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# Contest (1)

# sol.cpp

```
#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())
using 11 = long long;
using pii = pair<int, int>;
using vi = vector<int>;
auto& operator<<(auto&, pair<auto, auto>);
auto operator<<(auto& o, auto x) -> decltype(x.end(), o) {
  for (int i = 0; auto y : x) \circ << ", " + !i++ * 2 << y;
  return o << "}";</pre>
auto& operator<<(auto& o, pair<auto, auto> x) {
 return o << "(" << x.first << ", " << x.second << ")"; }
void __print(auto... x) { ((cerr << " " << x), ...) << endl; }</pre>
#define debug(x...) cerr << "[" #x "]:", __print(x)
#define debug(...) 2137
#endif
int main() {
 cin.tie(0)->sync_with_stdio(0);
```

## .vimrc

```
set nu et ts=2 sw=2
filetype indent on
syntax on
colorscheme habamax
hi MatchParen ctermfg=66 ctermbg=234 cterm=underline
nnoremap ; :
nnoremap : ;
inoremap {<cr> {<cr>}<esc>0 <bs>
```

## Makefile

```
CXXFLAGS=-std=c++20 -Wall -Wextra -Wshadow
 g++ $(CXXFLAGS) -fsanitize=address, undefined -g -DLOCAL \
      sol.cpp -o sol
fast: sol.cpp
  g++ $(CXXFLAGS) -02 sol.cpp -o fast
```

```
#!/bin/bash
for((i=1;i>0;i++)) do
  echo "$i"
  echo "$i" | ./gen > int
 diff -w <(./sol < int) <(./slow < int) || break
hash.sh
cpp -dD -P -fpreprocessed | tr -d '[:space:]' | md5sum |cut -c-6
.bashrc
```

# Struktury danych (2)

#include <ext/pb\_ds/tree\_policy.hpp>

```
OrderedSet.h
Opis: s.find_by_order(k) i s.order_of_key(k).
Czas: \mathcal{O}(\log n)
#include <ext/pb_ds/assoc_container.hpp>
```

```
using namespace gnu pbds;
template<class T>
using ordered_set = tree<T, null_type, less<T>, rb_tree_tag,
                        tree_order_statistics_node_update>;
```

## LineSet.h

};

alias rm='trash'

alias mv='mv -i'

alias cp='cp -i'

3

Opis: Znajduje maksimum funkcji liniowych online. Dla doubli inf = 1/.0, div(a,b) = a/b. Czas:  $\mathcal{O}(\log n)$ 

```
struct line {
 mutable 11 k, m, p;
 bool operator<(const line& o) const { return k < o.k; }</pre>
 bool operator<(11 x) const { return p < x; }</pre>
struct line_set : multiset<line, less<>>> {
 static const 11 inf = LLONG MAX;
 ll div(ll a, ll b) {
   return a / b - ((a ^ b) < 0 && a % b); }
 bool isect(iterator x, iterator y) {
   if (y == end()) return x \rightarrow p = inf, 0;
   if (x->k == y->k) x->p = x->m > y->m ? inf : -inf;
    else x->p = div(y->m - x->m, x->k - y->k);
   return x->p >= y->p;
 void add(ll k, ll m) {
   auto z = insert(\{k, m, 0\}), y = z++, x = y;
   while (isect(y, z)) z = erase(z);
   if (x != begin() \&\& isect(--x, y)) isect(x, y = erase(y));
   while ((y = x) != begin() && (--x)->p >= y->p)
     isect(x, erase(y));
 ll querv(ll x) {
   assert(!emptv());
   auto 1 = *lower bound(x);
   return 1.k * x + 1.m;
```

