# VC-DENSITY IN AN ADDITIVE REDUCT OF p-ADIC NUMBERS

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ABSTRACT. Aschenbrenner et. al. computed a bound  $\operatorname{vc}(n) \leq 2n-1$  for the VC density function in the field of p-adic numbers, but it is not known to be optimal. I investigate a certain P-minimal additive reduct of the field of p-adic numbers and use a cell decomposition result of Leenknegt to compute an optimal bound  $\operatorname{vc}(n) = n$  for that structure.

VC density was introduced into model theory in [1] by Aschenbrenner, Dolich, Haskell, MacPherson, and Starchenko as a natural notion of dimension for definable families of sets in NIP theories. In a NIP theory T we can define the vc-function

$$vc_T = vc : \mathbb{N} \longrightarrow \mathbb{N}$$

where vc(n) measures the worst-case complexity of families of definable sets in an n-dimensional space. The simplest possible behavior is vc(n) = n for all n. For  $T = Th(\mathbb{Q}_p)$ , the paper [1] computes an upper bound for this function to be 2n-1, and it is not known whether it is optimal. This same bound would hold in any reduct of the field of p-adic numbers, so one may expect that the simplified structure of the reduct would allow a better bound. In [2], Leenknegt provides a cell decomposition result for a certain P-minimal additive reduct of the field p-adic numbers. Using this result, in this paper we improve the bound for the VC function, showing that in Leenknegt's structure vc(n) = n.

Explain organization of this paper, notation

#### 1. VC-DIMENSION AND VC-DENSITY

**Definition 1.1.** Throughout this section we work with a collection  $\mathcal{F}$  of subsets of a set X. We call the pair  $(X, \mathcal{F})$  a set system.

- Given a subset A of X, we define the set system  $(A, A \cap \mathcal{F})$  where  $A \cap \mathcal{F} = \{A \cap F\}_{F \in \mathcal{F}}$ .
- For  $A \subset X$  we say that  $\mathcal{F}$  shatters A if  $A \cap \mathcal{F} = \mathcal{P}(A)$ .

**Definition 1.2.** We say  $(X, \mathcal{F})$  has VC-dimension n if the largest subset of X shattered by  $\mathcal{F}$  is of size n. If  $\mathcal{F}$  shatters arbitrarily large subsets of X, we say that  $(x, \mathcal{F})$  has infinite VC-dimension. We denote the VC-dimension of  $(X, \mathcal{F})$  by  $VC(\mathcal{F})$ .

**Note 1.3.** We may drop X from the previous definition, as it VC-dimension doesn't depend on the base set and is determined by  $(\bigcup \mathcal{F}, \mathcal{F})$ .

This allows us to distinguish between well behaved set systems of finite VC-dimension which tend to have good combinatorial properties and poorly behaved set systems with infinite VC dimension.

Another natural combinatorial notion is that of a dual system:

**Definition 1.4.** For  $a \in X$  define  $X_a = \{F \in \mathcal{F} \mid a \in F\}$ . Let  $\mathcal{F}^* = \{X_a\}_{a \in X}$ . We define  $(\mathcal{F}, \mathcal{F}^*)$  as the <u>dual system</u> of  $(X, \mathcal{F})$ . The VC-dimension of the dual system of  $(X, \mathcal{F})$  is referred to as the <u>dual VC-dimension</u> of  $(X, \mathcal{F})$  and denoted by  $VC^*(\mathcal{F})$ . (As before, this notion doesn't depend on X.)

**Lemma 1.5.** A set system has finite VC-dimension if and only if its dual system has finite VC-dimension. More precisely

$$VC^*(\mathcal{F}) \le 2^{1+VC(\mathcal{F})}.$$

For a more refined notion we look at the traces of our family on finite sets:

**Definition 1.6.** Define the shatter function  $\pi_{\mathcal{F}} \colon \mathbb{N} \longrightarrow \mathbb{N}$  and the <u>dual shatter function</u>  $\pi_{\mathcal{F}}^* \colon \mathbb{N} \longrightarrow \mathbb{N}$  of  $\mathcal{F}$  by

$$\pi_{\mathcal{F}}(n) = \max\{|A \cap \mathcal{F}| \mid A \subset X \text{ and } |A| = n\}$$

 $\pi_{\mathcal{F}}^*(n) = \max \{ \text{number of atoms in Boolean algebra generated by } B \mid B \subset \mathcal{F}, |B| = n \}$ 

Note that the dual shatter function is precisely the shatter function of the dual system:  $\pi_{\mathcal{F}}^* = \pi_{\mathcal{F}^*}$ 

A simple upper bound is  $\pi_{\mathcal{F}}(n) \leq 2^n$  (same for the dual). If VC-dimension is infinite then clearly  $\pi_{\mathcal{F}}(n) = 2^n$  for all n. Conversely we have the following remarkable fact:

**Theorem 1.7** (Sauer-Shelah '72). If the set system  $(X, \mathcal{F})$  has finite VC-dimension d then  $\pi_{\mathcal{F}}(n) \leq \binom{n}{\leq d}$  where  $\binom{n}{\leq d} = \binom{n}{d} + \binom{n}{d-1} + \ldots + \binom{n}{1}$ .

