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Estimating Differential Quantities using Polynomial fitting of Osculating Jets

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Abstract: This paper addresses the pointwise estimation of differential properties of a smooth manifold *S* —a curve in the plane or a surface in 3*D*— assuming a point cloud sampled over *S* is provided. The method consists of fitting the local representation of the manifold using a jet, and either interpolation or approximation. A jet is a truncated Taylor expansion, and the incentive for using jets is that they encode all local geometric quantities —such as normal, curvatures, extrema of curvature. On the way to using jets, the question of estimating differential properties is recasted into the more general framework of multivariate interpolation/approximation, a well-studied problem in numerical analysis. On a theoretical perspective, we prove several convergence results when the samples get denser. For curves and surfaces, these results involve asymptotic estimates with convergence rates depending upon the degree of the jet used. For the particular case of curves, an error bound is also derived. To the best of our knowledge, these results are among the first ones providing accurate estimates for differential quantities of order three and more. On the algorithmic side, we solve the interpolation/approximation problem using Vandermonde systems. Experimental results for surfaces of \mathbb{R}^3 are reported. These experiments illustrate the asymptotic convergence results, but also the robustness of the methods on general Computer Graphics models.

Key-words: Meshes, Point Clouds, Differential Geometry, Interpolation, Approximation.

Estimation des Quantités Différentielles par Ajustement Polynomial des Jets Osculateurs

Résumé : Ce rapport concerne l'estimation locale des propriétés différentielles d'une variété lisse S –une courbe dans le plan ou une surface en 3D– à partir d'un nuage de points échantillonnés sur S. La méthde consiste à ajuster la représentation locale de la variété par un jet, en interpolant ou approximant. Un jet est un développement de Taylor tronqué, et l'intérêt des jets est qu'ils codent toutes les quantités géométriques locales –telles que la normale, les courbures, les extrema de courbure. Avec l'utilisation des jets, le problème d'estimation des quantités différentielles est placé dans le cadre plus général de l'interpolation/approximation multivariée, un sujet classique d'analyse numérique. Sur le plan théorique, nous donnons des résultats de convergence lorsque l'échantillonnage est raffiné. Pour les courbes et surfaces, ces résutats sont des estimations asymptotiques avec des vitesses de convergence fonction du degré du jet utilisé. Pour le cas des courbes, une majoration d'erreur est aussi fournie. A notre connaissance, ces résutats sont parmi les premiers fournissant des estimations précises pour les quantités différentielles d'ordre trois et plus. Sur le plan algorithmique, nous traitons le problème d'interpolation/approximation avec des systèmes de Vandermonde. Des résultats expérimentaux pour les surfaces de R3 sont analysés. Ces expérimentations illustrent les résultats de convergence asymptotique, ainsi que la robustesse de la méthode sur des modèles de *Computer Graphics.*

Mots-clés : Maillages, Nuages de Points, Géométrie différentielle, Interpolation, Approximation.

1 Introduction

1.1 Estimating differential quantities

Several applications from Computer Vision, Computer Graphics, Computer Aided Design or Computational Geometry require estimating local differential quantities. Example such applications are surface segmentation, surface smoothing / denoising, surface reconstruction, shape design. In any case, the input consists of a point cloud or a mesh. Most of the time, estimating first and second order differential quantities, that is the tangent plane and curvature-related quantities, is sufficient. However, applications involving shape analysis $[H+99, Por01]$ require estimating third order differential quantities.

For first to third order differential quantities, a wealth of different estimators can be found in the vast literature of applied geometry [Pet01]. Most of these are adaptations to the discrete setting of smooth differential geometry results. For example, several definitions of normals, principal directions and curvatures over a mesh can be found in [Tau95, CW00]. Ridges of polyhedral surfaces as well as cuspidal edges of the focal sets are computed in [WB01]. Geodesics and discrete versions of the Gauss-Bonnet theorem are considered in [PS98].

A striking fact about estimation of second order differential quantities —using conics and quadrics— - is that the classification of Euclidean conics/quadrics is never mentioned. Another prominent feature is that few contributions address the question of the accuracy of these estimates or that of their convergence when the mesh or the sample points get denser. The question of convergence is one prime importance since estimates do not always asymptotically behave as one would expect. For example, it is proved in[BCM02] that the angular defect of triangulations does not in general provide information on the Gauss curvature of the underlying smooth surface.

The following are provably good approximation results. In [AB99], an error bound is proved on the normal estimate to a smooth surface sampled according to a criterion involving the skeleton. The surface area of a mesh and its normal vector field versus those of a smooth surface are considered in [MT01]. Asymptotic estimates for the normal and the Gauss curvature of a sampled surface for several methods are given in [MW00b]. In particular, a degree two interpolation is analyzed. Based upon the normal cycle and restricted Delaunay triangulations, an estimate for the second fundamental form of a surface is developed in [CSM03].

Deriving provably good differential operators is the goal pursued in this paper. To motivate our guideline and before presenting our contributions, we raise the following question. Second order differential properties for plane curves are almost always investigated using the osculating circle, while principal curvatures of surfaces are almost always computed using osculating paraboloids. Why not osculating parabolas for curves and osculating ellipsoids or hyperboloids for surfaces? Before answering this question and to clarify the presentation, we recall some fundamentals.

1.2 Curves and surfaces, height functions and jets

It is well known [dC76, Spi99] that any regular embedded smooth 1 curve or surface can be locally written as the graph of a univariate or bivariate function with respect to any *z* direction that does not belong to the tangent space. We shall call such a function a *height function*. Taking an order *n* Taylor expansion of the height function over a curve yields:

$$
f(x) = J_{B,n}(x) + O(x^{n+1}),
$$
\n(1)

with

$$
J_{B,n}(x) = B_0 + B_1 x + B_2 x^2 + B_3 x^3 + \dots + B_n x^n.
$$
 (2)

Similarly for a surface:

$$
f(x, y) = J_{B,n}(x, y) + O(||(x, y)||^{n+1}),
$$
\n(3)

with

$$
J_{B,n}(x,y) = \sum_{k=1}^{n} H_{B,k}(x,y), H_{B,k}(x,y) = \sum_{j=0}^{k} B_{k-j,j} x^{k-j} y^{j}.
$$
 (4)

Borrowing to the jargon of singularity theory [BG92], the truncated Taylor expansion $J_{B,n}(x)$ or $J_{B,n}(x, y)$ is called a degree *n* jet, or *n*-jet. Since the differential properties of a *n*-jet match those of its defining curve/surface up to order *n*, the jet is said to have a *n* order contact with its defining curve or surface. This also accounts for the term *osculating jet* —although osculating ² was initially meant for 2-jets. The degree *n*-jet of a curve involves $n + 1$ terms. For a surface, since there are $i + 1$ monomials of degree *i*, the *n*-jet involves $N_n = 1 + 2 + \cdots + (n+1) = (n+1)(n+2)/2$ terms. Notice that when *z* direction used is aligned with the normal vector to the curve/surface, one has $B_1 = 0$ 0 or $B_{10} = B_{01} = 0$. The osculating *n*-jet encloses differential properties of the curve/surface up to order *n*, that is any differential quantity of order *n* can be computed from the *n*-jet. In particular, the tangent space can be computed from the 1-jet, the curvature related information can be obtained from the 2-jet, locating ridges require coefficients of the 3-jet, and so on. To clarify the presentation, we summarize as follows:

Definition. 1 *For a curve or surface:*

- *• given a coordinate system, the* osculating *n*-jet *is the Taylor expansion of the height function truncated at order n,*
- *• the osculating n-jet is* principal *if the linear terms vanish in this coordinate system (i.e. the ordinate axis is the normal direction of the manifold),*
- *• an* osculating conic/quadric *is a conic/quadric whose 2-jet matches that of the curve/surface (independently of a given coordinate system),*
- *• an osculating conic/quadric is* degenerate *if it is the graph of its 2-jet,*

¹ Regular means that the tangent space has dimension one/two for a curve/surface everywhere. Embedded forbids selfintersections. Smooth means as many times differentiable as we need, typically C^3 or C^4 .

