mathbf{m}}$.
Homogeneous 3-D world point: $\widetilde{\mathrm{M}}=\left[\begin{array}{llll}X & Y & 0 & 1\end{array}\right]^{T}$.
Projection:

$$
\begin{gathered}
s \widetilde{\mathbf{m}}=\mathbf{A}\left[\begin{array}{ll}
\mathbf{R} & \mathbf{t}
\end{array}\right] \widetilde{\mathrm{M}}= \\
{\left[\begin{array}{lll}
\alpha & \gamma & u_{0} \\
0 & \beta & v_{0} \\
0 & 0 & 1
\end{array}\right]\left[\begin{array}{llll}
\mathbf{r}_{1} & \mathbf{r}_{2} & \mathbf{r}_{3} & \mathbf{t}
\end{array}\right]\left[\begin{array}{llll}
X & Y & 0 & 1
\end{array}\right]^{T}=} \\
\\
\mathbf{A}\left[\begin{array}{lll}
\mathbf{r}_{1} & \mathbf{r}_{2} & \mathbf{t}
\end{array}\right]\left[\begin{array}{lll}
X & Y & 1
\end{array}\right]^{T}
\end{gathered}
$$

## Homography, constraints

UNIVERSITEIT
VAN
AMSTERDAM

An homography $\mathbf{H}$ can be estimated between known points on the calibration object and the measured world points.
$\mathbf{H}=\left[\begin{array}{lll}\mathbf{h}_{1} & \mathbf{h}_{2} & \mathbf{h}_{3}\end{array}\right]=\lambda \mathbf{A}\left[\begin{array}{lll}\mathbf{r}_{1} & \mathbf{r}_{2} & \mathbf{t}\end{array}\right]$

## Homography, constraints

An homography $\mathbf{H}$ can be estimated between known points on the calibration object and the measured world points.
$\mathbf{H}=\left[\begin{array}{lll}\mathbf{h}_{1} & \mathbf{h}_{2} & \mathbf{h}_{3}\end{array}\right]=\lambda \mathbf{A}\left[\begin{array}{lll}\mathbf{r}_{1} & \mathbf{r}_{2} & \mathbf{t}\end{array}\right]$
We demand:
C1: $\quad \mathbf{r}_{1}^{T} \mathbf{r}_{2}=0$
( $\mathbf{r}_{1}, \mathbf{r}_{2}$ orthogonal),
C2: $\quad \mathbf{r}_{1}^{T} \mathbf{r}_{1}=\mathbf{r}_{2}^{T} \mathbf{r}_{2}$
( $\mathbf{r}_{1}, \mathbf{r}_{2}$ have same length).

## Homography, constraints

An homography $\mathbf{H}$ can be estimated between known points on the calibration object and the measured world points.
$\mathbf{H}=\left[\begin{array}{lll}\mathbf{h}_{1} & \mathbf{h}_{2} & \mathbf{h}_{3}\end{array}\right]=\lambda \mathbf{A}\left[\begin{array}{lll}\mathbf{r}_{1} & \mathbf{r}_{2} & \mathbf{t}\end{array}\right]$
We demand:
C1: $\quad \mathbf{r}_{1}^{T} \mathbf{r}_{2}=0$
( $\mathbf{r}_{1}, \mathbf{r}_{2}$ orthogonal),
C2: $\quad \mathbf{r}_{1}^{T} \mathbf{r}_{1}=\mathbf{r}_{2}^{T} \mathbf{r}_{2} \quad\left(\mathbf{r}_{1}, \mathbf{r}_{2}\right.$ have same length $)$.

