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Computing Nearest Gcd with Certification

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Abstract

A bisection method, based on exclusion and inclusion tests, is used to address the nearest univariate gcd problem formulated as a bivariate real minimization problem of a rational fraction.

The paper presents an algorithm, a first implementation and a complexity analysis relying on Smale's α -theory. We report its behavior on an illustrative example.

1 Introduction

1.1 A minimization problem

As usual, we denote by $\mathbf{R}_d[z]$ and $\mathbf{C}_d[z]$ the vector spaces of univariate polynomials of degree less or equal to d , with coefficients in \mathbf{R} and \mathbf{C} . Let $\mathbf{e}(z) = (e_0(z), \dots, e_d(z))^T$ be a polynomial basis of $\mathbf{C}_d[z]$. We denote by $f = (f_0, \dots, f_d)^T$ the coefficients of the polynomial $f(z) = \sum_{k=0}^d f_k e_k(z) \in \mathbf{R}_d[z]$. The 2-norm of f is $\|f\|^2 = \sum_{k=0}^d f_k \bar{f}_k$.

We address the nearest gcd problem, i.e. the following minimization problem:

Given two degree d polynomials $f(z)$ and $g(z)$, find two degree d polynomials $p(z)$ and $q(z)$ with a non trivial gcd, which are solutions of the minimization problem

$$\min_{\substack{p, q \in \mathbf{C}_d[z] \\ \text{Resultant}(p, q) = 0}} \|f - p\|^2 + \|g - q\|^2. \quad (1)$$

N.K. Karmarkar and Y.N. Lakshman in [14] reduced that problem to another minimization problem :

$$\min_{z \in \mathbf{C}} \frac{f(z)\overline{f(z)} + g(z)\overline{g(z)}}{\sum_{k=0}^d e_k(z)\overline{e_k(z)}}. \quad (2)$$

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Moreover, a global minimum z_0 of (2) is a root of a nearest gcd and determines the polynomials p and q of (1) by the formulas:

$$\begin{aligned} p(z) &= f(z) - \frac{f(z_0)}{\mathbf{e}(z_0)^T \mathbf{e}(z_0)} \overline{\mathbf{e}(z_0)}^T \mathbf{e}(z) \\ q(z) &= g(z) - \frac{g(z_0)}{\mathbf{e}(z_0)^T \mathbf{e}(z_0)} \overline{\mathbf{e}(z_0)}^T \mathbf{e}(z) \end{aligned}$$

with $\mathbf{e}(z) = (e_0(z), \dots, e_d(z))^T$.

We observe that with the change of variables $z = X + iY$, the problem amounts to minimize a rational bivariate function

$$F(X, Y) := \frac{N(X, Y)}{D(X, Y)}.$$

So we focus on the resolution of this task.

1.2 Our approach

To compute the global minima of $F(X, Y)$, we propose a bisection algorithm, based on an exclusion test and an inclusion test (see below), applied simultaneously to :

- 1- the polynomial $P_\mu(X, Y) = N(X, Y) - \mu D(X, Y)$
- 2- the system $G(X, Y) = (G_1(X, Y), G_2(X, Y))$ of the numerators of the gradient $\nabla F(X, Y)$.

We will also use the following notations.

- 1- $A_F := \{(\bar{x}_i, \bar{y}_i) \mid i = 1, \dots, N\}$ is the set of global minima of $F(X, Y)$.
- 2- Z_H is the set of zeroes of a polynomial function $H(X, Y)$ and $d(x, y, Z_H)$ is the Euclidean distance from $(x, y) \in \mathbf{R}^2$ to Z_H . The function $H(X, Y)$ will be either $P_\mu(X, Y)$ or $G_i(X, Y)$, $i = 1, 2$.
- 3- S or $S(x, y, r)$ is a square centered at (x, y) and of radius r .

Our bisection algorithm will iteratively update a list of retained squares and a list of approximate global minima of F . We prove a complexity result in Section 3.

We follow the approach initiated by S. Smale and his co-worker in a series of papers (see e.g. [5] and the references therein) relying on their celebrated α and γ -theorems.