²From the latin osculare, to kiss.

• an osculating conic/quadric is principal *if its 2-jet is principal.*

Degenerate osculating conics/quadrics are specific curves and surfaces since:

Theorem. 1 *[Ber87, Chapter 15] There are 9 Euclidean conics and 17 Euclidean quadrics.*

Observation. 1 *The degenerate osculating conics to a smooth curve are parabola or lines. The degenerate osculating quadrics to a smooth surface are either paraboloids (elliptic, hyperbolic), parabolic cylinders, or planes.*

Principal degenerate osculating conics and quadrics are therefore respectively 2 out of 9 conics and 4 out of 17 quadrics. Degenerate stands for the fact that the quadratic forms these conics/quadrics are defined with do not have full ranks.

Principal degenerate osculating conics and quadrics are related to the so-called Monge form of the curve or surface, that is the local Taylor expansion of the curve/surface in the Monge coordinate system. The Monge coordinate system of a curve is defined by its tangent and normal vectors. For a surface, the Monge coordinate system is such that the *z* axis is aligned with the normal and the *x*, *y* axis are aligned with the principal directions. In this particular system, the height function is called the Monge form, and letting k_1, k_2 stand for the principal curvatures, one has ³:

$$
f(x, y) = \frac{1}{2}(k_1x^2 + k_2y^2) + hot
$$
 (5)

From these observations, the question we ended paragraph 1.1 with can now be answered. By theorem 1 and observation 1, using a general conic/quadric or a principal degenerate one to approximate a curve or a surface does not make any difference. In both case and up to order two, the local differential properties of the curve/surface, degenerate conic/quadric, or full rank conic/quadric are identical. As an example, consider Figure $1(a,b,c)$. Figure $1(a)$ features a curve and its osculating circle 4 . In (b), the osculating circle is replaced by the principal osculating parabola —whose symmetry axis is the normal to *C* and whose curvature matches that of *C*. At last, in (c) a general parabola locally approximates *C*. Its symmetry axis is not aligned with the normal to *C*.

Summarizing our discussions so far, the rationale for fitting the *n*-jet of a curve/surface is that this polynomial contains all the differential information up to order *n*.

³ *hot* stands for *higher order terms* in the Taylor expansion.

⁴ A point worth noticing is the relative position of *C* and its osculating circle: the former usually crosses the latter at the intersection point. To see why, just take an order three Taylor expansion and check the relative positions of the curve and the circle!

Figure 1: A curve and (a) its osculating circle (special case of osculating conic), (b) its principal degenerate conic, (c) an degenerate osculating conic

1.3 Interpolation, approximation and related variations

Our methodology to retrieve differential quantities consists of fitting the osculating jet. The following variations need to be discussed in order to state our contributions precisely. The case of curves and surfaces being tantamount, our descrption focuses on surfaces. Assume we are given a set of *N* points $p_i(x_i, y_i, z_i)$, $i = 1, ..., N$ in the neighborhood of a given point *p* on the surface processed. Point *p* itself may or may not be one of the *N* samples, and one can assume without loss of generality that *p* is located at the origin of the coordinate system used.

Interpolation versus approximation. Interpolating consists of finding a polynomial that fits exactly a set of data points. In our case and following Equation (3), let *B* index a coefficient of the jet of the surface, and *A* index a coefficient of the jet sought ⁵. We aim at finding a *n*-jet $J_{A,n}$ such that

$$
f(x_i, y_i) = J_{B,n}(x_i, y_i) + O(||(x_i, y_i)||^{n+1}) = J_{A,n}(x_i, y_i), \quad \forall i = 1, ..., N.
$$
 (6)

Approximation, on the other hand, gives up on exactness, that is the graph of the jet sought may not contain the sample points. We shall focus on *least-square* approximation, which consists of minimizing the sum of the square errors between the value of the jet and that of the function. The quantity to be minimized is therefore

$$
\sum_{i=1}^{N} (J_{A,n}(x_i, y_i) - f(x_i, y_i))^2.
$$
 (7)

The two problems can actually be written in the same matrix form. To see why, let us write the jets in the polynomial basis consisting of monomials $x^i y^j$. Example other basis that could be used are the Bezier-Bernstein basis or the Newton basis. We use the monomials since this basis is convenient

⁵As a mnemonic, the reader may want to remind that index *A* stands for the *Answer* to the fitting problem.

for the asymptotic analysis but also the design of effective algorithms. Denote *A* be the *Nn*-vector of the coefficients of the jet sought, that is

$$
A = (A_{0,0}, A_{1,0}, A_{0,1}, \dots, A_{0,n})^t.
$$

Denote *B* the *N*-vector of the ordinates, i.e. with $z_i = f(x_i, y_i)$,

$$
B = (z_1, z_2, \dots, z_N)^t = (J_{B,n}(x_i, y_i) + O(||(x_i, y_i)||^{n+1}))_{i=1,\dots,N}.
$$

Equations (6) and (7) yield the following $N \times N_n$ Vandermonde matrix

$$
M = (1, x_i, y_i, x_i^2, \dots, x_i y_i^{n-1}, y_i^n)_{i=1,\dots,N}.
$$
\n(8)

For the interpolation case, the number of points matches the number of parameters, so that matrix *M* is square and Eq. (6) can be written as $MA = B$. For the approximation case, *M* is a rectangular $N \times N_n$ matrix, and Eq. (7) is summarized as $\min ||MA - B||_2$.

Choosing between interpolation and approximation depends upon the problem tackled. For noisy datasets, approximation is the method of choice. Otherwise, the alternative depends of the relative values of the number of model parameters versus the number of available points. If the two match one-another, a natural choice is interpolation. In any case, fitting yields a linear system, so that numerical issues arise. Facing these difficulties is the topic of section 2.

Mesh or meshless methods. An important difference between local geometry estimation algorithms is whether or not they require some topological information —typically the connectivity of a mesh. Mesh-based methods are usually faster. Meshless techniques are more general and better suited for noisy datasets. A difficulty of the latter methods, however, is to select the relevant points used to perform the estimates. While one can always resort to heuristics of the *k*-nearest-neighbors type, user defined parameters should be avoided. This issue is addressed in section 5.

