

# Universidad Complutense de Madrid

# #define int long long

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Contest (1)		
template.cpp		14 li

```
#include <bits/stdc++.h>
```

```
using namespace std;
#define rep(i, a, b) for(int i = a; i < (b); ++i)
\#define all(x) begin(x), end(x)
#define sz(x) (int)(x).size()
#define int long long
typedef pair<int, int> pii;
typedef vector<int> vi;
signed main() {
 cin.tie(0)->sync_with_stdio(0);
 cin.exceptions(cin.failbit);
```

#### troubleshoot.txt

Write a few simple test cases if sample is not enough. Are time limits close? If so, generate max cases. Is the memory usage fine? Could anything overflow? Make sure to submit the right file.

Wrong answer:

Print your solution! Print debug output, as well. Are you clearing all data structures between test cases Can your algorithm handle the whole range of input? Read the full problem statement again. Do you handle all corner cases correctly? Have you understood the problem correctly? Any uninitialized variables?

Any overflows? Confusing N and M, i and j, etc.?

Are you sure your algorithm works? What special cases have you not thought of?

Are you sure the STL functions you use work as you

Go through the algorithm for a simple case.

```
Add some assertions, maybe resubmit.
Create some testcases to run your algorithm on.
```

```
Go through this list again.
Explain your algorithm to a teammate.
Ask the teammate to look at your code.
Go for a small walk, e.g. to the toilet.
Is your output format correct? (including whitespace)
Rewrite your solution from the start or let a teammate
    do it.
Runtime error:
Have you tested all corner cases locally?
Any uninitialized variables?
Are you reading or writing outside the range of any
    vector?
Any assertions that might fail?
Any possible division by 0? (mod 0 for example)
Any possible infinite recursion?
Invalidated pointers or iterators?
Are you using too much memory?
Debug with resubmits (e.g. remapped signals, see
    Various).
Time limit exceeded:
Do you have any possible infinite loops?
What is the complexity of your algorithm?
Are you copying a lot of unnecessary data? (References)
How big is the input and output? (consider scanf)
Avoid vector, map. (use arrays/unordered_map)
What do your teammates think about your algorithm?
Memory limit exceeded:
What is the max amount of memory your algorithm should
Are you clearing all data structures between test cases
```

# Strings (2)

#### KMP.h

52 lines

4

**Description:** kmp[i] = The length of the longest non trivial suffix that ends at position i and coincides with a prefix of s.

Time:  $\mathcal{O}(n)$ 

```
vi kmp(const string& s) {
 vi res(sz(s));
 rep(i, 1, sz(s)) {
   int k = res[i - 1];
   while (k > 0 \&\& s[k] != s[i]) k = res[k - 1];
   res[i] = k + (s[k] == s[i]);
 return res; }
```

### Zfunction.h

**Description:** zfun[i] = The length of the longest non trivial prefix that starts at position i and coincides with a prefix of s. zfun[0] = 0. Time:  $\mathcal{O}(n)$ 

```
vi zfun(const string& s) {
```

```
vi z(sz(s)); int l = 0;
rep(i, 1, sz(s)) {
 z[i] = \max(\min(z[i-1], z[1] + 1 - i), OLL);
 while (i+z[i] < sz(s) \&\& s[z[i]] == s[i+z[i]]) z[i] ++;
 if(z[i] + i > z[1] + 1) 1 = i;
return z; }
```

# Number theory (3)

# Modular arithmetic

#### ModInverse.h

**Description:** Pre-computation of modular inverses. Assumes LIM < mod and that mod is a prime.

```
const 11 mod = 1000000007, LIM = 200000;
11* inv = new 11[LIM] - 1; inv[1] = 1;
rep(i,2,LIM) inv[i] = mod - (mod / i) * inv[mod % i] %
```

#### ModPow.h

```
const 11 mod = 1000000007; // faster if const
11 modpow(ll b, ll e) {
 ll ans = 1;
  for (; e; b = b * b % mod, e /= 2)
   if (e & 1) ans = ans \star b % mod;
  return ans;
```

# ModLog.h

**Description:** Returns the smallest x > 0 s.t.  $a^x = b \pmod{m}$ , or -1if no such x exists. modLog(a,1,m) can be used to calculate the order of

#### Time: $\mathcal{O}(\sqrt{m})$

ll modLog(ll a, ll b, ll m) { 11 n = (11) sqrt(m) + 1, e = 1, f = 1, j = 1;unordered map<11, 11> A; while  $(j \le n \& \& (e = f = e * a % m) != b % m)$ A[e \* b % m] = j++;if (e == b % m) return j;  $if (\underline{\hspace{0.5cm}} gcd(m, e) == \underline{\hspace{0.5cm}} gcd(m, b))$ rep(i,2,n+2) if (A.count(e = e \* f % m)) return n \* i - A[e]; return -1:

### ModSum.h

**Description:** Sums of mod'ed arithmetic progressions.

modsum(to, c, k, m) =  $\sum_{i=0}^{to-1} (ki+c)\%m$ . divsum is similar but for floored division.

**Time:**  $\log(m)$ , with a large constant.

typedef unsigned long long ull;

```
ull sumsq(ull to) { return to /2 * ((to-1) | 1); }
ull divsum(ull to, ull c, ull k, ull m) {
  ull res = k / m * sumsq(to) + c / m * to;
  k %= m; c %= m;
  if (!k) return res;
  ull to2 = (to * k + c) / m;
  return res + (to - 1) * to2 - divsum(to2, m-1 - c, m,
       k);
ll modsum(ull to, ll c, ll k, ll m) {
  c = ((c \% m) + m) \% m;
  k = ((k % m) + m) % m;
  return to \star c + k \star sumsq(to) - m \star divsum(to, c, k,
      m);
```

#### ModMulLL.h

 $7.2 \cdot 10^{18}$ .

**Time:**  $\mathcal{O}(1)$  for modmul,  $\mathcal{O}(\log b)$  for modpow

```
typedef unsigned long long ull;
ull modmul(ull a, ull b, ull M) {
 ll ret = a * b - M * ull(1.L / M * a * b);
  return ret + M * (ret < 0) - M * (ret >= (11)M);
ull modpow(ull b, ull e, ull mod) {
 ull ans = 1;
  for (; e; b = modmul(b, b, mod), e /= 2)
   if (e & 1) ans = modmul(ans, b, mod);
  return ans;
```

# ModSart.h

**Description:** Tonelli-Shanks algorithm for modular square roots. Finds x s.t.  $x^2 = a \pmod{p}$  (-x gives the other solution).

**Time:**  $\mathcal{O}(\log^2 p)$  worst case,  $\mathcal{O}(\log p)$  for most p

```
"ModPow.h"
                                                      24 lines
ll sgrt(ll a, ll p) {
  a \% = p; if (a < 0) a += p;
  if (a == 0) return 0;
 assert (modpow(a, (p-1)/2, p) == 1); // else no
      solution
  if (p % 4 == 3) return modpow(a, (p+1)/4, p);
  // a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8
      == 5
  11 s = p - 1, n = 2;
  int r = 0, m;
  while (s % 2 == 0)
   ++r, s /= 2;
  while (modpow(n, (p-1) / 2, p) != p-1) ++n;
 11 x = modpow(a, (s + 1) / 2, p);
  ll b = modpow(a, s, p), g = modpow(n, s, p);
  for (;; r = m)  {
   11 t = b;
   for (m = 0; m < r && t != 1; ++m)
```

```
t = t * t % p;
if (m == 0) return x;
ll qs = modpow(q, 1LL \ll (r - m - 1), p);
q = qs * qs % p;
x = x * qs % p;
b = b * q % p;
```

#### PrimRoot.h

**Description:** Given an integer n it return an integer g that  $\{q^k:k\in\mathbb{N}_0\}=\mathbb{Z}_n*$ . The primitive Root is also called generator of the group  $\mathbb{Z}_n$ \*. This code is only valid when n is a prime number. if n is  $p^k$  and g is a primitive root of p:  $q^p \equiv q \pmod{p^2} \iff q + p \text{ is a primitive root of } n$  $g^p \not\equiv g \pmod{p^2} \iff g \text{ is a primitive root of } n$ if n is  $2 \cdot p^k$  and q is a primitive root of  $p^k$ :

