# Factoring Polynomials via Polytopes

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# ABSTRACT

We introduce a new approach to multivariate polynomial factorisation which incorporates ideas from polyhedral geometry, and generalises Hensel lifting. Our main contribution is to present an algorithm for factoring bivariate polynomials which is able to exploit to some extent the sparsity of polynomials. We give details of an implementation which we used to factor randomly chosen sparse and composite polynomials of high degree over the binary field.

# **Categories and Subject Descriptors**

I.1.2 [Computing Methodologies]: Symbolic and Algebraic Manipulation—Algebraic Algorithms

# **General Terms**

algorithms, experimentation

# Keywords

multivariate polynomial, factorisation, Newton polytope

# 1. INTRODUCTION

Factoring polynomials is a fundamental problem in algebra and number theory and it is a basic routine in all major computer algebra systems. There is an extensive literature on this problem — we refer the reader to the references in [6]. Most of these papers deal with dense polynomials, two notable exceptions being [7, 9]. These two papers reduce sparse polynomials with more than two variables to bivariate or univariate polynomials which are then treated as dense polynomials. It is still open whether there is an efficient algorithm for factoring sparse bivariate or univariate polynomials. Our goal in this paper is to study sparse bivariate polynomials using their connection to integral polytopes.

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Newton polytopes of multivariate polynomials reflect to a certain extent the sparsity of polynomials and they carry a lot of information about the factorisation patterns of polynomials as demonstrated in our recent work [2, 3]. In [5], we deal with irreducibility of random sparse polynomials. In this paper our focus is on the more difficult problem of factoring sparse polynomials. We do not solve this problem completely. However, our approach is a practical new method which generalises Hensel lifting; its running time will in general improve upon that of Hensel lifting and sparse bivariate polynomials can often be processed significantly more quickly. As with Hensel lifting, it has an exponential worst-case running time. Also, our method does not work for all polynomials, but only for those that are squarefree on certain subsets of the edges of their Newton polytopes (see Theorem 7).

Here is a brief outline of the paper. In Section 2 we present a brief introduction to Newton polytopes and their relation to multivariate polynomials, and in Section 3 we state our central problem. Section 4 contains an outline of our method, and highlights the theoretical problems we need to address. The main theorem underpinning our method is proved in Section 6, after a key geometric lemma in Section 5. Section 7 contains a concise description of the algorithm. Finally in Section 8 we present a small example, as well as details of our computer implementation of the algorithm. We believe the main achievements of this paper are the theoretical results in Section 6, and the high degree polynomials we have factored using the method, as presented in Subsection 8.2.

## 2. OSTROWSKI'S THEOREM

This paper considers polynomial factorisation over a field  $\mathbb{F}$  of arbitrary characteristic. We denote by  $\mathbb{N}$  the *non-negative* integers, and  $\mathbb{Z}$ ,  $\mathbb{Q}$  and  $\mathbb{R}$  the integers, rationals and reals.

Let  $\mathbb{F}[X_1, X_2, \ldots, X_n]$  be the ring of polynomials in n variables over the field  $\mathbb{F}$ . For any vector  $e = (e_1, \ldots, e_n)$  of non-negative integers define  $X^e := X_1^{e_1} \ldots X_n^{e_n}$ . Let  $f \in \mathbb{F}[X_1, \ldots, X_n]$  be given by

$$f := \sum_{e} a_e X^e \tag{1}$$

where the sum is over finitely many points e in  $\mathbb{N}^n$ , and  $a_e \in \mathbb{F}$ . The Newton polytope of f, Newt(f), plays an essential role in all that follows. It is the polytope in  $\mathbb{R}^n$  obtained as the convex hull of all exponents e for which the corresponding coefficient  $a_e$  is non-zero. It has integer vertices,

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since all the e are integral points; we call such polytopes *integral*. Given two polytopes Q and R their *Minkowski sum* is defined to be the set

$$Q + R := \{ q + r \, | \, q \in Q, r \in R \}.$$

When Q and R are integral polytopes, so is Q+R. If we can write an integral polytope P as a Minkowski sum Q+R for integral polytopes Q and R then we call this an *(integral)* decomposition. The decomposition is trivial if Q or R has only one point. For an example of nontrivial decomposition, see the figures in Section 8 where Newt(f) = Q + R. The motivating theorem behind our investigation is (see [2]):

THEOREM 1 (OSTROWSKI). Let 
$$f, g, h \in \mathbb{F}[X_1, \dots, X_n]$$
.  
If  $f = gh$  then  $Newt(f) = Newt(g) + Newt(h)$ .

An immediate result of this theorem relates to testing polynomial irreducibility: In the simplest case in which the polytope does not decompose, one immediately deduces that the polynomial must be irreducible. This was explored in [2, 3, 5], in particular a quasi-polynomial time algorithm is presented in [5] for finding all the decompositions of any given integral polytope in a plane. In this paper, we address the more difficult problem: Given a decomposition of the polytope, how can we recover a factorisation of the polynomial whose factors have Newton polytopes of that shape, or show that one does not exist?

In the remainder of the paper, we restrict our attention to bivariate polynomials, and f always denotes a bivariate polynomial in the ring  $\mathbb{F}[x, y]$ . For  $e = (e_1, e_2) \in \mathbb{N}^2$ , we redefine the notation  $X^e$  to mean  $x^{e_1}y^{e_2}$ . We shall retain the term "Newton polytope" for the polygon Newt(f) to avoid confusion with other uses of the term "Newton polygon".