One or two stages methods. Fitting a 2-jet requires estimating the tangent plane and the curvature related information. These steps can be carried out sequentially or simultaneously. Following the guideline of [SZ90], most of the methods already mentioned proceed sequentially. The provably good algorithm we propose proceeds simultaneously. Along its analysis, we also provide theoretical results on the accuracy of sequential methods.

1.4 Contributions

Jet interpolation and approximation. Consider Eqs. (6) and (7). We expect $J_{A,n}$ and $J_{B,n}$ to be equivalent in some sense. To specify this, we shall study the convergence properties of the coefficients of $J_{A,n}$ when the points p_i s converge to p . More precisely, assume that the coordinates of the p_i s are given by $p_i(x_i = a_i h, y_i = b_i h, z_i = f(x_i, y_i))$. Parameters a_i and b_i are arbitrary real numbers, while h specifies that the p_i s uniformly tend to the origin. We actually expect

$$
A_{ij} = B_{ij} + O(r(h)).
$$

Function $r(h)$ describes the convergence rate or the precision of the fitting, and the main contribution of this paper is to quantify $r(h)$ for interpolation and approximation methods. As we shall see, interpolation or approximation of the same degree yield the same convergence rate. The difficulties posed are also similar and are essentially to deal with singular matrices.

Relationship to previous theoretical contributions. The theoretical results we are aware of are twofold. First, in [MW00a, Lemma 4.1], a degree two interpolation is used and analyzed. We deal with jets or arbitrary degree, for interpolation and approximation. Second, convergence results on the coefficients of the Lagrange interpolation polynomial versus the Taylor expansion of a function are proved in [Coa66]. Our results match those, but we also analyze the approximation case. Most importantly, we give constructive proofs and develop the corresponding algorithms.

It should be noticed that we are *not* concerned here with the convergence of the Lagrange interpolation polynomial to the height function on a whole given set. This problem requires specific conditions on the function and the position of the points, as illustrated by the Runge divergence phenomena [LS86, Chapter 2]. Therefore, our study is not to be confused with global fitting such as piecewise polynomial fitting encountered in CAD.

2 Interpolation, approximation, numerical issues

In this section, we recall the fundamentals of the fitting methods used, namely interpolation and approximation, together with the numerical issues arising from the resolutions.

2.1 Interpolation

The interpolation fitting is based upon the Lagrange interpolation, that is the construction of a polynomial constrained to fit a set of data points. Although this problem is classical for the univariate case, the multivariate case is still an active research field from both the theoretical and computational points of view. We briefly review the univariate and multivariate basics of Lagrange interpolation.

Univariate Lagrange interpolation. Let $X = \{x_0, \ldots, x_n\}$ be $n + 1$ distinct real values, the socalled *nodes*. Then, for any real function f , there is a unique polynomial P of degree n so that $P(x_i) = f(x_i)$, $\forall i = 0, \dots, n$. Polynomial *P* is called the *Lagrange interpolation polynomial* of *f* at the nodes *X*. For any choice of distinct nodes, this polynomial exists and is unique, and in that case the Lagrange interpolation is said to be *poised*.

Multivariate Lagrange interpolation. Consider now the following bivariate problem. Let Π*ⁿ* be the subspace of bivariate polynomials of total degree equal or less than *n*, whose dimension is $N_n = \binom{n+2}{n}$, and let $X = \{x_1, \ldots, x_N\}$ consist of $N = N_n$ values in \mathbb{R}^2 called nodes. (Notice that *N* is exactly the number of monomials found in the jet of Equation 3.) The Lagrange interpolation problem is said to be poised for *X* if for any function $f : \mathbb{R}^2 \to \mathbb{R}$, there exists a unique polynomial P in Π_n so that $P(x_i) = f(x_i)$, $\forall i = 1, ..., N$. It is intuitive and well known that this problem is poised iff the set of nodes X is not a subset of any algebraic curve of degree at most n , or equivalently the Vandermonde determinant formed by the interpolation equations does not vanish. As noticed in [SX95], the set of nodes for which the problem is not poised has measure zero, hence it is almost always poised.

However let us illustrate non-poised cases and almost non-poised or degenerate ones. Consider the two quadrics $q_1(x, y) = 2x + x^2 - y^2$ and $q_2(x, y) = x^2 + y^2$, whose intersection curve *I* projects in the (x, y) plane to the conic $C(x, y) = 0$ with $C(x, y) = x - y^2$ —Figure 2. If one tries to interpolate a height function using points on *I*, uniqueness of the interpolant is not achieved since any quadric in the pencil of q_1 and q_2 goes through *I*. A similar example featuring the four one-ring and one two-ring neighbors of a point p is depicted on figure 3. Notice that being able to figure out such configurations is rather a strength than a weakness of the method since a surface is sought and, the amount of information available does not determine uniquely this surface. A first fundamental difference between the univariate and multivariate cases is therefore the critical issue of choosing nodes so that the interpolation is poised.

In the particular case where the points lies on a regular square grid of the plane, the geometry of the configuration leads to the following remarks. On one hand, a non-poised degree *n* interpolation occurs if the points lies on *n* lines, since they define an algebraic curve of degree *n*. One the other hand, triangular lattices yield poised problems for every degree. These results and further extensions can be found in [GS00] and references therein.

2.2 Least square approximation

It is well known that the minimization problem of Eq. (7) has a unique solution iff the matrix *M* is of maximum rank N_n . In that case, the minimum value ρ is called the *residual* of the system, that is $\rho = \min ||MA - B||_2$. The important issue is again the rank of the matrix *M*. In terms of the relative values of N versus N_n , using too many points certainly smoothes out geometric features, but also makes rank deficient matrices less likely.

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2.3 Numerical Issues

The difficulties of solving linear and least-squares systems consist of dealing with rank-deficient matrices. We now discuss these issues in more detail. Distances between matrices and matrix norms refer to the Euclidean norm.

Singular systems and condition numbers. To quantify degeneracies, we resort to a Singular Value Decomposition (SVD) [GvL83]. Denote $\sigma_n, \ldots, \sigma_1$ the singular values of *M* sorted in decreasing order. It is well known that the least singular value of *M* is the distance from *M* to rank deficient matrices. The singular values also characterizes the sensitivity of the problem, that is the way errors on the input data induce errors on the computed solution. Notice that errors refer to the uncertainty attached to the input data and not to the rounding errors inherent to the floating point calculations. In our case, input data are the sample points, so that errors are typically related to the acquisition system —e.g. a laser range scanner.

