Computation of quasi-discrete Hankel transforms of integer order for propagating optical wave fields

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The method originally proposed by Yu *et al.* [Opt. Lett. **23**, 409 (1998)] for evaluating the zero-order Hankel transform is generalized to high-order Hankel transforms. Since the method preserves the discrete form of the Parseval theorem, it is particularly suitable for field propagation. A general algorithm for propagating an input field through axially symmetric systems using the generalized method is given. The advantages and the disadvantages of the method with respect to other typical methods are discussed. © 2004 Optical Society of America

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1. INTRODUCTION

In propagation of optical beams through systems with cylindrical symmetry, one encounters the need for the numerical computation of the pth-order Hankel transform (HT) and the inverse Hankel Transform (IHT):

$$f_2(\nu) = 2\pi \int_0^\infty f_1(r) J_p(2\pi\nu r) r dr,$$
 (1a)

$$f_1(r) = 2\pi \int_0^\infty f_2(\nu) J_p(2\pi r\nu) \nu \mathrm{d}\nu, \qquad (1\mathrm{b})$$

where J_p is the *p*th-order Bessel function of the first kind, r is the radial coordinate, and ν is the spatial frequency. Integrals (1) are especially difficult to compute because of the oscillatory behavior of the Bessel function and the infinite length of the interval. Since the seminal work by Siegman in 1977,¹ a number of algorithms for the numerical evaluation of the HT have been reported in the literature, for both zero-order HTs²⁻¹¹ and high-order HTs.¹²⁻²¹ Unfortunately, the performance of a method for computing the HT is highly dependent on the function to be transformed, and thus it is difficult to determine the optimal algorithm for a given function. In optics, we often deal with problems where the HT and the IHT need to be computed thousands of times starting from a known analytical expression for the input function. In this case, an efficient and fast numerical HT algorithm is needed. A useful numerical evaluation of methods for calculating HTs was provided recently by Markham and Conchello.¹¹

In 1998, Yu *et al.*¹⁰ developed a zero-order quasidiscrete HT method that approximates the input function by a Fourier–Bessel series over a finite integration interval. This led to a symmetric transformation matrix for the HT and the IHT that satisfies the discrete form of the Parseval theorem. Since energy conservation holds after successive applications of the HT and the IHT, this method has considerable application in iterative processes. The Yu method was not included by Markham and Conchello in their analysis¹¹; however, we have found tangible advantages of the Yu method over other typical methods.

The purpose of this work is twofold. First, we present an algorithm that extends the zero-order HT method of Yu^{10} to higher orders. We will denote this generalized method as the *p*th-order quasi-discrete HT (pQDHT). Second, we show the application of the pQDHT method to the field propagation through axially symmetric systems. The advantages and the disadvantages of the pQDHT method with respect to other common methods are discussed.

2. DERIVATION OF THE METHOD

To derive the algorithm, we first assume that the space domain $f_1(r)$ extends over a limited region $0 \le r \le R$. This follows from the fact that the functions f_1 representing physical fields are actually zero outside a disk whose radius is conveniently chosen equal to R. Thus, from a physical point of view, this assumption is reasonable; from a computational point of view, it is necessary. Since the HT and the IHT have indeed the same mathematical form [see Eqs. (1)], it is reasonable to assume that the frequency domain $f_2(\nu)$ is distributed over a limited band 0 $\leq \nu \leq V$ as well. Physically, this means that the spacebandwidth product RV of the transform remains finite, as required by numerical limitations. As pointed out by Siegman,¹ the evaluation of Eqs. (1) in a finite range instead of the actual limits $(0, \infty)$ is an inherent source of error in the computation of the HT and the IHT when numerical techniques are used. Therefore we have that

$$f_1(r \ge R) = 0, \qquad f_2(\nu \ge V) = 0, \qquad (2)$$

where R and V are the truncated radii of the spatial variable and the frequency variable, respectively.