Thus the systems with a finite VC-dimension are precisely the systems where the shatter function grows polynomially. Define VC-density to be the degree of that polynomial:

**Definition 1.8.** Define vc-density and dual vc-density of  $\mathcal{F}$  as

$$\operatorname{vc}(\mathcal{F}) = \limsup_{n \to \infty} \frac{\log \pi_{\mathcal{F}}(n)}{\log n} \in \mathbb{R}^{\geq 0} \cup \{+\infty\}$$

$$\mathrm{vc}^*(\mathcal{F}) = \limsup_{n \to \infty} \frac{\log \pi_{\mathcal{F}}^*(n)}{\log n} \in \mathbb{R}^{\geq 0} \cup \{+\infty\}$$

Generally speaking a shatter function that is bounded by a polynomial doesn't itself have to be a polynomial. Proposition 4.12 in [1] gives an example of a shatter function that grows like  $n \log n$  (so it has VC-density 1).

So far the notions that we have defined are purely combinatorial. We now adapt VC-dimension and VC-density to the model theoretic context.

**Definition 1.9.** Work in a structure M. Fix a finite collection of formulas  $\Phi(x,y) = \{\phi_i(x,y)\}.$ 

- For  $\phi(x,y) \in \mathcal{L}(M)$  and  $b \in M^{|y|}$  let  $\phi(M^{|x|},b) = \{a \in M^{|x|} \mid \phi(a,b)\} \subseteq M^{|x|}$ .
- Let  $\Phi(M^{|x|}, M^{|y|}) = \{\phi_i(M^{|x|}, b) \mid \phi_i \in \Phi, b \in M^{|y|}\} \subseteq \mathcal{P}(M^{|x|}).$
- Let  $\mathcal{F}_{\Phi} = \Phi(M^{|x|}, M^{|y|})$  giving a set system  $(M^{|x|}, \mathcal{F}_{\Phi})$ .
- Define <u>VC-dimension</u> of  $\Phi$ , VC( $\Phi$ ) to be the dual VC-dimension of  $(M^{|x|}, \mathcal{F}_{\Phi})$ .
- Define VC-density of  $\Phi$ ,  $vc(\Phi)$  to be the dual VC-density of  $(M^{|x|}, \mathcal{F}_{\Phi})$ .

We will also refer to the VC-density and VC-dimension of a single formula  $\phi$  viewing it as a one element collection  $\{\phi\}$ .

Counting atoms of a Boolean algebra in a model theoretic setting corresponds to counting types, so it is instructive to rewrite the shatter function in terms of types.

### Definition 1.10.

$$\pi_{\Phi}(n) = \max \{ \text{number of } \Phi \text{-types over } B \mid B \subset M, |B| = n \}$$

## Lemma 1.11.

$$\operatorname{vc}(\Phi) = degree \ of \ polynomial \ growth \ of \ \pi_{\Phi}(n) = \limsup_{n \to \infty} \frac{\log \pi_{\Phi}(n)}{\log n}$$

One can check that the shatter function and hence VC-dimension and VC-density of a formula are elementary notions, so they only depend on the first-order theory of the structure.

NIP theories are a natural context for studying VC-density. In fact we can take the following as the definition of NIP:

## **Definition 1.12.** Define $\phi$ to be NIP if it has finite VC-dimension.

[?] shows that in a general combinatorial context, VC-density can be any real number in  $0 \cup [1, \infty)$ . Less is known if we restrict our attention to NIP theories. Proposition 4.6 in [1] gives examples of formulas that have non-integer rational VC-density in an NIP theory, however it is open whether one can get an irrational VC-density in this context.

In general, instead of working with a theory formula by formula, we can look for a uniform bound for all formulas:

**Definition 1.13.** For a given NIP structure M, define the <u>vc-function</u>

$$\operatorname{vc}^{M}(n) = \sup \{ \operatorname{vc}(\phi(x, y)) \mid \phi \in \mathcal{L}(M), |x| = n \}$$

As before this definition is elementary, so it only depends on the theory of M. We omit the superscript M if it is understood from the context. One can easily check the following bounds:

**Lemma 1.14** (Lemma 3.22 in [1]).

$$vc(1) \ge 1$$

$$vc(n) \ge n vc(1)$$

However, it is not known whether the second inequality can be strict or even whether  $vc(1) < \infty$  implies  $vc(n) < \infty$ .

