

3.18 Nutrient cycling and nutrient limitation in Ecosim

Ecosim uses a very simple strategy to represent nutrient cycling and potential nutrient limitation of primary production rates. It is assumed that at any instant in time the system has a total nutrient concentration N_T , which is partitioned between nutrient 'bound' in biomass versus free in the environment (accessible to plants for nutrient uptake). That is, T is represented as the sum $N_T = \sum_i ?_i B_i + N_f$, where $?_i$ is (fixed) nutrient content per unit of pool i biomass, and N_f is free nutrient concentration. Then assuming that N_T varies as $dN_T/dt = I - vN_T$, where I is total inflow rate to the system from all nutrient loading sources and v is total loss rate from the system due to all loss agents (volume exchange, sedimentation, export in harvests, etc.), and that v is relatively large, N_T is approximated in Ecosim by the (possibly moving) equilibrium value $N_T = I/v$.

Changes in nutrient loading can be simulated by assigning a time forcing function number to N_T on the [Ecosim parameters](#) form, in which case N_T is calculated as $N_T = f_t N_{T0}$ where N_{T0} is the Ecopath base estimate of N_T (at the start of each simulation) and f_t is a time multiplier ($f_t = 1$ implies Ecopath base value of N_T) supplied by the user the same as any other time forcing function. Note that under the moving equilibrium assumption, changes in f_t can be viewed as caused by either changes in input rate I or nutrient loss rate v .

The Ecopath base estimate N_{T0} of total nutrient is entered by specifying the base free nutrient proportion $pf = N_f / N_{T0}$ on entry to Ecosim (also on the [Ecosim parameters](#) form: *Time dynamic (Ecosim) > Input > Ecosim parameters*), from which the Ecosim initialization can calculate N_{T0} as simply $N_{T0} = \sum_i ?_i B_i / (1 - pf)$. Note here that the units of nutrient concentration are contained in the per-biomass relative nutrient concentrations $?_i$ and these need not be specified in any particular absolute units. During each simulation, N_f is varied dynamically by setting it equal at any time to $N_T - \sum_i ?_i B_i$ so that accumulation of nutrient in any biomass pool(s) can reduce free nutrient available to promote primary production.

Primary production rates for producer pools j are linked to free nutrient concentration during each simulation through assumed Michaelis-Menten uptake relationships of the form $P/B_j = P/B_{max,j} N_f / (K_j + N_f)$, where the parameters $P/B_{max,j}$ and K_j are calculated as part of the Ecosim initialization using input estimates by the user of the ratios $P/B_{max,j} / P/B_{Ecopath,j}$ (Ecosim [Group Info](#) form). The Michaelis constant K_j is set so that $P/B_j = P/B_{Ecopath,j}$ when N_f is at the initial concentration determined by $N_T - \sum_i ?_i B_i$ when all B_i are at Ecopath base values). The user can increase sensitivity to changes in nutrient concentration (make P/B_j more variable with changes in N_T and N_f) by increasing the input $P/B_{max,j} / P/B_{Ecopath,j}$ ratio.

The default free nutrient proportion pf is set at unity, which causes N_f to be virtually constant over time (and hence P/B_j 's to be virtually independent of nutrient concentration changes). Thus to "turn on" nutrient limitation effects, you must set a lower value for pf , (e.g., 0.3) on the Ecosim parameters form.

Users should be aware that this simple approach to accounting for nutrient limitation can interact with the numerical method used to simulate very fast phytoplankton dynamics over time, to cause numerical instability or "chattering" in the values of phytoplankton biomass. This happens mainly in cases where p_i is low, so that N_f is initially small. Then any biomass decline in the system (e.g. due to decline in zooplankton biomass) results in a relatively large increase in N_f , which can cause an over-response in the calculated phytoplankton biomass(es) B_j , which then drives N_f to near zero, which in turn causes too large a decrease in calculated B_j for the next monthly Ecosim time step.

Chattering can be reduced by using the Runge-Kutta integration option and/or higher pf settings. Improved numerical integration procedures should allow us to avoid the problem entirely in future Ecosim versions, but at present the computational cost of avoiding the problem by 'brute force' (shorter simulation time steps) would be prohibitively expensive of computer time.

Note further that the single free nutrient concentration N_f is linked to all primary producer groups in the model (through the uptake kinetics- P/B relationships), implying competition among all plant types in the model for free nutrients. This can cause major shifts in primary production structure over time, e.g. between benthic and pelagic primary production and between grazeable and non-grazeable algal types.