# 3. PARTIAL FACTORISATIONS

Let Newt(f) = Q + R be a decomposition of the Newton polytope of f into integral polytopes in the first quadrant. For each lattice point  $q \in Q$  and  $r \in R$  we introduce indeterminates  $g_q$  and  $h_r$ . The polynomials g and h are then defined as

$$\begin{array}{rcl} g & := & \sum_{q \in Q} g_q X^q \\ h & := & \sum_{r \in R} h_r X^r. \end{array}$$

We call g and h the *generic* polynomials given by the decomposition Newt(f) = Q + R. Let #Newt(f) denote the number of lattice points in Newt(f). The equation f = ghdefines a system of #Newt(f) quadratic equations in the coefficient indeterminates obtained by equating coefficients of each monomial  $X^e$  with  $e \in Newt(f)$  on both sides. The aim is to find an efficient method of solving these equations for field elements. Our approach, motivated by Hensel lifting, is to assume that, along with the decomposition of the Newton polytope, we are given appropriate factorisations of the polynomials defined along its edges. This "boundary factorisation" of the polynomial is then "lifted" into the Newton polytope, and the coefficients of the possible factors q and h revealed in successive layers. Unfortunately, to describe the algorithm properly we shall need a considerable number of elementary definitions — the reader may find the figures in Section 8.1 useful in absorbing them all.

Let S be a subset of Newt(f). An S-partial factorisation of f is a specialisation of a subset of the indeterminates  $g_q$ and  $h_r$  such that for each lattice point  $s \in S$  the coefficients of monomials  $X^s$  in the polynomials gh and f are equal field elements. (A specialisation is just a substitution of field elements in place of indeterminates.) The case S = Newt(f) is equivalent to a factorisation of f in the traditional sense, and we will call this a *full factorisation*. Now suppose we have an S-partial factorisation and an S'-partial factorisation. Suppose further  $S \subseteq S'$  and the indeterminates specialised in the S-partial factorisation have been specialised to the same field elements as the corresponding ones in the S'-partial factorisation. Then we say the S'-partial factorisation extends the S-partial factorisation. We call this extension proper if S' has strictly more lattice points than S.

Let  $\operatorname{Edge}(f)$  denote the set of all edges of  $\operatorname{Newt}(f)$ . Each edge  $\delta \in \operatorname{Edge}(f)$  is viewed as directed so that  $\operatorname{Newt}(f)$  lies on the left hand side of the edge, and this directed edge can be defined by an affine function  $\ell$  as follows. Suppose the edge  $\delta$  is from  $(u_1, v_1)$  to  $(u_2, v_2)$ , vertices with integral coordinates;  $(u_1, v_1)$  is called the starting vertex of the edge. Let  $d = \operatorname{gcd}(u_2 - u_1, v_2 - v_1)$ ,  $u_0 = (u_2 - u_1)/d$ , and  $v_0 = (v_2 - v_1)/d$ . Then  $(u_0, v_0)$  represents the direction of  $\delta$  and the integral points on  $\delta$  are of the form

$$(u_1, v_1) + i(u_0, v_0), \quad 0 \le i \le d.$$

Let  $(\nu_1, \nu_2) := (-v_0, u_0)$ , a rotation of  $(u_0, v_0)$  by 90 degrees counter clockwise, and let  $\eta = v_0 u_1 - u_0 v_1$ . Define

 $\ell(e) = \nu_1 e_1 + \nu_2 e_2 + \eta$ , for  $e = (e_1, e_2) \in \mathbb{R}^2$ .

Then  $\ell$  has the property that  $\ell(e) \geq 0$  for each point  $e \in$ Newt(f), with the equation holding iff  $e \in \delta$ , that is, Newt(f) lies in the positive side of the line  $\ell = 0$ . We call this function  $\ell$  the primitive affine function associated with  $\delta$ , denoted by  $\ell_{\delta}$ .

The function  $\ell_{\delta}$  has another nice property: Since  $gcd(\nu_1, \nu_2) = 1$ , there exist integers  $\zeta_1$  and  $\zeta_2$  such that  $\zeta_1\nu_1 + \zeta_2\nu_2 = 1$ , and they are unique under the requirement that  $0 \leq \zeta_2 < \nu_1$ . Define the change of variables

$$z := x^{\nu_2} y^{-\nu_1}$$
 and  $w := x^{\zeta_1} y^{\zeta_2}$ . (2)

Then any monomial of the form  $x^{e_1}y^{e_2}$  can be written as  $x^{e_1}y^{e_2} = z^{i_1}w^{i_2}$ , where

$$egin{pmatrix} i_1\ i_2 \end{pmatrix} = egin{pmatrix} \zeta_2 & -\zeta_1\ 
u_1 & 
u_2 \end{pmatrix} egin{pmatrix} e_1\ e_2 \end{pmatrix}.$$

Its inverse transform is

$$\begin{pmatrix} e_1 \\ e_2 \end{pmatrix} = \begin{pmatrix} \nu_2 & \zeta_1 \\ -\nu_1 & \zeta_2 \end{pmatrix} \begin{pmatrix} i_1 \\ i_2 \end{pmatrix}.$$

This change of variables has the nice property that when  $(e_1, e_2)$  moves along the direction  $(u_0, v_0)$  of the edge  $\delta$ , then the exponent of w remain constant (as  $i_2 = \ell_{\delta}(e_1, e_2) - \eta$ ), while the exponent of z strictly increases (by  $1 = \zeta_2 u_0 - \zeta_1 v_0$  for each increment of  $(u_0, v_0)$ ).

For each  $\delta \in \text{Edge}(f)$ , there exists a unique pair of faces (either edges or vertices)  $\delta'$  and  $\delta''$  of Q and R, respectively, such that  $\delta = \delta' + \delta''$ . One can also easily show that there exists a unique integer  $c_{\delta}$  such that

$$\delta' = \{e \in Q \mid \ell_{\delta}(e) = c_{\delta} \}$$
  
$$\delta'' = \{e \in R \mid \ell_{\delta}(e) = -c_{\delta} + \eta \}$$

where  $\eta$  is the constant coefficient of  $\ell_{\delta}$ .