To quantity the previous arguments, we resort to the *conditioning* or *condition number* of the system [GvL83, Hig96]. The conditioning is defined as a magnification factor which relates the afore-mentioned errors by the following rule $Error$ on solution = $Error$ on input \times conditioning. Denote

$$
\kappa_2(M) = ||M||_2||M^{-1}||_2 = \sigma_n/\sigma_1
$$

the condition number of the matrix *M*. The conditioning in the two cases are respectively given by

$$
\begin{cases}\n\text{linear square system:} & \kappa_2(M), \\
\text{least square system:} & \kappa_2(M) + \kappa_2(M)^2 \rho \text{ with } \rho = ||MX - B||_2 \text{ the residual.} \n\end{cases} \tag{9}
$$

The following theorem provides precise error bounds:

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Theorem. 2 *Suppose X and* \widetilde{X} *are the solutions of the problems*

$$
\begin{cases}\n\text{linear square system:} & MX = B \text{ and } (M + \Delta M)\tilde{X} = B + \Delta B, \\
\text{least square system:} & \min||MX - B||_2 \text{ and } \min||((M + \Delta M)\tilde{X} - (B + \Delta B)||_2,\n\end{cases}\n\tag{10}
$$

with ε *a* positive real value such that $||\Delta M||_2 \leq \varepsilon ||M||_2$, $||\Delta B||_2 \leq \varepsilon ||B||_2$, and $\varepsilon \kappa_2(M) < 1$. Then *one has:*

$$
\frac{||X - X||_2}{||X||_2} \le \varepsilon \text{ conditioning.} \tag{11}
$$

In practice, if the conditioning is of order 10^a and the relative error on the input is $\varepsilon \approx 10^{-b}$ — with $\epsilon \kappa_2(M) < 1$, then the relative error on the solution is of order 10^{a-b} .

Pre-conditioning the Vandermonde system. As already discussed, a convenient way to solve Eqs. (6) and (7) consists of using the basis of monomials. One ends up with the Vandermonde matrix of Eq. (8), that can be solved with usual methods of linear algebra. Unfortunately, Vandermonde systems are known to be ill-conditioned due to the change of magnitude of the terms. We therefore pre-condition so as to improve the condition number. Assuming the x_i s, y_i s are of order *h*, the pre-conditioning consists of performing a column scaling by dividing each monomial $x_i^k y_i^l$ by h^{k+l} .

The new system is $M'Y = MDY = B$ with *D* the diagonal matrix $D = (1, h, h, h^2, \ldots, h^n, h^n)$, so that the solution *A* of the original system is $A = DY$. The condition number used in the sequel is precisely $\kappa(M')$. (Notice it has the geometric advantage to be invariant under homothetic transformations of the input points.) Then the accuracy of the result can be estimated a posteriori, and almost degenerate cases hight-lighted by large conditioning.

Alternatives for the interpolation case. An alternative to the Vandermonde system consists of using the basis of Newton polynomials. Resolution of the system can be done using divided differences [Sau95], a numerically accurate yet instable method [Hig96].

3 Surfaces

3.1 Problem addressed

Let *S* be a surface and *p* be a point of *S*. Without loss of generality, we assume *p* is located at the origin and we aim at investigating differential quantities at *p*. Consider the height function *f* given by Equation (3) in any coordinate system whose *z* axis is not in the tangent plane. We shall interpolate *S* by a bivariate *n*-jet $J_{A,n}(x, y)$ whose graph is denoted *Q*.

The normal to a surface given by Equation (3) is

$$
n_S = (-B_{10}, -B_{01}, 1)^t / \sqrt{1 + B_{10}^2 + B_{01}^2}.
$$
\n(12)

In order to characterize curvature properties, we resort to the Weingarten map *A* of the surface also called the shape operator, that is the tangent map of the Gauss map. (Recall that the second fundamental form *II* and *A* satisfy $I(v) = \langle A(v), v \rangle$ for any vector *v* of the tangent space.) The principal curvatures and principal directions are the eigenvalues(eigenvectors) of *A*, and the reader is referred to [dC76, Section 3.3]. If the *z* axis is aligned with the normal, the linear terms of Equation (3) vanish, and the second fundamental form reduces to the Hessian of the height function. Further α implifications are obtained in the Monge coordinate system, where $I = Id_2$, the Hessian is a diagonal matrix, and the principal curvatures are given by $2B_{20}$ and $2B_{02}$.

3.2 Polynomial fitting of the height function

We begin by an approximation result on the coefficients of the height function. We focus on the convergence rate given by the value of the exponent of parameter *h*.

Proposition. 1 *A poised polynomial interpolation or a polynomial approximation of degree n based* upon N points $p_i(x_i, y_i, z_i)$ whose abscissa are $x_i = O(h), y_i = O(h)$ estimates the coefficients of *degree k of the Taylor expansion of f to accuracy* $O(h^{n-k+1})$:

$$
A_{k-j,j} = B_{k-j,j} + O(h^{n-k+1}) \quad \forall k = 0, ..., n \ \forall j = 0, ..., k.
$$

Moreover, if the origin is one of the p_i s and interpolation is used, then $A_{0,0} = B_{0,0} = 0$.

Proof. [Proof of Prop. 1, interpolation case.] Using the notations introduced in section 1.3, we aim at solving the linear system $MA = B$, with $D = det M \neq 0$ since we assume the interpolation problem is poised. We apply Cramer's rule to solve it:

$$
A_{k-j,j} = \det(1, x_i, y_i, \dots, x_i^{k-j+1} y_i^{j-1}, J_{B,n}(x_i, y_i) + O(||(x_i, y_i)||^{n+1}), x_i^{k-j-1} y_i^{j+1}, \dots, y_i^{n})|_{i=1,\dots,N}/D.
$$

By linear combination of the columns of the numerator, all monomials of $J_{B,n}(x_i, y_i)$ disappear except $B_{k-j, j} x_i^{k-j} y_i^j$:

$$
A_{k-j,j} = \det(1, x_i, y_i, \dots, x_i^{k-j+1} y_i^{j-1}, B_{k-j,j} x_i^{k-j} y_i^j + O(||(x_i, y_i)||^{n+1}), x_i^{k-j-1} y_i^{j+1}, \dots, y_i^n))_{i=1,\dots,N}/D
$$

Eventually, splitting the numerator yields

 \sim

$$
A_{k-j,j} = B_{k-j,j} + O(h^{n-k+1}).
$$

For the particular case where the origin is one of the samples, notice that the equation involving point *p* is $A_{0,0} = f(0,0) = 0$. \square

Proof. [*Proof of Prop. 1, approximation case.]* Consider the least-square system $\min |MA - B||_2$. With the assumption than M is of rank N_n , the approximation is equivalent to the inversible linear system $^tMMA = ^tMB$. With the notation $\Sigma_l^k = \sum_{i=1}^N x_i^k y_i^l$ and the assumption $||(x_i, y_i)|| = O(h)$:

$$
{}^{t}MB = \begin{pmatrix} B_{0,0}\Sigma_{0}^{0} + B_{1,0}\Sigma_{0}^{1} + B_{0,1}\Sigma_{1}^{0} + \dots + B_{0,n}\Sigma_{n}^{0} + O(h^{n+1}) \\ B_{0,0}\Sigma_{0}^{1} + \dots + B_{0,n}\Sigma_{n}^{1} + O(h^{n+2}) \\ \vdots \\ B_{0,0}\Sigma_{n}^{0} + \dots + O(h^{n+1}) \end{pmatrix}
$$

$$
{}^{t}MM = \begin{pmatrix} \Sigma_{0}^{0} & \Sigma_{0}^{1} & \Sigma_{1}^{0} & \dots & \Sigma_{n}^{0} \\ \Sigma_{0}^{1} & \Sigma_{0}^{2} & \Sigma_{1}^{1} & \dots & \Sigma_{n}^{1} \\ \Sigma_{1}^{0} & \Sigma_{1}^{1} & \Sigma_{2}^{0} & \dots & \Sigma_{n+1}^{0} \\ \vdots & \vdots & \vdots & \vdots & \vdots \\ \Sigma_{n}^{0} & \Sigma_{n}^{1} & \Sigma_{n}^{0} & \dots & \Sigma_{2n}^{0} \end{pmatrix}
$$

