SPECTRAL METHOD FOR 3-DIMENSIONAL POISSON'S EQUATION IN CYLINDRICAL COORDINATES WITH REGULAR BOUNDARIES*

Yuri K. Batygin

Stanford Linear Accelerator Center, Stanford University, Stanford, CA 94309, USA

Abstract

Method of calculation of space charge field of the beam using an expansion of space charge potential and space charge distribution as Fourier-Bessel series is discussed. Coefficients of series are connected by an algebraic equation, which substantially simplifies solution of the problem. Efficiency and accuracy of the method are discussed. Suggested method is effective in multidimensional problems of study of intense chargedparticle beams.

1 INTRODUCTION

Accurate calculation of space charge potential of the beam is important for precise simulation of high current beam dynamics in particle accelerators. The most popular approaches combine finite difference methods and Fast Fourier Transforms where Poisson's equation is substituted by a finite-difference equation and resulted matrix equation is solved to find the grid function of potential [1]. In spectral methods solution of the problem is expressed as analytical functions [2], [3]. It provides more exact solution and simplifies numerical algorithm. Numerical solution in this case is a Fourier spectrum, which can be easily verified. However, spectral methods are usually slow.

2 MATHEMATICAL ALGORITHM

Consider periodic distribution of a space charge density of the beam in perfect conducting pipe of radius a (see Fig. 1). Space charge potential of the beam is calculated via solution of the Poisson's equation:

$$\frac{1}{r}\frac{\partial}{\partial r}\left(r\frac{\partial\Phi}{\partial r}\right) + \frac{1}{r^2}\frac{\partial^2\Phi}{\partial \varphi^2} + \frac{\partial^2\Phi}{\partial z^2} = -\frac{\rho(r,\varphi,z)}{\varepsilon_o} , \qquad (1)$$

with Dirichlet boundary conditions for potential Φ at the surface of conducting pipe and periodic conditions in longitudinal direction z:

$$\Phi (a, \phi, z) = 0, \qquad (2)$$

$$\Phi(\mathbf{r}, \boldsymbol{\varphi}, \mathbf{z}) = \Phi(\mathbf{r}, \boldsymbol{\varphi}, \mathbf{z} + \mathbf{L}) \quad . \tag{3}$$

Let us express unknown potential $\Phi(r, \phi, z)$ and space charge distribution $\rho(r, \phi, z)$ as Fourier-Bessel series:

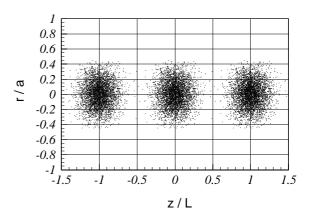


Fig. 1. Periodic distribution of charged particles.

$$\Phi = \sum_{k=-K}^{K} \sum_{n=-N}^{N} \sum_{m=1}^{M} \overline{\Phi}_{kmn} J_{n}(\nu_{nm}\frac{r}{a}) e^{-i(\frac{2\pi kz}{L} + n\phi)}, \quad (4)$$
$$\rho = \sum_{k=-K}^{K} \sum_{n=-N}^{N} \sum_{m=1}^{M} \overline{\rho}_{kmn} J_{n}(\nu_{nm}\frac{r}{a}) e^{-i(\frac{2\pi kz}{L} + n\phi)}, \quad (5)$$

where $J_n(x)$ is the Bessel function of the order n; v_{nm} is the m-th root of the equation $J_n(x) = 0$. Coefficients of space charge potential expansion in Eq. (5) are obtained from inverse Fourier-Bessel transformation:

$$\overline{\rho}_{kmn} = \frac{\int_{0}^{a} \int_{0}^{L} \int_{0}^{2\pi} \rho(r,\phi,z) J_{n}(\nu_{nm} \frac{r}{a}) e^{i(\frac{2\pi kz}{L} + n\phi)} r dr d\phi dz}{\pi a^{2} L J_{n+1}^{2}(\nu_{nm})}.$$
 (6)

Eq. (6) is obtained using orthogonality of Bessel functions

$$\int_{0}^{1} J_{m}(\mu_{i}x) J_{m}(\mu_{k}x) x \, dx = \begin{cases} 0, & i \neq k \\ \frac{1}{2} [J_{m+1}(\mu_{i})]^{2}, & i = k \end{cases}$$
(7)

After substitution of the expansions (4), (5) into Poisson's equation (1), the coefficients of potential, $\overline{\Phi}_{kmn}$, and space charge distribution, $\overline{\rho}_{kmn}$, are connected by the algebraic equation:

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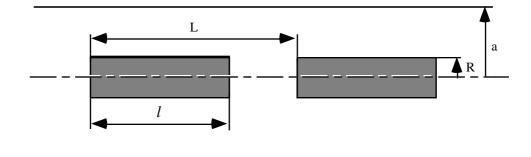


Fig. 2. Periodic train of uniformly charged cylindrical bunches.

$$\overline{\Phi}_{kmn} = \frac{\overline{\rho}_{kmn}}{\epsilon_0 [(\frac{2\pi k}{L})^2 + (\frac{\nu_{nm}}{a})^2]} .$$
(8)

Algorithm of space charge potential calculation is performed as 3 steps:

- calculation of coefficients of space charge distribution, Eq. (6),
- calculation of coefficients of space charge potential, Eq. (8),
- calculation of space charge potential, Eq. (4).

In special cases the problem can be simplified. For example, if the region is limited in longitudinal direction by conductive surfaces, it is enough to use expression:

$$\Phi = \sum_{k=1}^{K} \sum_{n=-N}^{N} \sum_{m=1}^{M} \overline{\Phi}_{kmn} J_n(\nu_{nm}\frac{r}{a}) e^{-in\phi} \sin(\frac{\pi kz}{L}),$$
⁽⁹⁾

analogously for $\rho(r, \phi, z)$.

3 CONVERGENCE OF ALGORITHM

To estimate convergence of the series (4), (5), let us calculate potential of the periodic sequence of uniformly charged bunches of radius R < a, length l < L and charge density ρ_o (see Fig. 2). For that problem the coefficients, Eq. (6), are calculated analytically:

$$\rho_{\rm kmo} = \rho_0 \frac{4}{\upsilon_{\rm om}} \left(\frac{R}{a}\right) \left(\frac{l}{L}\right) \left[\frac{J_1(\upsilon_{\rm om}\frac{R}{a})}{J_1^2(\upsilon_{\rm om})}\right] \left[\frac{\sin\left(\frac{\pi kl}{L}\right)}{\frac{(\pi kl)}{L}}\right]. \quad (10)$$

Analysis shows that the values of ρ_{kmo} decrease as ~1/(\sqrt{m} k). Substitution of Eq.(10) into Eqs. (8), (4) gives expression for potential of the train of cylindrical bunches:

$$\Phi(\mathbf{r}, \mathbf{z}) = \sum_{k=1}^{\infty} \sum_{m=1}^{\infty} \frac{4 \rho_0}{\varepsilon_0 \upsilon_{om} [(\frac{2\pi k}{L})^2 + (\frac{\upsilon_{om}}{a})^2]}$$

$$(\frac{R}{a})(\frac{l}{L}) \frac{J_1(\upsilon_{om} \frac{R}{a})}{J_1^2(\upsilon_{om})} \frac{\frac{\sin(\frac{\pi kl}{L})}{L}}{(\frac{\pi kl}{L})} J_0(\upsilon_{om} \frac{r}{a}) \cos(\frac{2\pi kz}{L}) .$$
(11)

Coefficients Φ_{kmo} decrease in this case as

$$\overline{\Phi}_{\rm kmo} \sim \frac{1}{m \, k \, (m^2 + k^2)} \, . \tag{12}$$

Example shows that the expression for potential is fast convergent. Realistic beam usually has a smooth distribution of space charge density. It is expected, that in practice the small number of coefficients in series, Eqs. (4), (5) is enough to obtain the required accuracy.

4 NUMERICAL EXAMPLE

Suggested algorithm was tested for the problem of space charge potential of the axial - symmetric beam. Calculations were done for test function

$$\Phi(\mathbf{r}, \mathbf{z}) = [1 - (\frac{\mathbf{r}}{a})^4] (\cos \frac{\pi \mathbf{z}}{L})^4 , \qquad (13)$$

which satisfies the boundary conditions, Eqs. (2), (3). The corresponding value of space charge is given by:

$$\frac{\rho(\mathbf{r}, \mathbf{z})}{\varepsilon_{o}} = 16 \frac{\mathbf{r}^{2}}{a^{4}} \left(\cos \frac{\pi \mathbf{z}}{L}\right)^{4} + \left(\frac{2\pi}{L}\right)^{2} \left[1 - \left(\frac{\mathbf{r}}{a}\right)^{4}\right] \left[\left(\cos \frac{\pi \mathbf{z}}{L}\right)^{4} - \frac{3}{4} \left(\cos \frac{2\pi \mathbf{z}}{L}\right)^{2}\right] .$$
 (14)