Let  $\Gamma \subseteq \text{Edge}(f)$ , and let  $K = (k_{\gamma})_{\gamma \in \Gamma}$  be a vector of positive integers, one for each edge  $\gamma \in \Gamma$ . Define

 $\operatorname{Newt}(f)|_{\Gamma,K} := \{ e \in \operatorname{Newt}(f) \mid 0 \le \ell_{\gamma}(e) < k_{\gamma} \text{ for } \gamma \in \Gamma \}.$ 

This defines a strip along the interior of Newt(f), or a union of such strips. We denote by  $Q|_{\Gamma,K}$  and  $R|_{\Gamma,K}$  the subsets of Q and R respectively given by

$$\begin{aligned} Q|_{\Gamma,K} &:= \{ e \in Q \,|\, 0 \le \ell_{\delta}(e) < k_{\delta} + c_{\delta} \text{ for } \delta \in \Gamma \} \\ R|_{\Gamma,K} &:= \{ e \in R \,|\, 0 \le \ell_{\delta}(e) < k_{\delta} - c_{\delta} + \eta \text{ for } \delta \in \Gamma \}. \end{aligned}$$

Once again these denote strips along the inside of Q and R whose sum contains the strip Newt $(f)|_{\Gamma,K}$  in Newt(f), as indicated by Figures 2, 3 and 4.

We now come to the main definition of this section.

DEFINITION 2. A  $Newt(f)|_{\Gamma,K}$ -factorisation is called a  $(\Gamma, K; Q, R)$ -factorisation if the following two properties hold:

- Exactly the indeterminate coefficients of g and h indexed by lattice points in  $Q|_{\Gamma,K}$  and  $R|_{\Gamma,K}$ , respectively, have been specialised.
- Let K' = (k'<sub>γ</sub>)<sub>γ∈Γ</sub> be a vector of positive integers with k'<sub>γ</sub> ≥ k<sub>γ</sub> for all γ ∈ Γ, with the inequality strict for at least one γ. Then not all of the indeterminate coefficients of g indexed by lattice points in Q|<sub>Γ,K'</sub> have been specialised.

The second property will be used only once, in the proof of Lemma 8. In most instances Q, R and  $\Gamma$  will be clear from the context. If so we will omit them and refer simply to a K-factorisation. Furthermore, if K is the all ones vector, denoted (<u>1</u>), of the appropriate length indexed by elements of some set  $\Gamma$ , then we call this a ( $\Gamma; Q, R$ )-boundary factorisation. We shall simplify this to partial boundary factorisation or (<u>1</u>)-factorisation when  $\Gamma$ , Q and R are evident from the context. This special case will be the "lifting off" point for our algorithm.

The central problem we address is

**Problem 3** Let  $f \in \mathbb{F}[x, y]$  have Newton polytope Newt(f)and fix a Minkowski decomposition Newt(f) = Q + R where Q and R are integral polygons in the first quadrant. Suppose we have been given a  $(\Gamma; Q, R)$ -boundary factorisation of f for some set  $\Gamma \subseteq \text{Edge}(f)$ . Construct a full factorisation of f which extends it, or show that one does not exist.

Moreover, one wishes to solve the problem in time bounded by a small polynomial function of #Newt(f).

## 4. THE POLYTOPE METHOD

#### 4.1 An outline of the method

We now give a basic sketch of our polytope factorisation method for bivariate polynomials. Throughout this section  $\Gamma$  will be a fixed subset of  $\operatorname{Edge}(f)$  and  $\operatorname{Newt}(f) = Q + R$  a fixed decomposition. We shall need to place certain conditions on  $\Gamma$  later on, but for the time being we will ignore them. Since  $\Gamma, Q$  and R are fixed we shall use our abbreviated notation when discussing partial factorisations.

We begin with  $K = (\underline{1})$  the all-ones vector of the appropriate length and a K-factorisation (partial boundary factorisation). Recall this is a partial factorisation in which exactly the coefficients in the sets  $Q|_{\Gamma,K}$  and  $R|_{\Gamma,K}$ , subsets of points on the boundaries of Q and R, have been specialised.

At each step of the algorithm we either show that no full factorisation of f exists which extends this partial factorisation, and halt, or that at most one can exist, and we find a

new K'-factorisation with  $K' = (k'_{\delta})$  such that

$$\sum_{\delta \in \Gamma} k'_{\delta} > \sum_{\delta \in \Gamma} k_{\delta}.$$

(Usually the sum will be incremented by just one.) Iterating this procedure either we halt after some step, in which case we know that no factorisation of f exists which extends the original partial boundary factorisation, or we eventually have  $Q \subseteq Q|_{\Gamma,K}$ , say. At that point all of the indeterminates in one of our partial factors have been specialised, and we may check to see if we have found a factor by division.

Note that in the situation in which Newt(f) is just a triangle with vertices (0, n), (n, 0) and (0, 0) for some n, our method reduces to the standard Hensel lifting method for bivariate polynomial factorisation. As such, our "polytope method" is a natural generalisation of Hensel lifting from the case of "generic" dense polynomials to arbitrary, possibly sparse, polynomials.

## 4.2 Hensel lifting equations

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In this section we derive the basic equations which are used in our algorithm.

Let f be as in (1). For any  $\delta \in \text{Edge}(f)$  recall that  $\ell_{\delta}$  is the associated primitive affine function. For  $i \geq 0$  we define

$$f_i^{\delta} := \sum_{\ell_{\delta}(e)=i} a_e X^e.$$

Thus  $f_i^{\delta}$  is just the polynomial obtained from f by removing all terms whose exponents do not lie on the "ith translate of the supporting line of  $\delta$  into the polytope Newt(f)". We call the polynomials  $f_0^{\delta}$  edge polynomials.

Given the decomposition Newt(f) = Q + R let  $\delta'$  and  $\delta''$ denote the unique faces of Q and R which sum to  $\delta$ . As before assume  $\ell_{\delta}(\delta') = c_{\delta}$  and  $\ell_{\delta}(\delta'') = -c_{\delta} + \eta$ . Let g and h denote generic polynomials with respect to Q and R. For  $i \geq 0$  define

$$g_i^{\delta} := \sum_{q \in Q, \, \ell_{\delta}(q) = c_{\delta} + i} g_q X^q$$
$$h_i^{\delta} := \sum_{r \in R, \, \ell_{\delta}(r) = -c_{\delta} + \eta + i} h_r X^r.$$

Once again  $g_i^{\delta}$  and  $h_i^{\delta}$  are obtained from g and h by considering only those terms which lie on particular lines. The next result is elementary but fundamental.