We know:
$\mathbf{h}_{1}=\lambda \mathbf{A} \mathbf{r}_{1} \quad \rightarrow \quad \mathbf{r}_{1}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{1}$
$\mathbf{h}_{2}=\lambda \mathbf{A} \mathbf{r}_{2} \quad \rightarrow \quad \mathbf{r}_{2}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{2}$

## Homography, constraints

An homography $\mathbf{H}$ can be estimated between known points on the calibration object and the measured world points.
$\mathbf{H}=\left[\begin{array}{lll}\mathbf{h}_{1} & \mathbf{h}_{2} & \mathbf{h}_{3}\end{array}\right]=\lambda \mathbf{A}\left[\begin{array}{lll}\mathbf{r}_{1} & \mathbf{r}_{2} & \mathbf{t}\end{array}\right]$
We demand:
C1: $\quad \mathbf{r}_{1}^{T} \mathbf{r}_{2}=0$
( $\mathbf{r}_{1}, \mathbf{r}_{2}$ orthogonal),
C2: $\quad \mathbf{r}_{1}^{T} \mathbf{r}_{1}=\mathbf{r}_{2}^{T} \mathbf{r}_{2} \quad\left(\mathbf{r}_{1}, \mathbf{r}_{2}\right.$ have same length $)$.

We know:
$\mathbf{h}_{1}=\lambda \mathbf{A} \mathbf{r}_{1} \quad \rightarrow \quad \mathbf{r}_{1}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{1}$
$\mathbf{h}_{2}=\lambda \mathbf{A} \mathbf{r}_{2} \quad \rightarrow \quad \mathbf{r}_{2}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{2}$
So the constraints are:
C1: $\quad \mathbf{h}_{1}^{T} \mathbf{A}^{-T} \mathbf{A}^{-1} \mathbf{h}_{2}=0$,
C2: $\quad \mathbf{h}_{1}^{T} \mathbf{A}^{-T} \mathbf{A}^{-1} \mathbf{h}_{1}=\mathbf{h}_{2}^{T} \mathbf{A}^{-T} \mathbf{A}^{-1} \mathbf{h}_{2}$.

## Closed-form solution using constraints $\mathbf{1 / 4}$

UNIVERSITEIT
VAN

Using the constraints, we can first find $\mathbf{A}$, followed by $\mathbf{R}$ and $\mathbf{t}$. Let

$$
\begin{aligned}
& \mathbf{B}=\mathbf{A}^{-T} \mathbf{A}^{-1}=\left[\begin{array}{lll}
B_{11} & B_{12} & B_{13} \\
B_{12} & B_{22} & B_{23} \\
B_{13} & B_{23} & B_{33}
\end{array}\right] \\
& =\left[\begin{array}{ccc}
\frac{1}{\alpha^{2}} & -\frac{\gamma}{\alpha^{2} \beta} & \frac{v_{0} \gamma-u_{0} \beta}{\alpha^{2} \beta} \\
-\frac{\gamma}{\alpha^{2} \beta} & \frac{\gamma^{2}}{\alpha^{2} \beta^{2}}+\frac{1}{\beta^{2}} & -\frac{\gamma\left(v_{0} \gamma-u_{0} \beta\right)}{\alpha^{2} \beta^{2}}-\frac{v_{0}}{\beta^{2}} \\
\frac{v_{0} \gamma-u_{0} \beta}{\alpha^{2} \beta} & -\frac{\gamma\left(v_{0} \gamma-u_{0} \beta\right)}{\alpha^{2} \beta^{2}}-\frac{v_{0}}{\beta^{2}} & \frac{\left(v_{0} \gamma-u_{0} \beta\right)^{2}}{\alpha^{2} \beta^{2}}+\frac{v_{0}^{2}}{\beta^{2}}+1
\end{array}\right] .
\end{aligned}
$$

This allows to solve for $\alpha, \beta$, etc.