In section 2 we will briefly review some of these notions, recall the definitions of α, β, γ , we also set $\gamma(G, A_F) := \max_{(x, y) \in A_F} \gamma(G; x, y)$. After that, we provide a precise quantitative definition for a point (x, y) to be, in our setting, an approximate global minimum of $F(x, y)$. Then thanks to the γ -theorem, the approximation is sufficiently good to imply that a Newton iteration converges quadratically towards a global minimum.

Theorem 1.1 *Assume that all the elements of A_F are regular points of the system $G(x, y) = 0$ and are contained in some square $S_0 := S(x_0, y_0, r_0)$.*

Let $C_d := 1 + \frac{\sqrt{2}}{\log 2}(4d - 2)$ and $J := \left\lceil \log_2 \frac{3r_0 C_d \gamma(G, A_F)}{K \delta_0} \right\rceil$, where δ_0 is the smallest positive root of $u - \frac{13 - 3\sqrt{17}}{4}(2u^2 - 4u + 1)^2$.

Denote by R_j the set of retained squares at step j . Then, for $j \geq J$,

- 1- *All the points of R_j are approximate global minima of $F(x, y)$. That is to say the Newton iteration applied to G from a point in R_j converges quadratically to a global minimum of F .*
- 2- *The number of exclusion tests is bounded by $1 + j N K^2$.*

3- When d tends to infinity J belongs to

$$O\left(\frac{d \log(d^3)}{\min_{(x,y) \in A_F} d_F(DG(x,y)^{-1}, \Sigma)} + \log K\right),$$

where d_F is the Frobenius distance and Σ is the set of singular matrices.

When Problem 1 has a unique solution the number of exclusions steps belongs to

$$O\left(\frac{N K^2 d^3 \log^3(d)}{\min_{(x,y) \in A_F} d_F(DG(x,y)^{-1}, \Sigma)} + N K^2 d^3 \log K\right).$$

where

- $V_\epsilon(G)$ is the tubular neighborhood of the zero set $G(x,y) = 0$;
- N is the maximal number of connected components of $U_\epsilon := V_\epsilon(G_1) \cap V_\epsilon(G_2)$, for all $\epsilon > 0$;
- for all $\epsilon > 0$, ϵK is bounding the radius of a connected component of U_ϵ .

Remark:

1. $\min_{(x,y) \in A_F} d_F(DG(x,y)^{-1}, \Sigma)$ is a constant but this gives information about the condition number of the system. That is why we keep this constant in the big O notation.
2. $\#A_F \leq \sum_{(p,q) \text{ solution of Problem 1}} \deg gcd(p,q)$.
Thus if Problem 1 has a unique solution then $\#A_F \leq \deg gcd(p,q) \leq d$.
3. In this paper, in order to apply α and γ -theorems, we suppose that all the elements of A_F are regular points of the system $G(x,y) = 0$. We can avoid this hypothesis, then instead of α and γ -theorems we have to use the tools developed in [11].

1.3 Other approaches in the literature

The nearest GCD problem has been studied with different approaches and with other formulations by many authors.

1.3.1 Algebraic approach via Euclid's algorithm, resultant and subresultant

The first papers, see for example Brown [6], on the complexity of Euclid's algorithm only works with exact coefficients. Then the numerical case has been successively considered by Schönhage [21], Noda and Sasaki [18], Corless, Gianni, Trager and Watt [7], Hribernic and Stetter [12], Emiris, Galligo and Lombardi [10], Beckermann, Labahn [2]. These authors consider the so-called near GCD or the ϵ -GCD problem. The singular value decomposition was first applied to the Sylvester matrix, in [7], and later applied to the subresultant matrices, in [9, 10] in order to get a certification a condition (depending on the level of accuracy) is satisfied and an efficient implementation [20].

1.3.2 Padé approximation and structured matrices approach

See the book Bini and Pan [3] and bibliography within, and more recently Boito, Bini [4].