LEMMA 4. Let  $f \in \mathbb{F}[x, y]$  and Newt(f) = Q + R be an integral decomposition with corresponding generic polynomials g and h. Let Edge(f) denote the set of edges of Newt(f) and  $\delta \in Edge(f)$ . The system of equations in the coefficient indeterminates of g and h defined by equating monomials on both sides of the equality f = gh has the same solutions as the system of equations defined by the following:

$$f_0^{\delta} = g_0^{\delta} h_0^{\delta}, \text{ and } g_0^{\delta} h_k^{\delta} + h_0^{\delta} g_k^{\delta} = f_k^{\delta} - \sum_{j=1}^{k-1} g_j^{\delta} h_{k-j}^{\delta} \text{ for } k \ge 1.$$
(3)

Thus any specialisation of coefficient indeterminates which is a solution of equations (3) will give a full factorisation of f. PROOF. In the equation f = gh gather together on each side all monomials whose exponent vectors lie on the same translate of the line supporting  $\delta$ .

These are precisely the equations which are used in Hensel lifting to try and reduce the non-linear problem of selecting a specialisation of the coefficients of g and h to give a factorisation of f, to a sequence of linear systems which may be recursively solved. We recall precisely how this is done, as our method is a generalisation.

We begin with a specialisation of the coefficients in the polynomials  $g_0^{\delta}$  and  $h_0^{\delta}$  which gives a factorisation of the polynomial  $f_0^{\delta}$ . Equation (3) with k = 1 gives a linear system for the indeterminate coefficients of  $g_1^{\delta}$  and  $h_1^{\delta}$ . In the special case in which standard Hensel lifting applies this system may be solved uniquely, and thus a unique partial factorisation of f is defined which extends the original one. This process is iterated for k > 1 until all the indeterminate coefficients in one of the generic polynomials have been specialised, at which stage one checks whether a factor has been found by division.

The problem with this method is that in general there may *not* be a unique solution to each of the linear systems encountered. There will be a unique solution in the dense bivariate case mentioned at the end of Section 4.1, subject to a certain coprimality condition. General bivariate polynomials may be reduced to ones of this form by randomisation, but the substitutions involved destroy the sparsity of the polynomial. Our approach avoids this problem, although again is not universal in its applicability. As explained earlier, our method extends a *special* kind of partial boundary factorisation of f, rather than just the factorisation of one of its edges. In this way uniqueness in the bivariate case is restored.

## 5. A GEOMETRIC LEMMA

This section contains a geometric lemma which ensures our method can proceed in a unique way at each step provided we start with a special type of partial boundary factorisation. We begin with a key definition.

DEFINITION 5. Let  $\Lambda$  be a set of edges of a polygon P in  $\mathbb{R}^2$  and r a vector in  $\mathbb{R}^2$ . We say that  $\Lambda$  dominates P in direction r if the following two properties hold:

- P is contained in the Minkowski sum of the set Λ and the infinite line segment rℝ<sub>≥0</sub> (the positive hull of r). Call this sum Mink(Λ, r).
- Each of the two infinite edges of Mink(Λ, r) contains exactly one point of P.

Thus  $\operatorname{Mink}(\Lambda, r)$  comprises a region bounded by the interior strip between its two infinite edges and the edges in  $\Lambda$ . This definition is illustrated in Figure 1 where  $\Lambda$  consists of all the bold edges on the boundary indicated by T.

We will call  $\Lambda$  an *irredundant* dominating set if it is a dominating set (in direction r, say) that does not strictly contain any other dominating set (in direction r). The edges in an irredundant dominating set are necessarily connected. For a polygon P in  $\mathbb{R}^2$  which is not a single point, it is obvious that there exists at least one irredundant dominating set: pick a direction r which is not a slope of any edge, and starting with the dominating set of all edges delete edges as required.



Figure 1: Dominating set of edges

The next lemma is at the heart of our algorithm.

LEMMA 6. Let P be an integral polygon and  $\Lambda$  an irredundant dominating set of edges of P. Suppose  $\Lambda_1$  is a polygonal line segment in P such that each edge of  $\Lambda_1$  is parallel to some edge of  $\Lambda$ . If  $\Lambda_1$  is different from  $\Lambda$  then  $\Lambda$  has at least one edge that has strictly more lattice points than the corresponding edge of  $\Lambda_1$ .

The lemma is illustrated in Figure 1, where T denotes the union of the edges in  $\Lambda$  and  $T_1$  the union of the line segments in  $\Lambda_1$ .

Before proving this lemma we make one more definition. We define a map  $\pi_r$  onto the orthogonal complement  $\langle r \rangle^{\perp} := \{s \in \mathbb{R}^2 \mid (s \cdot r) = 0\}$  of the vector r as follows:

$$\pi_r(v) = v - \left(\frac{v \cdot r}{r \cdot r}\right) r.$$

We call this projection by r, and we have that  $\pi_r(P) = \pi_r(\Lambda)$ . Notice that if  $e_1$  and  $e_2$  are adjacent edges in an irredundant dominating set, then the length of the projection by r of the polygonal line segment  $e_1e_2$  is just the sum of the lengths of the projections by r of the individual edges  $e_1$  and  $e_2$ . For otherwise, we would have, say,  $\pi_r(e_1) \subseteq \pi_r(e_2)$  and hence  $r\mathbb{R}_{\geq 0} + e_1 \subseteq r\mathbb{R}_{\geq 0} + e_2$ . Thus the edge  $e_1$  would be redundant, a contradiction. The same is true if we replace  $e_1$  and  $e_2$  by any adjacent line segments parallel to them — we still obtain an "additivity" in the lengths, which shall be used in the proof of the lemma.