**Time:** Assuming the generalized Riemann hypothesis,  $\mathcal{O}(\log^8 n)$ 

q is odd  $\iff$  q is a primitive root of n

q is even  $\iff q + p^k$  is a primitive root of n

```
"ModPow.h"
                                                      19 lines
int generator (int p) {
    vector<int> fact;
    int phi = p-1, n = phi;
    for (int i=2; i*i<=n; ++i)
        if (n % i == 0) {
            fact.push_back (i);
            while (n \% i == 0) n /= i;
    if (n > 1)
        fact.push_back (n);
    for (int res=2; res<=p; ++res) {</pre>
        bool ok = true;
```

 $\textbf{3.1.1} \quad \overset{\text{for (size-}}{\textbf{bigital-}} \overset{\text{i=0;}}{\textbf{poot}} \overset{\text{(<fact.size()}}{\text{(res, phi / fact[i], p)}} \overset{\text{(& ok; ++i)}}{\text{(= i)}}$ Given an integer n randra base b. We call the digital root,  $db_b(n)$ , of n the sum of its digits in the base b:  $db_b(n) = 1 + ((n-1) \mod (b-1))$ 

```
1. db_b(x+y) = db_b(db_b(x) + db_b(y))
```

2.  $db_b(xy) = db_b(db_b(x)db_b(y))$ 

```
3. db_b(x-y) \equiv db_b(x) - db_b(y) \pmod{b-1}
```

# Primality

MillerRabin.h

Description: Deterministic Miller-Rabin primality test. Guaranteed to work for numbers up to  $7 \cdot 10^{18}$ ; for larger numbers, use Python and extend A randomly.

**Time:** 7 times the complexity of  $a^b \mod c$ .

```
"ModMulLL.h"
bool MillerRabin (unsigned int n) { // returns true if n
     is prime, else returns false.
 unsigned int r = __builtin_ctzll(n-1); int d = n >> r
  for (int a: {2, 3, 5, 7, 11, 13, 17, 19, 23, 29, 31,
       37, 41}) {
   if (n == a) return true;
    unsigned int x = modpow(a, d, n), res = !(x == 1)|
         x == n - 1);
    for (int i = 1; i < r; i++) x = (__int128)x*x%n, res
        &= (x!=n-1);
    if (res) return false;
 return (n>=2);
```

#### Factor.h

**Description:** Pollard-rho randomized factorization algorithm. Returns prime factors of a number, in arbitrary order (e.g.  $2299 \rightarrow \{11, 19, 11\}$ ). **Time:**  $\mathcal{O}(n^{1/4})$ , less for numbers with small factors.

```
"ModMulLL.h", "MillerRabin.h"
ull pollard(ull n) {
 ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
 auto f = [\&](ull x) \{ return modmul(x, x, n) + i; \};
 while (t++ % 40 | | _{gcd}(prd, n) == 1) {
   if (x == y) x = ++i, y = f(x);
   if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd
   x = f(x), y = f(f(y));
  return __gcd(prd, n);
vector<ull> factor(ull n) {
 if (n == 1) return {};
 if (isPrime(n)) return {n};
 ull x = pollard(n);
 auto l = factor(x), r = factor(n / x);
 l.insert(l.end(), all(r));
 return 1;
```

#### Divisibility 3.3

euclid.h

**Description:** Finds two integers x and y, such that  $ax + by = \gcd(a, b)$ . If you just need gcd, use the built in  $\_gcd$  instead. If a and b are coprime, then x is the inverse of  $a \pmod{b}$ . 5 lines

```
ll euclid(ll a, ll b, ll &x, ll &y) {
 if (!b) return x = 1, y = 0, a;
```

```
11 d = euclid(b, a % b, y, x);
return y -= a/b * x, d;
```

#### CRT.h

Description: Chinese Remainder Theorem.

crt(a, m, b, n) computes x such that  $x\equiv a\pmod m$ ,  $x\equiv b\pmod n$ . If |a|< m and |b|< n, x will obey  $0\le x< \mathrm{lcm}(m,n)$ . Assumes  $mn<2^{62}$ .

Time:  $\log(n)$ 

# 3.3.1 Bézout's identity

$$ax + by = d$$

If (x, y) is one solution, then all solutions are given by

$$\left(x + \frac{kb}{\gcd(a,b)}, y - \frac{ka}{\gcd(a,b)}\right), \quad k \in \mathbb{Z}$$

# 3.4 Fractions

#### ContinuedFractions.h

**Description:** Given N and a real number  $x \ge 0$ , finds the closest rational approximation p/q with  $p,q \le N$ . It will obey  $|p/q - x| \le 1/qN$ . For consecutive convergents,  $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$ .  $(p_k/q_k$  alternates between > x and < x.) If x is rational, y eventually becomes  $\infty$ ; if x is the root of a degree 2 polynomial the a's eventually become cyclic. **Time:**  $\mathcal{O}(\log N)$ 

```
typedef double d; // for N \sim 1e7: long double for N \sim 1
                 e9
pair<11, 11> approximate(d x, 11 N) {
       11 \text{ LP} = 0, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1, 10 = 1
       for (;;) {
               ll lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q :
                           a = (ll) floor(y), b = min(a, lim),
                           NP = b*P + LP, NQ = b*Q + LQ;
               if (a > b) {
                        // If b > a/2, we have a semi-convergent that
                        // better approximation; if b = a/2, we *may*
                                         have one.
                         // Return {P, Q} here for a more canonical
                                          approximation.
                        return (abs(x - (d)NP / (d)NQ) < abs(x - (d)P / (
                                         d)0))?
```

```
make_pair(NP, NQ) : make_pair(P, Q);
}
if (abs(y = 1/(y - (d)a)) > 3*N) {
    return {NP, NQ};
}
LP = P; P = NP;
LQ = Q; Q = NQ;
}
```

#### FracBinarySearch.h

**Description:** Given f and N, finds the smallest fraction  $p/q \in [0,1]$  such that f(p/q) is true, and  $p,q \leq N$ . You may want to throw an exception from f if it finds an exact solution, in which case N can be removed.

```
Usage: fracBS([](Frac f) { return f.p>=3*f.q; }, 10);
// {1,3}
```

Time:  $\mathcal{O}(\log(N))$ 

```
struct Frac { ll p, q; };
template < class F >
Frac fracBS(F f, ll N) {
 bool dir = 1, A = 1, B = 1;
 Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search
      (0, N)
 if (f(lo)) return lo;
 assert(f(hi));
   ll adv = 0, step = 1; // move hi if dir, else lo
    for (int si = 0; step; (step *= 2) >>= si) {
     Frac mid{lo.p * adv + hi.p, lo.g * adv + hi.g};
      if (abs(mid.p) > N || mid.q > N || dir == !f(mid)
        adv -= step; si = 2;
   hi.p += lo.p \star adv;
   hi.q += lo.q * adv;
    dir = !dir;
    swap(lo, hi);
   A = B; B = !!adv;
 return dir ? hi : lo;
```

# 3.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

```
a = k \cdot (m^2 - n^2), \ b = k \cdot (2mn), \ c = k \cdot (m^2 + n^2),
```

with m > n > 0, k > 0,  $m \perp n$ , and either m or n even.

# 3.6 Primes

p=962592769 is such that  $2^{21}\mid p-1$ , which may be useful. For hashing use 970592641 (31-bit number), 31443539979727 (45-bit), 3006703054056749 (52-bit).  $pi(10^6)=78498$ .

Primitive roots exist modulo any prime power  $p^a$ , except for p=2, a>2, and there are  $\phi(\phi(p^a))$  many. For p=2, a>2, the group  $\mathbb{Z}_{2a}^{\times}\cong\mathbb{Z}_2\times\mathbb{Z}_{2a-2}$ .

