# A MULTISCALE FEATURE EXTRACTION APPROACH FOR 3D RANGE IMAGES 

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#### Abstract

This paper presents a multiscale feature extraction technique for 3D range images. Optimized and improved 3D Gaussian filtering and saliency map computation have been conceived jointly to the exploitation of connectivity relationships naturally induced by the 2D acquisition grid. The proposed algorithmic and implementation solutions guarantee good repeatability of detected features on different views and demonstrated superior computational performance compared to other known approaches.


## 1. INTRODUCTION

Nowadays, there is a growing interest in 3D shape acquisition and analysis because of their broad range of application fields (industrial, entertainment, security, medical, cultural heritage,...). In order to acquire the 3D shape of an object, a stereo scan system is usually employed to obtain a set of 3D range images (RI). If no information about relative displacements between scanner and object is available, the different views will not share a common reference system. In order to reconstruct the object shape, it is thus necessary to accurately align each view of the acquisition dataset through a registration procedure. A way to estimate the relative displacement between two views consists into the identification of a set of corresponding points taken from each view and the computation of an optimal rototranslation matrix out of them. Brute force approaches to solve this problem become impractical as RI size increases. In order to speed-up this procedure, a common technique is to reduce the number of candidates by automatically extracting a small set of feature points from both views and searching for correspondences between them. The problem then becomes how to select feature points that, with high repeatability rate, appear in same positions on overlapping areas of different views, i.e. taken from different angle directions. Feature extraction repeatability performance are important within a feature based model acquisition pipeline in that it heavily influences subsequent steps toward automated view alignment for 3D model acquisition [1]. Different solutions have been proposed for 3D feature extraction. In [2]

[^0]points are selected based on curvature criteria. In [3], for each point of the RI, the volume portion of the object inscribed into a sphere surrounding the point is estimated and only uncommon values are selected. Other solutions, like the presented one, are based on multiscale analyses [4], [5], [6] and are somewhat inspired to the Lowe's SIFT approach [7]. Although differing in algorithmic and implementation aspects, a typical sequence of processing steps for feature point extraction is:

- calculate the normals to the RI points (preliminary step)
- calculate N Gaussian filtered (scaled) versions of the RI
- derive N-1 saliency maps from the filtered data
- identify local maxima on each saliency map

Following this approach, we present some improving solutions for feature detection on 3D range data and we measure their performance in terms of feature repeatability and computation speed. Our main contributions are 1) a new directionconstrained Gaussian-like filtering, and 2) naturally induced neighborhood exploitation for computational speed-up.

## 2. FEATURE EXTRACTION ON 3D RANGE IMAGES

In a RI, each acquired pixel (usually on a image sensing grid $\mathbf{x}=(w, h))$ is associated to a depth or distance information (according to the modeled acquisition geometry). Some scanners produce true 3D localization of the point in the acquisition coordinate system. In our formalism a range image is a map $R I: I\left(\mathbb{Z}^{2}\right) \rightarrow R\left(\mathbb{R}^{3}\right)$, where the domain $I$ is a regular rectangular grid (usually corresponding to the CCD matrix), and the co-domain $R$ corresponds to the set of 3D points belonging to the target object. Not all points in $I$ have a valid corresponding point $\mathbf{p}_{\mathbf{x}} \in \mathrm{R}$ (this is due to the nature of the acquisition in terms of target object shape and/or measure range limitations) so a valid point subset $I_{v} \subseteq I$ can be defined. Moreover, although the 2D index grid is uniformly distributed, 3D points cannot be considered uniformly distributed over the 3D space. This is because point density is inversely proportional to the angle existing between the acquisition direction, and the normal associated to each point. Considering $R$ as a surface, its point distribution presents
holes in correspondence to occlusions, i.e. parts of the target object that are not directly visible by the incident rays. Despite these facts (non-valid regions, non-uniform density and holes) the regularity of the 2D grid is a useful structuring resource that can be used as a naturally induced connectivity lattice to be exploited in 3D point processing and to determine practical definitions to various data processing operators (e.g. filtering, interpolation/subsampling, normal computation,...) and speed-up their computation. We exploit this connectivity in all the steps of our method, and we propose original implementations of Gaussian filtering and saliency map calculations on $R I \mathrm{~s}$.