#### 2. P-ADIC NUMBERS

The field of p-adic numbers is often studied in the language of Macintyre  $\mathcal{L}_{Mac} = \{0, 1, +, -, \cdot, |, P_n\}$ . which is a language of fields together with unary predicates  $\{P_n\}_{n\in\mathbb{N}}$  interpreted in  $\mathbb{Q}_p$  by

$$P_n x \leftrightarrow \exists y \ y^n = x$$

and a divisibility relation where a|b holds when val  $a \leq \text{val } b$ .

Note that  $P_n \setminus \{0\}$  is a multiplicative subgroup of  $\mathbb{Q}_p$  with finitely many cosets.

**Theorem 2.1** (Macintyre '76). The  $\mathcal{L}_{Mac}$ -structure  $\mathbb{Q}_p$  has quantifier elimination.

There is also a cell decomposition result.

**Definition 2.2.** Define <u>n-cell</u> recursively. 0-cells are points in  $\mathbb{Q}_p$ . An n+1-cell is a subset of  $\mathbb{Q}_p^{n+1}$  of the following form:

$$\{(x,t) \in \mathbb{Q}_p \times D \mid \operatorname{val} a_1(x) \square_1 \operatorname{val}(t-c(x)) \square_2 \operatorname{val} a_2(x), t-c(x) \in \lambda P_n \}$$

where D is an n-cell,  $a_1(x), a_2(x), c(x)$  are  $\emptyset$ -definable,  $\square$  is  $<, \le$  or no condition, and  $\lambda \in \mathbb{Q}_p$ .

**Theorem 2.3** (Denef '84). Any subset of  $\mathbb{Q}_p$  defined by a  $\mathcal{L}_{Mac}$ -formula  $\phi(x,t)$  with |t|=1 and |x|=n decomposes into a finite union of n+1-cells.

In [1], Aschenbrenner, Dolich, Haskell, Macpherson, and Starchenko show that this structure has  $vc(n) \leq 2n - 1$ , however it is not known whether this bound is optimal.

In [2], Leenknegt analyzes the reduct of p-adic numbers to the language

$$\mathcal{L}_{aff} = \left\{0, 1, +, -, \{\bar{c}\}_{c \in \mathbb{Q}_p}, |, \{Q_{m,n}\}_{m,n \in \mathbb{N}}\right\}$$

where  $\bar{c}$  is a scalar multiplication by c, a|b stands for val  $a \leq \text{val } b$ , and  $Q_{m,n}$  is a unary predicate

$$Q_{m,n} = \bigcup_{k \in \mathbb{Z}} p^{km} (1 + p^n \mathbb{Z}_p).$$

Note that  $Q_{m,n}$  is a subgroup of the multiplicative group of  $\mathbb{Q}_p$  with finitely many cosets. One can check that the extra relation symbols are definable in the  $\mathcal{L}_{Mac}$ -structure  $\mathbb{Q}_p$ . The paper [2] provides a cell decomposition result with the following cells:

**Definition 2.4.** A 0-cell is a point in  $\mathbb{Q}_p$ . An n+1-cell is a subset of  $\mathbb{Q}_p^{n+1}$  of the following form:

$$\{(x,t) \in K \times D \mid \operatorname{val} a_1(x) \square_1 \operatorname{val} (t-c(x)) \square_2 \operatorname{val} a_2(x), t-c(x) \in \lambda Q_{m,n} \}$$

where D is an n-cell called the <u>base</u> of the cell,  $a_1(x), a_2(x), c(x)$  are linear polynomials,  $\square$  is < or no condition, and  $\lambda \in \mathbb{Q}_p$ .

**Theorem 2.5** (Leenknegt '12). Any formula  $\phi(x,t)$  in  $(\mathbb{Q}_p, \mathcal{L}_{aff})$  with |t| = 1 and |x| = n decomposes into a union of n + 1-cells.

Moreover, [2] shows that  $(\mathbb{Q}_p, \mathcal{L}_{aff})$  is a P-minimal reduct, that is the one-dimensional definable sets of  $(\mathbb{Q}_p, \mathcal{L}_{aff})$  coincide with the one-dimensional definable sets in the full structure  $(\mathbb{Q}_p, \mathcal{L}_{Mac})$ .