# 3.7 Important Functions

# 3.7.1 Prime counting function

PrimeCounting.h

**Description:** Given an integer it gives you  $\pi(n)$ .

Time:  $\mathcal{O}\left(n^{3/4}\right)$ 

19 line

```
int count_primes(int n) {
    auto f = [&] (int n) {return 1;}; // (f(ab) = f(a)f(ab))
    auto pref = [&](int n) {return n;}; // should
        return sum_{-}\{i=1..n\} f(i)
    vector<int> v; v.reserve((int)sqrt(n) * 2 + 20);
    int sq, k = 1;
    for (; k * k \le n; k++) v.push_back(k);
    sq = --k;
    if (k * k == n) k--;
    for (; k \ge 1; k--) v.push_back(n / k);
    vector<int> s(v.size());
    for (int i = 0; i < s.size(); i++) s[i] = pref(v[i</pre>
        1) - 1;
    auto geti = [\&] (int x) {return (x<=sq? (int)x-1:(
        int) (v.size() - (n / x)));};
    for (int p = 2; p * p <= n; p++)
        if (s[p-1] != s[p-2])
            for (int i = (int)v.size() - 1; p*p <= v[i]
                 && i >= 0; --i)
                s[i] = (s[geti(v[i] / p)] - s[p-2]) *
                     f(p);
    return s.back();
```

### 3.7.2 Möbius function

$$\mu(n) = \begin{cases} 0 & n \text{ is not square free} \\ 1 & n \text{ has even number of prime factors} \\ -1 & n \text{ has odd number of prime factors} \end{cases}$$

$$g(n) = \sum_{1 \le m \le n} f\left(\left\lfloor \frac{n}{m} \right\rfloor\right) \Leftrightarrow f(n) = \sum_{1 \le m \le n} \mu(m)g\left(\left\lfloor \frac{n}{m} \right\rfloor\right)$$
$$(f * g)(n) = \sum_{d \in \mathcal{D}} f(d)g\left(\frac{n}{d}\right)$$

```
1. Associativity: (f * q) * h = f * (q * h)
```

2. Commutativity: f \* q = q \* f

3. Distributive: f \* (q + h) = f \* q + f \* h

4. Identity element:  $f * \epsilon = \epsilon * f = f$ 

5. Möbius inversion:  $f * 1 = q \iff q * \mu = f$ 

The most important relations are:

```
\epsilon = 1 * \mu
                                              \sigma_1 = \varphi * \sigma_0
\Omega = 1_{\mathcal{P}} * 1
                                     \varphi * 1 = Id
\sigma_k = Id_k * 1 \sigma_0^3 * 1 = (\sigma_0 * 1)^2
\omega = 1_{\mathbb{P}} * 1
```

Where  $\mathcal{P}$  is the set of prime powers and  $\mathbb{P}$  is the set of prime numbers.

# 3.7.3 Multiplicative functions

LinearSieve.h

Description: Linear Sieve for prime numbers. Can be used for precomputing multiplicative functions

```
Time: \mathcal{O}(n)
void sieve () {
 //asignar 1
 for (int i=2; i <= N; i++) {
   if (lp[i] == 0) { // i es primo
      lp[i] = i; pr.push_back(i); // asignar primo
    for (int j = 0; i * pr[j] <= N; j++) {
      lp[i*pr[j]] = pr[j];
      if (i%pr[j] == 0) {/*asignar multiplo*/;break;}
      else{/*asignar coprimo*/}
```

# PrefixSumOpt.h

**Description:** Let f the multiplactive function to compute its prefix sum. Let q and c multiplicative functions so that f \* q = c and both 3 can be computed in constant time.

```
Time: \mathcal{O}\left(n^{2/3}\right)
```

```
unordered_map<int, int> mem;
int calc (int x) {
 if (x <= 0) return 0;
 if (x \le th) return p f(x);
 if (mem.count(x) != 0) return mem[x];
 int ans = 0;
 for (int i = 2, la; i \le x; i = la + 1) {
   la = x/(x/i);
   ans += (p_g(la)-p_g(i-1))*calc(x / i);
```

```
return mem[x] = (p_c(x) - ans)/p_g(1);
```

# Numerical (4)

# Polynomials and 4.1 recurrences

```
Polynomial.h
```

```
struct Poly {
 vector<double> a;
 double operator()(double x) const {
   double val = 0;
   for (int i = sz(a); i--;) (val *= x) += a[i];
   return val:
 void diff() {
   rep(i, 1, sz(a)) a[i-1] = i*a[i];
   a.pop_back();
 void divroot(double x0) {
    double b = a.back(), c; a.back() = 0;
    for (int i=sz(a)-1; i--;) c = a[i], a[i] = a[i+1] *x0
        +b, b=c;
   a.pop_back();
};
```

### PolyRoots.h

```
Description: Finds the real roots to a polynomial.
```

```
Usage: polyRoots(\{\{2, -3, 1\}\}, -1e9, 1e9\}) // solve x^2-3x+2
```

```
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
```

```
"Polynomial.h"
vector<double> polyRoots(Poly p, double xmin, double
   xmax) {
 if (sz(p.a) == 2) \{ return \{-p.a[0]/p.a[1]\}; \}
 vector<double> ret:
 Poly der = p;
 der.diff();
 auto dr = polyRoots(der, xmin, xmax);
 dr.push back(xmin-1);
 dr.push_back(xmax+1);
 sort(all(dr));
 rep(i, 0, sz(dr)-1) {
   double l = dr[i], h = dr[i+1];
   bool sign = p(1) > 0;
   if (sign ^ (p(h) > 0)) {
     rep(it,0,60) { // while (h - l > 1e-8)
       double m = (1 + h) / 2, f = p(m);
       if ((f \le 0) ^ sign) l = m;
       else h = m;
```

```
ret.push_back((1 + h) / 2);
return ret;
```

# PolyInterpolate.h

**Description:** Given n points (x[i], y[i]), computes an n-1-degree polynomial p that passes through them:  $p(x) = a[0] * x^0 + ... + a[n-1] * x^{n-1}$ . For numerical precision, pick  $x[k] = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1$ . Time:  $\mathcal{O}(n^2)$ 

```
typedef vector<double> vd;
vd interpolate(vd x, vd y, int n) {
 vd res(n), temp(n);
 rep(k, 0, n-1) rep(i, k+1, n)
   y[i] = (y[i] - y[k]) / (x[i] - x[k]);
  double last = 0; temp[0] = 1;
  rep(k, 0, n) rep(i, 0, n) {
   res[i] += y[k] * temp[i];
    swap(last, temp[i]);
    temp[i] -= last * x[k];
 return res;
```

#### BerlekampMassey.h

Time:  $\mathcal{O}(N^2)$ 

**Description:** Recovers any *n*-order linear recurrence relation from the first 2n terms of the recurrence. Useful for guessing linear recurrences after brute-forcing the first terms. Should work on any field, but numerical stability for floats is not guaranteed. Output will have size  $\leq n$ . **Usage:** berlekampMassey( $\{0, 1, 1, 3, 5, 11\}$ ) //  $\{1, 2\}$ 

```
"../number-theory/ModPow.h"
vector<ll> berlekampMassey(vector<ll> s) {
  int n = sz(s), L = 0, m = 0;
 vector<ll> C(n), B(n), T;
  C[0] = B[0] = 1;
```

```
11 b = 1;
rep(i, 0, n) \{ ++m;
  ll d = s[i] % mod;
  rep(j, 1, L+1) d = (d + C[j] * s[i - j]) % mod;
  if (!d) continue;
  T = C; 11 coef = d * modpow(b, mod-2) % mod;
  rep(j, m, n) C[j] = (C[j] - coef * B[j - m]) % mod;
  if (2 * L > i) continue;
  L = i + 1 - L; B = T; b = d; m = 0;
C.resize(L + 1); C.erase(C.begin());
for (ll& x : C) x = (mod - x) % mod;
return C:
```

# LinearRecurrence.h

**Description:** Generates the k'th term of an n-order linear recurrence  $S[i] = \sum_j S[i-j-1]tr[j]$ , given  $S[0... \ge n-1]$  and tr[0...n-1]. Faster than matrix multiplication. Useful together with Berlekamp–Massey.