# Grafy (3)

# 3.1 Przepływy

#### Dinic.h

 $\mathbf{Opis}$ : Dinic ze skalowaniem. Należy ustawić zakres it w flow zgodnie z U. Czas:  $\mathcal{O}(nm \log U)$ 

```
struct dinic {
  struct edge {
    int to, rev;
    11 cap;
 };
 vi lvl, ptr, q;
  vector<vector<edge>> adj;
  dinic(int n) : lvl(n), ptr(n), q(n), adj(n) {}
  void add_edge(int u, int v, ll cap, ll rcap = 0) {
    int i = sz(adj[u]), j = sz(adj[v]);
    adj[u].push_back(\{v, j + (u == v), cap\});
    adj[v].push_back({u, i, rcap});
 11 dfs(int v, int t, ll f) {
    if (v == t || !f) return f;
    for (int& i = ptr[v]; i < sz(adj[v]); i++) {</pre>
      edge& e = adj[v][i];
      if (lvl[e.to] == lvl[v] + 1)
       if (ll p = dfs(e.to, t, min(f, e.cap))) {
          e.cap -= p, adj[e.to][e.rev].cap += p;
          return p;
    return 0;
  11 flow(int s, int t) {
    11 f = 0; q[0] = s;
    for (int it = 29; it >= 0; it--) do {
      lvl = ptr = vi(sz(q));
      int qi = 0, qe = lvl[s] = 1;
      while (qi < qe && !lvl[t]) {
        int v = q[qi++];
        for (edge e : adj[v])
          if (!lvl[e.to] && e.cap >> it)
            q[qe++] = e.to, lvl[e.to] = lvl[v] + 1;
      while (ll p = dfs(s, t, LLONG MAX)) f += p;
    } while (lvl[t]);
    return f;
};
```

## GomorvHu.h

Opis: Tworzy drzewo gdzie min cut to minimum na ścieżce.

Czas:  $\mathcal{O}(n)$  przepływów

```
struct edge { int u, v; ll w; };
vector<edge> gomory_hu(int n, const vector<edge>& ed) {
 vector<edge> t; vi p(n);
 rep(i, 1, n) {
   dinic d(n);
   for (edge e : ed) d.add_edge(e.u, e.v, e.w, e.w);
   t.push_back({i, p[i], d.flow(i, p[i])});
   rep(j, i + 1, n) if (p[j] == p[i] \&\& d.lvl[j]) p[j] = i;
 return t;
```

MCMF.h

## MCMF Matching SCC TwoSat RDSU DMST

```
Czas: \mathcal{O}(Fm \log n)
#include <ext/pb_ds/priority_queue.hpp>
const 11 INF = 2e18;
struct MCMF {
  struct edge {
   int from, to, rev;
   11 cap, cost;
  vector<vector<edge>> adj;
  vector<ll> dst, pi;
  __gnu_pbds::priority_queue<pair<11, int>> q;
  vector<decltype(q)::point_iterator> it;
  vector<edge*> p;
  MCMF(int _n) : n(_n), adj(n), pi(n), p(n) {}
  void add edge(int u, int v, ll cap, ll cost) {
   int i = sz(adj[u]), j = sz(adj[v]);
    adj[u].push_back({u, v, j + (u == v), cap, cost});
    adj[v].push_back(\{v, u, i, 0, -cost\});
 bool path(int s, int t) {
    dst.assign(n, INF); it.assign(n, q.end());
    q.push(\{dst[s] = 0, s\});
    while (!q.empty()) {
     int u = q.top().second; q.pop();
     for (edge& e : adj[u]) {
       11 d = dst[u] + pi[u] + e.cost - pi[e.to];
       if (e.cap && d < dst[e.to]) {</pre>
         dst[e.to] = d, p[e.to] = &e;
         if (it[e.to] == q.end())
            it[e.to] = q.push({-dst[e.to], e.to});
          el se
            q.modify(it[e.to], {-dst[e.to], e.to});
    rep(i, 0, n) pi[i] = min(pi[i] + dst[i], INF);
   return pi[t] != INF;
  pair<11, 11> flow(int s, int t, 11 cap) {
    11 f = 0, c = 0;
    while (f < cap && path(s, t)) {
     11 d = cap - f;
     for (edge* e = p[t]; e; e = p[e->from])
       d = min(d, e->cap);
     for (edge* e = p[t]; e; e = p[e->from])
       e->cap -= d, adj[e->to][e->rev].cap += d;
     f += d, c += d * pi[t];
    return {f, c};
```

Opis: MCMF z Dijkstrą. Jeżeli są ujemne krawędzie to przed puszczeniem

flow w pi trzeba policzyć najkrótsze ścieżki z s.