I am able to compute the vc-function for this structure

**Theorem 2.6.** Theorem (B.)  $(\mathbb{Q}_p, \mathcal{L}_{aff})$  has vc(n) = n.

### 3. Key Lemmas and Definitions

To show that  $\operatorname{vc}(n) = n$  it suffices to bound  $\operatorname{vc}(\phi) \leq |x|$  for every formula  $\phi(x;y)$ . Fix such a formula  $\phi(x;y)$ . Instead of working with it directly, we simplify it using quantifier elimination. Quantifier elimination result can be easily obtained from cell decomposition:

**Lemma 3.1.** Any formula  $\phi(x;y)$  in  $(\mathbb{Q}_p, \mathcal{L}_{aff})$  can be written as a boolean combination of formulas from the following collection

$$\Psi(x;y) = \{ \text{val}(p_i(x) - c_i(y)) < \text{val}(p_j(x) - c_j(y)) \}_{i,j \in I} \cup \{ p_i(x) - c_i(y) \in \lambda_k Q_{m,n} \}_{i \in I, k \in K}$$

where I, K are finite index sets, each  $p_i$  is a linear polynomial in x without a constant term, each  $c_i$  is a linear polynomial in y, and  $\lambda_k \in \mathbb{Q}_p$ .

Proof. Let l = |x| + |y|. Apply cell decomposition theorem to  $\phi(x; y)$  to obtain  $\mathscr{D}^l$ , a collection of l-cells. Let  $\mathscr{D}^{l-1}$  be a collection l-1 of bases of cells in  $\mathscr{D}^l$ . Similarly, construct by induction  $\mathscr{D}^i$  for each  $0 \leq j < l$ , where  $\mathscr{D}_j$  is a collection of j-cells which are the bases of cells in  $\mathscr{D}_{j+1}$ . Let  $\mathscr{D} = \bigcup \mathscr{D}_j$ . Choose n, m large enough to cover all n', m' that come up in the cells for  $Q_{n',m'}$ . Choose  $\lambda_k$  to go over all the cosets of  $Q_{n,m}$ . Let  $q_i(x,y)$  enumerate all of the polynomials  $a_1(\bar{x}), a_2(\bar{x}), t-c(\bar{x})$  that show up in the cells of  $\mathscr{D}$ . Those are all polynomials of degree  $\leq 1$  in variables

x,y. We can split each of them as  $q_i(x,y) = p_i(x) - c_j(y)$  where the constant term goes into  $c_j$ . This gives us the appropriate finite collection of formulas  $\Psi$ . From cell decomposition it is easy to see that when a,a' have the same  $\Psi$ -type, then they would have they have the same  $\phi$ -type. Thus  $\phi$  can be written as a boolean combination of formulas from  $\Psi$ .

**Lemma 3.2.** If  $\phi$  can be written as a Boolean combination of formulas from  $\Psi$  then

$$\operatorname{vc}(\Psi) \le n \implies \operatorname{vc}(\phi) \le n$$

*Proof.* If a, a' have the same  $\Psi$ -type over B, then they have the same  $\phi$ -type over B, where B is some parameter set. Therefore the number of  $\phi$ -types is bounded by the number of  $\Psi$ -types. The bound follows from Lemma 1.11.

Therefore to show that  $\operatorname{vc}(\phi) \leq |x|$ , it suffices to bound  $\operatorname{vc}(\Psi) \leq |x|$ . More precisely, it is sufficient to show that if there is a parameter set B of size N then the number of  $\Psi$ -types over B is  $O(N^{|x|})$ . Fix such a parameter set B and work with it from now on. We will compute a bound for the number of  $\Psi$ -types over B.

Consider a set  $T = \{c_i(b) \mid b \in B, i \in I\} \subset \mathbb{Q}_p$ . In this definition B is the parameter set that we fixed and  $c_i(b)$  come from collection of formulas  $\Psi$  from the quantifier elimination above. View T as a tree as follows:

**Definition 3.3.** • For  $c \in \mathbb{Q}_p$ ,  $\alpha \in \mathbb{Z}$  define a <u>ball</u>

$$B(c,\alpha) = \{c' \in \mathbb{Q}_p \mid \operatorname{val}(c' - c) \le \alpha\}.$$

• Define a collection of balls  $\mathscr{B} = \{B(t_1, \operatorname{val}(t_1 - t_2))\}_{t_1, t_2 \in T}$ . An <u>interval</u>  $(B_1, B_2)$  is a set  $B_1 \backslash B_2$  for  $B_1, B_2 \in \mathscr{B}$  with  $B_1 \supset B_2$  and no balls from  $\mathscr{B}$  in between. We also define an interval  $(-\infty, B)$  as a set  $\mathbb{Q}_p \backslash B$  for a ball  $B(c, v) \in \mathscr{B}$  with the smallest valuation v of all the balls in  $\mathscr{B}$ . Note that there are at most 2|T| = 2N|I| = O(N) different intervals and they partition  $\mathbb{Q}_p$ .

