# **Diversification Improves Interpolation**

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# **ABSTRACT**

We consider the problem of interpolating an unknown multivariate polynomial with coefficients taken from a finite field or as numerical approximations of complex numbers. Building on the recent work of Garg and Schost, we improve on the best-known algorithm for interpolation over large finite fields by presenting a Las Vegas randomized algorithm that uses fewer black box evaluations. Using related techniques, we also address numerical interpolation of sparse polynomials with complex coefficients, and provide the first provably stable algorithm (in the sense of relative error) for this problem, at the cost of modestly more evaluations. A key new technique is a randomization which makes all coefficients of the unknown polynomial distinguishable, producing what we call a diverse polynomial. Another departure from most previous approaches is that our algorithms do not rely on root finding as a subroutine. We show how these improvements affect the practical performance with trial implementations.

#### **Categories and Subject Descriptors**

F.2.1 [Analysis of Algorithms and Problem Complexity]: Numerical Algorithms and Problems—*Computations on polynomials*; G.1.1 [Numerical Analysis]: Interpolation; G.4 [Mathematical Software]: Algorithm Design and Analysis

#### **General Terms**

Algorithms, Theory, Performance

#### **Keywords**

Sparse polynomials, symbolic-numeric, interpolation

# 1. INTRODUCTION

Polynomial interpolation is a long-studied and important problem in computer algebra and symbolic computation.

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Given a way to evaluate an unknown polynomial and an upper bound on its degree, the interpolation problem is to determine a representation for the polynomial. In *sparse* interpolation, we are also given an upper bound on the number of nonzero terms in the unknown polynomial, and the output is returned in the sparse (also *lacunary* or *supersparse*) representation, wherein only nonzero terms are explicitly stored.

Applications of sparse interpolation include the manipulation and factorization of multivariate polynomials and system solving (see, e.g., Canny et al. (1989); Kaltofen and Trager (1990); Díaz and Kaltofen (1995, 1998); Javadi and Monagan (2007, 2009). With the advent of hybrid symbolic-numeric algorithms for (systems of) multivariate polynomials with approximate coefficients, we find applications of approximate sparse interpolation, in particular for solving nonlinear systems of equations (see, e.g., Sommese et al. (2001, 2004); Stetter (2004)) and factoring approximate multivariate polynomials (see, e.g., Kaltofen et al. (2008)).

Sparse interpolation is also a non-trivial generalization of the *polynomial identity testing* problem, determining if an unknown polynomial given by an algebraic circuit is identically zero. A relevant result in our setting is Bläser et al. (2009); see also the recent survey by Saxena (2009).

Here we examine the sparse interpolation problem in two settings which have received recent attention: when the coefficients are elements of finite fields (particularly large finite fields, over which we have no choice) and when they are approximations to complex numbers. We give improvements over the state of the art in both cases, and demonstrate our new algorithms in practice with an implementation in C++.

#### 1.1 Problem definition

A multivariate polynomial  $f \in \mathsf{F}[x_1,\ldots,x_n]$  with coefficients in a field  $\mathsf{F}$  is t-sparse for some  $t \in \mathbb{N}$  if f has at most t nonzero terms; that is, f can be written

$$f = \sum_{i=1}^{t} c_i x_1^{e_{i1}} x_2^{e_{i2}} \cdots x_n^{e_{in}}$$

for coefficients  $c_i \in \mathsf{F}$  and exponent tuples  $(e_{i1},\ldots,e_{in}) \in \mathbb{N}^n$  for  $1 \leq i \leq t$ . If each  $e_{ij} < d$ , then the size of this representation is O(t) field elements plus  $O(tn \log d)$  bits. We seek algorithms which are polynomial-time in this size.

Let  $f \in \mathsf{F}[x_1,\ldots,x_n]$ . A black box for f is a function which takes as input a vector  $(a_1,\ldots,a_n) \in \mathsf{F}^n$  and produces  $f(a_1,\ldots,a_n) \in \mathsf{F}$ . The cost of the black box is the number of operations in  $\mathsf{F}$  required to evaluate it at a given input.

Clausen et al. (1991) showed that, if only evaluations over F are allowed, then for some instances at least  $\Omega(n^{\log t})$  black

box probes are required. Hence if we seek polynomial-time algorithms, we must extend the capabilities of the black box. To this end, Díaz and Kaltofen (1998) introduced the extended domain black box which is capable of evaluating  $f(b_1, \ldots, b_n) \in E$  for any  $(b_1, \ldots, b_n) \in E^n$  where E is any extension field of F. That is, we can change every black box operation to work over an extension field, usually paying an extra cost proportional to the size of the extension.

Motivated by the case of black boxes that are division-free algebraic circuits, we will use the following model which we believe to be fair and cover all previous relevant results. Here  $\mathsf{M}(m)$  is the number of field operations required to multiply two univariate polynomials with degrees less than m, and O(m) represents any function bounded by  $m(\log m)^{O(1)}$ .

DEFINITION 1.1. Let  $f \in \mathsf{F}[x_1, \ldots, x_n]$  and  $\ell > 0$ . A remainder black box for f with size  $\ell$  is a procedure which, given any monic square-free polynomial  $g \in \mathsf{F}[y]$  with  $\deg g = m$ , and any  $h_1, \ldots, h_n \in \mathsf{F}[y]$  with each  $\deg h_i < m$ , produces  $f(h_1, \ldots, h_n)$  rem g using at most  $\ell \cdot \mathsf{M}(m)$  operations in  $\mathsf{F}$ .

This definition is general enough to cover the algorithms we know of over finite fields, and we submit that the cost model is fair to the standard black box, extended domain black box, and algebraic circuit settings. The model also makes sense over complex numbers as well, as we will see.

# 1.2 Interpolation over finite fields

We first summarize previously known univariate interpolation algorithms when F is a finite field with q elements and identify our new contributions. For now, let  $f \in \mathbb{F}_q[x]$  have degree less than d and sparsity t. We will assume we have a remainder black box for f with size  $\ell$ . Since field elements can be represented with  $O(\log q)$  bits, a polynomial-time algorithm will have cost polynomial in  $\ell$ , t,  $\log d$ , and  $\log q$ .

For the dense output representation, one can use the classical method of Newton/Waring/Lagrange to interpolate in  $O^{\tilde{}}(\ell d)$  time (von zur Gathen and Gerhard, 2003, §10.2).