PROOF. We assume that  $\Lambda$  dominates P in the direction r as shown in Figure 1. Let  $\delta_1, \dots, \delta_k$  be the edges in  $\Lambda$  and  $\delta'_1, \dots, \delta'_k$  the corresponding edges of  $\Lambda_1$ . Let  $n_i$  be the number of lattice points on  $\delta_i$ , and  $m_i$  that on  $\delta'_i$ ,  $1 \le i \le k$ . We want to show that  $n_i > m_i$  for at least one  $i, 1 \le i \le k$ . Suppose otherwise, namely

$$n_i \le m_i, \quad 1 \le i \le k. \tag{4}$$

We derive a contradiction by considering the lengths of  $\Lambda$ and  $\Lambda_1$  on the projection by  $\pi_r$ . Note that if  $m_i = 0$  for some *i* then certainly  $n_i > m_i$  and we are done; thus we may assume that  $m_i \ge 1$  for all *i*.

First, certainly  $\pi(\Lambda_1) \subseteq \pi(\Lambda)$  as  $\Lambda$  is a dominating set. Since  $\Lambda_1$  is different from  $\Lambda$ , their corresponding end points must not coincide. Hence at least one end point of  $\Lambda_1$  will not be on the infinite edges in the direction r. Hence  $\pi_r(\Lambda_1)$ lies completely inside  $\pi_r(\Lambda)$ , so has length strictly shorter than  $\pi_r(\Lambda)$ .

Now for  $1 \leq i \leq k$  let  $\epsilon_i$  be the length of the projection of a primitive line segment on  $\delta_i$  (which means that the line segment has both end points on lattice points but no lattice points in between). Certainly  $\epsilon_i \geq 0$ . Since the end points of  $\delta_i$  are lattice points, the length of  $\pi_r(\delta_i)$  is exactly  $(n_i - 1)\epsilon_i$ for  $1 \leq i \leq k$ , hence  $\pi_r(\Lambda)$  has length  $\sum_{i=1}^k (n_i - 1)\epsilon_i$ . (Here we need the fact that the dominating set is irredundant, to give us the necessary "additivity" in the lengths.) For  $\delta'_i$ , since it is parallel to  $\delta_i$ , the projected length of a primitive line segment on it is also  $\epsilon_i$ . Hence the length of  $\pi_r(\Lambda_1)$  is at least  $\sum_{i=1}^k (m_i - 1)\epsilon_i$  and from (4) we know that

$$\sum_{i=1}^{k} (m_i - 1)\epsilon_i \ge \sum_{i=1}^{k} (n_i - 1)\epsilon_i$$

This contradicts our previous observation that  $\pi_r(\Lambda_1)$  is strictly shorter than  $\pi_r(\Lambda)$ . The lemma is proved.

# 6. THE MAIN THEOREM

Let  $\Gamma$  be an irredundant dominating set of Newt(f). We call a ( $\Gamma$ ; Q, R)-boundary factorisation of f a dominating edges factorisation relative to  $\Gamma, Q$  and R. A coprime dominating edges factorisation is a ( $\Gamma$ ; Q, R)-boundary factorisation with the property that for each  $\delta \in \Gamma$  the edge polynomials  $g_0^{\delta}$  and  $h_0^{\delta}$  are coprime, up to monomial factors. (In other words, they are coprime as Laurent polynomials.)

We are now ready to state our main theoretical result.

THEOREM 7. Let  $f \in \mathbb{F}[x, y]$  and Newt(f) = Q + R be a fixed Minkowski decomposition, where Q and R are integral polygons in the first quadrant. Let  $\Gamma$  be an irredundant dominating set of Newt(f) in direction r, and assume that Q is not a single point or a line segment parallel to  $r\mathbb{R}_{\geq 0}$ . For any coprime dominating edges factorisation of f relative to  $\Gamma, Q$  and R, there exists at most one full factorisation of f which extends it, and moreover this full factorisation may be found or shown not to exist in time polynomial in #Newt(f).

We shall prove this theorem inductively through the next two lemmas.

LEMMA 8. Let f, Q, R and  $\Gamma$  be as in the statement of Theorem 7. Suppose we are given a K-factorisation of f, where  $K = (k_{\delta})_{\delta \in \Gamma}$  (more specifically, a  $(\Gamma, K; Q, R)$  - factorisation). For each  $\delta \in \Gamma$ , denote by  $\delta'$  the face of Qsupported by  $\ell_{\delta} - c_{\delta}$ . There exists  $\delta \in \Gamma$  with the following properties

- The face  $\delta'$  is an edge (rather than a vertex).
- The number of unspecialised coefficients of g<sup>δ</sup><sub>kδ</sub> is nonzero but strictly less than the number of integral points on δ'.

All the unspecialised terms of g<sup>δ</sup><sub>kδ</sub> have exponents being consecutive integral points on the line defined by ℓ<sub>δ</sub> = (c<sub>δ</sub> + k<sub>δ</sub>).

PROOF. Let  $\overline{Q}$  be the polygon

$$Q := \{ r \in Q \mid \ell_{\delta}(r) \ge c_{\delta} + k_{\delta} \text{ for all } \delta \in \Gamma \}.$$

Note that the lattice points in  $\bar{Q}$  correspond to unspecialised coefficients of g. Let  $\Lambda$  denote the set of edges  $\delta \in \Gamma$  of Newt(f) such that the functional  $\ell_{\delta} - c_{\delta}$  supports an edge of Q (rather than just a vertex). Note that  $\Lambda \neq \emptyset$ , for otherwise Q must be a single point or a line segment in direction r, contradicting our assumption. We denote the edge by  $\delta'$ , and write  $\bar{\delta}$  for the face of  $\bar{Q}$  supported by  $\ell_{\delta} - (c_{\delta} + k_{\delta})$ . Note that each  $\overline{\delta}$  contains at least one lattice point. (This follows from the second property in Definition 2.) Certainly,  $\bar{\delta}$  is parallel to  $\delta'$  for each  $\delta \in \Lambda$ , and the edge sequence  $\{\bar{\delta}\}_{\delta\in\Lambda}$ , forms a polygonal line segment in Q. Since  $\Gamma$  is an irredundant dominating set for Newt(f), the set of edges  $\{\delta'\}_{\delta \in \Lambda}$  is an irredundant dominating set for Q. By Lemma 6, there is at least one edge  $\delta \in \Lambda$ , such that  $\delta'$  has strictly more lattice points than  $\overline{\delta}$ . This edge  $\delta$  has the required properties. This completes the proof.  $\Box$ 