1.3.3 Rootfinding and cluster root approach

Pan [19] and, more recently, Zeng [23] use root finding and least squares methods.

1.3.4 Optimization approach

The resolution of problem 2 is the main and the most "time-consuming" step of the method propose in [14] that we aim to improve. The authors rely on techniques from Arnon-McCallum [1] and Manocha-Demmel [15]. The second paper is based on resultant for expressing the intersection of two curves and on numerical computation of eigenvalues and eigenvectors by QR iterations. Therefore the expected running time is in $O(p^3)$ where p is the product of degrees of two curves, hence the complexity of the algorithm is at least in $O(d^6)$.

Kaltofen and his co-workers [13] determine approximate GCDs from methods based on structured total least square (STLN). The STLN is an iterative method of the family of Gauss-Newton methods. The authors describe its application to the case of the Sylvester matrix associated to the input polynomials f and g , then show its interest and efficiency by producing the results of experiments. However, since the starting point of their Gauss-Newton like method is not precised, the method can diverge. This is an important drawback.

More recently, Nie-Demmel-Gu [17] use a sum of squares (SOS) technique, the resolution relies on semi definite programming (SDP). For the nearest GCD problem, this yields linear matrix inequalities (LMI) whose size s is $O((d+2)(d+1))$ whose complexity of resolution by projective algorithm is in $O(s^3)$ [16]. So, one ends up with a complexity in $O(d^6)$.

2 The proposed bisection method

2.1 Principles

We propose a bisection method, described below, to approximate the global minima of the function $F(X, Y)$ defined in the introduction. *In the next section, a procedure for computing an initial square is given* ; so we suppose here that a square $S(x_0, y_0, r_0)$ containing all the global minima is known.

The bisection method is based on an exclusion test and an inclusion test. An exclusion test, denoted hereafter by $E(F, S)$ or $E^+(F, S)$, is defined on the set of squares and returns true if the function F has no zero in the square S or false if it might have a zero. Examples of exclusion tests are provided in section 3.

An inclusion test $I(G, S)$ is needed to numerically prove the existence of a local minimum: it takes $G = (G_1, G_2)$ and a square S then returns true if there exists a ball $B(x^*, y^*, r^*)$, containing S , which contains one zero of G , in that case it also returns (x^*, y^*, r^*) , otherwise it returns false.

Our definition of approximate minima is based on γ -theorem and α -theorem of Smale [22], [5], [11], applied to the system $G(x, y)$. We first introduce some quantities and the corresponding classical notations:

- $\beta := \beta(G; x, y) = \|DG(x, y)^{-1}G(x, y)\|$
- $\gamma := \gamma(G; x, y) = \sup_{k \geq 2} \left(\frac{1}{k!} \|DG(x, y)^{-1}D^k G(x, y)\| \right)^{1/(k-1)}$
- $\alpha := \alpha(G; x, y) = \beta\gamma$.
- $\gamma(G; A_F) = \max_{(x, y) \in A_F} \gamma(G; x, y)$.

Smale's γ -theorem states:

Theorem 2.1 *Let (x^*, y^*) be a zero of G and suppose that $DG(x^*, y^*)$ is invertible. If*

$$\|(x, y) - (x^*, y^*)\| \gamma(G, x^*, y^*) \leq \frac{3 - \sqrt{7}}{2}$$

then the Newton iteration from (x, y) converges quadratically to (x^, y^*) .*

This leads to the following definition, motivated by the γ -theorem [5]:

Definition 2.2 *The point (x, y) is an approximate global minimum of $F(x, y)$ if*

$$d((x, y), A_F) \leq \frac{3 - \sqrt{7}}{2\gamma(G, A_F)}.$$

Under this condition, the γ -theorem asserts that the Newton iteration from any point in the ball $B(x^*, y^*, \frac{3 - \sqrt{7}}{2\gamma(G, A_F)})$ where $(x^*, y^*) \in A_F$, converges quadratically towards (x^*, y^*) .

The following result gives a sufficient condition for a point to be a good starting point for the Newton iteration. This result is called α -theorem.

