



Moscow IPT

Ctrl-XD

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Numerical (1)

1.1 Newton’s method

To compute $B = \frac{1}{A}$ modulo x^m : define $B_1 = \{inv(A[0])\}$ and $B_{2n} = B_n(2 - A \cdot B_n)$.

To compute $B = \sqrt{A}$ modulo x^m : define $B_1 = \{\sqrt{A[0]}\}$ and $B_{2n} = \frac{B_n}{2} + \frac{A}{2B_n}$.

To compute $B = \log(1 + xA)$ modulo x^m : $B = \int \frac{(1+xA)'}{1+xA}$.

If T is EGF for some objects then $C = -\log(1 - T) = \sum_{k=1}^{+\infty} \frac{T^k}{k}$ is EGF for cycles of them.

To compute $B = e^{xA}$ modulo x^m : define $B_1 = \{1\}$ and $B_{2n} = B_n(1 + A - \log B_n)$.

If T is EGF for some objects then $F = e^T = \sum_{k=0}^{\infty} \frac{T^k}{k!}$ is EGF for their unordered combinations.

In general case you have equation $P(B, x) = 0$. (f.e. $\log B - A(x) = 0$).

The transition is $B_{2n-\alpha} = B_n - \frac{P(B,x)}{P'_B(B,x)}$. (if $P'_B(B,x)$ is divisible by x^α).

FFT

Description: Applies the discrete Fourier transform to a sequence of numbers modulo MOD.
Time: $\mathcal{O}(n \log n)$

```
int rev[N], root[N];

void init(int n) {
    static int last_init = -1;
    if (n == last_init) return;
    last_init = n;
    for (int i = 1; i < n; ++i) {
        rev[i] = (rev[i >> 1] >> 1) | (i & 1) * (n >> 1);
    }
    const int root_n = binpow(ROOT, (MOD - 1) / n);
    for (int i = 0, cur = 1; i < n / 2; ++i) {
        root[i + n / 2] = cur;
        cur = mul(cur, root_n);
    }
    for (int i = n / 2 - 1; i >= 0; --i) {
        root[i] = root[i << 1];
    }
}
```

```
    }
}

void dft(int* f, int n, bool inverse = false) {
    init(n);
    for (int i = 0; i < n; ++i) {
        if (i < rev[i]) swap(f[i], f[rev[i]]);
    }
    for (int k = 1; k < n; k <= 1)
        for (int i = 0; i < n; i += (k < 1))
            for (int j = 0; j < k; ++j) {
                int z = mul(f[i + j + k], root[j + k]);
                f[i + j + k] = add(f[i + j], MOD - z);
                f[i + j] = add(f[i + j], z);
            }
    if (inverse) {
        reverse(f + 1, f + n);
        const int inv_n = inv(n);
        for (int i = 0; i < n; ++i) f[i] = mul(f[i], inv_n);
    }
}
```

Middle product

Description: Calculates middle-product of two arrays using Tellegen’s principle.
Time: $\mathcal{O}(n \log n)$

```
vector<int> mult(const vector<int>& a, const vector<int>& b) {
    int n = sz(a), m = sz(b), k = 1;
    while (k < n) k <= 1;
    fill_n(fft1, k, 0), fill_n(fft2, k, 0);
    copy(all(a), fft1), copy(all(b), fft2);
    dft(fft1, k, true), dft(fft2, k);
    for (int i = 0; i < k; ++i) fft1[i] = mul(fft1[i], fft2[i]);
    dft(fft1, k);
    return {fft1, fft1 + n - m + 1};
}
```

Berlekamp-Massey

Description: Returns the polynomial of a recurrent sequence of order n from the first $2n$ terms.
Usage: `berlekamp_massey({0, 1, 1, 3, 5, 11}) // {1, -1, -2}`
Time: $\mathcal{O}(n^2)$

```
vector<int> berlekamp_massey(vector<int> s) {
    int n = sz(s), L = 0, m = 0;
    vector<int> c(n), b(n), t;
    c[0] = b[0] = 1;
```

```

int eval = 1;
for (int i = 0; i < n; ++i) {
    m++;
    int delta = 0;
    for (int j = 0; j <= L; ++j) {
        delta = add(delta, mul(c[j], s[i - j]));
    }
    if (delta == 0) continue;
    t = c;
    int coef = mul(delta, inv(eval));
    for (int j = m; j < n; ++j) {
        c[j] = sub(c[j], mul(coef, b[j - m]));
    }
    if (2 * L > i) continue;
    L = i + 1 - L, m = 0, b = t, eval = delta;
}
c.resize(L + 1);
return c;
}

```

Flows (2)

Dinitz

Description: Finds maximum flow using Dinitz algorithm.

