

1. GENERAL RELATIONSHIPS AND TECHNIQUES

The extra gain with respect to the multidimensional Cooley–Tukey method is that *there are no twiddle factors between p -primary pieces corresponding to different primes p .*

The case where \mathbf{N} is not diagonal has been examined by Guessoum & Mersereau (1986).

1.3.3.3.2.3. Nesting of Winograd small FFTs

Suppose that the CRT has been used as above to map an n -dimensional DFT to a μ -dimensional DFT. For each $\kappa = 1, \dots, \mu$ [κ runs over those pairs (i, j) such that $v(i, j) > 0$], the Rader/Winograd procedure may be applied to put the matrix of the κ th 1D DFT in the **CBA** normal form of a Winograd small FFT. The full DFT matrix may then be written, up to permutation of data and results, as

$$\bigotimes_{\kappa=1}^{\mu} (\mathbf{C}_{\kappa} \mathbf{B}_{\kappa} \mathbf{A}_{\kappa}).$$

A well known property of the tensor product of matrices allows this to be rewritten as

$$\left(\bigotimes_{\gamma=1}^{\mu} \mathbf{C}_{\gamma} \right) \times \left(\bigotimes_{\beta=1}^{\mu} \mathbf{B}_{\beta} \right) \times \left(\bigotimes_{\alpha=1}^{\mu} \mathbf{A}_{\alpha} \right)$$

and thus to form a matrix in which the *combined* pre-addition, multiplication and post-addition matrices have been *precomputed*. This procedure, called *nesting*, can be shown to afford a reduction of the arithmetic operation count compared to the row–column method (Morris, 1978).

Clearly, the nesting rearrangement need not be applied to all μ dimensions, but can be restricted to any desired subset of them.

1.3.3.3.2.4. The Nussbaumer–Quandalle algorithm

Nussbaumer’s approach views the DFT as the evaluation of certain polynomials constructed from the data (as in Section 1.3.3.2.4). For instance, putting $\omega = e(1/N)$, the 1D N -point DFT

$$X^*(k^*) = \sum_{k=0}^{N-1} X(k) \omega^{k^*k}$$

may be written

$$X^*(k^*) = Q(\omega^{k^*}),$$

where the polynomial Q is defined by

$$Q(z) = \sum_{k=0}^{N-1} X(k) z^k.$$

Let us consider (Nussbaumer & Quandalle, 1979) a 2D transform of size $N \times N$:

$$X^*(k_1^*, k_2^*) = \sum_{k_1=0}^{N-1} \sum_{k_2=0}^{N-1} X(k_1, k_2) \omega^{k_1^*k_1 + k_2^*k_2}.$$

By introduction of the polynomials

$$Q_{k_2}(z) = \sum_{k_1} X(k_1, k_2) z^{k_1}$$

$$R_{k_2^*}(z) = \sum_{k_2} \omega^{k_2^*k_2} Q_{k_2}(z),$$

this may be rewritten:

$$X^*(k_1^*, k_2^*) = R_{k_2^*}(\omega^{k_1^*}) = \sum_{k_2} \omega^{k_2^*k_2} Q_{k_2}(\omega^{k_1^*}).$$

Let us now suppose that k_1^* is coprime to N . Then k_1^* has a unique inverse modulo N (denoted by $1/k_1^*$), so that multiplication by k_1^* simply permutes the elements of $\mathbb{Z}/N\mathbb{Z}$ and hence

$$\sum_{k_2=0}^{N-1} f(k_2) = \sum_{k_2=0}^{N-1} f(k_1^*k_2)$$

for any function f over $\mathbb{Z}/N\mathbb{Z}$. We may thus write:

$$X^*(k_1^*, k_2^*) = \sum_{k_2} \omega^{k_1^*k_2^*k_2} Q_{k_1^*k_2}(\omega^{k_1^*})$$

$$= S_{k_1^*k_2}(\omega^{k_1^*})$$

where

$$S_{k^*}(z) = \sum_{k_2} z^{k^*k_2} Q_{k_2}(z).$$

Since only the value of polynomial $S_{k^*}(z)$ at $z = \omega^{k_1^*}$ is involved in the result, the computation of S_{k^*} may be carried out modulo the unique cyclotomic polynomial $P(z)$ such that $P(\omega^{k_1^*}) = 0$. Thus, if we define:

$$T_{k^*}(z) = \sum_{k_2} z^{k^*k_2} Q_{k_2}(z) \text{ mod } P(z)$$

we may write:

$$X^*(k_1^*, k_2^*) = T_{k_1^*k_2^*}(\omega^{k_1^*})$$

or equivalently

$$X^* \left(k_1^*, \frac{k_2^*}{k_1^*} \right) = T_{k_2^*}(\omega^{k_1^*}).$$

For N an odd prime p , all nonzero values of k_1^* are coprime with p so that the $p \times p$ -point DFT may be calculated as follows: (1) form the polynomials

$$T_{k_2^*}(z) = \sum_{k_1} \sum_{k_2} X(k_1, k_2) z^{k_1 + k_2^*k_2} \text{ mod } P(z)$$

- for $k_2^* = 0, \dots, p - 1$;
- (2) evaluate $T_{k_2^*}(\omega^{k_1^*})$ for $k_1^* = 0, \dots, p - 1$;
- (3) put $X^*(k_1^*, k_2^*/k_1^*) = T_{k_2^*}(\omega^{k_1^*})$;
- (4) calculate the terms for $k_1^* = 0$ separately by