### 2.1. Normal extraction

Our feature extraction algorithm requires the computation of normal vectors at each point location $\mathbf{p}_{\mathbf{x}} \in R$. Given a $\mathbf{p}_{\mathbf{x}}$, its normal is computed by calculating the difference vectors $\mathbf{d}=\mathbf{p}_{\mathbf{j}}-\mathbf{p}_{\mathbf{x}}$ between $\mathbf{p}_{\mathbf{x}}$ and its 8 neighbor points $\mathbf{p}_{\mathbf{j}}$ in the 2D index grid. By assigning an index $k$, ranging from 1 to 8, for each vector $\mathbf{d}$, associated in a anti-clockwise order (see Fig.1), we define the normal $\hat{\mathbf{n}}_{\mathbf{x}}$ as:

$$
\hat{\mathbf{n}}_{\mathbf{x}}=\frac{\sum_{k=1}^{8} \hat{\mathbf{d}}_{k} \times \hat{\mathbf{d}}_{j}}{8} \quad j= \begin{cases}k+2 & k<=6 \\ k-6 & k>6\end{cases}
$$

Fig. 1. $\hat{\mathbf{d}}$ selection for (a) $k=1$; (b) $k=2$; (c) $k=8$

### 2.2. 3D Gaussian filtering

An important step for our algorithm requires to extract a set of N Gaussian filtered versions $G(n)$ of the input range image, with $n=[1, N]$. In order to do so, N 3D Gaussian filters with progressively increasing kernel dimension $\sigma_{n}$ are applied to each point constituting the original range image, varying its position as follows:

$$
\begin{equation*}
\mathbf{p}_{\mathbf{x}}^{f}(n)=\frac{\sum_{\mathbf{p}_{\mathbf{j} \in \eta_{\mathbf{x}}\left(2 \sigma_{n}\right)}} \mathbf{p}_{\mathbf{j}} \cdot e^{-\frac{\left\|\mathbf{p}_{\mathbf{x}}-\mathbf{p}_{\mathbf{j}}\right\|^{2}}{2 \cdot \sigma_{n}^{2}}}}{\sum_{\mathbf{p}_{\mathbf{j} \in \eta_{\mathbf{x}}\left(2 \sigma_{n}\right)}} e^{-\frac{\left\|\mathbf{p}_{\mathbf{x}}-\mathbf{p}_{\mathbf{j}}\right\|^{2}}{2 \cdot \sigma_{n}^{2}}}} \tag{2}
\end{equation*}
$$

where $\sigma_{n}$ represents the $n_{t h}$ kernel radius, while $\eta_{\mathbf{x}}\left(2 \sigma_{n}\right)$ identifies the points within a distance $2 \sigma_{n}$ from $\mathbf{p}_{\mathbf{x}}$.


Fig. 2. $\mathbf{p}_{\mathbf{x}}^{f}(n)$ projected over $\hat{\mathbf{n}}_{\mathbf{x}}$ direction

The set of $\mathbf{p}_{\mathbf{x}}^{f}(n)$, as calculated in (2), could be directly used (as done in [5] and [6]) as filtered data. However, we do not yet consider it as such. This is because, due to the above described characteristics of $R$, border points (on validity region or hole boundaries) and points lying between two regions with different normal directions would suffer from shrinking and position biasing distortions respectively. These distortions cannot be considered negligible, since most feature points are located over highest curvature regions and therefore they are prone to position biasing due to point density disparity over $\eta_{\mathbf{x}}\left(2 \sigma_{n}\right)$. In addition, shrinking distortion would affect border areas and would lead to artifacts in the saliency computation, as we will see. In order to avoid such distortions, we propose to project $\mathbf{p}_{\mathbf{x}}^{f}(n)$ over the normal direction $\hat{\mathbf{n}}_{\mathbf{x}}$, as follows:

$$
\begin{equation*}
G_{\mathbf{x}}(n)=\mathbf{p}_{\mathbf{x}}+\left\langle\mathbf{p}_{\mathbf{x}}^{f}(n)-\mathbf{p}_{\mathbf{x}}, \hat{\mathbf{n}}_{\mathbf{x}}\right\rangle \hat{\mathbf{n}}_{\mathbf{x}} \tag{3}
\end{equation*}
$$

This is equivalent to force $\mathbf{p}_{\mathbf{x}}^{f}(n)$ to move along the axis associated to $\hat{\mathbf{n}}_{\mathbf{x}}$, therefore reducing the distortion caused by non-uniform points distribution over $\eta_{\mathbf{x}}\left(2 \sigma_{n}\right)$, as depicted in Fig.2.
We define $G(n)$, the set of points $G_{\mathbf{x}}(n), \forall \mathbf{x} \in \mathrm{RI}$, as the filtered version of the range image after having applied a Gaussian filter of radius $\sigma_{n}$, as defined in (3). As the kernel radius $\sigma_{n}$ increases, details which size is smaller than $\sigma_{n}$ disappear from $G(n)$. Once the $G(n)$ has been calculated, its normal vectors are recomputed as indicated in section 2.1.
As usually done in multiscale approaches, when kernel size doubles, a subsampling factor of two is applied to the image, also to reduce the computational burden. Here subsampling is performed on $I$, thus in a very simple and fast way.