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

```

struct Edge {
    int to, cap, flow;
};
vector<Edge> E;
vector<int> gr[N];

int n;
int d[N], ptr[N];

bool bfs(int v0 = 0, int cc = 1) {
    fill(d, d + n, -1);
    d[v0] = 0;
    vector<int> q{v0};
    for (int st = 0; st < sz(q); ++st) {
        int v = q[st];
        for (int id : gr[v]) {
            auto [to, cp, fl] = E[id];
            if (d[to] != -1 || cp - fl < cc) continue;
            d[to] = d[v] + 1;

```

```

            q.emplace_back(to);
        }
    }
    return d[n - 1] != -1;
}

int dfs(int v, int flow, int cc = 1) {
    if (v == n - 1 || !flow) return flow;
    for (; ptr[v] < sz(gr[v]); ++ptr[v]) {
        auto [to, cp, fl] = E[gr[v][ptr[v]]];
        if (d[to] != d[v] + 1 || cp - fl < cc) continue;
        int pushed = dfs(to, min(flow, cp - fl), cc);
        if (pushed) {
            int id = gr[v][ptr[v]];
            E[id].flow += pushed;
            E[id ^ 1].flow -= pushed;
            return pushed;
        }
    }
    return 0;
}

ll dinitz() {
    ll flow = 0;
    for (int c = INF; c > 0; c >>= 1) {
        while (bfs(0, c)) {
            fill(ptr, ptr + n, 0);
            while (int pushed = dfs(0, INF, c))
                flow += pushed;
        }
    }
    return flow;
}

```

MCMF

Description: Finds Minimal Cost Maximal Flow.

```

struct Edge {
    ll to, f, c, w;
};

vector<Edge> E;
vector<int> gr[N];

void add_edge(int u, int v, ll c, ll w) {
    gr[u].push_back(sz(E));

```

```

    E.emplace_back(v, 0, c, w);
    gr[v].push_back(sz(E));
    E.emplace_back(u, 0, 0, -w);
}

pair<ll, ll> mcmf(int n) {
    vector<ll> dist(n);
    vector<ll> pr(n);
    vector<ll> phi(n);
    auto dijkstra = [&] {
        fill(all(dist), INF);
        dist[0] = 0;
        priority_queue<pair<ll, int>, vector<pair<ll, int>>,
            greater<>> pq;
        pq.emplace(0, 0);
        while (!pq.empty()) {
            auto [d, v] = pq.top();
            pq.pop();
            if (d != dist[v]) continue;
            for (int idx : gr[v]) {
                if (E[idx].c == E[idx].f) continue;
                int to = E[idx].to;
                ll w = E[idx].w + phi[v] - phi[to];
                if (dist[to] > d + w) {
                    dist[to] = d + w;
                    pr[to] = idx;
                    pq.emplace(d + w, to);
                }
            }
        }
    };

    ll total_cost = 0, total_flow = 0;
    while (true) {
        dijkstra();
        if (dist[n - 1] == INF) break;
        ll min_cap = INF;
        int cur = n - 1;
        while (cur != 0) {
            min_cap = min(min_cap, E[pr[cur]].c - E[pr[cur]].f);
            cur = E[pr[cur] ^ 1].to;
        }
        cur = n - 1;
        while (cur != 0) {
            E[pr[cur]].f += min_cap;
            E[pr[cur] ^ 1].f -= min_cap;

```

```

        total_cost += min_cap * E[pr[cur]].w;
        cur = E[pr[cur] ^ 1].to;
    }
    total_flow += min_cap;
    for (int i = 0; i < n; ++i) {
        phi[i] += dist[i];
    }
}

return {total_flow, total_cost};
}

```

Number Theory (3)

Extended GCD

Description: Finds two integers x and y , such that $ax + by = \gcd(a, b)$.

```

ll exgcd(ll a, ll b, ll &x, ll &y) {
    if (!b) return x = 1, y = 0, a;
    ll d = euclid(b, a % b, y, x);
    return y -= a/b * x, d;
}

```

CRT

Description: Chinese Remainder Theorem. $\text{crt}(a, m, b, n)$ computes x s.t. $x \equiv a \pmod m, x \equiv b \pmod n$.

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

```

ll crt(ll a, ll m, ll b, ll n) {
    if (n > m) swap(a, b), swap(m, n);
    ll x, y, g = exgcd(m, n, x, y);
    assert((a - b) % g == 0); // else no solution
    x = (b - a) % n * x % n / g * m + a;
    return x < 0 ? x + m * n / g : x;
}

```

Combinatorics (4)

4.1 Derangements

Number of n -permutations that none of the elements appears in their original position.

$$D_n = (n - 1)(D_{n-1} + D_{n-2}) = nD_{n-1} + (-1)^n \approx \frac{n!}{e}$$

4.2 Burnside’s lemma

Given a group G of symmetries and a set Ω , the number of elements of Ω up to symmetry equal

$$\frac{1}{|G|} \sum_{g \in G} N(g)$$

where $N(g)$ is the number of elements fixed by g ($g(x) = x$).