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 Let *D* be $det({}^{t}MM)$, applying the Cramer's rule gives:

$$
A_{j,k-j} = \det \begin{pmatrix} \Sigma_0^0 & \cdots & B_{0,0} \Sigma_0^0 + B_{1,0} \Sigma_0^1 + B_{0,1} \Sigma_1^0 + \cdots + B_{0,n} \Sigma_n^0 + O(h^{n+1}) & \cdots & \Sigma_n^0 \\ \Sigma_0^1 & \cdots & B_{0,0} \Sigma_0^1 + \cdots + B_{0,n} \Sigma_n^1 + O(h^{n+2}) & \cdots & \Sigma_n^1 \\ \Sigma_1^0 & \cdots & B_{0,0} \Sigma_1^0 + \cdots + O(h^{n+2}) & \cdots & \Sigma_{n+1}^0 \\ \vdots & \vdots & & \vdots & \vdots \\ \Sigma_n^0 & \cdots & B_{0,0} \Sigma_n^0 + \cdots + O(h^{n+n+1}) & \cdots & \Sigma_{2n}^0 \end{pmatrix} / D
$$

Linear combinations of the columns gives:

$$
A_{j,k-j} = \det \begin{pmatrix} \Sigma_0^0 & \cdots & B_{j,k-j} \Sigma^{j,k-j} + O(h^{n+1}) & \cdots & \Sigma_n^0 \\ \Sigma_0^1 & \cdots & B_{j,k-j} \Sigma^{j+1,k-j} + O(h^{n+2}) & \cdots & \Sigma_n^1 \\ \Sigma_1^0 & \cdots & B_{j,k-j} \Sigma^{j,k-j+1} + O(h^{n+2}) & \cdots & \Sigma_{n+1}^0 \\ \vdots & \vdots & \vdots & \vdots & \vdots \\ \Sigma_n^0 & \cdots & B_{j,k-j} \Sigma^{j,k-j+n} + O(h^{n+n+1}) & \cdots & \Sigma_{2n}^0 \end{pmatrix} / D
$$

The numerator of this formula splits due to the multi-linearity of the determinant, and noticing that $\Sigma_{l}^{k} = O(h^{k+l})$ gives $A_{j,k-j} = B_{j,k-j} + O(h^{n-k+1})$. \square

Using the previous proposition, the order of accuracy of a differential quantity is linked to the degree of the interpolant and the order of this quantity. More precisely:

Theorem. 3 *A polynomial fitting of degree n estimates any k th -order differential quantity to accuracy* $O(h^{n-k+1})$ *. In particular:*

- the coefficients of the unit normal vector are estimated with accuracy $O(h^n)$, and so is the *angle between the normal and the estimated normal.*
- *• the coefficients of the first, second fundamental form and shape operator are approximated* with accuracy $O(h^{n-1})$, and so are the principal curvatures and directions.

Proof. It is easily checked that the formula corresponding to the geometric quantities are C^1 functions of the coefficients. The result follows from proposition 1 and lemma 1 —see appendix 8. \Box

The previous theorem generalizes [MW00a, Lemma 4.1] where 2-jet interpolations only are studied. The $O(h^n)$ bound on the normal should also be compared to the normal estimate of the normal vector using specific Voronoi centers called poles considered in [AB99]. The error bound proved there is equivalent to 2ε with ε the sampling density of the surface. Setting $h = \varepsilon$ lfs and assuming lfs is bounded from above, the estimation stemming from a polynimial fitting therefore yields more accurate results for the tangent plane, and also provides information on higher order quantities. From an algorithmic perspective, and according to proposition 2, it is sufficient to perform the fitting in any coordinate system.

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3.3 Accuracy of the fitting and relationship to the coordinate system

Proposition 1 involves any admissible coordinate system. From a practical standpoint, we would like to perform the fitting in some arbitrary coordinate system, compute the normal and principal directions, and infer the height function in the Monge coordinate system. Notice that this is not completely straightforward since rewriting a height function in a different coordinate system results in an implicit equation for the corresponding surface —we had $z = f(x, y)$, but x, y, z become linear combinations of the new variables x' , y' , z' . The following proposition shows that calculations can indeed be performed in this order, and that no accuracy is lost along the process. The constructive proof uses the implicit function theorem and is omitted.

Proposition. 2 Let $J_{A,n}$ be the n-jet estimating $J_{B,n}$ in a coordinate system D, and let Q be the surface associated with $J_{A,n}.$ Let D be another coordinate system, $J_{B,n}$ ($J_{A,n}$) the n-jet of S (Q) in D . If $A_{k-j,j} = B_{k-j,j} + O(h^{n-k+1})$, then $A'_{k-j,j} = B'_{k-j,j} + O(h^{n-k+1})$.

3.4 Influence of normal accuracy on higher order estimates

Following the guideline initiated in [SZ90] several algorithms first estimate the normal to the surface and then proceed with Hessian of the height function. We analyze the error incurred by the latter as a function of the accuracy on the former. We denote θ the angle between the normal n_S to the surface and the normal n_Q estimated by the two-stages method. In order to simplify calculations, we assume that n_Q is aligned with the *z*-axis and n_S is in the (x, z) -plane, so that $f(x, y) = B_{10}x +$ $B_{20}x^2 + B_{11}xy + B_{02}y^2 + O(||(x, y)||^3)$, with $B_{10} = -\tan\theta$. Expressed in the same coordinate system, the interpolant —a 2-jet to simplify calculations— reads as $J_{A,2}(x, y) = A_{20}x^2 + A_{11}xy + A_{02}y^2$. The proof of the following can be found in appendix 9.

Proposition. 3 *If a small error* θ *is done on the estimated normal, a 2-jet interpolation give the Gauss curvature with a linear error wrt* θ*:*

$$
k_Q - k_S = \theta O(h^{-1}) + O(h) + O(\theta^2).
$$

For a fixed *h*, the curvature error is a linear function of the angle between the normals. The term $\theta O(h^{-1})$ shows that if θ is fixed, the smaller *h* the worse the accuracy. Hence estimating the normal deserves specific care.

4 Plane Curves

All the results proved for surfaces in the previous section can also be proved for curve, and we omit them. Instead, for the interpolation case, we prove an error bound between the coefficients of the curve and those of the jet.

