The Two-Scale Approximation for Nonlinear Wave-Wave Interactions

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Why did we switch from 2nd Generation spectral models to 3rd Generation spectral models?

To allow detailed balance calculations of source terms in order to be able to treat complex situations.

The nonlinear interaction source term is critical to this balance.

If this term is incorrect, all the others must be tuned to compensate. This is not physics --- this is tuning.

Objective of this paper

Examine the accuracy of the DIA and its relationship to the full integral.

Introduce a new approximation method for nonlinear interaction source term with improved accuracy, while retaining "similar" computational efficiency



Coastal wave spectra can be very complex and contain directional shear! As can wave spectra in hurricanes/turning winds/slanting fetch/limited fetch

Even "simple" spectra can have relatively complex shapes!



Currituck Sound Mean Directional Distribution Function all long-fetch cases (327), u₁₀ > 7 m/s centered on wind direction, grouped by f/f_p



FIG. 10. The lobe angle (a, c) and the lobe ratio (b, d) of the bimodal distribution. In (a) and (b), the directional resolution is degraded to have a uniform resolution as done for the Fourier decomposition procedure. Solid curves: from polynomial fitting (coefficients listed in Table 2), and dashed curves: computed by $D_{kFFTA}(\theta)$. In (c) and (d), the directional resolution is not degraded. The dashed-and-dotted curves are computed from Eqs. (20) and (21). Numerical results of Banner and Young (1994) on the effect of dissipation functions are shown with stars: quadratic, triangles: cubic, and square: quartic frequency dependence. Quasi-steady wave field.





<u>General 4-wave interaction integral for rate of change of action density at k_1 </u> Involves action densities at 4 wavenumbers, the coupling coefficient (C) which is also a function of the wavenumbers, and a Jacobian tranform for moving from <u>k</u> to (s,n) space

$$\frac{\partial n(\underline{k}_1)}{\partial t} = \iint T(\underline{k}_1, \underline{k}_3) d\underline{k}_3$$

$$T(\underline{k}_1, \underline{k}_3) = \iint [n_1 n_3 (n_4 - n_2) + n_2 n_4 (n_3 - n_1)] C(\underline{k}_1, \underline{k}_2, \underline{k}_3, \underline{k}_4) \left| \frac{\partial W}{\partial \underline{n}} \right|^{-1} d\underline{s}$$

The integral for 4-wave interactions can be written as an integral of the transfers from one wavenumber into another, where 2 other wave numbers are free to assumed continuous values along loci determined by the resonant conditions:

$$\underline{k}_{4} = \underline{k}_{1} + \underline{k}_{2} - \underline{k}_{3} \qquad \qquad \vec{k}_{1} - \vec{k}_{3} = \vec{k}_{4} - \vec{k}_{2}$$

$$\underline{k}_{2} = \underline{k}_{2} \left(s, \underline{k}_{1}, \underline{k}_{3} \right) \longrightarrow \text{DIA} \qquad \qquad as$$

$$Constraint \qquad \qquad \vec{k}_{2} \rightarrow \vec{k}_{1}$$

$$W = \omega_{1} + \omega_{2} - \omega_{3} - \omega_{4}$$

DIA

Where s is determined

from W=0 solution

Represents subset of points for which two wavenumbers are (close to) equal. Essentially gives the interactions along Phillips' "figure-8" curve.

How are these two sets of interactions related?



There is one point along each s-locus for which k2=k1, up to the limit of the maximum k3 value in the Figure 8 resonance pattern. The Full Integral considers contributions over the 2-dimensions above. The DIA only samples along the line.









What happens if you violate this limit?



Comparison of a DIA calculation and an actual full-integral, finite-depth calculation for the case of kph=0.7 (JONSWAP spectrum with a peak period of 10 seconds in depth of 10.5 meters).

Since the DIA only samples a slice from the complete integral, and is tuned to fit a specific spectrum, how well does it work for a range of peakedness typical of wave generation conditions?