## Skojarzenia

### Matching.h

Opis: Dinic uproszczony do znajdywania największego skojarzenia. Czas:  $\mathcal{O}(m\sqrt{n})$ 

```
struct matching {
  vector<vi> adi;
 vi l, r, lvl;
  matching (int n, int m) : adj(n), l(n, -1), r(m, -1) {}
```

```
void add_edge(int u, int v) { adj[u].push_back(v); }
 bool dfs(int u) {
    int t = exchange(lvl[u], -1) + 1;
    for (int v : adj[u])
     if (r[v] == -1 || (lvl[r[v]] == t && dfs(r[v])))
        return 1[u] = v, r[v] = u, 1;
    return 0;
 int match() {
   int n = sz(1), ans = 0; vi q(n);
    for (int s = 0, t = 0;; s = t = 0) {
     lvl = vi(n); bool f = 0;
      rep(i, 0, sz(1)) if (l[i] == -1) lvl[i] = 1, q[t++] = i;
      while (s < t) {
       int u = q[s++];
       for (int v : adj[u]) {
         int x = r[v];
         if (x == -1) f = 1;
         else if (!lvl[x]) lvl[x] = lvl[u] + 1, q[t++] = x;
     if (!f) break;
     rep(i, 0, n) if (l[i] == -1) ans += dfs(i);
    return ans;
};
```

## 3.3 **DFS**

### SCC.h

Opis: Znajduje SCC w kolejności topologicznej. Czas:  $\mathcal{O}(n+m)$ 

```
struct SCC {
 int n, t = 0, cnt = 0;
 vector<vi> adi;
 vi val, p, st;
 SCC(int _n) : n(_n), adj(n), val(n), p(n, -1) {}
 void add_edge(int u, int v) { adj[u].push_back(v); }
 int dfs(int u) {
   int low = val[u] = ++t; st.push_back(u);
    for (int v : adj[u]) if (p[v] == -1)
     low = min(low, val[v] ?: dfs(v));
    if (low == val[u]) {
     for (int x = -1; x != u;)
       p[x = st.back()] = cnt, st.pop_back();
     cnt++:
   return low;
 void build() {
   rep(i, 0, n) if (!val[i]) dfs(i);
   rep(i, 0, n) p[i] = cnt - 1 - p[i];
};
```

#### TwoSat.h Opis: 2-SAT. Czas: $\mathcal{O}(n+m)$

```
struct two sat {
 int n;
 vector<pii> ed;
 vector<bool> b;
 two_sat(int _n) : n(_n) {}
 int add_var() { return n++; }
```

```
void either(int x, int y) {
    x = max(2 * x, -1 - 2 * x), y = max(2 * y, -1 - 2 * y);
    ed.push_back({x, y}); }
  void implies(int x, int y) { either(~x, y); }
  void must(int x) { either(x, x); }
  void at_most_one(const vi& v) {
    if (sz(v) <= 1) return;</pre>
    int cur = ~v[0];
    rep(i, 2, sz(v)) {
      int nxt = add var();
      either(cur, ~v[i]); either(cur, nxt);
      either(~v[i], nxt); cur = ~nxt;
    either(cur, ~v[1]);
 bool solve() {
    SCC scc(2 * n);
    for (auto [u, v] : ed)
      scc.add_edge(u ^ 1, v), scc.add_edge(v ^ 1, u);
    scc.build(); b.resize(n, 1);
    rep(i, 0, n) {
      if (scc.p[2 * i] == scc.p[2 * i + 1]) return 0;
      if (scc.p[2 * i] < scc.p[2 * i + 1]) b[i] = 0;
    return 1;
};
```

2

## 3.4 **DSU**

#### RDSU.h

Opis: DSU z rollbackami.