• Define a collection of balls  $\mathscr{B}' = \mathscr{B} \cup \{B(c_{i_1}(b), \operatorname{val}(c_{i_2}(b) - c_{i_3}(b)))\}_{i_1, i_2, i_3 \in I, b \in B}$ . An <u>sub-interval</u> is defined the same as an interval except using collection  $\mathscr{B}'$  instead of  $\mathscr{B}$ . Sub-intervals refine intervals, and there are at most  $2|T| + |B| \cdot |I|^3 = O(N)$  many of them.

**Definition 3.4.** Suppose  $a \in \mathbb{Q}_p$  lies in an interval  $B(t_L, \alpha_L) \setminus B(t_U, \alpha_U)$ .

- Define <u>T-valuation</u> of a to be  $T-val(a) = val(a t_U)$ .
- Define floor of a to be  $F(a) = \alpha_L$ .

**Definition 3.5.** Suppose  $a_1, a_2 \in \mathbb{Q}_p$  lie in our tree in the same interval  $B(t_L, \alpha_L) \setminus B(t_U, \alpha_U)$ . We say that  $a_i$  is close to boundary if  $|\operatorname{T-val}(a_i) - \alpha_L| \le m$  or  $|\operatorname{T-val}(a_i) - \alpha_U| \le m$ . Otherwise we say that it is far from boundary.

**Definition 3.6.** We say  $a_1, a_2$  have the same <u>interval type</u> if one of the following holds:

- Both  $a_1, a_2$  are far from boundary and  $a_1 t_U, a_2 t_U$  are in the same  $Q_{m,n}$  coset.
- Both  $a_1, a_2$  are close to boundary and  $val(a_1 a_2) > T-val(a_1) + n = T-val(a_2) + n$ .

The following lemma is an adaptation of lemma 7.4 in [1].

**Lemma 3.7.** For n, m there exists  $D = D(n, m) \in \mathbb{Z}$  such that for any  $x, y, a \in \mathbb{Q}_p$  if

$$val(x-c) = val(y-c) < val(x-y) - D$$

then x - c, y - c are in the same coset of  $Q_{n,m}$ .

*Proof.* Define that  $a, b \in \mathbb{Q}_p$  are similar if val a = val b and

$$a \upharpoonright [\operatorname{val} a, \operatorname{val} a + (m+n)] = b \upharpoonright [\operatorname{val} b, \operatorname{val} b + (m+n)]$$

If a, b are similar then

$$a \in Q_{n,m} \leftrightarrow b \in Q_{n,m}$$

Moreover for any  $\lambda \in \mathbb{Q}_p$ , if a, b are similar we would also have  $a/\lambda, b/\lambda$  are similar. Thus if a, b are similar, then they belong in the same coset of  $Q_{n,m}$ . If we pick D = n + m then conditions of the lemma force x - c, y - c to be similar.

The following construction is along the lines of lemmas 7.3, 7.5 of [1].

**Definition 3.8.** For two balls  $B(a, \alpha)$ ,  $B(b, \beta)$  let  $\gamma = \min(\alpha, \beta, \operatorname{val}(a - b))$ . Define the distance between those two balls to be  $|\alpha - \gamma| + |\beta - \gamma|$ . In  $\mathbb{Q}_p$  value group is discrete and residue field is finite, so there are finitely many balls at a fixed distance from a given ball. Near balls of  $B(a, \alpha)$  are defined to be balls with distance  $\mathcal{D}$  from  $B(a, \alpha)$ . Enumerate those as:

$$B_1(a,\alpha), B_2(c,\alpha), \dots B_{N_D}(a,\alpha)$$

Near balls partition the space

$$\{b \in \mathbb{Q}_p \mid |\operatorname{val}(a-b) - \alpha| \le D\}$$

**Lemma 3.9.** Suppose  $c_1, c_2 \in \mathbb{Q}_p^{|x|}$  satisfy the following three conditions

- For all  $i \in I$   $p_i(c_1)$  and  $p_i(c_2)$  are in the same sub-interval.
- For all  $i \in I$   $p_i(c_1)$  and  $p_i(c_2)$  have the same interval type.
- For all  $i, j \in I$ ,  $\operatorname{T-val}(p_i(c_1)) > \operatorname{T-val}(p_j(c_1))$  iff  $\operatorname{T-val}(p_i(c_2)) > \operatorname{T-val}(p_j(c_2))$ .

Then  $c_1, c_2$  have the same  $\Psi$ -type over B.