As the spectrum approaches full development 3 the DIA progressively overpredicts the amount of energy transferred to the forward face of the spectrum - requiring other terms in the detailed 2.5 balance to try to compensate for this transfer. 2 ^{1.5} 1 0.5 0 5 2 3 4 6 7 0 8 Y The ratio of the maximum value within the positive lobe predicted by the DIA to the maximum value predicted by the full integral,

as a function of JONSWAP spectral peakedness parameter, y.



Same as previous figure, defining a ratio for the largest negative values within the negative lobe, as a function of JONSWAP spectral peakedness parameter, γ .

Problem with DIA – The basis is not the integral that we are trying to estimate.

We need a new approximation that;

- conserves constants of motion (action, energy, momentum)
- has its basis in the correct integral (important for complex cases)
- retains the number of degrees of freedom in the modeled spectrum
- is not limited to $k_p h \ge 1$
- is much more efficient than the full integral

Basis for Two-Scale Approximation

$$N^{3} = \hat{n}_{1}\hat{n}_{3}(\hat{n}_{4} - \hat{n}_{2}) + \hat{n}_{2}\hat{n}_{4}(\hat{n}_{3} - \hat{n}_{1}) + n'_{1}n'_{3}(n'_{4} - n'_{2}) + n'_{2}n'_{4}(n'_{3} - n'_{1}) + \hat{n}_{1}\hat{n}_{3}(n'_{4} - n'_{2}) + \hat{n}_{2}\hat{n}_{4}(n'_{3} - n'_{1}) + n'_{1}n'_{3}(\hat{n}_{4} - \hat{n}_{2}) + n'_{2}n'_{4}(\hat{n}_{3} - \hat{n}_{1}) + \hat{n}_{1}n'_{3}(\hat{n}_{4} - \hat{n}_{2}) + \hat{n}_{2}n'_{4}(\hat{n}_{3} - \hat{n}_{1}) + n'_{1}\hat{n}_{3}(\hat{n}_{4} - \hat{n}_{2}) + n'_{2}\hat{n}_{4}(\hat{n}_{3} - \hat{n}_{1}) + \hat{n}_{1}n'_{3}(n'_{4} - n'_{2}) + n'_{2}\hat{n}_{4}(\hat{n}_{3} - n'_{1}) + \hat{n}_{1}n'_{3}(n'_{4} - n'_{2}) + \hat{n}_{2}n'_{4}(n'_{3} - n'_{1}) + n'_{1}\hat{n}_{3}(n'_{4} - n'_{2}) + n'_{2}\hat{n}_{4}(n'_{3} - n'_{1}) +$$

$$n = \hat{n} + n'$$

$$S_{nl}(f,\theta) = B + L + X$$

Line 1 contains interactions for only B Line 2 contains interactions for only L Lines 3-8 contain cross-interactions between B and L

This approximation to the full integral would be exact if all terms were retained.

The fundamental idea here is to capture the broad-scale distribution of energy parametrically and to allow "local" differences to be treated as shown below. Terms that are neglected tend to contribute in a +/- sense around s.

This could be a DIA form or a
diffusion operator, but we would
lose considerable accuracy.
$$\frac{\partial n_1}{\partial t} = B + \iint N^3_* C \left| \frac{\partial W}{\partial n} \right|^{-1} ds \ k_3 d\theta_3 dk_3 + \dots$$
$$N^3_* \text{ terms neglect terms containing } n'_2 \text{ and } n'_4 \text{ - retain } \hat{n}_2 \text{ and } \hat{n}_4$$

$$N_*^3 = \hat{n}_2 \hat{n}_4 (n'_3 - n'_1) + n'_1 n'_3 (\hat{n}_4 - \hat{n}_2) + \hat{n}_1 n'_3 (\hat{n}_4 - \hat{n}_2) + n'_1 \hat{n}_3 (\hat{n}_4 - \hat{n}_2)$$

Note: X is typically 2-3 times larger than L or B. This is why linear sums (neural networks, EOF's, etc.) do not work well for Snl estimation.