4.1 Problem addressed

Let *C* be a curve, and consider the height function *f* following Equation (1) in any coordinate system whose *y* axis is not tangent to the curve and, whose origin is on the curve (this implies that $B_0 = 0$). We shall fit *C* by a *n*-jet $J_{A,n}(x)$ whose graph is denoted *Q*. As already mentioned, there are $n + 1$ unknown coefficients A_i s, we assume N data points $P_i(x_i = a_i h, y_i = f(x_i))$ are given, where $N = n + 1$ for interpolation fitting. Notice again that parameter *h* specifies the uniform convergence of these data points to the origin. The fitting equations are:

$$
y_i = f(x_i) = J_{B,n}(x_i) + O(x_i^{n+1}) = J_{A,n}(x_i).
$$

Since curve *C* is given by Equation (1), the normal and the curvature of *C* at the origin are given by

$$
n_C = (-B_1, 1)^t / \sqrt{1 + B_1^2}, \ k_C = 2B_2 / (1 + B_1^2)^{\frac{3}{2}}.
$$
 (13)

Moreover, in the Monge coordinate system $-B_1 = 0$, these expressions simplify to $n_C = (0, 1)^t$ and $k_C = 2B_2$.

4.2 Error bounds for the interpolation

The equivalent of Prop. 1 for curves gives the magnitude of the accuracy of the interpolation. We can actually be more precise and provide precise error bounds depending upon the function interpolated as well as the relative position of the sample points used. The proof of the following can be found in appendix 10.

Proposition. 4 Consider a degree n ($n \ge 2$) interpolation problem for a curve $y = f(x)$. Let ε be a positive number so that the interpolation point abscissa lie in the interval $[-\varepsilon,\varepsilon].$ Let c be a positive *constant so that* $\sup_{x \in [-\varepsilon, \varepsilon]} |f^{(n+1)}(x)| \leq c$ \leq *c.* At last, let $d \leq 1$ be defined by $\min_{i \neq j} |x_i - x_j| = 2\varepsilon d/n$. (*d* is a measure of the minimum distance between the interpolation points.) Then for $k = 0, ..., n$:

$$
|A_k - B_k| \le \varepsilon^{(n-k+1)} c \left(\frac{n}{2d}\right)^{\frac{n(n-1)}{2}}.
$$

Here is an application of the previous result. Let θ denote the angle between the normal and the estimated normal. We have $\sin(\theta) = ||n_Q \wedge n_C|| = |A_1 - B_1| / \sqrt{(1 + A_1^2)(1 + B_1^2)} \le |A_1 - B_1|$. It $|A_1 - B_1|$. It is found that

$$
\theta \leq \arcsin(\ \varepsilon^n c\left(\frac{n}{2d}\right)^{\frac{n(n+1)}{2}}).
$$

Therefore, the coordinate system minimizing the error in the worst case is the one with respect to which the pairwise differences between the abscissa of the sample pointsis maximized. Generalizing this result for the curvature involves cumbersome calculations.

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5 Algorithm

The fitting algorithm to estimate the differential properties at a point *p* consists of (i)collecting the points used for the fitting. Recall that a *n*-jet involves $N_n = (n+1)(n+2)/2$ coefficients, so that when interpolating (approximating) we assume $N = N_n$ ($N > N_n$). (ii)solving the linear system (iii)recovering the differential properties. We examine in turn the details of these three steps.

5.1 Collecting *N* **neighbors**

The mesh case. Although no topological information is required by the fitting method, the connectivity information of a mesh can be used as follows. We sequentially visit the one-ring neighbors, two-ring neighbors, and so on until *N* points have been collected. Let R_1, \ldots, R_k be the *k* rings of neighbors necessary to collect N neighbors. All the points of the $k-1$ first rings are used. The complement up to *N* points is chosen arbitrarily out of the *k*th ring.

The point-cloud case. The normal at *p* is first estimated, and the neighbors of *p* are further retrieved from a power diagram in the estimated tangent plane [BF02] —a provably good procedure if the samples are dense enough. If the number of neighbors collected is less than *N*, we recursively collect the neighbors of the neighbors.

Collecting the points therefore boils down to estimating the tangent plane. One solution is to construct the Voronoi diagram of the point set and use these Voronoi vertices called *poles* [AB99, FR02]. Poles yield an accurate estimate of the normal vector but require a global construction.

An alternative is to resort to the algorithm of section 3, and solve a degree one interpolation problem —which requires three points and is well poised as soon as the three points are not collinear. Geometrically, the closer the three points to being aligned, the more unstable the tangent plane estimate. To see how one can get around this difficulty, denote *q* the nearest neighbor of *p*. Also, let *r* be the sample point so that the circum-radius r_{circ} of the triangle *pqr* is minimum. The estimated normal at *p* is the normal to the plane through *pqr*. Intuitively, minimizing the circum-radius *r circ* prevents two difficulties: on one hand triangles with a large angle (near to π) exhibit a large circum-circle and are discarded; on the other hand, triangles involving a third point *r* which is not a *local* neighbor of *p* cannot minimize *r circ* and are also discarded. A more formal argument advocating the choice of the triangle with minimum *r circ* is provided in [She02], where it is shown that the worst error on the approximation of the gradient of a bivariate function by a linear interpolant precisely involves r_{circ} .

5.2 Solving the fitting problem

The next stage consists of choosing the *z* direction to work with. Since the tangent plane has not been estimated, we use a principal component analysis to compute a rough estimate of the normal with the neighboring points. The polynomial fitting can be done in any coordinate system whose z axis is not tangent to the surface. Hence at least one of the three axis of the world coordinate system matches this requirement. A natural choice is to select the coordinate axis whose angle with the rough estimated normal is minimum. For these coordinates, we fill the Vandermonde matrix. The

matrix is further scaled as explained in section 2.3, with *h* the average value of the norms $|| (x_i, y_i) ||$. The corresponding system is solved using a Singular Value Decomposition. Practically, we use the SVD of the Gnu Scientific Library, available from http://sources.redhat.com/gsl.

As pointed out in section 2.3, the instability of the system is provided by the condition number. Whenever degenerate configurations are detected, one can proceed as follows. For the approximation strategy, one can either keep the same degree and increase the number of points used, or reuse the same points with a lower degree. These changes are likely to provide a non singular matrix *M*. In the worst-case, a degree one fitting must be possible since then only three linearly independent points are required! For the interpolation, things are a bit more involved since reducing the interpolation degree requires discarding some points. Selecting the subset yielding the best conditioning is a challenging problem [Las99, Hig96]. Notice also that for the approximation case, one can always retrieve a solution from an under-constrained least-square problem by choosing, e.g., the solution vector of least norm.