Usage: linearRec( $\{0, 1\}$ ,  $\{1, 1\}$ , k) // k'th Fibonacci number

Time:  $\mathcal{O}\left(n^2 \log k\right)$ 

26 lines

```
typedef vector<ll> Poly;
ll linearRec(Polv S, Polv tr, ll k) {
 int n = sz(tr);
 auto combine = [&](Poly a, Poly b) {
   Poly res(n \star 2 + 1);
   rep(i, 0, n+1) rep(j, 0, n+1)
     res[i + j] = (res[i + j] + a[i] * b[j]) % mod;
    for (int i = 2 * n; i > n; --i) rep(j,0,n)
      res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]
          ]) % mod;
   res.resize(n + 1);
   return res;
 };
 Poly pol(n + 1), e(pol);
 pol[0] = e[1] = 1;
 for (++k; k; k /= 2) {
   if (k % 2) pol = combine(pol, e);
   e = combine(e, e);
 11 \text{ res} = 0;
 rep(i,0,n) res = (res + pol[i + 1] * S[i]) % mod;
 return res:
```

# 4.2 Optimization

a = x1; x1 = x2; f1 = f2;x2 = a + r\*(b-a); f2 = f(x2);

#### GoldenSectionSearch.h

} else {

**Description:** Finds the argument minimizing the function f in the interval [a,b] assuming f is unimodal on the interval, i.e. has only one local minimum and no local maximum. The maximum error in the result is eps. Works equally well for maximization with a small change in the code. See TernarySearch.h in the Various chapter for a discrete version.

```
Usage: double func (double x) { return 4+x+.3*x*x; } double xmin = gss(-1000,1000,func);

Time: \mathcal{O}(\log((b-a)/\epsilon))

double gss(double a, double b, double (*f) (double)) { double r = (sqrt(5)-1)/2, eps = 1e-7; double x1 = b - r*(b-a), x2 = a + r*(b-a); double f1 = f(x1), f2 = f(x2); while (b-a > eps) if (f1 < f2) { //change to > to find maximum b = x2; x2 = x1; f2 = f1; x1 = b - r*(b-a); f1 = f(x1);
```

```
}
return a;
```

### HillClimbing.h

Description: Poor man's optimization for unimodal functions.

```
typedef array<double, 2> P;

template<class F> pair<double, P> hillClimb(P start, F
    f) {
    pair<double, P> cur(f(start), start);
    for (double jmp = 1e9; jmp > 1e-20; jmp /= 2) {
        rep(j,0,100) rep(dx,-1,2) rep(dy,-1,2) {
            P p = cur.second;
            p[0] += dx*jmp;
            p[1] += dy*jmp;
            cur = min(cur, make_pair(f(p), p));
        }
    }
    return cur;
}
```

#### Integrate.h

**Description:** Simple integration of a function over an interval using Simpson's rule. The error should be proportional to  $h^4$ , although in practice you will want to verify that the result is stable to desired precision when epsilon changes.

### IntegrateAdaptive.h

Description: Fast integration using an adaptive Simpson's rule.

```
Usage: double sphereVolume = quad(-1, 1, [](double x) { return quad(-1, 1, [&](double y) { return quad(-1, 1, [&](double z) { return x*x + y*y + z*z < 1; \});});}
```

```
return x*x + y*y + z*z < 1; });});

typedef double d;
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6

template <class F>
d rec(F& f, d a, d b, d eps, d S) {
  d c = (a + b) / 2;
  d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
  if (abs(T - S) <= 15 * eps || b - a < 1e-10)
    return T + (T - S) / 15;
  return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2);
}
template < class F>
```

```
d quad(d a, d b, F f, d eps = 1e-8) {
  return rec(f, a, b, eps, S(a, b));
}
```

#### Simplex.h

for (;;) {

**Description:** Solves a general linear maximization problem: maximize  $c^Tx$  subject to  $Ax \leq b, x \geq 0$ . Returns -inf if there is no solution, inf if there are arbitrarily good solutions, or the maximum value of  $c^Tx$  otherwise. The input vector is set to an optimal x (or in the unbounded case, an arbitrary solution fulfilling the constraints). Numerical stability is not guaranteed. For better performance, define variables such that x=0 is viable.

```
Usage: vvd A = {{1,-1}, {-1,1}, {-1,-2}};
vd b = {1,1,-4}, c = {-1,-1}, x;
T val = LPSolver(A, b, c).solve(x);
```

**Time:**  $\mathcal{O}(NM * \#pivots)$ , where a pivot may be e.g. an edge relaxation.  $\mathcal{O}(2^n)$  in the general case.

```
typedef double T; // long double, Rational, double +
    mod < P > \dots
typedef vector<T> vd;
typedef vector<vd> vvd;
const T eps = 1e-8, inf = 1/.0;
#define MP make_pair
\#define ltj(X) if (s == -1 \mid \mid MP(X[j], N[j]) < MP(X[s], N[
    s])) s=j
struct LPSolver {
 int m. n:
 vi N, B;
 vvd D;
 LPSolver(const vvd& A, const vd& b, const vd& c) :
    m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2), vd(n+2)) {
      rep(i, 0, m) rep(j, 0, n) D[i][j] = A[i][j];
      rep(i,0,m) \{ B[i] = n+i; D[i][n] = -1; D[i][n+1] 
          = b[i];
      rep(j, 0, n) \{ N[j] = j; D[m][j] = -c[j]; \}
      N[n] = -1; D[m+1][n] = 1;
  void pivot(int r, int s) {
    T *a = D[r].data(), inv = 1 / a[s];
    rep(i, 0, m+2) if (i != r \&\& abs(D[i][s]) > eps) {
      T *b = D[i].data(), inv2 = b[s] * inv;
      rep(j, 0, n+2) b[j] = a[j] * inv2;
      b[s] = a[s] * inv2;
    rep(j,0,n+2) if (j != s) D[r][j] \star= inv;
    rep(i,0,m+2) if (i != r) D[i][s] *= -inv;
    D[r][s] = inv;
    swap(B[r], N[s]);
 bool simplex(int phase) {
   int x = m + phase - 1;
```

```
int s = -1;
      rep(j, 0, n+1) if (N[j] != -phase) ltj(D[x]);
      if (D[x][s] >= -eps) return true;
      int r = -1;
      rep(i,0,m) {
        if (D[i][s] <= eps) continue;</pre>
        if (r == -1 || MP(D[i][n+1] / D[i][s], B[i])
                     < MP(D[r][n+1] / D[r][s], B[r])) r
      if (r == -1) return false:
      pivot(r, s);
  T solve(vd &x) {
    int r = 0:
    rep(i, 1, m) if (D[i][n+1] < D[r][n+1]) r = i;
    if (D[r][n+1] < -eps) {
      pivot(r, n);
      if (!simplex(2) \mid \mid D[m+1][n+1] < -eps) return -
      rep(i, 0, m) if (B[i] == -1) {
        int s = 0;
        rep(j,1,n+1) ltj(D[i]);
        pivot(i, s);
    bool ok = simplex(1); x = vd(n);
    rep(i, 0, m) if (B[i] < n) x[B[i]] = D[i][n+1];
    return ok ? D[m][n+1] : inf;
};
```

# 4.3 Matrices

#### Determinant.h

**Description:** Calculates determinant of a matrix. Destroys the matrix. **Time:**  $\mathcal{O}\left(N^3\right)$ 

```
double det(vector<vector<double>>& a) {
  int n = sz(a); double res = 1;
  rep(i,0,n) {
    int b = i;
    rep(j,i+1,n) if (fabs(a[j][i]) > fabs(a[b][i])) b =
        j;
    if (i != b) swap(a[i], a[b]), res *= -1;
    res *= a[i][i];
    if (res == 0) return 0;
    rep(j,i+1,n) {
        double v = a[j][i] / a[i][i];
        if (v != 0) rep(k,i+1,n) a[j][k] -= v * a[i][k];
    }
}
return res;
}
```

#### IntDeterminant.h

**Description:** Calculates determinant using modular arithmetics. Modulos can also be removed to get a pure-integer version.

