Vertical discretization

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Introduction

- this talk will be rather theoretical, showing some problems which must be faced when one attempts to solve set of PDEs numerically
- the only practical consequence will be that in NH integration (LNHDYN=.T.) proper vertical discretization must be used:

&NAMDYN NDLNPR=1,

- anyway, setup will complain if you forget
- so if you are tired ...

Equations to be discretized (1)

 basically these are Euler equations for perfect gas (multi-phasic case is not considered here), formulated in terrain following mass based η-coordinate:

$$\frac{\partial \pi}{\partial z} = -\rho g \qquad \pi_T \equiv 0$$

$$\pi(x, y, \eta, t) = A(\eta) + B(\eta)\pi_S(x, y, t)$$

$$\eta_T = 0 \qquad A(0) = 0 \qquad B(0) = 0$$

$$\eta_S = 1 \qquad \Rightarrow \qquad A(1) = 0 \qquad B(1) = 1$$

• for simplicity, following NH prognostic variables will be used:

$$\mathcal{P} \equiv \frac{p - \pi}{\partial w} - \text{scaled NH pressure departure}$$
$$d \equiv \frac{\partial w}{\partial z} - \text{true vertical divergence}$$

Equations to be discretized (2)

• with this choice of NH prognostic variables, system of Euler equations reads:

Equations to be discretized (3)

• system is closed with following diagnostic relations:

$$D_{3} = \nabla \cdot \mathbf{v} + X + d$$

$$X = \frac{\pi(1+\mathcal{P})}{RT} \cdot \frac{\partial \mathbf{v}}{\partial \pi} \cdot \nabla \phi$$

$$Z = \frac{\pi(1+\mathcal{P})}{RT} \cdot \frac{\partial \mathbf{v}}{\partial \pi} \cdot \nabla (gw)$$

$$gw = gw_{S} + \int_{\eta}^{1} \frac{mRT}{\pi(1+\mathcal{P})} d \, d\eta'$$

$$\phi = \phi_{S} + \int_{\eta}^{1} \frac{mRT}{\pi(1+\mathcal{P})} \, d\eta'$$

$$\dot{\pi} = \mathbf{v} \cdot \nabla \pi - \int_{0}^{\eta} \nabla \cdot (m\mathbf{v}) \, d\eta'$$

$$m\dot{\eta} = B \int_{0}^{1} \nabla \cdot (m\mathbf{v}) \, d\eta - \int_{0}^{\eta} \nabla \cdot (m\mathbf{v}) \, d\eta'$$

Basic principles of discretization

- at first glance it might seem that discretization is an easy task, just with many arbitrary choices to be done
- this is not true, such careless approach would most probably lead to unstable, in better case imprecise model
- in reality, discretization should be done in such way that it preserves as many continuous properties as possible
- this task is not trivial, since:
 - 1. no discretization scheme can preserve all continuous properties
 - 2. often it is not obvious which continuous properties are the important ones

Choice of vertical grid (1)

- in current ALADIN-NH, vertical discretization is done using finite difference approach
- vertically staggered grid of Lorenz type is used
- atmosphere is divided into *L* layers, numbered from top to bottom
- 3D prognostic variables are defined inside layers on so called full levels $1, \ldots, L$
- fluxes and vertical velocities are defined on layer interfaces or half levels $\tilde{0},\ldots,\tilde{L}$
- half level $\tilde{0}$ is the top boundary, half level \tilde{L} is the earth surface

Choice of vertical grid (2)

• schematically, vertical grid looks like this:



Introduction of continuous vertical operators

• following vertical operators can be identified in continuous system $(\Psi \text{ is arbitrary function of } \eta)$:

$$\mathcal{G}\Psi = \int_{\eta}^{1} \frac{m}{\pi} \Psi \, \mathrm{d}\eta'$$
$$\mathcal{S}\Psi = \frac{1}{\pi} \int_{0}^{\eta} m \Psi \, \mathrm{d}\eta'$$
$$\mathcal{N}\Psi = \frac{1}{\pi_{S}} \int_{0}^{1} m \Psi \, \mathrm{d}\eta$$
$$\mathcal{L}\Psi = \pi \frac{\partial^{2}}{\partial \pi^{2}} (\pi \Psi)$$

 their semi-implicit (SI) counterparts play a key role in derivation of structure equation and in elimination of variables used by Helmholtz solver

