

Structureless global bundle adjustment in MicMac

Les grandes lignes

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1 Relative orientations **NO_AllOri2Im**

The code launches **TestLib NO_AllOri2Im** (cf. Alg. 1) for a pattern of images. Inside the program, it iterates over all images and launches **TestLib NO_Ori2Im** for all connected images (cf. Alg. 2 and Fig. 1). The main method that calculates the relative orientation between an image pair is detailed in Alg. 4.

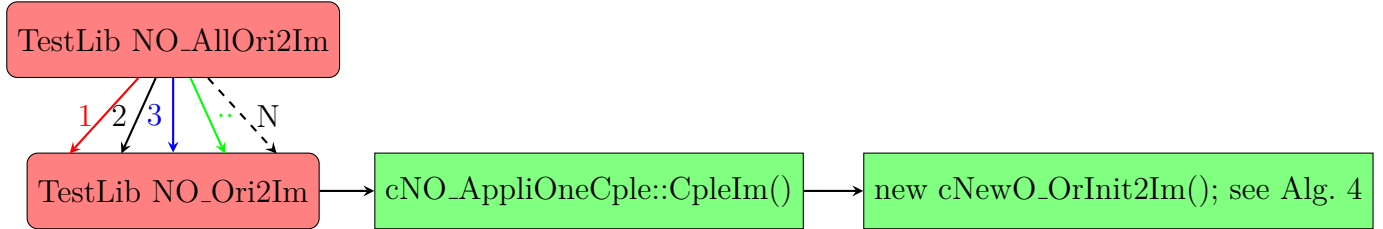


Figure 1: The relative orientation workflow. 1– N signify nodes (i.e. images) in the connectivity graph. Computation is done in parallel.

1.1 Main classes

- `cNO_AppliOneCple` – manager, prepares data for `cNewO_OrInit2Im`, contains the image pair and their names, multiple tie-pts structure, a directory/file manager ...
- `cNewO_OrInit2Im` – class managing the calculation of a relative orientation
- `cXml_O2IComputed` – saves the relative orientation result

1.2 Algorithms

Algorithm 1 **TestLib NO_AllOri2Im** def in `TestAllNewOriImage_main`

aVIm – vector of images in pattern
aHom – homologous points

```

                                ▷ Launch rel ori caculation for each image pair
for  $aK = 0; aK < aVIm \rightarrow size(); aK++$  do
    aVH – vector of overlapping images (with common tie-pts)
    for  $aKH = 0; aKH < aVH \rightarrow size(); aKH++$  do
        aIm1 = aVIm[aK]
        aIm2 = aVH[aKH]
        TestLib NO_Ori2Im aIm1 aIm2
    end for
end for
  
```

Algorithm 2 TestLib NO_Ori2Im def in TestNewOriImage_main

BenchNewFoncRot()

▷ Preparation of data, i.e. create multiple tie-pts structures, read orientations if InOri
cNO_AppliOneCple anAppli(argc, argv)

▷ Calculate relative orientation

*cNewO_OrInit2Im * aCple = anAppli.CpleIm();*

▷ Save result to xml

cXml_Ori2Im&aXml = aCple->XmlRes()

MakeFileXML(aXml, anAppli.NameXmlOri2Im(true))

MakeFileXML(aXml, anAppli.NameXmlOri2Im(false));

Algorithm 3 anAppli.CpleIm() def in NewOri/cNewO_CpleIm.cpp

▷ Initialise an object of class *cNewO_OrInit2Im*

return new *cNewO_OrInit2Im(mGenOri, mQuick, mIm1, mIm2, &mMergeStr, mTestSol, mRotInOri)*

Algorithm 4 *cNewO_OrInit2Im* Initial orientation of an image pair

Creates *PackReduit* which is a subset of 500 tie-pts chosen heuristically
Save photogrammetric points for the image pair, i.e. HomFloatSym.dat
Initialise the *cXml_Ori2Im* for the image pair

if *GpsIsInit* **then**

 Do something

end if

Calculate a homography between 2d points with *cElHomographie :: RobustInit*

Save results (including residuals, overlap hom, 2d ellipses) to *cXml_O2IComputed* class,
object *aXCmp*

▷ Test if the scene is not locally planar

ElRotation3D * aRP = TestOriPlanePatch()

if aRP **then**

 AmeliorerSolLinear(*aRP,)

end if

▷ Test the essential matrix with minimum pts and RANSAC (8pts algo)

ElRotation3D aMRR = TestcRanscMinimMatEss()

AmeliorerSolLinear(aMRR,);

▷ Test the essential matrix with more pts and RANSAC (8pts algo)

ElRotation3D aRR = RansacMatriceEssentielle()

▷ Test the essential matrix - classical test, L1+L2, no RANSAC

ElRotation3D aR = (aL2 ? mPackPStd.MepRelPhysStd(1.0,true) : mPackStdRed.MepRelPhysStd(1.0,false)) ;

▷ Test global homography, robust estimation

cResMepRelCoplan aRMC = ElPackHomologue::MepRelCoplan(...

...

Save the new homography in aXCmp.HomWithR()

▷ Test pure rotation of the camera

cResMepCoc aRCoc= MEPCoCentrik(...

...

Save the rotation in aXCmp.RPure().Ori()

▷ Adjust the solution in a few iterations (until now direct linear solution)

ElRotation3D aSol = aBundle→OneIterEq(...