The algorithm of Ben-Or and Tiwari (1988) for sparse polynomial interpolation, with improvements by Kaltofen and Yagati (1989), can be adapted to arbitrary finite fields. Unfortunately, these algorithms require t discrete logarithm computations in  $\mathbb{F}_q^*$ , whose cost is small if the field size q is chosen carefully (as in Kaltofen (2010)), but not in general. The best approach for arbitrary q seems to be Pollard's "kangaroo" algorithm (1978), using the fact that the discrete logs must fall in the range  $[0, \ldots, d-1]$ . The total cost of this method is  $O(t\ell + t^2 + t\sqrt{d})$ .

The current study builds most directly on the work of Garg and Schost (2009), who gave the first polynomial-time algorithm for sparse interpolation over an arbitrary finite field. Their algorithm works roughly as follows. For very small primes p, use the black box to compute f modulo  $x^p-1$ . A prime p is a "good prime" if and only if all terms of f are distinct modulo  $x^p-1$ . We do this for p in the range of roughly  $O(t^2 \log d)$  until there are sufficient good primes to recover the unique symmetric polynomial over  $\mathbb{Z}[y]$  whose roots are the exponents of nonzero terms in f. We then factor this polynomial to find those exponents, and correlate with any good prime image to determine the coefficients. The total cost is  $O\tilde{}(\ell t^4 \log^2 d)$  field operations. Using randomization, it is easy to reduce this to  $O\tilde{}(\ell t^3 \log^2 d)$ .

The coefficients of the symmetric integer polynomial in Garg & Schost's algorithm are bounded by  $O(d^t)$ , much

larger than the O(d) size of the exponents ultimately recovered. Our new algorithm over finite fields of size at least  $\Omega(t^2d)$  avoids evaluating the symmetric polynomial and finding its roots entirely. As a result, we reduce the total number of required evaluations and develop a randomized algorithm with cost  $O^-(\ell t^2 \log^2 d)$ , which is roughly quadratic in the input and output sizes. Since this can be deterministically verified in the same time, our algorithm (as well as the randomized version of Garg & Schost) is of the Las Vegas type.

These results are summarized in Table 1, where we assume in all cases that the field size q is "large enough". In the table, the "probe degree" refers to the degree of g in each evaluation of the remainder black box as defined above.

Two techniques are available to adapt a univariate interpolation algorithm to the multivariate case. The first is Kronecker substitution: given a remainder black box for an unknown  $f \in \mathsf{F}[x_1,\ldots,x_n]$  with max degree less than d, we can easily construct a remainder black box for the univariate polynomial  $f(x,x^d,x^{d^2},\ldots,x^{d^{n-1}}) \in \mathsf{F}[x]$ , whose terms correspond one-to-one with terms of f. The cost is that of the univariate algorithm with sparsity t and degree  $d^n$ .

The other method for constructing a multivariate interpolation algorithm is due to Zippel (1990). The technique is inherently probabilistic of the Monte Carlo type and works variable-by-variable, at each step solving a number of transposed Vandermonde systems and making O(t) calls to a univariate interpolation algorithm. Say this underlying algorithm uses  $\rho(d,t)$  black box evaluations of degree  $\Delta(d,t)$  and  $\psi(d,t)$  other field operations to interpolate univariate polynomials with sparsity t and degrees less than d. Then the resulting algorithm using Zippel's method has total cost

$$O(nt\psi(d,t) + \ell nt\rho(d,t)\Delta(n,t))$$

to interpolate an n-variate t-sparse polynomial with max degree less than d. Zippel (1990) used the dense algorithm for univariate interpolation; using Ben-Or and Tiwari's algorithm instead was studied by Kaltofen and Lee (2003).

Grigoriev et al. (1990) give a parallel algorithm with small depth but which is not competitive in our model due to the large number of processors required. A practical parallel version of Ben-Or and Tiwari's algorithm has been developed by Javadi and Monagan (2010). Kaltofen et al. (1990) and Avendaño et al. (2006) present modular algorithms for interpolating polynomials with rational and integer coefficients, but their methods do not seem to apply to finite fields.

# 1.3 Approximate polynomial interpolation

Recently, a number of numerically-focussed sparse interpolation algorithms have been presented. The algorithm of Giesbrecht et al. (2009) is a numerical adaptation of Ben-Or and Tiwari (1988), which samples f at O(t) randomly chosen roots of unity  $\omega \in \mathbb{C}$  on the unit circle. In particular,  $\omega$  is chosen to have (high) order at least the degree, and a randomization scheme is used to avoid clustering of nodes which will cause dramatic ill-conditioning. A relatively weak theoretical bound is proven there on the randomized conditioning scheme, though experimental and heuristic evidence suggests it is much better in practice. Cuyt and Lee (2008) adapt Rutishauser's qd algorithm to alleviate the need for bounds on the partial degrees and the sparsity, but still evaluate at high-order roots of unity. Approximate sparse rational function interpolation is considered by Kaltofen and Yang (2007) and Kaltofen et al. (2007), using the Structured

|                  | Probes          | Probe degree    | Computation cost     | Total cost                    | Type          |
|------------------|-----------------|-----------------|----------------------|-------------------------------|---------------|
| Dense            | d               | 1               | $O^{}(d)$            | $O^{}(\ell d)$                | deterministic |
| Ben-Or & Tiwari  | O(t)            | 1               | $O(t^2 + t\sqrt{d})$ | $O(\ell t + t^2 + t\sqrt{d})$ | deterministic |
| Garg & Schost    | $O(t^2 \log d)$ | $O(t^2 \log d)$ | $O(t^4 \log^2 d)$    | $O(\ell t^4 \log^2 d)$        | deterministic |
| Randomized G & S | $O(t \log d)$   | $O(t^2 \log d)$ | $O(t^3 \log^2 d)$    | $O(\ell t^3 \log^2 d)$        | Las Vegas     |
| Ours             | $O(\log d)$     | $O(t^2 \log d)$ | $O(t^2 \log^2 d)$    | $O(\ell t^2 \log^2 d)$        | Las Vegas     |

**Table 1:** Sparse univariate interpolation over large finite fields, with black box size  $\ell$ , degree d, and t nonzero terms

Total Least Norm (STLN) method and, in the latter, randomization to improve conditioning. Approximate sparse interpolation is also considered for integer polynomials by Mansour (1995), where a polynomial-time algorithm is presented in quite a different model from ours. In particular the evaluation error is absolute (not relative) and the complexity is sensitive to the bit length of the integer coefficients.