LEMMA 9. Let f, Q, R and  $\Gamma$  be as in the statement of Theorem 7. Suppose we are given a K-factorisation of f, where  $K = (k_{\delta})_{\delta \in \Gamma}$ . Moreover, assume this factorisation extends a coprime dominating edges factorisation, i.e., the polynomials  $g_{\delta}^{0}$  and  $h_{\delta}^{0}$  are coprime up to monomial factors for all  $\delta \in \Gamma$ . Then there exists  $\delta \in \Gamma$  such that the coefficients of  $g_{k_{\delta}}^{k}$  are not all specialised, but they may be specialised in at most one way consistent with equations (3). This specialisation may be computed in time polynomial in #Newt(f).

PROOF. The basic idea of proof is to first transform the bivariate equation (3) into equations of univariate polynomials determined by the individual edges, then to determine the existence or uniqueness of solutions.

Select  $\delta \in \Gamma$  such that the properties in Lemma 8 hold. Let  $n_{\delta}$  and  $m_{\delta}$  be the number of integral points on the edges  $\delta'$  and  $\bar{\delta}$  respectively, where  $\delta'$  and  $\bar{\delta}$  are defined as in the proof of Lemma 8. Thus we have  $m_{\delta} < n_{\delta}$  and  $m_{\delta} \geq 1$ . With the notation from Section 3, write  $\ell_{\delta}(e_1, e_2) = \nu_1 e_1 + \nu_2 e_2 + \eta$ , where  $\nu_1$  and  $\nu_2$  are coprime.

Let z and w be new variables. Using the transform (2), any monomial of the form  $x^{e_1}y^{e_2}$  can be written as

$$x^{e_1}y^{e_2} = z^{i_1}w^{i_2} (5)$$

where

$$i_1 = e_1\zeta_2 - e_2\zeta_1, i_2 = e_1\nu_1 + e_2\nu_2 = \ell_\delta(e_1, e_2) - \eta.$$

Every monomial in  $g_i^{\delta}$  is of the form  $x^{e_1}y^{e_2}$  where  $\ell_{\delta}(e_1, e_2) = c_{\delta} + i$ . Let the monomials s and t be the terms of g and h respectively whose exponents vectors are the starting vertices of the faces of Q and R defined by  $\ell_{\delta} - c_{\delta}$  and  $\ell_{\delta} + c_{\delta} - \eta$ , respectively. Thus we have  $g_i^{\delta}(z, w) = sw^i G_i(z)$  for some univariate Laurent polynomial  $G_i(z)$ . Similarly  $h_i^{\delta}(z, w) = tw^i H_i(z)$  and  $f_i^{\delta}(z, w) = stw^i F_i(z)$ , where  $H_i(z)$  and  $F_i(z)$  are univariate Laurent polynomials. With this construction,  $G_0(z), H_0(z)$  and  $F_0(z)$  have non-zero constant term and are "ordinary polynomials", i.e., contain no negative powers of

z. For  $i < k_{\delta}$  all of the coefficients in the polynomials  $G_i(z)$ and  $H_i(z)$  have been specialised. Moreover  $G_0(z)$  is of degree  $n_{\delta}$ , and all but  $m_{\delta}$  of the coefficients of  $G_{k_{\delta}}(z)$  have been specialised. (By "degree" of a Laurent polynomial we mean the difference in the exponents of the highest and lowest terms, if the polynomial is non-zero, and  $-\infty$  otherwise). Equations (3) with this change of variables may be written as  $F_0(z) = G_0(z)H_0(z)$ , and for k > 1

$$G_k(z)H_0(z) + G_0(z)H_k(z) = F_k(z) - \sum_{j=1}^{k-1} G_j(z)H_{k-j}(z).$$

We know that all of the coefficients of  $G_i(z)$  and  $H_i(z)$  have been specialised for  $0 \le i < k_{\delta}$  in such a way as to give a solution to  $F_0 = G_0 H_0$  and the first  $k_{\delta} - 1$  equations above. Thus we need to try and solve

$$G_{k_{\delta}}H_{0} + G_{0}H_{k_{\delta}} = F_{k_{\delta}} - \sum_{j=1}^{k_{\delta}-1} G_{j}H_{k_{\delta}-j}.$$
 (6)

for the unspecialised indeterminate coefficients of  $G_{k\delta}$  and  $H_{k\delta}.$ 

We first compute using Euclid's algorithm ordinary polynomials U(z) and V(z) such that

$$V(z)H_0(z) + U(z)G_0(z) = 1$$

where  $\deg_z(U(z)) < \deg_z(H_0(z))$  and  $\deg_z(V(z)) < \deg_z(G_0(z))$ . (Note that  $G_0(z)$  and  $H_0(z)$  are coprime since we have a coprime partial boundary factorisation.) Any solution  $G_{k_{\delta}}$  of Equation (6) must be of the form

$$G_{k_{\delta}} = \{ V(F_{k_{\delta}} - \sum_{j=1}^{k_{\delta}-1} G_j H_{k_{\delta}-j}) \bmod G_0 \} + \varepsilon G_0 \qquad (7)$$

for some Laurent polynomial  $\varepsilon(z)$  with undetermined coefficients.

