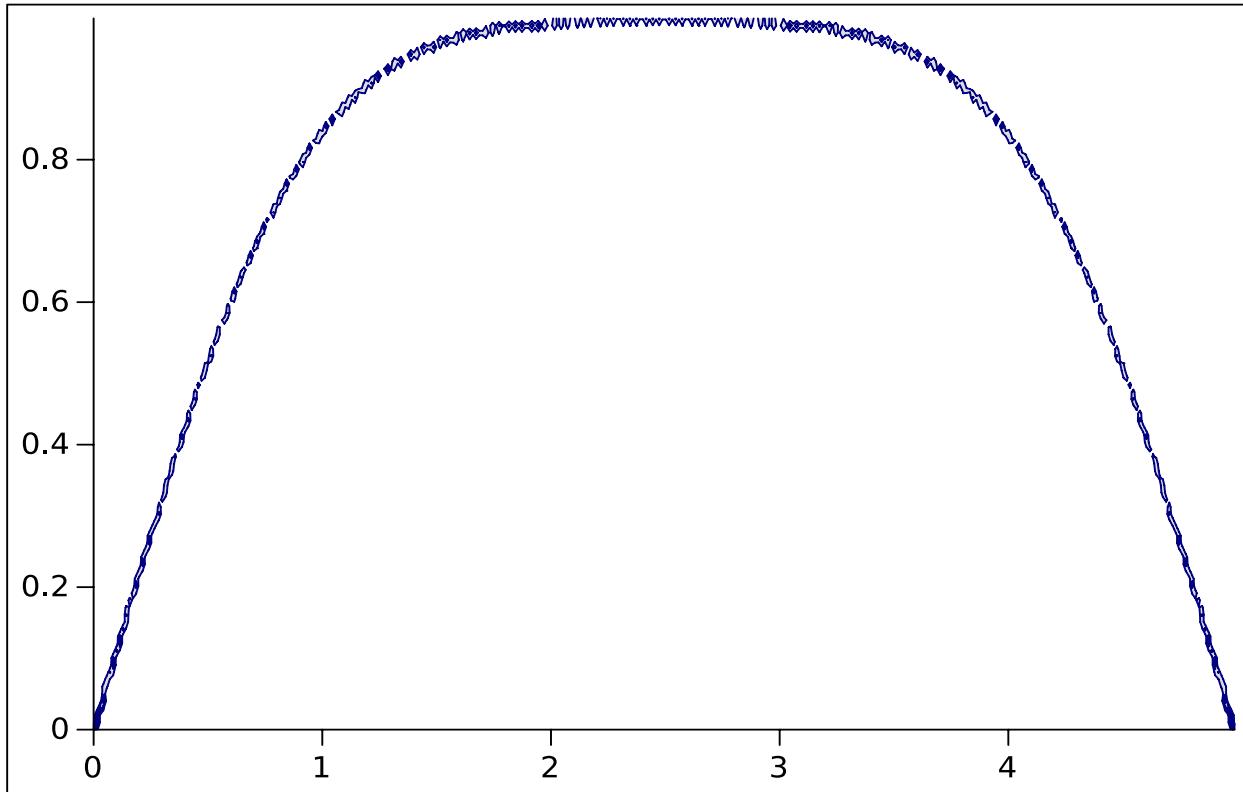


NUMERICAL Calculation of potential on a charged disk



```
#include <iostream>
#include <cmath>
#include <stdlib.h>

using namespace std;

int main() {

    const float Lx = 5.0, Ly = 5.0; // Disk diameter
    const float h = 0.01; // Grid spacing
    const int Nx = Lx / h;
    const int Ny = Ly / h;
    const int dx = Nx / 2, dy = Ny / 2; // Center of grid

    float phi1[Nx][Ny], phi2[Nx][Ny];
```

```

// Ù...Ù,Ø¬Ø§Ø±Ø¬Ù‡ÙŒ Ø§Ù^Ù„ÙŒÙ‡Ù¾Ø¤Ø§Ù+Ø³ÙŒÙ„
for (int x = 0; x < Nx; ++x) {
    for (int y = 0; y < Ny; ++ay) {
        phi1[x][y] = 1.0; // Ù...Ù,Ø¬Ø§Ø±Ø¬Ù‡ÙŒ Ù‡Ù...Ù‡Ù+Ù,Ø§Ø· Ø··Ù‡ 1
        phi2[x][y] = 1.0; // Ù...Ù,Ø¬Ø§Ø±Ø¬Ù‡ÙŒ Ù‡Ù...Ù‡Ù+Ù,Ø§Ø· Ø··Ù‡ 1
    }
}

```

// Ø§Ø¹Ù...Ø§Ù„ Ø’Ø±Ø§ÙŒØ· Ù...Ø±Ø²ÙŒ Ø¬ÙŒØ±ÙŒÙ©Ù„Ù‡ (Ù¾Ø¤Ø§Ù+Ø³ÙŒÙ„ Ø«Ø§Ø·Ø¤
ØµÙØØ± Ø¬Ø±Ù...Ø±Ø²Ù‡Ø§)

```
for (int x = 0; x < Nx; ++x) {
```

```
    phi1[x][0] = 0.0;
```

```
    phi2[x][0] = 0.0;
```

```
    phi1[x][Ny-1] = 0.0;
```

```
    phi2[x][Ny-1] = 0.0;
```

```
}
```

```
for (int y = 0; y < Ny; ++y) {
```

```
    phi1[0][y] = 0.0;
```

```
    phi2[0][y] = 0.0;
```

```
    phi1[Nx-1][y] = 0.0;
```

```
    phi2[Nx-1][y] = 0.0;
```

```
}
```

// Ø±Ù^Ø’ Ø¢Ø±Ø§Ù...â€ŒØ··Ø’ÙŒ Ø··Ø±Ø§ÙŒ Ù...ØØ§Ø³Ø··Ù‡Ù¾Ø¤Ø§Ù+Ø³ÙŒÙ„

```
for (int iter = 0; iter < 1000000; ++iter) {
```

```
    for (int x = 1; x < Nx-1; ++x) {
```

```
        for (int y = 1; y < Ny-1; ++y) {
```

```
            phi2[x][y] = 0.25 * (phi1[x+1][y] + phi1[x-1][y] + phi1[x][y+1] + phi1[x][y-1]);
```

```
    }

}

for (int x = 1; x < Nx-1; ++x) {
    for (int y = 1; y < Ny-1; ++y) {
        phi1[x][y] = phi2[x][y];
    }
}

// ØºØ®ÙŒØ±Ù‡ Ù†Ø¤Ø§ÙŒØ¬ Ø¬Ø± ÙØØ§ÙŒÙ„
ofstream out("pp15.dat");
for (int x = 0; x < Nx; ++x) {
    for (int y = 0; y < Ny; ++y) {
        out << x * h << "\t" << y * h << "\t" << phi1[x][y] << endl;
    }
}
out.close();
}
```