The numerics of hydrostatic structured-grid coastal ocean models: state of the art and future perspectives

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Abstract

The state of the art of the numerics of hydrodynamic non-hydrostatic structured-grid coastal ocean models is reviewed here. First, some fundamental differences in the hydrodynamics of the coastal ocean, such as the large surface elevation variation compared to the mean water depth, are contrasted against large scale ocean dynamics. Then the hydrodynamic equations as they are used in coastal ocean models as well as in large scale ocean models are presented, including parameterisations for turbulent transports. As steps towards discretisation, coordinate transformations and vertical and horizontal discretisations based on a finite-volume approach are discussed with focus on the specific requirements for coastal ocean models. As in large scale ocean models, splitting of internal and external modes is essential also for coastal ocean models, but specific care is needed when wetting an drying of intertidal flats is included. As one obvious characteristics of coastal ocean models, open boundaries occur and need to be treated in a way that correct model forcing from outside is transmitted to the model domain without reflecting waves from the inside. Here, also new developments in two-way nesting are presented. Single processes such as internal inertia-gravity waves, advection and turbulence closure models are discussed with focus on the coastal scales. Some overview on existing hydrostatic structured-grid coastal ocean models is given, including their extensions towards non-hydrostatic models. Finally, an outlook on future perspectives is made.

Keywords: coastal ocean models, hydrostatic models, adaptive coordinates, mode splitting, wetting and drying, two-way nesting

1. Introduction

This review paper presents the state of the art of hydrodynamic coastal ocean modelling, with focus on structured-grid hydrostatic models. Here, we define the coastal ocean as ranging from estuaries and shallow coastal areas to the shelf break region. On the small scale end, near-shore and coastal engineering scales are not resolved, but parameterised, if needed. On the large scale end, the dynamics of the ocean basins will not be explicitly treated, since their effects are introduced by means of lateral boundary conditions.

The coastal ocean as defined here, is strongly affected by the superposition of direct human impacts and climate change and variability. Direct human impacts include for example eutrophication, fishery, large scale constructions (such as offshore wind farms), dredging and dumping of material, discharges (e.g., of cooling water) and pollution. Aspect of natural and anthropogenic climate change and variability impacting on the coastal ocean are relative sea level rise (global and local), global warming, changes of the water cycle (precipitation, evaporation, run-off), and changed wind forcing. These local, regional and global forcings may lead to severe changes in the coastal environment, such as regime shifts in biogeochemistry, coastal erosion, flooding of low-laying coastal regions, increase of oxygen minimum zones, invasion of alien species, and many more. Changes are strong in many coastal regions around the world, but specifically serious in the Arctic where the warming is strongly accelerated.

To understand the changing coastal oceans, numerical modelling is one of the most effective tools, allowing for reconstructions of historic states, hindcasts and analysis of the dynamics in the last decades, short term forecasts of coastal ocean states, as well as for coastal climate projections and possible future scenarios. One core requirement for the predictability of such model exercises is the way how basic hydrodynamic processes are reproduced.

Main aspects which need to be resolved are exchanges of matter (such as salt, heat, oxygen, nutrients and particulate matter) in vertical and lateral directions. In the lateral direction, the coast-to-ocean exchange is mediated by estuarine dynamics, tidal dispersion, tidal fronts, and meso-scale and submeso-scale dynamics. In the vertical, critical interfaces play a crucial role, such as the sediment-water interface, the *clines* (lutocline, halocline, thermocline, redoxcline), the surface and bottom wave-affected layers as well as the air-sea interface. Internal waves play a special role for linking horizontal and vertical scales. All these features need to be properly resolved.

In this paper, we discuss specific requirements for the hydrodynamic cores of coastal ocean models in terms of driving biogeochemical and suspended sediment models. This will include consequences for interfacing with models for other compartments such as atmospheric models, surface wave models, sea ice models, hydrological models and bottom sediment models, without explicitly discussing such models itself. Also issues of data assimilation will be excluded here. Scales where resolution of non-hydrostatic pressure contributions become relevant will be parameterised such that we restrict the discussions on hydrostatic models. Furthermore, we restrict ourselves to the class of structured-grid models, since the numerical methods in unstructured-grid modelling differ significantly from the structured-grid methods. Due to their flexible grids, unstructured-grid models are known to be specifically suitable for investigating multi-scale dynamics in the ocean. We will here discuss how also structured-grid models combined with nesting strategies are able to resolve multiple scales.

2. Characteristics of coastal ocean dynamics

The coastal ocean has a number of characteristics which discriminate it from the global ocean in a way that specific physical and numerical treatment is required. In the following, these characteristics are expressed in terms of characteristic horizontal and vertical length scales H and L, respectively, and the time scale T.

2.1. Relatively thick mixed layers

In temperate shelf seas, the surface and bottom mixed layers may occupy a considerable fraction of the water depth. In summer, when stratification shows a maximum, they may still cover 60%-80% of the water column, while in autumn and winter they may completely merge into a vertically homogeneous water column (Bolding et al., 2002). In contrast to this, surface and bottom mixed layers in ocean basins, which may have a comparable absolute thickness to shelf seas may cover only a small portion of the total depth of several thousands of metres (Martin (1985)).

With H being the local time-averaged water depth, and D_s and D_b denoting surface and bottom mixed layer thickness, respectively, this characteristics of coastal seas can be expressed as

$$\frac{D_s + D_b}{H} = \mathcal{O}(1) \tag{1}$$

The major consequence of this is that turbulence closure models are needed which resolve the vertical structure of the mixed layers including the interaction of vertical stratification and mixing. For the relatively thick bottom boundary layers, another implication is that the bed friction generates relatively strong dissipation of kinetic energy in the water column. The latter may be enhanced in shallow water, when the surface wave length becomes comparable to the water depth, such that the orbital wave motion adds to the bed stress. Due to the dynamic importance of the bottom boundary layer it is required that it is vertically well-resolved everywhere. This means for the generally strongly variable bottom bathymetry of the coastal ocean that *bottom-following vertical coordinates* are mandatory (see section 4.2).

2.2. Relatively high surface elevation amplitude

The sea surface elevation variability due to tides and storm surges is strongly enhanced in the shallow coastal ocean, adding an important nonlinearity to the dynamics. When denoting the standard deviation of the surface elevation as $\sigma(\eta)$, this characteristics of the coastal ocean could be formulated as

$$\frac{\sigma(\eta)}{H} = \mathcal{O}(1). \tag{2}$$

First of all, coastal ocean models need to predict the temporally and spatially varying free surface elevation η, rather than applying the rigid-lid approximation which large scale ocean models originally used in order to allow for large time steps (Bryan, 1969). Furthermore, the relatively high surface elevation variation requires the use of *surface-following vertical coordinates* to ensure sufficient vertical resolution of the near surface region.

An extreme case occurs when the surface elevation variability is larger than the mean water depth, such that the sea bed is falling dry at times:

$$\frac{\sigma(\eta)}{H} > 1. \tag{3}$$

In such cases, wetting and drying algorithms must be implemented to prevent unphysical negative water depths (see e.g., Burchard et al., 2004; Warner et al., 2013, and section 5).

2.3. Relatively strong horizontal and vertical gradients

By definition, all global riverine input is discharged into coastal seas. Since the exchange between the coasts and the interior of the adjacent shelf

seas is limited, strong gradients build up in the coastal ocean. Salinity may vary from river concentration (almost zero) to ocean concentration within a few kilometres in the horizontal. Strong stratification may be generated in the vicinity of freshwater discharge such that also vertical gradients may be strong. In addition to the salinity, also other tracers such a nutrients or suspended matter discharged by rivers may show strong gradients. Let c be a relevant tracer (e.g., salinity), then the fact that horizontal or vertical gradients may be relative strong in coastal seas can be expressed as

$$\frac{\Delta c}{|\partial_{[x,y]} c|} \ll L, \quad \frac{\Delta c}{|\partial_z c|} \ll H, \tag{4}$$

with $\partial_{[x,y]}$ denoting horizontal gradients and ∂_z denoting vertical gradients.

2.4. Relatively short time scale

The relevant time scale T in the coastal ocean is typically much shorter than in the global ocean. The key time scales in the coastal ocean are tidal, inertial, diurnal or synoptic, i.e. $\mathcal{O}$(1 day), whereas large scale ocean studies generally concentrate on time scales reaching from seasonal to millenial or beyond. The latter can be achieved by parameterising effects of tides, inertial oscillations or storms (see, e.g., Eden et al. (2014)). In estuaries and coastal regions, complex dynamical changes will occur during a tidal cycle with strong tidal asymmetries in mixing and secondary circulation, and substantial impact on transport e.g. of salt and suspended matter (see, MacCready and Geyer (2010)). One example for this is the strain-induced periodic stratification (SIPS, see Simpson et al. (1990)) due to horizontal density gradients as often observed in regions of freshwater influence (see, e.g., Rippeth et al., 2001). Interior mixing in shelf seas is often caused by inertial shear caused by episodic storm events combined with coastal coupling (Craig (1989); van Haren et al. (1999); Knight et al. (2002)), a process which cannot be parameterised in the coastal ocean due to its complex bathymetry and coastline. Therefore, different model physics e.g. in terms of turbulence closure models (see the discussion at the end of section 3.2.1) requiring different numerical schemes are needed for coastal ocean models.

2.5. Relatively small spatial scales

Probably the most obvious characteristics of the coastal ocean is its limited spatial extent:

$$L \ll r_{Earth}, \tag{5}$$

with the Earth's radius r_{Earth}. Even if the entire coastal ocean is considered as one single global object (see concepts by Holt et al. (2009)), large scale global or regional drivers need to be transferred to the coastal ocean. For the remote meteorological forcing (weather systems) this is typically done by one-way or two-way nesting to a regional atmospheric model which itself is coupled to a global model. For the remote oceanic impact (tides, surges, planetary waves, etc.) this has to be done via lateral boundary conditions coming from external models. Therefore, the problem of *open boundary conditions (OBCs) and model nesting* is characteristic for coastal ocean models (as well as for the basin-scale ocean models), see the discussions in Section 6.

3. Basic equations

3.1. Momentum and tracer equations

Coastal and large scale ocean models typically use a common set of dynamic equations, which is based on the Reynolds-averaged primitive equations including hydrostatic and Boussinesq approximations. Detailed derivations of these equations are e.g. found in Griffies and Adcroft (2008).

These equations are derived from the Navier-Stokes equations on the rotating Earth by first adopting the Boussinesq approximation which includes the use of a constant reference density ρ_0 (instead of the potential density ρ) in all terms except in the gravity term. This has the consequence that mass conservation is replaced by volume conservation. In a next step, the effects of small-scale turbulent fluctuations on the larger scales are considered in a parameterised way without the necessity to numerically resolve them. This procedure assumes an ensemble averaging of the Navier-Stokes equations resulting in the Reynolds-averaged Navier-Stokes equations which now include unknown turbulent stresses which need to be parameterised (see section 3.2). Finally, a scaling analysis assuming a small aspect ratio (vertical scales are much smaller than horizontal scales) results in the degeneration of the vertical component of the momentum equation to the hydrostatic balance, neglecting vertical acceleration, the non-linear and the stress divergence terms and the effects of Earth rotation (see the discussion in Kanarska et al., 2007; Klingbeil and Burchard, 2013). The hydrostatic approximation includes the traditional approximation for the Coriolis terms, neglecting the horizontal contributions of Earth rotation in the horizontal momentum equations. The

dynamic equations are here for simplicity expressed in local Cartesian coordinates (see section 4 for spherical and curvilinear coordinates). After inserting the Boussinesq approximation into the mass conservation equation, the incompressibility condition remains:

$$\partial_x u + \partial_y v + \partial_z w = 0, \tag{6}$$

with the space coordinate vector (x, y, z) (eastwards, northwards, upwards), and the three-dimensional velocity vector (u, v, w).

Applying the incompressibility condition (6) and the hydrostatic condition (with pressure p and gravitational acceleration $g = 9.81\,\mathrm{m\,s^{-2}}$),

$$\partial_z p = -g\rho, \tag{7}$$

(remainder of the vertical component of the Navier-Stokes equations) into the horizontal momentum equations, the resulting dynamic equations are of the following form:

$$\begin{aligned} &\partial_t u + \partial_x\left(u^2\right) + \partial_y\left(uv\right) + \partial_z\left(uw\right) + \partial_{\tilde{x}} J^{\mathrm{u}}_{\tilde{x}} + \partial_{\tilde{y}} J^{\mathrm{u}}_{\tilde{y}} + \partial_z J^{\mathrm{u}}_z - fv \\ &\quad = -\frac{1}{\rho_0}\partial_x p_a - \frac{g}{\rho_0}\partial_x \int_z^\eta \rho \,\mathrm{d}\xi \\ &\quad = -\frac{1}{\rho_0}\partial_x p_a - g\partial_x \eta + \int_z^\eta \partial_x b \,\mathrm{d}\xi \end{aligned} \tag{8}$$

and

$$\begin{aligned} &\partial_t v + \partial_x\left(vu\right) + \partial_y\left(v^2\right) + \partial_z\left(vw\right) + \partial_{\tilde{x}} J^{\mathrm{v}}_{\tilde{x}}(u,v) + \partial_{\tilde{y}} J^{\mathrm{v}}_{\tilde{y}}(u,v) + \partial_z J^{\mathrm{v}}_z + fu \\ &\quad = -\frac{1}{\rho_0}\partial_y p_a - \frac{g}{\rho_0}\partial_y \int_z^\eta \rho \,\mathrm{d}\xi \\ &\quad = -\frac{1}{\rho_0}\partial_y p_a - g\partial_y \eta + \int_z^\eta \partial_y b \,\mathrm{d}\xi, \end{aligned} \tag{9}$$

with time t, the local Coriolis frequency $f = 2\Omega\sin(\phi)$ (with the angular speed of Earth rotation $\Omega = 7.29 \cdot 10^{-5}\mathrm{s}^{-1}$ and the latitude ϕ), the surface elevation η, and the surface pressure p_a. $J^{\mathrm{u}}_{[\tilde{x},\tilde{y},z]}$ and $J^{\mathrm{v}}_{[\tilde{x},\tilde{y},z]}$ are shear stress vectors along the local plain spanned by the vector $(\tilde{x}, \tilde{y})$ which might be

along geopotentials ($\tilde{x} = x$, $\tilde{y} = y$), or along general vertical coordinates or along isopycnals. This simplifying notation is used to avoid the introduction of rotated diffusion tensors and to facilitate the practical numerical implementation. The terms on the left hand sides of (8) and (9) are the tendency term, the three components of the advection terms (written in so-called integral form), the stress divergence terms, and the Coriolis term. The terms on the right hand side form the horizontal pressure gradient components $\partial_x p/\rho_0$ and $\partial_y p/\rho_0$. The formulation in the last line splits the pressure gradient into the external pressure gradient (composed of the atmospheric pressure gradient and the surface slope) and the internal pressure gradient, where $b = -(g/\rho_0)(\rho - \rho_0)$ is the buoyancy. The last lines of (8) and (9) result from assuming $\rho(\eta)/\rho_0 \approx 1$ which is consistent with the Boussinesq approximation. In (8) and (9), the flux form for the advective terms is used, which results for any quantity q from the incompressibility condition (6):

$$u\partial_x q + v\partial_y q + w\partial_z q = \partial_x(uq) + \partial_y(vq) + \partial_z(wq), \tag{10}$$

which is the preferred form for coastal models in contrast to large-scale models sometimes based on the vector-invariant form for nonlinear terms, which allow an integral conservation of potential enstrophy and/or total energy (e.g. Arakawa and Lamb, 1977) (see section 7.1).

Equations (8) and (9) are valid for $-H(x,y) < z < \eta(x,y,t)$ (with the bottom coordinate $z = -H$). Dynamic boundary conditions at the free surface are

$$J_z^{[\mathrm{u,v}]}\big|_{z=\eta} = \frac{\tau^{\mathrm{s}}_{[\mathrm{x,y}]}}{\rho_0}, \tag{11}$$

with the surface stress vector $(\tau^{\mathrm{s}}_{\mathrm{x}}, \tau^{\mathrm{s}}_{\mathrm{y}})$, which is calculated from meteorological data using bulk formulae (e.g., Kondo (1975); Fairall et al. (1996)). At the bottom, the dynamic boundary conditions are the no-slip conditions

$$[u, v]\big|_{z=-H} = 0. \tag{12}$$

For numerical reasons, however, generally an equivalent flux boundary condition is applied,

$$J_z^{[\mathrm{u,v}]}\big|_{z=-H} = \frac{\tau^{\mathrm{b}}_{[\mathrm{x,y}]}}{\rho_0}, \tag{13}$$

with the bed stress vector $(\tau^{\mathrm{b}}_{\mathrm{x}}, \tau^{\mathrm{b}}_{\mathrm{y}})$ reconstructed from a logarithmic velocity profile (see Section 7.5).

Kinematic surface and bottom boundary conditions are:

$$w\big|_{z=\eta} = \partial_t \eta + u\big|_{z=\eta} \partial_x \eta + v\big|_{z=\eta} \partial_y \eta + (E - P), \tag{14a}$$

$$w\big|_{z=-H} = -u\big|_{z=-H} \partial_x H - v\big|_{z=-H} \partial_y H, \tag{14b}$$

with the evaporation rate E and the precipitation rate P. River run-off can be treated in various ways, e.g. as lateral boundary condition.

There is a large variety of lateral boundary conditions, which will be further discussed in section 6.

Within the Boussinesq approximation, the budget equation for internal energy reduces to a conservation equation for potential temperature θ:

$$\partial_t \theta + \partial_x (u\theta) + \partial_y (v\theta) + \partial_z (w\theta) + \partial_{\tilde{x}} J^{\theta}_{\tilde{x}} + \partial_{\tilde{y}} J^{\theta}_{\tilde{y}} + \partial_z J^{\theta}_z = \frac{1}{C_p \rho_0} \partial_z I, \tag{15}$$

with the horizontal and vertical turbulent fluxes of θ given by $J^{\theta}_{[\tilde{x},\tilde{y},z]}$, the specific heat capacity of water at constant pressure, C_p, and irradiance in water, I. The positive vertical gradient of irradiance depends on the light absorption model applied. A classical model would be the model by Paulson and Simpson (1977), but often more complex models taking into account water colour, suspended matter and biogeochemical components are used. In (15), the heating due to viscous dissipation of turbulent kinetic energy is neglected, since in comparison to other terms in the local heat budget of the ocean this is a very small term (in contrast to conditions in the higher atmosphere). The surface boundary condition for the turbulent flux of potential temperature is formulated as

$$J^{\theta}_z\big|_{z=\eta} = \frac{Q_s + Q_l + Q_b}{C_p \rho_0}, \tag{16}$$

with the sensitive heat flux Q_s (due to the air-sea temperature difference), the latent heat flux Q_l (due to evaporation), and the long-wave back radiation Q_b. Note that upward (positive) fluxes of potential temperature lead to cooling at the surface. For the parameterisation of these heat fluxes various models are available, with the bulk formulae by Kondo (1975) and Fairall et al. (1996) being the most common ones for coastal applications.

The conservation equation for the salinity S is of similar form as the temperature conservation equation, but without source term:

$$\partial_t S + \partial_x (uS) + \partial_y (vS) + \partial_z (wS) + \partial_{\tilde{x}} J^{\mathrm{S}}_{\tilde{x}} + \partial_{\tilde{y}} J^{\mathrm{S}}_{\tilde{y}} + \partial_z J^{\mathrm{S}}_z = 0. \tag{17}$$

The diffusive salinity flux at the surface is given as

$$J_z^{\mathrm{S}}\big|_{z=\eta} = -(E-P)S, \tag{18}$$

see Beron-Vera et al. (1999) for details. The salinity unit which should be used in all ocean models is the absolute salinity which indicates the mass fraction of salt in seawater in [g/kg], defined according to the TEOS-10 standard (IOC, SCOR and IAPSO (2010)). The potential density is calculated by means of an empirical equation of state as function of potential temperature θ, absolute salinity S and pressure p:

$$\rho = \rho(\theta, S, p). \tag{19}$$

A recent and accurate formulation for Boussinesq models has been provided by Roquet et al. (2015).

3.2. Turbulence closure modelling

Turbulence closure modelling involves complex theories which cannot be discussed here. For an overview, the reader is referred to text books (Kantha and Clayson, 2000b; Griffies, 2004; Cushman-Roisin and Beckers, 2011; Olbers et al., 2012). In most ocean models, vertical and horizontal turbulent fluxes of scalars are parameterised by a down-gradient approximation of the form

$$J^c_{[\tilde{x},\tilde{y}]} = -A_c \partial_{[\tilde{x},\tilde{y}]} c, \tag{20}$$

$$J^c_z = -K_c \partial_z c, \tag{21}$$

using horizontal and vertical eddy diffusivities A_c and K_c for any scalar c. With this, the turbulent fluxes are directed towards the lower concentrations, as postulated for molecular fluxes in ideal gases.