We rearrange (7) as

$$G_{k_{\delta}} - \{V(F_{k_{\delta}} - \sum_{j=1}^{k_{\delta}-1} G_j H_{k_{\delta}-j}) \bmod G_0\} = \varepsilon G_0 \qquad (8)$$

Let the degree in z of the Laurent polynomial on the lefthand side of this equation be d. Now the degree of the polynomial  $G_0(z)$  as a Laurent polynomial (and an ordinary polynomial) is  $n_{\delta} - 1$ . If  $d < n_{\delta} - 1$  then we must have d = 0. In other words, (7) has a unique solution, namely that with  $\varepsilon = 0$ . Otherwise  $d \ge n_{\delta} - 1$  and the degree in z of  $\varepsilon(z)$  as a Laurent polynomial is  $d - (n_{\delta} - 1)$ . Hence in this case we need to also solve for the  $d - n_{\delta} + 2$  unknown coefficients of  $\varepsilon(z)$ . We know that all but  $m_{\delta}$  coefficients of  $G_{k_{\delta}}$  have already been specialised, and these unspecialised ones are adjacent terms. Hence exactly  $(d + 1) - m_{\delta}$  coefficients on the lefthand side of (8) have been specialised, which are adjacent lowest and highest terms. By assumption we have that  $m_{\delta} < n_{\delta}$ , and hence  $(d + 1) - m_{\delta} \ge d - n_{\delta} + 2$ .

All of the coefficients of the righthand side of Equation (8) have been specialised, except those of the unknown polynomial  $\varepsilon(z)$ . On the lefthand side all but the middle  $m_{\delta}$  coefficients have been specialised. This defines a pair of triangular systems from which one can either solve for the coefficients of  $\varepsilon$  uniquely, or show that no solution exists (this may happen when  $n_{\delta} > m_{\delta} + 1$ ). We describe precisely how this is done: Suppose that exactly r of the lowest terms on the lefthand side have been specialised, and hence also  $(d+1) - (m_{\delta} + r)$ of the highest terms. We can solve uniquely for the r lowest terms of  $\varepsilon(z)$  using the triangular system defined by considering coefficients of the powers  $z^a, z^{a+1}, \ldots, z^{a+r-1}$  on both sides of Equation (7), where  $z^a$  is the lowest monomial occurring on the lefthand side. One may also solve for the coefficients of the  $(d+1) - (m_{\delta} + r)$  highest powers uniquely using a similar triangular system. (Note that to ensure the triangular systems each have unique solutions we use here the fact that the constant term of  $G_0$  is non-zero, and the polynomial is of degree exactly  $n_{\delta} - 1$ .) Noticing that  $(d+1) - (m_{\delta} + r) + r = (d+1 - m_{\delta}) \ge d - n_{\delta} + 2$ , we see that all the coefficients of  $\varepsilon$  have been accounted for. However, if  $d+1-m_{\delta} > d-n_{\delta}+2$  (i.e.  $n_{\delta} > m_{\delta}+1$ ) there will be some "overlap", and the two triangular systems might not have a common solution. In this case there can be no solution to the Equation (7). If an  $\varepsilon(z)$  does exist which satisfies Equation (8) then the remaining coefficients of  $G_{k_{\delta}}$  can now be computed uniquely. Having computed the only possible solution of (7) for  $G_{k_{\delta}}$  we can substitute this into Equation (6) and recover  $H_{k_{\delta}}$  directly. More precisely compute

$$\frac{(F_{k_{\delta}} - \sum_{j=1}^{k_{\delta}-1} G_j H_{k_{\delta}-j}) - G_{k_{\delta}} H_0}{G_0}.$$
(9)

If its coefficients match with the known coefficients of  $H_{k\delta}$ then we have successfully extended the partial factorisation; otherwise we know no extension exists.

These computations can be done in time quadratic in the degree of the largest polynomial occurring in the above equations. Since all polynomials are Newton polytopes which are line segments lying within Newt(f) this is certainly quadratic in #Newt(f). (In fact, the running time is most closely related to the length of the side  $n_{\delta}$  from which we are performing the lifting step.) This completes the proof.  $\Box$ 

Theorem 7 may now be proved in a straightforward manner: Specifically, one first shows that for any partial factorisation extending a coprime dominating edges factorisation, there exists at most one full factorisation extending it, and this may be efficiently found. This is proved by induction on the number of unspecialised coefficients in the partial factorisation using Lemma 9. Theorem 7 then follows easily as a special case.

# 7. THE ALGORITHM

We now gather everything together and state our algorithm. We shall present it in an unadorned form, omitting detail on how to perform the more straightforward subroutines.

#### Algorithm 10

Input: A polynomial  $f \in \mathbb{F}[x, y]$ . Output: A factorisation of f or "failure".

Step A: Compute an irredundant dominating set  $\Gamma$  of Newt(f). For this choice of  $\Gamma$ , compute all coprime ( $\Gamma$ ; Q, R)-boundary factorisations of f, i.e., coprime partial boundary factorisations relative to the summands Q and R and the dominating set  $\Gamma$ . Here Q and R range over the summand pairs of Newt(f).

Step B: By repeatedly applying the method in the proof of Lemma 9, lift each coprime dominating edges factorisation of f as far as possible. If any of these lift to a full factorisation

output this factorisation and halt. If none of them lift to a full factorisation then output "failure".

Step A can be accomplished using a summand finding algorithm, an algorithm for finding dominating sets, and a univariate polynomial factorisation algorithm. A detailed description of these stages of the algorithm is given in the report [1]. For now, we just note that the summand finding algorithm is just a minor modification of the summand counting algorithm given in [3, Algorithm 17].

The algorithm is certainly correct, for it fails except when it finds a factor using the equations in Lemma 9. On the running time, using Theorem 7 lifting from each coprime dominating edges factorisation can be done in time polynomial (in fact cubic) in #Newt(f). However, although one can find such a dominating edges factorisation efficiently, the number of them may be exponential in the degree. In practice we recommend that a relative small number of dominating edges factorisations are tried before the polynomial is randomised and one resorts to other "dense polynomial" techniques.