$$\frac{\partial n_{1}}{\partial t} = B + \left(\frac{k}{k_{0}}\right)^{19/2} \left\langle \begin{pmatrix} \left(\frac{\beta}{\beta_{0}}\right) \iint (\hat{n}_{1}n'_{3} + n'_{1}\hat{n}_{3} + n'_{1}n'_{3})\Lambda(\hat{n}_{2} - \hat{n}_{4}, \underline{k}_{1}, k_{*}, \theta_{*}, \mathbf{x}_{1}, \dots, \mathbf{x}_{n}) k_{*}d\theta_{*}dk_{*} \right. \\ \left. + \left(\frac{\beta}{\beta_{0}}\right)^{2} \iint (n'_{1} - n'_{3})\Lambda(\hat{n}_{2}\hat{n}_{4}, \underline{k}_{1}, k_{*}, \theta_{*}, \mathbf{x}_{1}, \dots, \mathbf{x}_{n}) k_{*}d\theta_{*}dk_{*} \right\rangle$$

where

$$\Lambda(\hat{n}_{2} - \hat{n}_{4}, \underline{k}_{1}, k_{*}, \theta_{*}, \mathbf{x}_{1}, ..., \mathbf{x}_{n}) = \bigoplus C |\frac{\partial W}{\partial n}|^{-1} (\hat{n}_{4} - \hat{n}_{2}) ds$$

$$\Lambda(\hat{n}_{2}\hat{n}_{4}, \underline{k}_{1}, k_{*}, \theta_{*}, \mathbf{x}_{1}, ..., \mathbf{x}_{n}) = \bigoplus C |\frac{\partial W}{\partial n}|^{-1} \hat{n}_{2}\hat{n}_{4} ds$$

$$\frac{\partial n_{1}}{\partial t} = \sum_{\delta f_{i}} \sum_{\delta \theta_{j}} \mu_{ij\bar{s}p} N^{3}_{*_{p}} + \mu_{ij\bar{s}d} N^{3}_{*_{D}}$$

$$\overline{\text{Pumping}} \quad \overline{\text{Diffusion}}$$

Note: This has no free parameters. All coefficients are determined from the full integral. Still conserves action, energy, and momentum since each elemental transfer is conservative.

Note: computer time is about 3x that of DIA

Approximation used here is a simple parameterization of B based on:

1 *dof* only (Peakedness)

Normalized spectra stratified by inverse wave age based on data from Currituck Sound Bering Sea Atlantic Ocean









Comparison of TSA (dashed) and DIA (dot-dash) to full integral (solid) for finite-depth case ($k_ph = 0.7$).

Example of actual spectrum with crude parameterization. In this case, the spectral shape is highly variable in terms of its angular distribution and the peak shape is not well approximated.





Comparison of the L+X contributions to the equivalent contributions in the full integral provides a good measure of the effectiveness of the second-scale introduction in the TSA

The TSA

- conserves constants of motion (action, energy, momentum)
- has its basis in the correct integral (important for complex cases)
- retains the number of degrees of freedom as the modeled spectrum
- is not limited to kph \geq 1

AND

• is much more efficient than the full integral

Number of mathematical operations is

about 3xDIA for 1 quadraplet (about 1.5 if two sets of q's are used) BUT: DIA's instability makes it unstable for moderate time steps!

about 1/250 of the time of the Full Integral

CONCLUSIONS

- DIA has extreme difficulty in reproducing SnI since it does not represent the full set of 4-wave interactions.
- Although DIA is calibrated to provide similar energy transfer to the Low-frequency region of the spectrum, the calibration is only locally valid (near γ=3.3).
- The parametric extension of the DIA to shallow water does not capture essential elements of nonlinear energy fluxes and their effects on spectral shape in coastal waters
- The TSA appears to be a more accurate alternative to the DIA for both deep water and shallow water cases
- The initial TSA can be extended to improve the B-scale treatment which should improve its accuracy for operational applications.