The functions $f_1(r)$ and $f_2(\nu)$ may be expanded in terms of the *p*th-order Bessel series,²²

$$f(r) = \sum_{m=1}^{\infty} c_{pm} J_p \left(\alpha_{pm} \frac{r}{a} \right), \qquad 0 \le r \le a, \qquad (3)$$

where α_{pm} is the *m*th root of the *p*th Bessel function and the coefficients c_{pm} are determined with

$$c_{pm} = \frac{1}{a^2 J_{p+1}^2(a_{pm})} \int_0^a f(r) J_p \left(\alpha_{pm} \frac{r}{a} \right) r \mathrm{d}r.$$
(4)

If we evaluate the radius r at values $\alpha_{pn}/(2\pi V)$ and the frequency ν at $\alpha_{pm}/(2\pi R)$ and replace ∞ with a finite number N, Eqs. (1) are approximated to some level of accuracy by the discrete sums

$$f_{2}\left(\frac{\alpha_{pm}}{2\pi R}\right) = \frac{1}{\pi V^{2}} \sum_{n=1}^{N} \frac{f_{1}(\alpha_{pn}/(2\pi V))}{J_{p+1}^{2}(\alpha_{pn})} J_{p}\left(\frac{\alpha_{pn}\alpha_{pm}}{S}\right),$$
(5a)

$$f_1\left(\frac{\alpha_{pn}}{2\pi V}\right) = \frac{1}{\pi R^2} \sum_{m=1}^{N} \frac{f_2(\alpha_{pm}/(2\pi R))}{J_{p+1}^2(\alpha_{pm})} J_p\left(\frac{\alpha_{pn}\alpha_{pm}}{S}\right),$$
(5b)

where $S \equiv 2 \pi R V$. Equations (5) can be rewritten in a symmetric form by defining the vectors

$$F_{2}(m) = f_{2}\left(\frac{\alpha_{pm}}{2\pi R}\right) |J_{p+1}(\alpha_{pm})|^{-1}V,$$
 (6a)

$$F_{1}(n) = f_{1}\left(\frac{\alpha_{pn}}{2\pi V}\right) |J_{p+1}(\alpha_{pn})|^{-1}R;$$
 (6b)

thus Eqs. (5) reduce to

$$F_2(m) = \sum_{n=1}^{N} T_{mn} F_1(n), \qquad (7a)$$

$$F_1(n) = \sum_{m=1}^{N} T_{nm} F_2(m),$$
 (7b)

where

$$T_{mn} = \frac{2J_p(\alpha_{pn}\alpha_{pm}/S)}{|J_{p+1}(\alpha_{pn})||J_{p+1}(\alpha_{pm})|S}$$
(8)

defines the elements of an $N \times N$ transformation matrix **T**. From Eqs. (2) and (6), it is evident that $F_1(n > N) = F_2(m > N) = 0$ for $S \ge \alpha_{p,N+1}$; thus only the first $N \times N$ elements in the matrix **T** are relevant.

3.7

There are some mathematical properties of the transformation matrix \mathbf{T} to be discussed here. First, note that \mathbf{T} is a real square symmetric matrix, and thus \mathbf{T} is equal to its transpose \mathbf{T}^{T} . To preserve self-consistency, the matrix \mathbf{T} needs to be unitary, which means that an input function returns to itself after one HT and one IHT, i.e., $\mathbf{TT} = \mathbf{I}_N$, where \mathbf{I}_N is the identity matrix. Now, we see in Eq. (8) that \mathbf{T} is, in fact, a function of the continuous parameter S; thus it is still necessary to determine the optimum S that satisfies the equation $\mathbf{T}^2(S) - \mathbf{I}_N = 0$. In this respect, Yu *et al.*¹⁰ found numerically for the zero-order HT that if $S = \alpha_{0,N+1}$, then \mathbf{T} is very close to being a unitary matrix and gets closer as N increases. We have corroborated that this result holds for the *p*th-order HTs, i.e., $S = \alpha_{p,N+1}$.