▷ Triangulate tie-pts, calculate 3D ellipses, and save

... aXCmp.Elips() = anElips3D

Keep track of computational time with *cXml_O2ITiming* & *aTiming*

Algorithm 5 *cElRotation3D * TestOriPlanePatch* def in *photogramphgr_mep_patch_plane.cpp*
to do

Algorithm	6	cElRotation3D	*	TestcRanscMinimMatEss	def	in
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photogramphgr_mep_patch_plane.cpp

to do

Algorithm	7	cElRotation3D	*	RansacMatriceEssentielle	def	in
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photogramphgr_mep_patch_plane.cpp

to do

Algorithm 8	cElRotation3D	*	MEPCoCentrik	def in	<i>photogramphgr_mep_patch_plane.cpp</i>
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to do

2 Generate point hom per triplet **NO_AllImTriplet**

The goal of this program is to gather information about tie-points of multiplicity 3, i.e. tie-points with track length 3 and thus visible in all images of a triplet. The program *TestLib NO_AllImTriplet* launches the *CPP_GenAllImP3()* function the defined in *cNewO_PointsTriples.cpp*. Inside the *CPP_GenAllImP3()*, for each image within the called pattern the following programme is executed in parallel *TestLib NO_OneImTriplet*.

Then, the *TestLib NO_OneImTriplet*, calls the *CPP_GenOneImP3()* function (see Alg. 9) defined in *cNewO_PointsTriples.cpp*. Inside this function, it creates an object of class *cAppli_GenPTripleOneImage*, and launches this class' member function – *GenerateTriplets()* (cf. Fig. 10).

2.1 Main classes

- *cAppli_GenPTripleOneImage*

2.2 Algorithms

Algorithm 9 *CPP_GenOneImP3* def in *cNewO_PointsTriples.cpp*

▷ Create object of class *cAppli_GenPTripleOneImage* (
in constructor initialise *mVCams* which contains all images connected to the current image;
the first element of *mVCam* is the current/master image
cAppli_GenPTripleOneImage anAppli(argc,argv);
anAppli.GenerateTriplets();

▷ Launch triplet point generation, see Alg. 10

Algorithm 10 *cAppli_GenPTripleOneImage* :: *GenerateTriplets()* def in *cNewO_PointsTriples.cpp*

mVCams – vector of all images
▷ For every two images connected to the current image (i.e. *mVCams*[0]), calculate the triplet points
for *aK* = 0; *aK* < *mVCams* → *size()*; *aK* ++ **do**
 for *aK* = 0; *aK* < *mVCams* → *size()*; *aK* ++ **do**
 GenerateTriplet(*aKC1*,*aKC2*);
 end for
end for

Algorithm 11 *cAppli_GenPTripleOneImage* :: *GenerateTriplet(in,int)* def in *cNewO_PointsTriples.cpp*

▷ Load all tie-pts for an image pair
▷ Collect tie-pts with multiplicity 3
▷ Save the triplet points to a file
mNM – > *WriteTriplet(aName3, aVP1Exp, aVP2Exp, aVP3Exp, aVNb)*;

3 Generate triplets **NO_GenTripl**

The goal here is to build a set of optimal triplet pairs from the existing relative orientations. As is illustrated schematically in Fig. 2 the **TestLib NO_GenTripl** launches **GenTriplet_main** (see Alg. 12) for a pattern of images. Then, for each edge of the epipolar graph, the Alg. 15 is called.