## Closed-form solution using constraints 2/4

If we reshuffle the six unique elements of B into a vector $\mathbf{b}=\left[B_{11}, B_{12}, B_{22}, B_{13}, B_{23}, B_{33}\right]$,
we can rewrite both constraints as
$\mathbf{h}_{i}^{T} \mathbf{B h}_{j}=\mathbf{v}_{i j}^{T} \mathbf{b}$,
where

$$
\begin{aligned}
\mathbf{v}_{i j}= & {\left[h_{i 1} h_{j 1}, h_{i 1} h_{j 2}+h_{i 2} h_{j 1}, h_{i 2} h_{j 2},\right.} \\
& \left.h_{i 3} h_{j 1}+h_{i 1} h_{j 3}, h_{i 3} h_{j 2}+h_{i 2} h_{j 3}, h_{i 3} h_{j 3}\right]^{T},
\end{aligned}
$$

ultimately resulting in

$$
\left[\begin{array}{c}
\mathbf{v}_{12}^{T} \\
\left(\mathbf{v}_{11}-\mathbf{v}_{22}\right)^{T}
\end{array}\right] \mathbf{b}=0
$$

## Closed-form solution using constraints 3/4

Next, stack all the equations from $n$ measurements (estimated homographies) of the plane ('checkerboard'):

$$
\mathbf{V b}=0
$$

where $\mathbf{V}$ is a $2 n \times 6$ matrix. Solve as usual using the $S V D$.

## Closed-form solution using constraints 4/4

UNIVERSITEIT
VAN
AMSTERDAM

Once $\mathbf{A}$ is known, we can obtain $\mathbf{r}_{1}, \mathbf{r}_{2}$ and $\mathbf{t}$ :

$$
\begin{aligned}
& \mathbf{r}_{1}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{1} \\
& \mathbf{r}_{2}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{2} \\
& \mathbf{t}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{3}
\end{aligned}
$$

## Closed-form solution using constraints 4/4

UNIVERSITEIT
VAN

Once $\mathbf{A}$ is known, we can obtain $\mathbf{r}_{1}, \mathbf{r}_{2}$ and t :

$$
\begin{aligned}
& \mathbf{r}_{1}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{1} \\
& \mathbf{r}_{2}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{2} \\
& \mathbf{t}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{3}
\end{aligned}
$$

Now Zhang says
$\mathbf{r}_{3}=\mathbf{r}_{1} \times \mathbf{r}_{2}$,
and use SVD to make matrix $\mathbf{R}$ orthogonal, i.e., $\mathbf{R}=\mathbf{U V}^{T}$.

## Closed-form solution using constraints 4/4

Once $\mathbf{A}$ is known, we can obtain $\mathbf{r}_{1}, \mathbf{r}_{2}$ and $\mathbf{t}$ :

$$
\begin{aligned}
& \mathbf{r}_{1}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{1}, \\
& \mathbf{r}_{2}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{2} \\
& \mathbf{t}=\lambda^{-1} \mathbf{A}^{-1} \mathbf{h}_{3}
\end{aligned}
$$

Now Zhang says
$\mathbf{r}_{3}=\mathbf{r}_{1} \times \mathbf{r}_{2}$,
and use SVD to make matrix $\mathbf{R}$ orthogonal, i.e.,
$\mathbf{R}=\mathbf{U V}^{T}$.
I say:
Make $\mathbf{r}_{1}, \mathbf{r}_{2}$ orthogonal in least-squares sense.
The compute $\mathbf{r}_{3}=\mathbf{r}_{1} \times \mathbf{r}_{2}$.
Is simpler and boils down to the same thing.

## Radial distortion

Using the camera intrinsics and extrinsics undistorted coordinates of points (corners on the checkerboard) can be approximated. These is used to solve for $k_{1}, k_{2}$ :

$$
\left[\begin{array}{cc}
\left(u-u_{0}\right)\left(x^{2}+y^{2}\right) & \left(u-u_{0}\right)\left(x^{2}+y^{2}\right)^{2} \\
\left(v-v_{0}\right)\left(x^{2}+y^{2}\right) & \left(v-v_{0}\right)\left(x^{2}+y^{2}\right)^{2}
\end{array}\right]\left[\begin{array}{l}
k_{1} \\
k_{2}
\end{array}\right]=\left[\begin{array}{l}
\breve{u}-u \\
\breve{v}-v
\end{array}\right] .
$$