Our new algorithm for approximate sparse interpolation is presented in Section 4. We provide the first algorithm which is provably numerically stable without heuristics or conjectures. We define an " $\epsilon$ -approximate black box" as one which evaluates an unknown t-sparse target polynomial  $f \in \mathbb{C}[x]$ , of degree d, with relative error at most  $\epsilon > 0$ . Our goal is to build a t-sparse polynomial g such that  $||f - g||_2 \le \epsilon ||f||_2$ . A bound on the degree and sparsity of the target polynomial, as well as  $\epsilon$ , must also be provided. After demonstrating that the problem is well-posed, we adapt our variant of the Garg and Schost (2009) algorithm for the approximate case, prove it is numerically accurate in terms of the relative error of the output, and analyze its cost. Our algorithm typically requires  $O(t^2 \log^2 d)$  evaluations at primitive roots of unity of order  $O(t^2 \log d)$  (as opposed to order d in previous approaches). We guarantee that it finds a t-sparse polynomial g such that  $||g-f||_2 \leq 2\epsilon ||f||_2$ . An experimental demonstration of the numerical robustness is given in Section 5.

Note that all these works evaluate the polynomial only on the unit circle. This is necessary because we allow and expect f to have very large degree, which would cause a catastrophic loss of precision at data points of non-unit magnitude. Similarly, we assume that the complex argument of evaluation points is exactly specified, which is again necessary because any error in the argument would be exponentially magnified by the degree.

# 2. SPARSE INTERPOLATION FOR GENERIC FIELDS

Here and for the remainder, we focus on interpolating an unknown t-sparse univariate polynomial  $f \in \mathsf{F}[x]$  with degree less than d. As discussed above, these algorithms can be adapted to multivariate interpolation using Kronecker substitution or Zippel's method.

Assume a fixed, unknown  $f \in \mathsf{F}[x]$  with sparsity at most t and degree at most d. We will use a remainder black box for f to evaluate  $f \operatorname{rem}(x^p-1)$  for small primes p. We say p is a "good prime" if the sparsity of  $f \operatorname{rem}(x^p-1)$  is the same as that of f itself — that is, none of the exponents are equivalent modulo p.

The following lemma shows the size of primes required to randomly choose good primes with high probability.

LEMMA 2.1. Let  $f \in \mathsf{F}[x]$  be a t-sparse polynomial with degree d, and let  $\lambda = \max \left( 21, \lceil \frac{5}{3}t(t-1) \ln d \rceil \right)$ . A prime

chosen at random in the range  $\lambda, \ldots, 2\lambda$  is a good prime for f with probability at least 1/2.

PROOF. Let  $e_1, \ldots, e_t$  be the exponents of nonzero terms in f. If p is a bad prime, then p divides  $(e_j - e_i)$  for some i < j. Each  $e_j - e_i \le d$ , so there can be at most  $\log_{\lambda} d = \ln d / \ln \lambda$  primes that divide each  $e_j - e_i$ . There are exactly  $\binom{t}{2}$  such pairs of exponents, so the total number of bad primes is at most  $(t(t-1) \ln d)/(2 \ln \lambda)$ .

From Rosser and Schoenfeld (1962, Corollary 3 to Theorem 2), the total number of primes in the range  $\lambda, \ldots, 2\lambda$  is at least  $3\lambda/(5 \ln \lambda)$  when  $\lambda \geq 21$ , which is at least  $t(t-1) \ln d / \ln \lambda$ , at least twice the number of bad primes.  $\square$ 

Now observe an easy case for the sparse interpolation problem. If a polynomial  $f \in F[x]$ , has all coefficients distinct; that is,  $f = \sum_{1 \le i \le t} c_i x^{e_i}$  and  $c_i = c_j \Rightarrow i = j$ , then we say f is diverse. To interpolate a diverse polynomial  $f \in F[x]$ , we first follow the method of Garg and Schost (2009) by computing  $f \operatorname{rem}(x^{p_i}-1)$  for "good primes"  $p_i$  such that the sparsity of  $f \operatorname{rem}(x^{p_i} - 1)$  is the same as that of f. Since f is diverse,  $f \operatorname{rem}(x^{p_i} - 1)$  is also diverse and in fact each modular image has the same set of coefficients. Using this fact, we avoid the need to construct and subsequently factor the symmetric polynomial in the exponents. Instead, we correlate like terms based on the (unique) coefficients in each modular image, then use simple Chinese remaindering to construct each exponent  $e_i$  from its image modulo each  $p_i$ . This requires only  $O(\log d)$  remainder black box evaluations at good primes, gaining a factor of t improvement over the randomized version of Garg and Schost (2009).

In the following sections, we will show how to choose an  $\alpha \in \mathsf{F}$  so that  $f(\alpha x)$  — which we can easily construct a remainder black box for — is diverse. With such a procedure, Algorithm 1 gives a Monte Carlo algorithm for interpolation over a general field.

Theorem 2.2. With inputs as specified, Algorithm 1 correctly computes the unknown polynomial f with probability at least  $1-\mu$ . The total cost in field operations (except for step 8) is

$$O\left(\ell \cdot \left(\frac{\log D}{\log T + \log\log D} + \log\frac{1}{\mu}\right) \cdot \mathsf{M}\left(T^2 \log D\right)\right).$$

PROOF. The for loop on line 3 searches for the true sparsity t and a single good prime  $\varrho$ . Since each prime p in the given range is good with probability at least 1/2 by Lemma 2.1, the probability of failure at this stage is at most  $\mu/3$ .

The for loop on line 12 searches for and uses sufficiently many good primes to recover the exponents of f. The product of all the good primes must be at least D, and since each prime is at least  $\lambda$ , at least  $(\ln D)/(\ln \lambda)$  good primes are required.