For the vertical turbulent momentum fluxes, an equivalent formulation is generally used:

$$J_z^{[\mathrm{u,v}]} = -K_{\mathrm{M}} \partial_z [u, v], \tag{22}$$

with the vertical eddy viscosity K_{M}. For the parameterisation of the horizontal turbulent momentum fluxes, additional invariance to vector rotation

needs to be taken into account (Blumberg and Mellor, 1987):

$$\begin{aligned} J^{\mathrm{u}}_{\tilde{x}}(u) &= -2A_{\mathrm{M}}\partial_{\tilde{x}}u, \\ J^{\mathrm{u}}_{\tilde{y}}(u,v) &= -A_{\mathrm{M}}\left(\partial_{\tilde{y}}u + \partial_{\tilde{x}}v\right), \\ J^{\mathrm{v}}_{\tilde{x}}(u,v) &= -A_{\mathrm{M}}\left(\partial_{\tilde{y}}u + \partial_{\tilde{x}}v\right), \\ J^{\mathrm{v}}_{\tilde{y}}(v) &= -2A_{\mathrm{M}}\partial_{\tilde{y}}v, \end{aligned} \tag{23}$$

with the horizontal eddy viscosity A_{M}.

3.2.1. Small-scale closure models

For the parameterisation of vertical eddy viscosity and diffusivity, two basic approaches are commonly used in ocean models.

One of them is based on direct algebraic computation of the eddy parameters from bulk properties (boundary fluxes, mixed layer depth, bulk Richardson number) of the surface and bottom mixed layers, the so-called K-profile parameterisation (KPP) which has been originally proposed for large-scale ocean models by Large et al. (1994). In addition to the down-gradient fluxes, also counter-gradient flux components (for example for convectively unstable boundary layers) are included in the KPP model. Many extensions of the original KPP model have been proposed (see, e.g., Smyth et al., 2002; Durski et al., 2004; Uchiyama et al., 2010) to extend its applicability also to coastal ocean applications (see, e.g., Blaas et al., 2007). The KPP model assumes that the boundary layer turbulence is in equilibrium with the surface or bottom fluxes, respectively. For large scale ocean models, this might often be an adequate approximation. In many coastal model applications however, the deviation from this equilibrium is a first-order process which cannot be neglected.

The other closure approach, which in contrast to the KPP model considers high temporal and spatial variability in boundary layers, is based on a dynamic equation for the turbulent kinetic energy (TKE), k, which can be derived from the Navier-Stokes equations by means of Reynolds-averaging (Wilcox, 1998; Burchard, 2002a):

$$\partial_t k + \partial_x\left(uk\right) + \partial_y\left(vk\right) + \partial_z\left(wk\right) - \partial_z\left(\frac{K_{\mathrm{M}}}{Sc_k}\frac{\partial k}{\partial z}\right) = P + B - \varepsilon, \tag{24}$$

with the shear production, $P = K_{\mathrm{M}} M^2$ including the vertical shear squared, M^2, and the buoyancy production, $B = -K_c N^2$, with the buoyancy frequency (or Brunt-Väisälä frequency) squared, N^2, and the dissipation rate, ε. The shear production P transfers kinetic energy lost from the mean flow due to vertical stress divergence to the turbulent kinetic energy, whereas the buoyancy production B exchanges energy between the mean potential energy to the turbulent kinetic energy. Note, that in (24) the down-gradient approximation has been adopted with the turbulent Schmidt number Sc_k, which is of order unity. Horizontal turbulent fluxes are generally neglected for the TKE budget (Delhez et al., 1999). This may also be the case for the advective terms, but in tidally energetic regimes, such a neglect might lead to significant errors or instabilities. In addition to the TKE, also the length scale of the energy containing eddies, l, needs to be approximated to obtain an estimate for the vertical eddy viscosity and eddy diffusivity. Once k and l are known, the vertical eddy coefficients can be calculated as

$$K_{\mathrm{M}} = S_{\mathrm{M}} k^{1/2} l, \qquad K_c = S_c k^{1/2} l, \tag{25}$$

with the non-dimensional stability functions S_{M} and S_c. These stability functions can be parameterised by means of second-moment closures based on Reynolds-averaging the Navier-Stokes equations (Mellor and Yamada, 1982; Canuto et al., 2001) or based on simpler empirical relations (Rodi, 1987; Burchard and Baumert, 1995). The dissipation rate ε can be calculated from l via the cascading relation

$$\varepsilon = c_\varepsilon \frac{k^{3/2}}{l} \tag{26}$$

with the empirical constant c_ε.

Surface and bottom boundary conditions are typically derived from the law of the wall, resulting in k being proportional to the surface or bed stress, respectively, and the length sale to be proportional to the respective roughness length. Equivalent flux conditions, which proved to be numerically more stable have been proposed (Burchard and Petersen, 1999). To parameterise effects of breaking surface waves, injection of TKE at the surface is applied (Burchard, 2001). Effects of Langmuir Circulation on the dynamics of the surface mixed layer may be parameterised as well (see, e.g., Harcourt, 2014). Additional frictional effects of surface waves on wave drag in shallow water can also be considered in the bottom boundary condition (Mellor, 2002).

Various methods are used in ocean modelling to calculate the length scale l. In one-equation turbulence closure models (the one equation is for the

computation of the TKE), l is calculated from combinations of geometric length scales (such as the distance from the surface or the bottom) or from hydrodynamic length scales (such as the Ozmidov scale), see e.g., Mellor and Yamada (1982); Gaspar et al. (1990). More common in coastal ocean models are however two-equation turbulence closure models which in addition to the TKE equation calculate a budget equation for a length-scale related quantity such as for the dissipation rate ε (k-ε models, see Rodi, 1987; Burchard and Baumert, 1995), for the turbulent frequency $\omega = \varepsilon/k$ (k-ω models, see Wilcox, 1998; Umlauf et al., 2003), the quantity kl (k-kl models, see Mellor and Yamada, 1982), or the generic length-scale related property $k^m l^n$ with real numbers m and n (the generalised length-scale model, see Umlauf and Burchard, 2003). When properly calibrated, all these two-equation models perform similarly (Burchard et al., 1998; Umlauf and Burchard, 2003; Warner et al., 2005).

Two-equation turbulence closure models including algebraic second-moment closures as outlined here (see also the review by Umlauf and Burchard, 2005) have proven in many applications as a good compromise between computational efficiency and realistic representation of relevant processes in the coastal ocean.

There is a number of reasons, why such second-moment turbulence closure models are a necessary feature of coastal ocean models. First of all, as discussed in section 2.1, a large portion of water columns in the coastal ocean (if not the entire water column) are covered by surface or bottom boundary layers. These boundary layers have a vertical structure which locally influences the mixing coefficients. Therefore, boundary layer models reacting on local forcing instead of surface or bottom fluxes only are required. Secondly, relevant times scale in the coastal ocean are much shorter than in the global ocean (see section 2.4). Therefore, the turbulence closure models need to react on correct time scales to rapid time changes in surface or bottom stress instead of instantaneouly adjusting the eddy coefficients in the entire mixed layers such as in the KPP model. Such requirements are only fulfilled by carefully calibrated second-moment turbulence closure models.

There are a few global ocean models using low-level second-moment closures (see, e.g., Oschlies (2002); Danabasoglu et al. (2014)), which typically are based on the Gaspar et al. (1990) one-equation model, where the turbulent length scale l is algebraically computed. Two-equation turbulence closure models are generally too complex for global ocean models. This is mainly due to the large baroclinic time steps of global models (about 1 hour)

and the coarse resolution in boundary layers. Increasing spatial and temporal resolution of large-scale ocean models in order to allow for the inclusion of two-equation turbulence closure models would result in addionally high computational costs which probably might not pay off on the temporal and spatial scales relevant for global ocean simulations. Therefore, in large-scale ocean models, the KPP model (see above) or other models with reduced dynamics (see, e.g., Jackson et al. (2008)) are used.

3.2.2. Meso-scale and submeso-scale closure models

Whereas in ocean models a consistent theoretical framework for the vertical turbulent fluxes exists, such a rigorous closure for the horizontal turbulent fluxes due to meso-scale and submeso-scale eddies is still missing. In the stratified ocean, the physically most relevant direction of horizontal mixing is along isopycnal surfaces. Since this is however numerically demanding in non-isopycnal models (see the discussion by Griffies et al., 1998; Beckers et al., 1998), in many model applications horizontal mixing is simply carried out along coordinate surfaces.

Two simplistic common approaches used in the coastal ocean modelling community are to either neglect horizontal turbulent fluxes (and then rely on the numerical fluxes due to monotone, upstream-biased advection schemes, see section 7) or to set A_M and A_c to constant values. In the latter case, the proper choice of the horizontal exchange coefficients is then determined by sensitivity studies. A physically more sound, resolution-dependent parameterisation of horizontal momentum fluxes has been proposed by Smagorinsky (1963). For incompressible fluids under the (quasi)hydrostatic pressure approximation the horizontal eddy viscosity can be derived as

$$A_\mathrm{M} = C_\mathrm{M}(\Delta x\,\Delta y)\sqrt{(\partial_{\bar{x}}u)^2 + \tfrac{1}{2}\left(\partial_{\bar{y}}u + \partial_{\bar{x}}v\right)^2 + (\partial_{\bar{y}}v)^2}, \tag{27}$$

with the horizontal grid sizes Δx and Δy, and with the empirical parameter C_M. The latter has been estimated by Holt and James (2006) in a shelf sea application as $C_\mathrm{M} \approx 0.2$. The same approach as (27) can also be applied for horizontal tracer diffusivities (see, e.g., Gräwe et al., 2015), using a different value for the empirical parameter.

For large scale ocean models which do not resolve meso-scale eddies, other closures have been proposed, such as the Gent and McWilliams (1990) closure for isopycnal mixing or the dynamic meso-scale eddy closure by Eden and Greatbatch (2008). However, in coastal ocean models eddies are at least

partially resolved, such that these closures cannot directly be used in such models. On the other hand, theories for submeso-scale dynamics are not yet mature and general enough to derive applicable eddy closures.

4. Coordinate transformations and spatial discretisation

The set of governing equations, consisting of the continuity equation (6), the hydrostatic momentum equations (8) and (9), and the prognostic equations for temperature (15) and salinity (17), so far has been formulated in a local Cartesian reference frame at the Earth surface for simplicity. However, the transformation to coordinate systems that align with the geometry of the domain is often advantageous.

Most coastal ocean models support orthogonal curvilinear coordinates in the horizontal to offer to some extent the alignment with coastlines and an increased horizontal resolution towards the coast as required by the typically strong near-coastal horizontal gradients, see (4). The corresponding basis transformation introduces additional metric coefficients and terms that also offer the formulation in spherical coordinates, see Sec. 4.3 for details. For clarity, the governing equations in the remainder of this review will be kept in horizontal Cartesian coordinates.

In the vertical the alignment with the boundaries (free surface and bottom topography) and proper resolution (of e.g. the boundary layers and pycnoclines) is even more critical than in the horizontal. A general description that supports various types of vertical coordinates is given by the s-coordinates of Kasahara (1974). The general vertical coordinate $s(x, y, z, t)$ must be strictly monotone in z in order to guarantee a unique inverse transformation $z(x, y, s, t)$ and a well-defined Jacobian $\partial_s z = (\partial_z s)^{-1}$ for the transformation of the governing equations. For isopycnal coordinates $s = \rho$ this precludes the representation of unstable stratification. The horizontal coordinates and time do not change under the transformation and, if necessary for the purpose of clarity, derivatives with respect to constant s and z will be distinguished by $(\cdot)_s$ and $(\cdot)_z$, respectively. With the Lagrangian derivative $\mathrm{D}/\mathrm{D}t$ it can easily be shown, that the grid-related vertical velocity, defined as the difference of the fluid's vertical velocity and the vertical velocity of the coordinate

surfaces w^{grid}, is given by

$$w^{\mathrm{s}} = w - w^{\mathrm{grid}}$$
$$= \left[\frac{\mathrm{D}}{\mathrm{D}t} - \left(\frac{\mathrm{D}}{\mathrm{D}t}\right)_s\right] z\left(x,y,s,t\right) = \partial_s z \left[\frac{\mathrm{D}}{\mathrm{D}t} - \left(\frac{\mathrm{D}}{\mathrm{D}t}\right)_s\right] s = \partial_s z \frac{\mathrm{D}s}{\mathrm{D}t}. \quad (28)$$

4.1. Layer-integrated equations

The vertical discretisation of the transformed space into $k_{\max}$ layers with interfaces at fixed levels $s_{k+1/2}$ (independent of x, y, t and with $k \in [0, k_{\max}]$) is equivalent to a discretisation of the physical space into $k_{\max}$ layers with interfaces at variable levels $z_{k+1/2}\left(x,y,t\right) = z\left(x,y,s_{k+1/2},t\right)$ and $z_{k+1/2} > z_{k-1/2}$, see Burchard and Petersen (1997).

With layer heights $h_k(x,y,t) = \int_{s_{k-1/2}}^{s_{k+1/2}} \partial_s z \mathrm{d}s = z_{k+1/2} - z_{k-1/2}$, layer-averaged prognostic quantities $\varphi_k(x,y,t) = h_k^{-1} \int_{z_{k-1/2}}^{z_{k+1/2}} \varphi \,\mathrm{d}z$, for $k \in [1, k_{\max}]$ and $\varphi \in \{u, v, \theta, S\}$, and $w^{\mathrm{s}}_{k+1/2}\left(x,y,t\right) = w^{\mathrm{s}}\left(x,y,z_{k+1/2},t\right)$, the layer-integration of (6), (8), (9), (15) and (17) yields:

$$\partial_t h_k + \partial_x\left(h_k u_k\right) + \partial_y\left(h_k v_k\right) + \left(w^{\mathrm{s}}_{k+1/2} - w^{\mathrm{s}}_{k-1/2}\right) = 0, \quad (29)$$

$$\partial_t\left(h_k u_k\right) + \partial_x\left(h_k u_k u_k\right) + \partial_y\left(h_k v_k u_k\right) + \left(w^{\mathrm{s}}_{k+1/2}\widetilde{u}_{k+1/2} - w^{\mathrm{s}}_{k-1/2}\widetilde{u}_{k-1/2}\right)$$
$$= f h_k v_k - g h_k \partial_x \eta - h_k \tfrac{1}{\rho_0}\partial_x p_a + \left(\tau_{\mathrm{x},k+1/2} - \tau_{\mathrm{x},k-1/2}\right)$$
$$+ h_k \widetilde{F}^{\mathrm{h}}_{\mathrm{x}}(h_k, u_k, v_k) + h_k F^{\mathrm{IP}}_{\mathrm{x},k}, \quad (30)$$

$$\partial_t\left(h_k v_k\right) + \partial_x\left(h_k u_k v_k\right) + \partial_y\left(h_k v_k v_k\right) + \left(w^{\mathrm{s}}_{k+1/2}\widetilde{v}_{k+1/2} - w^{\mathrm{s}}_{k-1/2}\widetilde{v}_{k-1/2}\right)$$
$$= -f h_k u_k - g h_k \partial_y \eta - h_k \tfrac{1}{\rho_0}\partial_y p_a + \left(\tau_{\mathrm{y},k+1/2} - \tau_{\mathrm{y},k-1/2}\right) \quad (31)$$
$$+ h_k \widetilde{F}^{\mathrm{h}}_{\mathrm{y}}(h_k, u_k, v_k) + h_k F^{\mathrm{IP}}_{\mathrm{y},k},$$

$$\partial_t\left(h_k \theta_k\right) + \partial_x\left(h_k u_k \theta_k\right) + \partial_y\left(h_k v_k \theta_k\right) + \left(w^{\mathrm{s}}_{k+1/2}\widetilde{\theta}_{k+1/2} - w^{\mathrm{s}}_{k-1/2}\widetilde{\theta}_{k-1/2}\right)$$
$$= -\partial_x\left(h_k J^{\theta}_{\tilde{x},k}\right) - \partial_y\left(h_k J^{\theta}_{\tilde{y},k}\right) - \left(J^{\theta}_{z,k+1/2} - J^{\theta}_{z,k-1/2}\right) \quad (32)$$
$$+ \frac{1}{C_p \rho_0}\left(I_{k+1/2} - I_{k-1/2}\right),$$

$$\partial_t\left(h_k S_k\right)+\partial_x\left(h_k u_k S_k\right)+\partial_y\left(h_k v_k S_k\right)+\left(w^{\mathrm{s}}_{k+1/2}\widetilde{S}_{k+1/2}-w^{\mathrm{s}}_{k-1/2}\widetilde{S}_{k-1/2}\right)$$

$$=-\partial_x\left(h_k J^{\mathrm{S}}_{\tilde{x},k}\right)-\partial_y\left(h_k J^{\mathrm{S}}_{\tilde{y},k}\right)-\left(J^{\mathrm{S}}_{z,k+1/2}-J^{\mathrm{S}}_{z,k-1/2}\right). \tag{33}$$

In (29)-(33) the interfacial values $\widetilde{\varphi}_{k+1/2}$ with $\varphi\in\{u,v,\theta,S\}$ are approximated by the vertical advection scheme, see Sec. 7.1. The salinity of freshwater fluxes is zero, i.e. $\widetilde{S}_{k_{\max}+1/2}=\widetilde{S}_{1/2}=0$. According to (28) the kinematic boundary conditions (14a) and (14b) transform to $w^{\mathrm{s}}_{k_{\max}+1/2}=E-P$ and $w^{\mathrm{s}}_{1/2}=0$.

In (30) and (31) the internal pressure gradient force is given by

$$F^{\mathrm{IP}}_{[x,y],k}=\frac{1}{h_k}\int_{z_{k-1/2}}^{z_{k+1/2}}\left(\partial_{[x,y]}\right)_z\int_z^{\eta} b\mathrm{d}\xi\mathrm{d}z, \tag{34}$$

and is approximated in terms of the layer-averaged buoyancies b_k (see, e.g. Shchepetkin and McWilliams, 2003 and Sec. 7.2).

Considering the dynamic boundary conditions (11) and (13), the vertical shear stresses are defined as

$$\tau_{[\mathrm{x,y}],k+1/2}=\begin{cases}\tau^{\mathrm{s}}_{[\mathrm{x,y}]}(u_{k_{\max}},v_{k_{\max}}), & k=k_{\max}\\ K_{\mathrm{M},k+1/2}\dfrac{[u,v]_{k+1}-[u,v]_k}{z_{k+1}-z_k}, & k\in[1,k_{\max}-1]\\ \tau^{\mathrm{b}}_{[\mathrm{x,y}]}(h_1,u_1,v_1), & k=0\end{cases}, \tag{35}$$

with centre positions $z_k=\frac{1}{2}\left(z_{k-1/2}+z_{k+1/2}\right)$.

The lateral contributions of the stress tensor are approximated according to Mellor and Blumberg (1985) and Kantha and Clayson (2000a):

$$\widetilde{F}^{\mathrm{h}}_{[\mathrm{x,y}]}(h,u,v)=-\frac{1}{h}\left(\partial_x\left(hJ^{[\mathrm{u,v}]}_{\tilde{x}}(u,v)\right)+\partial_y\left(hJ^{[\mathrm{u,v}]}_{\tilde{y}}(u,v)\right)\right), \tag{36}$$

with the diffusive fluxes $J^{[\mathrm{u,v}]}_{[\tilde{x},\tilde{y}]}$ given by (23) and all gradients calculated along layer k. In case of diffusion along directions not aligned with constant s, additional terms due to non-diagonal elements of the diffusion tensor would occur (Griffies (2004)).

In (32) and (33) $J^{c}_{[\tilde{x},\tilde{y}],k}$ denotes the lateral turbulent fluxes (again with all gradients calculated along layer k), $J^{c}_{z,k\pm1/2}$ the vertical turbulent fluxes at the layer interfaces, with $J^{\mathrm{S}}_{k_{\max}+1/2}=0$ and $I_{k\pm1/2}$ the interfacial irradiance.