Two important properties of continuous vertical operators

- SI scheme is based on linear model which uses resting background state with flat orography, constant temperature T^* and constant surface pressure π^*_S
- for formulation of SI scheme, following two relations between continuous vertical operators are crucial:

$$\mathcal{G}^* + \mathcal{S}^* - \mathcal{G}^* \mathcal{S}^* = \mathcal{N}^* \tag{1}$$

$$\mathcal{L}^* \left[\mathcal{S}^* \mathcal{G}^* - \frac{c_p}{c_v} \mathcal{G}^* - \frac{c_p}{c_v} \mathcal{S}^* \right] = \frac{R}{c_v} \mathcal{I}$$
(2)

 \mathcal{I} - identity operator ($\mathcal{I}\Psi = \Psi$)

• these relations should hold also in discretized case

Discretized form of vertical operators (1)

• vertical functions A, B and hydrostatic pressures π are primarily defined on half levels \tilde{l} :

$$\pi_{\tilde{l}} = A_{\tilde{l}} + B_{\tilde{l}}\pi_{\tilde{L}}$$

• pressure difference across the layer l is given by formula:

$$\delta \pi_l = \pi_{\tilde{l}} - \pi_{\tilde{l}-1}$$

• integral operators \mathcal{G} , \mathcal{S} , \mathcal{N} applied on full level quantity Ψ are discretized as:

$$(\mathbf{G}\Psi)_{l} = \sum_{k=l+1}^{L} \Psi_{k} \delta_{k} + \Psi_{l} \alpha_{l}$$
$$(\mathbf{S}\Psi)_{l} = \frac{1}{\pi_{l}} \sum_{k=1}^{l-1} \Psi_{k} \delta \pi_{k} + \Psi_{l} \beta_{l}$$
$$(\mathbf{N}\Psi)_{l} = \frac{1}{\pi_{\tilde{L}}} \sum_{k=1}^{L} \Psi_{k} \delta \pi_{k}$$

Discretized form of vertical operators (2)

• symbols δ_l , α_l and β_l denote logarithmic pressure thickness of layer l, resp. its lower and upper part:



- it is not supposed that $\delta_l = \alpha_l + \beta_l$
- vertical laplacian *L* is discretized in most compact form possible, i.e. as 3-diagonal matrix:

$$(\mathbf{L}\Psi)_l = a_l \Psi_{l-1} + b_l \Psi_l + c_l \Psi_{l+1}$$

Fulfilling the constraints (1)

- the task is to find such expressions for quantities α_l^* , β_l^* , δ_l^* , π_l^* , a_l^* , b_l^* , c_l^* , which will respect constraints (1) and (2) in discretized case
- requirement that vertically discretized system respects both constraints (1) and (2) turned to be too strong, in fact it cannot be respected by any reasonable choice
- however, it is possible to satisfy weaker requirement:

$$\mathbf{G}^* + \mathbf{S}^* - \mathbf{G}^* \mathbf{S}^* = \mathbf{N}^*$$
$$\mathbf{L}^* \left[\mathbf{S}^* \mathbf{G}^* - \frac{c_p}{c_v} \mathbf{G}^* - \frac{c_p}{c_v} \mathbf{S}^* \right] = \frac{R}{c_v} \mathbf{T}^*$$

 T^* – 3-diagonal smoothing operator (sum of each row is 1)

- \bullet appearance of smoothing operator T^{\ast} means that constraint (2) is fulfilled only approximately in discretized system
- \bullet anyway, $T^*\Psi \to \Psi$ as vertical mesh size tends to zero

Fulfilling the constraints (2)

- in order to maintain stability, solution found for SI quantities must be extended also to non-linear model
- the procedure is straightforward (it is enough to remove stars), bottom boundary treatment for vertical laplacian L being the only exception
- desired results for full levels $2, \ldots, L$ are:

$$\pi_l = \sqrt{\pi_{\tilde{l}-1}\pi_{\tilde{l}}} \qquad \delta_l = \frac{\delta\pi_l}{\pi_l} \qquad \alpha_l = \beta_l = 1 - \sqrt{\frac{\pi_{\tilde{l}-1}}{\pi_{\tilde{l}}}}$$

• full level 1 is a special case:

$$\delta_1 = 2 + \frac{c_v}{R} \qquad \pi_1 = \frac{\delta \pi_1}{\delta_1} \qquad \alpha_1 = \beta_1 = 1$$