Theorem 2.3 If $\alpha < (13 - 3\sqrt{17})/4$ then G has one and only one zero in the open ball $B(x, y, \sigma(x, y))$ with

$$\sigma(x, y) = \frac{1 + \alpha - \sqrt{1 - 6\alpha + \alpha^2}}{4\gamma} \leq \frac{2 - \sqrt{2}}{2\gamma}.$$

and the Newton iteration from (x, y) converges to (x^*, y^*) . Furthermore, we have

$$\|(x^*, y^*) - N_G^k(x, y)\| \leq \frac{5 - \sqrt{17}}{4\gamma} \left(\frac{1}{2}\right)^{2^k - 1},$$

where $N_G^k(x, y)$ is the k -th iterate of the Newton iteration applied to G with starting point (x, y) .

The following result gives a radius r^* of a ball centered in (x^*, y^*) such that every point in $B(x^*, y^*, r^*)$ satisfy the hypothesis of the α -theorem.

Proposition 2.4 Let (x^*, y^*) be a zero of G and suppose that $DG(x^*, y^*)$ is invertible. Let δ_0 be the smallest positive root of $u - \frac{13 - 3\sqrt{17}}{4}\Psi(u)^2$, where $\Psi(u) = 2u^2 - 4u + 1$. If

$$\|(x, y) - (x^*, y^*)\|\gamma(G, x^*, y^*) \leq \delta_0$$

then

$$\alpha(G, x, y) \leq \frac{13 - 3\sqrt{17}}{4}.$$

Remark: $0.07 \leq \delta_0 \leq 0.08$.

Proof. We use the following inequality, see [22]

$$\alpha(G, x, y) \leq \frac{u}{\Psi(u)^2},$$

where $u = \|(x, y) - (x^*, y^*)\|\gamma(G, x^*, y^*)$.

Furthermore, we have $u/\Psi(u)^2 \leq (13 - 3\sqrt{17})/4$ for $u \in [0; \delta_0]$ and this gives the desired result. \square

In our complexity study we will need a bound on $\gamma(G, A_F)$. The following proposition will be useful.

Proposition 2.5 Let $\deg G = \max(\deg G_1, \deg G_2)$, $(x, y) \in S(0, 0, r_0)$, $G_k(X, Y) = \sum_{l=0}^{\deg G_k} \sum_{i+j=l} G_{k;i,j} X^i Y^j$,
 $\|G\|_B^2 = \sum_{k=1,2} \sum_{l=0}^{\deg G_k} \sum_{i+j=l} G_{k;i,j} \bar{G}_{k;i,j} \frac{i!j!}{\deg G_k!}$,

$$\gamma(G; x, y) \leq \|G\|_B \|DG(x, y)^{-1}\| (\deg G)^2 (1 + 2r_0^2)^{(\deg G - 2)/2}$$

Proof. We set some notations:

$$\|(x, y)\|_1 = (1 + x^2 + y^2)^{1/2},$$

$\Delta(a_i)$ is the diagonal matrix with coefficients a_i .

Now, we use the following bound, see [22, Proposition 3],

$$\gamma(G; x, y) \leq \frac{\mu(G; x, y) \deg G^{3/2}}{2\|(x, y)\|_1},$$

where ,

$$\mu(G; x, y) = \|G\|_B \|DG(x, y)^{-1}\| \Delta(\deg(G_i)^{1/2} \|(x, y)\|_1^{\deg G_i - 1})\|.$$

We use the following classical inequality $\|AB\| \leq \|A\| \|B\|$ to conclude. \square

Let a threshold $\epsilon > 0$ be given, the output of this bisection method will be a set (eventually empty) $Z = \{(x_i^*, y_i^*, r_i^*, \mu_i^*)\}$, such that :

- 1- the (x_i^*, y_i^*) 's are approximate global minima of $F(x, y)$ and $\mu_i^* = F(x_i^*, y_i^*)$.
- 2- the ball $B(x_i^*, y_i^*, r_i^*)$ contains one and only one zero of $G(x, y)$.
- 3- If $i \neq j$ then $B(x_j^*, y_j^*, r_j^*) \cap B(x_i^*, y_i^*, r_i^*) = \emptyset$ and $|\mu_i^* - \mu_j^*| < \epsilon$.

