# Algorithms for matrix groups

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angle \leq \operatorname{GL}(d,R)$  where R is a ring; usually finite field  $\operatorname{GF}(q)$ 

Goal: efficient algorithms, for their study, which are both theoretically and practically effective.

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# Why do we care?

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- Invariant theory: irreducible representations, Kronecker products, tensor-induced representations.
- Energy levels of systems of identical particles: irreducible representations of classical groups

#### Two $d \times d$ matrices A and B Cost of $A \times B$ using conventional algorithm is $O(d^3)$ .

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Where do we notice improvements? Perhaps for  $d \ge 100$ .

Given  $G \leq \operatorname{GL}(d, \mathbb{Z})$ , and  $x \in \operatorname{GL}(d, \mathbb{Z})$ : is  $x \in G$ ?

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Given  $G \leq \operatorname{GL}(d, \mathbb{Z})$ , and  $x \in \operatorname{GL}(d, \mathbb{Z})$ : is  $x \in G$ ? Mihailova (1958): membership problem is undecidable for  $d \geq 4$ .

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Discrete log problem  $F = GF(q), \ \omega \in F$  primitive. Given  $\alpha \in F$ , determine k so that  $\alpha = \omega^k$ .

No polynomial-time algorithm known.

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Babai & Beals (1999):

#### Theorem

If the set of primes dividing a multiplicative upper-bound B for |g| is known, then the precise value of |g| can be determined in polynomial time.

• Compute a "good" multiplicative upper bound E for |g|.

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$$m(x) = \prod_{i=1}^t f_i(x)^{m_i}$$

where deg( $f_i$ ) =  $d_i$  and  $\beta = \lceil \log_p \max m_i \rceil$ .

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 $E = \operatorname{lcm}(q^{d_i} - 1) \times p^{\beta}$ |g| divides E.

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If t = 1, then compute  $g^{p'_1}$  for  $j = 1, 2, \ldots, \alpha_1$ .

Otherwise write E = uv where u, v are coprime and have approximately same number of distinct prime factors.

Now  $g^u$  has order k say, dividing v; and  $g^k$  has order  $\ell$  say, dividing u.

The order of g is  $k\ell$ .

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Implementations in both GAP and Magma use databases of factorisations of numbers of the form  $q^i - 1$ , prepared as part of the Cunningham Project.

If we just know E, then we can learn in polynomial time the *exact* power of 2 (or of any specified prime) which divides |g|.

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Now we compute  $h = g^b$ , and determine (by powering) its order which divides  $2^m$ .

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Many algorithms are **randomised**: use random search in G to find elements having prescribed property  $\mathcal{P}$ .

#### Example

- Characteristic polynomial having factor of degree > d/2.
- Order divisible by prescribed prime.

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Common feature: algorithms depend on detailed analysis of **proportion** of elements of finite simple groups satisfying  $\mathcal{P}$ .

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To find element satisfying  $\mathcal{P}$  by random search with a probability of failure less than given  $\epsilon \in (0, 1)$ : choose a sample of uniformly distributed random elements in G of size at least  $[-\log_e(\epsilon)]k$ .

## Challenge Problem II: Generate random elements

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Walk is repeated  $O(\log |G|)$  times.

Final list S of  $O(\log |G|)$  elements input to construction phase.

$$g_1^{\epsilon_1}\cdots g_m^{\epsilon_m}$$

where  $S = \{g_1, \ldots, g_m\}$  and  $\epsilon_i \in \{0, 1\}$  (chosen independently).

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Cost per random element is  $O(\log |G|)$ .

## CLMNO (1995): Product replacement algorithm

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Input: ordered list of generators  $[g_1, \ldots, g_m]$  for G.

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• Select at random i, j where  $1 \le i, j \le m$ .

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- Multiply r by g<sub>i</sub>.

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Basic step repeated a number, say t, of times.

Now to obtain random element: execute basic operation once, and return r as random element.