## Algorithm 1: Generic interpolation

```
Input: \mu \in \mathbb{R}_{>0}, T, D, q \in \mathbb{N}, and a remainder black
                   box for unknown T-sparse f \in F[x] with
      Output: t \in \mathbb{N}, e_1, \dots, e_t \in \mathbb{N}, and c_1, \dots, c_t \in \mathsf{F} such that f = \sum_{1 \leq i \leq t} c_i x^{e_i}
  1 t \leftarrow 0
  2 \lambda \leftarrow \max\left(21, \left\lceil \frac{5}{3}T(T-1)\ln D\right\rceil\right)
  3 for \lceil \log_2(3/\mu) \rceil primes p \in \{\lambda, \dots, 2\lambda\} do
            Use black box to compute f_p = f(x) \operatorname{rem}(x^p - 1)
            if f_p has more than t terms then
  6
                 t \leftarrow \text{sparsity of } f_p
  8 \alpha \leftarrow element of F s.t. \Pr[f(\alpha x) \text{ not diverse}] < \mu/3
  9 g_{\varrho} \leftarrow f(\alpha x) \operatorname{rem}(x^{\varrho} - 1)
10 c_1, \ldots, c_t \leftarrow \text{nonzero coefficients of } g_{\varrho}
11 e_1, \ldots, e_t \leftarrow 0
12 for \lceil 2 \ln(3/\mu) + 4(\ln D)/(\ln \lambda) \rceil primes p \in \{\lambda, \dots, 2\lambda\}
            Use black box to compute g_p = f(\alpha x) \operatorname{rem}(x^p - 1)
13
           if g_p has exactly t nonzero terms then
14
15
                 for i = 1, ..., t do Update e_i with exponent of
                c_i in g_p modulo p via Chinese remaindering
16 for i=1,\ldots,t do c_i \leftarrow c_i \alpha^{-e_i}
17 return f(x) = \sum_{1 \leq i \leq t} c_i x^{e_i}
```

Let  $n = \lceil 2 \ln(3/\mu) + 4(\ln D)/(\ln \lambda) \rceil$  be the number of primes sampled in this loop, and  $k = \lceil (\ln D)/(\ln \lambda) \rceil$  the number of good primes required. We can derive that

$$\exp\left(\frac{-2\left(n/2-k\right)^2}{n}\right) < \frac{\mu}{3}.$$

Then from Hoeffding's Inequality (1963), the probability of encountering fewer than k good primes is less than  $\mu/3$ .

Therefore the total probability of failure is at most  $\mu$ . For the cost analysis, the dominating cost will be the modular black box evaluations in the last for loop. The number of evaluations in this loop is  $O(\log(1/\mu) + (\log D)/(\log \lambda))$ , and each evaluation has cost  $O(\ell \cdot M(\lambda))$ . Since

$$(\log D)/(\log \lambda) \in \Theta((\log D)/(\log T + \log\log D)),$$

the complexity bound is correct as stated.  $\ \ \Box$ 

In case the bound T on the number of nonzero terms is very bad, we could choose a smaller value of  $\lambda$  based on the true sparsity t before line 8, improving the cost of the remainder of the algorithm.

In addition, as our bound on possible number of "bad primes" seems to be quite loose, a more efficient approach in practice would be to replace the for loop on line 12 with one that starts with a prime much smaller than  $\lambda$  and incrementally searches for the next larger primes until the product of all good primes is at least D. We could choose the lower bound to start searching from based on lower bounds on the birthday problem. That is, assuming (falsely) that the exponents are randomly distributed modulo p, start with the least p that will have no exponents collide modulo p with high probability. This would yield an algorithm more sensitive to the true bound on bad primes, but unfortunately gives a worse formal cost analysis.

# 3. SPARSE INTERPOLATION OVER FINITE FIELDS

# 3.1 Diversification

For a prime power q, we use  $\mathbb{F}_q$  to denote the finite field with q elements. In order to use Algorithm 1 over  $\mathbb{F}_q[x]$ , we must find an  $\alpha$  so that  $f(\alpha x)$  is diverse. A surprisingly simple trick works: evaluating  $f(\alpha x)$  for a random nonzero  $\alpha \in \mathbb{F}_q$ .

THEOREM 3.1. For  $q \geq T(T-1)D$  and any T-sparse polynomial  $f \in \mathbb{F}_q[x]$  with deg f < D, if  $\alpha$  is chosen uniformly at random from  $\mathbb{F}_q^*$ , the probability that  $f(\alpha x)$  is diverse is at least 1/2.

PROOF. Let  $t \leq T$  be the exact number of nonzero terms in f, and write  $f = \sum_{1 \leq i \leq t} c_i x^{e_i}$ , with nonzero coefficients  $c_i \in \mathbb{F}_q^*$  and  $e_1 < e_2 < \cdots < e_t$ . So the ith coefficient of  $f(\alpha x)$  is  $c_i \alpha^{e_i}$ .

If  $f(\alpha x)$  is not diverse, then we must have  $c_i \alpha^{e_i} = c_j \alpha^{e_j}$  for some  $i \neq j$ . Therefore consider the polynomial  $A \in \mathbb{F}_q[y]$  defined by

$$A = \prod_{1 \le i < j \le t} (c_i y^{e_i} - c_j y^{e_j}).$$

We see that  $f(\alpha x)$  is diverse if and only if  $A(\alpha) \neq 0$ , hence the number of roots of A over  $\mathbb{F}_q$  is exactly the number of unlucky choices for  $\alpha$ .

The polynomial A is the product of exactly  $\binom{t}{2}$  binomials, each of which has degree less than D. Therefore

$$\deg A < \frac{T(T-1)D}{2},$$

and this also gives an upper bound on the number of roots of A. Hence  $q-1 \geq 2 \deg A$ , and at least half of the elements of  $\mathbb{F}_q^*$  are not roots of A, yielding the stated result.  $\square$ 

Using this result, given a black box for f and the exact sparsity t of f, we can find an  $\alpha \in \mathbb{F}_q$  such that  $f(\alpha x)$  is diverse by sampling random values  $\alpha \in \mathbb{F}_q$ , evaluating  $f(\alpha x)$  rem  $x^p-1$  for a single good prime p, and checking whether the polynomial is diverse. With probability at least  $1-\mu$ , this will succeed in finding a diversifying  $\alpha$  after at most  $\lceil \log_2(1/\mu) \rceil$  iterations. Therefore we can use this approach in Algorithm 1 with no effect on the asymptotic complexity.