Fulfilling the constraints (3)

• vertically discretized laplacian applied on scaled NH pressure departure \mathcal{P} reads (l = 2, ..., L - 1):

$$(\mathbf{L}\mathcal{P})_{l} = \left[\pi \frac{\partial}{\partial \pi} \left(\frac{\partial(\pi \mathcal{P})}{\partial \pi}\right)\right]_{l} = \frac{1}{\delta_{l}} \left[\left(\frac{\partial(\pi \mathcal{P})}{\partial \pi}\right)_{\tilde{l}} - \left(\frac{\partial(\pi \mathcal{P})}{\partial \pi}\right)_{\tilde{l}-1}\right]$$
$$\left(\frac{\partial(\pi \mathcal{P})}{\partial \pi}\right)_{\tilde{l}} = \frac{\pi_{l+1}\mathcal{P}_{l+1} - \pi_{l}\mathcal{P}_{l}}{\pi_{l+1} - \pi_{l}}$$

• treatment of full level 1 must be consistent with elastic top boundary condition $p_T = 0$ and mass coordinate specification $\pi_T = 0$:

$$(\mathbf{L}\mathcal{P})_{1} = \frac{1}{\delta_{1}} \left[\left(\frac{\partial(\pi\mathcal{P})}{\partial\pi} \right)_{\tilde{1}} - \left(\frac{\partial(\pi\mathcal{P})}{\partial\pi} \right)_{\tilde{0}} \right]$$
$$\left(\frac{\partial(\pi\mathcal{P})}{\partial\pi} \right)_{\tilde{0}} = \frac{\pi_{1}\mathcal{P}_{1} - (\pi\mathcal{P})_{\tilde{0}}}{\pi_{1} - \pi_{\tilde{0}}} = \frac{\pi_{1}\mathcal{P}_{1} - (p_{T} - \pi_{T})}{\pi_{1} - \pi_{T}} = \mathcal{P}_{1}$$

• treatment of full level L will be explained in separate talk devoted to bottom boundary condition

Discretization of term $\partial \mathbf{v} / \partial \pi \cdot \nabla \Psi$

- quantity $\partial \mathbf{v} / \partial \pi \cdot \nabla \Psi$ appears in X and Z terms, where Ψ is either geopotential or gw
- discretization on full levels is done using per partes rule:

$$\begin{aligned} \frac{\partial \mathbf{v}}{\partial \pi} \cdot \nabla \Psi &= \frac{\partial}{\partial \pi} (\mathbf{v} \cdot \nabla \Psi) - \mathbf{v} \cdot \frac{\partial}{\partial \pi} \nabla \Psi \\ \left(\frac{\partial \mathbf{v}}{\partial \pi} \cdot \nabla \Psi \right)_l &= \frac{\mathbf{v}_{\tilde{l}} \cdot \nabla \Psi_{\tilde{l}} - \mathbf{v}_{\tilde{l}-1} \cdot \nabla \Psi_{\tilde{l}-1}}{\delta \pi_l} - \mathbf{v}_l \cdot \frac{\nabla \Psi_{\tilde{l}} - \nabla \Psi_{\tilde{l}-1}}{\delta \pi_l} = \\ &= \frac{(\mathbf{v}_{\tilde{l}} - \mathbf{v}_l) \cdot \nabla \Psi_{\tilde{l}} - (\mathbf{v}_l - \mathbf{v}_{\tilde{l}-1}) \cdot \nabla \Psi_{\tilde{l}-1}}{\delta \pi_l} \end{aligned}$$

• half level velocity $\mathbf{v}_{\tilde{l}}$ is determined by interpolation $(\tilde{l} = \tilde{1}, \dots \tilde{L} - 1)$:

$$\mathbf{v}_{\tilde{l}} = \varepsilon_l \mathbf{v}_l + (1 - \varepsilon_l) \mathbf{v}_{l+1} \qquad \mathbf{v}_{\tilde{0}} = \mathbf{v}_1$$
$$\varepsilon_l = \frac{\delta_{l+1} - \alpha_{l+1}}{\delta_{l+1} - \alpha_{l+1} + \alpha_l} \qquad \mathbf{v}_{\tilde{L}} = \mathbf{v}_L$$