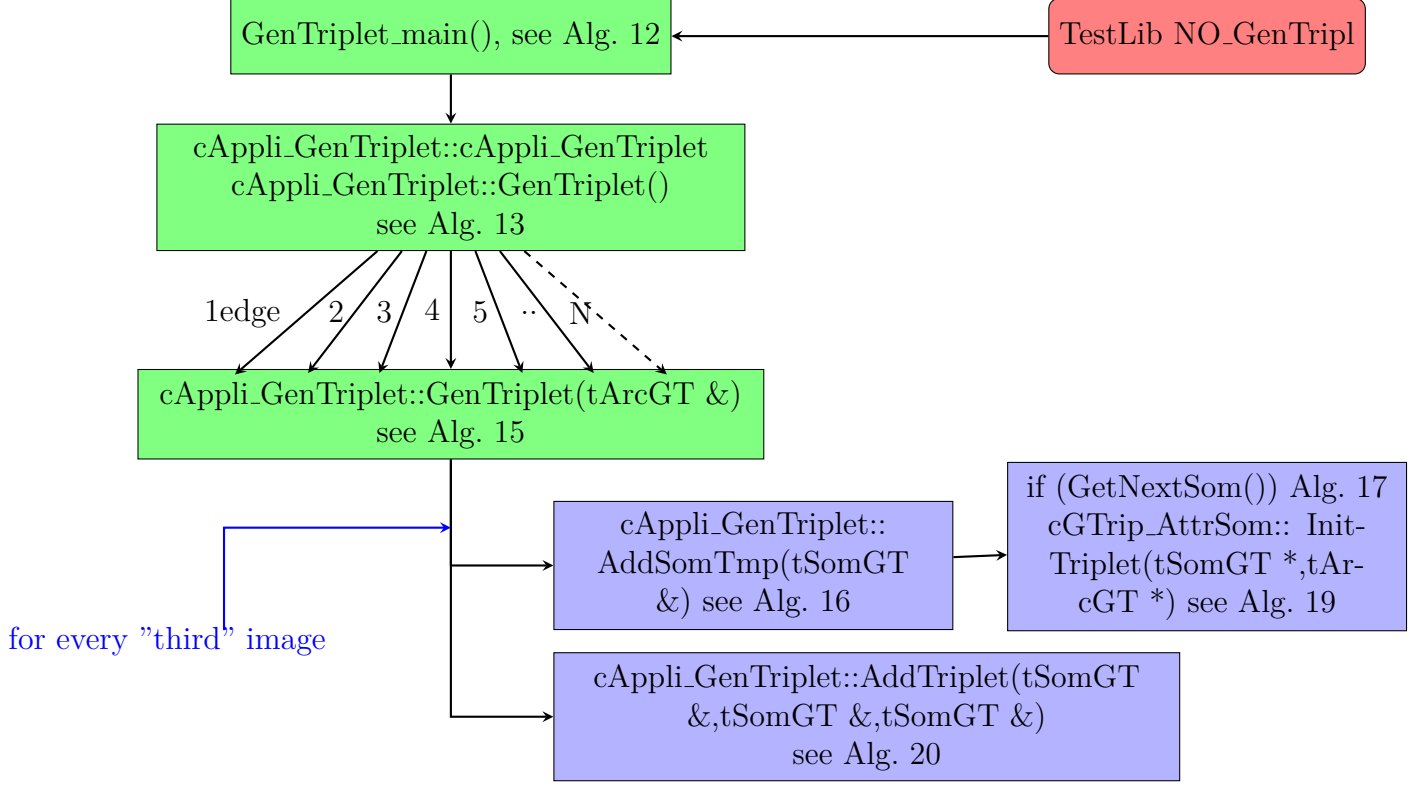


Figure 2: The triplet generation workflow. The enumerated arrows launch the **GenTriplet** method for each edge (i.e. a pair of nodes and the edge connecting them) of the epipolar graph. The methods in purple are launched for each "third" image attached to each edge of the graph.

3.1 Main classes, some typedefs and variables

- *cAppli_GenTriplet*
- *cGTrip_AttrSom* – class corresponding to a triplet node (its orientation parameters etc)
- *cGTrip_AttrASym* – ?
- *cGTrip_AttrArc* – class containing relative rotation corresponding to an edge
- *cResTriplet* – class containing *cXml_Ori3ImInit* (xml save)
- *cTripletInt* – typedef to *cTplTriplet < int >* containing ids of images within a triplet
- *tSomGT* – a node of the graph
- *tArcGT* – an edge of the graph
- *tGrGT* – a graph
- * *mHautBase* – height, used e.g. in the B to H ratio calcul

- * *mMapTriplets* – contains all triplets (for each triplet it stores the identifiers of nodes defining the node and the resulting orientations)
- * *mTopoTriplets* – contains a list of all accepted triplets (xml of type *XmlTopoTriplet*)
- * *mTriOfCple* – stores edges (i.e. 2 images) and a list of "third" images for each edge
- * *mVSomVois* – a vector of nodes that are neighbouring with an edge in consideration (i.e., they're outside that edge)
- * *mVSomEnCourse* – vector of "third" images that are currently connected to an edge
- * *mVSomSelected* – vector containing the selected triplets

3.2 Algorithms

$$|C_2 + \lambda \cdot C_3 \quad U_3 \quad U_1| = 0 \quad (1)$$

$$\lambda = -\frac{|C_2 \quad U_3 \quad U_1|}{|C_3 \quad U_3 \quad U_1|} \quad (2)$$

Figure 3: Computation of orientations of a triplet consistent with relative orientations. See also Alg. 19

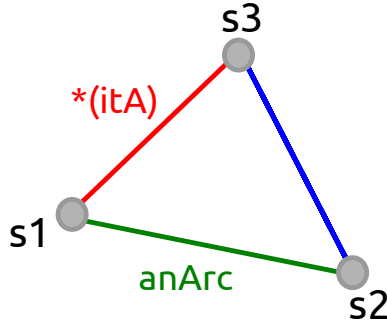


Figure 4: A triplet defined by 3 nodes and three edges, illustration of the Alg. 15. The green edge is the edge entering the *GenTriplet(tArcGT&anArc)* method. If s3 exists, there is an edge between s1 and s3. If (*mGrT.edge_s1s2(anArc.s2(), aS3)*), the blue edge connecting s2 and s3 exists, and therefore a complete triplet is defined.

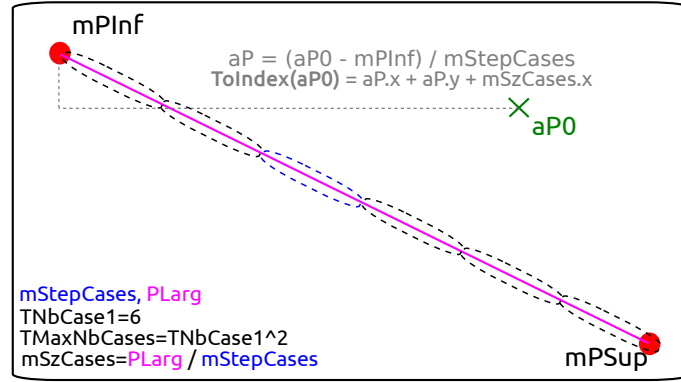


Figure 5: Illustration of the tie-pts indexing method implemented in $ToIndex(const Pt2df \&aP0)$.

Algorithm 12 *GenTriplet_main* def in cNewO_OldGenTriplets.cpp

cAppli_GenTriplet anAppli(argc,argv);

▷ Initialise the object class

anAppli.GenTriplet();

▷ Call the GenTriplet()

Algorithm 13 *cAppli_GenTriplet* :: *cAppli_GenTriplet* constructor def in cNewO_OldGenTriplets.cpp

▷ Some member variables

mVecAllSom - a vector of all nodes

mMapS - a map of graph nodes and the respective image names

mGrT - the epipolar graph

tSubGrGT SubAll allows to iterate over a subgraph of edges attached to a node

▷ Initialise the nodes of the graph

mVecAllSom, mMapS

▷ Initialise the edges of the graph

for Each stereo pair **do**

 Get respective nodes from the map of nodes (aS1,aS2)

 Get the relative orientation from XML file

 Test the edge from aS1 to aS2 ?