#### 3.2 Verification

So far, Algorithm 1 over a finite field is probabilistic of the Monte Carlo type; that is, it may give the wrong answer with some controllably-small probability. To provide a more robust Las Vegas probabilistic algorithm, we require only a fast way to check that a candidate answer is in fact correct. To do this, observe that given a modular black box for an unknown T-sparse  $f \in \mathbb{F}_q[x]$  and an explicit T-sparse polynomial  $g \in \mathbb{F}_q[x]$ , we can construct a modular black box for the 2T-sparse polynomial f-g of their difference. Verifying that f=g thus reduces to the well-studied problem of deterministic polynomial identity testing.

The following algorithm is due to Bläser et al. (2009) and provides this check in essentially the same time as the interpolation algorithm; we restate it in Algorithm 2 for completeness and to use our notation.

#### Algorithm 2: Verification over finite fields

**Input**:  $T, D, q \in \mathbb{N}$  and modular black box for unknown T-sparse  $f \in \mathbb{F}_q[x]$  with deg  $f \leq D$ Output: ZERO iff f is identically zero

1 for the least  $(T-1)\log_2 D$  primes p do

Use black box to compute  $f_p = f \operatorname{rem}(x^p - 1)$ 

if  $f_p \neq 0$  then return NONZERO

4 return ZERO

Theorem 3.2. Algorithm 2 works correctly as stated and uses at most

$$O(\ell T \log D \cdot \mathsf{M} (T \log D \cdot (\log T + \log \log D)))$$

field operations.

PROOF. See (Bläser et al., 2009, Theorem 14).  $\square$ 

This provides all that we need to prove the main result of this section:

Theorem 3.3. Given  $q \geq T(T-1)D+1$ , any  $T, D \in \mathbb{N}$ , and a modular black box for unknown T-sparse  $f \in \mathbb{F}_q[x]$ with deg  $f \leq D$ , there is an algorithm that always produces the correct polynomial f and with high probability uses only  $O^{\sim}(\ell T^2 \log^2 D)$  field operations.

PROOF. Use Algorithms 1 and 2 with  $\mu = 1/2$ , looping as necessary until the verification step succeeds. With high probability, only a constant number of iterations will be necessary, and so the cost is as stated.  $\square$ 

For the small field case, when  $q \in O(T^2D)$ , the obvious approach would be to work in an extension E of size  $O(\log T + \log D)$  over  $\mathbb{F}_q$ . Unfortunately, this would presumably increase the cost of each evaluation by a factor of  $\log D$ , potentially dominating our factor of T savings compared to the randomized version of Garg and Schost (2009) when the unknown polynomial has very few terms and extremely high degree.

In practice, it seems that a much smaller extension than this is sufficient in any case to make each  $gcd(e_i - e_i, q - 1)$ small compared to q-1, but we do not yet know how to prove any tighter bound in the worst case.

# **ALGORITHMS FOR APPROXIMATE** SPARSE INTERPOLATION

In this section we consider the problem of interpolating an approximate sparse polynomial  $f \in \mathbb{C}[x]$  from evaluations on the unit circle. We will generally assume that f is t-sparse:

$$f = \sum_{1 \le i \le t} c_i x^{e_i} \text{ for } c_i \in \mathbb{C} \text{ and } e_1 < \dots < e_t = d.$$
 (4.1)

We require a notion of size for such polynomials, and define the coefficient 2-norm of  $f = \sum_{0 \le i \le d} f_i x^i$  as

$$||f||_2 = \sqrt{\sum_{0 \le i \le d} |f_i|^2}.$$

The following identity relates the norm of evaluations on the unit circle and the norm of the coefficients. As in Section 2, for  $f \in \mathbb{C}[x]$  is as in (4.1), we say that a prime p is a good prime for f if  $p \nmid (e_i - e_j)$  for all  $i \neq j$ .

LEMMA 4.1. Let  $f \in \mathbb{C}[x]$ , p a good prime for f, and  $\omega \in \mathbb{C}$  a pth primitive root of unity. Then

$$||f||_2^2 = \frac{1}{p} \sum_{0 \le i \le n} |f(\omega^i)|^2.$$

See Giesbrecht and Roche (2010, Theorem 2.9).

We can now formally define the approximate sparse univariate interpolation problem.

Definition 4.2. Let  $\epsilon > 0$  and assume there exists an unknown t-sparse  $f \in \mathbb{C}[x]$  of degree at most D. An  $\epsilon$ approximate black box for f takes an input  $\xi \in \mathbb{C}$  and produces a  $\gamma \in \mathbb{C}$  such that  $|\gamma - f(\xi)| \leq \epsilon |f(\xi)|$ .

The approximate sparse univariate interpolation problem is then as follows: given  $D, T \in \mathbb{N}$  and  $\delta \geq \epsilon > 0$ , and an  $\epsilon$ approximate black box for an unknown T-sparse polynomial  $f \in \mathbb{C}[x]$  of degree at most D, find a T-sparse polynomial  $g \in \mathbb{C}[x]$  such that  $||f - g||_2 \le \delta ||g||_2$ .

The following theorem shows that t-sparse polynomials are well-defined by good evaluations on the unit circle.

Theorem 4.3. Let  $\epsilon > 0$  and  $f \in \mathbb{C}[x]$  be a t-sparse polynomial. Suppose there exists a t-sparse polynomial  $g \in \mathbb{C}[x]$ such that for a prime p which is good for f and f - g, and pth primitive root of unity  $\omega \in \mathbb{C}$ , we have

$$|f(\omega^i) - g(\omega^i)| \le \epsilon |f(\omega^i)| \quad \text{for } 0 \le i < p.$$

Then  $||f - g||_2 \le \epsilon ||f||_2$ . Moreover, if  $g_0 \in \mathbb{C}[x]$  is formed from g by deleting all the terms not in the support of f, then  $||f - g_0||_2 \le 2\epsilon ||f||_2$ .

Proof. Summing over powers of  $\omega$  we have

$$\sum_{0 \leq i < p} |f(\omega^i) - g(\omega^i)|^2 \leq \epsilon^2 \sum_{0 \leq i < p} |f(\omega^i)|^2.$$

Thus, since p is a good prime for both f-g and f, and using Lemma 4.1,  $p \cdot ||f - g||_2^2 \le \epsilon^2 \cdot p \cdot ||f||_2^2$  and  $||f - g||_2 \le \epsilon ||f||_2$ . Since  $g - g_0$  has no support in common with f,

$$||g - g_0||_2 \le ||f - g||_2 \le \epsilon ||f||_2$$
.