### 2.3. Saliency maps calculation

Once the $G(n)$ have been computed, $\mathrm{N}-1$ saliency maps can be derived. A saliency map is a 2D array of difference values, obtained by pairwise subtraction of $G(n)$ at different kernel sizes. In a multiscale framework, a pairwise subtraction between $G(n)$ retains only the details comprised between the two scale limits (which can be roughly thought as cut-off frequencies), in other words it highlights features which dimension is comprised between two kernel sizes, say $\sigma_{n}$ and
$\sigma_{n+1}$. Thanks to the shrinking artifact avoidance guaranteed by (3), straightforward pointwise subtraction at different scales can be performed correctly. We then calculate saliency maps $S(n)$ as follows:

$$
\begin{equation*}
S(n)=|G(n)-G(n+1)| \cdot\left\langle\hat{\mathbf{n}}_{\mathbf{x}}(n), \hat{\mathbf{n}}_{\mathbf{x}}(n+1)\right\rangle \tag{4}
\end{equation*}
$$

Unlike [4] and [5], the factor $\left\langle\hat{\mathbf{n}}_{\mathbf{x}}(n), \hat{\mathbf{n}}_{\mathbf{x}}(n+1)\right\rangle$ has been introduced in order to give more importance to points for which the normal direction does not change after the filtering process, thus restricting saliency distribution to more specific areas of the range image. A similar approach is employed in [6], where the correction factor takes into account $\hat{\mathbf{n}}_{\mathbf{x}}(n)$.

### 2.4. Maximum extraction

Once each saliency map $S(n)$ has been determined, saliency maximums are extracted as follows:

$$
\begin{equation*}
\operatorname{sMax}(n)=\left\{\mathbf{x} \mid S_{\mathbf{x}}(n)>S_{\mathbf{j}}(n), \forall \mathbf{j} \in \eta_{\mathbf{x}}\left(2 \sigma_{n+1}\right)\right\} \tag{5}
\end{equation*}
$$

In order to generate the maximum list for a saliency map, an iterative search is performed. Once the greatest valid saliency value for $S(n)$ is found, no other maximum may be selected within a neighborhood region of $\eta_{\mathbf{x}}\left(2 \sigma_{n+1}\right)$. This is because the greatest detail size that can be detected within $S(n)$ is $\sigma_{n+1}$. Such constraints on the distance between maxima guarantee the absence of overlap between features.
Each maximum is further tested through two conditions before being selected as feature point. At first, its neighborhood must be defined. In order to avoid selecting features which descriptor would result calculated out of few neighbor points, maximums that possess less than $50 \%$ of the estimated points for $\eta_{\mathbf{x}}\left(2 \sigma_{n}\right)$ are discarded. The second check is about position stability. Saliency maps may slightly vary due to changes of acquisition direction, therefore maximums detected over regions with similar saliency values may vary their position significantly. In order to avoid such points, we only retain maximums for which the saliency value is much greater than its neighboring points, and it is not located over saliency ridges. An example of saliency map with superimposed feature points at a given scale is shown in Fig.3.

## 3. RESULTS

To assess the contribution of each step and the overall performance of the proposed feature extraction method, four different implementations have been tested, depending on which version of filtering and saliency computation have been used. The first version, we called basic implementation, reflects what proposed in [5], i.e. Gaussian filtering performed as in (2) and saliency computed as pairwise difference between filtered versions of data. The second version only implements the improved version of Gaussian filter (3); the third version only implements the saliency correction (4). Al last, the


Fig. 3. A saliency map with superimposed feature points.


Fig. 4. Datasets: (a) Angels, (b) Dolphin, (c) Teeth.
fourth version implements both the proposed improvements. The comparison has been performed over three pre-aligned range image datasets which are shown in Fig. 4 (where different colors have been used for each view). All $R I$ s have been acquired with a commercial structured-light scanner (Optical Rev-Eng LE, produced by Open Technologies srl, Italy). All the range images used for our tests have a resolution of 1280 $\times 1024$, which means that $I_{v}$ can have up to over 1.3 million points. Prior to the first Gaussian filtering, each range image is down-sampled by a factor 2 , thus reducing initial dimensions to $640 \times 512$. In our implementation, three $G(n)$ have been calculated, with $\sigma_{n}=2 \sigma_{n-1}$, while the $\sigma_{1}$ is set to be eight times the average interdistance between points in $R$. Although, according to the SIFT methodology [7], the implemented software can work with any number of scaling level and $\sigma$ progression, we experimentally found the above parameters to give the best compromise between repeatability and computational performances for the considered datasets. Repeatability performance evaluations have been executed as follows: from each range image of a dataset, its features are extracted, then for each pair of consecutive range images (according to a predetermined order), features in the overlapping area are counted; subsequently, both feature sets are compared in order to find correspondences (feature points are considered correspondent if their position differ in less than $\frac{\sigma}{3}$ ). At last, feature repeatability is computed as the number of features that have a correspondence relation divided by the