 Add the edge to the graph mGrT

end for

Algorithm 14 *cAppli_GenTriplet* :: *GenTriplet()* def in cNewO_OldGenTriplets.cpp

▷ Iterate over edges

for All edges between nodes in mVecAllSom **do**

 Call GenTriplet(*itA);

end for

Algorithm 15 *cAppli_GenTriplet* :: *GenTriplet(tArcGT&anArc)* def in
cNewO_OldGenTriplets.cpp

▷ Only symmetric edges are dealt with

if (!anArc.attr().IsDirASym()) return;

▷ Load tie-pts for the stereo pair

▷ Get the distance between the two most distant points (mPInf,mPSup) in the pair model;
return if it is too large

▷ Some variables

mCurPMed – median point calculated on all tie-pts of a pair
mCurS1 = & (anArc.s1()); – the aS1 in Fig. 4
mCurS2 = & (anArc.s2()); – the aS2 in Fig. 4

▷ Initialise the triplets. See Fig. 4 for illustration

for all edges attached to mCurS1/aS1 **do**
 Get the second node (**itA*).s2() for an edge in consideration and call it aS3 of the triplet
 if there is an edge between aS3 and aS2 of the anArc **then**
 call *AddSomTmp(aS3)*, see Alg. 16
 end if
end for

▷ Iterate over all connected images that are found in mVSomEnCourse and
select the "best" ones. The number of selected triplets is defined by *TQuickNbMaxTriplet*
and *TStdNbMaxTriplet*. See Alg. 17

while aSom = GetNextSom() **do**
 *AddTriplet(*aSom, mCurArc- > s1(), mCurArc- > s2());* see Alg. 20
end while

Algorithm 16 *cAppli_GenTriplet* :: *AddSomTmp(tSomGT&aS)* def in
cNewO_OldGenTriplets.cpp

mVSomVois.push_back(&aS);

▷ If an approx orientation of triplet is possible, add the "third" image to the current nodes

if *thenInitTriplet* (see Alg. 19)
 mVSomEnCourse.push_back(&aS);
end if

Algorithm 17 *cAppli_GenTriplet :: GetNextSom()*, update the mVSomSelected; def in cNewO_OldGenTriplets.cpp

▷ Leave if your objective is reached, i.e. there are no "third" images in the current nodes or if the number of selected "third" images is already bigger than the *aNbMaxTriplet*

```

if(mVSomEnCourse.empty())return0;
if(int(mVSomSelected.size()) > aNbMaxTriplet)return0;

    ▷ Get the best GainGlob for all available "third" images and store it in:
aGainMax, aRes, aIndexRes

    ▷ Remove the best image from mVSomEnCourse
mVSomEnCourse.erase(mVSomEnCourse.begin() + aIndexRes);

    ▷ And save it in the selected images
mVSomSelected.push_back(aRes);

    ▷ Update the cost/gain of the remaining images given the best result in aRes
for all images in mVSomEnCourse do
    tSomGT * aSom = mVSomEnCourse[aK];
    aSom->attr().UpdateCost(aSom, aRes); see Alg. 18
end for

```

Algorithm 18 *cGTrip_AttrSom :: UpdateCost(tSomGT * aSomThis, tSomGT * aSomSel)*, def in cNewO_OldGenTriplets.cpp

▷ Calculates the new gain as a function of b sur h between aSom and the aSomSel being the best node

```

aNewGain = mAppli->GainBSurH(aSomThis, aSomSel);

    ▷ Get Pds and Dens that will help to define the new global gain/cost
int * aPdsGlob = mAppli->Pds();
int * aDens2 = aSomSel->attr().mDens;

    ▷ Re-define the gain accordingly:

for All "cases" (here 6*6) do
    int aD2 = aDens2[aK];
    ElSetMin(mGain[aK], mDens[aK] * ( aNewGain * aD2 + TQuant*(TQuant-aD2)));
    mGainGlob += mGain[aK] * aPdsGlob[aK];
end for

```

Algorithm 19 *cGTrip_AttrSom* :: *InitTriplet*(*tSomGT* * *aSom*, *tArcGT* * *anA12*) def in *cNewO_OldGenTriplets.cpp*

▷ Some variables

anA13 – edge from *s1* to *s3* (see Fig. 4)
anA23 – edge from *s2* to *s3* (see Fig. 4)
aR21 – affine rotation (Rot et tr) corresponding to edge between *s1* and *s2*
aR31 – affine rotation (Rot et tr) corresponding to edge between *s1* and *s3*
aR31Bis – affine rotation (Rot et tr) corresponding to edge between *s1* and *s3* calculated from *aR21* and *aR32*
aR32 – affine rotation (Rot et tr) corresponding to edge between *s2* and *s3*
aVP1, aVP2, aVP3 – vectors containing tie-pts visible in the triplet
mNb – vector that contains information about tie-pts in an indexed form

▷ Calculate orientations consistent within a triplet (origin in *s1*) using the known relative orientations. We search for a scale factor λ for which the three directions stemming from three images (i.e. nodes) and corresponding to a tie-point will intersect in 3D. We use the coplanarity condition as the mathematical model to calculate it. Its determinant shall be null for the solution to exist. Because we will have as many λ s as there is tie-points, the final result is the median of all results. Since the multiple points are not really considered in this approach, it will be ill-defined for three colinear perspective center.