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

```
struct RDSU {
 vi e; vector<pii> st;
 RDSU(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find(e[x]); }</pre>
  int time() { return sz(st); }
  void rollback(int t) {
    for (int i = time(); i-- > t;)
      e[st[i].first] = st[i].second;
    st.resize(t);
 bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push back({a, e[a]});
    st.push_back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
```

#### 3.5 Drzewa

#### DMST.h

```
Opis: Skierowane MST z r. Zwraca {-1, {}} gdy nie istnieje.
Czas: \mathcal{O}(m \log m)
```

```
struct edge { int a, b; ll w; };
struct node {
  edge kev;
  node *1 = 0, *r = 0;
  11 delta = 0;
  void prop() {
   kev.w += delta;
   if (1) 1->delta += delta;
   if (r) r->delta += delta;
   delta = 0:
  edge top() { prop(); return key; }
node* merge(node* a, node* b) {
 if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a:
void pop(node*\& a) { a->prop(); a = merge(a->1, a->r); }
pair<11, vi> dmst(int n, int r, const vector<edge>& g) {
 RDSU uf(n);
  vector<node*> heap(n);
  for (edge e : g) heap[e.b] = merge(heap[e.b], new node{e});
  11 \text{ res} = 0;
  vi seen(n, -1), path(n), par(n);
  seen[r] = r;
  vector<edge> Q(n), in(n, \{-1, -1, -1\}), comp;
  deque<tuple<int, int, vector<edge>>> cycs;
  rep(s, 0, n) {
    int u = s, qi = 0, w;
    while (seen[u] < 0) {</pre>
     if (!heap[u]) return {-1, {}};
      edge e = heap[u]->top();
     heap[u]->delta -= e.w, pop(heap[u]);
     Q[qi] = e, path[qi++] = u, seen[u] = s;
      res += e.w, u = uf.find(e.a);
      if (seen[u] == s) {
       node* cyc = 0;
       int end = qi, time = uf.time();
        do cyc = merge(cyc, heap[w = path[--qi]]);
        while (uf.join(u, w));
       u = uf.find(u), heap[u] = cyc, seen[u] = -1;
        cycs.push_front({u, time, {&Q[qi], &Q[end]}});
    rep(i, 0, qi) in[uf.find(Q[i].b)] = Q[i];
  for (auto& [u, t, cmp] : cycs) {
   uf.rollback(t);
    edge inedge = in[u];
    for (auto& e : cmp) in[uf.find(e.b)] = e;
   in[uf.find(inedge.b)] = inedge;
  rep(i, 0, n) par[i] = in[i].a;
  return {res, par};
```

# Matma (4)

## 4.1 Arytmetyka modularna

GCD.h

 $\mathbf{Opis:}\ \mathrm{Rozszerzony}\ \mathrm{algorytm}\ \mathrm{Euklidesa}.$ 

```
Czas: \mathcal{O}(\log \min(a, b))
```

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

### CRT.h

 $\mathbf{Opis:}\,$  Chińskie twierdzenie o resztach.

Czas:  $\mathcal{O}(\log \min(m, n))$ 

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

#### ModMul.h

**Opis:** Mnożenie i potęgowanie dwóch long longów modulo. Jest to wyraźnie szybsze niż zamiana na \_\_int128.

```
using ull = uint64_t;
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 >= (ll)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;
}
```

#### ModInt.h

```
template<int M, int R>
struct mod {
 static const int MOD = M, ROOT = R;
  mod(11 y = 0) : x(y % M) { x += (x < 0) * M; }
  mod operator+= (mod o) {
   if ((x += 0.x) >= M) x -= M;
    return *this; }
  mod operator-= (mod o) {
    if ((x -= 0.x) < 0) x += M;
    return *this; }
  mod operator*=(mod o) {
   x = 111 * x * o.x % M;
   return *this: }
  mod operator/=(mod o) { return (*this) *= o.inv(); }
  friend mod operator+(mod a, mod b) { return a += b; }
  friend mod operator-(mod a, mod b) { return a -= b; ]
  friend mod operator*(mod a, mod b) { return a *= b; }
  friend mod operator/(mod a, mod b) { return a /= b; }
  auto operator<=>(const mod&) const = default;
  mod pow(ll n) const {
   mod a = x, b = 1;
   for (; n; n /= 2, a *= a) if (n & 1) b *= a;
   return h:
 mod inv() const { return pow(M - 2); }
using mint = mod<998244353, 3>;
```