Since the determinant of unitary matrices is indeed unity, it is possible to use this result to verify the accuracy of the matrix $\mathbf{T}(\alpha_{p,N+1})$ for a given N. By using numerical evaluation, we have verified that $|\det[\mathbf{T}(\alpha_{p,N+1})]$ $-1| < 10^{-8}$, 10^{-9} , and 10^{-11} for N = 50, 200, and 500, respectively. These tolerances are enough for practical purposes and can be used as criteria to orient the reader in the number of sampling points N. If a more accurate matrix **T** is desired, it could be improved (once N is defined) by finding the root of the nonlinear equation $\det[\mathbf{T}(S)] - 1 = 0$ in the neighborhood of $S = \alpha_{p,N+1}$. Note that because **T** is a real unitary matrix, its rows (and columns) form a set of orthonormal vectors that constitute a basis for the radial and frequency N-dimensional spaces.²²

An important feature of the pQDHT method is that it is energy preserving, i.e., satisfies the Parseval theorem:

$$\int_{0}^{\infty} |f_{1}(r)|^{2} r \mathrm{d}r = \int_{0}^{\infty} |f_{2}(\nu)|^{2} \nu \mathrm{d}\nu.$$
(9)

To prove this, we substitute Eqs. (5) into the last equation, and after some algebra we arrive at

$$\sum_{m=1}^{N} \frac{|f_2(\alpha_{pm}/(2\pi R))|^2}{2\pi^2 R^2 J_{p+1}^2(\alpha_{pm})} = \sum_{n=1}^{N} \frac{|f_1(\alpha_{pn}/(2\pi V))|^2}{2\pi^2 V^2 J_{p+1}^2(\alpha_{pn})},$$
(10)

which can be rewritten in the discrete form of the Parseval theorem by using Eqs. (6), namely,

$$\frac{2}{S^2} \sum_{n=1}^{N} |F_1(n)|^2 = \frac{2}{S^2} \sum_{m=1}^{N} |F_2(m)|^2.$$
(11)

The conservation of energy can also be inferred from the fact that the norm of an input vector F_1 that is linearly mapped onto an output vector F_2 is invariant when the transformation matrix **T** is unitary.

3. NUMERICAL TESTS

To test the numerical accuracy of the pQDHT method, we consider two examples. All codes were written in Matlab software, mainly because of the existence of a wealth of built-in mathematical functions.

For the first example, we choose the sinc function $f_1(r) = \sin(2\pi\gamma r)/2\pi\gamma r$, whose analytical *p*th-order HT is given by

$$f_{2}(\nu) = \frac{\nu^{p} \cos(p \pi/2)}{2 \pi \gamma \sqrt{\gamma^{2} - v^{2}} (\gamma + \sqrt{\gamma^{2} - v^{2}})^{p}}, \qquad 0 \le \nu < \gamma$$
(12a)

$$=\frac{\sin[p \arcsin(\gamma/\nu)]}{2\pi\gamma\sqrt{\nu^2-\gamma^2}}, \quad \nu > \gamma.$$
(12b)

This function was also considered in Ref. 14 and provides a useful benchmark for our routines. Figures 1(a) and 1(c) depict the analytical and computed HTs of the sinc function with $\gamma = 5$ for orders p = 1 and 4. We used N = 256 points to sample the radius in the interval (0, 3). We can appreciate that the calculated values are quite close to the exact values. Since we must truncate

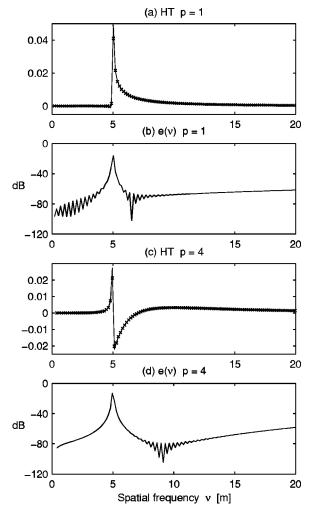


Fig. 1. Exact and computed Hankel transforms of the sinc function with $\gamma = 5$ for (a) p = 1 and (c) p = 4. Solid curves denote exact transforms, and crosses denote the calculated transforms. Dynamic errors are shown in (b) and (d).