for all almost tie-pts in *aVP1* (*aStep* defines the subsample) **do**

▷ Get normalised vector directions

aU31 – direction corresponding to tie-pt in *s3* calculated from *R31*
aU2 – direction corresponding to tie-pt in *s2*
aU1 – direction corresponding to tie-pt in *s1*
aU31Bis – direction corresponding to tie-pt in *s3* calculated from *R31Bis*

aVL13, aVL23 are vectors containing all the results calculated using two combinations

▷ Calculate intersection of the points using *s1, s2*

aPInt12 = *InterSeg*(*aC1, aC1+aU1, aC2, aC2+aU2, OkI*);

▷ Impose that *s3* is found at the intersection of line *s1 s3* and on the bundle stemming from *aPInt12* and going towards *U31*. Save result to *aVLByInt*

CoordInterSeg(*aC1, aC3, aPInt12, aPInt12+aU31, OkI, p, q*);

▷ Calculate the perspective center of *s3* from *aVL13* and *aVL23* and *aVLByInt*

Pt3dr *aC31* = *aC3* / *MedianeSup*(*aVL13*);
Pt3dr *aC32* = *aC2* + *aV3Bis* * *MedianeSup*(*aVL23*);
mC3 = (*aC31* + *aC32*) / 2.0;
Pt3dr *aC3I* = *aC3* * *MedianeSup*(*aVLByInt*);

▷ The final solution for the perspective center is the one from multiple points (i.e. from the inverse *aVLByInt*). As far as the rotation, the average result from *R31* et *R31Bis* is taken (orthonormality is imposed with nearest rotation.)

mC3 = *aC3I*;
mM3 = *NearestRotation*((*aR31.Mat()* + *aR31Bis.Mat()*)*0.5);

end for

▷ Calculate gain and density

$aGain1$, b sur h between $s1$ and $s3$
 $aGain2$, b sur h between $s2$ and $s3$
 $aGain = ElMin(aGain1, aGain2);$

▷ Inside this method, the mNb vector is updated. The size of the vector is equal to $TMaxNbCases$ and it encodes indexed position in an image. To update the mNb , each tie-pt is first indexed with `ToIndex()` (see Fig. 3.2). The output of the indexing is an int from 0 to $TMaxNbCases$ and it indicates the position within the mNb vector that will be incremented.

$InitNb(aVP1);$

▷ Using the mNb we now calculate gain scores for each *case element* of the indexed image. The gain depends on points' density in a given *case element*

for all element cases i.e. $mSzCases.x * mSzCases.y$, see Fig. 3.2 **do**

$mDens[aK]$ – the density per case

$mGain[aK]$ – gain per case

$mGainGlob$ – global gain

end for

Algorithm 20 *cAppli.GenTriplet* :: *AddTriplet*($tSomGT\&, tSomGT\&, tSomGT\&$) def in *cNewO_OldGenTriplets.cpp*

▷ Get the corresponding nodes

$aA1, aA2, aA3$

▷ Normalise the orientations so that the first node has an identity rotation matrix and the bases between the first and the second node is equal to 1. To do that, multiply orientations of nodes $aA2$ and $aA3$ by the inverse of $aA1$; then, divide the orientations by the length of the base between $aA1$ and $aA2$.

$aR1Inv = aA1.R3().inv();$

$aR1 = aR1Inv * aA1.R3();$

$aR2 = aR1Inv * aA2.R3();$

$aR3 = aR1Inv * aA3.R3();$

$double aD = euclid(aR2.tr());$

$aR2 = ElRotation3D(aR2.tr()/aD, aR2.Mat(), true);$

$aR3 = ElRotation3D(aR3.tr()/aD, aR3.Mat(), true);$

▷ Iterate over all (triplet) tie-pts. For each tie-pt in each image get its direction with *AddSegOfRot*. The directions are stored in vectors aW . Intersect the directions in 3D with *InterSeg*. Get the residuals in images, store the mean value in $aVRes$ vector.

▷ Check if this triplet already exists in the *mMapTriplets*. If it does and the median residual of the current triplet is larger than that of the triplet already stored in the map, it returns from the *AddTriplet* method. Otherwise it continues.

▷ Save. Update the *mMapTriplets*, the *mTopoTriplets* and the *mTriOfCple*.

4 Optimisation I NO_AllImOptTrip

The goal of this program is to optimise the calculated triplet orientations by (i) filtering tie-pts, (ii) testing with *TestOPA()* algorithm based on resection and *TestTomasiKanade()* algorithm suited for long focal lengths. Finally, bundle adjustment *SolveBundle3Image* is called for each triplet. Schematic workflow of the algorithms is presented in Fig. 6.