## 4.2 Liczby pierwsze

### MillerRabin.h

Opis: Test pierwszości Millera-Rabina.

```
bool prime(ull n) {
   if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;
   ull A[] = {2, 325, 9375, 28178, 450775, 9780504, 1795265022},
        s = _builtin_ctzll(n - 1), d = n >> s;
   for (ull a : A) {
      ull p = modpow(a % n, d, n), i = s;
      while (p != 1 && p != n - 1 && a % n && i--)
            p = modmul(p, p, n);
      if (p != n - 1 && i != s) return 0;
   }
   return 1;
}
```

#### ronarumno.n

Opis: Algorytm faktoryzacji rho Pollarda.

Czas:  $\mathcal{O}(n^{1/4})$ 

```
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 = q;
    x = f(x), y = f(f(y));
  }
  return __gcd(prd, n);
}
void factor(ull n, map<ull, int>& cnt) {
  if (n == 1) return;
  if (prime(n)) { cnt[n]++; return; }
  ull x = pollard(n);
  factor(x, cnt); factor(n / x, cnt);
}
```

## 4.3 Wielomiany

#### NTT.h

Czas:  $\mathcal{O}((n+m)\log(n+m))$ 

```
template<class T>
void ntt(vector<T>& a, bool inv) {
 int n = sz(a); vector<T> b(n);
 for (int i = n / 2; i; i /= 2, swap(a, b)) {
   T w = T(T::ROOT).pow((T::MOD - 1) / n * i), m = 1;
   for (int j = 0; j < n; j += 2 * i, m *= w) rep(k, 0, i) {
     T u = a[j + k], v = a[j + k + i] * m;
     b[j / 2 + k] = u + v, b[j / 2 + k + n / 2] = u - v;
 if (inv) {
   reverse(1 + all(a));
   T z = T(n).inv(); rep(i, 0, n) a[i] *= z;
template<class T>
vector<T> conv(vector<T> a, vector<T> b) {
 int s = sz(a) + sz(b) - 1, n = 1 << li>lg(2 * s - 1);
 a.resize(n); ntt(a, 0); b.resize(n); ntt(b, 0);
 rep(i, 0, n) a[i] *= b[i];
 ntt(a, 1); a.resize(s);
 return a;
```

```
Czas: \mathcal{O}((n+m)\log(n+m))
template<class T>
vector<T> mconv(const auto& x, const auto& y) {
  auto con = [&] (const auto& v) {
   vector<T> w(sz(v)); rep(i, 0, sz(v)) w[i] = v[i].x;
    return w; };
  return conv(con(x), con(y));
template<class T>
vector<T> conv3(const vector<T>& a, const vector<T>& b) {
 using m0 = mod<754974721, 11>; auto c0 = mconv<m0>(a, b);
  using m1 = mod<167772161, 3>; auto c1 = mconv<m1>(a, b);
  using m2 = mod<469762049, 3>; auto c2 = mconv<m2>(a, b);
  int n = sz(c0); vector<T> d(n); m1 r01 = m1(m0::MOD).inv();
  m2 r02 = m2 (m0::MOD).inv(), r12 = m2 (m1::MOD).inv();
   int x = c0[i].x, y = ((c1[i] - x) * r01).x,
       z = (((c2[i] - x) * r02 - y) * r12).x;
   d[i] = (T(z) * m1::MOD + y) * m0::MOD + x;
  return d;
```

 $\mbox{\bf Opis:}$ NTT z Garnerem. Działa dla  $n+m \leq 2^{24}$  i  $c_k < 5 \cdot 10^{25}.$ 

## **4.4** Inne

#### SameDiv.h

**Opis:** Dzieli przedział  $[1,\infty)$  na przedziały równości dzielenia. Czas:  $\mathcal{O}(\sqrt{n})$ 

```
vector<1l> same_floor(ll n) {
   vector<1l> v;
   for (ll i = 1; i <= n; i = n / (n / i) + 1) v.push_back(i);
   return v.push_back(n + 1), v;
}
vector<1l> same_ceil(ll n) {
   vector<1l> v;
   for (ll i = 1, j; i < n; i = (n + j - 2) / (j - 1)) {
      j = (n + i - 1) / i;
      v.push_back(i);
   }
   return v.push_back(n), v;
}</pre>
```