The algorithm will always succeed when one starts with a dominating set  $\Gamma$  of Newt(f) such that the polynomials  $f_0^{\delta}, \delta \in \Gamma$ , are all squarefree (up to a monomial factor). Precisely, if the algorithm outputs "failure" one knows that in fact the polynomial f is irreducible, and otherwise the algorithm will output a factor. One might call polynomials for which such sets exist nice. This algorithm should be compared with the standard method of factoring "nice" polynomials using Hensel lifting. Precisely, in the literature a bivariate polynomial of total degree n which is squarefree upon reduction modulo y is often called "nice". The standard Hensel lifting algorithm will factor "nice" bivariate polynomials, on average very quickly [4], although in exponential time in the worst case. Notice a "nice" polynomial would be one whose Newton polytope has "lower boundary" a single edge of length n which is squarefree. The above algorithm factors not just these polynomials, but also any polynomials which have a "squarefree dominating set". (The algorithm also includes as a special case that given in Wan [10], where one "lifts downward" from the edge joining (n, 0) and (0, n).

# 8. EXAMPLES AND IMPLEMENTATION

## 8.1 Example

Suppose we want to factor the following binary polynomial  $f = x^{12} + x^{19} + (x^{10} + x^{11} + x^{13})y + (x^8 + x^9 + x^{12} + x^{17})y^2 + x^7y^3 + (x^4 + x^{11})y^4 + (x^2 + x^5 + x^{10})y^5 + y^6 + x^{10}y^8 + (x^8 + x^{11})y^9 + x^6y^{10} + x^9y^{12} + x^{15}y^{16}$ , with Newton polytope pictured in Figure 2, where a star indicates a nonzero term of f. Newt(f) is found to have three non-trivial decompositions, and eight irredundant dominating sets. None of these sets have edge polynomials which are all squarefree; however, fortunately we are still able to lift successfully from one of the coprime partial boundary factorisations. Specifically, consider the decomposition Newt(f) = Q + R, where Q and R are as in Figures 3 and 4, with generic polynomials g and h. The dominating edges of Newt(f) which allow a coprime edge factorisation are the two on the lower boundary of the polygon. Details of the lifting process are given in [1]. For now, we just note that the lines drawn in the interior of the polygons in Figures 3 and 4 indicate the first few layers

of coefficients which are revealed during the lifting, and the lines in the interior of Newt(f) the known coefficients of f which are used to do this.



Figure 2: Newton polytope of f



Figure 3: Newton polytope Q of the generic polynomial g

## 8.2 Implementation

We have developed a preliminary implementation of the algorithm with the aim of demonstrating how it would work for bivariate polynomials over  $\mathbb{F}_2$ . The work was carried out at the Oxford University Supercomputing Centre (OSC) on the Oswell machine, using an UltraSPARC III processor running at about 122 Mflop/s and with 2 GBytes of memory. The implementation was written using a combination of C and Magma programs, and was divided into three phases. In the first phase, the input polynomial is read and its Newton



Figure 4: Newton polytope R of the generic polynomial h

polytope computed using the asymptotically fast Graham's algorithm for computing convex hulls [8]. In that phase we also compute all irredundant dominating sets, and output the edge polynomials. In the second phase, a Magma program invokes a univariate factorisation algorithm to perform the partial boundary factorisations, and the results are directed into the third phase program. In this last phase, a search for coprime dominating edges factorisations is performed, and when appropriate, the lifting process is started. The polynomial arithmetic was performed using classical multiplication and division, and the triangular systems were solved using dense Gaussian elimination over  $\mathbb{F}_2$ .

We generated a number of random experiments with total degree reaching d = 2000. In all these cases, the input polynomial f was constructed by multiplying two random polynomials q and h of degree d/2 each with a given number of non-zero terms. Specifically, for each polynomial the given number of exponent vectors  $(e_1, e_2)$  were chosen uniformly at random subject to  $0 \leq e_1 + e_2 \leq d/2$ . These vectors always included ones of the form  $(e_1, 0), (0, e_2)$  and  $(e_3, (d/2) - e_3)$  to ensure the polynomial was of the correct degree and had no monomial factor. As the polynomials chosen were sparse the corresponding Newton polytopes had very few edges. In all these cases, the components of edge vectors of Newt(f) had a very small gcd, so that the edges had few integral points and consequently the polygon itself had very few summands. Table 1 gives the running times of the total factorisation process to find at least one non-trivial factor involving all three phases described above. Here s is the number of non-zero terms of the input polynomial f; #Newt(f), #Newt(g), and #Newt(h) are the total number of lattice points in Newt(f), Newt(g) and Newt(h) respectively; and t is the total running time in seconds. The actual polynomials f, g and h in each of the cases are listed in [1].

## 9. CONCLUSION

In this paper we have investigated a new approach for bivariate polynomial factorisation based on a study of New-

Table 1: Run time data for random experiments.

d	s	#Newt $(f)$	$\#\operatorname{Newt}(g)$	#Newt $(h)$	t
50	14	561	166	50	2.3
100	16	2234	472	222	11.6
500	15	52940	12758	11282	21.5
1000	30	206461	28582	56534	42.9
2000	28	848849	133797	132932	410.7

ton polytopes and generalised Hensel lifting. In standard Hensel lifting, one lifts a factorisation from a single edge, and uniqueness can be ensured by randomising the polynomial to enforce coprimality conditions and by making sure the edge being lifted from is sufficiently long. This randomisation is by substitution of linear forms which destroys the sparsity of the input polynomial. Our main theoretical contribution is to show how uniqueness may be ensured in the bivariate case, only under certain coprimality conditions, and without restrictions on the lengths of the edges. For certain classes of sparse polynomials, namely those whose Newton polytopes have few Minkowski decompositions, this gives a practical new approach which greatly improves upon Hensel lifting. As with Hensel lifting, our method has an exponential worst-case running time; however, we demonstrated the practicality of our algorithm on several randomly chosen composite and sparse binary polynomials of high degree.

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