Discretization of term $\nabla(gw)$

- quantity $\nabla(gw)$ appears in Z term
- discretization on half level \tilde{l} is straightforward:

$$gw_{\tilde{l}} = gw_{\tilde{L}} + \sum_{k=l+1}^{L} (RT)_k d_k \delta_k$$

$$\downarrow$$

$$\nabla(gw_{\tilde{l}}) = \nabla(gw_{\tilde{L}}) + \sum_{k=l+1}^{L} \left[\nabla(RT)_{k} d_{k} \delta_{k} + (RT)_{k} \nabla(d_{k}) \delta_{k} + (RT)_{k} d_{k} \nabla \delta_{k} \right]$$

Discretization of term $\partial(\pi \mathcal{P})/\partial \pi$

- except from vertical laplacian L, quantity $\partial(\pi P)/\partial \pi$ appears in pressure gradient term, where it is needed at full levels
- discretization is done in natural way:

$$\left(\frac{\partial(\pi\mathcal{P})}{\partial\pi}\right)_{l} = \frac{\pi_{\tilde{l}}\mathcal{P}_{\tilde{l}} - \pi_{\tilde{l}-1}\mathcal{P}_{\tilde{l}-1}}{\delta\pi_{l}}$$

• scaled NH pressure departure $\mathcal{P}_{\tilde{L}}$ at half levels is determined by simple averaging $(\tilde{l} = \tilde{1}, \dots, \tilde{L} - 1)$:

$$\mathcal{P}_{\tilde{l}} = \frac{1}{2} (\mathcal{P}_l + \mathcal{P}_{l+1}) \qquad \mathcal{P}_{\tilde{0}} = \mathcal{P}_1 \\ \mathcal{P}_{\tilde{L}} = \mathcal{P}_L$$

Discretization of term $\dot{\eta}\partial\Psi/\partial\eta$

- when eulerian advection is used, quantity $\dot{\eta}\partial\Psi/\partial\eta$ must be evaluated on full levels, Ψ being any prognostic variable
- discretization of vertical advection term is dictated by requirement of total energy conservation, which follows from per partes rules:

$$\int_{0}^{1} m\dot{\eta} \frac{\partial\Psi}{\partial\eta} \,\mathrm{d}\eta = -\int_{0}^{1} \Psi \frac{\partial}{\partial\eta} (m\dot{\eta}) \,\mathrm{d}\eta$$
$$\int_{0}^{1} m\dot{\eta} \Psi \frac{\partial\Psi}{\partial\eta} \,\mathrm{d}\eta = -\int_{0}^{1} \frac{\Psi^{2}}{2} \cdot \frac{\partial}{\partial\eta} (m\dot{\eta}) \,\mathrm{d}\eta$$

• in discretized case, they can be satisfied by choice:

$$\left(\dot{\eta}\frac{\partial\Psi}{\partial\eta}\right)_{l} = \frac{(m\dot{\eta})_{\tilde{l}}(\Psi_{l+1} - \Psi_{l}) + (m\dot{\eta})_{\tilde{l}-1}(\Psi_{l} - \Psi_{l-1})}{2\delta\pi_{l}}$$

• this formula is applicable also for l = 1 and l = L, since $\dot{\eta}_{\tilde{0}} = \dot{\eta}_{\tilde{L}} = 0$

Discretization of term $\nabla \pi / \pi$ (1)

 conservation of net angular momentum in global atmosphere is a consequence of continuous identity:

$$\int_0^1 \frac{mRT}{p} \nabla p \, \mathrm{d}\eta = \int_0^1 (\phi - \phi_S) \nabla \frac{\partial p}{\partial \eta} \, \mathrm{d}\eta$$

• its discrete counterpart can be satisfied by following choice:

$$\left(\frac{\nabla p}{p}\right)_{l} = \frac{1}{1 + \mathcal{P}_{l}} \cdot \frac{1}{\delta \pi_{l}} \left[\alpha_{l} \nabla(\delta p_{l}) + \delta_{l} \sum_{k=1}^{l-1} \nabla(\delta p_{k}) \right]$$

• application of this formula in pressure gradient term led to unstable behaviour due to incompatible $\nabla \mathcal{P}$ discretization in linear SI system and non-linear model

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 application of this formula in pressure gradient term led to unstable behaviour due to incompatible ∇P discretization in linear SI system and non-linear model ⇒ requirement of net angular momentum conservation is in conflict with model stability