2.2 Computation of an initial square

Lemma 1 in [14] gives us a bound for the initial square:

Lemma 2.6 *Let F be the rational bivariate function corresponding to the approximate gcd problem. Let $(x, y) \in A_F$ then $\|(x, y)\| \leq 5 \max(\|f\|^2, \|g\|^2)$*

2.3 Sketch of algorithm

The algorithm consists of an initialization followed by a while loop with an internal for loop. We call step k , the k^{th} step of the while loop.

Algorithm 2.1: Approximate gcd

Input: $F = N/D$, $G = (G_1, G_2)$, an initial square S_0 and a threshold $\epsilon > 0$, as described above.

◊ Create a set of squares $L := \{S_0\}$, a set of solutions $Z := \emptyset$ and a value (the minimum to be updated) $\mu := +\infty$.

◊ While L is not empty do

- Compute $\delta = \min F(x_i, y_i)$ where the (x_i, y_i) 's are the centers of squares of L .
- If $\delta < \mu$ then $\mu := \delta$.
- For each square S of L perform the exclusion tests $E^+(P_\mu, S)$, $E(G_1, S)$ and $E(G_2, S)$. If at least one of these exclusion tests is true then remove S from L ; else, perform an inclusion test $I(G, S)$.
- If it returns false then divide S in 4 equal squares; else an approximate local minimum (x^*, y^*) is provided, it is the unique zero of $G(x, y)$ in the ball $B(x^*, y^*, r^*)$.
 - If $\mu^* = F(x^*, y^*) > \mu + \epsilon$ then remove S ; else add (x^*, y^*, r^*) to Z and update the set Z as follows:
 - If $\mu^* = F(x^*, y^*) < \mu$ then put $\mu := \mu^*$. For each element $(x_i^*, y_i^*, r_i^*, \mu_i^*)$ of Z do
 - * If $\mu < \mu_i^*$ and $|\mu - \mu_i^*| > \epsilon$ then remove the solution $(x_i^*, y_i^*, r_i^*, \mu_i^*)$ from Z
 - * If for any element $(x_i^*, y_i^*, r_i^*, \mu_i^*)$ of Z , $|\mu - \mu_i^*| < \epsilon$ and (x^*, y^*) is not in the ball $B(x_i^*, y_i^*, r_i^*)$ then add (x^*, y^*, r^*) to Z .

Output: The set Z of approximate global minima of $F(x, y)$.

Proposition 2.7 *Assume there exists $\epsilon > 0$ such that for any square $S(x, y, \epsilon) \subset S_0$ we have :*

- 1- *the inclusion test is true if $S(x, y, \epsilon)$ contains a zero of G .*
- 2- *the exclusion test is true if $S(x, y, \epsilon)$ does not contain a zero of G .*

Then we have :

- 1- *The previous algorithm stops.*
- 2- *Let μ_k be the value of μ at the end of step k . If the inclusion test is not used in the previous algorithm then $\lim_{k \rightarrow \infty} \mu_k = \min_{(x, y) \in S_0} F(x, y)$.*

Proof. The point 1 holds by construction under these assumptions. For the 2, if the inclusion test is not used in the while loop, the squares containing a global minim is divided into four squares and the algorithm doesn't stop. Moreover the sequence of the μ'_k 's decreases and the sequence of the radius r_k of the retained squares decreases strictly. Hence from the continuity of F , we obtain $\lim_{k \rightarrow \infty} \mu_k = \min_{(x, y) \in S_0} F(x, y)$. ◻

Proposition 2.8 *Assume there exists $\epsilon > 0$ such that for all square $S(x, y, \epsilon) \subset S_0$ we have :*

- 1- *the inclusion test is true if $S(x, y, \epsilon)$ contains a zero of G .*

2- the exclusion test is true if $S(x, y, \epsilon)$ does not contain a zero of G .