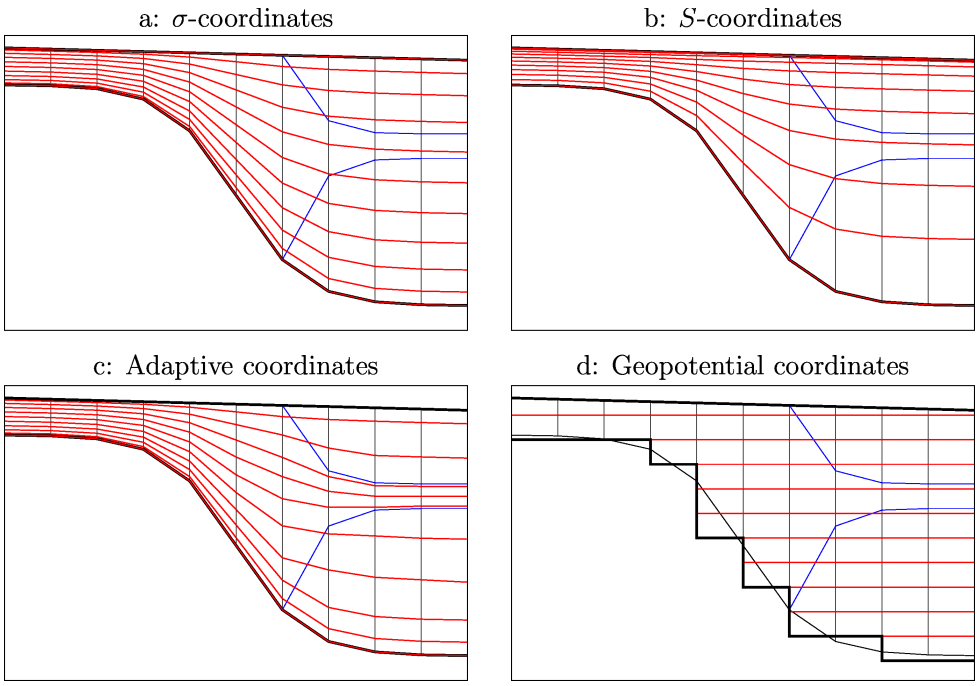

Figure 1: Sketch showing a vertical transect through a coastal ocean model domain including four different vertical coordinates. The blue lines indicate two selected isopycnals representing a tidal front with a stratified shelf sea region to the right and a well-mixed coastal region to the left. The vertical black lines indicate the horizontal discretisation, the red lines show the vertical coordinate surfaces and the bold black lines show the bottom and surface representation of the coordinates. a: σ-coordinates with zooming towards the surface and the bottom; b: S-coordinates with high near-surface resolution; c: Adaptive coordinates with increased resolution in the stratified region; d: Geopotential coordinates without partial cells at the bottom.

4.2. Examples for vertical coordinates

The almost arbitrary choice of the transformation function $s = s(x, y, z, t)$ or equivalently of the moving layer interfaces $z_{k+1/2}(x, y, t)$ allows for a high degree of freedom for the vertical layout which can be exploited to improve the model performance.

4.2.1. Geopotential coordinates

Geopotential (or z-) coordinates (see Figure 1d) are based on globally defined z-levels $z_{k+1/2}$ which do not depend on location or time, except for the uppermost coordinate level $z_{k_{\max}+1/2} = \eta$ which is identical to the free surface. Vertical resolution is typically increased near the surface to properly reproduce the dynamics of the surface mixed layer. Since for too fine near-surface vertical resolution the moving and sloping free surface might intersect with one or more coordinate levels, the uppermost geopotential coordinate level $z_{k_{\max}-1/2}$ is chosen such that it is below the minimum surface elevation. For models of shelf seas with a large maximum tidal range (see Sec. 2.2) this would result in a large mean thickness of the upper layer of $\mathcal{O}(10\text{m})$, such that near-surface processes are not well-resolved (see, e.g., Mathis et al. (2015)). For simulations of shallow coastal regions including drying and flooding this would mean that all areas with intertidal flats are not vertically resolved. One solution to this problem has been proposed by Casulli and Cheng (1992) and Burchard and Baumert (1998): Layers at the surface are locally added or removed when the surface layer becomes too thick or too thin, respectively. This problem can also be solved by applying so-called z^*-coordinates (Adcroft and Campin (2004)), where geopotential coordinates are obtained for $\eta = 0$ everywhere, and all layer thicknesses are then locally multiplied by $(\eta + H)/H$, such that the layers are squeezed or expanded with the changing water depth. Another disadvantage of geopotential models near the surface (which is not solved by the above mentioned treatment) is that the tidal elevation triggers a large oscillating vertical velocity relative to the fixed coordinate levels which generates strong numerical mixing due to the vertical advection schemes, see Sec. 7.1.

At the bottom geopotential layers will be intersecting with the sloping bottom. This leads to a stair-case type of discretisation of bottom slopes which leads to low accuracy in resolving bottom parallel currents such as dense bottom currents (Ezer and Mellor (2004)). A partial solution to this problem has been proposed by Beckmann and Döscher (1997) by coupling an empirical terrain-following bottom boundary layer model to the geopotential

model grid, allowing for bottom-parallel down-slope flow. Another approach to improve representation of along-bottom flow has been proposed by Adcroft et al. (1997): to introduce *shaved cells* at the bottom to better fit the bottom topography. Geopotential coordinates over sloping topography have one other inherent problem: they do not allow a high vertical resolution near the bottom. A resolution of the bottom layer of 1 m would require equidistant geopotential coordinates over the entire water column, which is often computationally not affordable.

Given all these disadvantages of geopotential coordinates they can generally not be recommended for coastal ocean simulations. There is however one potential problem of vertical coordinates which is not relevant for geopotential coordinates: the internal pressure gradient problem, see Sec. 7.2. Therefore, models with geopotential coordinates are still applied for problems with strong stratification such as the Baltic Sea (Wang et al. (2015)).

The numerical treatment of geopotential coordinates fits into the concept of general vertical coordinates by formally allowing layers of zero thickness at the bottom and, in the case of removable surface layers, also at the surface.

4.2.2. Terrain-following coordinates

A transformation function of the form $z(s) - \eta = S(x, y, s, t)D(x, y, t)$ with the local water depth $D = \eta + H$, the stretching function S being strictly monotone in s, $S\left(s_{1/2}\right) = -1$ and $S\left(s_{k_{\max}+1/2}\right) = 0$ generates terrain-following coordinates with $z_{1/2} = -H$ and $z_{k_{\max}+1/2} = \eta$. In addition to the free surface being a coordinate surface (as for the z^*-coordinates), also the bottom is a coordinate surface. Thus, in every watercolumn $k_{\max}$ layers are active and masking of layers is not necessary. A classical option for coastal ocean model simulations are σ-coordinates defined by $S = s = \sigma$ (Nihoul et al. (1986); Blumberg and Mellor (1987)), see Figure 1a. They result in layer thicknesses proportional only to the local water depth, which is a disadvantage especially in deeper waters (where layers may become too thick and thus resolution in critical parts of the water column too coarse) or above steep bottom slopes (where coordinate slopes are relatively steep throughout the water column, see Deleersnijder and Beckers (1992)). One other problem of terrain-following coordinates is the intersection of sloping coordinate surfaces with isopycnals. For the calculation of the internal pressure gradient, this results in a balance of two large terms which typically results in large truncation errors and thus artificial currents (see Sec. 7.2 for a detailed discussion.

Some of these problems can be reduced by using non-equidistant σ-levels $\sigma_{k+1/2}$ (equivalent to non-linear stretching functions $S(s)$) and modified stretching functions $S(x, y, s)$ that rescale the near-surface layer heights proportional to a fixed critical water depth instead of to the local water depth (Song and Haidvogel, 1994; Shchepetkin and McWilliams, 2009), see Figure 1b.

An alternative transformation function that maintains high resolution at the surface or at the bottom independently of the local water depth has been developed by Burchard and Petersen (1997) and Burchard and Bolding (2002).

On order to obtain higher accuracy, hybridisations between s- and z-coordinates have been proposed, leading to less steep coordinate slopes than for s coordinates and smaller bottom steps than for z-coordinates, see Gerdes (1993); Burchard and Petersen (1997) for details.

4.2.3. Arbitrary Lagrangian-Eulerian (ALE) coordinates

For geopotential and terrain-following coordinates the transformation function $z(s)$ is explicitly prescribed. The temporal evolution of the corresponding interface positions $z_{k+1/2}(x, y, t)$ and layer heights $h_k(x, y, t)$ is fully determined by a prognostic equation for the sea surface elevation $\eta(x, y, t)$ (see Sec. 5.1). In this case the continuity equation (29) does not serve as a prognostic equation for h_k, but as a diagnostic equation for the grid-related vertical velocity $w^{\mathrm{s}}_{k+1/2}$. Following Adcroft and Hallberg (2006), this is the Eulerian treatment of the vertical direction (EVD).

In contrast, the Lagrangian treatment of the vertical direction (LVD) is based on a Lagrangian vertical coordinate, formally defined by $\frac{\mathrm{D}s}{\mathrm{D}t} = 0$ (Starr, 1945). In this case the grid-related vertical velocity vanishes $w^{\mathrm{s}} = 0$, see (28), and the continuity equation (29) serves as a prognostic equation for the layer heights h_k.

Numerical mixing (induced by truncation errors of discrete advective fluxes; see e.g. Klingbeil et al., 2014) is significantly reduced in Lagrangian vertical coordinates because of the missing advective transport through the coordinate surfaces. However, a pure Lagrangian movement of the layers is prone to grid distortion and possibly vanishing layers, which complicates the numerical treatment.

Therefore, a combination of LVD and EVD gains increasingly popularity. Originally Hirt et al. (1974) defined the Arbitrary Lagrangian-Eulerian (ALE) method by a Lagrangian mesh movement (LVD step) followed by

an optional readjustment of the mesh associated with grid-related advective transports (EVD step). The advection during the EVD step can be replaced by instantaneous "remapping" (monotone and conservative interpolation, see e.g. White and Adcroft, 2008) which relaxes the strict CFL constraint to a less severe Lipschitz condition.

Another approach is to consider Lagrangian tendencies in the calculation of the layer heights used in EVD algorithms. Separable terrain-following coordinates of the form $z - \eta = S(s)D(x, y, t)$ inherently imply a Lagrangian treatment of merely barotropic motions (Shchepetkin and McWilliams, 2009). Non-separable terrain-following coordinates with time-independent stretching functions $S(x, y, s)$ do not exactly treat barotropic motions in a Lagrangian way, but also significantly reduce the corresponding grid-related vertical transports by scaling the layer heights according to $h_k = \frac{D}{H} h_k^{(\eta=0)}$. The same holds for the z^*-coordinates of Adcroft and Campin (2004).

Recently, Leclair and Madec (2011) proposed to also treat fast oscillating baroclinic currents in a Lagrangian way by introducing the so-called $\tilde{z}$-coordinate.

4.2.4. Adaptive coordinates

In order to further optimise the vertical grid layout in 3D models, Hofmeister et al. (2010) presented adaptive coordinates, see Figure 1c. These do not only consider Lagrangian tendencies, but also a zooming of layers towards boundaries, stratification and shear. This zooming is based on a vertical diffusion equation for the interface positions,

$$\partial_t z_{k+1/2} - \left(D^{\mathrm{grid}}_{k+1}(z_{k+3/2} - z_{k+1/2}) - D^{\mathrm{grid}}_{k}(z_{k+1/2} - z_{k-1/2}) \right) = 0, \qquad (37)$$

with boundary conditions $z_{1/2} = -H$ and $z_{k_{\max}+1/2} = \eta$ (Burchard and Beckers, 2004). A vertically constant layer "diffusion" coefficient D^{grid} (unit s^{-1}) tends to generate equidistant layers, whereas locally increased coefficients (e.g. proportional to stratification and shear) increase the local concentration of layer interfaces and thus the vertical resolution. Improved vertical resolution decreases the truncation errors of vertical advective fluxes and thus reduces vertical numerical mixing. The layer distribution in adjacent water columns is coupled by lateral adjustment of layer heights and interface positions. In addition, an optional isopycnal tendency aligns the layers with isopycnals and thus avoids the spurious diapycnal mixing of tracers due

to explicit and numerical along-layer mixing in the model. Furthermore, an isopycnal alignment of the layers reduces internal pressure gradient errors (see Sec. 7.2). It should be noted, that pure isopycnal coordinates are not appropriate for coastal ocean models, because of the relatively large mixed layers according to (1). Gräwe et al. (2015) demonstrated the clear advantages of adaptive coordinates compared to σ-coordinates in a realistic state-of-the-art model application with different dynamical regimes (tidal/non-tidal parts, seasonal thermocline / permanent halocline).

4.3. Horizontal discretisation

The layer-integrated equations (29) - (33) include horizontal divergence terms for advection and diffusion which still need to be discretised. Independently from the vertical coordinate specification, a horizontal coordinate transformation is carried out. For the simplest case, no coordinate transformation is needed, such that horizontal coordinates remain Cartesian, with typically x denoting the coordinate to the right and y denoting the coordinate orthogonally to it in a right-hand system. Such Cartesian coordinates are consistent with UTM coordinates. This however limits model applications to domains being very small compared to the Earth's radius, since the model domain is assumed to be planar. For larger domains (such as shelf seas), a transformation into spherical coordinates is carried out, with coordinates λ (pointing eastwards) and ϕ (pointing northwards), which is consistent with geographical coordinates. The transformation to spherical coordinates results in geometric coefficients by which terms in the dynamic equations are weighted (see, e.g., the text book by Kantha and Clayson (2000a)). To support multi-scale applications also with regular coordinates (and not only with unstructured-grid models), orthogonal curvi-linear coordinates have been introduced (Haidvogel et al. (1991)), having enough flexibility to focus into certain parts of the domain such as estuaries or coastal regions. Curvi-linear coordinates may be referenced to Cartesian or to spherical coordinates, depending on the scale of the problem, but the resulting grids (after discretisation) are of the same type. It should be noted that Cartesian and spherical coordinates can be regarded as special cases of orthogonal curvi-linear coordinates, such that models which include the latter do automatically offer the numerics for the other two types. Note that also non-orthogonal curvilinear coordinates are in use, to further increase flexibility in spatial resolution, but those have not become common in coastal ocean modelling (Kantha and Clayson (2000a)).

After the horizontal transformation, a horizontal discretisation is carried out, by dividing the two coordinates of the Cartesian, spherical or orthogonal curvi-linear coordinates into not necessarily equidistant spatial increments. This leads to a tiling of the horizontal model domain by grid cells which may be rectangles (Cartesian coordinates), trapezoids (spherical coordinates) or general quadruples (orthogonal curvi-linear coordinates). Note that angles between edges are orthogonal only for Cartesian coordinates. Each grid cell has thus four corners and four edges. Grid cells which include mainly land areas are then masked out.

The layer-integrated equations (29) - (33) are then area-integrated over these cells, such that in combination with the vertical layer-integration an integration over so-called finite volumes is obtained. By means of this volume integration, divergence terms are transformed to differences of interface-integrated advective and diffusive fluxes, such that advection and diffusion are discretised in a conservative way. Non-conservative source and sink terms appear as volume-integrals. These finite volumes have 8 corners, 12 edges and 6 interfaces, see Fig. 2. With $\Delta x_{i,j}$ and $\Delta y_{i,j}$ denoting mean length and width of the volumes and $h_{i,j,k}$ denoting the mean height, these volumes are defined as $\mathcal{V}_{i,j,k} = \Delta x_{i,j} \Delta y_{i,j} h_{i,j,k}$. With $\Delta x_{i,j\pm 1/2}$ and $\Delta y_{i\pm 1/2,j}$ denoting the length and width at the lateral edges, the vertical interfaces of the finite volume have the areas $A_{i\pm 1/2,j,k} = \Delta y_{i\pm 1/2,j} h_{i\pm 1/2,j,k}$ and $A_{i,j\pm 1/2,k} = \Delta x_{i,j\pm 1/2} h_{i,j\pm 1/2,k}$. With these notations, the volume-integral of the layer height equation (29) reads as

$$
\begin{aligned}
&\Delta x_{i,j} \Delta y_{i,j} \partial_t h_{i,j,k} \\
&\quad + \left(\Delta y_{i+1/2,j} h_{i+1/2,j,k} u_{i+1/2,j,k} - \Delta y_{i-1/2,j} h_{i-1/2,j,k} u_{i-1/2,j,k} \right) \\
&\quad + \left(\Delta x_{i,j+1/2} h_{i,j+1/2,k} v_{i,j+1/2,k} - \Delta x_{i,j-1/2} h_{i,j-1/2,k} v_{i,j-1/2,k} \right) \\
&\quad + \Delta x_{i,j} \Delta y_{i,j} \left(w^{\mathrm{s}}_{i,j,k+1/2} - w^{\mathrm{s}}_{i,j,k-1/2} \right) = 0,
\end{aligned}
\tag{38}
$$

with $u_{i\pm 1/2,j,k}$, $v_{i,j\pm 1/2,k}$, and $w^{\mathrm{s}}_{i,j,k\pm 1/2}$, denoting area-averaged velocities at the respective interfaces. Tracer mass conservation requires the finite-volume

discretisation of e.g. the salinity equation of the following form:

$$
\begin{aligned}
&\Delta x_{i,j}\Delta y_{i,j}\partial_t h_{i,j,k}S_{i,j,k} \\
&\quad +\big(\Delta y_{i+1/2,j}h_{i+1/2,j,k}u_{i+1/2,j,k}\widetilde{S}_{i+1/2,j,k} \\
&\qquad -\Delta y_{i-1/2,j}h_{i-1/2,j,k}u_{i-1/2,j,k}\widetilde{S}_{i-1/2,j,k}\big) \\
&\quad +\big(\Delta x_{i,j+1/2}h_{i,j+1/2,k}v_{i,j+1/2,k}\widetilde{S}_{i,j+1/2,k} \\
&\qquad -\Delta x_{i,j-1/2}h_{i,j-1/2,k}v_{i,j-1/2,k}\widetilde{S}_{i,j-1/2,k}\big) \\
&\quad +\Delta x_{i,j}\Delta y_{i,j}\big(w^{\mathrm{s}}_{i,j,k+1/2}\widetilde{S}_{i,j,k+1/2}-w^{\mathrm{s}}_{i,j,k-1/2}\widetilde{S}_{i,j,k-1/2}\big)= \\
&\quad -\Delta y_{i+1/2,j}\left(h_{i+1/2,j,k}J^{\mathrm{S}}_{\tilde{x},i+1/2,j,k}\right)+\Delta y_{i-1/2,j}\left(h_{i-1/2,j,k}J^{\mathrm{S}}_{\tilde{x},i-1/2,j,k}\right) \\
&\quad -\Delta x_{i,j+1/2}\left(h_{i,j+1/2,k}J^{\mathrm{S}}_{\tilde{y},i,j+1/2,k}\right)+\Delta x_{i,j-1/2}\left(h_{i,j-1/2,k}J^{\mathrm{S}}_{\tilde{y},i,j-1/2,k}\right) \\
&-\Delta x_{i,j}\Delta y_{i,j}\left(J^{\mathrm{S}}_{z,i,j,k+1/2}-J^{\mathrm{S}}_{z,i,j,k-1/2}\right),
\end{aligned}
\tag{39}
$$

with the interfacial salinity values $\widetilde{S}_{i\pm1/2,j,k}$, $\widetilde{S}_{i,j\pm1/2,k}$ and $\widetilde{S}_{i,j,k\pm1/2}$ and the volume averaged salinity $S_{i,j,k}$. For constant tracers and zero turbulent tracer fluxes the tracer equation must recover (38).

The natural spatial discretisation which results from the volume-integration carried out in (38) and (39) suggests to locate the discrete velocity points $u_{i\pm1/2,j,k}$, $v_{i,j\pm1/2,k}$, and $w^{\mathrm{s}}_{i,j,k\pm1/2}$ at the centres of the 6 interfaces of the finite volume. The natural location for the volume-averaged salinity $S_{i,j,k}$ (and consequently also of the potential temperature $\theta_{i,j,k}$ and the buoyancy $b_{i,j,k}$) would then be the centre of the finite volume, see Fig. 2. This would be in accordance with the classical Arakawa C-grid (Arakawa and Lamb (1977)), where velocity points and *pressure points* (salinity, potential temperature, buoyancy, and thus also pressure) are staggered in space. Such a grid layout guarantees that (38) and (39) can be discretised at second order in space without any interpolations of velocity components. For the volume-integrated u- and v-velocity equations (not shown here; note that for the w-velocity no dynamic equation has to be integrated) separate (staggered)

finite volumes have to be constructed for which transport velocities have to be interpolated to the respective interfaces (which has to be done for any staggered or non-staggered grid layout). Turbulent quantities (such as vertical eddy viscosity/diffusivity and turbulent kinetic energy) are best located at the w-points, such that also for the turbulence equations specific finite volumes centred around the w-points need to be constructed. There is one disadvantage of the fully staggered grid layout: The respectively other horizontal velocity component for the calculation of the Coriolis terms needs to be interpolated from four adjacent velocity points in order to be second-order in space. A grid layout which avoids this by collocating u- and v-components at the corners of the finite volume is given by the B-grid (which still is staggered with respect to vertical velocity and pressure points). Therefore large scale ocean models for which an accurate spatial discretisation is essential typically use the B-grid, whereas coastal ocean models have traditionally been written as C-grid models (see also Sec. 8.1). An excellent overview over the discretisation of the linearised shallow water equations and its accuracies in A- (all discrete points collocated), B- (only horizontal velocity points collocated), C- (all staggered), and D-grids (as C-grid, but with u- and v-velocity locations exchanged) is given by Beckers and Deleersnijder (1993).

5. Choices for mode splitting and wetting & drying

An essential feature of coastal ocean models is the consideration of a temporally and spatially variable free surface, see (2) in section 2.2. Due to the presence of a free surface, fast surface gravity waves are part of the numerical solution, and drying and flooding in shallow areas of the domain is possible. Both issues require special numerical techniques that will be outlined in the following two subsections.