Time:  $\mathcal{O}\left(N^3\right)$ 

```
const ll mod = 12345;
ll det(vector<vector<ll>>& a) {
  int n = sz(a); ll ans = 1;
  rep(i,0,n) {
    rep(j,i+1,n) {
    while (a[j][i] != 0) { // gcd step
        ll t = a[i][i] / a[j][i];
        if (t) rep(k,i,n)
            a[i][k] = (a[i][k] - a[j][k] * t) % mod;
        swap(a[i], a[j]);
        ans *= -1;
    }
}
ans = ans * a[i][i] % mod;
if (!ans) return 0;
}
return (ans + mod) % mod;
}
```

#### SolveLinear.h

**Description:** Solves A\*x=b. If there are multiple solutions, an arbitrary one is returned. Returns rank, or -1 if no solutions. Data in A and b is lost.

Time:  $\mathcal{O}\left(n^2m\right)$ 

```
typedef vector<double> vd;
const double eps = 1e-12;
int solveLinear(vector<vd>& A, vd& b, vd& x) {
 int n = sz(A), m = sz(x), rank = 0, br, bc;
 if (n) assert(sz(A[0]) == m);
 vi col(m); iota(all(col), 0);
 rep(i,0,n) {
   double v, bv = 0;
   rep(r,i,n) rep(c,i,m)
     if ((v = fabs(A[r][c])) > bv)
       br = r, bc = c, bv = v;
   if (bv <= eps) {
      rep(j,i,n) if (fabs(b[j]) > eps) return -1;
      break;
   swap(A[i], A[br]);
   swap(b[i], b[br]);
   swap(col[i], col[bc]);
   rep(j,0,n) swap(A[j][i], A[j][bc]);
   bv = 1/A[i][i];
   rep(j,i+1,n) {
     double fac = A[j][i] * bv;
     b[j] = fac * b[i];
     rep(k, i+1, m) A[j][k] = fac*A[i][k];
   rank++;
```

```
x.assign(m, 0);
for (int i = rank; i--;) {
   b[i] /= A[i][i];
   x[col[i]] = b[i];
   rep(j,0,i) b[j] -= A[j][i] * b[i];
}
return rank; // (multiple solutions if rank < m)
}</pre>
```

#### SolveLinear2.h

**Description:** To get all uniquely determined values of x back from SolveLinear, make the following changes:

#### SolveLinearBinary.h

**Description:** Solves Ax = b over  $\mathbb{F}_2$ . If there are multiple solutions, one is returned arbitrarily. Returns rank, or -1 if no solutions. Destroys A and b.

```
Time: \mathcal{O}(n^2m)
```

34 lines

```
typedef bitset<1000> bs;
int solveLinear(vector<bs>& A, vi& b, bs& x, int m) {
 int n = sz(A), rank = 0, br;
  assert(m \le sz(x));
 vi col(m); iota(all(col), 0);
  rep(i,0,n) {
    for (br=i; br<n; ++br) if (A[br].any()) break;</pre>
    if (br == n) {
      rep(j,i,n) if(b[j]) return -1;
      break;
    int bc = (int)A[br]._Find_next(i-1);
    swap(A[i], A[br]);
    swap(b[i], b[br]);
    swap(col[i], col[bc]);
    rep(j, 0, n) if (A[j][i] != A[j][bc]) {
     A[j].flip(i); A[j].flip(bc);
    rep(j,i+1,n) if (A[j][i]) {
     b[j] ^= b[i];
     A[j] ^= A[i];
    rank++;
 x = bs();
 for (int i = rank; i--;) {
    if (!b[i]) continue;
    x[col[i]] = 1;
```

```
rep(j, 0, i) b[j] ^= A[j][i];
return rank; // (multiple solutions if rank < m)
```

#### MatrixInverse.h

**Description:** Invert matrix A. Returns rank; result is stored in A unless singular (rank < n). Can easily be extended to prime moduli; for prime powers, repeatedly set  $A^{-1} = A^{-1}(2I - AA^{-1}) \pmod{p^k}$  where  $A^{-1}$  starts as the inverse of A mod p, and k is doubled in each step.

Time:  $\mathcal{O}\left(n^3\right)$ 

```
int matInv(vector<vector<double>>& A) {
 int n = sz(A); vi col(n);
 vector<vector<double>> tmp(n, vector<double>(n));
 rep(i, 0, n) tmp[i][i] = 1, col[i] = i;
 rep(i,0,n) {
   int r = i, c = i;
   rep(j,i,n) rep(k,i,n)
     if (fabs(A[j][k]) > fabs(A[r][c]))
       r = j, c = k;
   if (fabs(A[r][c]) < 1e-12) return i;</pre>
   A[i].swap(A[r]); tmp[i].swap(tmp[r]);
   rep(j,0,n)
      swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c
          ]);
    swap(col[i], col[c]);
   double v = A[i][i];
    rep(j,i+1,n) {
      double f = A[j][i] / v;
     A[j][i] = 0;
     rep(k, i+1, n) A[j][k] -= f*A[i][k];
     rep(k,0,n) tmp[j][k] -= f*tmp[i][k];
   rep(j,i+1,n) A[i][j] /= v;
   rep(j,0,n) tmp[i][j] /= v;
   A[i][i] = 1;
 for (int i = n-1; i > 0; --i) rep(j,0,i) {
   double v = A[i][i];
   rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
 rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
 return n;
```

### Tridiagonal.h

**Description:** x = tridiagonal(d, p, q, b) solves the equation system

$$\begin{pmatrix} & b_0 \\ & b_1 \\ & b_2 \\ & b_3 \\ & \vdots \\ & b_{n-1} \end{pmatrix} = \begin{pmatrix} d_0 & p_0 & 0 & 0 & \cdots & 0 \\ q_0 & d_1 & p_1 & 0 & \cdots & 0 \\ 0 & q_1 & d_2 & p_2 & \cdots & 0 \\ \vdots & \vdots & \ddots & \ddots & \ddots & \vdots \\ 0 & 0 & \cdots & q_{n-3} & d_{n-2} & p_{n-2} \\ 0 & 0 & \cdots & 0 & q_{n-2} & d_{n-1} \end{pmatrix} \begin{pmatrix} x_0 \\ x_1 \\ x_2 \\ x_3 \\ \vdots \\ x_{n-1} \end{pmatrix}$$

This is useful for solving problems on the type

$$a_i = b_i a_{i-1} + c_i a_{i+1} + d_i, \ 1 \le i \le n,$$

where  $a_0, a_{n+1}, b_i, c_i$  and  $d_i$  are known. a can then be obtained from

$$\begin{aligned} \{a_i\} &= \operatorname{tridiagonal}(\{1,-1,-1,...,-1,1\}, \{0,c_1,c_2,...,c_n\},\\ \{b_1,b_2,...,b_n,0\}, \{a_0,d_1,d_2,...,d_n,a_{n+1}\}). \end{aligned}$$

Fails if the solution is not unique.

If  $|d_i| > |p_i| + |q_{i-1}|$  for all i, or  $|d_i| > |p_{i-1}| + |q_i|$ , or the matrix is positive definite, the algorithm is numerically stable and neither tr nor the check for diag[i] == 0 is needed.