the input function, the calculated transform suffers from the Gibbs phenomenon, particularly near the discontinuity; however, it can be reduced with proper windowing (apodizing). Now, a suitable measure of accuracy of the numerical method is provided by the dynamic error

$$e(\nu) = 20 \log_{10} \left[\frac{|f_2(\nu) - f_2^*(\nu)|}{\max|f_2^*(\nu)|} \right].$$
 (13)

This error function compares the difference between exact $[f_2(\nu)]$ and estimated $[f_2^*(\nu)]$ transform samples with the maximum value of the HT. The dynamic error is given in decibels and plotted on a [0, -120 dB] scale in subplots 1(b) and 1(d). The error is greatest around $\nu = \gamma$, i.e., near the point where the transform exhibits the discontinuity. Beyond this point, the error remains below -60 dB.

Successive application of the HT and the IHT to the input function f_1 will presumably return the same function. We quantify the approximation error in the numerical computation by taking the average absolute difference between the exact $[f_1(r)]$ and retrieved $[f_1^*(r)]$ functions:

$$\epsilon_1 = \frac{1}{N} \sum_{j=1}^{N} |f_1 - f_1^*|. \tag{14}$$

For the fourth-order HT and IHT of the sinc function, we obtain $\epsilon_1 \sim 10^{-10} (N = 100)$, $\epsilon_1 \sim 10^{-12} (N = 200)$, and $\epsilon_1 \sim 10^{-14} (N \ge 300)$ over the whole range of the radial variable.

In the second example, we compare numerically the pQDHT with respect to the well-assessed procedure of the quasi-fast Hankel transform (QFHT) introduced by Siegman.¹ To this end, we choose as test function the generalized version of the top-hat function, namely, $f_1(r) = r^p[H(r) - H(r - a)]$ with a > 0, where H(r) is the step function H(r) = 1 for $r \ge 0$ and H(r) = 0 elsewhere, whose analytical *p*th-order HT is given by

$$f_2(\nu) = a^{p+1} \frac{J_{p+1}(2\pi a\,\nu)}{\nu}.$$
(15)

For the calculations, we take p = 4 and a = 1 and sample the function $f_1(r)$ in the range $r \in (0, 2)$ for N = 512 and 1024. The pQDHT requires the selection of two free parameters (R and N), whereas the QFHT needs the selection of four free parameters (typically, R, N, V, and $K_2 =$ minimum allowed number of "points per cycle" in the frequency space). Since the QFHT has an inherent problem of lower-end correction, it requires a careful selection of the parameters¹; thus, to minimize the missing contribution due to this problem, we have chosen V = 20 and $K_2 = 2$.

The errors and the CPU times (two HT calculations) for the pQDHT and the QFHT are compared in Table 1. The error ϵ_2 corresponds to the average absolute difference between the exact $[f_2(\nu)]$ and computed $[f_2^*(\nu)]$ transformed functions, and the error ϵ_1 is computed with Eq. (14). For the retrieved function $f_1^*(r)$, the results obtained with the pQDFT are several orders of magnitude more precise than those produced by the standard QFHT. Conversely, the QFHT is faster, but it has lower accuracy for the calculated points. This circumstance can be explained by considering that one of the main advantages of the QFHT and of similar algorithms based on the fast Fourier transform (FFT) is that they require the storage of the kernel in the form of a vector instead of a square matrix. The number of complex multiplications for a QFHT is $4N \log_2 2N + 2N$ (see Refs. 2 and 7), and the number of real multiplications for a pQDHT is N^2 . Because one complex multiplication is almost equal to three real multiplications, it is obvious that the pQDHT is more efficient than the QFHT for the same accuracy.