Then we have :

1- The previous algorithm stops.

2- Let $\mu(\epsilon)$ be the value of μ at the end of the algorithm with input ϵ . Then $\lim_{\epsilon \rightarrow 0} \mu(\epsilon) = \min_{(x,y) \in S_0} F(x, y)$.

Proof. The point 1 holds by construction under these assumptions. For point 2, the radius $r(\epsilon)$ of the retained squares in the last iteration decreases. Hence from the continuity of F , we obtain $\lim_{\epsilon \rightarrow 0} \mu(\epsilon) = \min_{(x,y) \in S_0} F(x, y)$. \square

3 Complexity analysis

3.1 Exclusion test

Let H be a polynomial in $\mathbf{R}[X, Y]$ and denote by $D^k H(x, y)$ the homogeneous part of degree k of the Taylor expansion of H at the point (x, y) .

Let $S(x_0, y_0, r_0)$ be a square. To define an exclusion function $E(H, S)$, we rely on the following expression and lemma:

$$M_H(x_0, y_0, r_0) = |H(x_0, y_0)| - \sum_{k \geq 1} \frac{\|D^k H(x_0, y_0)\|}{k!} r_0^k.$$

Lemma 3.1 *If $M_H(x_0, y_0, r_0) > 0$ then the closed square $\bar{S}(x_0, y_0, r_0)$ does not contain any zero of the polynomial $H(x, y)$.*

Proof. The proof follows from Taylor formula and a simple inequality. \square

A key to analyse the complexity of the algorithm of section 2 is the following lemma, see [8].

Lemma 3.2 *Let $H \in \mathbf{R}[X, Y]$ be a polynomial of degree e , consider the associated algebraic variety $Z_H = \{(x, y) \in \mathbf{R}^2 : H(x, y) = 0\}$.*

Let $L_e = \frac{2^{1/e} - 1}{\sqrt{2}}$ and $m_H(x, y)$ be the function implicitly defined by $M_H(x, y, m_H(x, y)) = 0$.

Then $m_H(x, y)$ is related to the distance $d(x, y, Z_H)$ by the following inequalities:

$$L_e \cdot d(x, y, Z_H) \leq m_H(x, y) \leq d(x, y, Z_H).$$

The exclusion tests to be used for the algorithm of section 2 are defined for a polynomial P by:

1- $E(P, S)$ is true if $M_P(x, y, r) > 0$,

2- $E^+(P, S)$ is true if $M_P(x, y, r) > 0$ and $P(x, y) > 0$.

Since the degree of P_μ is $2d$ and the degree of the G_i 's is $4d - 2$, we get: if $P_\mu(x, y) > 0$ then $m_{P_\mu}(x, y) \geq L_{2d} \cdot d(x, y, Z_{P_\mu})$ and $m_{G_i}(x, y) \geq L_{4d-2} \cdot d(x, y, Z_{G_i})$.

Remark: $\deg G_i \leq 4d - 2$ because the coefficient of the term of degree $4d - 1$ is $2df_d^2 + 2dg_d^2 - 2df_d^2 - 2dg_d^2 = 0$.

Putting these facts together, we can state the proposition:

Proposition 3.3 *Let $H(X, Y) = P_\mu(X, Y)$ or $H(X, Y) = G_i(X, Y)$, $i = 1, 2$. Let $S := S(x, y, r)$. The logical relation*

$$E(H, S) \text{ is true means } m_H(x, y) > r,$$

so S does not contain any zero and is excluded in the bisection algorithm. Otherwise

$$E(H, S) \text{ is false means } m_H(x, y) \leq r,$$

and S may contain zeros. In this case S will be divided into four squares each of them with a radius $r/2$.