Proof. There are two kinds of formulas in  $\Psi$  (see Lemma 3.1). First we show that  $d_1, d_2$  agree on formulas of the form  $p_i(x) - c_i(y) \in \lambda_k Q_{m,n}$ . It is enough to show that for every  $i \in I, b \in B$  we have  $p_i(d_1) - c_i(b), p_i(d_2) - c_i(b)$  are in the same  $Q_{m,n}$ -coset. Fix such i, b. For brievety let  $a = p_i(d_1), a' = p_i(d_2)$  and  $Q = Q_{m,n}$ . We want to show that  $a - c_i(b), a' - c_i(b)$  are in the same Q-coset.

Suppose a is in one of the near balls. As a' has the same interval type, it has to be in the same near ball. By definition of the near ball we then have  $val(a - c_i(b)) = val(a' - c_i(b)) < val(a - a') - D$ . Thus by Lemma 3.7 we have  $a - c_i(b), a' - c_i(b)$  in the same Q-coset.

Now, suppose both a, a' aren't in any near balls. Label their interval as  $B(c_L, \alpha_L) \setminus B(c_U, \alpha_U)$ . Then we have

$$\alpha_L + D < \text{val}(a - c_U) < \alpha_U - D$$

$$\alpha_L + D < \text{val}(a' - c_U) < \alpha_U - D$$

as otherwise one (both) of them would be in one of the near balls. We have either  $\operatorname{val}(c_U - c_i(b)) \geq \alpha_U$  or  $\operatorname{val}(c_U - c_i(b)) \leq \alpha_L$  as otherwise it would contradict the definition of an interval.

Suppose it is the first case  $val(c_U - c_i(b)) \ge \alpha_U$ . Then

$$val(a - c_i(b)) = val(a - c_U) < \alpha_U - D \le val(c_U - c_i(b)) - D$$

so by Lemma 3.7 we have  $a-c_i(b)$ ,  $a-c_U$  are in the same Q-coset. By a parallel argument we have  $a'-c_i(b)$ ,  $a'-c_U$  are in the same Q-coset. As we are assuming a, a' have the same tree type it implies that  $a-c_U, a'-c_U$  are in the same Q-coset. Thus by transitivity we get that  $a-c_i(b)$ ,  $a'-c_i(b)$  are in the same Q-coset.

For the second case, suppose  $val(c_U - c_i(b)) \leq \alpha_L$ . Then

$$\operatorname{val}(a - c_i(b)) = \operatorname{val}(c_U - c_i(b)) \le \alpha_L < \operatorname{val}(a - c_U) - D$$

so by Lemma 3.7 we have  $a - c_i(b)$ ,  $c_U - c_i(b)$  are in the same Q-coset. By a parallel argument we have  $a' - c_i(b)$ ,  $c_U - c_i(b)$  are in the same Q-coset. Thus by transitivity we get that  $a - c_i(b)$ ,  $a' - c_i(b)$  are in the same Q-coset.

Next, we need to show that  $d_1, d_2$  agree on formulas of the form  $\operatorname{val}(p_i(x) - c_i(y)) < \operatorname{val}(p_j(x) - c_j(y))$  (see Lemma 3.1).

This gives us an upper bound on the number of types - there are at most |I|! many choices for the order of T-val, O(N) many choices for the interval for each  $p_i$ , and K many choices for the interval type for each  $p_i$ , giving a total of  $O(N^{|I|}) \cdot K^{|I|} \cdot |I|! = O(N^{|I|})$  many types. This implies  $\operatorname{vc}(\Psi) \leq |I|$ . The biggest contribution to this bound are the choices among the O(N) many intervals for each  $p_i$  with  $i \in I$ . Are all of those choices realized? Intuitively there are |x| many variables and |I| many equations, so once we choose an interval for |x| many  $p_i$ 's, the interval for the rest should be determined. This would give the required  $\operatorname{vc}(\Psi) \leq |x|$  bound. The next section outlines this proof formally.

### 4. Main Proof

**Lemma 4.1.** Suppose we have a finite collection of vectors  $\{\vec{p}_i\}_{i\in I}$  with each  $\vec{p}_i \in \mathbb{Q}_p^{|x|}$ . Suppose  $J \subset I$  and  $i \in I$  satisfy

$$\vec{p}_i \in \operatorname{span} \left\{ \vec{p}_j \right\}_{j \in J}$$
,

and we have  $\vec{x} \in \mathbb{Q}_p^{|x|}, \alpha \in \mathbb{Z}$  with

$$\operatorname{val}(\vec{p_j} \cdot \vec{x}) > \alpha \text{ for all } j \in J$$

Then

$$\operatorname{val}(\vec{p_i} \cdot \vec{x}) > \alpha - \gamma$$

for some  $\gamma \in \mathbb{N}$ . Moreover  $\gamma$  can be chosen independently from  $J, j, \vec{x}, \alpha$  depending only on  $\{\vec{p}_i\}_{i \in I}$ .