5.1. Mode splitting

Within the EVD-treatment (see Section 4.2.3) the layer-integrated continuity equation (29) serves as a diagnostic equation for the grid-related vertical velocities $w^{\mathrm{s}}_{k+1/2}$, and the layer heights h_k are determined in terms of the total water depth D and the free surface elevation η. In this case a prognostic equation for the free surface elevation must be derived from the vertical sum of the layer-integrated continuity equation (29) and consideration of the kinematic boundary conditions (14a) and (14b):

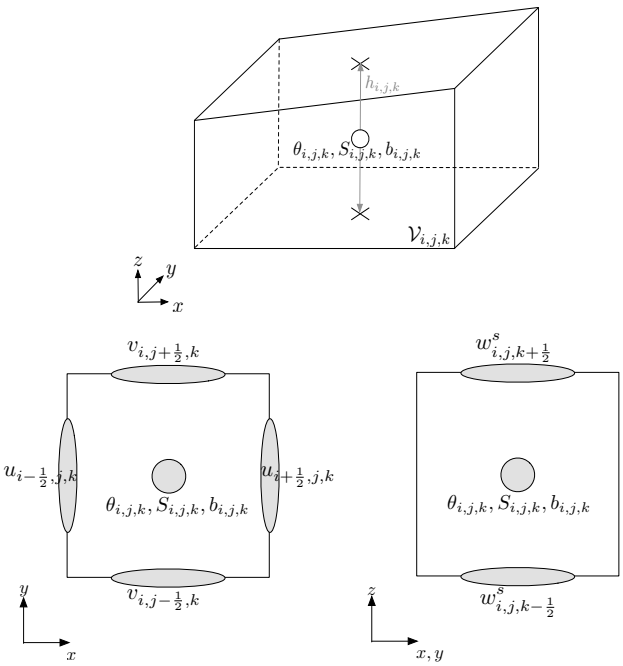

Figure 2: Control volume for cell-centered variables (top), C-grid staggering showing the location of the dependent variables in the horizontal (bottom, left) and in the vertical (bottom, right).

$$\partial_t(\eta{+}H) = \partial_t D = \partial_t \left(\sum_{k=1}^{k_{\max}} h_k \right) = -\partial_x \left(\sum_{k=1}^{k_{\max}} h_k u_k \right) - \partial_y \left(\sum_{k=1}^{k_{\max}} h_k v_k \right) - (E{-}P). \tag{40}$$

The direct numerical integration of the set (40) and (29)–(33) is possible. However, within an explicit temporal discretisation the time step is strongly constrained by the celerity of the fastest surface gravity waves. Alternatively, an implicit discretisation offers larger time steps (Casulli and Cheng, 1992), but also requires the inversion of a penta-diagonal system for the free surface elevation.

Since surface gravity waves are solely represented by the barotropic mode, the computational costs can be significantly reduced by splitting the governing equations into the barotropic and remaining baroclinic mode. Usually the slightly depth-dependent barotropic mode is approximated by the depth-integrated equations (also called external mode), given by (40) and the vertical sum of the layer-integrated momentum equations (30) and (31):

$$\partial_t D + \partial_x \left(D\bar{u} \right) + \partial_y \left(D\bar{v} \right) = -(E - P), \tag{41a}$$

$$\begin{aligned} &\partial_t \left(D\bar{u} \right) + \partial_x \left(D\bar{u}\bar{u} \right) + \partial_y \left(D\bar{v}\bar{u} \right) \\ &= fD\bar{v} - gD\partial_x \eta - D\tfrac{1}{\rho_0}\partial_x p_a + \tau_{\mathrm{x}}^{\mathrm{s}}(\bar{u},\bar{v}) - \tau_{\mathrm{x}}^{\mathrm{b}}(D,\bar{u},\bar{v}) + D\widetilde{F}_{\mathrm{x}}^{\mathrm{h}}(D,\bar{u},\bar{v}) + S_{\mathrm{x}}, \end{aligned} \tag{41b}$$

$$\begin{aligned} &\partial_t \left(D\bar{v} \right) + \partial_x \left(D\bar{u}\bar{v} \right) + \partial_y \left(D\bar{v}\bar{v} \right) \\ &= -fD\bar{u} - gD\partial_y \eta - D\tfrac{1}{\rho_0}\partial_y p_a + \tau_{\mathrm{y}}^{\mathrm{s}}(\bar{u},\bar{v}) - \tau_{\mathrm{y}}^{\mathrm{b}}(D,\bar{u},\bar{v}) + D\widetilde{F}_{\mathrm{y}}^{\mathrm{h}}(D,\bar{u},\bar{v}) + S_{\mathrm{y}}, \end{aligned} \tag{41c}$$

with the water depth,

$$D \equiv \sum_{k=1}^{k_{\max}} h_k, \tag{42}$$

the barotropic transports,

$$[D\bar{u}, D\bar{v}] \equiv \sum_{k=1}^{k_{\max}} [h_k u_k, h_k v_k], \tag{43}$$

and the barotropic-baroclinic interaction terms:

$$
\begin{aligned}
S_{[\mathrm{x,y}]} = \quad & \partial_x \left(D\bar{u}[\bar{u},\bar{v}] \right) && - \sum_{k=1}^{k_{\max}} \partial_x \left(h_k u_k[u_k,v_k] \right) \\
& + \partial_y \left(D\bar{u}[\bar{u},\bar{v}] \right) && - \sum_{k=1}^{k_{\max}} \partial_y \left(h_k u_k[u_k,v_k] \right) \\
& - D\widetilde{F}^{\mathrm{h}}_{[\mathrm{x,y}]}(D,\bar{u},\bar{v}) && + \sum_{k=1}^{k_{\max}} h_k \widetilde{F}^{\mathrm{h}}_{[\mathrm{x,y}]}(h_k,u_k,v_k) \\
& - \tau^{\mathrm{s}}_{[\mathrm{x,y}]}(\bar{u},\bar{v}) && + \tau^{\mathrm{s}}_{[\mathrm{x,y}]}(u_{k_{\max}},v_{k_{\max}}) \\
& + \tau^{\mathrm{b}}_{[\mathrm{x,y}]}(D,\bar{u},\bar{v}) && - \tau^{\mathrm{b}}_{[\mathrm{x,y}]}(h_1,u_1,v_1) \\
& && + \sum_{k=1}^{k_{\max}} h_k F^{\mathrm{IP}}_{[\mathrm{x,y}],k}. \qquad (44)
\end{aligned}
$$

The remaining 3D equations (30)–(33), as well as the turbulence closure equations (section 3.2) and other tracer equations such as biogeochemical model equations represent the internal (baroclinic) mode. With (42), (43) and (44) the splitting into external and internal mode equations is exact. But, as mentioned above, it does not exactly separate the slightly depth-dependent fast barotropic dynamics from the slow baroclinic ones. Therefore, the time-stepping of the internal mode might still be constrained by some fast dynamics.

A careful stability analysis of the linearised (flat-bottom) version of the external mode (41a)–(41c),

$$
\partial_t \eta = -H \left(\partial_x \bar{u} + \partial_y \bar{v} \right); \quad \partial_t \bar{u} = -g\partial_x \eta + f\bar{v}; \quad \partial_t \bar{v} = -g\partial_y \eta - f\bar{u}, \qquad (45)
$$

has been carried out by Beckers and Deleersnijder (1993). For an explicit discretisation on an equidistant C-grid, the stability criterion would be $\Delta t \leq \Delta x / \left(2gH\right)^{1/2}$, which means for $\Delta x = 1000\mathrm{m}$ and a depth of $H = 200\mathrm{m}$ a maximum timestep of $\Delta t \approx 5\mathrm{s}$. Compared to that, the stability criterion for discretising the internal mode is typically given by the advection terms. For an explicit one-stage two-level discretisation (Durran, 2013) of the linear advection equation the stability criterion is $\Delta t \leq \Delta x / u_{\max}$, which would for a typical maximum velocity of $u_{\max} = 1\mathrm{m\,s^{-1}}$ give a maximum timestep of $\Delta t = 200\mathrm{s}$, which is a factor of about 40 larger. Non-linear systems retaining all terms tend to have comparable timestep criteria.

If the surface elevation and the barotropic transports are not integrated prognostically as part of the internal mode, but provided by the external mode, an explicit temporal discretisation of the internal mode is not constrained by the celerity of the fastest surface gravity waves anymore, but only by the much slower internal motions. In this case the internal mode can be integrated explicitly with a timestep which, for coastal oceans, is $\mathcal{O}(10^1)$ times larger than that constrained by the celerity of shallow water surface gravity waves. Since this increased timestep is usually already almost required to resolve the dynamical processes in e.g. tidally dominated coastal oceans, the overhead of an implicit treatment of the 3D internal mode equations would be by far not balanced by the possible slight further increase of the timestep. Therefore, the prognostic integration of the internal mode is usually performed explicitly.

The prognostic integration of the external mode can be either carried out implicitly with the same timestep as used for the explicit integration of the internal mode, or explicitly within a series of smaller timesteps constrained by the celerity shallow water surface gravity waves. Both methods have their own merits and complications. Such a different numerical treatment of the two modes is called mode splitting and has first been introduced by Madala and Piacsek (1977).

5.1.1. Semi-implicit external mode computation

Let (41a)–(41c) being solved simultaneously, with the transports $D\bar{u}$ and $D\bar{v}$ in (41a) being temporally weighted such that they are between the levels of the old and the new surface elevation, and equivalently, with η in (41b) and (41c) being temporally between the old and new transports. then, a semi-implicit time stepping scheme for the external mode results. With the weighting parameter $0 \leq \gamma \leq 1$, this can be expressed as $X^{n+\gamma} = (1-\gamma)X^n + \gamma X^{n+1}$ for any quantity X which is treated semi-implicitly (with superscripts denoting the number of time step). For $\gamma = 0$, the scheme would be fully explicit, for $\gamma = 1$ the scheme would be fully implicit and for $\gamma = 0.5$, the Crank-Nicolson scheme is obtained. The schemes are more stable (and more implicit) the more the bias is towards the new time step, i.e., the larger γ is. This procedure gives a linear system of equations with a matrix of dimension $N \times N$ with N denoting the number of free-surface grid boxes. Depending on the treatment of the Coriolis terms (explicit or semi-implicit), and when calculating non-linear terms explicitly, each row of the matrix has 5 or 9 non-zero entries, such that the matrix is very sparse. One of the challenges

for solving these large systems of equations is to find accurate (typically iterative) solvers which are efficient also on massively parallel high-performance computers. An early example for a semi-implicit coastal ocean model including semi-implicit treatment of the Coriolis term has been presented by Backhaus (1985). This semi-implicit scheme is unconditionally stable for $\gamma \geq 0.5$.

To avoid the requirement of solving large linear systems of equations due to the semi-implicit treatment of the external mode, a directional-split method had been introduced by Peaceman and Rachford (1955), in which in an alternating manner the external mode is solved along one and then along the other coordinate direction. This method is called Alternating Directions Implicit (ADI). Many models in coastal engineering had adopted this approach (see, e.g., Pietrzak et al., 2002). While ADI is easier to handle computationally than a semi-implicit treatment without directional splitting, additional errors due to the splitting have to be considered.

The advantage of the semi-implicit treatment is the relatively easy coupling between the external and the internal mode, since both are computed with the same time step. A disadvantage of implicit schemes is the demanding computation of large and sparse linear system on parallel computers. While the Crank-Nicolson scheme is energy-conserving (but still sensitive to instabilities arising from the discretisation of other terms), a larger degree of implicitness ($\gamma \rightarrow 1$) stabilises the scheme, a procedure which however for larger time steps might become too dissipative (Walters et al. (2009)). In tidal or tsunami simulations this sets a practical limitation to the overall model time step.

5.1.2. Explicit external mode computation

In order to provide a reasonable free surface elevation and barotropic transports for updating the internal mode from stage ^{n} to stage $^{n+1}$, the external mode is integrated explicitly within a subcycle of $M_{\max}$ timesteps from stage ^{n,0} to stage $^{n,M_{\max}}$. The timestep of the external mode $\triangle t_{2\mathrm{D}}$ is constrained by the celerity of shallow water surface gravity waves, whereas the internal mode is integrated with a larger internal timestep $\triangle t_{3\mathrm{D}} = M \triangle t_{2\mathrm{D}}$ with $M \leq M_{\max}$.

It is easy and accurate to discretise the external mode with an explicit-in-time solver. In the following a time-staggered two-level scheme is outlined. Staggering in time between the surface elevations and the barotropic transports would still guarantee second-order in time accuracy. This would be

achieved by subsequently calculating the surface elevation and the barotropic transports with using the newly updated transports and elevations on the right hand sides, see section 7.3. In the following the focus will be on volume and tracer mass conservation. Therefore, the discretisation of the momentum equations will be not presented in detail. A time-staggered discretisation of the free-surface equation (41a) is given by:

$$\frac{\eta^{n,m+1}-\eta^{n,m}}{\triangle t_{2\mathrm{D}}}+\partial_x\left(D\bar{u}\right)^{n,m+1/2}+\partial_y\left(D\bar{v}\right)^{n,m+1/2} = -\left(E-P\right)^{n,m+1/2} \quad (46)$$

Following Shchepetkin and McWilliams (2005) a filtered surface elevation is provided to the internal mode,

$$\langle\eta\rangle^{n+1}\equiv\sum_{m=0}^{M_{\max}}a_m\eta^{n,m}, \quad (47)$$

with normalised filter weights $\sum_{m=0}^{M_{\max}}a_m=1$. If the initial surface elevation of each external mode cycle is reset to $\eta^{n,0}\equiv\langle\eta\rangle^n$, the compatible surface elevation equation resolved by the internal mode can be derived from (46) as

$$\frac{\langle\eta\rangle^{n+1}-\langle\eta\rangle^{n}}{\triangle t_{3\mathrm{D}}}+\partial_x\left\langle\langle D\bar{u}\rangle\right\rangle^{n+1/2}+\partial_y\left\langle\langle D\bar{v}\rangle\right\rangle^{n+1/2} = -\left\langle\langle E-P\rangle\right\rangle^{n+1/2} \quad (48)$$

with

$$\left\langle\langle\psi\rangle\right\rangle^{n+1/2}\equiv\sum_{m=0}^{M_{\max}-1}b_m\psi^{n,m+1/2} \quad (49)$$

and $b_m\equiv\frac{1}{M}\sum_{m'=m+1}^{M_{\max}}a_{m'}$. The general filter applied in (47) offers different types of filtering. Deleersnijder (1993) proposed to provide the instantaneous surface elevation $\langle\eta\rangle^{n+1}=\eta^{n,M}$ and averaged barotropic transports $\left\langle\langle[D\bar{u},D\bar{v}]\rangle\right\rangle^{n+1/2}=\sum_{m=0}^{M-1}\left[D\bar{u},D\bar{v}\right]^{n,m+1/2}$ to the internal mode. This treatment corresponds to $M_{\max}=M$, $a_m=\delta_{mM}$, $b_m=1$ and involves no reset of the external quantities, i.e. $\eta^{n+1,0}=\eta^{n,M}$ and $\left[D\bar{u},D\bar{v}\right]^{n+1,0-1/2}=\left[D\bar{u},D\bar{v}\right]^{n,M-1/2}$. In order to stabilise the retained fast dynamics within the

internal mode and to prevent aliasing errors Griffies (2007) suggested to provide an averaged surface elevation $\langle\eta\rangle^{n+1} = \frac{1}{2M+1}\sum_{m=0}^{2M}\eta^{n,m}$ to the internal mode, obtained by $M_{\max} = 2M$, $a_m = \frac{1}{2M+1}$ and $b_m = \frac{2M-m}{M(2M+1)}$. Non-uniform filter weights, centered around $m = M < M_{\max}$, were presented by Shchepetkin and McWilliams (2005) to offer more frequency-selective filtering.

Volume conservation requires that the layer-integrated continuity equation (29) in the internal mode,

$$\frac{h_k^{n+1} - h_k^n}{\triangle t_{3\mathrm{D}}} + \partial_x \left(h_k u_k\right)^{n+1/2} + \partial_y \left(h_k v_k\right)^{n+1/2} + \left(w_{k+1/2}^{\mathrm{s},n+1/2} - w_{k-1/2}^{\mathrm{s},n+1/2}\right) = 0, \tag{50}$$

recovers (48) when vertically summed. Thus

$$\sum_{k=1}^{k_{\max}} h_k^{n+1} \equiv \langle\eta\rangle^{n+1} + \langle H\rangle^{n+1} = \langle D\rangle^{n+1}, \tag{51a}$$

$$\sum_{k=1}^{k_{\max}} \left[h_k u_k, h_k v_k\right]^{n+1/2} \equiv \langle\langle [D\bar{u}, D\bar{v}]\rangle\rangle^{n+1/2} \tag{51b}$$

$$w_{k_{\max}+1/2}^{\mathrm{s},n+1/2} \equiv \langle\langle E - P\rangle\rangle^{n+1/2}. \tag{51c}$$

Compatibility condition (51b) can be obtained by shifting the velocity profiles obtained from the baroclinic momentum equations (30) and (31). It should be noted again, that for tracer mass conservation the discrete tracer equations must recover (50) for constant tracers. Therefore, the shifting of velocity profiles must be done before the grid-related vertical velocity is diagnosed from (50) and the tracer equations are integrated. At the end of each internal timestep, the interaction-terms (44) can be updated in terms of $\langle\langle [D\bar{u}, D\bar{v}]\rangle\rangle^{n+1/2}$, $\left[h_k u_k, h_k v_k\right]^{n+1/2}$ and $F_{[\mathrm{x,y}],k}^{\mathrm{IP},n+1}$. Since these interaction-terms are held constant over the next external subcycle, integrating the barotropic transports from $(D\bar{u})^{n+1,0-1/2}$ to $(D\bar{u})^{n+1,M_{\max}-1/2}$, a time lag is introduced.

5.2. Drying and flooding

In many coastal areas intertidal flats are an important topographic feature. During high water those areas are inundated and during ebb tide they may fall dry, approaching zero water depth $D \to 0$, but never becoming negative. To understand how this is reflected in the mathematical formulations for

coastal ocean dynamics, it is instructive to reformulate the depth-integrated continuity equation (41a) into an advection equation for the water depth:

$$\partial_t D + \bar{u}\partial_x D + \bar{v}\partial_y D = -\left(\partial_x \bar{u} + \partial_y \bar{v}\right) D - (E - P). \tag{52}$$

The conservative form (41a) shows that total water volume is conserved. The advective form (52) demonstrates positivity of the water depth, because its total derivative depends on a sink term which vanishes for decaying water depth. The divergence stays bounded, because for frictionally dominated regimes the depth-integrated momentum equations (41b) and (41c), and the quadratic bed friction (77) guarantee bounded depth-mean velocities. The evaporation E which reduces the water depth must be numerically treated such that it vanishes for vanishing water depth.

Various ways have been proposed in coastal ocean models to guarantee the two major properties (i) volume conservation (for incompressible flow) according to (41a) and (ii) positivity according to (52), also numerically.

It should be straight-forward to directly discretise one of the equations. Usually (41a) is used as a basis. In contrast, for a vertically integrated model, Stelling and Duinmeijer (2003) successfully applied a first-order upstream scheme to discretise the advection equation (52), and later Zijlema et al. (2011) postulated that positivity could also be achieved using a Total Variation Diminishing (TVD) scheme with the MUSCL limiter (van Leer (1979)).

Given that the spatial discretisation is for a C-grid (section 4.3), and in order to conserve volume locally and globally, the transports $\bar{u}D$ and $\bar{v}D$ need to be identical to the sum of the layer-integrated transports used in (29). This consistency of local and global volume conservation is also essential for the conservation of momentum and tracers. Therefore, a numerical recalculation of $\bar{u}D$ and $\bar{v}D$ in (41a) would require a correction to the layer-integrated transports calculated prognostically in (30) and (31). This is the general principle of conservative drying and flooding: to manipulate the layer-integrated momentum equations in a physically sound and reasonable way, and to use the same transports in (41a).

There are various approaches to this, all of them leaving a thin film of water in dry cells, where the dynamics is manipulated in some way to avoid zero or negative water depth. In early coastal ocean models (Flather and Heaps, 1975; Backhaus, 1976, 1985; Jungclaus and Backhaus, 1994), a number of rules were defined on how to reduce fluxes out of a grid cell when

the application of these fluxes would lead to negative water depth (or the undercut of a prescribed minimum water depth). Also the wetting and drying schemes of Oey (2005) and Warner et al. (2013) are based on this principle. The method works reliably, but to guarantee positive water depth, it has to be taken into account that reduction or shut-off of fluxes around one grid cell would change the net fluxes into the neighboring cells, such that an iterative procedure might become necessary. This problem has been partially solved by Warner et al. (2013), by shutting off all transports around a cell when the water depth at the old time step is below the critical depth, see also Defne and Ganju (2015) who simulated a large lagoonal back-barrier estuary at the US east coast including extensive intertidal flats. Moreover, the direct correction of fluxes might in some models lead to spatially and temporally oscillating patterns of dry and wet cells. Furthermore, shutting on and off transports instantaneously could lead to complications in turbulence closure models, since the bed stress which largely determines the eddy viscosity profiles (via the turbulent kinetic energy and the macro length scale of turbulence, see section 3.2.1) would be suddenly switched on or shut off regularly.