4.3 Stirling numbers (first kind)

Unsigned Stirling numbers of the first kind $c_{n,k}$ is the number of permutations of n elements with k cycles as well as the coefficient on x^k in the expansion $x(x+1)(x+2)\dots(x+(n-1))$.

Signed Stirling numbers of the first kind $s_{n,k}$ is the coefficient on x^k in the expansion $x(x-1)(x-2)\dots(x-(n-1))$.

$$c_{0,0} = s_{0,0} = 1 \quad c_{k,0} = s_{k,0} = 0 \quad c_{n,k} = s_{n,k} = 0 \text{ for } k > n$$

$$c_{n,k} = c_{n-1,k-1} + (n-1)c_{n-1,k}$$

$$s_{n,k} = s_{n-1,k-1} - (n-1)s_{n-1,k}$$

$$s_{n,k} = (-1)^{n+k} c_{n,k}$$

$$\text{EGF: } \sum_{n=0}^\infty \sum_{k=0}^n s_{n,k} \frac{x^n}{n!} y^k = (1+x)^y$$

$$\text{EGF: } \sum_{n=k}^\infty s_{n,k} \frac{x^n}{n!} = \frac{(\log(1+x))^k}{k!}$$

4.4 Stirling numbers (second kind)

The Stirling numbers of the second kind $S(n,k)$, count the number of ways to partition a set of n labelled objects into k nonempty unlabelled subsets.

$$S(n,n) = 1 \text{ for } n \geq 0 \quad S(0,n) = S(n,0) = 0 \text{ for } n > 0$$

$$S(n+1,k) = k \cdot S(n,k) + S(n,k-1)$$

$$S(n,k) = \sum_{t=0}^k \frac{(-1)^{k-t} t^n}{(k-t)! t!}$$

$$\text{EGF: } \sum_{n=0}^\infty \sum_{k=0}^n S(n,k) \frac{x^n}{n!} y^k = e^{y(e^x-1)}$$

$$\text{EGF: } \sum_{n=k}^\infty S(n,k) \frac{x^n}{n!} = \frac{(e^x-1)^k}{k!}$$

4.5 Bell numbers

Bell number B_n is the number of partitions of n labeled elements.

$$B_0 = B_1 = 1$$

$$B_n = \sum_{k=0}^{n-1} C_n^k B_k = \sum_{k=0}^n S(n,k)$$

$$\text{EGF: } \sum_{n=1}^\infty \frac{B_n}{n!} x^n = e^{e^x-1}$$

4.6 Narayana numbers

Narayana number $N(n,k)$ is the number of correct bracket sequences with length $2n$ and exactly k distinct nestings. Also the number of unlabeled ordered rooted trees with $n+1$ vertices and k leaves.

$$N(n,k) = \frac{1}{n} C_n^k C_n^{k-1}$$

4.7 Labeled unrooted trees

Every tree on n vertices has unique sequence of $n-2$ integers from $\{1\dots n\}$ associated with the tree.

Vertex with degree d appears in sequence $d-1$ times.

On n vertices: n^{n-2} .

With degrees d_1, d_2, \dots, d_n : $\frac{(n-2)!}{(d_1-1)! \dots (d_n-1)!}$.

Strings (5)

KMP

Description: Calculates prefix function and Z-function of the given string.

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

```
vector<int> pi(const string& s) {
    vector<int> p(sz(s));
    for (int i = 1; i < sz(s); ++i) {
        int g = p[i - 1];
        while (g && s[i] != s[g]) g = p[g - 1];
        p[i] = g + (s[i] == s[g]);
    }
    return p;
}

vector<int> zf(const string& s) {
    vector<int> z(sz(s));
    int l = -1, r = -1;
    for (int i = 1; i < sz(s); ++i) {
        z[i] = i >= r ? 0 : min(r - i, z[i - 1]);
        while (i + z[i] < sz(s) && s[i + z[i]] == s[z[i]]) ++z[i];
        if (i + z[i] > r) l = i, r = i + z[i];
    }
    return z;
}
```