# Teksty (5)

## 5.1 Podstawy

#### KMP.h

Opis: p[k] – nadjłuższy ścisły sufiks s[0, k] który jest prefiksem s. Czas:  $\mathcal{O}(n)$ 

```
vi kmp(const string& s) {
  vi p(sz(s));
  rep(i, 1, sz(s)) {
    int g = p[i - 1];
    while (g && s[i] != s[g]) g = p[g - 1];
    p[i] = g + (s[i] == s[g]);
  }
  return p;
}
```

```
Z.h Opis: f[k] – nadjłuższy prefiks s[k, n) który jest prefiksem s. Czas: O(n)
```

```
vi z (const string& s) {
  int n = sz(s), l = -1, r = -1;
  vi f(n); f[0] = n;
  rep(i, l, sz(s)) {
   if (i < r) f[i] = min(r - i, f[i - 1]);
   while (i + f[i] < n && s[i + f[i]] == s[f[i]]) f[i]++;
   if (i + f[i] > r) l = i, r = i + f[i];
  }
  return f;
```

#### Manacher.h

Opis: p[1] [k] – środek w k, p[0] [k] – środek między k – 1 a k. Czas:  $\mathcal{O}(n)$ 

```
array<vi, 2> manacher(const string& s) {
  int n = sz(s);
  array<vi, 2> p = {vi(n + 1), vi(n)};
  rep(z, 0, 2) for (int i = 0, 1 = 0, r = 0; i < n; i++) {
    int t = r - i + !z;
    if (i < r) p[z][i] = min(t, p[z][1 + t]);
    int L = i - p[z][i], R = i + p[z][i] - !z;
    while (L >= 1 && R + 1 < n && s[L - 1] == s[R + 1])
        p[z][i]++, L--, R++;
    if (R > r) 1 = L, r = R;
  }
  return p;
}
```

#### Duval.h

**Opis:** Rozkłada słowo na nierosnący ciąg podsłów mniejszych od swoich wszystkich nietrywialnych sufiksów.

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

```
vi duval(const string& s) {
  int n = sz(s); vi f;
  for (int i = 0; i < n;) {
    int j = i + 1, k = i;
    for (; j < n && s[k] <= s[j]; j++) {
      if (s[k] < s[j]) k = i;
      else ++k;
    }
  for (; i <= k; i += j - k) f.push_back(i);
  }
  return f.push_back(n), f;
}</pre>
```

## 5.2 Struktury sufiksowe

#### SuffixArrav.h

**Opis:** Zawiera pusty sufiks. lcp[k] – najdłuższy wspólny prefiks k-1 i k. Czas:  $\mathcal{O}(n \log n)$ 

```
struct suffix_array {
    vi sa, lcp;
    suffix_array(const string& s, int lim = 128) {
        int n = sz(s) + 1, k = 0, a, b;
        vi x(all(s) + 1), y(n), ws(max(n, lim)), rank(n);
        sa = lcp = y, iota(all(sa), 0);
        for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
            p = j, iota(all(y), n - j);
            rep(i, 0, n) if (sa[i] >= j) y[p++] = sa[i] - j;
            fill(all(ws), 0);
```

# Geometria (6)

## 6.1 Podstawy

#### Point.h

Opis: Podstawowy szablon do geometrii.

```
using D = long double;
const D EPS = 1e-91;
int sgn(D x) \{ return (x > EPS) - (x < -EPS); \}
struct P {
 D x, y;
  P operator+(P o) const { return {x + o.x, y + o.y}; }
  P operator-(P o) const { return {x - o.x, y - o.y}; }
  P operator*(D a) const { return {x * a, y * a}; }
  P operator/(D a) const { return {x / a, y / a}; }
  \verb"auto operator"<=>(P \circ) \verb"const" \{
    return pair(sgn(x - o.x), sgn(y - o.y)) <=> pair(0, 0); }
  bool operator==(P o) const {
    return !sgn(x - o.x) && !sgn(y - o.y); }
D cross(P a, P b) { return a.x * b.y - a.y * b.x; }
D cross(P p, P a, P b) { return cross(a - p, b - p); }
D dot(P a, P b) { return a.x * b.x + a.y * b.y; }
D dot(P p, P a, P b) { return dot(a - p, b - p); }
D abs2(P a) { return a.x * a.x + a.v * a.v; }
D abs(P a) { return sqrt(abs2(a)); }
auto& operator<<(auto& o, P a) {</pre>
  return o << "(" << a.x << ", " << a.y << ")"; }
```