In some models (e.g., Casulli and Cattani, 1994) the water depth at the transport points of a C-grid is chosen as the minimum depth between the adjacent pressure points, with the consequence that velocities feel the shallow water and bed friction strongly reduces the volume transport. This however might lead to significantly delayed drying and flooding processes (Burchard and Bolding, 2002).

In order to avoid explicit switching on or off of transports in drying cells, Burchard et al. (2004) (see also Burchard and Bolding, 2002) proposed the following measures when calculating the layer-integrated momentum equations:

- A critical water depth D_{crit}, and a minimum water depth, D_{min} are defined with $D_{\mathrm{crit}} > D_{\mathrm{min}}$ (see fig. 3). Typical values would be $D_{\mathrm{crit}} = 0.2\,\mathrm{m}$ and $D_{\mathrm{min}} = 0.05\,\mathrm{m}$. To simplify the dynamics in very shallow water with $D < D_{\mathrm{crit}}$, many terms in the layer-integrated momentum equations are reduced by a factor linearly decreasing between 1 (for $D = D_{\mathrm{crit}}$) and 0 (for $D = D_{\mathrm{min}}$), which basically reduces the local dynamics to an external pressure gradient - friction balance. Such a balance supports reduction of local transports, since the friction coefficient calculated from the law of the wall in the lowest grid box increases towards infinity for the water depth converging towards zero, see equa-

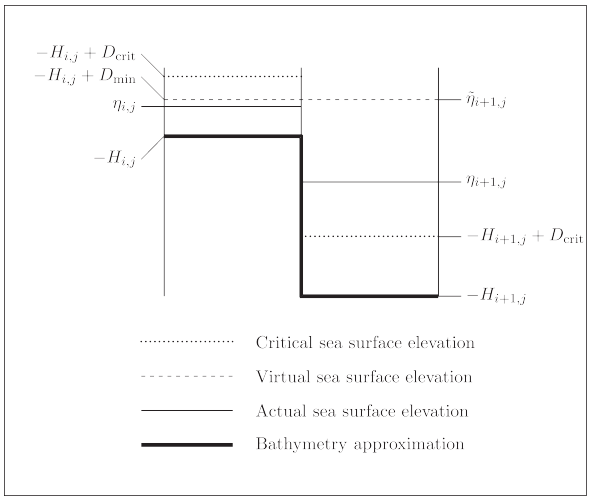

Figure 3: Sketch explaining the geometry of bathymetry and surface elevation approximations as it is defined in GETM. This figure has been modified from Burchard et al. (2004).

tion (76).

- Additionally, when the surface elevation in one grid box is below the bottom coordinate plus the minimum depth in any adjacent cell (which often happens at steep bathymetry in shallow water and relatively coarse horizontal resolution, see fig. 3, where $\eta_{i+1,j} < -H_{i,j} + D_{\mathrm{min}}$), a virtual surface elevation is defined from which the external pressure gradient $(\partial_x \eta, \partial_y \eta)$ is calculated. In that case, $\tilde{\eta}_{i+1,j} = \max\{\eta_{i+1,j}, -H_{i,j} + D_{\mathrm{min}}\}$, and the surface elevation gradient is calculated as $(\tilde{\eta}_{i+1,j} - \eta_{i,j})/\Delta x$. In the extreme case shown in fig. 3 water volume would actually flow against the elevation gradient, because $\eta_{i,j} < -H_{i,j} + D_{\mathrm{min}}$.

- Despite the above measures, direct reduction or shut-off of transports surrounding a grid cell may still be necessary at singular locations. To stabilise the drying and flooding algorithm, increasing the minimum and critical water depth is often a preferred option.

Using these algorithms Gräwe et al. (2016) could carry out a stable and multi-annual simulation of the entire Wadden Sea of the South-Eastern North Sea at a horizontal resolution of 200 m.

6. Treatment of open boundary conditions and model nesting

Coastal ocean models focus on limited areas of interest and thus data are needed to force the models at open boundaries, see Section 2.5. The data can typically come from an external source (e.g. climatology, coarser resolution run on a wider domain) following a one-way (coarse → fine) approach. In comparison, in a two-way approach, the external data (a numerical model) feel the local coastal solution through a feedback term. Local and external models generally differ on several key points among them grid resolution (in space and time) and physics. Across the interface, inflow and outflow can occur. In the case of inflow conditions, the quality of the local solution is strongly dependent on the quality of external data while in the case of outflow, the main issue is to prevent wave reflection at the interface. A two-way approach helps in maintaining some consistency between local and external solutions, at least when the physics of the two models are close. In both one-way and two-way approaches, the design of the boundary conditions must be carefully done. Due to the complexity of the physics, the scale heterogeneity in space and time, ideal boundary conditions are not achievable and a reasonable choice relies on a compromise between accuracy and robustness. Section (6.1) briefly introduces the general framework. Section (6.2) treats the problem of model forcing in a one-way approach (open boundary conditions). Section (6.3) introduces two-way approaches. Some considerations on discrete approximations are given in Section (6.4).

6.1. General framework

We consider a domain Ω_{loc} embedded in a domain Ω_{ext} as shown in Figure 4. The local and external solutions, ϕ_{loc} and ϕ_{ext}, are integrated according to

$$\partial_t \phi_{\text{loc}} + L_{\text{loc}}(\phi_{\text{loc}}) = f_{\text{loc}} \text{ in } \Omega_{\text{loc}}, \quad \partial_t \phi_{\text{ext}} + L_{\text{ext}}(\phi_{\text{ext}}) = f_{\text{ext}} \text{ in } \Omega_{\text{ext}}.$$

L_{loc} and L_{ext} are spatial differential operators, f_{loc} and f_{ext} model forcings. ϕ is the vector of prognostic state variables. The models interact through a transition area Γ that, depending on the boundary conditions, may be restricted to a single interface γ.

Let Δx_{loc} and Δx_{ext} (resp. Δt_{loc} and Δt_{ext})) be the grid (resp. time) steps on Ω_{loc} and Ω_{ext}. Today, the availability of global/regional operational or reanalysis products allows to consider the case where typical values of time/grid

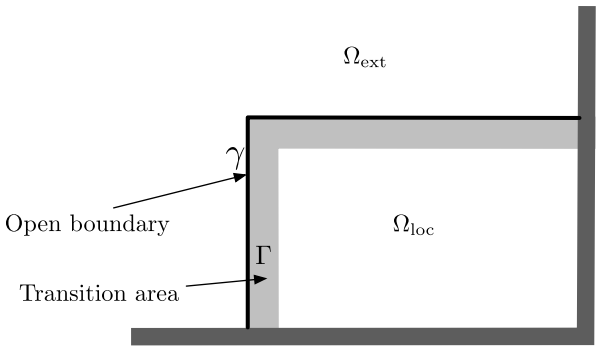

Figure 4: The local domain Ω_{loc} embedded in the exterior domain Ω_{ext}. Ω_{loc} has an artificial interface γ and possibly a transition area Γ on which boundary conditions are prescribed.

steps obey:

$$\frac{\Delta x_{\text{ext}}}{\Delta x_{\text{loc}}} \approx \frac{\Delta t_{\text{ext}}}{\Delta t_{\text{loc}}} \approx 2 - 10.$$

The model interactions across the interface γ and/or within the transition area Γ are described in the next Sections.

6.2. Open boundary conditions (OBC) and one-way interaction

We start with a description of the two main classes of OBC methods: relaxation methods which aim to relax the local solution towards the external solution within the transition area Γ and radiation methods which propagate out the information to prevent waves reflection. Simple radiation methods are appropriate only for a wave equation, the extension to general hyperbolic systems using a characteristic variables approach is explained in (6.2.2). Advanced open boundary conditions like Perfectly Matched Layer (PML) and Absorbing Boundary Conditions (ABC) (see Nataf, 2013) will not be discussed here since their derivations require very simplified problems and they have still to be validated in realistic applications. In (6.2.4), we discuss the specificities of local mesh refinement when the local computational grid is nested within a coarser resolution grid and also mention the impact

of variabilities at interfaces. For a review of open boundary conditions, the reader is referred to Palma and Matano (2000); Blayo and Debreu (2005).

6.2.1. Relaxation and radiation methods:

Relaxation methods. Basic Dirichlet (a.k.a. clamped) boundary conditions do not perform well (except for certain basic equations): this abrupt forcing generates inconsistencies and discontinuities near the interface γ. The flow relaxation scheme (FRS) (Davies, 1976) relaxes the interior solution towards external data in the transition area Γ near the boundary and can be written under the following relaxation (or nudging) form (Martinsen and Engedahl, 1987) (considering ϕ as a scalar variable and omitting the model forcing):

$$\partial_t \phi_{\text{loc}} + L_{\text{loc}}(\phi_{\text{loc}}) = -\frac{1}{\tau}(\phi_{\text{loc}} - \phi_{\text{ext}}) \quad \text{in } \Omega_{\text{loc}}, \tag{53}$$

The right hand side of (53) is the relaxation term. τ is a relaxation time scale. $\lambda = \dfrac{1}{\tau}$ is a positive function, which decreases away from the lateral boundary γ and vanishes outside of Γ. At the discrete level, the right hand side of (53) can be solved implicitly to remove associated stability constraints. It is difficult to prescribe the size of the transition area Γ and the variation of λ within Γ, since they depend on the space and time scales of the phenomena (Kållberg, 1977). Typical profiles of λ are based on quadratically or exponentially decreasing functions of the distance to the boundary γ.

Due to the smallness of the coastal model domain, it may also happen that the large scales of the solution are better reproduced in the external domain Ω_{ext}. It is particularly the case when the external model makes use of data assimilation techniques (reanalysis product or operational forecast). In this situation, there is a potential benefit to relax the large scales of the local solution to the ones computed by the external model while the small scales features are still governed by the coastal model evolution. To this aim, spectral nudging methods (Waldron et al., 1996) can be viewed as a FRS scheme which acts on the whole local domain on specified (large) scales. It can be written under the following form:

$$\partial_t \phi_{\text{loc}} + L_{\text{loc}}(\phi_{\text{loc}}) + \Lambda(\phi_{\text{loc}} - \phi_{\text{ext}}) = 0 \quad \text{in } \Omega_{\text{loc}}, \tag{54}$$

where the operator Λ is now a low pass filter acting on the whole Ω_{loc} domain. The expression of Λ may be based on Fourier transform of $(\phi_{\text{loc}}, \phi_{\text{ext}})$ (with relaxation coefficients fixed in the Fourier space), on Gaussian correlation

kernels and more simply on iterative Laplacian smoothing. When applied to coastal modelling, the nudging term is usually reduced near the coast where the external model may have a coarser representation of the coastline.

Finally, most of the time, the FRS scheme is associated to a sponge layer that damps the small scales near the boundary using increased viscosity/diffusivity ν:

$$\partial_t \phi_{\text{loc}} + L_{\text{loc}}(\phi_{\text{loc}}) = -\frac{1}{\tau}(\phi_{\text{loc}} - \phi_{\text{ext}}) + \nu_\Gamma \, \Delta\phi_{\text{loc}} \quad \text{ in } \Omega_{\text{loc}}. \tag{55}$$

Radiation methods. In the case of outflow conditions, the FRS scheme will only behave well (and in particular will prevent waves reflection) when the external data is already an accurate representation of the solution. In other cases, radiation methods can be used to propagate out the disturbances. The basic radiation methods solve the Sommerfeld radiation condition (56) at the interface,

$$\partial_t \phi_{\text{loc}} + c \, \partial_n \phi_{\text{loc}} = 0 \qquad \text{on } \gamma, \tag{56}$$

and corresponds to the transport of ϕ through γ at the phase speed projected to the boundary, c (∂_n denotes a derivative perpendicular to and towards the outside of teh boundary). Obviously this boundary condition is only exact for a wave propagation problem and when c is known. In practice, c $\left(= -\partial_t \phi_{\text{loc}} \left(\partial_n \phi_{\text{loc}}\right)^{-1}\right)$ is estimated at the discrete level using internal values (Orlanski, 1976), leading to different implementations, including, in two dimensions, variants taking into account tangential derivatives (Raymond and Kuo, 1984). Besides the evaluation of the phase speed c, the relevancy of this type of boundary condition for non-monochromatic and/or dispersive waves is questionable.

The radiation condition is valid only for outflow conditions (e.g. $c > 0$ at an eastern boundary). In other cases, relaxation methods as described in the previous paragraph are more appropriate. Radiation and relaxation methods can be combined to give:

$$\partial_t \phi_{\text{loc}} + c \, \partial_n \phi_{\text{loc}} = -\frac{\phi_{\text{loc}} - \phi_{\text{ext}}}{\tau_\gamma} \qquad \text{on } \gamma. \tag{57}$$

where τ_γ, the relaxation time scale, has different values $\tau_{\text{in}}, \tau_{\text{out}}$ according to inflow ($\tau_\gamma = \tau_{\text{in}}$) or outflow ($\tau_\gamma = \tau_{\text{out}}$) conditions (Marchesiello et al., 2001). During inflow conditions c is set to zero and τ_{in} is small leading to a strong relaxation towards external solution. τ_{out} is large but not infinity in order to

prevent the internal solution to drift for external data. Note that even if (57) is written at the interface γ, it is most of the time associated to a relaxation term active on the whole transition area Γ. For consistency, the boundary value of the relaxation time scale τ defined in (53) should be greater than τ_{out}.

6.2.2. Characteristic methods

As mentioned previously, radiation methods are only valid when the temporal evolution of the state variable is governed by a wave equation. They can be extended to more general cases by looking at the hyperbolic parts of the equations. By definition hyperbolic systems can be decomposed into a series of wave propagation problems whose solutions are the characteristic variables. The problem is then well posed if the incoming characteristics variable are specified and the outgoing characteristic extrapolated from internal values (Blayo and Debreu, 2005).

Considering a one-dimensional version of the linearised barotropic equation (45)

$$\partial_t \eta = -H \partial_x \bar{u}, \quad \partial_t \bar{u} = -g \partial_x \eta,$$

the two characteristics are given by

$$w_\pm = \bar{u} \pm \sqrt{\frac{g}{H}}\, \eta$$

and propagate with speeds $c_\pm = \pm\sqrt{gH}$ (i.e. $\partial_t w_\pm + c_\pm \partial_x w_\pm = 0$). At open boundaries, the incoming characteristics are prescribed ($w_{\text{loc}} = w_{\text{ext}}$) and the outgoing characteristic are computed from internal values. The computation of outgoing characteristic variables at the boundary can be based on simple extrapolation techniques or on the application of upwind biased schemes. It is well known that the specification of the incoming characteristics of the barotropic equations is equivalent to the Flather boundary condition (Flather, 1976) (which can also be obtained by a Sommerfeld radiation condition for $\bar{u}$, with $c = \sqrt{gH}$, and a 1D version of the continuity equation). At an eastern boundary where $w_- = \bar{u} - \sqrt{\frac{g}{H}}\, \eta$ must be prescribed, the Flather boundary condition reads

$$\bar{u} = \bar{u}_{\text{ext}} + \sqrt{\frac{g}{H}}\, (\eta - \eta_{\text{ext}}) \tag{58}$$

and where η is given by another boundary condition (e.g. $\partial_x \eta = 0$). The use of the full characteristic method gives instead:

$$\bar{u} = \frac{1}{2}\left(w_{+,\mathrm{int}} + w_{-,\mathrm{ext}}\right) = \frac{1}{2}\left(\bar{u}_{\mathrm{ext}} + \bar{u}_{\mathrm{int}}\right) + \frac{1}{2}\sqrt{\frac{g}{H}}\left(\eta_{\mathrm{int}} - \eta_{\mathrm{ext}}\right) \tag{59}$$

and

$$\eta = \frac{1}{2}\sqrt{\frac{H}{g}}\left(w_{+,\mathrm{int}} - w_{-,\mathrm{ext}}\right) = \frac{1}{2}\left(\eta_{\mathrm{ext}} + \eta_{\mathrm{int}}\right) + \frac{1}{2}\sqrt{\frac{H}{g}}\left(\bar{u}_{\mathrm{int}} - \bar{u}_{\mathrm{ext}}\right), \tag{60}$$

where $w_{+,\mathrm{int}} = \bar{u}_{\mathrm{int}} + \sqrt{\frac{g}{H}}\eta_{\mathrm{int}}$ is computed from internal values. The Flather type boundary condition (58), or some of its variants, is used in most of the coastal ocean models where it is assumed that surface gravity waves are indeed dominant for the specification of the boundary conditions. If $\bar{u}_{\mathrm{ext}}$ is not known but only η_{ext} (e.g. tidal conditions), a reduced model (e.g. geostrophic balance) may be used to deduce $\bar{u}_{\mathrm{ext}}$.

6.2.3. 3D models

The 3D primitive equation models support the propagation of internal gravity waves whose characteristics can be obtained through a vertical normal mode decomposition. Setting 3D boundary conditions on the original state variables without taking into account these propagations renders the problem ill-posed (Oliger and Sundström, 1978). A remedy can be to add a small vertical viscosity in the hydrostatic balance equations (7) (Témam and Tribbia, 2003). Characteristics-based methods can formally be extended to 3D-based on a vertical mode decomposition (Jensen, 1998; Blayo and Debreu, 2005) resulting into a series of shallow water models. However the number of assumptions used in their derivations (in particular the linearisation around an horizontally constant background density field and a constant background state velocity field, typically the barotropic solution) make them difficult to apply in practice. In addition, the staggering of the grid variables on an Arakawa C-grid add some difficulties to the discrete implementation of a combination of variables at different grid locations, thus introducing a number of additional assumptions. That is why, in 3D primitive equation models, the preceding relaxation/radiation boundary conditions are usually applied on each of the prognostic variables $(\eta, \bar{u}, \bar{v}, u, v, T, S)$, with a Flather-like boundary condition for the barotropic part.

6.2.4. Local mesh refinement (one-way nesting)

One-way nested models can use any of the preceding boundary conditions with parameters adjusted according to the space and time refinement factors which are here assumed to be relatively small. One-way nesting may be performed online (Penven et al., 2006) or offline (Mason et al., 2010). The key element for a choice between these two approaches is the spatio-temporal variability of the boundary conditions. When this variability is high, online nesting, with high frequency / high resolution exchanges between local and external domains, is more efficient. Indeed storage and performance issues limit the frequency/resolution of the outputs made during the coarse resolution run. However in an offline approach, the obvious advantage is that the area covered by the local (high resolution) domain does not have to be known in advance, i.e. before running the external (coarse resolution) model.

6.3. Two-way interaction

We consider now a two-way interaction between the external and local models. The external model feels the local model by a feedback of the local solution. We differentiate between traditional two-way nesting algorithms where the local domain corresponds to a local grid refinement of the external domain and more general two-way coupling algorithms.
We will not detail here the interpolation/update operators between the two models. The reader is referred to Debreu and Blayo (2008) for their design, in particular in terms of scale selectivity and conservation.

6.3.1. Two-way nesting

Two-way nesting algorithms have been reviewed in Debreu and Blayo (2008). In this paragraph, we describe strongly coupled models which eventually can be viewed as a single model with an abrupt change of resolution at the high/coarse resolution grids interface γ.

Grid interactions. In two-way nesting applications, the feedback from the high to the coarse resolution grid prevents the two solutions to drift. In the limit of equal local and external grid resolutions ($L_{\text{ext}} = L_{\text{loc}}$) the original solution (without nesting) should be recovered. This implies a strong criterion on the forcing of the local model: the boundary conditions have to act on the difference between external and local solutions. We will call this a consistency condition. Typically, the sponge layer in (55) should now be written

$$\partial_t \phi_{\text{loc}} + \ldots = \nu_\Gamma \, \Delta(\phi_{\text{loc}} - \phi_{\text{ext}}) \quad \text{in } \Omega_{\text{loc}}, \tag{61}$$

and if radiation conditions are used, (57) should be replaced by:

$$\partial_t(\phi_{\rm loc} - \phi_{\rm ext}) + c\,\partial_n(\phi_{\rm loc} - \phi_{\rm ext}) = -\frac{\phi_{\rm loc} - \phi_{\rm ext}}{\tau}. \tag{62}$$

In a sense, in these equations, the temporal and spatial scales of the external solutions are not assumed to be much more larger than those of the local domain so that the simplification occurring in (55) and (57) cannot be made. Condition (62) was used in Perkins et al. (1997).

Elliptic systems and/or time splitting. The discrete versions of $L_{\rm ext}$ and $L_{\rm loc}$ may involve the solution of an elliptic system. This occurs for example when an implicit treatment of the surface gravity waves is applied, see Section (5.1.1), or if a nonhydrostatic kernel is involved (leading to the solution of a Poisson system), see Section 8.2. In that case, special care must be taken in order to get to correct solution (i.e. continuity of the solution and its derivatives). For being properly treated, the implicit solver has to be transformed into a multi-resolution solver which includes the grid interactions (e.g., paragraph 2.2.1 of Debreu and Blayo (2008) and Haley and Lermusiaux (2010); Cailleau et al. (2008)).