5.3 Retrieving differential quantities

We have already mentioned how to compute the normal. For the second order information, we compute the Weingarten map of the surface [dC76, Section 3.3]. Its eigenvalues (eigenvectors) provide the principal curvatures (directions) of the surface. For a parameterized surface given as a height function, one ends up with the formula given on Table 1. Notice that a basis of the tangent space associated to the parameterization $X(u, v) = (u, v, h(u, v))$ consists of the two vectors $X_u = (1, 0, h_u)^t$ and $X_v = (0, 1, h_v)^t$. A Gram-Schmidt orthonormalization of the basis $\{X_u, X_v\}$ gives another basis *Y*, *Z*} of the tangent space. The diagonalization of the symmetric matrix representing the Weingarten map in the basis $\{Y, Z\}$ provides the expression of the principal curvature directions with respect to the ${Y, Z}$ orthonormal basis. Note that the sign of principal curvatures and hence the definition of minimal and maximal directions rely on the orientation of the normal. As long as our experimental study is performed on meshes of oriented surfaces, it is straightforward to find a global and coherent orientation of the normals.

$$
E = 1 + a_1^2
$$

\n
$$
F = a_2 a_1
$$

\n
$$
G = 1 + a_2^2
$$

\n
$$
e = \frac{2a_3}{\sqrt{a_1^2 + 1 + a_2^2}}
$$

\n
$$
f = \frac{a_4}{\sqrt{a_1^2 + 1 + a_2^2}}
$$

\n
$$
A^t = -\begin{pmatrix} e & f \\ f & g \end{pmatrix} \begin{pmatrix} E & F \\ F & G \end{pmatrix}^{-1}
$$

\n
$$
B = \frac{2a_5}{\sqrt{a_1^2 + 1 + a_2^2}}
$$

Table 1: Computing the matrix A of the Weingarten map of $h(u, v) = a_1 u + a_2 v + a_3 u^2 + a_4 u v + a_5 v^2$ in the basis $\{X_u, X_v\}$

6 Experimental study

We present results along two lines. First, we illustrate the convergence theorems proved in section 3. Second, we present illustrations on standard computer graphics models.

6.1 Convergence estimates on a graph

Setup. We illustrate the convergence properties with the smooth height fields $f(u, v) = 0.1e^{2u+v-v^2}$ and $g(u, v) = 4u^2 + 2v^2$ defined over the parametric domain $(u, v) \in [-0.5, 0.5]^2$ —see Figs. 4 and 5. At selected points on the graphs of *f* and *g*, we study the angle between the normals —more precisely its sine $sin(n, \tilde{n})$, and the relative errors on principal curvatures. The values output by our algorithm are compared against the exact values computed analytically with arbitrary precision under Maple, and we report both average and maximum errors over samples points. More precisely, the graph of f or g are sampled with points $p_i(x_i, y_i, f(x_i, y_i))$ where the (x_i, y_i) s lies on a randomly perturbed triangulated square grid of side *h*. The triangulation is randomly perturbed to avoid simple degenarate configurations such as points lying on lines. The perturbation for the point (u, v) of the regular grid is the point (x, y) with $x = u + \delta h$, $y = v + \delta' h$ and δ, δ' random numbers in [0,0.9]. The connectivity of the graph is that of the triangulated grid. Notice also that since calculations are carried out on a surface patch parameterized over a square grid, the direction choosen for the polynomial fitting is the *z* direction near the origin, and either the the *x* or *y* directions at the periphery of the domain.

The convergence properties are illustrated (i)with respect to the discretization step *h* of the grid —for a given fitting degree *n* (ii)with respect to the fitting degree *n* —for a given discretization step *h*. We compare the convergence properties of the interpolation and approximation schemes, for fitting degrees ranging from one to nine. To quantify the observations, notice that according to theorem 3, the error δ on a k^{th} -order differential quantity is $O(h^{n-k+1})$, hence

$$
\delta \approx c \, h^{n-k+1} \Leftrightarrow \log(1/\delta) \approx \log(1/c) + (n - k + 1)\log(1/h) \tag{14}
$$

$$
\Leftrightarrow \frac{\log(1/\delta)}{\log(1/h)} \approx \frac{\log(1/c)}{\log(1/h)} + (n - k + 1). \tag{15}
$$

Convergence wrt to *h***.** To highlight the convergence properties with respect to the size of the neighborhood, we consider sequences of meshes with $h \to 0$, more precisely *h* ranges from 2^{-2} to 2^{-6} . The results for *f* and *g* being alike, we focus on the exponential *f*.

Curves of Figs. 6 to 11 show the average convergence behavior as the size of the neighborhood decreases. For a given degree and following equation (14), curves of figures 6 to 11 should be lines of slope $(n - k + 1)$. The behavior is more regular for the approximation case, and the estimate is also better: a gain of about a digit can be observed between *kmax* estimates of figures 10 and 11.

Convergence wrt to the interpolation degree. For the convergence wrt to the interpolation degree —with *h* fixed, we present results for the polynomial *g*. Conclusions are similar for *f* , but it should however be noticed that since the graph of *g* is more curvy than that of *f* , a finer grid is required. (The higher the degree the more points required by the fitting ... and we compute local quantities!) To be precise, we ran experiments with $h = 2^{-5}$ for f and $h = 2^{-7}$ for g.

Curves of figures 12 and 13 show the convergence as a function of the degree of the fitted polynomial for a fixed neighborhood size. According to Eq. (15), curves of these figures should be lines of unit slope, with a vertical shift of one unit between normal and curvatures errors since curvature is a 2^{nd} order quantity whereas normal is 1^{st} order. The gap between the average values and the maximal values is greater for interpolation than for approximation. The other charts provide the conditioning and the least singular value. Interpolation fitting is allways more ill-conditioned than approximation, and closer to a degenerate problem (the least singular value is the distance of the matrix system to singular matrices). The particular case of a degree 7 approximation reveals to be badly conditioned due to the regular connectivity of the mesh used to find the neighbors: there is only one more point than for the degree 7 interpolation fitting.

6.2 Illustrations

We depict differential informations on several models. When principal directions are displayed, blue and red respectively correspond to k_{min} and k_{max} —that is $k_{min} \le k_{max}$ —, assuming the surface normal points to the outside. To display patches of osculating *n*-jets, it is sufficient to select a rectangular domain in parameter space, sample it with a grid, and plot the corresponding mesh.

Consider the mesh models of the elliptic paraboloid $z = 2x^2 + y^2$ —16k points, Fig. 18—, and the surface of revolution $z = 0.1 \sin(10\sqrt{(x^2 + y^2)})$ —8k points, Fig. 19. The arrangement of curvature lines provides informations on umbilical points —where principal directions are not defined since $k_{min} = k_{max}$. On the paraboloid, it is easy to follow curvature lines and see how they turn around an umbilic. The surface of revolution provides an example of two parabolic lines (where the principal curvature k_{max} vanishes), that is a curve along which the Gauss curvature K_{Gauss} vanishes. This specific line splits the surface into elliptic ($K_{Gauss} > 0$) and hyperbolic regions ($K_{Gauss} < 0$). This model also illustrates a line of umbilical points where minimum and maximum principal directions swap each over.

For a standard example from Computer Graphics, consider the Michelangelo's David of Fig. 20. On this model of 95922 pts, the principal curvatures provide meaningful information for shape perception ⁶. To finish up, we illustrate the robustness of the method. Figure 21 displays random patches on the Mechanic model, a 12,500 points model reconstructed from the output of a range scanner. In spite of the coarse sampling, patches and principal directions provide faithful information. In a similar vein, approximation fitting with large neighborhoods Fig. 22 features a noisy triangulation of a graph. In spite of the severe level of noise, surface patches *average* the available information. On Fig. 23, a noisy triangulation of an ellisoid, 15k points, principal directions are enough precise to recognize an umbilic.

⁶See also [HCV52, p197] as well as [HGY⁺99].