## Radial distortion

Using the camera intrinsics and extrinsics undistorted coordinates of points (corners on the checkerboard) can be approximated. These is used to solve for $k_{1}, k_{2}$ :
$\left[\begin{array}{cc}\left(u-u_{0}\right)\left(x^{2}+y^{2}\right) & \left(u-u_{0}\right)\left(x^{2}+y^{2}\right)^{2} \\ \left(v-v_{0}\right)\left(x^{2}+y^{2}\right) & \left(v-v_{0}\right)\left(x^{2}+y^{2}\right)^{2}\end{array}\right]\left[\begin{array}{l}k_{1} \\ k_{2}\end{array}\right]=\left[\begin{array}{c}\breve{u}-u \\ \breve{v}-v\end{array}\right]$.
These equations are stacked $\left(\mathbf{D}\left[\begin{array}{ll}k_{1} & k_{2}\end{array}\right]^{T}=\mathbf{d}\right)$ and we solve least squares $\left[\begin{array}{ll}k_{1} & k_{2}\end{array}\right]^{T}=\left(\mathbf{D}^{T} \mathbf{D}\right)^{-1} \mathbf{D}^{T} \mathbf{d}$.

Then iterate both algorithm (internal+external, radial) until convergence.

## Maximum likelihood estimation

UNIVERSITEIT
VAN
Amsterdam

Optimize: use Levenberg-Marquardt to find minimum of
$\sum_{i=1}^{n} \sum_{j=1}^{m}\left\|\mathbf{m}_{i j}-\breve{\mathbf{m}}\left(\mathbf{A}, k_{1}, k_{2}, \mathbf{R}_{i}, \mathbf{t}_{i}, \mathrm{M}_{i}\right)\right\|^{2}$
( $n$ images, $m$ points per image)

All done . . .

## Notable experimental results

UNIVERSITEIT
VAN
AMSTERDAM

- Using three different images, the results are pretty good. Results keep getting better with more images.


## Notable experimental results

- Using three different images, the results are pretty good. Results keep getting better with more images.
- 45 degree angle between image plane and checkerboard seems to give best result. Loss of precision in corner detection was not taken into account (simulated data).


## Notable experimental results

- Using three different images, the results are pretty good. Results keep getting better with more images.
- 45 degree angle between image plane and checkerboard seems to give best result. Loss of precision in corner detection was not taken into account (simulated data).
- Systematic non-planarity of checkerboard has more effect than random noise (duh).


## Notable experimental results

- Using three different images, the results are pretty good. Results keep getting better with more images.
- 45 degree angle between image plane and checkerboard seems to give best result. Loss of precision in corner detection was not taken into account (simulated data).
- Systematic non-planarity of checkerboard has more effect than random noise (duh).
- Cylindrical non-planarity is worse than spherical non-planarity (cylindrical more common in practice?).


## Notable experimental results

- Using three different images, the results are pretty good. Results keep getting better with more images.
- 45 degree angle between image plane and checkerboard seems to give best result. Loss of precision in corner detection was not taken into account (simulated data).
- Systematic non-planarity of checkerboard has more effect than random noise (duh).
- Cylindrical non-planarity is worse than spherical non-planarity (cylindrical more common in practice?).
- Even with systematic non-planarity, results still usable.


## Notable experimental results

- Using three different images, the results are pretty good. Results keep getting better with more images.
- 45 degree angle between image plane and checkerboard seems to give best result. Loss of precision in corner detection was not taken into account (simulated data).
- Systematic non-planarity of checkerboard has more effect than random noise (duh).
- Cylindrical non-planarity is worse than spherical non-planarity (cylindrical more common in practice?).
- Even with systematic non-planarity, results still usable.
- Error in compute sensor center seems not to have too much effect in 3-D reconstruction.