3.2 Inclusion test

Our inclusion test is based on Smale's α -theory. The test is true if :

$$1- \alpha(G; x, y) < \frac{14 - 3\sqrt{17}}{4}.$$

$$2- S(x, y, r) \subset B(x^*, y^*, r^*) \text{ where } (x^*, y^*) \text{ is obtained from } (x, y) \text{ by } \log_2 \left(\log_2 \left(\frac{5 - \sqrt{17}}{4\gamma\epsilon} \right) + 1 \right)$$

Newton iterations and satisfies $d((x^*, y^*), A_F) \leq \epsilon$ and $r^* = \sigma(x^*, y^*)$, see Theorem 2.3.

3.3 Proof of theorem 1.1

Consider a retained square $S := S(x_l, y_l, \frac{r_0}{2^k})$ at step k . We have $\mu := \mu_k$ and $P_\mu(x_l, y_l) \geq 0$.

If $E(P_\mu, S)$ is false and $E(G_i, S)$ is false, it means, by Proposition 3.3, that $P_\mu(x_l, y_l) = 0$ or $m_{P_\mu}(x_l, y_l) \leq \frac{r_0}{2^k}$ and $m_{G_i}(x_l, y_l) \leq \frac{r_0}{2^k}$.

From Lemma 3.2, we deduce that

$$\text{if } P_\mu(x_l, y_l) > 0 \text{ then } L_{2d} \cdot d(x_l, y_l, Z_{P_\mu}) \leq m(x_l, y_l) \leq \frac{r_0}{2^k}.$$

Hence if $P_\mu(x_l, y_l) > 0$, then for each $(x, y) \in S$ we have

$$d(x, y, Z_{P_\mu}) \leq \|(x, y) - (x_l, y_l)\| + d(x_l, y_l, Z_{P_\mu}) \leq \sigma_k r_k$$

where $\sigma_k = (1 + 1/L_{2d})$ and $r_k := \frac{r_0}{2^k}$. In the same way,

$$\begin{aligned} d(x, y, Z_{G_i}) &\leq \|(x, y) - (x_l, y_l)\| + d(x_l, y_l, Z_{G_i}) \\ &\leq \omega_k r_k, \quad i = 1, 2. \end{aligned}$$

where $\omega_k = (1 + 1/L_{4d-2})$. For $H(x, y) \in \mathbf{R}[x, y]$ and $\epsilon > 0$, let

$$\begin{aligned} V_\epsilon(H) &= \{(x, y) \in \mathbf{R}^2, \text{st. } \exists (u, v) \in \mathbf{C}^2, H(u, v) = 0, \\ &\quad d((x, y), (u, v)) < \epsilon\}. \end{aligned}$$

$V_\epsilon(H)$ is called the tubular neighborhood of the solution set of $H(x, y) = 0$.

By the previous inequalities, if $P_\mu(x_l, y_l) > 0$ then we have

$$S \subset U_k := V_{\sigma_k}(P_\mu) \cap V_{\omega_k}(G_1) \cap V_{\omega_k}(G_2).$$

Notice that $A_F \subset \bigcap_{k \geq 0} U_k$. Let $\kappa_k r_k$ be half the maximal diameter of a connected component U_k . The area of U_k is bounded by $\pi \nu_k \kappa_k^2 r_k^2$ where ν_k is the number of connected components of U_k . For k big enough, this number of connected components is the number of real roots of $G_1(x, y) = G_2(x, y) = 0$ with $P_\mu(x, y) > 0$. We denote by $N < \infty$ the maximum of all the ν_k .

When k tends to infinity, the connected components of U_k tend to the real roots of $G_1(x, y) = G_2(x, y) = 0$ with $P_\mu(x, y) > 0$ and κ_k tend $\sqrt{2(1 + \cos(\alpha))}$ where α is the angle between the tangent of the curves $G_1(x, y) = 0$ and $G_2(x, y) = 0$ at the root. Let K be a bound on all the possible κ_k .

Let J be the first J such that $K r_J \leq \frac{\delta_0}{\gamma(G, A_F)}$, then for $k \geq J$, all the points of the retained squares are approximated zeros of the set A_F , by Proposition 2.4 and Theorem 2.3. An upper bound for J is given in the theorem.