*Proof.* Fix i, J satisfying the conditions of the lemma. For some  $c_j \in \mathbb{Q}_p$  for  $j \in J$  we have

$$\vec{p_i} = \sum_{j \in J} c_j \vec{p_j},$$

hence

$$\vec{p}_i \cdot \vec{x} = \sum_{j \in J} c_j \vec{p}_j \cdot \vec{x}.$$

We have

$$\operatorname{val}(c_j \vec{p}_j \cdot \vec{x}) = \operatorname{val}(c_j) + \operatorname{val}(\vec{p}_j \cdot \vec{x}) > \operatorname{val}(c_j) + \alpha.$$

Let  $\gamma = \max(0, \min - \operatorname{val}(c_j))$ . Then we have

$$\operatorname{val}(c_{j}\vec{p}_{j}\cdot\vec{x}) > \alpha - \gamma \qquad \text{for all } j \in J$$

$$\sum_{j \in J} c_{j}\vec{p}_{j}\cdot\vec{x} > \alpha - \gamma$$

This shows that we can pick such  $\gamma$  for a given choice of i, J, but independent from  $\alpha, \vec{x}$ . To get a choice independent from i, J, go over all such eligible choices (i ranges over I and J ranges over subsets of I), pick  $\gamma$  for each, and then take the maximum of those values.

Alternative way to write  $p_i(x)$  is  $\vec{p}_i \cdot \vec{x}$ , where  $\vec{p}_i$  and  $\vec{x}$  are vectors in  $\mathbb{Q}_p^{|x|}$ . The lemma above is a general result, but we only use it applied to the vectors  $\vec{p}_i$  given by our collection of formulas  $\Psi$ .

**Definition 4.2.** For  $c \in \mathbb{Q}_p$  and  $\alpha, \beta \in \mathbb{Z}$  define  $c \upharpoonright [\alpha, \beta] \in (\mathbb{Z}/p\mathbb{Z})^{\beta - \alpha + 1}$  to be the record of the coefficients of c for the valuations between  $\alpha, \beta$ . More precisely write c in its power series form

$$c = \sum_{\gamma \in \mathbb{Z}} c_{\gamma} p^{\gamma} \text{ with } c_{\gamma} \in \mathbb{Z}/p\mathbb{Z}$$

Then  $c \upharpoonright [\alpha, \beta]$  is just  $(c_{\alpha}, c_{\alpha+1}, \dots c_{\beta})$ .

Fix  $\gamma$  corresponding to  $\{\vec{p_i}\}_{i\in I}$  according to Lemma 4.1.

**Definition 4.3.** Denote  $\mathbb{Z}/p\mathbb{Z}^{\gamma}$  as Ct.

**Definition 4.4.** Let  $f: \mathbb{Q}_p^{|x|} \longrightarrow \mathbb{Q}_p^I$  with  $f(\bar{c}) = (p_i(\bar{c}))_{i \in I}$ . Define the segment space Sg to be the image of f.

Given a tuple  $(a_i)_{i\in I}$  in the segment space look at the corresponding floors  $\{F(a_i)\}_{i\in I}$ . Those are ordered as elements of  $\mathbb{Z}$ . Partition the segment space by order type of  $\{F(a_i)\}$ . Work in a fixed partition Sg'. After relabeling we may assume that

$$F(a_1) \geq F(a_2) \geq \dots$$

Consider the (relabeled) sequence of vectors  $\vec{p}_1, \vec{p}_2, \dots, \vec{p}_I$ . There is a unique subset  $J \subset I$  such that all vectors with indices in J are linearly independent, and all vectors with indices outside of J are a linear combination of preceding vectors. For any index  $i \in I$  we call it independent if  $i \in J$  and we call it dependent otherwise.

Now, we define the following function

$$g: \operatorname{Sg}' \longrightarrow \operatorname{Bt}^I \times \operatorname{Pt}^J \times \operatorname{Ct}^{I-J}$$

Let  $\bar{a} = (a_i)_{i \in I} \in \operatorname{Sg}'$ . To define  $g(\bar{a})$  we need to specify where it maps  $\bar{a}$  in each individual component of the product.

For all  $a_i$  record its interval type  $\in$  Bt, giving the first component.

For  $a_j$  with  $j \in J$ , record the interval of  $a_j$ , giving the second component.

For the third component do the following computation. Pick  $a_i$  with i dependent. Let j be the largest independent index with j < i. Record  $a_i \upharpoonright [F(a_j) - \gamma, F(a_j)]$ .