#### AngleCmp.h

**Opis:** Sortuje punkty rosnąco po kącie z przedziału  $(-\pi,\pi]$ . Punkt (0,0) ma kat 0.

```
bool angle_cmp(P a, P b) {
   auto half = [](P p) { return sgn(p.y) ?: -sgn(p.x); };
   int A = half(a), B = half(b);
   return A == B ? sgn(cross(a, b)) > 0 : A < B;
}</pre>
```

#### LineDist.h

Opis: Najkrótsza odległość między punktem i prostą/odcinkiem.

```
D line_dist(P p, P a, P b) {
    return abs(cross(p, a, b)) / abs(b - a);
}
D seg_dist(P p, P a, P b) {
    if (sgn(dot(a, p, b)) <= 0) return abs(p - a);
    if (sgn(dot(b, p, a)) <= 0) return abs(p - b);
    return line_dist(p, a, b);
}</pre>
```

LineInter.h

## LineInter ConvexHull HullTangents

```
Opis: Punkt przecięcia prostych/odcinków.
pair<int, P> line_inter(P a, P b, P c, P d) {
  auto u = cross(b - a, d - c);
  if (sgn(u) == 0) // parallel
   return {-(!sgn(cross(a, b, c))), {}};
  auto p = cross(c, b, d), q = cross(c, d, a);
  return {1, (a * p + b * q) / u};
vector<P> seg_inter(P a, P b, P c, P d) {
  auto oa = cross(c, d, a), ob = cross(c, d, b),
       oc = cross(a, b, c), od = cross(a, b, d);
  if (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn(od) < 0) // strict
   return { (a * ob - b * oa) / (ob - oa) };
  set<P> s;
  if (!sgn(seg_dist(a, c, d))) s.insert(a);
  if (!sqn(seq_dist(b, c, d))) s.insert(b);
  if (!sgn(seg_dist(c, a, b))) s.insert(c);
  if (!sgn(seq_dist(d, a, b))) s.insert(d);
  return {all(s)};
```

# 6.2 Wielokąty

## ConvexHull.h

 $\mathbf{Opis:}$  Otoczka wypukła w kierunku CCW.

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

#### HullTangents.h

**Opis:** Znajduje punkty styczne różne od a. Wielokąt musi być CCW i  $n \geq 3$ . Punkt a nie może leżeć w ściśłym wnętrzu wielokąta.

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

```
pii tangents (const vector < P > & p, P a) {
  int n = sz(p), t[2];
  rep(it, 0, 2) {
    auto dir = [&](int i) {
     P u = p[i] - a, v = p[(i + 1) % n] - a;
     D c = cross(u, v);
     if (sgn(c) != 0) return sgn(c) < 0;
     if (sgn(dot(u, v)) <= 0) return true;</pre>
      return sgn(abs2(u) - abs2(v)) > 0;
    auto idir = [&](int i) { return dir(i) ^ it; };
    if (idir(0) && !idir(n - 1)) { t[it] = 0; continue; }
    int s[2] = \{0, n - 1\};
    while (s[1] - s[0] > 2) {
     int mid = (s[0] + s[1]) / 2, x = idir(mid);
     if (idir(s[x ^ 1]) == (x ^ 1)) {
       s[x] = mid;
      } else {
       bool b = sgn(cross(p[mid] - a, p[s[1]] - a)) < 0;
```

```
s[b ^ x ^ it ^ 1] = mid;
}

t[it] = s[0] + 1 + (idir(s[0] + 1) == 0);

return {t[0], t[1]};
}
```