Table 1. Comparison of the pQDHT and the QFTT

N	ϵ_2	ϵ_1	CPU Time (s)	V
pQDHT				
512	$1.3~ imes~10^{-3}$	$2.2~ imes~10^{-13}$	0.028	128.7
1024	$4.8~ imes~10^{-5}$	$2.7~ imes~10^{-14}$	0.071	256.7
QFHT				
512	$3.8~ imes~10^{-3}$	$1.3~ imes~10^{-2}$	0.015	20
1024	$1.9~ imes~10^{-4}$	$6.4~ imes~10^{-3}$	0.025	20

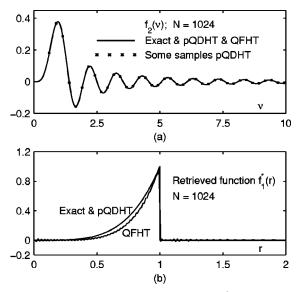


Fig. 2. (a) Analytical $[f_2(\nu)]$ and computed $[f_2^*(\nu)]$ transforms, (b) exact and retrieved functions $f_1^*(r)$. Both plots correspond to N = 1024 points sampled according to the criteria of the pQDHT and the QFHT.

In Fig. 2(a), we show the analytical $[f_2(\nu)]$ and computed $[f_2^*(\nu)]$ transforms in the interval $\nu \in [0, 10]$. The behavior of the retrieved functions $f_1^*(r)$ is depicted in Fig. 2(b). Note that the plot of $f_1^*(r)$ calculated with the pQDHT coincides with the exact function $f_1(r)$. Conversely, the observed Gibbs-phenomenon ripples in $f_1^*(r)$ computed with the QFHT result from the lower finite truncation in the transform domain (V = 20).

4. WAVE PROPAGATION WITH USE OF THE *p*th-ORDER QUASI-DISCRETE HANKEL TRANSFORM

In this section, we apply the pQDHT method to demonstrate its effectiveness for studying the propagation through axially symmetric systems. The general algorithm for propagating an input field $u_0(r)$ with the pQDHT method is summarized below. For clarity, overbarred variables denote column vectors. The operators \otimes and \oslash denote element-by-element multiplication and division of vectors, respectively, and * denotes the usual matrix product.

1. Choose the input parameters: p (order), R (maximum physical radius), and N (number of points).

2. Read and store the vector of roots of the *p*th-order Bessel function $\bar{\alpha} = [\alpha_1, ..., \alpha_N]$ and the (N + 1)th root α_{N+1} . See Refs. 23 and 24.

3. Calculate the radius vector $\bar{r} = \bar{\alpha}R/\alpha_{N+1}$, the radial frequency vector $\bar{v} = \bar{\alpha}/(2\pi R)$, and the limiting frequency $V = \alpha_{N+1}/(2\pi R)$.

4. Generate the symmetric $N \times N$ transformation matrix **T** [Eq. (8)] and the column vector $\overline{J} = |J_{p+1}(\overline{\alpha})|/R$.

5. Choose the propagating parameters: z_{max} (propagation distance), nz (number of steps), $\Delta z = z_{\text{max}}/(nz)$ (increment), and λ (wavelength).

6. Calculate the propagator²⁵

$$\overline{\text{prop}} = \exp(i2\pi\Delta z \sqrt{\lambda^{-2} - \bar{\nu}^2}).$$
(16)

- 7. Evaluate the input field $\bar{u}_0(r)$.
- 8. Calculate the HT of the input field: \overline{H} = $\mathbf{T} * [\overline{u}_0(r) \oslash \overline{J}].$
- 9. Propagating loop: for j = 1,...,nz; Propagation in Fourier space:

$$\bar{H} = \bar{H} \otimes \operatorname{prop}$$

Calculate and store the spatial field at this step:

$$\bar{u}_i = \mathbf{T} * (\bar{H} \oslash \bar{J})$$

end

Note that the matrix \mathbf{T} is computed only once when the algorithm is applied to propagation and the HT is reduced to a matrix-vector multiplication.