| Feat. extraction versions | Feature repeatability |  |  |
| :---: | :---: | :---: | :---: |
|  | Angels | Dolphin | Teeth |
| Basic implementation | $37 \%$ | $8 \%$ | $19 \%$ |
| Gaussian projection | $56 \%$ | $34 \%$ | $41 \%$ |
| Saliency projection | $35 \%$ | $11 \%$ | $19 \%$ |
| Proposed implementation | $\mathbf{5 7 \%}$ | $\mathbf{3 6 \%}$ | $\mathbf{4 1 \%}$ |

Table 1. Repeatability performance

| Dataset <br> name | RI <br> No. | Min-max <br> overlap | Avg points <br> $\left(\times \mathbf{1 0}^{\mathbf{3}}\right)$ | Exec time <br> [ms] |
| :---: | :---: | :---: | :---: | :---: |
| Angels | 8 | $20-60 \%$ | 1040 | 4823 |
| Dolphin | 20 | $20-90 \%$ | 406 | 1894 |
| Teeth | 8 | $40-90 \%$ | 408 | 1836 |

Table 2. Computational performance
number of features in the overlapping area. In Table1, feature repeatability performances are presented for each dataset and for each feature extraction version. Reported data are averaged over all scales and all couples of overlapping RIs. Our evaluation shows that the modified Gaussian filter improves repeatability performances by more than 20 percentage points with respect to the basic implementation of the Gaussian filter (used in the first and third version). Such improvement is due to the fact that the proposed filter does not introduce any shrinking effect usually associated to the Gaussian filtering, thus improving feature localization. The saliency correction does not yield as much, however a slight performance improvement is achieved through its introduction. It can be argued that such repeatability rates may not be considered as extraordinary. As a matter of fact, to our knowledge, no study about feature repeatability have been published. In [4], some feature repeatability results against range data subsampling and noise addition are presented, but they are not useful here for comparison. We deem our performances to be satisfactory because of two reasons. The first one is related to the image sets employed, since they present a heterogeneous grade of regularity, different feature types and dimensions, and usually big differences (borders, holes, point density) when acquired with different acquisition directions. Secondly, we deem such repeatability rates as entirely adequate for a robust subsequent search of correspondences, if suitable feature descriptors or signatures can be calculated and used to this end.
Computational performances are presented in Table 2, where we also describe datasets in terms of number of RI views per dataset, min-max overlap area between views pairs, and average number of points (i.e. $\left|I_{v}\right|$ ) per RI. Algorithms have been implemented in $\mathrm{C}++$, and executed on a PC with INTEL CORE2 CPU, 2.80 GHz each, and 4 GB of RAM. As the differences in computational time for the above four versions are negligible, we only give the processing time required by the proposed implementation, which is around 4.5 seconds for each million of points. This is quite an improvement with respect to [5] (around 5 minutes per million points), [4] (50
minutes per million points), [3] (around 5 minutes per million points) and [6] (around 7 minutes per million points), even considering the upgraded computational resources. Such performance boost is mainly due to the exploitation of neighborhood relationships defined on $I$. Secondly, the use of a fast subsampling, and some coding efficiency solutions on frequent and time-consuming operations make the rest (e.g. look-up tables for square root computing).

## 4. CONCLUSION

In this paper a multiscale method for 3D feature extraction on range images has been presented. This method, that can also be considered as a 3D extension of a SIFT-based approach, demonstrated to yield good robustness in terms of feature repeatability, and outstanding computational performance. The work is mainly characterized by the introduction of a modified version of Gaussian filtering of 3D range images and a weighted saliency map computation. They both contribute to improve feature repeatability. In addition, the direct exploitation of the 2D domain neighborhood helps the definition of the filter action and is determinant for the computational speed. Experimental performance confirm that our method is suitable to be used for automated multiple view registration and 3D model acquisition, where the degree of accuracy and obtainable automation strongly depends from the goodness of the feature extraction phases. The proposed solutions can be also useful for other 3D feature-based applications, such as content analysis and description of 3D objects, 3D object tracking, etc., where features robustness and computational speed are crucial aspects.

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[^0]:    *The first author performed the work while at Open Technologies srl, Italy