Time:  $\mathcal{O}(N)$ 

26 lines

35 lines

```
typedef double T;
vector<T> tridiagonal(vector<T> diag, const vector<T>&
    super,
   const vector<T>& sub, vector<T> b) {
 int n = sz(b); vi tr(n);
 rep(i, 0, n-1) {
   if (abs(diag[i]) < 1e-9 * abs(super[i])) { // diag[}
        i / == 0
     b[i+1] = b[i] * diag[i+1] / super[i];
      if (i+2 < n) b[i+2] = b[i] * sub[i+1] / super[i]
      diag[i+1] = sub[i]; tr[++i] = 1;
    } else {
      diag[i+1] -= super[i] * sub[i] / diag[i];
     b[i+1] = b[i] * sub[i] / diag[i];
 for (int i = n; i--;) {
   if (tr[i]) {
     swap(b[i], b[i-1]);
      diag[i-1] = diag[i];
     b[i] /= super[i-1];
    } else {
     b[i] /= diag[i];
     if (i) b[i-1] -= b[i] * super[i-1];
   }
 return b;
```

# Fourier transforms

FastFourierTransform.h

**Description:** fft(a) computes  $\hat{f}(k) = \sum_{x} a[x] \exp(2\pi i \cdot kx/N)$  for all k. N must be a power of 2. Useful for convolution: conv(a, b) = c, where  $c[x] = \sum a[i]b[x-i]$ . For convolution of complex numbers or more than two vectors: FFT, multiply pointwise, divide by n, reverse(start+1, end), FFT back. Rounding is safe if  $(\sum a_i^2 + \sum b_i^2) \log_2 N < 9 \cdot 10^{14}$  (in practice 10<sup>16</sup>; higher for random inputs). Otherwise, use NTT/FFT-

```
Time: O(N \log N) with N = |A| + |B| (~1s for N = 2^{22})
```

```
typedef complex<double> C;
typedef vector<double> vd;
void fft(vector<C>& a) {
```

```
int n = sz(a), L = 31 - _builtin_clz(n);
  static vector<complex<long double>> R(2, 1);
  static vector<C> rt(2, 1); // (^ 10% faster if
      double)
  for (static int k = 2; k < n; k *= 2) {
   R.resize(n); rt.resize(n);
    auto x = polar(1.0L, acos(-1.0L) / k);
    rep(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i
 vi rev(n):
  rep(i,0,n) \ rev[i] = (rev[i / 2] | (i \& 1) << L) / 2;
  rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);</pre>
  for (int k = 1; k < n; k *= 2)
   for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
      Cz = rt[j+k] * a[i+j+k]; // (25\% faster if hand-
      a[i + j + k] = a[i + j] - z;
      a[i + j] += z;
vd conv(const vd& a, const vd& b) {
 if (a.empty() || b.empty()) return {};
  vd res(sz(a) + sz(b) - 1);
  int L = 32 - \underline{\text{builtin\_clz}(\text{sz}(\text{res}))}, n = 1 << L;
 vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
  fft(in);
  for (C& x : in) x \star = x;
  rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
 rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4 * n);
 return res;
```

#### FastFourierTransformMod.h

Description: Higher precision FFT, can be used for convolutions modulo arbitrary integers as long as  $N \log_2 N \cdot \text{mod} < 8.6 \cdot 10^{14}$  (in practice 10<sup>16</sup> or higher). Inputs must be in [0, mod).

**Time:**  $\mathcal{O}(N \log N)$ , where N = |A| + |B| (twice as slow as NTT or FFT)

```
"FastFourierTransform.h"
                                                      22 lines
typedef vector<ll> vl;
template<int M> vl convMod(const vl &a, const vl &b) {
 if (a.empty() || b.empty()) return {};
 vl res(sz(a) + sz(b) - 1);
  int B=32-__builtin_clz(sz(res)), n=1<<B, cut=int(sqrt</pre>
  vector < C > L(n), R(n), outs(n), outl(n);
  rep(i, 0, sz(a)) L[i] = C((int)a[i] / cut, (int)a[i] %
  rep(i, 0, sz(b)) R[i] = C((int)b[i] / cut, (int)b[i] %
      cut);
  fft(L), fft(R);
  rep(i,0,n) {
    int j = -i \& (n - 1);
    outl[j] = (L[i] + conj(L[j])) * R[i] / (2.0 * n);
```

#### NumberTheoreticTransform FastSubsetTransform

```
outs[j] = (L[i] - conj(L[j])) * R[i] / (2.0 * n) /
       1i:
fft(outl), fft(outs);
rep(i, 0, sz(res)) {
  ll\ av = ll\ (real\ (outl[i]) + .5),\ cv = ll\ (imag\ (outs[i])
  11 \text{ bv} = 11(\text{imag}(\text{outl}[i]) + .5) + 11(\text{real}(\text{outs}[i]) + .5)
  res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
return res;
```

#### NumberTheoreticTransform.h

**Description:** ntt(a) computes  $\hat{f}(k) = \sum_{x} a[x]g^{xk}$  for all k, where  $q = \text{root}^{(mod-1)/N}$ . N must be a power of 2. Useful for convolution modulo specific nice primes of the form  $2^ab + 1$ , where the convolution result has size at most  $2^a$ . For arbitrary modulo, see FFTMod. conv (a, b) = c, where  $c[x] = \sum a[i]b[x-i]$ . For manual convolution: NTT the inputs, multiply pointwise, divide by n, reverse(start+1, end), NTT back. Inputs must be in [0, mod).

Time:  $\mathcal{O}(N \log N)$ 

```
"../number-theory/ModPow.h"
const 11 mod = (119 << 23) + 1, root = 62; // =
    998244353
// For p < 2^30 there is also e.g. 5 << 25, 7 << 26,
    479 << 21
// and 483 \ll 21 (same root). The last two are > 10^9.
typedef vector<ll> vl:
void ntt(vl &a) {
  int n = sz(a), L = 31 - \underline{builtin_clz(n)};
  static vl rt(2, 1);
  for (static int k = 2, s = 2; k < n; k \neq 2, s++) {
    rt.resize(n);
   ll z[] = \{1, modpow(root, mod >> s)\};
    rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod;
  rep(i,0,n) \ rev[i] = (rev[i / 2] | (i \& 1) << L) / 2;
  rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
  for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += 2 * k) rep(j, 0, k) {
      ll z = rt[j + k] * a[i + j + k] % mod, &ai = a[i
      a[i + j + k] = ai - z + (z > ai ? mod : 0);
      ai += (ai + z >= mod ? z - mod : z);
vl conv(const vl &a, const vl &b) {
 if (a.empty() || b.empty()) return {};
  int s = sz(a) + sz(b) - 1, B = 32 - \underline{builtin\_clz(s)},
      n = 1 << B;
  int inv = modpow(n, mod - 2);
  vl L(a), R(b), out(n);
  L.resize(n), R.resize(n);
  ntt(L), ntt(R);
  rep(i,0,n)
```

```
out[-i \& (n - 1)] = (ll)L[i] * R[i] % mod * inv %
ntt(out);
return {out.begin(), out.begin() + s};
```

#### FastSubsetTransform.h

**Description:** Transform to a basis with fast convolutions of the form  $c[z] = \sum_{z = x \oplus y} a[x] \cdot b[y],$  where  $\oplus$  is one of AND, OR, XOR. The size of a must be a power of two.

Time:  $\mathcal{O}(N \log N)$ 

void FST(vi& a, bool inv) { for (int n = sz(a), step = 1; step < n; step \*= 2) { for (int i = 0; i < n; i += 2 \* step) rep(i,i,i+ step) { int &u = a[j], &v = a[j + step]; tie(u, v) = inv ? pii(v - u, u) : pii(v, u + v); // ANDinv ? pii(v, u - v) : pii(u + v, u); // ORpii(u + v, u - v);if (inv) for (int& x : a) x /= sz(a): // XOR only vi conv(vi a, vi b) { FST(a, 0); FST(b, 0); rep(i, 0, sz(a)) a[i] \*= b[i];FST(a, 1); return a;

# Mathematics (5)

#### 5.1**Equations**

$$ax^2 + bx + c = 0 \Rightarrow x = \frac{-b \pm \sqrt{b^2 - 4ac}}{2a}$$

The extremum is given by x = -b/2a.

$$ax + by = e$$

$$cx + dy = f$$

$$\Rightarrow x = \frac{ed - bf}{ad - bc}$$

$$y = \frac{af - ec}{ad - bc}$$

In general, given an equation Ax = b, the solution to a variable  $x_i$  is given by

$$x_i = \frac{\det A_i'}{\det A}$$

where  $A'_i$  is A with the i'th column replaced by b.