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7 Conclusion

Estimating differential quantities is of prime importance in many applications from Computer Vision, Computer Graphics, Computer Aided Design or Computational Geometry. This importance accounts for the many different differential estimators one can find in the vast literature of applied geometry. Unfortunately, few of these have undergone a precise theoretical analysis. Another striking fact is that estimates of second order differential quantities are always computed using degenerate conics/quadrics without even mentioning the classification of Euclidean conics/quadrics.

The main contribution of the paper is to bridge the gap between the question of estimating differential properties of arbitrary order and multivariate interpolation and approximation. In making this connection, the use of jets —truncated Taylor expansions— is advocated. Precise asymptotic convergence rates are proved for curves and surfaces, both for the interpolation and approximation schemes. To the best of our knowledge, these results are among the first ones providing accurate estimates for differential quantities of order three and more. Experimental results for surfaces of \mathbb{R}^3 are reported. These experimentsillustrate the asymptotic convergence results, but also the robustness of the methods on general Computer Graphics models.

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8 Appendix: a useful lemma

In several occasions we derive differential quantities —unit normal vectors, curvatures— as a function *F* of the coefficients of the jet. The following lemma makes it easy to derive the precision on the quantity investigated if F is regular enough and if the some precision of the jet's coefficients is known.

Lemma. 1 Define a kth-order differential quantity —for a curve or a surface— as a C¹ function F of the of the coefficients of the k-jet of the height function. Also assume that a degree n fitting yields a precision $A_j = B_j + O(h^{n+1-j}), k = 1,...,n$ for a curve, and $A_{k-j,j} = B_{k-j,j} + O(h^{n-k+1}), k$ $1, \ldots, n, j = 0, \ldots, k$ *for a surface.*

Then, a polynomial fitting of degree n estimates a k th -order differential quantity to accuracy $O(h^{n-k+1}).$

Proof. The proof is the same for curves and surfaces, and we use the notations corresponding to a curve.

To begin with, we perform a substitution on the error terms in the expression of *F*:

$$
F((A_i)_{i=1,\ldots,k}) = F((B_i + O(h^{n-i+1}))_{i=1,\ldots,k}) = F((B_i + O(h^{n-k+1}))_{i=1,\ldots,k}).
$$

Since $F((B_i)_{i=1,\ldots,k})$ is a C^1 function, and denoting DF_p the differential of F at point p , an order one Taylor formula yields:

$$
F((B_i + O(h^{n-k+1}))_{i=1,\ldots,k}) = F((B_i)_{i=1,\ldots,k}) + DF_{(B_i + uO(h^{n-k+1}))_{i=1,\ldots,k}} (O(h^{n-k+1}), \ldots, O(h^{n-k+1}))^t, u \in]0,1[
$$

= $F((B_i)_{i=1,\ldots,k}) + O(h^{n-k+1}).$

9 Appendix to *Surfaces*

Proof. [Prop. 3] The system of equations for the interpolation is:

$$
A_{20}x_i^2h + A_{11}x_iy_i + A_{02}y_i^2 = B_{10}x_i + B_{20}x_i^2 + B_{11}x_iy_i + B_{02}y_i^2 + O(|(x_i, y_i)|^3) \quad i = 1, ..., 3.
$$

Let *D* be the determinant: $D = \det(x_i^2, x_i y_i, y_i^2)_{i=1,\dots,3}$, Cramer's rule gives:

$$
A_{20} = \det(B_{10}x_i + B_{20}x_i^2 + B_{11}x_iy_i + B_{02}y_i^2 + O(|(x_i, y_i)|^3), x_iy_i, y_i^2)_{i=1,\dots,3}/D
$$

= $B_{10}O(h^{-1}) + B_{20} + O(h).$

Similar calculations gives: $A_{11} = B_{10}O(h^{-1}) + B_{11} + O(h)$, and $A_{02} = B_{10}O(h^{-1}) + B_{02} + O(h)$. The Gaussian curvature of *Q* is then:

$$
k_Q = 4A_{20}A_{02} - A_{11}^2
$$

= $4(B_{10}O(h^{-1}) + B_{20} + O(h))(B_{10}O(h^{-1}) + B_{02} + O(h)) - (B_{10}O(h^{-1}) + B_{11} + O(h))^2$
= $4B_{20}B_{02} - B_{11}^2 + B_{10}O(h^{-1}) + B_{10}^2O(h^{-2}) + O(h)$
= $4B_{20}B_{02} - B_{11}^2 + \tan\theta O(h^{-1}) + \tan^2\theta O(h^{-2}) + O(h)$.

The Gaussian curvature of *S* is

$$
k_S = \frac{4B_{20}B_{02} - B_{11}^2}{(1 + B_{10}^2)^2} = \frac{4B_{20}B_{02} - B_{11}^2}{(1 + \tan^2 \theta)^2} = (4B_{20}B_{02} - B_{11}^2)\cos^4 \theta.
$$

Thus the error on the curvature is:

$$
k_Q - k_S = (4B_{20}B_{02} - B_{11}^2)\cos^4\theta + \tan\theta O(h^{-1}) + \tan^2\theta O(h^{-2}) + O(h)
$$

= $\theta O(h^{-1}) + O(h) + O(\theta^2)$.

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10 Appendix to *Curves*

Proof. [Proof of Prop. 4] The interpolation system leads to the $n + 1$ equations:

$$
J_{A,n}(x_i) = J_{B,n}(x_i) + \frac{c_i x_i^{n+1}}{(n+1)!}
$$

where $\forall i = 0, ..., n |c_i| \leq c$ $\leq c$. The usual calculations yield

$$
A_k - B_k = \det(1, x_i, \dots, x_i^{k-1}, \frac{c_i x_i^{n+1}}{(n+1)!}, x_i^{k+1}, \dots, x_i^n)_{i=0,\dots,n} / \det(1, x_i, x_i^2, x_i^3, \dots, x_i^n)_{i=0,\dots,n}
$$

In order to find an upper bound of the numerator, we use the inequality:

$$
|\det(a_{ij})_{i,j=0,...,n}| = |\sum_{\sigma \in S_{n+1}} \varepsilon(\sigma) \prod_{i=0}^{n} a_{i\sigma(i)}| \leq \sum_{\sigma \in S_{n+1}} \prod_{i=0}^{n} |a_{i\sigma(i)}| \leq (n+1)! \max_{\sigma \in S_{n+1}} \prod_{i=0}^{n} |a_{i\sigma(i)}|
$$

With $|x_i|$: $\leq \varepsilon$ and $|c_i| \leq c$ $\leq c$, this leads to:

$$
|\det(1,\ldots,x_i^{k-1},\frac{c_ix_i^{n+1}}{(n+1)!},x_i^{k+1},\ldots,x_i^n)_{i=0,\ldots,n}|\leq c\epsilon^{\frac{n(n+1)}{2}+(n-k+1)}
$$

The denominator can be computed with the Vandermonde formula:

$$
|\det(1, x_i, x_i^2, \ldots, x_i^n)_{i=0,\ldots,n}| = \prod_{j>i} |x_j - x_i| \ge \prod_{j>i} \frac{2\epsilon d}{n} \ge \left(\frac{2\epsilon d}{n}\right)^{\frac{n(n+1)}{2}}.
$$

The result is then the ratio of these bounds. \Box

11 Appendix to *Experimental study*

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