Similarly, when an explicit mode splitting scheme is employed (see Section 5.1), the models should be coupled at the barotropic time step instead of at the baroclinic time step (Debreu et al., 2012). This also allows to satisfy the consistency condition previously mentioned.

Vertical refinement. Most of the current two-way nested models apply only horizontal mesh refinements. Local vertical refinements (e.g. increase of resolution in surface and/or bottom boundary layers) are difficult to manage for several reasons. Vertical parameterizations (see Section 3.2.1) involve implicit solvers due to stability constraints and, so, as for horizontal implicit solvers (paragraph above) are more difficult to tackle on multiresolution grids. Coupling between barotropic and baroclinic modes (see Section 5.1) can also become cumbersome if only part of the water column is computed on the fine grid. Instead of a local vertical mesh refinement, a local change of vertical coordinate systems is achievable. In coastal applications, the capacity to change locally the vertical coordinate systems (4.2) can become crucial for some applications. Indeed, mesh refinement along coastal areas with very small water depths may require some local adaptation of the parameters of

the vertical coordinate systems (parameters of the terrain-following coordinates, number of layers, ...). In that case, vertical remapping has to be applied between coarse and fine grids.

Time refinement. In traditional applications, horizontal mesh refinement is most of the time coupled with time refinement in order to satisfy stability conditions. If the main stability constraint is given by external gravity waves propagation (cf 5.1), at a velocity of $\sqrt{gH}$, we may expect that, when refining the grid near the coast and so with smaller depths, the time step does not have to be reduced. This is certainly the case for some applications, but generally having the possibility to refine the time steps may become important when the actual time step is constrained by other phenomena (e.g. drying and flooding, advection ...).

6.3.2. Two-way coupling

Less strong local/external coupling may also be of interest, particularly when the space and time refinement factors are large and/or if the models L_{ext} and L_{loc} have different physics. In that case, two-way relaxation methods may also be of interest. The spectral nudging method introduced in Section 6.2.1 can be extended to two-way interactions. The smooth semi-prognostic method (Sheng et al., 2005) is an example of such a technique. More generally, grids interactions can be (weakly) incorporated in any assimilation scheme coupling at the same time the assimilation of observations and the grid interactions.

6.4. Discrete point of view

In addition to the preceding continuous formulations, a number of questions arise when taking into account the model discretisations.

6.4.1. Implementation of boundary conditions

We already mentioned the discretisation of the Sommerfeld radiation condition, including the sensitivity to the computation of the normal wave speed in 2D. The discrete implementation of boundary conditions has to deal with the time stepping algorithm and the grid staggering. As an example, Mason et al. (2010) proposed a stable implementation of a 2D characteristic method for the barotropic equations on a staggered C-grid. This paper introduces a stable treatment of the outgoing characteristics (computation of $\bar{u}_{\text{int}}$ and η_{int} in (59)).

6.4.2. Link with internal numerical schemes

More generally, the problem of open boundary conditions and model nesting is also fundamentally tied to the discrete internal schemes (e.g. implicit/explicit treatment, time integration scheme, discrete conservation properties, diffusive terms). All these aspects interact with an efficient implementation of the boundary conditions. The trivial example is the use of a first-order upwind biased scheme for a simple advection equation: the original equation $\partial_t \phi + c\partial_x \phi = 0$ is hyperbolic and thus requires only one boundary condition at inflow, otherwise over-specification occurs. However, the modified equation (i.e. including spatial truncation errors) is of parabolic nature $\partial_t \phi + c\partial_x \phi = \mu \partial_{xx} \phi + \mathcal{O}(\Delta x)^2, \mu = c\,\dfrac{\Delta x}{2}$ and allows for the specification of boundary conditions both at inflow and outflow boundaries. Indeed, in practice the numerical scheme will naturally take into account only external data at inflow. Similarly if a dispersive scheme is used to discretise the spatial derivative (e.g. second-order centered), radiation methods will not perform well even in the monochromatic wave case. This makes the derivation of a very general OBC difficult and may also explain why some boundary conditions work for some models and not for others.

A promising approach concerning the derivation of boundary conditions is the summation-by-parts approach (Svärd and Nordström, 2014) that can make the link between internal schemes and boundary conditions (possibly weekly imposed) according to a discrete energy estimate.

Sanity control. At the discrete level, some basic sanity controls can be made. This is particularly important for one- and/or two-way nesting. The consistency condition has been mentioned previously: assuming that the same numerical/physical choices are made for the local and external solution, a simulation with spatial and temporal refinement equal to one should lead to the same solution with or without nesting. This has several implications on the formulation of the boundary conditions, the choice of the number of ghost cells and the treatment of barotropic/baroclinic coupling.

The second sanity control is to verify the constancy preservation (see Section 7.1): in absence of sinks and sources, if a tracer is globally initialised to a constant value, the tracer should remain constant on both grid and with the same value. This has again several implications on the way interpolation/update schemes are constructed and the consistency of barotropic and baroclinic exchanges.

7. Numerical treatment of single terms and processes

This section discusses discrete algorithms that are commonly used for the solution of the governing equations formulated in Sec. 3 and 4. Originally, coastal ocean models were discretised using numerical methods loosely inherited from large scale ocean models, namely second-order in space and Leapfrog-in-time schemes (Blumberg and Mellor, 1987). However, over the years, the numerics of those models have evolved toward the use of more advanced discretisation techniques, more similar to techniques used in other fields of computational fluid dynamics, in order to accommodate to the specific constraints of coastal applications. Indeed, as discussed in Sec. 2, coastal ocean dynamics is characterised by strong vertical and horizontal gradients, long waves, rapid variations on short time and spatial scales as well as large dissipation and velocities near the bottom which all require specific numerical treatments. The diversity of methods in use in coastal models is large so that the discussion can only focus on techniques which are more widely used or are expected to become more widely used. Since the numerical requirements for each process can be rather different, this section considers separately sub-systems of the full system of primitive equations. However, it should not overshadow the fact that a numerical kernel must be designed as a whole, not just as an advection or a wave propagation equation since interferences between numerical methods used for different processes exist.

There often exists some confusion to properly characterise the class of discretisation techniques used in state-of-the-art oceanic coastal models. Since there are significant differences in truncation error between finite-difference and finite-volume advection-diffusion schemes (e.g. Leonard, 1995), clarifications are needed. As far as the tracer conservation is concerned, it is obvious that they must be interpreted in a finite-volume sense, meaning that for a given tracer q the discrete quantity $q_{i,j,k}$ is the average over a control volume $\mathcal{V}_{i,j,k}$ (Fig. 2). For the momentum equations, the generally adopted formulation is inherited from Lilly (1965) who introduced a discretisation approach which has the property to conserve energy in a discrete sense as long as second-order centered schemes are used. Momentum advection terms in Lilly (1965) can be viewed as a finite-volume discretisations provided that proper horizontal metric terms are added. However, not every term in the momentum equation can be cast in a flux divergence form (e.g. Coriolis term, curvilinear metric terms). For those terms momentum components should be interpreted in a finite-different sense. Note that the fact that tracer values

are cell-averaged implies that the pressure deduced from the hydrostatic relation is a point-wise value located on cell interfaces in the vertical. Generally speaking, discretisation techniques in most existing structured-grid oceanic coastal models can be classified as finite volume discretisations. The arrangement on the computational grid is shown in Fig. 2. We now give more details on the specifics of different terms.

7.1. Tracer and momentum advection

We consider here the advection equation written in a conservative way for a layer-averaged quantity q_k (q can either represent a tracer or horizontal momentum components when a flux-form of nonlinear terms is used) along with the continuity equation (29)

$$\begin{aligned} \partial_t(h_k q_k) + \partial_x(h_k u_k q_k) + \partial_y(h_k v_k q_k) + \delta_k(w^s q) &= 0, \\ \partial_t h_k + \partial_x(h_k u_k) + \partial_y(h_k v_k) + \delta_k(w^s) &= 0, \end{aligned} \tag{63}$$

using the notations introduced in Sec. 4 and the definition of a vertical difference located at the centre of the cell k, $\delta_k(a) = a_{k+1/2} - a_{k-1/2}$ for any quantity a. We retain the continuity equation here to emphasise that the advection equation must be conservative while satisfying a constancy preservation property (i.e. if q is initialised with a constant, it remains so). An overview of the time and space discretisations of (63) used in large and mesoscales oceanic models can be found in Lemarié et al. (2015). Broadly speaking, it is well known that classical even-ordered linear advection schemes tend to produce unphysical oscillations which can be controlled by an explicit scale-selective dissipation whose coefficients A_M and A_c must be tuned to ensure no build-up of grid scale noise. On the contrary, odd-ordered linear advection schemes are inherently dissipative near the grid-scale (e.g. App. B in Soufflet et al., 2016), but are not oscillation free with the exception of the first-order upwind scheme which can not, however, be considered seriously due to its unacceptable artificial diffusion. The linear compact implicit schemes (Lele, 1992) are a very attractive alternative but they require a matrix inversion which makes them tractable for vertical advection only. Note that an important constraint on the vertical advection is that the oceanic interior vertical mixing is exceptionally weak so that upwind schemes in models with quasi-Eulerian vertical coordinate or dissipative remapping schemes in models with an ALE coordinate (see section 4.2) must be avoided.

7.1.1. Design of numerical models

In practice, several strategies coexist in the coastal community. A first choice is to consider simple second-order centered schemes supplemented by an explicit diffusion and a non-dissipative time-stepping algorithm in order to ensure energy consistency of the whole code at a discrete level (Marsaleix et al., 2008). A second choice is to rely exclusively on standard high-order upwind schemes in the horizontal direction and low-dispersive centered advection schemes in the vertical (e.g. fourth-order compact schemes), without any explicit dissipation, combined with a high-order predictor-corrector time-stepping algorithm using a method of lines (e.g. Shchepetkin and McWilliams, 2005). In this case, it is accepted that strong gradients will generate overshoots and undershoots and that there is an uncontrolled sink of kinetic energy and tracer variance in the model caused by dissipative truncation errors. This last point is, however, generally not worrisome as long as a first-order upwind scheme is avoided. As an alternative to the method of lines, where space and time discretisations are considered separately, a coupled space-time approach[1] could be used for advective terms (e.g. Hundsdorfer and Trompert, 1994; Daru and Tenaud, 2004). This type of approach can be viewed as a conservative semi-Lagrangian approach where the characteristics are tracked in space-time. The resulting schemes are one-step schemes with interfacial fluxes depending on the local Courant number $\alpha^x = u_k \Delta t / \Delta x$. They have the property to be free of computational mode, exact for a Courant number $\alpha^x = 1$, and to be easily extended to arbitrary high-order of accuracy (e.g. Daru and Tenaud, 2004). Examples of such schemes are the first-order upwind scheme, the Lax-Wendroff scheme or the QUICKEST scheme which are respectively first, second and third order accurate in space and time. Whatever the strategy used (uncoupled or coupled space and time), it is generally preferable to prevent spurious numerical oscillations by using non-oscillatory schemes which are more computationally demanding but provide a way to ensure nonlinear stability. Since non-oscillatory schemes are expected to be increasingly used to simulate coastal flows, we provide additional comments on this point to highlight the associated delicacies.

[1] The coupled space and time approach is also known in the literature as flux-form semi-Lagrangian approach, as well as transient interpolation method, or direct space-time discretisation.

7.1.2. Nonlinear stability and flux correcting/limiting methods

Numerical methods with an order of accuracy larger than one are prone to numerical oscillations. A nonlinear stability criteria often considered is that the numerical solution must not oscillate without bounds. Using the notion of total variation (TV) as a measure of the overall amount of oscillations, numerical schemes have been constructed to ensure that the total variation is either bounded (TVB) or diminishing (TVD) (Harten, 1983; Shu, 1987). Since TVD schemes necessarily degenerate to first order accuracy near smooth extrema, they are prone to clipping errors manifested by the appearance of a flat plateau in the solution due to an excess of diffusion. To avoid this type of error, a less stringent constraint is the monotonicity-preservation (MP) which relaxes the TVD constraint near extrema (e.g. Suresh and Huynh, 1997; Daru and Tenaud, 2004). Flux correcting/limiting is generally used to impose the TVD, TVB or MP stability conditions (e.g. Durran, 2010). Historically, the flux limiting/correcting methods were derived in a coupled space and time framework using the second-order Lax-Wendroff scheme. They can be easily extended to higher-order accuracy using one-step coupled time-space approach of arbitrary high order (e.g. Daru and Tenaud, 2004). However if a method of lines is used the time and space dimensions are decoupled therefore the constraints on total variation must be imposed both in time and space. For this purpose, a family of Runge-Kutta (RK) schemes satisfying the TVD property has been derived by Shu (1988) and Gottlieb and Shu (1998). These RK schemes are TVD over a range of time-steps which is generally smaller than the stability range given by the more traditional CFL-type condition. Note that a consequence of the TVD, TVB and MP stability constraints is a better control of numerical oscillations hence those constraints could be used not only for momentum advection as a way to limit spurious mixing by reducing the grid Reynolds number (Ilicak et al., 2012), but also for the advection of tracers and layer thickness subject to monotonicity or positivity constraints. There is an abundant literature on the comparison of different flux limiting/correcting strategies, see Mohammadi-Aragh et al. (2015) for a recent study in the oceanic modelling context. Note that in multi-dimensions, the TVD, TVB or MP properties are generally applied independently in each spatial directions. It can be shown that in this case the resulting scheme does not strictly satisfy those properties anymore. Nonetheless this approach generally provides acceptable results.

7.1.3. Multidimensional aspects and stability constraints

The extension of one-dimensional advection schemes to multi-dimensions in the context of a method of lines is straight-forward since the necessary transverse terms are discretised in a transparent way. If a quasi-Eulerian vertical coordinate is used, the stability constraint in three dimensions reads

$$\alpha^x + \alpha^y + \beta\alpha^z \leq \alpha^\star \tag{64}$$

where $\alpha^\star$ is the CFL stability constraint associated to the numerical schemes used for the time and space discretisations, and β a coefficient arising from the fact that different advection schemes with different stability criteria may be used in the horizontal and vertical directions. Typical values of $\alpha^\star$ and β can be found in Lemarié et al. (2015). As far as the coupled space and time approach is concerned, a simple application of the one-dimensional fluxes in each spatial direction would result in a scheme that can be at most first-order accurate, whatever the order of accuracy of the 1D fluxes, and the scheme would be unconditionally unstable. To overcome this issue a directional splitting can be used to introduce the necessary transverse terms in the discretisation. A simple method to introduce transverse terms while preserving constancy is the splitting of Bott (2010). With this type of approach the stability constraint is given by

$$\max\left\{\alpha^x, \alpha^y, \alpha^z\right\} \leq 1$$

which is less stringent than (64).

For coastal applications, Shchepetkin (2015) has emphasised the severe stability constraint potentially imposed by vertical advection. Several ways can be used to mitigate this issue *(i)* use a semi-implicit advection scheme *(ii)* use a local time-step for vertical advective terms based on the maximum vertical Courant number in each water column *(iii)* use a quasi-Lagrangian vertical coordinate which is subject to a gentle Lipschitz stability condition rather than the usually severe CFL condition (e.g. White et al., 2009). This last remark applies to models using rezoning/remapping strategies because models using a transport algorithm instead of remapping (e.g. Hofmeister et al., 2010) will still be subject to a CFL type stability constraint.

7.2. Internal pressure gradient errors

Since in coastal ocean models the use of bottom-following coordinates is mandatory (see Sec. 2.1), the coordinate surfaces may intersect with isopycnal or geopotential surfaces at a relatively large angle when the bathymetry

is steep. This makes it significantly harder to accurately compute the internal horizontal pressure gradient since the pressure is not known along constant horizontal surfaces. Using the notations introduced in Sec. 4, for a generalised vertical coordinate s, the x-component of the internal horizontal pressure gradient (HPG) is expressed using the chain rule

$$\frac{1}{\rho_0}\left(\partial_x p\right)_z = \frac{1}{\rho_0}\left[\left(\partial_x p\right)_s - \partial_z p\left(\partial_x z\right)_s\right], \tag{65}$$

where p is the hydrostatic pressure deduced from the hydrostatic relation (7). The subscript s is denoting differentiation with respect to constant s. In an arbitrary general coordinate the two terms in rhs are nontrivial, of opposite sign and may be large compared to the order of magnitude of their difference, except for two special cases: For isopycnal coordinates, the first term vanishes. The second term in the rhs is sometimes referred to as compensating hydrostatic term which vanishes for geopotential coordinates ($s = z$). At a discrete level, unbalanced truncation errors in the computation of the two large terms in the rhs will interfere with each other and give rise to a possibly large pressure-gradient error leading to hydrostatic inconsistency manifested by unphysical spontaneous motions, even in the absence of lateral density gradients (i.e., $\rho = \rho(z)$). A very large body of literature is dedicated to this subject (e.g. Blumberg and Mellor, 1987; Mellor et al., 1994; McCalpin, 1994; Burchard and Petersen, 1997; Lin, 1997; Shchepetkin and McWilliams, 2003; Adcroft et al., 2008; Marsaleix et al., 2011; Berntsen, 2011) where different strategies have been followed to reformulate (65) in a way more convenient for the discretisation. Those strategies are generally motivated either by achieving a high-order accuracy to cancel the two terms in (65) in the case of a flat stratification (e.g. McCalpin, 1994; Shchepetkin and McWilliams, 2003) or by ensuring an energetic consistency (e.g. Marsaleix et al., 2008). In this latter case, the aim is to make sure in a discrete sense that the change of kinetic energy due to the work of pressure forces is balanced by the change of potential energy due to buoyancy forces. This balance occurs only with a linear equation of state and involves the discrete advection for density (e.g. Shchepetkin and McWilliams, 2003, App. A).

Mainly three strategies have been followed in practical applications to reformulate and discretise (65). *(i)* A first strategy is to directly use the hydrostatic assumption (6) to express p in terms of density integrals

$$\frac{1}{\rho_0}\left(\partial_x p\right)_z = \frac{g}{\rho_0}\left[\left(\partial_x I_\rho(z)\right)_s + \rho\left(\partial_x z\right)_s\right], \qquad I_\rho(z) = \int_z^\eta \rho(x,y,\xi)d\xi, \tag{66}$$

assuming a zero atmospheric pressure gradient at the surface. When discretising this form of pressure gradient, averaging of $I_\rho(z)$ in the vertical and ρ in the horizontal are needed since the only term computed at its natural position is $(\partial_x z)_s$. According to Shchepetkin and McWilliams (2003) this explains why this approach is often prone to large hydrostatic inconsistencies. However, high-order schemes have been proposed in McCalpin (1994) to compute the horizontal derivatives and horizontal averaging of ρ combined with a spectral representation in the vertical. Note that high-order methods in the vertical should be used with care in case of strongly stratified flows because density is generally not smooth on the grid scale, especially with coarse vertical resolution. *(ii)* a second and intuitive approach consists in computing the pressure (i.e. $I_\rho(z)$) and then to vertically interpolate the result at an horizontal surface of reference before differentiating (e.g. Lin, 1997; White et al., 2009). This is the preferred option in atmospheric models. In Lin (1997) a flux-form finite-volume approach based on this concept is adopted. *(iii)* the preferred approach for most oceanic models is to use a *density-Jacobian* discretisation. As done in (8) and (9), by combining the Leibniz rule and the hydrostatic relation we easily obtain that

$$\frac{1}{\rho_0}\left(\partial_x p\right)_z = g\frac{\rho(\eta)}{\rho_0}\partial_x \eta - \int_z^\eta \underbrace{\left[\left(\partial_x b\right)_s - \partial_{z'} b\left(\partial_x z'\right)_s\right]}_{\mathcal{J}_{x,z'}(b,z')} dz' \tag{67}$$

where the buoyancy b is defined in Sec. 3. Compared to the straight-forward approach (66) differentiation comes first and integration then. The density-Jacobian formulation clearly shows that if $s = \rho$ only one term is nonzero and the pressure gradient term can be computed accurately provided that proper care is taken to the equation of state (Adcroft et al., 2008). The Jacobian in (67) must be computed at cell corners, if we refer to Fig. 5 a standard second-order discretisation reads

$$\begin{aligned}\left(\mathcal{J}_{x,z}\right)_{i+1/2,k+1/2} &= \frac{1}{2\Delta x}\big\{(1+r_x)b_{i+1,k+1} + (1-r_x)b_{i+1,k} \\ &\qquad -(1-r_x)b_{i,k+1} - (1+r_x)b_{i,k}\big\}\end{aligned} \tag{68}$$

with $r_x = -\frac{z_{i+1,k+1/2} - z_{i,k+1/2}}{z_{i+1/2,k+1} - z_{i+1/2,k}}$ the grid slope. This discretisation of the Jacobian has been widely used in coastal models for years. It can be shown that the standard second-order Jacobian discretisation corresponds to a discrete

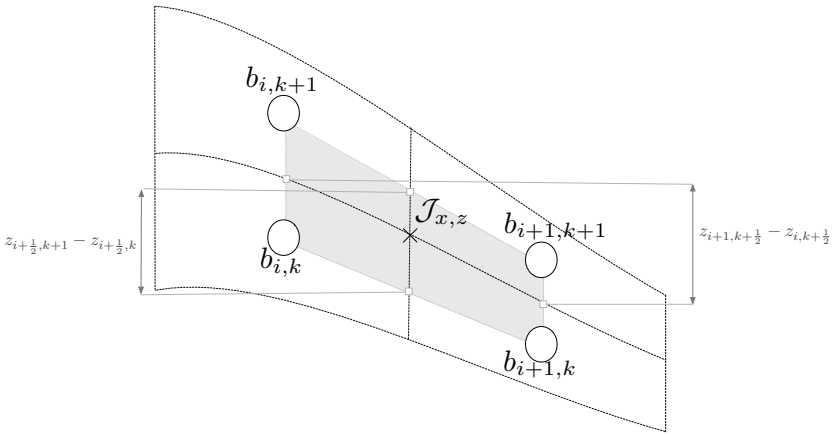

Figure 5: Elementary computational stencil of the standard second-order density Jacobian $\mathcal{J}_{x,s}$. The distances involved in the computation of the grid slope r_x are reported.

form of Green's theorem. Indeed, (68) can be written as the sum of four line integrals discretised using a simple trapezoidal rule. In an attempt to further reduce the pressure gradient errors, Shchepetkin and McWilliams (2003) have introduced a high-order (cubic) discretisation for those line integrals. In their work, care has been taken to the robustness of the scheme even in case of non-smooth density fields. The resultant scheme turned out to be a significant step forward compared to existing approaches. Note that besides the imbalance in the truncation errors in (65) there is an other source of pressure gradient error stemming from the use of a nonlinear equation of state including compressibility of seawater (Dukowicz, 2001). Ways to remove this source of error by rearranging the terms in the equation of state are presented in Shchepetkin and McWilliams (2003), Adcroft et al. (2008) or Marsaleix et al. (2011). This latter source of error is sometimes overlooked, however it must be stressed that the careless inclusion of compressibility terms into the equation of state would have significant dynamic consequences.