Thus

$$||f - g_0||_2 = ||f - g + (g - g_0)||_2$$
  

$$\leq ||f - g||_2 + ||g - g_0||_2 \leq 2\epsilon ||f||_2. \quad \Box$$

In other words, any t-sparse polynomial whose values are very close to f must have the same support except possibly for some terms with very small coefficients.

# Constructing an $\epsilon$ -approximate remainder black box

Assume that we have chosen a good prime p for a t-sparse  $f \in F[x]$ . Our goal in this subsection is a simple algorithm and numerical analysis to accurately compute  $f \text{ rem } x^p - 1$ .

Assume that  $f \operatorname{rem} x^p - 1 = \sum_{0 \le i < p} \bar{b_i} x^i$  exactly. For a primitive pth root of unity  $\omega \in \mathbb{C}$ , let  $V(\omega) \in \mathbb{C}^{p \times p}$  be the Vandermonde matrix built from the points  $1, \omega, \dots, \omega^{p-1}$ . Recall that  $V(\omega) \cdot (b_0, \dots, b_{p-1})^T = (f(\omega^0), \dots, f(\omega^{p-1}))^T$  and  $V(\omega^{-1}) = p \cdot V(\omega)^{-1}$ . Matrix vector product by such Vandermonde matrices is computed very quickly and in a numerically stable manner by the Fast Fourier Transform (FFT).

#### Algorithm 3: Approximate Remainder

Input: An  $\epsilon$ -approximate black box for the unknown t-sparse  $f \in \mathbb{C}[x]$ , and  $p \in \mathbb{N}$ , a good prime for fOutput:  $h \in \mathbb{C}[x]$  such that  $\|(f \operatorname{rem} x^p - 1) - h\|_2 \le \epsilon \|f\|_2.$ 1  $w \leftarrow (f(\omega^0), \dots, f(\omega^{p-1})) \in \mathbb{C}^p$  computed using the  $\epsilon$ -approximate black box for f2  $u \leftarrow (1/p) \cdot V(\omega^{-1})w \in \mathbb{C}^p$  using the FFT algorithm
3 return  $h = \sum_{0 \le i < p} u_i x^i \in \mathbb{C}[x]$ 

Theorem 4.4. Algorithm 3 works as stated, and produces an h such that  $\|(f \operatorname{rem} x^p - 1) - h\|_2 \le \epsilon \|f\|_2$ . It requires  $O(p \log p)$  floating point operations and p evaluations of the black box.

PROOF. Because f and f rem  $x^p-1$  have exactly the same coefficients (p is a good prime for f), they have exactly the same norm. The FFT in Step 2 is accomplished in  $O(p\log p)$  floating point operations. This algorithm is numerically stable since  $(1/\sqrt{p})\cdot V(\omega^{-1})$  is unitary. That is, assume  $v=(f(\omega_0),\ldots,f(\omega^{p-1}))\in\mathbb{C}^p$  is the vector of exact evaluations of f, so  $\|v-w\|_2 \le \epsilon \|v\|_2$  by the black box specification. Then, using the fact that  $\|v\|_2 = \sqrt{p} \|f\|_2$ ,

$$\begin{aligned} \left\| (f \operatorname{rem} x^{p-1}) - h \right\|_{2} &= \left\| \frac{1}{p} V(\omega^{-1}) v - \frac{1}{p} V(\omega^{-1}) w \right\|_{2} \\ &= \frac{1}{\sqrt{p}} \left\| \frac{1}{\sqrt{p}} V(\omega^{-1}) \cdot (v - w) \right\|_{2} = \frac{1}{\sqrt{p}} \left\| v - w \right\|_{2}, \end{aligned}$$

which is at most  $(\epsilon/\sqrt{p}) \|v\|_2 = \epsilon \|f\|_2$ .  $\square$ 

# 4.2 Creating $\epsilon$ -diversity

First, we extend the notion of polynomial diversity to the approximate case.

DEFINITION 4.5. Let  $f \in \mathbb{C}[x]$  be a t-sparse polynomial as in (4.1) and  $\delta \geq \epsilon > 0$  such that  $|c_i| \geq \delta \|f\|_2$  for  $1 \leq i \leq t$ . The polynomial f is said to be  $\epsilon$ -diverse if and only if every pair of distinct coefficients is at least  $\epsilon \|f\|_2$  apart. That is, for every  $1 \leq i < j \leq t$ ,  $|c_i - c_j| \geq \epsilon \|f\|_2$ .

Intuitively, if  $(\epsilon/2)$  corresponds to the machine precision, this means that an algorithm can reliably distinguish the coefficients of a  $\epsilon$ -diverse polynomial. We now show how to choose a random  $\alpha$  to guarantee  $\epsilon$ -diversity.

Theorem 4.6. Let  $\delta \geq \epsilon > 0$  and  $f \in \mathbb{C}[x]$  a t-sparse polynomial whose non-zero coefficients are of magnitude at least  $\delta \|f\|_2$ . If s is a prime satisfying s > 12 and

$$t(t-1) \le s \le 3.1 \frac{\delta}{\epsilon},$$

then for  $\zeta = \mathbf{e}^{2\pi \mathbf{i}/s}$  an s-PRU and  $k \in \mathbb{N}$  chosen uniformly at random from  $\{0, 1, \ldots, s-1\}$ ,  $f(\zeta^k x)$  is  $\epsilon$ -diverse with probability at least  $\frac{1}{2}$ .

PROOF. For each  $1 \leq i \leq t$ , write the coefficient  $c_i$  in polar notation to base  $\zeta$  as  $c_i = r_i \zeta^{\theta_i}$ , where each  $r_i$  and  $\theta_i$  are nonnegative real numbers and  $r_i \geq \delta \|f\|_2$ .