Discretization of term $\nabla \pi / \pi$ (2)

• situation can be solved by splitting term $\nabla p/p$ into hydrostatic and non-hydrostatic parts:

$$\frac{\nabla p}{p} = \frac{\nabla [\pi (1+\mathcal{P})]}{\pi (1+\mathcal{P})} = \frac{\nabla \pi}{\pi} + \frac{\nabla \mathcal{P}}{1+\mathcal{P}}$$

 angular momentum conservation is imposed only on hydrostatic part:

$$\begin{pmatrix} \overline{\nabla}\pi \\ \pi \end{pmatrix}_{l} = \frac{1}{\delta\pi_{l}} \left[\alpha_{l} \nabla(\delta\pi_{l}) + \delta_{l} \sum_{k=1}^{l-1} \nabla(\delta\pi_{k}) \right] =$$

$$= \frac{1}{\delta\pi_{l}} \left[\alpha_{l} \nabla(\pi_{\tilde{l}} - \pi_{\tilde{l}-1}) + \delta_{l} \nabla\pi_{\tilde{l}-1} \right] = \frac{1}{\delta\pi_{l}} \left[\frac{C_{l}}{\pi_{l}} + \delta B_{l} \right] \nabla\pi_{\tilde{L}}$$

$$C_{l} = A_{\tilde{l}} B_{\tilde{l}-1} - A_{\tilde{l}-1} B_{\tilde{l}} \qquad \delta B_{l} = B_{\tilde{l}} - B_{\tilde{l}-1}$$

• term $\nabla \mathcal{P}/(1+\mathcal{P})$ is discretized in natural way

Discretization of term $\nabla \phi$

- \bullet quantity $\nabla\phi$ occurs in pressure gradient term and in X term
- it is needed on both half and full levels:

$$\phi_{\tilde{l}} = \phi_{\tilde{L}} + \sum_{k=l+1}^{L} \frac{(RT)_k}{1 + \mathcal{P}_k} \delta_k \qquad \phi_l = \phi_{\tilde{l}} + \frac{(RT)_l}{1 + \mathcal{P}_l} \alpha_l$$

$$\Downarrow$$

$$\nabla \phi_{\tilde{l}} = \nabla \phi_{\tilde{L}} + \sum_{k=l+1}^{L} \left[\frac{\nabla (RT)_k}{1 + \mathcal{P}_k} \delta_k - \frac{(RT)_k \nabla \mathcal{P}_k}{(1 + \mathcal{P}_k)^2} \delta_k + \frac{(RT)_k}{1 + \mathcal{P}_k} \nabla \delta_k \right]$$
$$\nabla \phi_l = \nabla \phi_{\tilde{l}} + \frac{\nabla (RT)_l}{1 + \mathcal{P}_l} \alpha_l - \frac{(RT)_l \nabla \mathcal{P}_l}{(1 + \mathcal{P}_l)^2} \alpha_l + \frac{(RT)_l}{1 + \mathcal{P}_l} \nabla \alpha_l$$

Discretization of terms $\nabla \alpha_l$, $\nabla \delta_l$

• discretization of quantities $\nabla \alpha_l$, $\nabla \delta_l$ follows directly from their definitions (l = 2, ..., L):

$$\nabla \alpha_{l} = -\frac{1}{2} \cdot \frac{C_{l}}{\pi_{\tilde{l}} \pi_{\tilde{l}-1}} (1 - \alpha_{l}) \nabla \pi_{\tilde{L}} \qquad \nabla \alpha_{1} = 0$$

$$\nabla \delta_{l} = -\frac{1}{2} \cdot \frac{C_{l}}{\pi_{\tilde{l}} \pi_{\tilde{l}-1}} \cdot \frac{1 + (1 - \alpha_{l})^{2}}{1 - \alpha_{l}} \nabla \pi_{\tilde{L}} \qquad \nabla \delta_{1} = 0$$

$$C_{l} = A_{\tilde{l}} B_{\tilde{l}-1} - A_{\tilde{l}-1} B_{\tilde{l}}$$

Conclusions

- design of reliable vertical discretization is a difficult task
- only restricted set of continuous properties can be preserved by discretized scheme
- this set must be selected very carefully, since the wrong choices easily lead to model instability
- another source of problems can be inconsistency between linear SI system and non-linear model
- sometimes conservation constraints must be relaxed in favour of model stability