To see how the pQDHT method works in practice, we consider the focusing evolution of a trucated fourth-order Bessel beam $u(\mathbf{r}) = J_4(k_t r) \exp(i4\phi + ik_z z)$. The beam is normally incident on a circular lens with radius R = 4 mm and focal distance f = 0.5 m that is located at plane z = 0. The transverse and longitudinal wave numbers satisfy the relation $k^2 = (2 \pi / \lambda)^2 = k_t^2 + k_z^2$. For *p*th-order Bessel beams, it is known that the image in

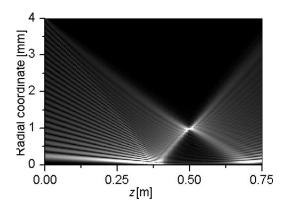


Fig. 3. Focusing evolution of a fourth-order Bessel beam along the plane (r, z). Physical parameters are chosen to produce a circular ring of radius 1 mm in the focal plane (at z = 0.5 m).

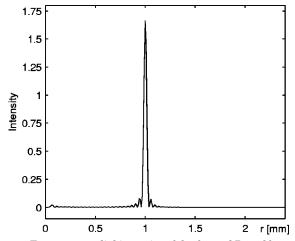


Fig. 4. Transverse radial intensity of the focused Bessel beam at the focal plane.

the focal plane corresponds to a circular ring of radius ρ = fk_t/k_z modulated by the angular spectrum $A(\phi)$ $\exp(ip\phi)$.²⁶ For an illumination at $\lambda = 632.8$ nm, we have chosen $k_t = 19858.32 \text{ m}^{-1}$ to ensure an image ring with radius $\rho = 1$ mm. The paraxial transmittance of the lens is given by $\exp[-ikr^2/(2f)]$. The propagation of the focused Bessel beam along the plane r-z is displayed in Fig. 3. This spatial evolution was obtained by sampling the radius at 256 points and calculating the field at 300 transverse planes evenly spaced through a length of 75 cm. All numerical calculations were completed within 7.11 s on a 1-GHz personal computer with 256-Mbyte RAM. In Fig. 4, we show the radial intensity of the focusing Bessel beam at z = 0.5 m. As expected, the plot exhibits an intense peak corresponding to a circular ring at the focal plane. The good agreement between these results and results reported in Ref. 26 (obtained with twodimensional FFT-based algorithms) shows that the pQDHT method is highly accurate, and it can thus be used for studying the wave propagation through axially symmetric systems.

5. CONCLUSIONS

In conclusion, we have proposed a discrete algorithm for computing general integer-order HTs and IHTs for a given input function. Our algorithm is a generalization of the original zero-order HT method proposed recently by Yu *et al.*¹⁰ In tests done with known transform pairs, agreement between exact and computed transforms was excellent. The main advantages and disadvantages of the pQDHT method are summarized as follows:

• The pQDHT method does not need either to interpolate the set of data points or to use very large zero padding, as is usual in the popular backprojection and sliceprojection methods (BP-SP)^{5,8,12–15} and methods based on the one-dimensional FFT through an exponential change of variables (FT-EXP).^{1–3,19,20} In this way, the accuracy of the method depends exclusively on the size of the transformation matrix **T** and is not affected by the selection of additional parameters or by the choice of interpolation method, quadrature scheme, or resampling procedure.

• By construction, the pQDHT method is energy preserving; i.e., the original field is retrieved after two successive applications of the HT. Other known methods (e.g., Refs. 1 and 8) have no simple retrieval expression. This property makes the pQDHT method very useful, particularly when it is applied to propagation.

• The pQDHT method transforms complex inputs in one step, contrary to the BP-SP and FT-EXP methods, which typically deal with the real and imaginary contributions separately (e.g., Refs. 4, 5, and 8).

• The computation of the transformation matrix \mathbf{T} may be time-consuming; however, tables and efficient algorithms for computing the zeros of the Bessel functions can be found in the numerical literature (see, e.g., Refs. 23 and 24). A text file with the first 3000 roots of the first ten orders of the Bessel *J* function is available by writing to the corresponding author. Moreover, the matrix \mathbf{T} needs to be computed only once in propagation applications.

We have shown how to apply the pQDHT method in optical field propagation. Since the HT and the IHT are reduced to a simple matrix-vector multiplication, the algorithm is very fast and efficient. The advantages outlined above should make the pQDHT method valuable in optical beam propagation as well as other optical and nonoptical applications.

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