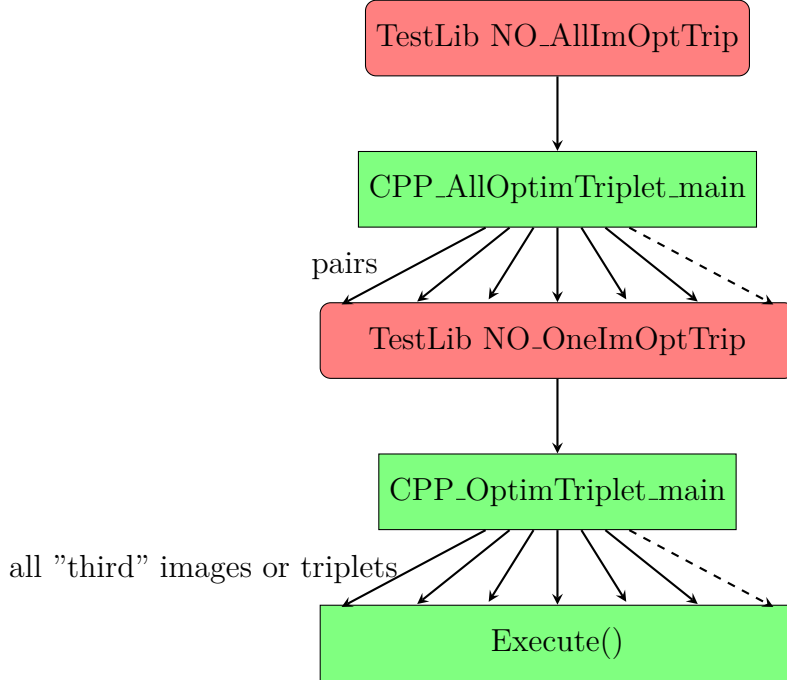


Figure 6: Triplet optimisation I workflow.

4.1 Main classes, some typedefs and variables

- *cAppliOptimTriplet* – manager

4.2 Algorithms

Algorithm 21 *CPP_AllOptimTriplet_main* constructor def in cNewO_OptimTriplet.cpp

- ▷ Iterate over all couples in *ListeCpleOfTriplets.xml* (must be coherent with pattern too!) and launch the *TestLibNO_OneImOptTrip* programme for each.
-

Algorithm 22 *CPP_OptimTriplet_main* constructor def in cNewO_OptimTriplet.cpp

- ▷ Call the object of the "managing" class
- ```
cAppliOptimTriplet anAppli(argc,argv,true);
```
-

---

**Algorithm 23** *cAppliOptimTriplet* :: *cAppliOptimTriplet* constructor def in  
cNewO\_OptimTriplet.cpp

---

▷ Check whether testing with Tomas-Kanade algo shall be done (specific to long focal  
lengths)

*mModeTTK*, *mWithTTK*

▷ For a pair of images, get a list of its "third" images forming a triplet with that couple.  
Launch:

*Execute()*

---

---

**Algorithm 24** *cAppliOptimTriplet :: Execute()* constructor def in cNewO\_OptimTriplet.cpp

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▷ Define the number of selected points. If Quick mode is on (true by default) 200 points are used, otherwise its 500 points.

QuickDefNbMaxSel=200, StdDefNbMaxSel=500;  
mNbMaxSel = mQuick ? QuickDefNbMaxSel : StdDefNbMaxSel;

mNbMaxInit = mQuick ? QuickNbMaxInit : StdNbMaxInit ;

▷ Calculate a mean focal length *mFoc*

▷ Load images to the vector *mIm*s. Create variables *mIm*1, *mIm*2, *mIm*3 for each image of the triplet. Load pairs of images to vector *mPairs*. Create variables *mP*12, *mP*13, *mP*23 for each couple of the triplet.

▷ Load tie-pts with the *VFullPtOf3()* method.

▷ "Attenuate" the number of selections with  $\frac{a \cdot b}{a+b}$  as shown below:

int aNbFull3 = *mIm*2- > *VFullPtOf3()*.size() ;  
int aNb3 = round\_up(CoutAttenuateTetaMax(aNbFull3,mNbMaxSel));  
mNbMaxSel = aNb3;

▷ Filter tie-pts according to set parameters

*mSel*3 = *IndPackReduit*(*mIm*2- > *VFullPtOf3()*, *mNbMaxInit*, *mNbMaxSel*);

▷ Filter tie-pts for Tomas-Kanade tests

*mTKNbMaxSel* = *ElMin*(*mTKNbMaxSel*, *aNbFull3*);  
*mTKSel*3 = *IndPackReduit*(*mIm*2- > *VFullPtOf3()*, *ElMin*(*aNbFull3*, 5 \*  
*mTKNbMaxSel*), *ElMin*(*aNbFull3*, *mTKNbMaxSel*));

▷ Update some variables for all triplet images

*mIm*s[*aK*]- > *SetReduce*(*mSel*3, *mTKSel*3);  
*mFullH*123.push\_back(&(mIm[s[*aK*]- > *VFullPtOf3()*));  
*mRedH*123.push\_back(&(mIm[s[*aK*]- > *VRedPtOf3()*));  
*mTK\_H*123.push\_back(&(mIm[s[*aK*]- > *VT\_K\_PtOf3()*));

▷ Calculate a global residual which is the mean residual of residuals calculated on pairs 12, 13 and 23

mBestResidu = ResiduGlob();

▷ TestOPA

**for** (int aKP=0 ; aKP<int(mPairs.size()) ; aKP++) **do**  
    TestOPA(\*(mPairs[aKP])); see Alg. 25  
**end for**

▷ TestTomasiKanade

TestTomasiKanade(); see Alg. 27

▷ Perform bundle adjustemnt on the result

*SolveBundle3Image*(*mFoc*, *aR*21, //*mIm*2- > *Ori*(), *aR*31, //*mIm*3- >  
*Ori*(), *aP*Med, *aB*OnH, *mRedH*123, *mP*12- > *RedHoms*(), *mP*13- >  
*RedHoms*(), *mP*23- > *RedHoms*(), *mP*ds3, *aParamSB3I*);

▷ Calculate and save the "ellipse"

*cXml\_Elips3D anElips3D*;  
*aXml.Elips*() = *anElips3D*; 16  
*MakeFileXML*(*aXml*, *aNameSauveXml*);

---



---

**Algorithm 25** *cAppliOptimTriplet :: TestOPA(cPairOfTriplet&aPair)* constructor def in cNewO\_OptimTriplet.cpp

---

▷ Load filtered tie-pts mRedH123 into aVP and images to aI.