# Recurrences

If  $a_n = c_1 a_{n-1} + \cdots + c_k a_{n-k}$ , and  $r_1, \ldots, r_k$  are distinct roots of  $x^k - c_1 x^{k-1} - \cdots - c_k$ , there are  $d_1, \ldots, d_k$  s.t.

$$a_n = d_1 r_1^n + \dots + d_k r_k^n.$$

Non-distinct roots r become polynomial factors, e.g.  $a_n = (d_1 n + d_2)r^n.$ 

#### 5.3 Trigonometry

$$\sin(v+w) = \sin v \cos w + \cos v \sin w$$
$$\cos(v+w) = \cos v \cos w - \sin v \sin w$$

$$\tan(v+w) = \frac{\tan v + \tan w}{1 - \tan v \tan w}$$
$$\sin v + \sin w = 2\sin\frac{v+w}{2}\cos\frac{v-w}{2}$$
$$\cos v + \cos w = 2\cos\frac{v+w}{2}\cos\frac{v-w}{2}$$

$$(V+W)\tan(v-w)/2 = (V-W)\tan(v+w)/2$$

where V, W are lengths of sides opposite angles v, w.

$$a\cos x + b\sin x = r\cos(x - \phi)$$
  
$$a\sin x + b\cos x = r\sin(x + \phi)$$

where  $r = \sqrt{a^2 + b^2}$ ,  $\phi = \operatorname{atan2}(b, a)$ .

#### 5.4 Geometry

# 5.4.1 Triangles

Side lengths: a, b, c

Semiperimeter: 
$$p = \frac{a+b+c}{2}$$
  
Area:  $A = \sqrt{p(p-a)(p-b)(p-c)}$ 

Circumradius: 
$$R = \frac{abc}{4A}$$

Inradius: 
$$r = \frac{A}{p}$$

Length of median (divides triangle into two equal-area triangles):  $m_a = \frac{1}{2}\sqrt{2b^2 + 2c^2 - a^2}$ 

Length of bisector (divides angles in two):

$$s_a = \sqrt{bc \left[ 1 - \left( \frac{a}{b+c} \right)^2 \right]}$$

UCM

Law of sines:  $\frac{\sin \alpha}{a} = \frac{\sin \beta}{b} = \frac{\sin \gamma}{c} = \frac{1}{2R}$ Law of cosines:  $a^2 = b^2 + c^2 - 2bc \cos \alpha$ 

Law of tangents:  $\frac{a+b}{a-b} = \frac{\tan \frac{\alpha + \beta}{2}}{\tan \frac{\alpha - \beta}{2}}$ 

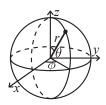
# 5.4.2 Quadrilaterals

With side lengths a, b, c, d, diagonals e, f, diagonals angle  $\theta$ , area A and magic flux  $F = b^2 + d^2 - a^2 - c^2$ :

$$4A = 2ef \cdot \sin \theta = F \tan \theta = \sqrt{4e^2f^2 - F^2}$$

For cyclic quadrilaterals the sum of opposite angles is 180°. ef = ac + bd, and  $A = \sqrt{(p-a)(p-b)(p-c)(p-d)}$ .

# 5.4.3 Spherical coordinates



$$\begin{aligned} x &= r \sin \theta \cos \phi & r &= \sqrt{x^2 + y^2 + z^2} \\ y &= r \sin \theta \sin \phi & \theta &= a \cos(z/\sqrt{x^2 + y^2 + z^2}) \\ z &= r \cos \theta & \phi &= a \tan(y, x) \end{aligned}$$

#### 5.5 Derivatives/Integrals

$$\frac{d}{dx}\arcsin x = \frac{1}{\sqrt{1-x^2}} \quad \frac{d}{dx}\arccos x = -\frac{1}{\sqrt{1-x^2}}$$

$$\frac{d}{dx}\tan x = 1 + \tan^2 x \quad \frac{d}{dx}\arctan x = \frac{1}{1+x^2}$$

$$\int \tan ax = -\frac{\ln|\cos ax|}{a} \quad \int x\sin ax = \frac{\sin ax - ax\cos ax}{a^2}$$

$$\int e^{-x^2} = \frac{\sqrt{\pi}}{2}\operatorname{erf}(x) \quad \int xe^{ax}dx = \frac{e^{ax}}{a^2}(ax-1)$$

Integration by parts:

$$\int_{a}^{b} f(x)g(x)dx = [F(x)g(x)]_{a}^{b} - \int_{a}^{b} F(x)g'(x)dx$$

#### 5.6 Sums

$$c^{a} + c^{a+1} + \dots + c^{b} = \frac{c^{b+1} - c^{a}}{c - 1}, c \neq 1$$

$$B_{n} = 1 - \sum_{k=0}^{n-1} {n \choose k} \frac{B_{k}}{n - k + 1}$$

$$\sum_{i=0}^{n-1} i^{p} = \frac{1}{p+1} \sum_{r=0}^{p} {p+1 \choose r} B_{r} n^{p+1-r}$$

$$1 + 2 + 3 + \dots + n = \frac{n(n+1)}{2}$$

$$1^{2} + 2^{2} + 3^{2} + \dots + n^{2} = \frac{n(2n+1)(n+1)}{6}$$

$$1^{3} + 2^{3} + 3^{3} + \dots + n^{3} = \frac{n^{2}(n+1)^{2}}{4}$$

$$1^{4} + 2^{4} + 3^{4} + \dots + n^{4} = \frac{n(n+1)(2n+1)(3n^{2} + 3n - 1)}{30}$$

$$1^{5} + 2^{5} + 3^{5} + \dots + n^{5} = \frac{n^{2}(n+1)^{2}(2n^{2} + 2n - 1)}{12}$$

$$1^{6} + 2^{6} + 3^{6} + \dots + n^{6} = \frac{(n^{3} + 2n^{2} + x)(3n^{4} + 6n^{3} - 3n + 1)}{42}$$

5.7 Series
$$e^{x} = 1 + x + \frac{x^{2}}{2!} + \frac{x^{3}}{3!} + \dots, (-\infty < x < \infty)$$

$$\ln(1+x) = x - \frac{x^{2}}{2} + \frac{x^{3}}{3} - \frac{x^{4}}{4} + \dots, (-1 < x \le 1)$$

$$\sqrt{1+x} = 1 + \frac{x}{2} - \frac{x^{2}}{8} + \frac{2x^{3}}{32} - \frac{5x^{4}}{128} + \dots, (-1 \le x \le 1)$$

$$\sin x = x - \frac{x^{3}}{3!} + \frac{x^{5}}{5!} - \frac{x^{7}}{7!} + \dots, (-\infty < x < \infty)$$

$$\cos x = 1 - \frac{x^{2}}{2!} + \frac{x^{4}}{4!} - \frac{x^{6}}{6!} + \dots, (-\infty < x < \infty)$$

#### Probability theory 5.8

Let X be a discrete random variable with probability  $p_X(x)$ of assuming the value x. It will then have an expected value (mean)  $\mu = \mathbb{E}(X) = \sum_{x} x p_X(x)$  and variance  $\sigma^2 = V(X) = \mathbb{E}(X^2) - (\mathbb{E}(X))^2 = \sum_x (x - \mathbb{E}(X))^2 p_X(x)$ where  $\sigma$  is the standard deviation. If X is instead continuous it will have a probability density function  $f_X(x)$ and the sums above will instead be integrals with  $p_X(x)$ replaced by  $f_X(x)$ .