Cylindrical region 0 < r < a, 0 < z < L, was covered by the grid with steps $h_r = a / N_r$, $h_z = L / N_z$, where N_r , N_z are

the number of grid points in r and z, respectively. Coefficients in space charge density expansion, Eq. (5), are expressed as :

$$\overline{\rho}_{kmo} = \frac{2}{J_1^2(\nu_{om}) N_r N_z}.$$

$$\sum_{s=0}^{N_z-1} \sum_{j=1}^{N_r-1} \rho_{sj}(\frac{r_j}{a}) J_0(\nu_{mo}\frac{r_j}{a}) \exp(i\frac{2\pi k}{L} z_s) \ , \eqno(15)$$

where the grid points are used:

$$z_s = s \cdot h_z$$
, $s = 0, 1, 2, \dots, N_z - 1$, (16)

$$\mathbf{r}_{j} = \mathbf{j} \cdot \mathbf{h}_{r}, \qquad \mathbf{j} = 0, 1, 2, \dots, N_{r} - 1.$$
 (17)

For z-expansion the Fast Fourier Transform was used. The number of coefficients in z-direction was equal to the number of grid points, $K = N_z$. The number of coefficients in radial direction M was varied. An averaged deviation of numerical solution, $\Phi(r_j, z_s)$, from the exact analytical solution, $\Phi(r, z)$, was calculated as an error δ :

$$\delta = \frac{1}{N_z N_r} \sqrt{\sum_{s=0}^{N_z-1} \sum_{j=1}^{N_r-1} [\Phi(r, z) - \Phi(r_j, z_s)]^2} .$$
(18)

Results of calculation are summarized in Table 1.

For comparison, the same problem was tested by the finite-difference method. Poisson's equation (1) was substituted by the finite-difference analog:

$$U_{k,j+1}(1+\frac{1}{2(j-1)}) - 2U_{k,j}(1+\frac{h_r^2}{h_z^2}) + U_{k,j-1}(1-\frac{1}{2(j-1)}) + U_{k+1,j}(\frac{h_r}{h_z})^2 + U_{k-1,j}(\frac{h_r}{h_z})^2 = -\frac{\rho_{k,j}}{\varepsilon_0}h_r^2.$$
(19)

Calculations started with Fourier expansion of unknown potential, $U_{k, j}$, in z-direction

$$U_{k,j} = \sum_{m=1}^{N_z} \overline{U}_m(j) \exp(-i \frac{2\pi (k-1) (m-1)}{N_z}) , \qquad (20)$$

Table 1. Error δ of the test problem.

	Spectral method			Finite -difference
Grid	M=5	M=10	M=40	method
16 x 16	$3 \cdot 10^{-4}$	3.10-4	3.10-	⁴ 3·10 ⁻⁴
32 x 32	$1 \cdot 10^{-4}$	4·10 ⁻⁵	3.10-5	⁵ 4·10 ⁻⁵
64 x 64	6·10 ⁻⁵	1.10^{-5}	4·10 ⁻⁶	6·10 ⁻⁶
128 x 128	3.10-5	5.10-6	5·10 ⁻⁷	1·10 ⁻⁶
256 x 256	1·10 ⁻⁵	2.10-6	i 1·10 ^{−7}	1.10-6

similar for space charge density $\rho_{k,j}$. Coefficients of Fourier expansion, $\overline{U}_m(j)$, were expressed via inverse Fourier transform:

$$\overline{U}_{m}(j) = \frac{1}{N_{z}} \sum_{k=1}^{N_{z}} U_{kj} \exp(i \frac{2\pi (k-1) (m-1)}{N_{z}})$$
, (21)

and similar for $\overline{\rho}_m(j)$. Substitution of expansion, Eq. (21), into the finite-difference analog of the Poisson's equation, Eq. (19), results in a three-diagonal matrix equation:

$$\alpha_{j} \overline{U}_{m}(j+1) + \beta_{j} \overline{U}_{m}(j) + \gamma_{j} \overline{U}_{m}(j-1) = w_{j} , \qquad (22)$$

which was solved utilizing the Gauss elimination method [1]. After that, the potential in grid points was calculated using Fourier series, Eq. (20).

Comparison of numerical results shows that spectral method gives better accuracy than the finite-difference one using relatively small number of harmonics M in radial direction.

5 REFERENCES

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