Suppose  $f(\zeta^k x)$  is not  $\epsilon$ -diverse. Then there exist indices  $1 \le i < j \le t$  such that

$$\left| r_i \zeta^{\theta_i} \zeta^{ke_i} - r_j \zeta^{\theta_j} \zeta^{ke_j} \right| \le \epsilon \|f\|_2.$$

Algorithm 4: Adaptive diversification

Input:  $\epsilon$ -approximate black box for f, known good prime p, known sparsity t

**Output**:  $\zeta, k$  such that  $f(\zeta^k x)$  is  $\epsilon$ -diverse, or FAIL

 $1 s \leftarrow 1, \delta \leftarrow \infty, f_p \leftarrow 0$ 

**2** while  $s \le t^2$  and  $\#\{coeffs \ c \ of \ f_s \ s.t. \ |c| \ge \delta\} < t \ do$ 

**3**  $s \leftarrow \text{least prime} \geq 2s$ 

4  $\zeta \leftarrow \exp(2\pi \mathbf{i}/s)$ 

5  $k \leftarrow \text{random integer in } \{0, 1, \dots, s-1\}$ 

6 Compute  $f_s = f(\zeta^k x) \operatorname{rem} x^p - 1$ 

7  $\delta \leftarrow$  least number s.t. all coefficients of  $f_s$  at least  $\delta$  in absolute value are pairwise  $\epsilon$ -distinct

8 if  $\delta > 2\epsilon$  then return FAIL

9 else return  $\zeta^k$ 

Because  $\min(r_i, r_j) \geq \delta \|f\|_2$ , the left hand side is at least  $\delta \|f\|_2 \cdot |\zeta^{\theta_i + ke_i} - \zeta^{\theta_j + ke_j}|$ . Dividing out  $\zeta^{\theta_j + ke_i}$ , we get

$$\left| \zeta^{\theta_i - \theta_j} - \zeta^{k(e_j - e_i)} \right| \le \frac{\epsilon}{\delta}.$$

By way of contradiction, assume that there exist distinct choices of k that satisfy the above inequality, say  $k_1, k_2 \in \{0, \ldots, s-1\}$ . Since  $\zeta^{\theta_i - \theta_j}$  and  $\zeta^{e_j - e_i}$  are a fixed powers of  $\zeta$  not depending on the choice of k, this means

$$\left| \zeta^{k_1(e_j - e_i)} - \zeta^{k_2(e_j - e_i)} \right| \le 2\frac{\epsilon}{\delta}.$$

Because s is prime,  $e_i \neq e_j$ , and we assumed  $k_1 \neq k_2$ , the left hand side is at least  $|\zeta - 1|$ . Observe that  $2\pi/s$ , the distance on the unit circle from 1 to  $\zeta$ , is a good approximation for this Euclidean distance when s is large. In particular, since s > 12,

$$\frac{\left|\zeta-1\right|}{2\pi/s} > \frac{\sqrt{2}\left(\sqrt{3}-1\right)/2}{\pi/6},$$

and therefore  $|\zeta - 1| > 6\sqrt{2}(\sqrt{3} - 1)/s > 6.2/s$ , which from the statement of the theorem is at least  $2\epsilon/\delta$ . This is a contradiction, and therefore the assumption was false; namely, there is at most one choice of k such that the i'th and j'th coefficients collide.

Then, since there are exactly  $\binom{t}{2}$  distinct pairs of coefficients, and  $s \geq t(t-1) = 2\binom{t}{2}$ ,  $f(\zeta^k x)$  is diverse for at least half of the choices for k.  $\square$ 

We note that the diversification which maps the polynomial f(x) to  $f(\zeta^k x)$  and back is numerically stable since  $\zeta$  is on the unit circle.

In practice, the previous theorem will be far too pessimistic. We therefore propose the method of Algorithm 4 to adaptively choose s,  $\delta$ , and  $\zeta^k$  simultaneously, given a good prime p.

Suppose there exists a threshold  $S \in \mathbb{N}$  such that for all primes s > S, a random sth primitive root of unity  $\zeta^k$  makes  $f(\zeta^k x)$   $\epsilon$ -diverse with high probability. Then Algorithm 4 will return a root of unity whose order is within a constant factor of S, with high probability. From the previous theorem, if such an S exists it must be  $O(t^2)$ , and hence the number of iterations required is  $O(\log t)$ .

Otherwise, if no such S exists, then we cannot diversify the polynomial. Roughly speaking, this corresponds to the situation that f has too many coefficients with absolute value

close to the machine precision. In this case, we can simply use the algorithm of Garg and Schost (2009) numerically, achieving the same stability but using a greater number of evaluations and bit operations. It is possible to establish an adaptive hybrid between our algorithm and that of Garg and Schost (2009) by making f as  $\epsilon$ -diverse as possible given our precision. The non-zero coefficients of f are clustered into groups which are not  $\epsilon$ -diverse (i.e., are within  $\epsilon ||f||_2$  of each other). We can use the symmetric polynomial reconstruction of Garg and Schost (2009) to extract the exponents within each group.

# 4.3 Approximate interpolation algorithm

We now plug our  $\epsilon$ -approximate remainder black box, and method for making f  $\epsilon$ -diverse, into our generic Algorithm 1 to complete our algorithm for approximate interpolation.

THEOREM 4.7. Let  $\delta > 0$ ,  $f \in \mathbb{C}[x]$  with degree at most D and sparsity at most T, and suppose all nonzero coefficients c of f satisfy  $|c| > \delta ||f||_2$ . Suppose also that  $\epsilon < 1.5\delta/(T(T-1))$ , and we are given an  $\epsilon$ -approximate black box for f. Then, for any  $\mu < 1/2$  we have an algorithm to produce a  $g \in \mathbb{C}[x]$  satisfying the conditions of Theorem 4.3. The algorithm succeeds with probability at least  $1-\mu$  and uses  $O^*(T^2 \cdot \log(1/\mu) \cdot \log^2 D)$  black box evaluations and floating point operations.

PROOF. Construct an approximate remainder black box for f using Algorithm 3. Then run Algorithm 1 using this black box as input. On step 8 of Algorithm 1, run Algorithm 4, iterating steps 5–7  $\lceil \log_2(3/\mu) \rceil$  times on each iteration through the while loop to choose a diversifying  $\alpha = \zeta^k$  with probability at least  $1 - \mu/3$ .

The cost comes from Theorems 2.2 and 4.4 along with the previous discussion and Theorem 4.6.  $\qed$ 

Observe that the resulting algorithm is Monte Carlo, but could be made Las Vegas by combining the finite fields zero testing algorithm discussed in Section 3.2 with the guarantees of Theorem 4.3.