The value of the parameter r_x in (68) is a good qualitative indicator of discretisation errors. For $r_x = 0$ the grid is flat and the pressure gradient term reduces to a single term while for $|r_x| = 1$ there is a 45 degrees slope between the grid and the horizontal direction, and (68) reduces to a simple differentiation of two terms in the diagonal. More generally, for $|r_x| \leq 1$ we expect the

discretisation to be fairly accurate since it is based on vertical interpolations, this is the so-called "hydrostatic consistency" situation (Fig. 5). Values of $|r_x|$ greater than 1 generally lead to a degradation of the accuracy due to the need of vertical extrapolation to compute the pressure gradient term. This corresponds to a situation of "hydrostatic inconsistency", see Haney (1991). A strong requirement to avoid the emergence of spurious motions is thus to keep the value of r_x at a reasonable level. To do so several ways are commonly used concurrently, *(i)* bathymetry smoothing (e.g. Martinho and Batteen, 2006) is systematically used in terrain-following coordinates to limit the maximum slopes *(ii)* the function responsible for the vertical placement of grid points is designed in such a way to minimise the departure of the vertical coordinate isosurfaces from the horizontal direction (e.g. Schär et al., 2002; Shchepetkin and McWilliams, 2009; Berntsen, 2011; Lemarié et al., 2012b) *(iii)* a quasi-Lagrangian vertical coordinate whose layer distribution in the rezoning step can be chosen to reduce grid slopes or to follow isopycnals (e.g. Burchard and Petersen, 1997; White et al., 2009; Hofmeister et al., 2010). Note that trying to confine the maximum grid slopes at the bottom to quickly obtain nearly flat coordinate isosurfaces when going toward the surface must be done with care in the sense that large pressure gradient errors at depth will affect the circulation in the upper layers anyway through baroclinic-barotropic coupling. Besides the approaches discussed so far, we could imagine alternative strategies. For example, the coordinate system could be determined using only the low-frequency variations in the bathymetry and high frequencies could be handled using an immersed boundary method (e.g. Lundquist et al., 2010). Other alternatives could be the use of well-balanced discretisations (Botta et al., 2004) or to express horizontal momentum using the covariant velocity components instead of the usual contravariant components (Weller and Shahrokhi, 2014). Eventually, as emphasised by Mellor et al. (1994), the interaction between pressure gradient and other terms in the equations should be further clarified, especially the interaction with active tracers advection. More specifically, the interplay between numerically-induced mixing of active tracers and HPG errors is expected to play a crucial role in the time-evolution of circulation errors (e.g. Lemarié et al., 2012b, Sec. 6.2). However, the importance of HPG errors in simulations of coastal currents is difficult to evaluate in practical applications. The competing effects of the model dynamics, its sensitivity to the horizontal/vertical grid resolution and discrete numerical algorithms, are tricky to untangle.

7.3. Inertia-gravity waves

Most of the work on the discretisation of inertia-gravity waves was mainly concerned with finding the most appropriate variable arrangement among the Arakawa grids A to E by looking at discrete dispersion relations (e.g. Mesinger and Arakawa, 1976). Here we restrict ourselves to the C-grid arrangement which is the preferred option in coastal models since the Rossby radius of deformation is generally properly resolved, with the exception of fjord-type systems such as the Baltic Sea, where very high resolution $< 1km$ is required (Holtermann et al., 2014). Typical space-time discretisation techniques can be found in Beckers and Deleersnijder (1993) or Lemarié et al. (2015) and are summarised below.

7.3.1. Inertial oscillations

In the absence of other terms but Coriolis terms in (30) and (31) we obtain the equations of free motion on a rotating plane

$$\partial_t(h_k u_k) = f h_k v_k; \quad \partial_t(h_k v_k) = -f h_k u_k. \tag{69}$$

A characteristic of the Coriolis acceleration is that it is directed perpendicular to the velocity direction. Hence, it does not play any role in the kinetic energy budget, which forms an important constraint for the discretisation (e.g. Arakawa and Lamb, 1977). In (69), the quantity $f h_k v_k$ must be given at u-points and $f h_k u_k$ at v-points. If (69) is semi-discretised in space, this constraint is trivially satisfied locally on an Arakawa B-grid, since u and v are collocated, but can only be satisfied globally on an Arakawa C-grid. Considering that the Coriolis frequency f is cell centered, the following discretisation of the rhs in (69)

$$(f h_k v_k)_{i+1/2,j} = \overline{f_{i,j} h_{i,j,k} \overline{v}^{\,x}_k}^{\,x}, \qquad (f h_k u_k)_{i,j+1/2} = \overline{f_{i,j} h_{i,j,k} \overline{u}^{\,y}_k}^{\,y}, \tag{70}$$

where $\overline{\cdot}^{\,m}$ is an averaging operator in the direction m ($m = x, y$), ensures that all the terms associated to the Coriolis force cancel when a kinetic energy equation is derived from (30) and (31) (Sec. III.C in Arakawa and Lamb, 1977). In (70), the horizontal variations of the curvature metric terms have been ignored for clarity. An alternative energetically-consistent discretisation providing a different logic for the cancellation of the terms when the kinetic energy budget is formed at the discrete level has been introduced in Espelid et al. (2000) and reads

$$(f h_k v_k)_{i+1/2,j} = f_{i+1/2,j} \overline{(h_k v_k)}^{\,w}_{i+1/2,j}; (f h_k u_k)_{i,j+1/2} = f_{i,j+1/2} \overline{(h_k u_k)}^{\,w}_{i,j+1/2}, \tag{71}$$

where $\overline{\cdot}^{w}$ is a weighted average defined as

$$\overline{(h_k v_k)}^{w}_{i+1/2,j} = \frac{w_{i+1/2,j,k}}{4} \sum_{n=0}^{n=1} \sum_{m=0}^{m=1} \frac{(hv)_{i+n,j-m+1/2,k}}{w_{i+n,j-m+1/2,k}};\ w_{i+1/2,j,k} = \sqrt{\frac{g h_{i+1/2,j,k}}{|f_{i+1/2,j}|}}$$

for each layer k, same rationale applies to $\overline{(h_k u_k)}^{w}_{i,j+1/2}$. Compared to (70), this formulation requires the Coriolis parameter to be defined at u and v points and allows more freedom in the way to define the thickness layer h_k at cell interfaces.

When the time dimension is discretised, the energy conservation property is less straight-forward to ensure. First, a simple forward Euler step to integrate (69) would be unconditionally unstable. The use of a (semi)-implicit scheme would remove stability constraints. This approach is essentially adopted in models which treat the gravity waves in a semi-implicit way (e.g. Walters et al., 2009). In general, for coastal applications, the baroclinic time-step is much smaller than the inertial period meaning that the Coriolis term is not expected to impose any limitation on the time-step. For this reason explicit schemes are generally preferred. To circumvent the stability issues of the Euler forward step, a Forward-Backward in Time (FBT) approach is often used (Bleck and Smith, 1990)

$$\begin{aligned} (h_k u_k)^{n+1}_{i+1/2,j} &= (h_k u_k)^{n}_{i+1/2,j} + \Delta t (f h_k v_k)^{n+s}_{i+1/2,j} \\ (h_k v_k)^{n+1}_{i+1/2,j} &= (h_k v_k)^{n}_{i+1/2,j} - \Delta t (f h_k u_k)^{n+1-s}_{i,j+1/2}, \end{aligned} \tag{72}$$

and is stable for $f\Delta t \leq 2$. In (72), s is either 0 if n is even or 1 otherwise. The major drawbacks of this approach are that it requires to switch at each time-step the order of integration of the u and v components and it is not energetically-consistent even if the spatial discretisation is. Other popular explicit schemes to integrate (69) include the third-order Adams-Bashforth scheme (stable for $f\Delta t \leq 0.72$) and the leapfrog scheme (stable for $f\Delta t \leq 1$) which both globally ensure that there is no false generation of kinetic energy since u and v are considered at the same moment in time.

7.3.2. Internal gravity waves

Elements about the temporal discretization of external gravity waves are introduced in Sec. 5.1 so that we focus here on internal gravity waves. The

numerical integration of those waves is generally studied on the following one-dimensional subsystem

$$\begin{aligned} \partial_t p_q + \rho_0 c_q^2 \partial_x u_q &= 0 \\ \rho_0 \partial_t u_q + \partial_x p_q &= 0, \end{aligned} \tag{73}$$

for a given vertical mode q ($q = 0$ corresponds to the barotropic mode) with c_q the gravity wave speed associated with q^{th} vertical mode. The details to derive (73) from the full three-dimensional equations can be found, e.g., in Lemarié et al. (2015). In (73), the term $\partial_x u_q$ in the first equation can be traced back to the horizontal divergence in the continuity equation and the term $\partial_x p_q$ in the second equation to the horizontal pressure gradient term in the primitive equations. The stability constraint imposed by internal gravity waves reads

$$\Delta t \sqrt{c_1^2 \left(\frac{1}{\Delta x^2} + \frac{1}{\Delta y^2} \right)} \leq \alpha^\star_{\text{igw}}$$

where $\alpha^\star_{\text{igw}}$ is the CFL stability criterion depending on space and time discretization used for internal waves and c_1 the speed of the fastest internal wave. Typical values for $\alpha^\star_{\text{igw}}$ can be found in Lemarié et al. (2015) and Tab. 1. Following Hallberg (1997) and Shchepetkin and McWilliams (2008), a general family of two-level schemes for the time integration of (73) in a forward manner is

$$\begin{aligned} p_q^\star &= p_q^n - c_q^2 \Delta t \; \partial_x (\rho_0 u_q^n) \\ \rho_0 u_q^\star &= \rho_0 u_q^n - \Delta t \; \partial_x \left(\vartheta p_q^\star + (1-\vartheta) p_q^n \right) \\ p_q^{n+1} &= p_q^n - c_q^2 \Delta t \; \partial_x \left(\varsigma \rho_0 u_q^\star + (1-\varsigma) \rho_0 u_q^n \right) \\ \rho_0 u_q^{n+1} &= \rho_0 u_q^n - \Delta t \; \partial_x \left((1-\varphi) \left[(1-\varepsilon) p_q^{n+1} + \varepsilon p_q^\star \right] + \varphi p^n \right) \end{aligned} \tag{74}$$

where the parameters ϑ, ς, ε and φ can be chosen to improve both accuracy and stability. As reported in Tab. 1, the general discretization (74) encompasses several well-known schemes presented in the literature. Among the various alternatives, by far the simpler strategy is the so-called forward-backward approach (e.g. Mesinger and Arakawa, 1976). Other alternatives include the time staggering of tracers and momentum (e.g. Backhaus, 1985) or the use of a larger stencil in time (Brown and Campana, 1978; Shchepetkin and McWilliams, 2005). Of interest for the coastal modeling community, Rueda et al. (2007) showed that the original implementation of the Casulli

	Forward-Backward	Casulli and Cattani (1994)	Rueda et al. (2007)	Shchepetkin and McWilliams (2008)
ϑ	-	-	$1/6$	$13/24$
ς	0	0	$1/2$	$1/4$
φ	0	$1/2$	$1/2$	$1/4$
ε	0	0	0	$5/18$
$\alpha^{\star}_{\text{igw}}$	1	unstable	1	$\sqrt{3}$
Order of accuracy	2^{nd}	1^{st}	3^{rd}	3^{rd}

Table 1: Parameter values (ϑ, ς, ε, φ) in the general two-level forward-time integration scheme (74) corresponding to different alternatives existing in the literature. The associated stability constraint $\alpha^{\star}_{\text{igw}}$ and order of accuracy are also given. For simplicity the parameters reported for the Shchepetkin and McWilliams (2008) scheme have been slightly modified compared to the original paper without compromising its properties.

and Cattani (1994) time-integration scheme with a semi-implicit treatment of the external gravity waves is baroclinically unstable for the integration of internal gravity waves unless a backward Euler scheme is used for the external mode. As far as the spatial discretisation is concerned, the overwhelming majority of oceanic models consider simple second-order schemes to compute $\partial_x u_q$ (i.e. the continuity equation) and $\partial_x p_q$ (i.e. the horizontal pressure gradient). However studies by Blayo (2000) and Demange et al. (2014) emphasised that higher-order approximations of both the pressure gradient term and the horizontal divergence in the continuity equation allow a significantly better representation of internal waves propagation, compared to standard algorithmic choices.

7.4. Orientation of diffusion tensor and numerically-induced mixing

In (8), (9), (15), and (17), the horizontal turbulent fluxes are oriented in a direction $(\widetilde{x},\widetilde{y})$ to be defined. In the case of a stratified flow, if the vertical coordinate $s \neq \rho$ it is well known that maintaining the diffusion tensor along the coordinate lines can be responsible for non-physical mixing of water masses and associated spurious baroclinic currents (e.g. Barnier et al., 1998). It is thus relatively common for non-isopycnic models to implement a rotation of the diffusion tensor in a direction non-aligned with the computational grid (e.g. Stelling and Van Kester, 1994; Griffies et al., 1998). However, the

implementation of such a rotation tensor turned out to be very delicate notably to avoid unphysical undershooting and overshooting and to maintain reasonable accuracy in case of large slopes, similarly to the HPG calculations (Beckers et al., 1998, 2000; Lemarié et al., 2012a). Since for a generalized coordinate it is expected that the slope between the isopycnal direction and the direction of the computational grid is larger than for a classical z coordinate, a good compromise is generally to rotate the explicit diffusion in the geopotential direction rather than the physically correct isopycnal direction. This allows to avoid delicacies associated with the computation of isopycnal slopes in unstratified regions and at the intersection with the surface and the bottom.

A non negligible amount of non-physical mixing of water masses can also be due to the advection operator. Indeed, as mentioned earlier, oceanic coastal models generally rely either on high-order linear upwind schemes or on flux correcting/limiting approaches to discretise horizontal advection. The associated truncation errors are dissipative. Much effort have been recently directed toward the quantification and/or modification of physical orientation of this mixing of numerical nature (e.g. Burchard and Rennau, 2008; Marchesiello et al., 2009; Ilicak et al., 2012; Klingbeil et al., 2014). For example, Marchesiello et al. (2009) showed that the dissipative truncation error associated to a third-order upwind scheme in a terrain-following coordinate model was greatly exceeding the small level of physically acceptable mixing observed in the oceanic interior. The physical results proved to be greatly improved by splitting the third-order upwind scheme between a centered and a diffusive part and when the diffusive part is rotated in the isopycnal direction (Marchesiello et al., 2009; Lemarié et al., 2012b). This approach, which is viable only for linear advection schemes, required a thorough analysis of the biharmonic isopycnal mixing operator to obtain a stable and sufficiently accurate discretisation for large slopes (Lemarié et al., 2012a). However, this methodology has some limitations: the rotated operator is derived under the small slope approximation which may no longer be valid in coastal applications (Lemarié et al., 2012b, Sec. 2.4.2) and it is unclear how the diffusion operator behaves in case of fast temporal variations of isopycnals. An alternative approach to have more control on the orientation of mixing for coastal applications is the use of quasi-Lagrangian approach (e.g. Hofmeister et al., 2010; Leclair and Madec, 2011). Besides the orientation of the diffusion tensor, Ilicak et al. (2012) emphasised that spurious diapycnal mixing can be aggravated if the grid Reynolds number is too large (i.e. if the velocity field

is not smooth on the grid scale). Useful diagnostics to help characterising the amount and/or orientation of spurious mixing can be found in Ilicak et al. (2012) and Klingbeil et al. (2014). Note that even if much of the attention is on the advection operator, a non-negligible amount of numerical mixing could also be associated to other terms in the equations (Soufflet et al., 2016).

7.5. Bottom friction

Due to the relatively thick bottom boundary layers in the coastal ocean, see (1), an accurate discretisation of bottom friction is required. The no-slip bottom boundary condition for u and v (12) is typically not directly discretised, because a vertical resolution on scales of millimetres would be required to obtain sufficient accuracy. Instead, it is assumed that the velocity profiles in the lowest grid box of an ocean model follow the law of the wall:

$$\left[u\left(z'\right), v\left(z'\right)\right] = \frac{\left[u_*^{\mathrm{b}}, v_*^{\mathrm{b}}\right]}{\kappa} \ln\left(\frac{z' + z_0^{\mathrm{b}}}{z_0^{\mathrm{b}}}\right), \tag{75}$$

with the bottom roughness length, z_0^{b}, the bottom friction velocity vector, $\left(u_*^{\mathrm{b}}, v_*^{\mathrm{b}}\right)$ and the distance from the wall, $z' = z + H$. With (75), the discrete bottom boundary condition has the form of a quadratic friction law

$$\begin{aligned} \frac{\tau_{[\mathrm{x,y}]}^{\mathrm{b}}(h,u,v)}{\rho_0} &= \left(\frac{\kappa}{\ln\left(\frac{h/2+z_0^{\mathrm{b}}}{z_0^{\mathrm{b}}}\right)}\right)^2 \cdot \left[u,v\right]\left(u^2+v^2\right)^{1/2} \\ &= c_d \cdot \left[u,v\right]\left(u^2+v^2\right)^{1/2}. \end{aligned} \tag{76}$$

In (76),

$$\tau_{[\mathrm{x,y}]}^{\mathrm{b}} = \rho_0 \left[u_*^{\mathrm{b}}, v_*^{\mathrm{b}}\right] \left(\left(u_*^{\mathrm{b}}\right)^2 + \left(v_*^{\mathrm{b}}\right)^2\right)^{1/2} \tag{77}$$

is the bed stress vector, c_d is the non-dimensional bottom drag coefficient, h is the thickness of the lowest grid box, and (u, v) is the velocity vector in the lowest grid box. Note that for the limit of $h \to 0$ we obtain $c_d \to \infty$ and thus retain the original no-slip condition (12). The numerically calculated bed stress vector calculated in (76) is then used in the bottom boundary condition (13). In situations when the near-bottom velocity profile deviates strongly from the law of the wall, e.g. when stratification due to high sediment concentrations or a wave-enhanced bottom boundary layer (or both) are

present, the vertical resolution needs to be sufficiently high. One other possibility would be to parameterise such unresolved processes (Burchard et al. (2008); Warner et al. (2008)). Specific care needs to be taken for the bottom (and surface) boundary conditions when turbulence closure models are used. As shown by Burchard and Petersen (1999) and Burchard et al. (2005b), flux boundary conditions for turbulent quantities provide much higher accuracy in reproducing the law of the wall than Dirichlet boundary conditions.