Aho-Corasick

Description: Builds Aho-Corasick

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

```
const int C = 26;
struct node {
    int nx[C], first = -1, suff = -1; //, zsuff = -1;
    vector<int> idx;
    node() {
        fill(nx, nx + C, -1);
    }
};

vector<node> t(1);
void add_word(string& s, int id) {
    int v = 0;
    for (char ch : s) {
        int x = ch - 'a';
        if (t[v].nx[x] == -1) {
            t[v].nx[x] = sz(t);
            t.emplace_back();
        }
        v = t[v].nx[x];
    }
```

```
    }
    t[v].idx.emplace_back(id);
}

void build_aho() {
    vector<pair<int, int>> q;
    for (int x = 0; x < C; ++x) {
        if (t[0].nx[x] == -1) {
            t[0].nx[x] = 0;
        } else {
            q.emplace_back(0, x);
        }
    }
    for (int st = 0; st < sz(q); ++st) {
        auto [par, x] = q[st];
        int a = t[par].nx[x];
        if (t[par].suff == -1) {
            t[a].suff = 0;
        } else {
            t[a].suff = t[t[par].suff].nx[x];
            // t[a].zsuff = t[t[a].suff].idx.empty() ? t[t[a].suff].zsuff : t[a].suff;
        }
        for (int y = 0; y < C; ++y) {
            if (t[a].nx[y] == -1) {
                t[a].nx[y] = t[t[a].suff].nx[y];
            } else {
                q.emplace_back(a, y);
            }
        }
    }
}
```

Suffix array

Description: Calculates suffix array, inverse suffix array and LCP array of the given string.

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

```
const int M = 1e5 + 10;
vector<int> sa, pos, lcp;

void suffix_array(string& s) {
    int n = sz(s);
    vector<int> c(n), cur(n);
    sa.resize(n), pos.resize(n), lcp.resize(n);
    for (int i = 0; i < n; ++i) {
        sa[i] = i, c[i] = s[i];
    }
```

```

sort(all(sa), [&](int i, int j) { return c[i] < c[j]; });
vector<int> cnt(M);
for (int k = 1; k < n; k <= 1) {
    fill(all(cnt), 0);
    for (int x : c) cnt[x]++;
    for (int i = 1; i < M; ++i) cnt[i] += cnt[i - 1];
    for (int i : sa) {
        int c2 = c[(i - k + n) % n] - 1;
        cur[cnt[c2]++] = (i - k + n) % n;
    }
    swap(cur, sa);
    int x = -1, y = -1, p = 0;
    for (int i : sa) {
        if (c[i] != x || c[(i + k) % n] != y) {
            x = c[i], y = c[(i + k) % n], p++;
        }
        cur[i] = p;
    }
    swap(cur, c);
}
for (int i = 0; i < n; ++i) pos[sa[i]] = i;
int l = 0;
for (int i = 0; i < n; ++i) {
    if (pos[i] == n - 1) {
        l = 0;
    } else {
        while (s[(i + 1) % n] == s[(sa[pos[i] + 1] + 1) % n])
            ++l;
        lcp[pos[i]] = l;
        l = max(0, l - 1);
    }
}
}

```

Minimal rotation

Description: Rotates the given string until it is lexicographically minimal, returns shift.

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

```

int min_rotation(string& s, int len) {
    s += s;
    int i = 0, ans = 0;
    while (i < len) {
        ans = i;
        int j = i + 1, k = i;
        while (j < len * 2 && s[k] <= s[j]) {
            if (s[k] < s[j]) {

```

```

                k = i;
            } else {
                k += 1;
            }
            j += 1;
        }
        while (i <= k) {
            i += j - k;
        }
    }
    s = s.substr(ans, len);
    return ans;
}

```

Miscellaneous (6)

Integrate

Description: Function integration over an interval using Simpson's rule. The error is proportional to h^4 .

```

double integrate(double a, double b, auto&& f, int n = 1000) {
    double h = (b - a) / 2 / n, rs = f(a) + f(b);
    for (int i = 1; i < n * 2; ++i) {
        rs += f(a + i * h) * (i & 1 ? 4 : 2);
    }
    return rs * h / 3;
}

```

Fractional binary search

Description: Finds the smallest fraction $p/q \in [0, 1]$ s.t. $f(p/q)$ is true and $p, q \leq N$.

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

```

struct frac { ll p, q; };

```

```

frac fracBS(auto&& f, ll N) {
    bool dir = true, A = true, B = true;
    frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N]
    if (f(lo)) return lo;
    assert(f(hi));
    while (A || B) {
        ll adv = 0, step = 1;
        for (int si = 0; step; (step *= 2) >>= si) {
            adv += step;
            frac mid{lo.p * adv + hi.p, lo.q * adv + hi.q};
            if (abs(mid.p) > N || mid.q > N || dir == !f(mid)) {

```

```
        adv -= step; si = 2;
    }
}
hi.p += lo.p * adv;
hi.q += lo.q * adv;
dir = !dir;
swap(lo, hi);
A = B; B = !!adv;
}
return dir ? hi : lo;
}
```