To compute an upper bound for the number of exclusion tests, we notice that at step k , U_k contains the union of the kept squares (whose exclusion test is false), their number is denoted by q_k . Since the area of this union is $q_k r_k^2$ and must be less or equal than the area of U_k , we get

$$q_k \leq \pi \nu_k \delta_k^2 \leq \pi N K^2.$$

Now, let p_k be the number of excluded squares at step k . As we know the relation $4q_{k-1} = p_k + q_k$ with $p_0 = 0, q_0 = 1$ holds, the number of exclusion tests until step j is bounded by

$$\sum_{k=0}^j p_k + q_k = 1 + 4 \sum_{k=0}^{j-1} q_k \leq 1 + j \pi N K^2.$$

The last part of Theorem 1.1 comes from Proposition 2.5 which gives: $\log \gamma$ belongs to

$$O\left(\frac{d^2 r_0^{4d-3}}{\min_{(x,y) \in A_F} d_F(DG(x,y)^{-1}, \Sigma)}\right). \quad \square$$

4 Example

We consider the following example from the work of D. Rupprecht [20]. In his Ph.D. thesis, he developed a technique which allowed him to certify the degree of an approximate gcd but only if the required precision ε belongs to some intervals. There are small gaps between these intervals and the numerical computation of an approximate gcd is sensitive.

In our formulation of the nearest gcd problem, we look for the nearest pair (p, q) of degree d with a non-trivial gcd.

$$\begin{aligned} f &= (x^2 - 1.001)(x^2 + 1.00000001)(x^3 + 2x^2 - 2.999999x + 1) \\ &= 1.0000000000000000x^7 + 2.0000000000000000x^6 - 3.0009989900000000x^5 \\ &\quad + 0.9980000200000000x^4 - 0.998000041009990x^3 - 2.003000010020000x^2 \\ &\quad + 3.002999029029990x - 1.001000010010000 \\ g &= (x^2 - 0.999) * (x^2 + 1.00000003) * (x^4 + 3 * x - 1.0000002) \\ &= 1.0000000000000000x^8 + 0.0010000300000000x^6 + 3.0000000000000000x^5 \\ &\quad - 1.999000229970000x^4 + 0.0030000900000000x^3 - 0.001000030200006x^2 \\ &\quad - 2.997000089910000x + 0.999000229770006 \end{aligned}$$

The nearest perturbed pair (p, q) with a non-trivial gcd that is computed is:

$$\begin{aligned} p &= (0.000000001537219 + 0.000000006148875i)x^8 \\ &\quad + (0.999999993851125 + 0.000000001537219i)x^7 \\ &\quad + (1.999999998462781 - 0.000000006148875i)x^6 \\ &\quad + (-3.000998983851125 - 0.000000001537219i)x^5 \\ &\quad + (0.998000021537219 + 0.000000006148874i)x^4 \\ &\quad + (-0.998000047158864 + 0.000000001537219i)x^3 \\ &\quad + (-2.003000011557218 - 0.000000006148874i)x^2 \\ &\quad + (3.002999035178864 - 0.000000001537219i)x \\ &\quad + (-1.001000008472781 + 0.000000006148874i) \\ q &= (1.000000000000000 + 0.000000008719619i)x^8 \\ &\quad + (-0.000000008719619 - 0.000000000000000i)x^7 \\ &\quad + (0.001000030000000 - 0.000000008719619i)x^6 \\ &\quad + (3.000000008719619 + 0.000000000000000i)x^5 \\ &\quad + (-1.999000229970001 + 0.000000008719619i)x^4 \\ &\quad + (0.003000081280381 - 0.000000000000000i)x^3 \\ &\quad + (-0.001000030200006 - 0.000000008719619i)x^2 \\ &\quad + (-2.997000081190381 + 0.000000000000000i)x \\ &\quad + (0.999000229770005 + 0.000000008719619i) \end{aligned}$$

The order of the perturbation is of the order 10^{-8} and the degree of the corresponding gcd is 1.

Note that in this example, even if f is of degree 7, we allow a perturbation of degree 8 (which is the degree of g).

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