**for** All tie-pts in aVP **do**

    Get direction of the point in aI1 and aI2 and store in aVA,aVB

        ▷ Intersect the two directions

*Pt3dranInter = InterSeg(aVA, aVB, OkI);*

    ▷ Store in aL32 the correspondence 3D - 2D for the "third" images aI3. Note that the 2D position is not a projection of the anInter.

*aL32.push\_back(Appar23(Pt2dr(aP3.x, aP3.y), anInter));*

**end for**

        ▷ Estimate orientation of aI3 from RANSAC-based resection

*ElRotation3DaR3 = aCSI.RansacOFPA(true, aNbRansac, aL32, &anEcart);*

    ▷ Normalise the orientations so that the base between image1 and image2 is equal to 1 and the rotation matrix of image1 is identity (same as in Alg. 20)

    ▷ Test the solution by comparing residuals of the new solution with the former residuals.  
    TestSol(aVR); see Alg. 26

---



---

**Algorithm 26** *cAppliOptimTriplet :: TestSol(const std :: vector < ElRotation3D > &aVR);* def in cNewO\_OptimTriplet.cpp

---

        ▷ Compute the residuals based on the aVR solution

*aResidu = ResiduGlob(aVR[0], aVR[1], aVR[2]);*

    ▷ If the current residual is smaller than the former best residual, update the orientations and the best residual

**if** *then* *aResidu < mBestResidu*

*mBestResidu = aResidu;*

**for** *do* *intaK = 0; aK < 3; aK ++*

*mIms[aK] -> SetOri(aVR[aK]);*

**end for**

**end if**

---



---

**Algorithm 27** *cAppliOptimTriplet :: TestTomasikanade();* def in cNewO\_OptimTriplet.cpp

---

        ▷ Orientate with Tomasi-Kanade algorithm

*std :: vector < ElRotation3D > aVR = OrientTomasikanade*

    (*aPrec, mTK\_H123, 3, 50, 1e - 5, (std :: vector < ElRotation3D > \*)NULL*);

        ▷ Test the new solution and accept if smaller residual

    TestSol(aVR);

---

## 5 Optimisation II NO\_SolInit3

The goal of this program is ...

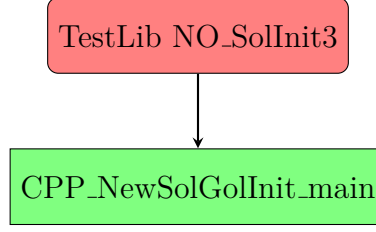


Figure 7: Triplet optimisation II workflow.

### 5.1 Main classes, some typedefs and variables

- *cAppli\_NewSolGolInit* – the manager class
- *cRandNParimiQ* – class for random selections
- *cNOSolIn\_AttrSom* – class corresponding to a node, defined in *SolInitNewOri.h*
- *cNOSolIn\_AttrArc* – class corresponding to an edge, defined in *SolInitNewOri.h*
- *cNOSolIn\_Triplet* – triplet solution
- *cLinkTripl* – it is an ordered *cNOSolIn\_Triplet*
- *typedef ElSomIterator < cNOSolIn\_AttrSom, cNOSolIn\_AttrArc > tItSNSI* – iterator over nodes
- *typedef ElArcIterator < cNOSolIn\_AttrSom, cNOSolIn\_AttrArc > tItANSI* – iterator over arcs

### 5.2 Algorithms

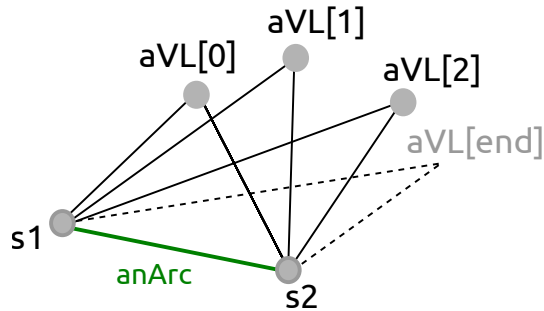


Figure 8: Illustration to Alg. 29. The coherence scores are calculated for all combinations of triplets, e.g. a couple of triplet  $s1 - s2 - aVL[0]$  and  $s1 - s2 - aVL[1]$ .

---

**Algorithm 28** *cAppli\_NewSolGolInit* :: *cAppli\_NewSolGolInit*; def in  
cNewO\_SolGlobInit.cpp

---

▷ Load the graph, i.e. all the nodes/images and the edges : mMapS, mGr and mV3

▷ Calculate an average coherence score for each triplet.