Expectation is linear:

$$\mathbb{E}(aX + bY) = a\mathbb{E}(X) + b\mathbb{E}(Y)$$

For independent X and Y,

$$V(aX + bY) = a^2V(X) + b^2V(Y).$$

# 5.8.1 Discrete distributions Binomial distribution

The number of successes in n independent yes/no experiments, each which yields success with probability p is  $Bin(n, p), n = 1, 2, ..., 0 \le p \le 1.$ 

$$p(k) = \binom{n}{k} p^k (1-p)^{n-k}$$

$$\mu = np, \, \sigma^2 = np(1-p)$$

Bin(n, p) is approximately Po(np) for small p.

#### First success distribution

The number of trials needed to get the first success in independent yes/no experiments, each which yields success with probability p is Fs(p), 0 .

$$p(k) = p(1-p)^{k-1}, k = 1, 2, \dots$$
  
$$\mu = \frac{1}{p}, \sigma^2 = \frac{1-p}{p^2}$$

# Poisson distribution

The number of events occurring in a fixed period of time t if these events occur with a known average rate  $\kappa$  and independently of the time since the last event is  $Po(\lambda), \lambda = t\kappa.$ 

$$p(k) = e^{-\lambda} \frac{\lambda^k}{k!}, k = 0, 1, 2, \dots$$
$$\mu = \lambda, \sigma^2 = \lambda$$

# 5.8.2 Continuous distributions Uniform distribution

If the probability density function is constant between a and b and 0 elsewhere it is U(a, b), a < b.

$$f(x) = \begin{cases} \frac{1}{b-a} & a < x < b \\ 0 & \text{otherwise} \end{cases}$$

$$\mu = \frac{a+b}{2}, \ \sigma^2 = \frac{(b-a)^2}{12}$$

# **Exponential distribution**

The time between events in a Poisson process is  $\operatorname{Exp}(\lambda)$ ,  $\lambda > 0$ .

$$f(x) = \begin{cases} \lambda e^{-\lambda x} & x \ge 0\\ 0 & x < 0 \end{cases}$$
$$\mu = \frac{1}{\lambda}, \, \sigma^2 = \frac{1}{\lambda^2}$$

#### Normal distribution

Most real random values with mean  $\mu$  and variance  $\sigma^2$  are well described by  $\mathcal{N}(\mu, \sigma^2)$ ,  $\sigma > 0$ .

$$f(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

If  $X_1 \sim \mathcal{N}(\mu_1, \sigma_1^2)$  and  $X_2 \sim \mathcal{N}(\mu_2, \sigma_2^2)$  then

$$aX_1 + bX_2 + c \sim \mathcal{N}(\mu_1 + \mu_2 + c, a^2\sigma_1^2 + b^2\sigma_2^2)$$

# 5.9 Markov chains

A Markov chain is a discrete random process with the property that the next state depends only on the current state. Let  $X_1, X_2, \ldots$  be a sequence of random variables generated by the Markov process. Then there is a transition matrix  $\mathbf{P} = (p_{ij})$ , with  $p_{ij} = \Pr(X_n = i | X_{n-1} = j)$ , and  $\mathbf{p}^{(n)} = \mathbf{P}^n \mathbf{p}^{(0)}$  is the probability distribution for  $X_n$  (i.e.,  $p_i^{(n)} = \Pr(X_n = i)$ ), where  $\mathbf{p}^{(0)}$  is the initial distribution.

# Stationary distribution

 $\pi$  is a stationary distribution if  $\pi = \pi \mathbf{P}$ . If the Markov chain is irreducible (it is possible to get to any state from any state), then  $\pi_i = \frac{1}{\mathbb{E}(T_i)}$  where  $\mathbb{E}(T_i)$  is the expected time between two visits in state i.  $\pi_j/\pi_i$  is the expected number of visits in state j between two visits in state i.

For a connected, undirected and non-bipartite graph, where the transition probability is uniform among all neighbors,  $\pi_i$  is proportional to node *i*'s degree.

# Ergodicity

A Markov chain is *ergodic* if the asymptotic distribution is independent of the initial distribution. A finite Markov chain is ergodic iff it is irreducible and *aperiodic* (i.e., the gcd of cycle lengths is 1).  $\lim_{k\to\infty} \mathbf{P}^k = \mathbf{1}\pi$ .

# Absorption

A Markov chain is an A-chain if the states can be partitioned into two sets **A** and **G**, such that all states in **A** are absorbing  $(p_{ii} = 1)$ , and all states in **G** leads to an absorbing state in **A**. The probability for absorption in state  $i \in \mathbf{A}$ , when the initial state is j, is  $a_{ij} = p_{ij} + \sum_{k \in \mathbf{G}} a_{ik} p_{kj}$ . The expected time until absorption, when the initial state is i, is  $t_i = 1 + \sum_{k \in \mathbf{G}} p_{ki} t_k$ .

# Techniques (A)

# techniques.txt

159 lines

Recursion Divide and conquer Finding interesting points in N  $\log$  N Algorithm analysis Master theorem Amortized time complexity Greedy algorithm Scheduling Max contiguous subvector sum Invariants Huffman encoding Graph theory Dynamic graphs (extra book-keeping) Breadth first search Depth first search \* Normal trees / DFS trees Dijkstra's algorithm MST: Prim's algorithm Bellman-Ford Konig's theorem and vertex cover Min-cost max flow Lovasz toggle Matrix tree theorem Maximal matching, general graphs Hopcroft-Karp Hall's marriage theorem Graphical sequences Floyd-Warshall Euler cycles Flow networks \* Augmenting paths \* Edmonds-Karp Bipartite matching Min. path cover Topological sorting Strongly connected components Cut vertices, cut-edges and biconnected components Edge coloring \* Trees Vertex coloring \* Bipartite graphs (=> trees) \* 3^n (special case of set cover) Diameter and centroid K'th shortest path Shortest cycle Dynamic programming Knapsack Coin change Longest common subsequence Longest increasing subsequence Number of paths in a dag Shortest path in a dag

Dynprog over intervals

Dynprog over subsets Dynprog over probabilities Dynprog over trees 3^n set cover Divide and conquer Knuth optimization Convex hull optimizations RMQ (sparse table a.k.a 2^k-jumps) Bitonic cycle Log partitioning (loop over most restricted) Combinatorics Computation of binomial coefficients Pigeon-hole principle Inclusion/exclusion Catalan number Pick's theorem Number theory Integer parts Divisibility Euclidean algorithm Modular arithmetic \* Modular multiplication \* Modular inverses \* Modular exponentiation by squaring Chinese remainder theorem Fermat's little theorem Euler's theorem Phi function Frobenius number Quadratic reciprocity Pollard-Rho Miller-Rabin Hensel lifting Vieta root jumping Game theory Combinatorial games Game trees Mini-max Games on graphs Games on graphs with loops Grundy numbers Bipartite games without repetition General games without repetition Alpha-beta pruning Probability theory Optimization Binary search Ternary search Unimodality and convex functions Binary search on derivative Numerical methods Numeric integration Newton's method Root-finding with binary/ternary search Golden section search Matrices Gaussian elimination

Exponentiation by squaring Sorting Radix sort Geometry Coordinates and vectors \* Cross product \* Scalar product Convex hull Polygon cut Closest pair Coordinate-compression Ouadtrees KD-trees All segment-segment intersection Sweeping Discretization (convert to events and sweep) Angle sweeping Line sweeping Discrete second derivatives Strings Longest common substring Palindrome subsequences Knuth-Morris-Pratt Tries Rolling polynomial hashes Suffix array Suffix tree Aho-Corasick Manacher's algorithm Letter position lists Combinatorial search Meet in the middle Brute-force with pruning Best-first (A\*) Bidirectional search Iterative deepening DFS / A\* Data structures LCA (2<sup>k</sup>-jumps in trees in general) Pull/push-technique on trees Heavy-light decomposition Centroid decomposition Lazy propagation Self-balancing trees Convex hull trick (wcipeg.com/wiki/Convex\_hull\_trick) Monotone queues / monotone stacks / sliding queues Sliding queue using 2 stacks Persistent segment tree

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