#### 5. IMPLEMENTATION RESULTS

We implemented our algorithms in C++ using GMP (gmplib.org) and NTL (www.shoup.net/ntl) for the exponent arithmetic.

In our timing results, "Determ" refers to the deterministic algorithm as stated in Garg and Schost (2009) and "Alg 1" is the algorithm we have presented here over finite fields, without the verification step. We also developed and implemented a more adaptive, Monte Carlo version of these algorithms, as briefly described at the end of Section 2. The basic idea is to sample modulo  $x^p - 1$  for just one prime  $p \in \Theta(t^2 \log d)$  that is good with high probability, then to search for much smaller good primes. This good prime search starts at a lower bound of order  $\Theta(t^2)$  based on the birthday problem, and finds consecutively larger primes until enough primes have been found to recover the symmetric polynomial in the exponents (for Garg & Schost) or just the exponents (for our method). The corresponding improved algorithms are referred to as "G&S MC" and "Alg 1++" below, respectively.

Table 2 summarizes some timings for these four algorithms over the finite field  $\mathbb{Z}/65521\mathbb{Z}$ . This modulus was chosen for convenience of implementation, although of course other

| $\log_2 D$ | T  | Determ | G&S MC | Alg 1 | Alg 1++ |
|------------|----|--------|--------|-------|---------|
| 12         | 10 | 3.77   | 0.03   | 0.03  | 0.01    |
| 16         | 10 | 46.82  | 0.11   | 0.11  | 0.08    |
| 20         | 10 | _      | 0.38   | 0.52  | 0.33    |
| 24         | 10 | _      | 0.68   | 0.85  | 0.38    |
| 28         | 10 | _      | 1.12   | 2.35  | 0.53    |
| 32         | 10 | _      | 1.58   | 2.11  | 0.66    |
| 12         | 20 | 37.32  | 0.15   | 0.02  | 0.02    |
| 16         | 20 | _      | 0.91   | 0.52  | 0.28    |
| 20         | 20 | _      | 3.5    | 3.37  | 1.94    |
| 24         | 20 | _      | 6.59   | 5.94  | 2.99    |
| 28         | 20 | _      | 10.91  | 10.22 | 3.71    |
| 32         | 20 | _      | 14.83  | 16.22 | 4.24    |
| 12         | 30 | _      | 0.31   | 0.01  | 0.01    |
| 16         | 30 | _      | 3.66   | 1.06  | 0.65    |
| 20         | 30 | _      | 10.95  | 6.7   | 3.56    |
| 24         | 30 | _      | 25.04  | 12.42 | 9.32    |
| 28         | 30 | _      | 38.86  | 19.36 | 13.8    |
| 32         | 30 | _      | 62.53  | 68.1  | 14.66   |
| 12         | 40 | _      | 0.58   | 0.01  | 0.02    |
| 16         | 40 | _      | 8.98   | 3.7   | 1.54    |
| 20         | 40 | _      | 30.1   | 12.9  | 8.42    |
| 24         | 40 | _      | 67.97  | 38.34 | 16.57   |
| 28         | 40 | _      | _      | 73.69 | 36.24   |
| 32         | 40 | _      | _      | _     | 40.79   |

Table 2: Finite Fields Algorithm Timings

| Noise          | Mean Error             | Median Error           | Max Error              |
|----------------|------------------------|------------------------|------------------------|
| 0              | $4.440\mathrm{e}{-16}$ | $4.402\mathrm{e}{-16}$ | $8.003\mathrm{e}{-16}$ |
| $\pm 10^{-12}$ | $1.113\mathrm{e}{-14}$ | $1.119\mathrm{e}{-14}$ | $1.179\mathrm{e}{-14}$ |
| $\pm 10^{-9}$  | $1.149\mathrm{e}{-11}$ | $1.191\mathrm{e}{-11}$ | $1.248\mathrm{e}{-11}$ |
| $\pm 10^{-6}$  | $1.145\mathrm{e}{-8}$  | $1.149\mathrm{e}{-8}$  | $1.281\mathrm{e}{-8}$  |

Table 3: Approximate Algorithm Stability

algorithms might be more efficient over this particularly small finite field. The timings are given in seconds of CPU time on a 64-bit AMD Phenom II 3.2GHz processor with  $512\mathrm{K}/2\mathrm{M}/6\mathrm{M}$  cache. Note that the numbers listed reflect the base-2 logarithm of the degree bound and the sparsity bound for the randomly-generated test cases.

The timings are mostly as expected based on our complexity estimates, and also confirm our suspicion that primes of size  $O(t^2)$  are sufficient to avoid exponent collisions. It is satisfying but not particularly surprising to see that our "Alg 1++" is the fastest on all inputs, as all the algorithms have a similar basic structure. Had we compared to the Ben-Or and Tiwari or Zippel's method, they would probably be more efficient for small sizes, but would be easily beaten for large degree and arbitrary finite fields as their costs are super-polynomial.

The implementation of the approximate algorithm uses machine double precision (IEEE), the C++ standard template library complex<double> type, and the popular FFTW package (www.fftw.org) for FFTs. Our stability results are summarized in Table 3. Each test case was randomly generated with degree at most 2<sup>20</sup> and at most 50 nonzero terms. We varied the precision as specified in the table and ran 10

tests in each range. Observe that the error in our results was often less than the  $\epsilon$  error on the evaluations themselves.

Both implementations are freely available for download at http://www.cs.uwaterloo.ca/~droche/diverse/.

## 6. CONCLUSIONS

We have shown how to use the idea of diversification to improve the complexity of sparse interpolation over large finite fields by a factor of t, the number of nonzero terms. We achieve a similar complexity for approximate sparse interpolation, and provide the first provably numerically stable algorithm for this purpose. Our experiments confirm these theoretical results.

Numerous open problems remain. A primary shortcoming of our algorithms is the quadratic dependence on t, as opposed to linear in the case of dense interpolation or even sparse interpolation in smaller or chosen finite fields using the Ben-Or and Tiwari algorithm. It seems that reducing this quadratic dependency will not be possible without a different approach, because of the birthday problem embedded in the diversification step. In the approximate case, a provably numerically stable algorithm for sparse interpolation with only O(t) probes is still an open question. And, while general backward error stability is not possible in the high degree case, it would be interesting in the case of low degree and many variables.

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