*EstimCoherenceMed*();

▷ Estimate robust rotations from available triplets

*EstimRotsArcsInit*();

▷ dd

*EstimCoheTriplet*();

▷ dd

*FilterTripletValide*();

▷ dd

*mTestTrip* → *CalcCoherFromArcs*(true);

▷ dd

*NumeroteCC*();

▷ dd

*CalculOrient*();

---

---

**Algorithm 29** *EstimCoherenceMed()*; def in cNewO\_SolGlobInit.cpp

---

```

 ▷ Calculate the number of triplet couples that share the same edge and store in:
 aNbTT

 ▷ Iterate over all nodes
 for tItSNSI anItS = mGr.begin(mSubAll); anItS.go_on(); anItS++ do
 tSomNSI * aS1 = &(*anItS);

 ▷ Iterate over all edges attached to node aS1
 for tItANSI anItA = aS1->begin(mSubAll); anItA.go_on(); anItA++ do

 tArcNSI & anArc = (*anItA); current edge ▷ Get vector of "third" images (triplets)
 for that edge
 std::vector< cLinkTripl > & aVL = anArc.attr().ASym()->Lnk3();

 ▷ Two iterations over all "thirds" to calculate some coherence scores between all available
 triplets for the anArc edge. See Fig. 8.
 for intaK1 = 0; aK1 < int(aVL.size()); aK1++ do
 for intaK2 = aK1 + 1; aK2 < int(aVL.size()); aK2++ do
 if aSel.GetNext() then i.e., only a number of calls are executed according to
 aSel(ElMin(aNbTT, NbMaxATT), aNbTT). The NbMaxATT = 100000
 cNOSolIn_Triple * aTri2 = aVL[aK2].m3;

 ▷ Checks coherence from the difference of coordinate systems in TriA and TriB
 aDCAB = DistCoherenceAtoB(&anArc, aTri1, aTri2); see Alg. 30
 aNb = ElMin(aTri1->Nb3(), aTri2->Nb3());
 aVPAB.push_back(Pt2df(aDCAB, aNb));

 ▷ Checks coherence from relative orientations
 aDC12 = DistCoherence1to2(&anArc, aTri1, aTri2); see Alg. 31
 aVP12.push_back(Pt2df(aDC12, aNb));
 end if
 end for
 end for
 end for

 ▷ Saves the median coherence for CohAB and Coh12
 mCoherMedAB = DefMedianPond(0, -1, aVPAB, 0);
 mCoherMed12 = DefMedianPond(0, -1, aVP12, &aKMed);

```

---

---

**Algorithm 30**     *DistCoherenceAtoB(tArcNSI \* anArc, cNOSolInTriplet \* aTriA, cNOSolInTriplet \* aTriB);* def in cNewO\_SolGlobInit.cpp

---

▷ Get orientations of the nodes s1 and s2 in the coordinate system of Triplet A and Triplet 2  
aR1A, aR2A, aR2A, aR2B

▷ Calculate transformations from the c system of tripletA to tripletB. Calculate twice from s1 and s2, then take a mean

$aR1AtoB = aR1B * aR1A.inv()$   
 $aR2AtoB = aR2B * aR2A.inv();$   
 $aMatA2B = NearestRotation((aR1AtoB.Mat() + aR2AtoB.Mat()) * 0.5);$

▷ Calculate the deviation of rotations calculated from s1 and s2 wrt the mean rotation

$aD1 = (aMatA2B - aR1AtoB.Mat()).L2();$   
 $aD2 = (aMatA2B - aR2AtoB.Mat()).L2();$

▷ Calculate the base between s1 and s2 and transform from c sys of tripletA to tripletB (aVA12). Calculate the base in c sys of triplet B (aVB12)

$aVA12 = aMatA2B * (aR2A.tr() - aR1A.tr());$   
 $aVB12 = aR2B.tr() - aR1B.tr();$

▷ Transform the deviation of rotations to distance measure

$aDistRot = sqrt(aD1 + aD2) * (2.0/3);$

▷ Calculate the B to H ratio as a harmonic mean (because of inverse of proportionality)  
 $aBOnH = MoyHarmonik(aTriA -> BOnH(), aTriB -> BOnH());$

▷ Compute deviation on the s1-s2 base calculated from two different triplets. Multiply by BtoH

$aDistTr = DistBase(aVA12, aVB12) * aBOnH;$

▷ The output is the sum of deviation on rotation and the base.

return  $aDistRot + aDistTr;$

---



---

**Algorithm 31**     *DistCoherence1to2(tArcNSI \* anArc, cNOSolInTriplet \* aTriA, cNOSolInTriplet \* aTriB);* def in cNewO\_SolGlobInit.cpp

---

▷ The *RotationC2toC1* return relative orientation of s2 wrt s1 (i.e.  $R1^{-1} \cdot R2$ ). Here, it is calculated twice for two triplets and it is passed to *DistanceRot* together with the B to H ratio harmonic mean.

return  $DistanceRot(RotationC2toC1(anArc, aTriA), RotationC2toC1(anArc, aTriB), MoyHarmonik(aTriA -> BOnH(), aTriB -> BOnH()));$  see Alg. 32.

---

---

**Algorithm 32** *DistanceRot(constElRotation3D&aR1, constElRotation3D&aR2, doubleaBSurH);*  
def in cNewO\_SolGlobInit.cpp

---

▷ Difference of rotations

$$aDif = aR1.Mat() - aR2.Mat();$$

▷ Calculate L2 on the difference

$$aDistRot = \text{sqrt}(aDif.L2());$$

▷ Calculate the euclidean difference between two bases and multiply by the B to H ratio.

$$aDistTr = \text{DistBase}(aR1.tr(), aR2.tr()) * aBSurH;$$

▷ Return the sum of DistRot and DistTr

---

---

**Algorithm 33** *EstimRotsArcsInit();*; def in cNewO\_SolGlobInit.cpp

---

---

**Algorithm 34** *EstimCohTriplets();*; def in cNewO\_SolGlobInit.cpp

---

---

**Algorithm 35** *FilterTripletsValide();*; def in cNewO\_SolGlobInit.cpp

---

## 6 Other

**Thresholds** are defined in *NewOri/NewOri.h*

**Graphs** are defined in *graphes/graphe.h*