**Lemma 4.5.** For  $\bar{a}, \bar{a}' \in \operatorname{Sg}'$  if  $g(\bar{a}) = g(\bar{a}')$  then  $a_i, a_i'$  have the same tree type for all  $i \in I$ .

*Proof.* For each i we show that  $a_i, a_i'$  are in the same interval and have the same interval type, so the conclusion follows by Lemma ??. Bt records the interval type of each element, so if  $g(\bar{a}) = g(\bar{a}')$  then  $a_i, a_i'$  have the same interval type for all  $i \in I$ . Thus it remains to show that  $a_i, a_i'$  lie in the same interval for all  $i \in I$ .

Suppose i is an independent index. Then by construction, Pt records the interval for  $a_i, a_i'$ , so those have to belong to the same interval. Now suppose i is dependent. Pick the largest j < i such that j is independent. We have  $F(a_i) \leq F(a_j)$  and  $F(a_i') \leq F(a_j')$ . Moreover  $F(a_j) = F(a_j')$  as they are mapped to the same interval (using the earlier part of the argument as j is independent).

**Claim 4.6.** val
$$(a_i - a_i') > F(a_j) - \gamma$$

*Proof.* Let  $\vec{x}, \vec{x}' \in \mathbb{Q}_p^{|x|}$  be some elements with

$$\vec{p}_k \cdot \vec{x} = a_k$$
 
$$\vec{p}_k \cdot \vec{x}' = a_k' \text{ for all } k \in I$$

It is always possible to do that as  $\bar{a}, \bar{a}' \in Sg'$ . Let J' be the set of the independent indices less than i. We have

$$\operatorname{val}(a_k - a_k') > F(a_k) \text{ for all } k \in J'$$

as for the independent indices  $a_k, a_k'$  lie in the same interval.

$$\operatorname{val}(a_k - a_k') > F(a_j)$$
 for all  $k \in J'$  by monotonicity of  $F(a_k)$ 

$$\operatorname{val}(\vec{p}_k \cdot \vec{x} - \vec{p}_k \cdot \vec{x}') > F(a_j) \text{ for all } k \in J'$$

$$\operatorname{val}(\vec{p}_k \cdot (\vec{x} - \vec{x}')) > F(a_j) \text{ for all } k \in J'$$

J' and i match the requirements of Lemma 4.1 so we conclude

$$\operatorname{val}(\vec{p}_i \cdot (\vec{x} - \vec{x}')) > F(a_j) - \gamma$$

$$\operatorname{val}(\vec{p}_i \cdot \vec{x} - \vec{p}_i \cdot \vec{x}') > F(a_j) - \gamma$$

$$\operatorname{val}(a_i - a_i')) > F(a_j) - \gamma$$

as needed, finishing the proof of the claim.

Additionally  $a_i, a'_i$  have the same image in Ct component, so we have

$$\operatorname{val}(a_i - a_i') > F(a_i)$$

As  $F(a_i) \leq F(a_j)$ ,  $a_i, a'_i$  have to lie in the same interval.

Corollary 4.7.  $\Psi(x,y)$  has VC-density  $\leq |x|$ 

Proof. Suppose we have  $c, c' \in \mathbb{Q}_p^{|x|}$  such that f(c), f(c') are in the same partition and g(f(c)) = g(f(c')). Then by the previous lemma  $p_i(c)$  has the same tree type as  $p_i(c')$  for all  $i \in I$ . Then by Lemma ?? c, c' have the same  $\Psi$ -type. Thus the number of possible  $\Psi$ -types is bounded by the size of the range of g times the number of possible partitions

(number of partitions) 
$$\cdot |Bt|^{|I|} \cdot |Pt|^{|J|} \cdot |Ct|^{|I-J|}$$

We have

 $|\operatorname{Bt}| = N_D + \text{number of cosets of } Q|\operatorname{Pt}| \leq N \cdot I^2 \text{ (the only component dependent on } N)$ 

$$|\operatorname{Ct}| = p^{\gamma}$$

and there are at most |I|! many partitions of Sg. This gives us a bound

$$|I|!\cdot|Bt|^{|I|}\cdot(N\cdot{|I|}^2)^{|J|}\cdot p^{\gamma|I-J|}=O(N^{|J|})$$

Every  $p_i$  is an element of a |x|-dimensional vector space, so there can be at most |x| many independent vectors. Thus we have  $|J| \leq |x|$  and the bound follows.  $\square$ 

Corollary 4.8. In the language  $\mathcal{L}_{aff}$  we have vc(n) = n.

*Proof.* Previous lemma implies that  $\operatorname{vc}(\phi) \leq \operatorname{vc}(\Psi) \leq |x|$ . As choice of  $\phi$  was arbitrary, this implies that VC-density of any formula is bounded by the arity of x.

## References

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