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 $Markov\ CHAIN:$  a discrete random process with a finite number of states and it satisfies the property that the next state depends only on the current state.

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#### Theorem

Let T be set of all m-tuples of generators of G. Then the algorithm constructs a Markov chain over state space T, and if m is at least twice the size of a minimal generating set of generators for G, this Markov chain is connected and aperiodic. Cost: after initialisation, two matrix multiplications.

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The random walk approaches a limiting distribution at exponential rate  $O((1 - \delta)^t)$  where t is number of steps taken.

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Diaconis & Saloff-Coste (1997, 1998):
 t = O(δ<sup>2</sup>(G, S) · m), where δ(G, S) is the maximal diameter for the Cayley graph of G wrt generating set S.
 Comparison of two Markov chains on different but related state spaces and combinatorics of random paths.

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- Pak (2001): Mixing time is polynomial. Multi-commodity flow technique.
- Lubotzky & Pak (2002):

Does the group of automorphisms of a free group of rank >3 have Kazhdan's property (T)? If so, then "graph of states" is well-behaved, giving excellent mixing time.

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## Permutation groups

Sims (1970, 1971): base and strong generating set (BSGS). G acts faithfully on  $\Omega = \{1, \ldots, n\}$ 

 $G_{\epsilon} = \{g \in G \mid \epsilon^g = \epsilon\}.$ 

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*Base*: sequence of points  $B = [\epsilon_1, \epsilon_2, \ldots, \epsilon_k]$  where  $G_{\epsilon_1, \epsilon_2, \ldots, \epsilon_k} = 1$ .

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*Base*: sequence of points  $B = [\epsilon_1, \epsilon_2, \dots, \epsilon_k]$  where  $G_{\epsilon_1, \epsilon_2, \dots, \epsilon_k} = 1$ . This determines chain of stabilisers

$$G = G^{(0)} \ge G^{(1)} \ge \cdots \ge G^{(k-1)} \ge G^{(k)} = 1,$$

where  $G^{(i)} = G_{\epsilon_1, \epsilon_2, \dots, \epsilon_i}$ . S strong generating set:  $G^{(i)} = \langle S \cap G^{(i)} \rangle$ 

#### Example

$$G = \langle (1,5,2,6), (1,2)(3,4)(5,6) \rangle$$
  

$$B = [1,3]$$
  

$$G > G_1 > G_{1,3} = 1$$
  

$$S = \{ (1,5,2,6), (1,2)(3,4)(5,6), (3,4) \}$$

$$|G^{(i)}:G^{(i+1)}| = \#B_i$$

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Variations underpin both theoretical and practical approaches to permutation group algorithms.

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Butler (1979): action of G on one-dimensional subspaces of V.

Murray & O'Brien (1995): heuristic algorithm to select base points. Neunhöffer et al. (2000s): use "helper subgroups" to construct large orbits Critical for success: index of one stabiliser in its predecessor.

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"Optimal" subgroup chain for GL(d, q)?

$$\operatorname{GL}(d,q) \ge q^{d-1}.\operatorname{GL}(d-1,q) \ge \operatorname{GL}(d-1,q) \ge \dots$$

**Leading index**:  $q^d - 1$ .

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Leading index:  $q^d - 1$ .

#### Example

Largest maximal subgroup  $2^{11}$ :  $M_{24} \le J_4$  index 173 067 389.

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- *G* preserves some natural linear structure associated with the action of *G* on *V*, and has normal subgroup related to this structure,
- or G is almost simple modulo scalars: T ≤ G/Z ≤ Aut(T) where T is simple.

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- **2** If  $N \lhd G$  exists, recognise N and G/N recursively, ultimately obtaining a composition series for the group.

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Then  $\phi$  :  $G \rightarrow S_r$  where r|d and  $N = \ker \phi$ .

COMPOSITIONTREE: exploits geometry to produce composition series for G, factors are **leaves** of tree.

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