7.6. Numerical treatment of turbulence closure equations

The discretisation of turbulence closure budget equations requires specific care, since turbulence quantities are often non-negative scalars such as the turbulent kinetic energy (TKE) k or its dissipation rate ε. Furthermore, these quantities often cover several orders of magnitude across fairly short time and spatial scales such as in oscillating tidal flow. The problem of a straight forward, explicit-in-time discretisation of the TKE equation is best demonstrated for the idealised case of freely decaying homogeneous turbulence. In this case the TKE budget equation (24) simplifies to the following form:

$$\partial_t k = -\varepsilon. \tag{78}$$

An explicit discretisation of (78) would be

$$\frac{k^{n+1}-k^n}{\Delta t} = -\varepsilon^n \quad \Rightarrow \quad k^{n+1} = k^n - \Delta t \varepsilon^n, \tag{79}$$

where n denotes a variable on the old time level and $n+1$ denotes a variable on the new time level. With (79) the time-step criterion for positivity of k^{n+1} is $\Delta t < \varepsilon^n/k^n$, which for example in tidal flow might require very small time steps during slack tide, when production is small and dissipation is high. A solution to this problem has been proposed by Patankar (1980):

$$\begin{aligned} \frac{k^{n+1}-k^n}{\Delta t} &= -\varepsilon^n \frac{k^{n+1}}{k^n} &&= -c_\varepsilon \frac{k^{n+1}\left(k^n\right)^{1/2}}{l^n} \\ \Rightarrow \quad k^{n+1} &= \frac{k^n}{1+\Delta t \frac{\varepsilon^n}{k^n}} &&= \frac{k^n}{1+\Delta t c_\varepsilon \frac{(k^n)^{1/2}}{l^n}}, \end{aligned} \tag{80}$$

which gives unconditionally positive results for k^{n+1}. Note that the second equalities in equations (80) are derived by using the cascading relation (26).

This so-called Patankar trick is used in almost all discretisations of turbulence closure budget equations, including those for other parameters such as the dissipation rate ε and the turbulence frequency ω, see Deleersnijder et al. (1997) and Burchard et al. (2005b) for details. The Patankar trick may also be applied to sink terms in biogeochemical models to guarantee positivity, but to obtain conservation of state variables, source terms must be treated accordingly (Burchard et al. (2005a)).

Another issue of the numerical treatment of turbulence closure models is the energy-consistency between mean flow and turbulence. Physically, the kinetic energy extracted from the mean flow due to vertical stress divergence is a source of turbulent kinetic energy transferred as shear production P. The same applies to the exchange of energy between mean potential energy and turbulence kinetic energy which is transferred as buoyancy production B. Burchard (2002b) shows how a consistent energy exchange can be obtained also on the numerical level.

8. Existing coastal ocean community models

Previous sections have discussed the various choices that have to be made when designing a numerical model with a focus on coastal applications. The present section gives an overview of the specifics of some existing hydrostatic and non-hydrostatic structured-grid models to illustrate the diversity of methods in use in coastal models and their evolution over the years. We have strived here to present a fair description of past and ongoing developments, however the reader must be warned that omissions and inaccuracies in the model description are inherent to this type of exercise.

8.1. Hydrostatic structured-grid models

In this subsection examples of structured-grid oceanic models used in practice to tackle research or engineering applications related to coastal environments are given, see Tab. 2 for a list of the most popular models historically used in this context. In this list we can find models originally designed for shallow seas applications (e.g. BOM, COHERENS, GETM, HAMSOM, Mars3D, MOHID), as well as mesoscale oceanic models, like NCOM, POLCOMS, POM or ROMS, adapted to handle nearshore and coastal scales (e.g. Oey, 2005; Uchiyama et al., 2010; Warner et al., 2013). More recently, global climate models (e.g. NEMO, Hycom) are also sometimes used to study shelf processes provided some adjustments in their numerical formulation (e.g.

Leclair and Madec, 2011; Lahaye et al., 2011). Finally, models specifically designed for commercial/engineering applications (e.g. Delft3D, ECOM-si, MIKE 3, TRIM3D) can also be used for real case studies (e.g. Cheng et al., 1993). As summarised in Tab. 3, the choice of discrete algorithms in oceanic models bear the imprints of their original target application and development time. In the 90's the situation was very clear: climate models were generally discretised on a B-grid using a geopotential coordinate under the rigid-lid assumption and leapfrog (LF) time-stepping (Griffies, 2004) while mesoscale models, largely inspired by POM, made use of a C-grid, a terrain-following coordinate and an explicit free surface. At that time coastal and engineering models were predominantly C-grid models based on a semi-implicit free surface and baroclinic foward in time methods influenced by Backhaus (1985) and Casulli and Cheng (1992). Nowadays, the distinction between the different class of models is less clear because newly developed models are generally designed to handle a larger range of spatio-temporal scales permitted by the advances in computational power and the maturity of nesting strategies (see Sec. 6). In early 2000's, the emergence of models like ROMS, GETM or Hycom came with a thorough rethinking of several model components like time-stepping algorithms (Shchepetkin and McWilliams, 2005), vertical coordinate (Bleck, 2002; Burchard and Beckers, 2004), turbulent closure schemes (Umlauf and Burchard, 2003), or open boundary conditions (Marchesiello et al., 2001) which increased further the diversity of models. Tab. 3 reflects the variety of models and modelling techniques that are presently used. However some general trends can be distinguished :

- Models originally developed with coastal applications in mind are characterized by 2-level (one-step) time-stepping with a coupled space and time treatment of advective terms. This choice combines simplicity and ease to implement monotonicity-preserving advection.

- The majority of models based on a coupled space and time approach favour the superbee slope limiter to impose the TVD property. This choice corresponds to the upper boundary of the TVD region and leads to the least numerical diffusion among the existing slope limiters (e.g. Cushman-Roisin and Beckers, 2011). However, this limiter tends to erroneously amplify the gradients. According to Durran (2010), the monotonized centered limiter is expected to provide superior results in terms of accuracy. Other flux-limiter methods currently used for practical applications is the ultimate quickest scheme (or equivalently the

so-called P2-PDM scheme). Note that the WENO or FCT approaches are still relatively marginally used.

- As motivated in Sec. 3.2.1, the overwhelming majority of coastal models rely on TKE-based vertical closure models. A dynamic equation for TKE is solved either in combination with a diagnostic mixing length or with a second dynamic equation for a length related quantity.

- The recently developed models are almost exclusively based on an explicit external mode computation which allows more accuracy in the computation of fast barotropic waves and is expected to provide enhanced scalability and performance on parallel computers.

- The vertical coordinate system is one of the aspect that focuses the most attention nowadays. The general trend is to allow more flexibility in the definition of the coordinate within the Arbitrary Lagrangian-Eulerian framework.

8.2. Non-hydrostatic models

Within the last decade several existing coastal ocean models were extended to abandon the hydrostatic pressure assumption (POM: Kanarska and Maderich (2003), BOM: Heggelund et al. (2004), ROMS: Kanarska et al. (2007), Symphonie: Auclair et al. (2011), GETM: Klingbeil and Burchard (2013)). Scaling analysis shows that the nonhydrostatic pressure contribution, formally defined by

$$\partial_z p_{\mathrm{nh}} = -\left(\partial_t w + \partial_x\left(uw\right) + \partial_y\left(vw\right) + \partial_z\left(ww\right) - f_{\mathrm{h}} u - F_{\mathrm{z}}\right), \qquad (81)$$

cannot be neglected when approaching sufficiently high spatial resolutions that promote processes with significant vertical accelerations (Marshall et al., 1997; Klingbeil and Burchard, 2013). Practical examples on coastal ocean scales are flows above irregular topography, high-frequency internal gravity waves or gravity plumes. In order to simulate these processes in a physically correct way, the vertical momentum balance must not be degenerated to (7) anymore.

For solving the full set of incompressible Navier-Stokes equations algorithms were implemented on top of the hydrostatic model kernels that do not require a complete rewrite of the existing kernels. These were usually based on either the classical projection method (Chorin, 1968; Témam, 1969)

Model	Acronym	website
BOM	Bergen Ocean Model	http://www.mi.uib.no/BOM/
COHERENS	COupled Hydrodynamical Ecological model for REgioNal Shelf seas	http://odnature.naturalsciences.be/coherens/
Delft-3D	Deltares flow	http://oss.deltares.nl/web/delft3d
ECOM-si	Estuarine, Coastal and Ocean Model (semi-implicit)	http://woodshole.er.usgs.gov/operations/modeling/ecomsi.html
GETM	General Estuarine Transport Model	http://www.getm.eu/
HAMSOM	Hamburg Shelf Ocean Model	https://wiki.zmaw.de/ifm/TO/Hamsom/modelhistory
ROMS	Regional Oceanic Modeling System	
↪ Rutgers		https://www.myroms.org/
↪ UCLA		http://research.atmos.ucla.edu/cesr/ROMS_page.html
↪ AGRIF		http://www.romsagrif.org/
Mars3D	Model for Applications at Regional Scale	http://wwz.ifremer.fr/mars3d/
MIKE 3	Danish Hydraulic Institute	https://www.mikepoweredbydhi.com/
MOHID	MOdelo HIDrodinâmico	http://www.mohid.com/
NCOM	Navy Coastal Ocean Model	http://www.dtic.mil/dtic/tr/fulltext/u2/a508063.pdf
POLCOMS	Proudman Oceanographic Laboratory Coastal Ocean Modelling System	http://cobs.noc.ac.uk/modl/polcoms/
POM	Princeton Ocean Model	http://www.ccpo.odu.edu/POMWEB/
Symphonie	Sirocco Ocean Model	http://sirocco.omp.obs-mip.fr/ocean_models/S-model
TRIM-3D	Tidal Residual, Intertidal Mudflat	
Hycom	HYbrid Coordinate Ocean Model	https://hycom.org/
NEMO	Nucleus for the European Modeling of the Ocean	http://www.nemo-ocean.eu/

Table 2: Summary of most significant oceanic models historically used for applications to three-dimensional coastal studies. The ROMS code has three variants referred to as ROMS-Rutgers, ROMS-UCLA, and ROMS-Agrif (see Shchepetkin and McWilliams, 2009, for more details). Hycom and NEMO are examples for large scale models which are also applied to the coastal ocean.

	Introd. ref.	Horz grid	Mode splitting	Internal Time-stepping	External Time-stepping	Mom. advec.	Tracer advec.	Vertical discret.	Turb. closure
BOM	Svendsen et al. (1996)	C	SPE	LF-AM3	LF-TR	TVD Superbee	TVD Superbee	s	MY2.5
COHERENS	Luyten et al. (1999)	C	SPI/SPE	2-level	-	TVD Superbee	TVD Superbee	s	k-ε
Delft-3D	Gerritsen et al. (2004)	C	SPI	ADI	ADI	UP3	UP3 w/ Forester fltr	z & s	k-ε
ECOM-si	Blumberg (1992)	C	SPI	2-level	θ-scheme	C2	C2	s	MY2.5
GETM	Burchard and Bolding (2002)	C	SPE	2-level	FB	TVD	TVD	adapt.	GOTM
HAMSOM	Backhaus (1985)	C	SPI	2-level	θ-scheme	upstream	upstream	σ	k-ε
ROMS (Rutgers)	Haidvogel et al. (2000)	C	SPE	LF-TR	LF-AM3	UP3	MPDATA	s	GLS
ROMS (UCLA)	Shchepetkin and McWilliams (2005)	C	SPE	LF-AM3	Gen FB	UP3	UP3	s	KPP
ROMS (Agrif)	Penven et al. (2006)	C	SPE	LF-AM3	Gen FB	UP3	WENO5Z	s	KPP
Mars3D	Lazure and Salomon (1991)	C	SPI	2-level	ADI	QK3	QK3	s	GLS
MIKE 3	Pietrzak et al. (2002)	C	SPI	2-level	ADI	Ultimate QK3	Ultimate QK3	s	k-ε
MOHID	Martins et al. (2001)	C	SPI	2-level	ADI	upstream	TVD Superbee	s	GOTM
NCOM	Martin (2000)	C	SPI	LF	θ-scheme	SUP3	FCT(UP3)	hyb z/s	MY2.5
POLCOMS	Holt and James (2001)	B	SPE	2-level	FB	PPM w/ limiters	PPM w/ limiters	σ	MY2.5
POM	Blumberg and Mellor (1978)	C	SPE	LF	LF	C2	MPDATA	s	MY2.5
Symphonie	Johns et al. (1992) Marsaleix et al. (1998)	C	SPE	LF	LF	C4	SUP3	s	TKE
TRIM-3D	Cheng et al. (1993)	C	SPI	2-level	θ-scheme	SL	TVD Van-Leer	-	GLS
Hycom	Bleck (2002)	C	SPE	LF	LF	Enstrophy-conserving	FCT(C2)	hyb z/s/ρ	KPP
NEMO	Madec et al. (1991)	C	SPE	LF	LF	SUP3	FCT(C2)	z & s	GLS

Table 3: Properties of some structured-grid coastal ocean models (note that the models might have many more choices than those listed here). Acronyms represent the following: FB=Forward-Backward, Gen = generalised, ADI = Alternating Direction Implicit, SPE=SPlit-Explicit, SPI=SPlit-Implicit, (S)UP3= (Split) third-order UPwind, QK3=third-order quickest, LF=Leapfrog, PPM=Piecewise Parabolic Method, TVD=Total Variation Diminishing, C2=second-order centered, C4=fourth-order centered, TR=trapezoidal, AM3=third-order Adams-Moulton interpolation, MPDATA=Multidimensional Positive Definite Advection Transport Algorithm.

or the pressure correction method (van Kan, 1986). Both methods require the solution of a Poisson equation for either the nonhydrostatic pressure (Kanarska and Maderich, 2003; Heggelund et al., 2004) or its correction (Kanarska et al., 2007; Auclair et al., 2011), which should be derived by inserting the discrete momentum equations into the discrete incompressibility condition to force a divergence-free final velocity field. The specific characteristics of coastal ocean models demand for some adaptations and care, e.g. for the consistent coupling of the free surface to the corrected nonhydrostatic velocities or the numerical treatment of the additional diagonals in the coefficient matrix that is neither symmetric nor positive-definite in general. Strategies to avoid the additional diagonals, only caused by the cross-terms of the vertical coordinate transformation, were presented by Heggelund et al. (2004) and Berntsen and Furnes (2005). Following Berntsen and Furnes (2005), Keilegavlen and Berntsen (2009) demonstrated that simulations with the nonhydrostatic pressure defined directly in σ-coordinates showed no significant deviations to results obtained with the pressure defined in z-coordinates. An alternative algorithm for weakly nonhydrostatic regimes (e.g. internal lee waves) that avoids the Poisson equation was presented by Klingbeil and Burchard (2013). They treat the nonhydrostatic pressure as the hydrostatic one, i.e. diagnose the time-lagged right-hand side of (81), integrate it vertically and insert it into the horizontal momentum equations.

9. Future perspectives

The present review of the numerics of structured-grid coastal ocean models gives an overview of the numerical methods available for todays coastal ocean models. There is no single model which combines all those methods. Thus, by just recombining existing methods, todays model platforms may be further improved.

But what are the future challenges in coastal ocean modelling? There are various trends which will drive further developments. One driver is certainly the ever ongoing growth of computational resources allowing for higher spatial resolution, better process representations, integration into regional Earth system models of growing complexity, application of more elaborate data assimilation schemes, and use of large ensemble simulations with high numbers of individual model simulations for better representation of model statistics.

Non-hydrostatic and Boussinesq assumption

The possible higher spatial resolution will in many coastal ocean scenarios allow for resolving non-hydrostatic processes. In cases where the horizontal resolution becomes substantially smaller than the water depth, hydrostatic models would loose their predictability when for example relatively short internal waves still propagate at their hydrostatic phase speed. Therefore, one challenge to the developers of coastal ocean models is to allow for efficient non-hydrostatic pressure calculation in situations where needed. As discussed in section 8.2, there are only a few of such examples among the structured-grid models. The challenge would therefore be that models automatically analyse whether in some regions the non-hydrostatic pressure contribution is relevant, and if so, the non-hydrostatic dynamics should be applied. On other (e.g., shallower) parts of the model domain it might be ignored. It should however be kept in mind that diagnosed non-hydrostatic pressure might also be a numerical artefact (Vitousek and Fringer, 2011).

Another possibility to consistently add non-hydrostatic effects, while avoiding the costly direct solution of a 3D Poisson equation for the pressure, is to relax both the hydrostatic and Boussinesq assumptions (Auclair et al., 2016). Since relaxing the Boussinesq assumption reintroduces the acoustic waves, an acoustic mode must be considered on top of the internal and external modes. This three-mode approach has the advantage to be completely explicit (i.e. local in space) and thus to have very good properties in terms of scalability. However, it raises new challenges to derive consistent open boundary conditions to couple models based on different modeling assumptions.

Physics-dynamics coupling

The development of numerical models generally dissociates the dynamical kernel, which handles the resolved scales, from the "physics", that account for under-resolved processes. However, both components are strongly coupled since there is constantly a transfer of information back and forth from the dynamics to the physics so that the dynamical core must not violate physical principles. Energy-consistency is certainly one important example of such coupling. As discussed in section 7.1, most advection schemes are not variance conserving, which means for the momentum budget violation of kinetic energy conservation. The rate of numerical dissipation can be diagnosed (e.g. Klingbeil et al., 2014) and it can be shown that more complex advection schemes limit the numerical dissipation (Mohammadi-Aragh et al., 2015). Another source of energy-conservation violation is due to the exchange

between mean kinetic energy and potential energy. This exchange is represented by the pressure-gradient term in the momentum equations and the advection terms of the potential temperature and salinity equations. Since those terms are discretised completely independently, energy-consistency is not given. Another issue is the energy consistency between loss of mean kinetic energy due to vertical stress divergence and production of turbulent kinetic energy (see section 7.6 and Burchard, 2002b). Generally better energy-consistency may also be achieved by more suitable choices for vertical coordinates. As discussed in section 4.2, generalised vertical coordinates give a high degree of freedom to construct an optimised vertical grid layout. This flexibility needs to be better exploited in the future to improve the performance of coastal ocean models.

It should be briefly mentioned that also the physics of coastal ocean models (as well as for large scale ocean models) is generally not energy-consistent, since e.g. the energy lost by horizontal stress divergence is not reintroduced into any other energy compartment of the model. This is an important issue which could however not be further discussed in this numerical review (see e.g., the discussion by Eden et al., 2014).

Multi-resolution strategies

One other aspect of ongoing discussions in the coastal ocean developers community is the optimal layout for horizontal grids. In both, structured-grid and unstructured-grid models, substantial progress has been made in the past. While unstructured-grid models have the better ability to solve multi-scale problems, the numerical cost per nominal grid point is up to one order of magnitude larger than for structured-grid models. It does therefore depend on the kind of problem to decide which type of model is better suited. For complex domains such as lagoon systems or estuarine networks, unstructured-grid models are the better choice while for classical coastal or shelf sea scenarios structured-grid models are more efficient. Structured-grid models have gained flexibility due to more flexible grids (such as curvilinear grids, see section 4.3 or two-way nesting, see section 6.3). Flexible multi-level nesting strategies reducing the number of masked grid points (land points) in the model domain might in the future help to better tackle multi-scale problems using structured grids. And there are efforts to increase the efficiency of unstructured-grid models. Both types of models - structured-grid models with flexible nesting and unstructured-grid models - might in the future support horizontal mesh-adaptivity to resolve moving or intermittent dynamic

features (as for example done in larger scale models to follow hurricanes). It depends on the success of such developments which type of model is gaining more ground in the future of coastal ocean modelling.

Coupling with other earth system compartments

Finally, stand-alone hydrodynamic coastal ocean models will always have a limited area of application. All such models are developed to also function as hydrodynamic core of coupled model systems. Most common is the coupling to biogeochemical models, where the hydrodynamic model provides advective and turbulent transports as well as the light regime and many other potential functional links between hydrodynamics and biogeochemistry. Without accurate physical and numerical representation of the hydrodynamics (e.g., in terms of the mixed-layer depth), biogeochemical models will have no predictability. Links between hydrodynamics and biogeochemistry may be complex. One tool to flexibly provide such a link is FABM (Framework of Biogeochemical Models, Bruggeman and Bolding (2014)).

Coupling to several other Earth system compartments is common, such as to statistical surface wave models, sea ice models, atmospheric models, benthic models, and hydrological models. Several comprehensive system models exist, such as COAWST (Warner et al., 2008), interactively coupling hydrodynamics (from ROMS) with surfaces waves, atmosphere, biogeochemistry, sediment transport and benthos to study complex coastal ocean scenarios in hindcast and sensitivity mode, or NEMO-Nordic (Gröger et al., 2015), coupling hydrodynamics (from NEMO) with atmosphere, sea ice, biogeochemistry, and hydrology for the purpose of regional climate downscaling and sensitivity scenarios. It is obvious that all coupling interfaces to other compartments require hydrodynamic process adaptations on the side of the hydrodynamic module, and that those need to be numerically discretised with care. It is expected that in the future, such complex coupled model systems will be much more widely used, such that quite some pressure to further develop hydrodynamic model will be generated.

Despite the substantial growth of computational resources, regional climate system models with run-times of typically more than a century and which maybe executed as large ensemble runs with many individual simulations will require high accuracy at relatively low resolution. Therefore the development of efficient and accurate numerical schemes at low resolution will be an important task also for the future.

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